



May 7, 2018

Buzz Oates Development and Construction Services
Attention: Troy Estacio
555 Capitol Mall, Ninth Floor
Sacramento, CA 95814

DRAFT WASTE MATERIALS REUSE WORKPLAN
WASTE UNIT 1 MATERIALS
MURPHY PARKWAY PROPERTY
APN 198-012-009
17100 Murphy Parkway
Lathrop, California
Brusca Project No. 137-002

INTRODUCTION

In accordance with your request, we have prepared this *Draft Workplan* for the onsite reuse of solid waste materials and associated soils that were excavated, stockpiled, tested, and evaluated at the subject property as a part of clean closure activities. The property consists of an approximate 48-acre, vacant parcel identified as APN 198-012-009 in Lathrop, San Joaquin County, California. Recent waste management work at the site included excavation/removal, stockpiling, sampling, and testing of a layer of solid waste materials (primarily broken glass) and associated soils that existed beneath a clean soil “cap” on the far northwesterly portion of the subject parcel. These wastes on the property have been referred to as Waste Unit 1 (WU1). The removed/stockpiled waste materials were sampled and tested to evaluate whether the materials can be reused onsite in engineered fills during upcoming site development.

The waste management work recently performed at the site is a component of the clean closure activities for the subject parcel in general accord with CCR Title 27, Section 21810 and with the guidance outlined in the Local Enforcement Agency (LEA) Advisory #16 published by CalRecycle. Advanced GeoEnvironmental (AGE) prepared a *Clean Closure Plan* for the property dated March 5, 2018.¹ An element of the *Clean Closure Plan* is an *Excavation and Material Management Plan (EMMP)* that was prepared by our firm and presented measures and procedures for the removal, management, sampling, testing, and evaluation of residual solid waste materials within WU1.² Our firm prepared *Addendum* to the *EMMP* dated March 22,

¹ AGE; “Clean Closure Plan, 17100 Murphy Parkway, Lathrop, San Joaquin County, California, PNA Parcel B (Portion of Solid Waste Information System [SWIS] number 39-CR-0022), APN 198-120-09”; March 5, 2018.

² Brusca Associates, Inc.; “Excavation and Material Management Plan, Murphy Parkway Property, APN 198-012-009, 17100 Murphy Parkway, Lathrop, San Joaquin County, California”; March 9, 2018.



2018 to slightly revise the verification soil sampling program (testing of soils beneath the area of removed wastes) as suggested by the Central Valley Regional Water Quality Control Board (CVRWQCB).³

This *Draft Workplan* includes: a general description of the site; background information; a description of the planned site development; information, laboratory data and discussion pertaining to stockpiled materials generated during the recent removal of solid waste materials from WU1; justification for onsite reuse of materials excavated from WU1; procedures for the reuse of these materials in onsite engineered fills; and verification/reporting procedures to ensure that the materials are reused in the manner described herein.

A *Vicinity Map* showing the location of the subject parcel is presented as Plate 1. A *Site Map* is presented as Plate 2 showing the layout of the entire parcel, the former area of capped wastes (WU1) on the far northwesterly portion of the site, the locations of stockpiled materials that were recently excavated and removed from the capped waste area, and the general layout of the planned large commercial/light-industrial building on the property. A detail map showing WU1 area and the locations of verification soil samples collected beneath the area of removed wastes is presented as Plate 3. A civil engineering drawing depicting additional details pertaining to the planned construction is presented as Plate 4. Analytical laboratory data for samples of the stockpiled materials from WU1, as well as for the verification soil samples collected beneath WU1, are presented on Tables I through VII; the laboratory reports and chain-of-custody documentation are included in Appendix A. The *Foundation Investigation* (geotechnical engineering report) for the site development project is presented in Appendix B.

SITE DESCRIPTION AND BACKGROUND

Site Location and Description

The approximate 48-acre subject property is located about 500 feet easterly of Murphy Parkway and about 1,000 feet northerly of D'Arcy Parkway in a commercial and industrial area of Lathrop, San Joaquin County, California. The property consists of an approximate 48-acre parcel identified as APN 198-012-009. The subject parcel is currently vacant/undeveloped land. The site is generally flat; however, two depressed areas are situated on the northwesterly portion of the site (bordering the WU1 area). Until recently, portions of the site were used for livestock grazing; wire fencing is in place on portions of the site for livestock enclosure purposes. A *Vicinity Map* showing the location of the property is presented as Plate 1, and a *Site Map* is presented as Plate 2.

Historical Site Activities

The 48-acre subject parcel historically was vacant farmland. In the early 1960s, a large glass manufacturing facility (currently referred to as former Pilkington North America [PNA] facility) was constructed directly northerly of the subject property. The subject 48-acre parcel (identified by APN 198-012-009) was associated with the former PNA facility, and, along with other offsite areas, was used for solid waste handling and waste water disposal related to the PNA facility.

³ Brusca Associates, Inc.; "Addendum, Excavation and Material Management Plan, Murphy Parkway Property, APN 198-012-009, 17100 Murphy Parkway, Lathrop, San Joaquin County, California"; March 22, 2018.



Solid wastes comprised largely of waste glass materials were handled within the northwesterly portion of the subject 48-acre site; the vast majority of waste glass materials historically handled and temporarily stored in this area were hauled from the property in the past to offsite disposal locations. Glass manufacturing activities at the former PNA facility were discontinued several years ago. In recent years, the subject 48-acre parcel generally has been unused, except for some livestock grazing.

Treated industrial and domestic waste water, as well as storm water, generated at the former PNA facility northerly of the subject property was disposed at that facility, and also on the subject 48-acre parcel via spray irrigation. The waste water discharge was performed under permit with the CVRWQCB. Because the subject 48-acre property is no longer intended for waste water discharge, the CVRWQCB is currently in the process of removing this area from the waste water discharge permit.

Past Onsite Waste Investigations and Remedial Work

It is indicated that the handling of glass wastes on the subject property ceased in the 1980s. Around 2012, the local oversight agency (the San Joaquin County Environmental Health Department [SJCEHD]) required the placement of a clean soil cap over the northwesterly portion of the property (WU1 area) where a thin layer of residual glass wastes was known to exist at the site surface. A *Soil Cover Workplan* was prepared in early 2013, and the capping work was implemented in late 2013.⁴ The area of the cap (WU1 area) is shown on Plate 2.

Starting in 2014, a substantial amount of investigation, waste characterization, and waste removal work was implemented on the subject parcel by AGE with direction and oversight by the SJCEHD. This work included:

- a significant number of explorations (test pits and borings) within the capped area and other site areas, including areas to the south and east of the capped area.
- a geophysical survey (involving ground penetrating radar and a magnetometer survey) on the westerly portion of the property to identify any unknown areas of disturbed materials/buried wastes. The geophysical and trenching work reportedly identified two areas (approximate 2,000 square feet each in plan area) of deeper disturbed materials/wastes along the far westerly margin of the property and outside of the cap area. These areas were identified as the South Waste Cell (SWC) and the North Waste Cell (NWC). The SWC and NWC waste materials subsequently were completely excavated and removed from the parcel for offsite landfill disposal.
- collection of samples of soil and waste materials from WU1, SWC, and NWC areas, as well as from other site areas for analytical laboratory testing.

AGE's investigative, remedial, and characterization work associated with solid wastes on the subject parcel was summarized in a *Site Assessment Summary Report* dated December 21, 2017.⁵

⁴ Antea Group; "Soil Cover Workplan, Former Cullet Staging Area, Pilkington North America, Inc., 500 East Louise Avenue, Lathrop, California"; May 22, 2013.

⁵ AGE; "Revised Site Assessment Summary Report, 17100 Murphy Parkway, Lathrop, San Joaquin County, California, Solid Waste Information System (SWIS) Number 39-CR-022 (APN 198-120-09); January 10, 2018.



PLANNED SITE DEVELOPMENT/CONSTRUCTION

Redevelopment of the subject property is planned to be initiated in the very near future. The planned development includes construction of an approximate 870,000 square-foot commercial/light-industrial structure intended for use for warehouse/distribution purposes. The layout of the planned development is shown on Plates 2 and 3. The planned structure will be of tilt-up concrete construction with a concrete slab-on-grade floor. Onsite areas exterior to the building will support asphaltic-concrete and concrete pavements, storm water basins, and very limited landscaping.

WU1 WASTE EXCAVATION, STOCKPILING, SAMPLING/TESTING, AND EVALUATION

Excavation, removal, and stockpiling of waste materials and associated soil within WU1 (capped area within the northwesterly portion of the subject parcel) was performed in April 2018 by Ramcon Engineering and Environmental Contracting. The work was overseen by a Certified Engineering Geologist from our firm; additionally, representatives of SJCEHD, the CVRWQCB, and CalRecycle observed the work. The work was performed in accordance with the agency-approved *EMMP* and associated *Addendum* referenced earlier in this *Workplan*.

Prior to removal of the waste materials from WU1, the clean soil cap materials were removed from this area and stockpiled along the northerly margin of the site. Subsequently, the underlying waste materials and associated soils were excavated with earthmoving equipment, removed from the WU1 area, and stockpiled directly to the south of the WU1 area on the property (see Plate 2). Five stockpiles of WU1 materials were created (identified as SP1 through SP5). Stockpiles SP1, SP2, SP3, and SP5 range in volume from about 1,000 to about 5,000 cubic yards. Stockpile SP4 is comprised predominantly of sizable cemented masses (boulder-size and smaller) of glass fragments in a fine-grained matrix; the volume of SP4 is quite small (less than 10 cubic yards). The total volume of stockpiled materials removed from WU1 is on the order of 13,000 to 14,000 cubic yards. The stockpiles were covered with plastic sheeting, and the plastic sheeting was secured with sandbags to prevent wind disturbance.

Per the provisions of the *EMMP*, our firm collected verification soil samples below the removed wastes in the WU1 area, and we collected samples of the stockpiled materials removed from WU1 for analytical laboratory testing. The sampling was conducted by a Certified Engineering Geologist from our firm and was overseen by regulatory agency representatives. A total of 49 spatially distributed, random/representative samples of the WU1 materials was collected. The samples were placed in laboratory-provided glass jars, immediately placed on ice, and transported to State-certified analytical laboratories under chain-of-custody for analysis for the following:

- Total Petroleum Hydrocarbons as gasoline, diesel, and motor oil by EPA Method 8015B
- SVOCs by EPA Method 8270C
- Polychlorinated Biphenyls (PCBs) by EPA Method 8082
- Title 22 Metals Total Threshold Limit (TTLC) by EPA Method 6010/200.7
- Dioxins and Furans by EPA Method E1613



The results of the above-listed analyses performed on stockpiled materials from WU1 are presented on summary data tables attached to this *Workplan*; summary data tables from the verification soil sampling beneath the area of removed wastes also are included. For reference, hazardous waste criteria and screening levels are included on the summary data tables. The laboratory reports and chain-of-custody documentation are presented in Appendix A.

As outlined in the *EMMP*, samples stockpiled materials from WU1 containing the highest concentrations of total metals were analyzed for soluble Title 22 metals by the Waste Extraction Method (WET) by EPA Method 1311 using deionized water (DI) as the extractant; this element of the testing program was required by the CVRWQCB for the purpose of evaluating potential future leaching impact to groundwater posed by the waste materials. A total of 11 samples was selected for the DI WET metals testing, including samples from each of the five stockpiles.

WU1 MATERIALS REUSE EVALUATION

Per the provisions of the approved *EMMP*, the laboratory data from the samples of stockpiled materials removed from WU1 have been evaluated to determine whether these materials are acceptable for onsite reuse in engineered fills during upcoming site development. As shown on Tables IV, VI, and VII, none of the 49 samples of the stockpiled materials from WU1 contained total metals, petroleum hydrocarbons, PCBs, SVOCs or dioxins/furans at concentrations above the commercial/industrial screening levels established in the *EMMP* (or in the case of arsenic, above the site-specific background concentration). Additionally, none of the results exceed hazardous waste criteria established in the California Code of Regulations (CCR) Title 22, Division 4.5. As such, these results indicate that the stockpiled materials from WU1 are acceptable for onsite reuse.

Materials reuse evaluation has included consideration of the potential for future impact to groundwater as a result of possible leaching of metals from the materials. As shown on Table V, the DI WET concentrations from the tested samples for the majority of the Title 22 metals do not exceed California Maximum Contaminant Level (MCL) drinking water concentrations. However, the concentrations of DI WET arsenic, chromium and lead concentrations detected in most of the samples exceed the MCL values. We note that several groundwater samples have been collected and tested beneath the WU1 area as a part of past site assessment work, and none of the groundwater samples from this area have contained lead or chromium at concentrations above their respective MCL values, despite the fact that waste materials were handled and stored in this area for several decades starting in the 1960s (and during operation of the PNA facility, much greater volumes of wastes were stored in this area than were removed during the recent waste management work). As such, it does not appear likely that the chromium and lead in the removed/stockpiled waste materials represents a significant threat to groundwater quality. It is noted that elevated concentrations of arsenic (above the MCL value) have been detected in groundwater at various locations beneath the subject property, including beneath the WU1 area. However, in general, the distribution of elevated concentrations of arsenic in groundwater beneath the site is somewhat erratic, and some of the past groundwater samples collected beneath WU1 in the past have not contained elevated concentrations of arsenic. It is known that naturally-occurring elevated concentrations of arsenic exist in groundwater in other areas of San Joaquin County, and the arsenic detected in groundwater beneath the subject property may also be naturally occurring.



To further evaluate the leaching potential of arsenic in the stockpiled materials removed from WU1, we have compared the estimated mass of the DI soluble arsenic contained in the stockpiled materials to a potential volume of receptor groundwater beneath the site. Based on an average DI WET arsenic concentration of 0.02 milligrams per Liter (mg/L), and recognizing that the DI WET test utilizes 50 grams of materials and 500 milliliters of water as an extractant, we calculate the total mass of DI soluble arsenic in the stockpiled materials to be approximately 4,120 grams. Assuming, that the entirety of this mass is leached to groundwater, a saturated zone void ratio of 0.5, and a twenty-foot depth of dispersion and diffusion within the saturated zone, the soluble arsenic mass in the stockpiled materials would represent 0.007 mg/L beneath the 48-acre property; this value is below the drinking water standard for arsenic (0.01 mg/L). This mass calculation likely substantially overestimates potential arsenic impact to groundwater because the DI WET data utilized is from the samples with the highest total metals results (the actual average DI WET arsenic concentration in the stockpiled materials likely is less than 0.02 mg/L), and the mass calculation assumes that all soluble arsenic in the stockpiled materials leaches to groundwater.

Despite the foregoing that suggests that the soluble arsenic, chromium and lead detected in the stockpiled materials removed from WU1 is not likely to represent a significant leaching threat to shallow groundwater quality, as a highly protective and conservative measure we propose specific procedures to be implemented during the onsite reuse of these materials to mitigate leaching potential. Specifically, we propose that the stockpiled materials removed from WU1 be placed beneath the concrete floor slab of the planned building on the property (and to a lesser degree, beneath planned concrete pavements) so that these materials are not subject to rainwater infiltration and related leaching potential. Construction plans indicate that the concrete slab-on-grade floor for the planned building on the property will be seven inches thick and reinforced with #4 reinforcing bars at 24 inches on center, each way; floor slab concrete will be designed to achieve a minimum 28-day unconfined compressive strength of 3,000 pounds per square inch (psi). The building will be completely surrounded by concrete and asphaltic-concrete pavements. Site pavements will be sloped to drain and onsite storm water will be directed toward drain inlets that discharge to the storm water system. As such, rainwater infiltration and associated leaching potential to materials directly below the future building's floor slab is considered to be negligible.

Once the building pad is constructed utilizing the materials removed from WU1 within the pad subgrade, a comparatively minor amount of these materials will subsequently be excavated during foundation and underground utility trench excavation within the building pad. We suggest that these materials be placed below concrete pavements that will adjoin the building (i.e. the planned concrete-paved dock areas northerly and southerly of the planned building). The concrete pavements are designed to be six inches thick and reinforced with #4 reinforcing bars at 24 inches on center, each way; pavement concrete will be designed to achieve a minimum 28-day unconfined compressive strength of 4,000 psi. The concrete pavements will substantially prevent rainwater infiltration and associated leaching potential to underlying materials.

We propose that the materials removed from WU1 be placed within the upper one-foot of the subgrades beneath the floor slab and concrete pavement sections mentioned above since these subgrades will be cement treated as a part of construction activities to enhance the engineering/strength properties of the subgrades. The cement treatment of these materials would further reduce the leaching potential of metals in the materials. It is well recognized that cement treatment of soils can mitigate leaching potential of heavy metal ions by providing both physical



resistance to leaching and by creating an elevated pH environment. The cement treatment process results in an impermeable monolith that substantially reduces surface area available for leaching. Moreover, the increased pH of the treated material (caused by alkali nature of Portland cement) limits the solubility of metals. As such, incorporating the materials from WU1 into floor slab and pavement subgrades that will be treated with cement affords an even greater reduction of the potential for leaching of soluble metals in these materials. Following the placement of the WU1 materials beneath the building floor slab (and to a lesser degree beneath concrete pavements) and the treatment of these materials with cement, the potential for leaching of metals from these materials to groundwater is considered to be negligible.

Specific measures for the placement and cement treatment of the materials are presented below.

PROCEDURES OF REUSE OF WU1 MATERIALS

Notification

At least one week prior to the onset of construction activities at the site that involves the relocation of the stockpiled materials from WU1 and the placement of these materials within building pad and pavement subgrades, the SJCEHD, the CVRWQCB, and CalRecycle will be notified of the work schedule.

Placement and Cement Treatment of WU1 Materials

All of the materials contained in the stockpiles of material removed from WU1 will be placed within the upper one foot of the building pad subgrade for the proposed approximate 870,000 square-foot concrete tilt up building planned to be constructed on the site. Materials placement and compaction should conform to the requirements of project plans and specifications, including the project *Foundation Investigation* prepared by Raney Geotechnical, Inc.; a copy of the *Foundation Investigation* is included in Appendix B of this *Workplan*.⁶

As described in the appended *Foundation Investigation*, the upper one foot of the building pad subgrade will be treated with at least five percent Portland cement, as determined by dry unit weight. The treated mixture will be moisture conditioned and compacted to at least 92 percent of the maximum dry density as determined by the ASTM D1557-02 test procedure.

Any materials generated during trimming of the treated building pad to final grade, and all materials generated during foundation and utility trench excavation within the upper one foot of the building pad should be placed beneath areas adjoining the building that are designated to be concrete paved. These materials will be placed within the upper one foot of the concrete pavement subgrades and will subsequently be cement treated. The project *Foundation Investigation* indicates that the pavement subgrades will be cement treated in a similar manner as described above for the building pad subgrade.

⁶ Raney Geotechnical, Inc.; "Foundation Investigation, Murphy Parkway Warehouse, Murphy Parkway near D'Arcy Parkway, Lathrop, California"; December 19, 2016.



This *Workplan* does not address physical hazards (such as those posed by the waste glass within the materials removed from WU1 on the property); protection against physical and construction hazards are the responsibility of future contractors/workers and other site personnel and are not included in the scope of this plan. It is recommended that all agencies, companies, and workers whose activities may result in exposure to the materials removed from WU1 (including waste glass materials and dust) develop health and safety plans specific to the site conditions and their activities, and that such measures consider/address potential inhalation hazards posed by glass dust, as well as cut/abrasion hazards posed by glass fragments.

Verification Observation and Reporting

A qualified environmental professional will make observations of site activities that involve the onsite transport, placement, and cement treatment of materials removed from WU1 to confirm that the procedures described herein are implemented. Additionally, earthwork quality control observations and testing will be performed by the project geotechnical engineering firm. The locations of placement of the materials removed from WU1 will be recorded on a scaled site plan. Upon completion, a report will be prepared summarizing the materials reuse activities

CLOSING

This *Draft Workplan* will be submitted to the SJCEHD, the CVRWQCB, and CalRecycle for review and approval prior to finalizing the *Workplan* and implementing the onsite reuse of the materials removed from WU1. If you have any questions or require additional information, please contact the undersigned at (916) 677-1470.

Sincerely,

BRUSCA ASSOCIATES, INC.

Joe Brusca
Principal Engineering Geologist
Certified Engineering Geologist No. 1948



JB:jb

Attachments: Plate 1, *Vicinity Map*
Plate 2, *Site Map*
Plate 3, *Verification Soil Sample Detail Map*
Plate 4, *Planned Development Drawing*

Table I, Summary of Total Metals Data – Verification Soil Samples
Table II, Summary of Petroleum Hydrocarbons Data – Verification Soil Samples



Table III, Summary of SVOCs, PCBs, and Dioxins/Furans Data – Verification
Soil Samples

Table IV, Summary of Total Metals Data – Stockpiled Waste Material Samples

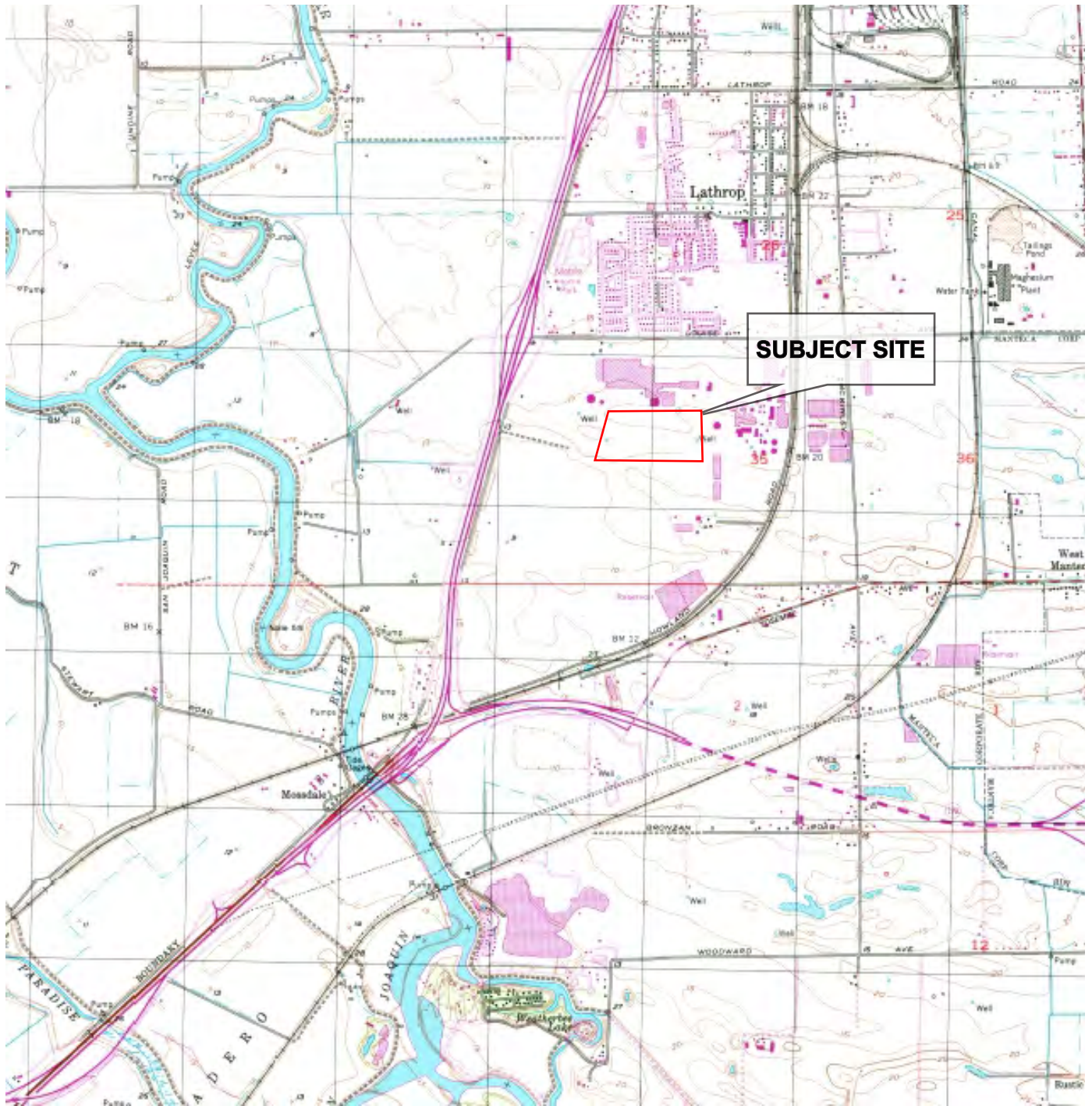
Table V, Summary of DI WET Metals Data – Stockpiled Waste Material Samples

Table VI, Summary of Petroleum Hydrocarbons Data – Stockpiled Waste
Material Samples

Table VII, Summary of SVOCs, PCBs, and Dioxins/Furans Data – Stockpiled
Waste Material

Appendix A: Laboratory Reports and Chain-of-Custody Documentation

Appendix B: *Foundation Investigation* report by Raney Geotechnical, Inc.



SOURCE: U.S.G.S. 7.5-minute Lathrop Quadrangle, California, 1996
Scale 1:24,000



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Environmental Engineering Geology
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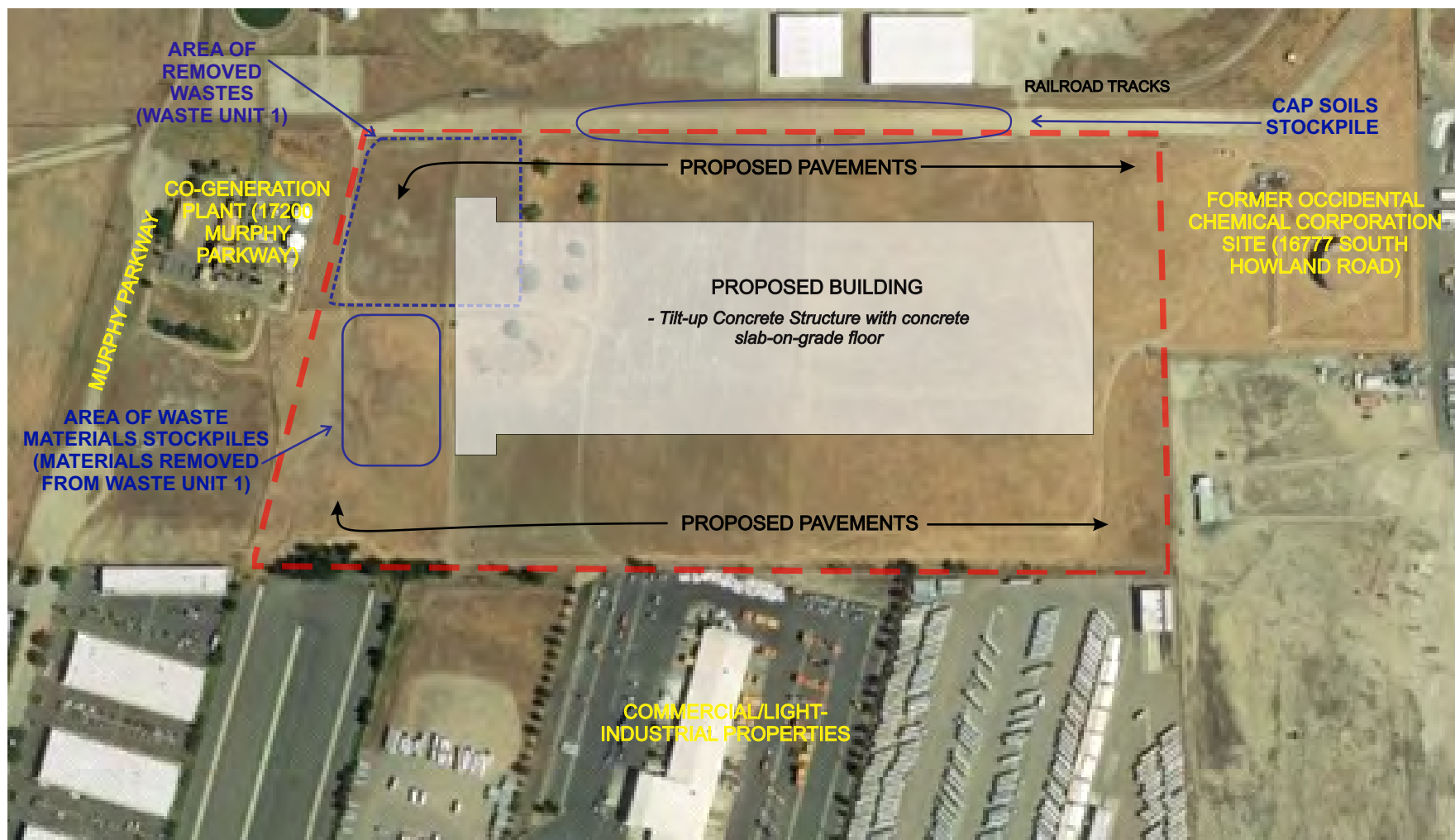
**MURPHY PARKWAY PROPERTY
WASTE MATERIALS REUSE
WORKPLAN**

APN 198-012-009
17100 Murphy Parkway
Lathrop, California

Brusca Project No. 137-002

VICINITY MAP

PLATE 1



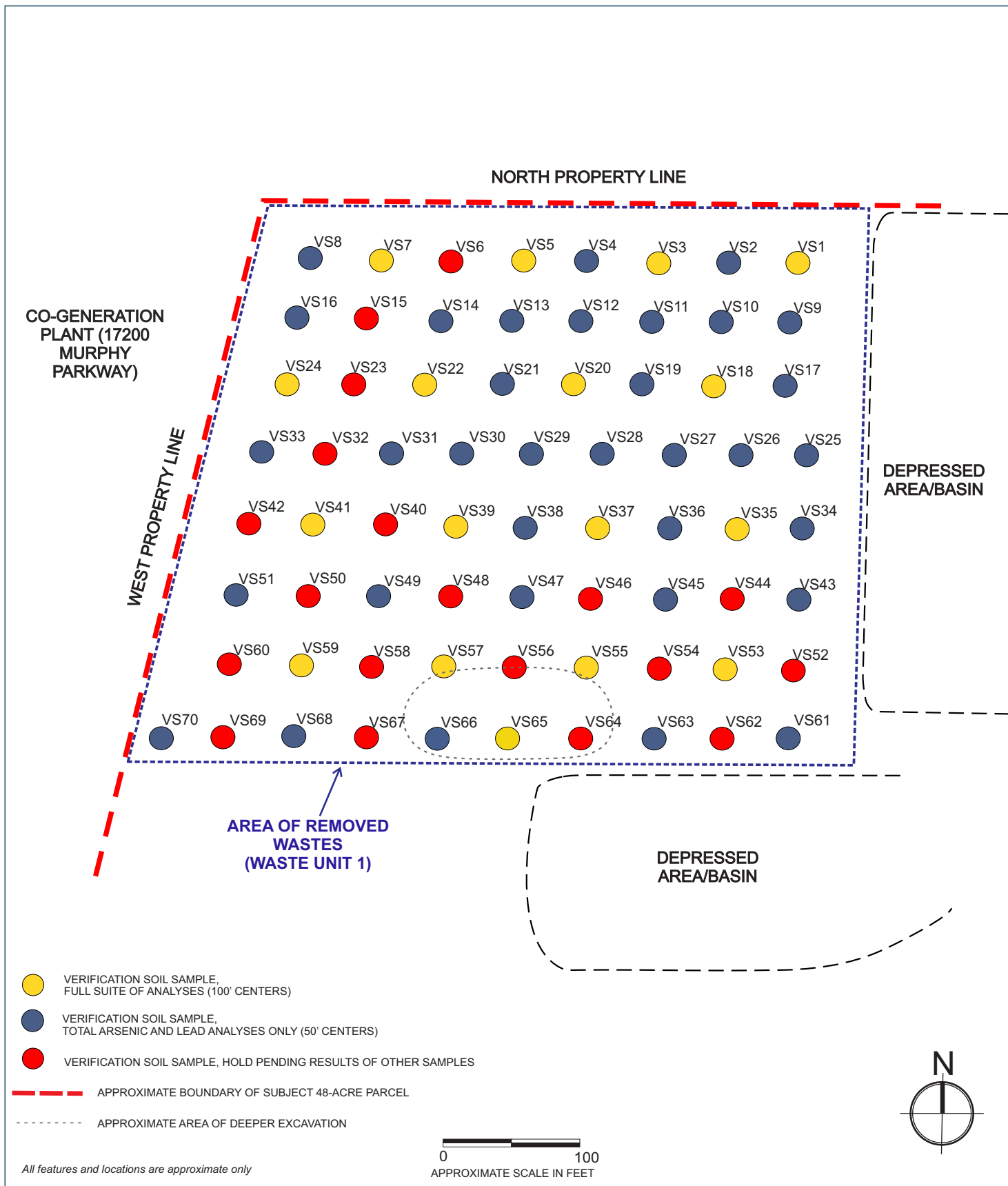
0 400
APPROXIMATE SCALE IN FEET

— — — — — APPROXIMATE BOUNDARY OF SUBJECT 48-ACRE PARCEL

All features and locations are approximate only

 Brusca Associates, Inc. Environmental Engineering Geology 1860 Sierra Gardens Drive, # 332 Roseville, CA 95661 ph (916) 677-1470 fax (916) 677-1471 BruscaAssociates.com	MURPHY PARKWAY PROPERTY WASTE MATERIALS REUSE WORK PLAN		SITE MAP
	APN 198-012-009 17100 Murphy Parkway Lathrop, California		
	Brusca Project No. 137-002		

PLATE 2



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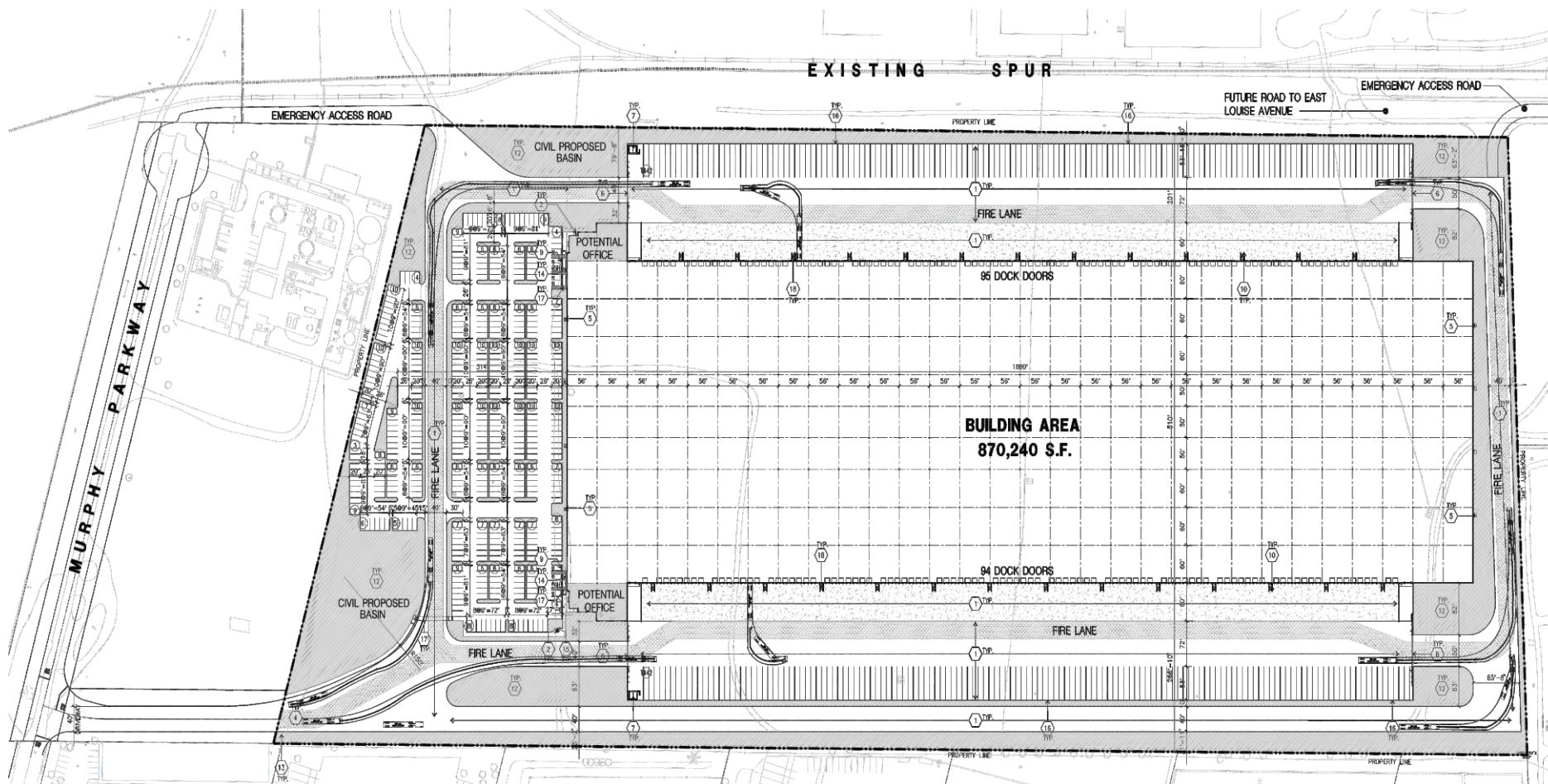
**MURPHY PARKWAY PROPERTY
WASTE MATERIALS REUSE WORK PLAN**

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**VERIFICATION SOIL
SAMPLE DETAIL MAP**

PLATE 3



0 300
APPROXIMATE SCALE IN FEET



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**MURPHY PARKWAY PROPERTY
WASTE MATERIALS REUSE
WORK PLAN**

**APN 198-012-009
17100 Murphy Parkway
Lathrop, California**

Brusca Project No. 137-002

**PLANNED
DEVELOPMENT
DRAWING**

PLATE 4

TABLE I
SUMMARY OF TOTAL METALS DATA
VERIFICATION SOIL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6010B/200.7																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
VS1	4/2/18	ND	ND	ND	49	ND	ND	5.2	5.0	3.3	ND	ND	5.8	ND	ND	19	21	ND
VS2	4/2/18			ND							ND							
VS3	4/2/18	ND	ND	ND	55	ND	ND	4.2	4.2	1.4	ND	ND	6.2	ND	ND	16	17	ND
VS4	4/2/18			ND							ND							
VS5	4/2/18	ND	ND	ND	47	ND	ND	4.8	4.0	ND	ND	ND	3.9	ND	ND	18	18	ND
VS7	4/2/18	ND	ND	ND	45	ND	ND	5.2	3.7	ND	ND	ND	4.1	ND	ND	17	19	ND
VS8	4/2/18			ND							ND							
VS9	4/11/18			ND							ND							
VS10	4/11/18			ND							ND							
VS11	4/11/18			ND							ND							
VS12	4/11/18			ND							ND							
VS13	4/11/18			ND							ND							
VS14	4/11/18			ND							ND							
VS16	4/11/18			ND							ND							
VS17	4/11/18			ND							ND							
VS18	4/11/18	ND	ND	ND	44	ND	ND	3.8	3.4	ND	ND	ND	3.8	ND	ND	13	18	ND
VS19	4/11/18			ND							ND							
VS20	4/11/18	ND	ND	ND	55	ND	ND	4.7	3.6	1.4	ND	ND	4.5	ND	ND	15	22	ND
VS21	4/11/18			ND							ND							
VS22	4/11/18	ND	ND	ND	72	ND	ND	4.3	4.0	1.6	ND	ND	4.4	ND	ND	16	20	ND
VS24	4/11/18	ND	ND	ND	52	ND	ND	4.2	3.8	3.1	ND	ND	4.9	ND	ND	16	19	ND
VS25	4/11/18			ND							ND							
VS26	4/11/18			ND							ND							

TABLE I
SUMMARY OF TOTAL METALS DATA
VERIFICATION SOIL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6010B/200.7																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
VS27	4/11/18			ND							ND							
VS28	4/11/18			ND							ND							
VS29	4/11/18			ND							ND							
VS30	4/11/18			ND							ND							
VS31	4/11/18			ND							ND							
VS33	4/11/18			ND							ND							
VS34	4/11/18			ND							ND							
VS35	4/11/18	ND	ND	ND	66	ND	ND	5.1	6.3	3.6	ND	ND	6.8	ND	ND	22	28	ND
VS36	4/11/18			ND							ND							
VS37	4/11/18	ND	ND	ND	77	ND	ND	5.9	6.4	4.1	ND	ND	7.5	ND	ND	18	26	ND
VS38	4/11/18			ND							ND							
VS39	4/11/18	ND	ND	ND	66	ND	ND	4.7	4.8	4.5	ND	ND	7.6	ND	ND	15	20	ND
VS41	4/11/18	ND	ND	ND	53	ND	ND	4.3	4.1	3.2	ND	ND	4.9	ND	ND	16	18	ND
VS43	4/11/18			ND							ND							
VS45	4/11/18			ND							ND							
VS47	4/11/18			ND							ND							
VS49	4/11/18			ND							ND							
VS51	4/11/18			ND							ND							
VS53	4/13/18	ND	ND	ND	80	ND	ND	7.1	6.6	6.8	ND	ND	8.8	ND	ND	23	28	ND
VS55	4/13/18	ND	ND	ND	65	ND	ND	7.7	5.8	4.1	ND	ND	8.2	ND	ND	24	27	ND
VS57	4/13/18	ND	ND	ND	49	ND	ND	6.6	4.7	3.0	ND	ND	5.8	ND	ND	21	32	ND
VS59	4/13/18	ND	ND	ND	62	ND	ND	6.0	5.0	3.5	ND	ND	7.2	ND	ND	20	22	ND
VS61	4/13/18			ND							ND							

TABLE I
SUMMARY OF TOTAL METALS DATA
VERIFICATION SOIL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY
17100 Murphy Parkway, Lathrop, San Joaquin County, California
Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6010B/200.7																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
VS63	4/13/18			ND							ND							
VS65	4/13/18	ND	ND	ND	26	ND	ND	3.7	2.7	1.9	ND	ND	3.4	ND	ND	11	11	ND
VS66	4/13/18			ND							ND							
VS68	4/13/18			ND							12							
VS70	4/13/18			ND							ND							
SCREENING LEVELS																		
Hazardous Waste Criteria²		500	500	500	10,000	75	100	2,500	8,000	2,500	1,000	3,500	2,000	100	700	2,400	5,000	20
DTSC SLs - residential³		N/A	390	0.11 ⁶	N/A	15	5.2	36,000	N/A	N/A	80	N/A	490	N/A	N/A	390	N/A	1.0
DTSC SLs - commercial³		N/A	1,500	0.36 ⁶	N/A	210	7.3	170,000	N/A	N/A	320	N/A	3,100	N/A	N/A	1,000	N/A	4.5
USEPA RSLs - residential⁴		31	390	0.68 ⁶	15,000	160	71	120,000	23	3,100	400	390	1,500	390	0.78	390	23,000	11
USEPA RSLs - commercial⁴		470	5,800	3.0 ⁶	220,000	2,300	980	1,800,000	350	47,000	800	5,800	22,000	5,800	12	5,800	350,000	46

Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg)

2. California Code of Regulations Title 22, Division 4.5

3. DTSC-Modified Screening Levels (DTSC-SLs); January 2018

4. USEPA Regional Screening Levels, TR=1E-06, HQ=1"; November 2017

5. N/A = screening value not established

6. Arsenic data for the project are compared to background concentration of 8.6 mg/kg rather than cited screening levels.

TABLE II
SUMMARY OF PETROLEUM HYDROCARBONS DATA
VERIFICATION SOIL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TOTAL PETROLEUM HYDROCARBONS EPA Method 8015C		
		C6-C12 (Gasoline Range)	C13-C28 (Diesel Range)	C29-C40 (Motor Oil Range)
VS1	4/2/18	ND	ND	ND
VS3	4/2/18	ND	ND	ND
VS5	4/2/18	ND	ND	ND
VS7	4/2/18	ND	950	2,100
VS18	4/11/18	ND	ND	ND
VS20	4/11/18	ND	ND	ND
VS22	4/11/18	ND	ND	ND
VS24	4/11/18	ND	220	310
VS35	4/11/18	ND	ND	ND
VS37	4/11/18	ND	ND	ND
VS39	4/11/18	ND	ND	ND
VS41	4/11/18	ND	ND	ND
VS53	4/13/18	ND	ND	ND
VS55	4/13/18	ND	21	26
VS57	4/13/18	ND	ND	ND
VS59	4/13/18	ND	ND	ND
VS65	4/13/18	ND	ND	ND
SCREENING LEVELS				
ESLs - residential²		740	230	11,000
ESLs - commercial²		3,900	1,100	140,000

Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg)

2. San Francisco Bay Regional Water Quality Control Board; "Environmental Screening Levels"; February 2016; Direct Exposure Human Health Levels; residential and commercial/industrial.

TABLE III
SUMMARY OF SVOCs, PCBs, AND DIOXINS/FURANS DATA
VERIFICATION SOIL SAMPLES
48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	PCBs EPA Method 8082			SEMI-VOLATILE ORGANIC COMPOUNDS EPA Method 8270C	DIOXINS/FURANS EPA Method 1613	
		PCB-1254	PCB-1260	Other PCBs		2,3,7,8-TCDD ⁶	TEQ ⁷
VS1	4/2/18	ND	ND	ND	ND	ND	0.00672
VS3	4/2/18	ND	ND	ND	ND	ND	0.00529
VS5	4/2/18	ND	ND	ND	ND	ND	0.00850
VS7	4/2/18	0.10	ND	ND	ND	ND	0.665
VS18	4/11/18	ND	ND	ND	ND	ND	0.0453
VS20	4/11/18	ND	ND	ND	ND	ND	0.252
VS22	4/11/18	ND	ND	ND	ND	ND	0.0430
VS24	4/11/18	ND	0.021	ND	ND	ND	1.05
VS35	4/11/18	ND	0.057	ND	ND	ND	0.158
VS37	4/11/18	ND	ND	ND	ND	ND	0.0529
VS39	4/11/18	ND	ND	ND	ND	ND	0.117
VS41	4/11/18	ND	ND	ND	ND	ND	0.215
VS53	4/13/18	ND	ND	ND	ND	ND	0.0123
VS55	4/13/18	ND	ND	ND	ND	ND	0.0501
VS57	4/13/18	ND	ND	ND	ND	ND	0.0622
VS59	4/13/18	ND	ND	ND	ND	ND	0.332
VS65	4/13/18	ND	ND	ND	ND	ND	0.139
SCREENING LEVELS							
Hazardous Waste Criteria²		50	50			1x10 ⁷	N/A
DTSC SLs - residential³		N/A	N/A			N/A	5
DTSC SLs - commercial³		N/A	N/A			N/A	22
USEPA RSLs - residential⁴		0.24	0.24			N/A	5
USEPA RSLs - commercial⁴		0.97	0.99			N/A	22

Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg) except for dioxin/furan results which are expressed in picograms per gram (pg/g) [parts per trillion]

2. California Code of Regulations Title 22, Division 4.5

3. DTSC-Modified Screening Levels (DTSC-SLs), January 2018, and DTSC HERO Note No. 2, April 2017 (preliminary remediation goals for dioxins/furans)

4. USEPA Regional Screening Levels, TR=1E-06, HQ=1"; November 2017

5. N/A = screening value not established

6. See laboratory report for other dioxin/furan compounds tested

7. TEQ = Toxic Equivalency calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs)

TABLE IV
SUMMARY OF TOTAL METALS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6010B/200.7																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
SP1-1	4/4/18	ND	ND	ND	54	ND	ND	5.4	3.7	7.8	22	ND	5.1	ND	ND	12	36	ND
SP1-2	4/4/18	ND	ND	ND	65	ND	ND	5.4	4.0	3.7	ND	ND	5.4	ND	ND	16	29	ND
SP1-3	4/4/18	ND	ND	ND	61	ND	ND	5.3	3.9	4.0	ND	ND	5.1	ND	ND	16	28	ND
SP1-4	4/4/18	ND	ND	ND	69	ND	ND	5.3	4.6	4.2	8.4	ND	5.9	ND	ND	17	32	ND
SP1-5	4/4/18	ND	ND	ND	41	ND	ND	4.6	3.0	11	45	ND	3.9	ND	ND	8.0	30	ND
SP1-6	4/4/18	ND	ND	ND	44	ND	ND	4.2	3.0	5.1	24	ND	4.2	ND	ND	11	24	ND
SP1-7	4/4/18	ND	ND	ND	60	ND	ND	5.3	4.2	5.1	11	ND	5.5	ND	ND	14	29	ND
SP1-8	4/4/18	ND	ND	ND	66	ND	ND	6.1	3.6	7.2	13	ND	5.1	ND	ND	15	46	ND
SP1-9	4/4/18	ND	ND	ND	67	ND	ND	5.7	4.1	5.3	ND	ND	5.6	ND	ND	17	35	ND
SP1-10	4/4/18	ND	ND	ND	64	ND	ND	5.5	3.9	5.8	ND	ND	5.8	ND	ND	16	28	ND
SP1-11	4/4/18	ND	ND	ND	54	ND	ND	5.4	3.7	6.1	13	ND	4.9	ND	ND	13	28	ND
SP1-12	4/4/18	ND	ND	ND	55	ND	ND	4.5	3.7	4.6	12	ND	4.9	ND	ND	14	26	ND
SP1-13	4/4/18	ND	ND	ND	52	ND	ND	4.6	3.6	6.0	22	ND	4.5	ND	ND	13	29	ND
SP1-14	4/4/18	ND	ND	ND	52	ND	ND	4.2	3.6	3.2	21	ND	3.9	ND	ND	12	21	ND
SP1-15	4/4/18	ND	ND	ND	52	ND	ND	6.0	3.9	5.7	18	ND	5.0	ND	ND	14	27	ND
SP1-16	4/4/18	ND	ND	ND	64	ND	ND	5.6	5.1	5.2	ND	ND	6.5	ND	ND	17	31	ND
SP1-17	4/4/18	ND	ND	ND	49	ND	ND	5.0	3.6	3.9	9.4	ND	4.6	ND	ND	13	23	ND
SP1-18	4/4/18	ND	ND	ND	62	ND	ND	5.2	3.9	4.4	13	ND	5.7	ND	ND	15	27	ND
SP1-19	4/4/18	ND	ND	ND	58	ND	ND	5.4	3.8	4.6	ND	ND	5.3	ND	ND	14	28	ND
SP1-20	4/4/18	ND	ND	ND	57	ND	ND	5.9	3.9	4.6	10	ND	5.3	ND	ND	16	30	ND
SP2-1	4/5/18	ND	ND	ND	58	ND	ND	6.2	4.9	5.2	13	ND	5.9	ND	ND	20	33	0.52
SP2-2	4/5/18	ND	ND	ND	55	ND	ND	5.5	4.7	3.5	11	ND	5.3	ND	ND	20	29	ND
SP2-3	4/5/18	ND	ND	ND	63	ND	ND	6.7	4.2	5.5	ND	ND	6.2	ND	ND	22	33	ND
SP2-4	4/5/18	ND	ND	ND	62	ND	ND	6.5	4.9	5.5	9.8	ND	6.6	ND	ND	22	30	ND
SP2-5	4/5/18	ND	ND	ND	59	ND	ND	6.2	4.4	5.6	17	ND	5.9	ND	ND	18	31	ND
SP2-6	4/5/18	ND	ND	ND	68	ND	ND	7.1	5.0	5.7	13	ND	7.5	ND	ND	24	33	ND
SP2-7	4/5/18	ND	ND	ND	58	ND	ND	6.0	4.6	4.1	ND	ND	6.4	ND	ND	22	24	ND
SP2-8	4/5/18	ND	ND	ND	61	ND	ND	5.5	4.4	6.5	11	ND	5.1	ND	ND	16	27	ND
SP2-9	4/5/18	ND	ND	ND	52	ND	ND	5.5	5.6	8.5	32	ND	9.5	ND	ND	17	28	ND
SP2-10	4/5/18	ND	ND	ND	61	ND	ND	6.1	4.6	4.1	ND	ND	6.2	ND	ND	21	26	ND
SP2-11	4/5/18	ND	ND	ND	58	ND	ND	6.7	4.4	6.5	17	ND	6.4	ND	ND	20	31	ND
SP2-12	4/5/18	ND	ND	ND	58	ND	ND	5.8	4.5	5.8	11	ND	5.4	ND	ND	17	29	ND

TABLE IV
SUMMARY OF TOTAL METALS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY
17100 Murphy Parkway, Lathrop, San Joaquin County, California
Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6010B/200.7																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
SP3-1	4/12/18	ND	ND	ND	61	ND	ND	5.8	4.5	4.5	ND	ND	6.0	ND	ND	17	30	ND
SP3-2	4/12/18	ND	ND	ND	72	ND	ND	5.6	4.7	8.0	ND	ND	6.6	ND	ND	17	33	ND
SP3-3	4/12/18	ND	ND	ND	41	ND	ND	4.4	3.5	5.0	20	ND	4.1	ND	ND	10	24	ND
SP3-4	4/12/18	ND	ND	ND	53	ND	ND	5.1	4.4	4.3	ND	ND	5.7	ND	ND	15	26	ND
SP3-5	4/12/18	ND	ND	ND	64	ND	ND	4.9	4.2	5.2	ND	ND	5.1	ND	ND	15	29	ND
SP3-6	4/12/18	ND	ND	ND	43	ND	ND	5.1	3.8	4.4	14	ND	4.5	ND	ND	11	23	ND
SP3-7	4/12/18	ND	ND	ND	43	ND	ND	4.6	3.8	3.3	10	ND	4.1	ND	ND	11	25	ND
SP3-8	4/12/18	ND	ND	ND	67	ND	ND	5.7	4.9	4.6	17	ND	6.6	ND	ND	17	26	ND
SP3-9	4/12/18	ND	ND	ND	59	ND	ND	5.9	5.6	4.8	ND	ND	6.6	ND	ND	18	29	ND
SP3-10	4/12/18	ND	ND	ND	57	ND	ND	5.7	4.9	4.6	ND	ND	5.7	ND	ND	14	28	ND
SP3-11	4/12/18	ND	ND	ND	56	ND	ND	5.0	4.4	4.0	ND	ND	5.7	ND	ND	15	28	ND
SP3-12	4/12/18	ND	ND	ND	38	ND	ND	3.8	3.9	7.6	38	ND	3.6	ND	ND	8.3	22	ND
SP4-1	4/12/18	ND	ND	ND	31	ND	ND	3.5	3.2	7.0	25	ND	2.9	ND	ND	7.8	21	ND
SP5-1	4/13/18	ND	ND	ND	59	ND	ND	8.6	4.3	4.4	ND	ND	6.2	ND	ND	20	33	ND
SP5-2	4/13/18	ND	ND	ND	63	ND	ND	6.4	5.3	4.9	ND	ND	7.2	ND	ND	20	30	ND
SP5-3	4/13/18	ND	ND	ND	61	ND	ND	7.0	4.5	8.6	ND	ND	6.4	ND	ND	18	30	ND
SP5-4	4/13/18	ND	ND	ND	64	ND	ND	6.7	4.9	5.2	ND	ND	7.2	ND	ND	20	31	ND
SCREENING LEVELS																		
Hazardous Waste Criteria²		500	500	500	10,000	75	100	2,500	8,000	2,500	1,000	3,500	2,000	100	700	2,400	5,000	20
DTSC SLs - residential³		N/A	390	0.11 ⁶	N/A	15	5.2	36,000	N/A	N/A	80	N/A	490	N/A	N/A	390	N/A	1.0
DTSC SLs - commercial³		N/A	1,500	0.36 ⁶	N/A	210	7.3	170,000	N/A	N/A	320	N/A	3,100	N/A	N/A	1,000	N/A	4.5
USEPA RSLs - residential⁴		31	390	0.68 ⁶	15,000	160	71	120,000	23	3,100	400	390	1,500	390	0.78	390	23,000	11
USEPA RSLs - commercial⁴		470	5,800	3.0 ⁶	220,000	2,300	980	1,800,000	350	47,000	800	5,800	22,000	5,800	12	5,800	350,000	46

Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg)

2. California Code of Regulations Title 22, Division 4.5

3. DTSC-Modified Screening Levels (DTSC-SLs); January 2018

4. USEPA Regional Screening Levels, TR=1E-06, HQ=1"; November 2017

5. N/A = screening value not established

6. Arsenic data for the project are compared to background concentration of 8.6 mg/kg rather than cited screening levels.

TABLE V
SUMMARY OF DI WET METALS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY
17100 Murphy Parkway, Lathrop, San Joaquin County, California
Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TITLE 22 METALS EPA Method 6020																
		Antimony	Silver	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Molybdenum	Nickel	Selenium	Thallium	Vanadium	Zinc	Mercury
SP1-1	4/4/18	0.002	0.005	0.030	0.48	0.001	0.002	0.069	0.045	0.12	0.30	0.021	0.06	ND	ND	0.26	0.44	0.0002
SP1-5	4/4/18	0.002	0.009	0.020	0.34	ND	0.002	0.060	0.039	0.20	0.36	0.016	0.045	ND	ND	0.16	0.35	0.0003
SP1-6	4/4/18	0.002	0.005	0.024	0.51	0.001	0.002	0.056	0.039	0.11	0.39	0.016	0.054	ND	ND	0.19	0.38	0.0002
SP1-8	4/4/18	0.001	0.004	0.027	0.52	0.001	0.001	0.057	0.037	0.085	0.14	0.024	0.053	ND	ND	0.24	0.37	0.0001
SP2-1	4/5/18	ND	0.005	0.017	0.45	0.001	ND	0.065	0.049	0.096	0.17	0.009	0.061	ND	ND	0.22	0.34	0.0003
SP2-6	4/5/18	ND	0.003	0.028	0.52	0.002	ND	0.054	0.044	0.099	0.11	0.007	0.073	ND	ND	0.25	0.32	0.0002
SP2-9	4/5/18	0.001	0.001	0.009	0.12	0.0004	0.0003	0.016	0.013	0.034	0.051	0.006	0.015	0.002	ND	0.080	0.068	0.0001
SP3-3	4/12/18	ND	0.003	0.013	0.34	ND	0.002	0.040	0.033	0.12	0.17	0.007	0.041	ND	ND	0.099	0.21	0.0003
SP3-12	4/12/18	ND	0.007	0.009	0.360	0.001	0.002	0.040	0.049	0.11	0.20	0.005	0.054	ND	ND	0.11	0.23	0.0001
SP4-1	4/12/18	ND	0.003	0.011	0.23	ND	0.002	0.033	0.031	0.11	0.17	0.007	0.028	ND	ND	0.075	0.25	0.00003
SP5-1	4/14/18	ND	ND	0.031	0.40	0.001	ND	0.047	0.032	0.055	0.046	0.007	0.047	ND	ND	0.20	0.20	ND
SCREENING LEVELS																		
MCLs ²		0.006	N/A	0.01	1.0	0.004	0.005	0.05	N/A	1.3	0.015	N/A	0.1	0.05	0.002	N/A	N/A	0.002

- Notes: 1. All concentrations expressed in milligrams per liter (mg/L)
2. Maximum Contaminant Levels (MCLs) for drinking water; California; July 2014
3. N/A = MCL not established

TABLE VI
SUMMARY OF PETROLEUM HYDROCARBONS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TOTAL PETROLEUM HYDROCARBONS		
		EPA Method 8015C		
		C6-C12 (Gasoline Range)	C13-C28 (Diesel Range)	C29-C40 (Motor Oil Range)
SP1-1	4/4/18	ND	47	100
SP1-2	4/4/18	ND	13	86
SP1-3	4/4/18	ND	47	68
SP1-4	4/4/18	ND	24	75
SP1-5	4/4/18	ND	96	52
SP1-6	4/4/18	ND	44	78
SP1-7	4/4/18	ND	88	92
SP1-8	4/4/18	ND	28	64
SP1-9	4/4/18	ND	44	28
SP1-10	4/4/18	ND	18	58
SP1-11	4/4/18	ND	75	87
SP1-12	4/4/18	ND	97	55
SP1-13	4/4/18	ND	71	74
SP1-14	4/4/18	ND	16	35
SP1-15	4/4/18	ND	170	240
SP1-16	4/4/18	ND	85	55
SP1-17	4/4/18	ND	81	40
SP1-18	4/4/18	ND	26	82
SP1-19	4/4/18	ND	39	51
SP1-20	4/4/18	ND	76	150
SP2-1	4/5/18	ND	62	59
SP2-2	4/5/18	ND	24	42
SP2-3	4/5/18	ND	18	55
SP2-4	4/5/18	ND	94	69
SP2-5	4/5/18	ND	51	46
SP2-6	4/5/18	ND	90	49
SP2-7	4/5/18	ND	60	57
SP2-8	4/5/18	ND	170	120
SP2-9	4/5/18	ND	370	380
SP2-10	4/5/18	ND	150	76
SP2-11	4/5/18	ND	110	68
SP2-12	4/5/18	ND	140	150
SP3-1	4/12/18	ND	29	29
SP3-2	4/12/18	ND	150	160
SP3-3	4/12/18	ND	51	37
SP3-4	4/12/18	ND	100	96
SP3-5	4/12/18	ND	32	15
SP3-6	4/12/18	ND	56	32

TABLE VI
SUMMARY OF PETROLEUM HYDROCARBONS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	TOTAL PETROLEUM HYDROCARBONS <i>EPA Method 8015C</i>		
		C6-C12 (Gasoline Range)	C13-C28 (Diesel Range)	C29-C40 (Motor Oil Range)
SP3-7	4/12/18	ND	39	19
SP3-8	4/12/18	ND	29	17
SP3-9	4/12/18	ND	85	82
SP3-10	4/12/18	ND	360	380
SP3-11	4/12/18	ND	26	33
SP3-12	4/12/18	ND	140	130
SP4-1	4/12/18	ND	170	220
SP5-1	4/13/18	ND	53	61
SP5-2	4/13/18	ND	70	58
SP5-3	4/13/18	ND	35	50
SP5-4	4/13/18	ND	17	56
SCREENING LEVELS				
ESLs - residential²		740	230	11,000
ESLs - commercial²		3,900	1,100	140,000

Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg)

2. San Francisco Bay Regional Water Quality Control Board; "Environmental Screening Levels"; February 2016; Direct Exposure Human Health Levels; residential and commercial/industrial.

TABLE VII
SUMMARY OF SVOCs, PCBs, AND DIOXINS/FURANS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	PCBs EPA Method 8020				SEMI-VOLATILE ORGANIC COMPOUNDS EPA Method 8270C										DIOXINS/FURANS EPA Method 1613	
		PCB -1016	PCB-1254	PCB-1260	Other PCBs	Pyrene	Benzo (a) anthracene	Benzo (b) fluoranthene	Benzo (k) fluoranthene	Benzo (a) pyrene	Chrysene	Fluoranthene	Indeno (1,2,3-cd) pyrene	Phenanthrene	Other SVOCs	2,3,7,8-TCDD ⁶	TEQ ⁷
SP1-1	4/4/18	ND	0.035	ND	ND	2.4	0.90	0.73	0.41	0.36	1.1	2.1	0.34	1.3	ND	0.375	6.49
SP1-2	4/4/18	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.668
SP1-3	4/4/18	ND	0.025	ND	ND	0.30	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.10
SP1-4	4/4/18	ND	0.013	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.510
SP1-5	4/4/18	ND	0.047	ND	ND	0.42	ND	ND	ND	ND	ND	0.36	ND	ND	ND	ND	4.88
SP1-6	4/4/18	ND	0.052	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	5.00
SP1-7	4/4/18	ND	0.043	ND	ND	0.49	ND	ND	ND	ND	ND	0.42	ND	ND	ND	ND	1.65
SP1-8	4/4/18	ND	0.032	ND	ND	0.36	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.93
SP1-9	4/4/18	ND	0.011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.139
SP1-10	4/4/18	ND	0.013	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.23
SP1-11	4/4/18	ND	0.061	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.94
SP1-12	4/4/18	ND	0.024	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.08
SP1-13	4/4/18	ND	0.043	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.89
SP1-14	4/4/18	ND	0.019	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.316
SP1-15	4/4/18	ND	0.110	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	3.20
SP1-16	4/4/18	ND	0.025	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.97
SP1-17	4/4/18	ND	0.018	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.429
SP1-18	4/4/18	ND	0.042	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.891
SP1-19	4/4/18	ND	0.021	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.84
SP1-20	4/4/18	ND	0.050	ND	ND	0.42	ND	ND	ND	ND	ND	0.34	ND	ND	ND	ND	3.62
SP2-1	4/5/18	ND	ND	0.029	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.77

TABLE VII
SUMMARY OF SVOCs, PCBs, AND DIOXINS/FURANS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	PCBs EPA Method 8020				SEMI-VOLATILE ORGANIC COMPOUNDS EPA Method 8270C										DIOXINS/FURANS EPA Method 1613	
		PCB -1016	PCB-1254	PCB-1260	Other PCBs	Pyrene	Benzo (a) anthracene	Benzo (b) fluoranthene	Benzo (k) fluoranthene	Benzo (a) pyrene	Chrysene	Fluoranthene	Indeno (1,2,3-cd) pyrene	Phenanthrene	Other SVOCs	2,3,7,8-TCDD ⁶	TEQ ⁷
SP2-2	4/5/18	0.066	ND	0.072	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	4.65
SP2-3	4/5/18	ND	0.036	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.73
SP2-4	4/5/18	ND	ND	0.077	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.96
SP2-5	4/5/18	ND	ND	0.032	ND	0.95	0.34	ND	ND	ND	0.44	0.88	ND	0.53	ND	ND	1.64
SP2-6	4/5/18	ND	ND	0.058	ND	1.2	0.49	0.41	ND	ND	0.57	1.1	ND	0.75	ND	ND	2.71
SP2-7	4/5/18	ND	ND	0.086	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.45
SP2-8	4/5/18	ND	ND	0.058	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.57
SP2-9	4/5/18	ND	ND	0.058	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	3.43
SP2-10	4/5/18	0.072	ND	0.045	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.08
SP2-11	4/5/18	ND	ND	0.026	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.75
SP2-12	4/5/18	ND	ND	0.061	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.21
SP3-1	4/12/18	ND	ND	ND	ND	0.65	0.30	0.30	ND	ND	0.34	0.73	ND	0.46	ND	ND	1.79
SP3-2	4/12/18	ND	0.020	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.62
SP3-3	4/12/18	ND	0.017	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.933
SP3-4	4/12/18	ND	0.015	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.522
SP3-5	4/12/18	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.76
SP3-6	4/12/18	ND	0.012	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.995
SP3-7	4/12/18	ND	0.012	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.823
SP3-8	4/12/18	ND	0.019	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.34
SP3-9	4/12/18	ND	0.016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.47	0.731
SP3-10	4/12/18	ND	0.018	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.492

TABLE VII
SUMMARY OF SVOCs, PCBs, AND DIOXINS/FURANS DATA
STOCKPILED WASTE MATERIAL SAMPLES

48-ACRE MURPHY PARKWAY PROPERTY

17100 Murphy Parkway, Lathrop, San Joaquin County, California

Brusca Reference No. 137-002

SAMPLE ID	DATE SAMPLED	PCBs EPA Method 8020				SEMI-VOLATILE ORGANIC COMPOUNDS EPA Method 8270C										DIOXINS/FURANS EPA Method 1613	
		PCB-1016	PCB-1254	PCB-1260	Other PCBs	Pyrene	Benzo (a) anthracene	Benzo (b) fluoranthene	Benzo (k) fluoranthene	Benzo (a) pyrene	Chrysene	Fluoranthene	Indeno (1,2,3-cd) pyrene	Phenanthrene	Other SVOCs	2,3,7,8-TCDD ⁶	TEQ ⁷
SP3-11	4/12/18	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.616
SP3-12	4/12/18	ND	0.034	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.79
SP4-1	4/12/18	0.17	0.11	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.91
SP5-1	4/13/18	ND	0.017	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.759
SP5-2	4/13/18	ND	0.014	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.670
SP5-3	4/13/18	ND	0.035	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.24
SP5-4	4/13/18	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.539
SCREENING LEVELS																	
Hazardous Waste Criteria²		50	50	50	50	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A		1x10 ⁷	N/A
DTSC SLs - residential³		N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A		N/A	5
DTSC SLs - commercial³		N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A		N/A	22
USEPA RSLs - residential⁴		4.1	0.24	0.24		1,800	1.1	1.1	11	0.11	110	2,400	1.1	N/A		N/A	5
USEPA RSLs - commercial⁴		27	0.97	0.99		23,000	21	21	210	2.1	2,100	30,000	21	N/A		N/A	22

- Notes: 1. All concentrations expressed in milligrams per kilogram (mg/kg) except for dioxin/furan results which are expressed in picograms per gram (pg/g) [parts per trillion]
2. California Code of Regulations Title 22, Division 4.5
3. DTSC-Modified Screening Levels (DTSC-SLs), January 2018, and DTSC HERO Note No. 2, April 2017 (preliminary remediation goals for dioxins/furans)
4. USEPA Regional Screening Levels, TR=1E-06, HQ=1"; November 2017
5. N/A = screening value not established
6. See laboratory report for other dioxin/furan compounds tested
7. TEQ = Toxic Equivalency calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs)

APPENDIX A – Laboratory Reports



25712 Commercentre Drive
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10 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/03/18 10:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
VS1	T181118-01	Soil	04/02/18 11:45	04/03/18 10:00
VS2	T181118-02	Soil	04/02/18 11:49	04/03/18 10:00
VS3	T181118-03	Soil	04/02/18 11:52	04/03/18 10:00
VS4	T181118-04	Soil	04/02/18 11:55	04/03/18 10:00
VS5	T181118-05	Soil	04/02/18 11:57	04/03/18 10:00
VS7	T181118-07	Soil	04/02/18 12:03	04/03/18 10:00
VS8	T181118-08	Soil	04/02/18 12:05	04/03/18 10:00

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

DETECTIONS SUMMARY

Sample ID: VS1

Laboratory ID: T181118-01

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	49	1.0	mg/kg	EPA 6010B	
Chromium	5.2	2.0	mg/kg	EPA 6010B	
Cobalt	5.0	2.0	mg/kg	EPA 6010B	
Copper	3.3	1.0	mg/kg	EPA 6010B	
Nickel	5.8	2.0	mg/kg	EPA 6010B	
Vanadium	19	5.0	mg/kg	EPA 6010B	
Zinc	21	1.0	mg/kg	EPA 6010B	

Sample ID: VS2

Laboratory ID: T181118-02

No Results Detected

Sample ID: VS3

Laboratory ID: T181118-03

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	55	1.0	mg/kg	EPA 6010B	
Chromium	4.2	2.0	mg/kg	EPA 6010B	
Cobalt	4.2	2.0	mg/kg	EPA 6010B	
Copper	1.4	1.0	mg/kg	EPA 6010B	
Nickel	6.2	2.0	mg/kg	EPA 6010B	
Vanadium	16	5.0	mg/kg	EPA 6010B	
Zinc	17	1.0	mg/kg	EPA 6010B	

Sample ID: VS4

Laboratory ID: T181118-04

No Results Detected

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Sample ID: VS5

Laboratory ID: T181118-05

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	47	1.0	mg/kg	EPA 6010B	
Chromium	4.8	2.0	mg/kg	EPA 6010B	
Cobalt	4.0	2.0	mg/kg	EPA 6010B	
Nickel	3.9	2.0	mg/kg	EPA 6010B	
Vanadium	18	5.0	mg/kg	EPA 6010B	
Zinc	18	1.0	mg/kg	EPA 6010B	

Sample ID: VS7

Laboratory ID: T181118-07

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	950	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	2100	10	mg/kg	EPA 8015B	
Barium	45	1.0	mg/kg	EPA 6010B	
Chromium	5.2	2.0	mg/kg	EPA 6010B	
Cobalt	3.7	2.0	mg/kg	EPA 6010B	
Nickel	4.1	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	19	1.0	mg/kg	EPA 6010B	
PCB-1254	100	10	ug/kg	EPA 8082	

Sample ID: VS8

Laboratory ID: T181118-08

No Results Detected

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS1

T181118-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040323	04/03/18	04/04/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		81.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	49	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.2	2.0	"	"	"	"	"	"	
Cobalt	5.0	2.0	"	"	"	"	"	"	
Copper	3.3	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	19	5.0	"	"	"	"	"	"	
Zinc	21	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040314	04/03/18	04/03/18	EPA 7471A Soil	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS1

T181118-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040318	04/03/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

119 %

35-140

"

"

"

"

Surrogate: Decachlorobiphenyl

139 %

35-140

"

"

"

"

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS1

T181118-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS1

T181118-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	58.5 %	15-121			"	"	"	"	
Surrogate: Phenol-d6	61.7 %	24-113			"	"	"	"	
Surrogate: Nitrobenzene-d5	82.7 %	21.3-119			"	"	"	"	
Surrogate: 2-Fluorobiphenyl	79.5 %	32.4-102			"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	73.3 %	18.1-105			"	"	"	"	
Surrogate: Terphenyl-d14	132 %	29.1-130			"	"	"	"	S-GC

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS2

T181118-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

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04/10/18 16:14

VS3

T181118-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040323	04/03/18	04/04/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		72.8 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	55	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.2	2.0	"	"	"	"	"	"	
Cobalt	4.2	2.0	"	"	"	"	"	"	
Copper	1.4	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	17	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040314	04/03/18	04/03/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Reported:
04/10/18 16:14

VS3

T181118-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040318	04/03/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		106 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		146 %	35-140		"	"	"	"	S-GC

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS3

T181118-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS3

T181118-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	66.0 %	15-121			"	"	"	"	
Surrogate: Phenol-d6	69.0 %	24-113			"	"	"	"	
Surrogate: Nitrobenzene-d5	90.2 %	21.3-119			"	"	"	"	
Surrogate: 2-Fluorobiphenyl	84.5 %	32.4-102			"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	77.8 %	18.1-105			"	"	"	"	
Surrogate: Terphenyl-d14	133 %	29.1-130			"	"	"	"	S-GC

SunStar Laboratories, Inc.

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Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS4

T181118-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS5

T181118-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040323	04/03/18	04/04/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: p-Terphenyl		71.9 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	47	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.8	2.0	"	"	"	"	"	"	
Cobalt	4.0	2.0	"	"	"	"	"	"	
Copper	ND	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	18	5.0	"	"	"	"	"	"	
Zinc	18	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040314	04/03/18	04/03/18	EPA 7471A Soil	

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Reported:
04/10/18 16:14

VS5

T181118-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040318	04/03/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		124 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		137 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS5

T181118-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS5

T181118-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	66.2 %	15-121			"	"	"	"
Surrogate: Phenol-d6	67.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	91.4 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	83.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	76.0 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	126 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS7

T181118-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040323	04/03/18	04/04/18	EPA 8015B	
C13-C28 (DRO)	950	10	"	"	"	"	"	"	
C29-C40 (MORO)	2100	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		87.6 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	45	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.2	2.0	"	"	"	"	"	"	
Cobalt	3.7	2.0	"	"	"	"	"	"	
Copper	ND	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	19	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040314	04/03/18	04/03/18	EPA 7471A Soil	

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Reported:
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VS7

T181118-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040318	04/03/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	100	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		107 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		106 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS7

T181118-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

VS7

T181118-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040221	04/04/18	04/05/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		55.5 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		59.7 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		81.5 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		78.4 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		62.0 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		92.8 %	29.1-130		"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
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Reported:
04/10/18 16:14

VS8

T181118-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8040313	04/03/18	04/03/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040323 - EPA 3550B GC

Blank (8040323-BLK1)

Prepared: 04/03/18 Analyzed: 04/04/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	71.0		"	100		71.0	65-135			

LCS (8040323-BS1)

Prepared & Analyzed: 04/03/18

C13-C28 (DRO)	500	10	mg/kg	495		100	75-125			
Surrogate: <i>p</i> -Terphenyl	79.8		"	99.0		80.6	65-135			

LCS Dup (8040323-BSD1)

Prepared & Analyzed: 04/03/18

C13-C28 (DRO)	540	10	mg/kg	505		106	75-125	7.88	20	
Surrogate: <i>p</i> -Terphenyl	82.4		"	101		81.6	65-135			

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Reported:
04/10/18 16:14

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040313 - EPA 3050B

Blank (8040313-BLK1)

Prepared & Analyzed: 04/03/18

Antimony	ND	3.0	mg/kg
Silver	ND	2.0	"
Arsenic	ND	5.0	"
Arsenic	ND	5.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	2.0	"
Chromium	ND	2.0	"
Cobalt	ND	2.0	"
Copper	ND	1.0	"
Lead	ND	3.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Lead	ND	3.0	"
Selenium	ND	5.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	ND	1.0	"

LCS (8040313-BS1)

Prepared & Analyzed: 04/03/18

Arsenic	103	5.0	mg/kg	100	103	75-125
Arsenic	103	5.0	"	100	103	75-125
Barium	104	1.0	"	100	104	75-125
Cadmium	103	2.0	"	100	103	75-125
Chromium	104	2.0	"	100	104	75-125
Lead	109	3.0	"	100	109	75-125
Lead	109	3.0	"	100	109	75-125

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Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040313 - EPA 3050B

Matrix Spike (8040313-MS1)

Source: T181119-01

Prepared & Analyzed: 04/03/18

Arsenic	90.6	5.0	mg/kg	100	1.36	89.3	75-125			
Arsenic	90.6	5.0	"	100	1.36	89.3	75-125			
Barium	161	1.0	"	100	66.0	95.4	75-125			
Cadmium	90.0	2.0	"	100	0.633	89.3	75-125			
Chromium	99.2	2.0	"	100	8.01	91.2	75-125			
Lead	95.6	3.0	"	100	ND	95.6	75-125			
Lead	95.6	3.0	"	100	2.80	92.8	75-125			

Matrix Spike Dup (8040313-MSD1)

Source: T181119-01

Prepared & Analyzed: 04/03/18

Arsenic	88.3	5.0	mg/kg	97.1	1.36	89.5	75-125	2.59	20	
Arsenic	88.3	5.0	"	97.1	1.36	89.5	75-125	2.59	20	
Barium	156	1.0	"	97.1	66.0	92.8	75-125	3.29	20	
Cadmium	87.9	2.0	"	97.1	0.633	89.9	75-125	2.30	20	
Chromium	96.9	2.0	"	97.1	8.01	91.6	75-125	2.36	20	
Lead	94.6	3.0	"	97.1	ND	97.5	75-125	1.04	20	
Lead	94.6	3.0	"	97.1	2.80	94.6	75-125	1.04	20	

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Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040314 - EPA 7471A Soil

Blank (8040314-BLK1)

Prepared & Analyzed: 04/03/18

Mercury	ND	0.10	mg/kg							

LCS (8040314-BS1)

Prepared & Analyzed: 04/03/18

Mercury	0.431	0.10	mg/kg	0.417		104	80-120			

Matrix Spike (8040314-MS1)

Source: T181119-01

Prepared & Analyzed: 04/03/18

Mercury	0.395	0.10	mg/kg	0.391	ND	101	75-125			

Matrix Spike Dup (8040314-MSD1)

Source: T181119-01

Prepared & Analyzed: 04/03/18

Mercury	0.426	0.10	mg/kg	0.417	ND	102	75-125	7.63	20	

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Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040318 - EPA 3550 ECD/GCMS

Blank (8040318-BLK1)

Prepared: 04/03/18 Analyzed: 04/06/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	11.0		"	9.80		112	35-140			
Surrogate: Decachlorobiphenyl	12.7		"	9.80		130	35-140			

LCS (8040318-BS1)

Prepared: 04/03/18 Analyzed: 04/06/18

PCB-1016	72.2	10	ug/kg	98.0		73.6	40-130			
PCB-1260	66.4	10	"	98.0		67.7	40-130			
Surrogate: Tetrachloro-meta-xylene	11.7		"	9.80		119	35-140			
Surrogate: Decachlorobiphenyl	14.0		"	9.80		142	35-140			S-GC

Matrix Spike (8040318-MS1)

Source: T181118-01

Prepared: 04/03/18 Analyzed: 04/06/18

PCB-1016	53.3	10	ug/kg	99.0	ND	53.8	40-130			
PCB-1260	58.4	10	"	99.0	ND	59.0	40-130			
Surrogate: Tetrachloro-meta-xylene	12.2		"	9.90		123	35-140			
Surrogate: Decachlorobiphenyl	16.5		"	9.90		166	35-140			S-GC

Matrix Spike Dup (8040318-MSD1)

Source: T181118-01

Prepared: 04/03/18 Analyzed: 04/06/18

PCB-1016	50.4	10	ug/kg	99.0	ND	50.9	40-130	5.54	30	
PCB-1260	50.6	10	"	99.0	ND	51.1	40-130	14.4	30	
Surrogate: Tetrachloro-meta-xylene	11.2		"	9.90		113	35-140			
Surrogate: Decachlorobiphenyl	12.4		"	9.90		125	35-140			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040221 - EPA 3550 ECD/GCMS

Blank (8040221-BLK1)

Prepared: 04/04/18 Analyzed: 04/05/18

Carbazole	ND	300	ug/kg
Aniline	ND	300	"
Phenol	ND	1000	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
2-Methylnaphthalene	ND	300	"
1-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040221 - EPA 3550 ECD/GCMS

Blank (8040221-BLK1)

Prepared: 04/04/18 Analyzed: 04/05/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
Pyridine	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
2,4,6-Trichlorophenol	ND	1000	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040221 - EPA 3550 ECD/GCMS

Blank (8040221-BLK1)

Prepared: 04/04/18 Analyzed: 04/05/18

Surrogate: 2-Fluorophenol	1990		ug/kg	3440		58.0	15-121			
Surrogate: Phenol-d6	2040		"	3440		59.4	24-113			
Surrogate: Nitrobenzene-d5	2900		"	3440		84.3	21.3-119			
Surrogate: 2-Fluorobiphenyl	2670		"	3440		77.7	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2500		"	3440		72.7	18.1-105			
Surrogate: Terphenyl-d14	4460		"	3440		130	29.1-130			

LCS (8040221-BS1)

Prepared: 04/02/18 Analyzed: 04/05/18

Phenol	1840	1000	ug/kg	3230		56.9	34-114			
2-Chlorophenol	1850	1000	"	3230		57.2	34-114			
1,4-Dichlorobenzene	1970	300	"	3230		61.0	34-114			
N-Nitrosodi-n-propylamine	2490	300	"	3230		77.3	30-110			
1,2,4-Trichlorobenzene	2180	300	"	3230		67.5	39-119			
4-Chloro-3-methylphenol	2620	1000	"	3230		81.1	50-130			
Acenaphthene	2460	300	"	3230		76.3	34-114			
Pentachlorophenol	1700	1000	"	3230		52.8	50-130			
Pyrene	2400	300	"	3230		74.4	30-110			
Surrogate: 2-Fluorophenol	1780		"	3230		55.0	15-121			
Surrogate: Phenol-d6	1870		"	3230		57.9	24-113			
Surrogate: Nitrobenzene-d5	2640		"	3230		81.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2480		"	3230		76.8	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2530		"	3230		78.6	18.1-105			
Surrogate: Terphenyl-d14	4310		"	3230		133	29.1-130			S-GC

Matrix Spike (8040221-MS1)

Source: T181108-01

Prepared: 04/02/18 Analyzed: 04/05/18

Phenol	2260	1000	ug/kg	3240	ND	70.0	34-114			
2-Chlorophenol	2240	1000	"	3240	ND	69.1	34-114			
1,4-Dichlorobenzene	2240	300	"	3240	ND	69.2	34-114			
N-Nitrosodi-n-propylamine	2660	300	"	3240	ND	82.3	30-110			
1,2,4-Trichlorobenzene	2510	300	"	3240	ND	77.4	39-119			
4-Chloro-3-methylphenol	2680	1000	"	3240	ND	82.9	50-130			
Acenaphthene	2680	300	"	3240	ND	82.8	34-114			
Pentachlorophenol	1660	1000	"	3240	ND	51.4	50-130			
Pyrene	2370	300	"	3240	37.5	72.0	30-110			
Surrogate: 2-Fluorophenol	2140		"	3240		66.2	15-121			
Surrogate: Phenol-d6	2270		"	3240		70.2	24-113			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040221 - EPA 3550 ECD/GCMS

Matrix Spike (8040221-MS1)

Source: T181108-01

Prepared: 04/02/18 Analyzed: 04/05/18

Surrogate: Nitrobenzene-d5	3040		ug/kg	3240		93.8	21.3-119			
Surrogate: 2-Fluorobiphenyl	2560		"	3240		79.1	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2670		"	3240		82.4	18.1-105			
Surrogate: Terphenyl-d14	4020		"	3240		124	29.1-130			

Matrix Spike Dup (8040221-MSD1)

Source: T181108-01

Prepared: 04/02/18 Analyzed: 04/05/18

Phenol	2090	1000	ug/kg	3360	ND	62.2	34-114	8.06	42	
2-Chlorophenol	2260	1000	"	3360	ND	67.2	34-114	0.852	40	
1,4-Dichlorobenzene	2240	300	"	3360	ND	66.8	34-114	0.00715	28	
N-Nitrosodi-n-propylamine	2770	300	"	3360	ND	82.7	30-110	4.11	38	
1,2,4-Trichlorobenzene	2460	300	"	3360	ND	73.4	39-119	1.65	28	
4-Chloro-3-methylphenol	2710	1000	"	3360	ND	80.8	50-130	1.07	42	
Acenaphthene	2660	300	"	3360	ND	79.3	34-114	0.696	31	
Pentachlorophenol	1780	1000	"	3360	ND	52.9	50-130	6.52	50	
Pyrene	2310	300	"	3360	37.5	67.7	30-110	2.57	31	
Surrogate: 2-Fluorophenol	2160		"	3360		64.3	15-121			
Surrogate: Phenol-d6	2290		"	3360		68.1	24-113			
Surrogate: Nitrobenzene-d5	3090		"	3360		92.2	21.3-119			
Surrogate: 2-Fluorobiphenyl	2620		"	3360		78.0	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2730		"	3360		81.3	18.1-105			
Surrogate: Terphenyl-d14	4160		"	3360		124	29.1-130			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/10/18 16:14

Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

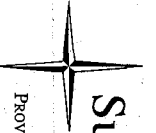
dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, Rosemead, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/2/18 Page: 1 of 1
Project Name: LATHROP 48-AC PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 718118 EDF #:

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD AND ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers
VS1	4/2/18	11:45	SOIL	JAR	X			X				X	X			01		1
VS2		11:49														02		1
VS3		11:52			X			X				X	X			03		1
VS4		11:53			X			X				X	X			04		1
VS5		11:57			X			X				X	X			05		1
VS6		12:00														06		1
VS7		12:03														07		1
VS8		12:05			X							X				08		1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/2/18, 2:37 PM</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/2/18 2:37 PM</u>																		
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/3/18 10:00</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/3/18 10:00</u>																		
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/3/18 10:00</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/3/18 10:00</u>																		
Total # of containers: <u>4,9</u>																		
Chain of Custody seals Y/N/NA: <u>Seals intact? Y/N/NA</u>																		
Received good condition/cold: <u>4,9</u>																		
Turn around time: <u>NORM</u>																		

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

COC 160903



25712 Commercentre Drive
Lake Forest, California 92630
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949.297.5027 Fax

19 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/12/18 10:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
VS9	T181234-01	Soil	04/11/18 11:00	04/12/18 10:00
VS10	T181234-02	Soil	04/11/18 11:02	04/12/18 10:00
VS11	T181234-03	Soil	04/11/18 11:03	04/12/18 10:00
VS12	T181234-04	Soil	04/11/18 11:05	04/12/18 10:00
VS13	T181234-05	Soil	04/11/18 11:07	04/12/18 10:00
VS14	T181234-06	Soil	04/11/18 11:09	04/12/18 10:00
VS16	T181234-08	Soil	04/11/18 11:14	04/12/18 10:00
VS17	T181234-09	Soil	04/11/18 11:20	04/12/18 10:00
VS18	T181234-10	Soil	04/11/18 11:22	04/12/18 10:00
VS19	T181234-11	Soil	04/11/18 11:26	04/12/18 10:00
VS20	T181234-12	Soil	04/11/18 11:29	04/12/18 10:00
VS21	T181234-13	Soil	04/11/18 11:31	04/12/18 10:00
VS22	T181234-14	Soil	04/11/18 11:34	04/12/18 10:00
VS24	T181234-16	Soil	04/11/18 11:37	04/12/18 10:00
VS25	T181234-17	Soil	04/11/18 11:55	04/12/18 10:00
VS26	T181234-18	Soil	04/11/18 11:57	04/12/18 10:00
VS27	T181234-19	Soil	04/11/18 11:58	04/12/18 10:00
VS28	T181234-20	Soil	04/11/18 12:00	04/12/18 10:00
VS29	T181234-21	Soil	04/11/18 12:03	04/12/18 10:00
VS30	T181234-22	Soil	04/11/18 12:04	04/12/18 10:00
VS31	T181234-23	Soil	04/11/18 12:06	04/12/18 10:00
VS33	T181234-25	Soil	04/11/18 12:10	04/12/18 10:00
VS34	T181234-26	Soil	04/11/18 12:41	04/12/18 10:00
VS35	T181234-27	Soil	04/11/18 12:43	04/12/18 10:00
VS36	T181234-28	Soil	04/11/18 12:45	04/12/18 10:00
VS37	T181234-29	Soil	04/11/18 12:48	04/12/18 10:00

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
VS38	T181234-30	Soil	04/11/18 12:50	04/12/18 10:00
VS39	T181234-31	Soil	04/11/18 12:52	04/12/18 10:00
VS41	T181234-33	Soil	04/11/18 12:57	04/12/18 10:00
VS43	T181234-35	Soil	04/11/18 13:14	04/12/18 10:00
VS45	T181234-37	Soil	04/11/18 13:17	04/12/18 10:00
VS47	T181234-39	Soil	04/11/18 13:21	04/12/18 10:00
VS49	T181234-41	Soil	04/11/18 13:26	04/12/18 10:00
VS51	T181234-43	Soil	04/11/18 13:30	04/12/18 10:00

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

DETECTIONS SUMMARY

Sample ID: VS9 **Laboratory ID:** T181234-01

No Results Detected

Sample ID: VS10 **Laboratory ID:** T181234-02

No Results Detected

Sample ID: VS11 **Laboratory ID:** T181234-03

No Results Detected

Sample ID: VS12 **Laboratory ID:** T181234-04

No Results Detected

Sample ID: VS13 **Laboratory ID:** T181234-05

No Results Detected

Sample ID: VS14 **Laboratory ID:** T181234-06

No Results Detected

SunStar Laboratories, Inc.

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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS16

Laboratory ID: T181234-08

No Results Detected

Sample ID: VS17

Laboratory ID: T181234-09

No Results Detected

Sample ID: VS18

Laboratory ID: T181234-10

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	44	1.0	mg/kg	EPA 6010B	
Chromium	3.8	2.0	mg/kg	EPA 6010B	
Cobalt	3.4	2.0	mg/kg	EPA 6010B	
Nickel	3.8	2.0	mg/kg	EPA 6010B	
Vanadium	13	5.0	mg/kg	EPA 6010B	
Zinc	18	1.0	mg/kg	EPA 6010B	

Sample ID: VS19

Laboratory ID: T181234-11

No Results Detected

Sample ID: VS20

Laboratory ID: T181234-12

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	55	1.0	mg/kg	EPA 6010B	
Chromium	4.7	2.0	mg/kg	EPA 6010B	
Cobalt	3.6	2.0	mg/kg	EPA 6010B	
Copper	1.4	1.0	mg/kg	EPA 6010B	
Nickel	4.5	2.0	mg/kg	EPA 6010B	
Vanadium	15	5.0	mg/kg	EPA 6010B	
Zinc	22	1.0	mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS21

Laboratory ID: T181234-13

No Results Detected

Sample ID: VS22

Laboratory ID: T181234-14

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Barium	72	1.0		mg/kg	EPA 6010B	
Chromium	4.3	2.0		mg/kg	EPA 6010B	
Cobalt	4.0	2.0		mg/kg	EPA 6010B	
Copper	1.6	1.0		mg/kg	EPA 6010B	
Nickel	4.4	2.0		mg/kg	EPA 6010B	
Vanadium	16	5.0		mg/kg	EPA 6010B	
Zinc	20	1.0		mg/kg	EPA 6010B	

Sample ID: VS24

Laboratory ID: T181234-16

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	220	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	310	10		mg/kg	EPA 8015B	
Barium	52	1.0		mg/kg	EPA 6010B	
Chromium	4.2	2.0		mg/kg	EPA 6010B	
Cobalt	3.8	2.0		mg/kg	EPA 6010B	
Copper	3.1	1.0		mg/kg	EPA 6010B	
Nickel	4.9	2.0		mg/kg	EPA 6010B	
Vanadium	16	5.0		mg/kg	EPA 6010B	
Zinc	19	1.0		mg/kg	EPA 6010B	
PCB-1260	21	10		ug/kg	EPA 8082	

Sample ID: VS25

Laboratory ID: T181234-17

No Results Detected

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS26

Laboratory ID: T181234-18

No Results Detected

Sample ID: VS27

Laboratory ID: T181234-19

No Results Detected

Sample ID: VS28

Laboratory ID: T181234-20

No Results Detected

Sample ID: VS29

Laboratory ID: T181234-21

No Results Detected

Sample ID: VS30

Laboratory ID: T181234-22

No Results Detected

Sample ID: VS31

Laboratory ID: T181234-23

No Results Detected

Sample ID: VS33

Laboratory ID: T181234-25

No Results Detected

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS34

Laboratory ID: T181234-26

No Results Detected

Sample ID: VS35

Laboratory ID: T181234-27

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Barium	66	1.0		mg/kg	EPA 6010B	
Chromium	5.1	2.0		mg/kg	EPA 6010B	
Cobalt	6.3	2.0		mg/kg	EPA 6010B	
Copper	3.6	1.0		mg/kg	EPA 6010B	
Nickel	6.8	2.0		mg/kg	EPA 6010B	
Vanadium	22	5.0		mg/kg	EPA 6010B	
Zinc	28	1.0		mg/kg	EPA 6010B	
PCB-1260	57	10		ug/kg	EPA 8082	

Sample ID: VS36

Laboratory ID: T181234-28

No Results Detected

Sample ID: VS37

Laboratory ID: T181234-29

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Barium	77	1.0		mg/kg	EPA 6010B	
Chromium	5.9	2.0		mg/kg	EPA 6010B	
Cobalt	6.4	2.0		mg/kg	EPA 6010B	
Copper	4.1	1.0		mg/kg	EPA 6010B	
Nickel	7.5	2.0		mg/kg	EPA 6010B	
Vanadium	18	5.0		mg/kg	EPA 6010B	
Zinc	26	1.0		mg/kg	EPA 6010B	

Sample ID: VS38

Laboratory ID: T181234-30

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS38

Laboratory ID: T181234-30

No Results Detected

Sample ID: VS39

Laboratory ID: T181234-31

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Barium	66	1.0		mg/kg	EPA 6010B	
Chromium	4.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.8	2.0		mg/kg	EPA 6010B	
Copper	4.5	1.0		mg/kg	EPA 6010B	
Nickel	7.6	2.0		mg/kg	EPA 6010B	
Vanadium	15	5.0		mg/kg	EPA 6010B	
Zinc	20	1.0		mg/kg	EPA 6010B	

Sample ID: VS41

Laboratory ID: T181234-33

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Barium	53	1.0		mg/kg	EPA 6010B	
Chromium	4.3	2.0		mg/kg	EPA 6010B	
Cobalt	4.1	2.0		mg/kg	EPA 6010B	
Copper	3.2	1.0		mg/kg	EPA 6010B	
Nickel	4.9	2.0		mg/kg	EPA 6010B	
Vanadium	16	5.0		mg/kg	EPA 6010B	
Zinc	18	1.0		mg/kg	EPA 6010B	

Sample ID: VS43

Laboratory ID: T181234-35

No Results Detected

Sample ID: VS45

Laboratory ID: T181234-37

No Results Detected

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Sample ID: VS47

Laboratory ID: T181234-39

No Results Detected

Sample ID: VS49

Laboratory ID: T181234-41

No Results Detected

Sample ID: VS51

Laboratory ID: T181234-43

No Results Detected

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Reported:
04/19/18 13:50

VS9

T181234-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Reported:
04/19/18 13:50

VS10
T181234-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Reported:
04/19/18 13:50

VS11
T181234-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Reported:
04/19/18 13:50

VS12

T181234-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Reported:
04/19/18 13:50

VS13

T181234-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Reported:
04/19/18 13:50

VS14
T181234-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS16
T181234-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS17

T181234-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS18
T181234-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		109 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	44	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	3.8	2.0	"	"	"	"	"	"	
Cobalt	3.4	2.0	"	"	"	"	"	"	
Copper	ND	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	13	5.0	"	"	"	"	"	"	
Zinc	18	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS18
T181234-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		97.4 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		78.2 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS18
T181234-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS18
T181234-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	43.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	47.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	50.7 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	63.6 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	74.4 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	67.4 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS19

T181234-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS20

T181234-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		102 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	55	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.7	2.0	"	"	"	"	"	"	
Cobalt	3.6	2.0	"	"	"	"	"	"	
Copper	1.4	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	22	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS20

T181234-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

105 %

35-140

"

"

"

"

Surrogate: Decachlorobiphenyl

98.8 %

35-140

"

"

"

"

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS20

T181234-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS20

T181234-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		46.1 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		49.8 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		50.7 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		63.6 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		75.0 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		69.3 %	29.1-130		"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS21

T181234-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS22

T181234-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	72	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.3	2.0	"	"	"	"	"	"	
Cobalt	4.0	2.0	"	"	"	"	"	"	
Copper	1.6	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	20	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS22

T181234-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

94.7 %

35-140

"

"

"

"

Surrogate: Decachlorobiphenyl

93.0 %

35-140

"

"

"

"

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS22

T181234-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS22

T181234-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	49.4 %	15-121			"	"	"	"
Surrogate: Phenol-d6	52.9 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	54.8 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	67.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	76.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	67.4 %	29.1-130			"	"	"	"

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Brusca Associates Inc.
PO Box 332
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS24

T181234-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	220	10	"	"	"	"	"	"	
C29-C40 (MORO)	310	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		109 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	52	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.2	2.0	"	"	"	"	"	"	
Cobalt	3.8	2.0	"	"	"	"	"	"	
Copper	3.1	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	19	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS24

T181234-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	21	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		80.2 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		81.8 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS24

T181234-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS24

T181234-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	47.3 %	15-121			"	"	"	"
Surrogate: Phenol-d6	55.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	54.9 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	72.9 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	84.6 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	71.4 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS25

T181234-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS26
T181234-18 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS27

T181234-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS28

T181234-20 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS29

T181234-21 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS30

T181234-22 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS31

T181234-23 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS33

T181234-25 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS34

T181234-26 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS35

T181234-27 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		94.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	66	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.1	2.0	"	"	"	"	"	"	
Cobalt	6.3	2.0	"	"	"	"	"	"	
Copper	3.6	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	22	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS35

T181234-27 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	57	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		108 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		101 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS35

T181234-27 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS35
T181234-27 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	51.1 %	15-121			"	"	"	"
Surrogate: Phenol-d6	55.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	56.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	68.9 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	78.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	73.4 %	29.1-130			"	"	"	"

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS36
T181234-28 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041304	04/13/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS37

T181234-29 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		105 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	77	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.9	2.0	"	"	"	"	"	"	
Cobalt	6.4	2.0	"	"	"	"	"	"	
Copper	4.1	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	18	5.0	"	"	"	"	"	"	
Zinc	26	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS37

T181234-29 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		93.7 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		97.2 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS37

T181234-29 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS37

T181234-29 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.8 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	62.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	75.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	84.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	77.9 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS38

T181234-30 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS39

T181234-31 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	66	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.7	2.0	"	"	"	"	"	"	
Cobalt	4.8	2.0	"	"	"	"	"	"	
Copper	4.5	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	20	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

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04/19/18 13:50

VS39

T181234-31 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		89.1 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		100 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS39

T181234-31 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS39

T181234-31 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		60.7 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		66.0 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		69.4 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		81.2 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		96.1 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		74.8 %	29.1-130		"	"	"	"	

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS41
T181234-33 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/12/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		110 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041307	04/13/18	04/13/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	53	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.3	2.0	"	"	"	"	"	"	
Cobalt	4.1	2.0	"	"	"	"	"	"	
Copper	3.2	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	18	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041308	04/13/18	04/13/18	EPA 7471A Soil	

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04/19/18 13:50

VS41

T181234-33 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041144	04/11/18	04/14/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

88.2 %

35-140

"

"

"

"

Surrogate: Decachlorobiphenyl

89.5 %

35-140

"

"

"

"

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS41

T181234-33 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS41
T181234-33 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041145	04/11/18	04/13/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	53.8 %	15-121			"	"	"	"
Surrogate: Phenol-d6	58.1 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	58.6 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	72.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	80.4 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	73.5 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS43

T181234-35 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS45

T181234-37 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Reported:
04/19/18 13:50

VS47

T181234-39 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

VS49

T181234-41 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Reported:
04/19/18 13:50

VS51

T181234-43 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	4.5	mg/kg	1	8041305	04/13/18	04/17/18	EPA 6010b	
Lead	ND	2.7	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041214 - EPA 3550B GC

Blank (8041214-BLK1)

Prepared: 04/12/18 Analyzed: 04/18/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	97.8		"	99.0		98.8	65-135			

LCS (8041214-BS1)

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	500	10	mg/kg	495		101	75-125			
Surrogate: <i>p</i> -Terphenyl	101		"	99.0		102	65-135			

Matrix Spike (8041214-MS1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	560	10	mg/kg	505	53	100	75-125			
Surrogate: <i>p</i> -Terphenyl	106		"	101		105	65-135			

Matrix Spike Dup (8041214-MSD1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	590	10	mg/kg	500	53	108	75-125	5.92	20	
Surrogate: <i>p</i> -Terphenyl	104		"	100		104	65-135			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041304 - EPA 3050B

Blank (8041304-BLK1)

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	ND	5.0	mg/kg
Lead	ND	3.0	"

LCS (8041304-BS1)

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	103	5.0	mg/kg	100	103	75-125
Lead	108	3.0	"	100	108	75-125

Matrix Spike (8041304-MS1)

Source: T181234-01

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	94.4	5.0	mg/kg	94.3	ND	100	75-125
Lead	94.7	3.0	"	94.3	0.0883	100	75-125

Matrix Spike Dup (8041304-MSD1)

Source: T181234-01

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	96.3	5.0	mg/kg	94.3	ND	102	75-125	1.96	20
Lead	101	3.0	"	94.3	0.0883	107	75-125	6.06	20

Batch 8041305 - EPA 3050B

Blank (8041305-BLK1)

Prepared: 04/13/18 Analyzed: 04/17/18

Arsenic	ND	5.0	mg/kg
Lead	ND	3.0	"

LCS (8041305-BS1)

Prepared: 04/13/18 Analyzed: 04/17/18

Arsenic	106	5.0	mg/kg	100	106	75-125
Lead	112	3.0	"	100	112	75-125

Matrix Spike (8041305-MS1)

Source: T181234-30

Prepared: 04/13/18 Analyzed: 04/17/18

Arsenic	100	5.0	mg/kg	95.2	ND	105	75-125
Lead	105	3.0	"	95.2	ND	110	75-125

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041305 - EPA 3050B

Matrix Spike Dup (8041305-MSD1)

Source: T181234-30

Prepared: 04/13/18 Analyzed: 04/17/18

Arsenic	100	5.0	mg/kg	91.7	ND	109	75-125	0.0163	20	
Lead	107	3.0	"	91.7	ND	116	75-125	1.55	20	

Batch 8041307 - EPA 3050B

Blank (8041307-BLK1)

Prepared & Analyzed: 04/13/18

Antimony	ND	3.0	mg/kg							
Silver	ND	2.0	"							
Arsenic	ND	5.0	"							
Barium	ND	1.0	"							
Beryllium	ND	1.0	"							
Cadmium	ND	2.0	"							
Chromium	ND	2.0	"							
Cobalt	ND	2.0	"							
Copper	ND	1.0	"							
Lead	ND	3.0	"							
Molybdenum	ND	5.0	"							
Nickel	ND	2.0	"							
Selenium	ND	5.0	"							
Thallium	ND	2.0	"							
Vanadium	ND	5.0	"							
Zinc	ND	1.0	"							

LCS (8041307-BS1)

Prepared & Analyzed: 04/13/18

Arsenic	94.3	5.0	mg/kg	100		94.3	75-125			
Barium	101	1.0	"	100		101	75-125			
Cadmium	99.6	2.0	"	100		99.6	75-125			
Chromium	101	2.0	"	100		101	75-125			
Lead	98.8	3.0	"	100		98.8	75-125			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041307 - EPA 3050B

Matrix Spike (8041307-MS1)	Source: T181234-10			Prepared & Analyzed: 04/13/18						
Arsenic	95.1	5.0	mg/kg	100	1.26	93.9	75-125			
Barium	137	1.0	"	100	44.3	92.6	75-125			
Cadmium	93.6	2.0	"	100	ND	93.6	75-125			
Chromium	98.9	2.0	"	100	3.82	95.0	75-125			
Lead	95.6	3.0	"	100	1.54	94.1	75-125			

Matrix Spike Dup (8041307-MSD1)	Source: T181234-10			Prepared & Analyzed: 04/13/18						
Arsenic	83.2	5.0	mg/kg	100	1.26	81.9	75-125	13.4	20	
Barium	125	1.0	"	100	44.3	80.8	75-125	9.07	20	
Cadmium	84.9	2.0	"	100	ND	84.9	75-125	9.84	20	
Chromium	89.0	2.0	"	100	3.82	85.2	75-125	10.5	20	
Lead	83.0	3.0	"	100	1.54	81.5	75-125	14.1	20	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041308 - EPA 7471A Soil

Blank (8041308-BLK1)

Prepared & Analyzed: 04/13/18

Mercury	ND	0.10	mg/kg

LCS (8041308-BS1)

Prepared & Analyzed: 04/13/18

Mercury	0.427	0.10	mg/kg	0.417	103	80-120

Matrix Spike (8041308-MS1)

Source: T181234-10

Prepared & Analyzed: 04/13/18

Mercury	0.399	0.10	mg/kg	0.417	ND	95.7	75-125

Matrix Spike Dup (8041308-MSD1)

Source: T181234-10

Prepared & Analyzed: 04/13/18

Mercury	0.399	0.10	mg/kg	0.417	ND	95.9	75-125	0.195	20

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041144 - EPA 3550 ECD/GCMS

Blank (8041144-BLK1)

Prepared: 04/11/18 Analyzed: 04/13/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	9.99		"	10.0		99.9	35-140			
Surrogate: Decachlorobiphenyl	10.9		"	10.0		109	35-140			

LCS (8041144-BS1)

Prepared: 04/11/18 Analyzed: 04/13/18

PCB-1016	85.4	10	ug/kg	100		85.4	40-130			
PCB-1260	66.4	10	"	100		66.4	40-130			
Surrogate: Tetrachloro-meta-xylene	8.79		"	10.0		87.9	35-140			
Surrogate: Decachlorobiphenyl	10.4		"	10.0		104	35-140			

LCS Dup (8041144-BSD1)

Prepared: 04/11/18 Analyzed: 04/13/18

PCB-1016	87.6	10	ug/kg	100		87.6	40-130	2.54	30	
PCB-1260	73.8	10	"	100		73.8	40-130	10.6	30	
Surrogate: Tetrachloro-meta-xylene	8.32		"	10.0		83.2	35-140			
Surrogate: Decachlorobiphenyl	10.5		"	10.0		105	35-140			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041145 - EPA 3550 ECD/GCMS

Blank (8041145-BLK1)

Prepared: 04/11/18 Analyzed: 04/13/18

Carbazole	ND	300	ug/kg
Aniline	ND	300	"
Phenol	ND	1000	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
1-Methylnaphthalene	ND	300	"
2-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041145 - EPA 3550 ECD/GCMS

Blank (8041145-BLK1)

Prepared: 04/11/18 Analyzed: 04/13/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
Pyridine	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
2,4,6-Trichlorophenol	ND	1000	"

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Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041145 - EPA 3550 ECD/GCMS

Blank (8041145-BLK1)

Prepared: 04/11/18 Analyzed: 04/13/18

Surrogate: 2-Fluorophenol	1400		ug/kg	3240		43.4	15-121			
Surrogate: Phenol-d6	1560		"	3240		48.1	24-113			
Surrogate: Nitrobenzene-d5	1750		"	3240		53.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2060		"	3240		63.8	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2450		"	3240		75.8	18.1-105			
Surrogate: Terphenyl-d14	2410		"	3240		74.5	29.1-130			

LCS (8041145-BS1)

Prepared: 04/11/18 Analyzed: 04/13/18

Phenol	1600	1000	ug/kg	3390		47.1	34-114			
2-Chlorophenol	1750	1000	"	3390		51.8	34-114			
1,4-Dichlorobenzene	1860	300	"	3390		55.0	34-114			
N-Nitrosodi-n-propylamine	2140	300	"	3390		63.3	30-110			
1,2,4-Trichlorobenzene	2100	300	"	3390		62.0	39-119			
4-Chloro-3-methylphenol	2340	1000	"	3390		69.1	50-130			
Acenaphthene	1860	300	"	3390		55.0	34-114			
Pentachlorophenol	2320	1000	"	3390		68.4	50-130			
Pyrene	1570	300	"	3390		46.2	30-110			
Surrogate: 2-Fluorophenol	1340		"	3390		39.6	15-121			
Surrogate: Phenol-d6	1550		"	3390		45.6	24-113			
Surrogate: Nitrobenzene-d5	1770		"	3390		52.4	21.3-119			
Surrogate: 2-Fluorobiphenyl	2070		"	3390		61.0	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2440		"	3390		72.0	18.1-105			
Surrogate: Terphenyl-d14	2350		"	3390		69.2	29.1-130			

LCS Dup (8041145-BSD1)

Prepared: 04/11/18 Analyzed: 04/13/18

Phenol	1670	1000	ug/kg	3400		49.1	34-114	4.50	42	
2-Chlorophenol	1830	1000	"	3400		53.8	34-114	4.28	40	
1,4-Dichlorobenzene	1910	300	"	3400		56.3	34-114	2.69	28	
N-Nitrosodi-n-propylamine	2220	300	"	3400		65.2	30-110	3.36	38	
1,2,4-Trichlorobenzene	2220	300	"	3400		65.4	39-119	5.60	28	
4-Chloro-3-methylphenol	2500	1000	"	3400		73.4	50-130	6.42	42	
Acenaphthene	2010	300	"	3400		59.1	34-114	7.55	31	
Pentachlorophenol	2530	1000	"	3400		74.3	50-130	8.67	50	
Pyrene	1780	300	"	3400		52.4	30-110	12.9	31	
Surrogate: 2-Fluorophenol	1410		"	3400		41.4	15-121			
Surrogate: Phenol-d6	1600		"	3400		47.1	24-113			

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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041145 - EPA 3550 ECD/GCMS

LCS Dup (8041145-BSD1)

Prepared: 04/11/18 Analyzed: 04/13/18

Surrogate: Nitrobenzene-d5	1800		ug/kg	3400		52.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2170		"	3400		63.7	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2630		"	3400		77.2	18.1-105			
Surrogate: Terphenyl-d14	2580		"	3400		76.0	29.1-130			

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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/19/18 13:50

Notes and Definitions

DET Analyte DETECTED
ND Analyte NOT DETECTED at or above the reporting limit
NR Not Reported
dry Sample results reported on a dry weight basis
RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



**SunStar
Laboratories**

Chain of Custody Record

25712 Commerce Centre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, Roseville, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/11/18 Page: 1 of 3
Project Name: LATHROP 4E-AC PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 7/87/154 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 PCBs 9002	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD AND ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers	
VS09	4/11/18	11:00	SOIL	JAR															
VS10		11:02																	
VS11		11:03																	
VS12		11:05																	
VS13		11:07																	
VS14		11:09																	
VS15		11:17																	
VS16		11:14																	
VS17		11:20																	
VS18		11:22																	
VS19		11:26																	
VS20		11:29																	
VS21		11:31																	
VS22		11:34																	
VS23		11:35																	
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/11/18 2:58 PM</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4/11/18 2:58</u>					Total # of containers					Notes				
Relinquished by: (signature) _____ Date / Time _____					Received by: (signature) _____ Date / Time _____					Chain of Custody seals X/N/A					Seals intact? Y/N/A				
Relinquished by: (signature) _____ Date / Time _____					Received by: (signature) _____ Date / Time _____					Received good condition/cold					Turn around time: <u>NOE WAK</u>				

COC 170171



**SunStar
Laboratories**

Chain of Custody Record

25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
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Project Manager: JOE BRUSCA

Date: 4/11/18 Page: 2 of 3
Project Name: LATHROP 4E-A PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 718294 EDF #:

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260	8260 + OXY	8260 BTEX, OXY only	8270	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD AND ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers
VS24	4/11/18	11:37	Soil	JAR	X			X				X	X					1
VS25		11:55																1
VS26		11:57																1
VS27		11:58																1
VS28		12:00																1
VS29		12:03																1
VS30		12:04																1
VS31		12:06																1
VS32		12:08																1
VS33		12:10																1
VS34		12:41																1
VS35		12:43																1
VS36		12:45																1
VS37		12:48																1
VS38		12:50																1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/11/18 2:58pm</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/11/18 2:58</u>																		
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 10:00</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 10:00</u>																		
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 10:00</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 10:00</u>																		
Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____																		

COC 170170



SunStar Laboratories

25712 Commerce Drive, Lake Forest, CA 92630
949-297-6020

Chain of Custody Record

REVISED

Client: BRUSCA ASSOCIATES INC.
Address: PO Box 332, Roseville, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/11/18 Page: 3 of 3
Project Name: LATHROP 4E-AC REENTRY
Collector: BRUSCA Client Project #: 137-002
Batch #: 7/10/2014 EDF #:

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 - SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 - PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Lens and Arsenic	Laboratory ID #	Comments/Preservative	Total # of containers
V3 39	4/11/18	12:52	SOIL	JAR	X	X	X	X	X	X	X	X	X	X	X	X	140LD	1
V3 40		12:53															140LD	1
V3 41		12:54															140LD	1
V3 42		12:55															140LD	1
V3 43		1:14 PM															140LD	1
V3 44		1:16 PM															140LD	1
V3 45		1:17 PM															140LD	1
V3 46		1:19 PM															140LD	1
V3 47		1:21 PM															140LD	1
V3 48		1:25 PM															140LD	1
V3 49		1:26 PM															140LD	1
V3 50		1:28 PM															140LD	1
V3 51		1:30 PM															140LD	1

Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time	Total # of containers	Chain of Custody seals	Seals intact?	Received good condition/cold	Turn around time:	Notes
<u>[Signature]</u>	<u>4/11/18 2:58 PM</u>	<u>[Signature]</u>	<u>4/11/18 2:58</u>					<u>NO</u>	
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time						
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time						

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

COC 170160



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

23 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/14/18 08:10. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
VS53	T181266-02	Soil	04/13/18 10:44	04/14/18 08:10
VS55	T181266-04	Soil	04/13/18 10:49	04/14/18 08:10
VS57	T181266-06	Soil	04/13/18 10:54	04/14/18 08:10
VS59	T181266-08	Soil	04/13/18 10:57	04/14/18 08:10
VS61	T181266-10	Soil	04/13/18 11:14	04/14/18 08:10
VS63	T181266-12	Soil	04/13/18 11:19	04/14/18 08:10
VS65	T181266-14	Soil	04/13/18 11:23	04/14/18 08:10
VS66	T181266-15	Soil	04/13/18 11:27	04/14/18 08:10
VS68	T181266-17	Soil	04/13/18 11:32	04/14/18 08:10
VS70	T181266-19	Soil	04/13/18 11:35	04/14/18 08:10

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

DETECTIONS SUMMARY

Sample ID: VS53

Laboratory ID: T181266-02

Reporting

Analyte	Result	Limit	Units	Method	Notes
Barium	80	1.0	mg/kg	EPA 6010B	
Chromium	7.1	2.0	mg/kg	EPA 6010B	
Cobalt	6.6	2.0	mg/kg	EPA 6010B	
Copper	6.8	1.0	mg/kg	EPA 6010B	
Nickel	8.8	2.0	mg/kg	EPA 6010B	
Vanadium	23	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	

Sample ID: VS55

Laboratory ID: T181266-04

Reporting

Analyte	Result	Limit	Units	Method	Notes
C13-C28 (DRO)	21	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	26	10	mg/kg	EPA 8015B	
Barium	65	1.0	mg/kg	EPA 6010B	
Chromium	7.7	2.0	mg/kg	EPA 6010B	
Cobalt	5.8	2.0	mg/kg	EPA 6010B	
Copper	4.1	1.0	mg/kg	EPA 6010B	
Nickel	8.2	2.0	mg/kg	EPA 6010B	
Vanadium	24	5.0	mg/kg	EPA 6010B	
Zinc	27	1.0	mg/kg	EPA 6010B	

Sample ID: VS57

Laboratory ID: T181266-06

Reporting

Analyte	Result	Limit	Units	Method	Notes
Barium	49	1.0	mg/kg	EPA 6010B	
Chromium	6.6	2.0	mg/kg	EPA 6010B	
Cobalt	4.7	2.0	mg/kg	EPA 6010B	
Copper	3.0	1.0	mg/kg	EPA 6010B	
Nickel	5.8	2.0	mg/kg	EPA 6010B	
Vanadium	21	5.0	mg/kg	EPA 6010B	
Zinc	32	1.0	mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Sample ID: VS59

Laboratory ID: T181266-08

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	62	1.0	mg/kg	EPA 6010B	
Chromium	6.0	2.0	mg/kg	EPA 6010B	
Cobalt	5.0	2.0	mg/kg	EPA 6010B	
Copper	3.5	1.0	mg/kg	EPA 6010B	
Nickel	7.2	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	22	1.0	mg/kg	EPA 6010B	

Sample ID: VS61

Laboratory ID: T181266-10

No Results Detected

Sample ID: VS63

Laboratory ID: T181266-12

No Results Detected

Sample ID: VS65

Laboratory ID: T181266-14

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Barium	26	1.0	mg/kg	EPA 6010B	
Chromium	3.7	2.0	mg/kg	EPA 6010B	
Cobalt	2.7	2.0	mg/kg	EPA 6010B	
Copper	1.9	1.0	mg/kg	EPA 6010B	
Nickel	3.4	2.0	mg/kg	EPA 6010B	
Vanadium	11	5.0	mg/kg	EPA 6010B	
Zinc	11	1.0	mg/kg	EPA 6010B	

Sample ID: VS66

Laboratory ID: T181266-15

No Results Detected

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Sample ID: VS68

Laboratory ID: T181266-17

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Lead	12	3.0		mg/kg	EPA 6010b	

Sample ID: VS70

Laboratory ID: T181266-19

No Results Detected

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS53

T181266-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	80	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	7.1	2.0	"	"	"	"	"	"	
Cobalt	6.6	2.0	"	"	"	"	"	"	
Copper	6.8	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	8.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	23	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS53

T181266-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		91.3 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		93.8 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS53

T181266-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS53

T181266-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	62.3 %	15-121			"	"	"	"
Surrogate: Phenol-d6	71.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	72.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	82.5 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	99.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	84.7 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS55

T181266-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	21	10	"	"	"	"	"	"	
C29-C40 (MORO)	26	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		104 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	65	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	7.7	2.0	"	"	"	"	"	"	
Cobalt	5.8	2.0	"	"	"	"	"	"	
Copper	4.1	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	8.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	24	5.0	"	"	"	"	"	"	
Zinc	27	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS55

T181266-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		89.9 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		93.0 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS55

T181266-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS55

T181266-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	51.4 %	15-121			"	"	"	"
Surrogate: Phenol-d6	62.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	64.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	76.7 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	91.4 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	85.7 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS57

T181266-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	49	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.6	2.0	"	"	"	"	"	"	
Cobalt	4.7	2.0	"	"	"	"	"	"	
Copper	3.0	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	21	5.0	"	"	"	"	"	"	
Zinc	32	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS57

T181266-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		96.3 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		101 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS57

T181266-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS57

T181266-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	62.5 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	62.2 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	72.2 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	86.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	83.8 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS59

T181266-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: p-Terphenyl		99.9 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	62	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.0	2.0	"	"	"	"	"	"	
Cobalt	5.0	2.0	"	"	"	"	"	"	
Copper	3.5	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	22	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS59

T181266-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		92.7 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		109 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS59

T181266-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS59

T181266-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	61.1 %	15-121			"	"	"	"
Surrogate: Phenol-d6	70.9 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	69.8 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	79.6 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	93.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	90.0 %	29.1-130			"	"	"	"

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04/23/18 10:43

VS61
T181266-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS63

T181266-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS65
T181266-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	ND	10	"	"	"	"	"	"	
C29-C40 (MORO)	ND	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		100 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	26	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	3.7	2.0	"	"	"	"	"	"	
Cobalt	2.7	2.0	"	"	"	"	"	"	
Copper	1.9	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	11	5.0	"	"	"	"	"	"	
Zinc	11	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

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04/23/18 10:43

VS65

T181266-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		98.5 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		113 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS65

T181266-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS65
T181266-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	56.7 %	15-121			"	"	"	"
Surrogate: Phenol-d6	68.1 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	67.2 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	81.1 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	97.1 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	92.1 %	29.1-130			"	"	"	"

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Reported:
04/23/18 10:43

VS66

T181266-15 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Manager: Joe Brusca

Reported:
04/23/18 10:43

VS68

T181266-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010b	
Lead	12	3.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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04/23/18 10:43

VS70

T181266-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6010B

Arsenic	ND	5.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010b	
Lead	ND	3.0	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041214 - EPA 3550B GC

Blank (8041214-BLK1)

Prepared: 04/12/18 Analyzed: 04/18/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	97.8		"	99.0		98.8	65-135			

LCS (8041214-BS1)

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	500	10	mg/kg	495		101	75-125			
Surrogate: <i>p</i> -Terphenyl	101		"	99.0		102	65-135			

Matrix Spike (8041214-MS1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	560	10	mg/kg	505	53	100	75-125			
Surrogate: <i>p</i> -Terphenyl	106		"	101		105	65-135			

Matrix Spike Dup (8041214-MSD1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	590	10	mg/kg	500	53	108	75-125	5.92	20	
Surrogate: <i>p</i> -Terphenyl	104		"	100		104	65-135			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041623 - EPA 3050B

Blank (8041623-BLK1)

Prepared & Analyzed: 04/16/18

Antimony	ND	3.0	mg/kg
Antimony	ND	3.0	"
Silver	ND	2.0	"
Arsenic	ND	5.0	"
Arsenic	ND	5.0	"
Barium	ND	1.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	2.0	"
Cadmium	ND	2.0	"
Chromium	ND	2.0	"
Cobalt	ND	2.0	"
Chromium	ND	2.0	"
Copper	ND	1.0	"
Cobalt	ND	2.0	"
Lead	ND	3.0	"
Copper	ND	1.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Lead	ND	3.0	"
Selenium	ND	5.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	ND	1.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Selenium	ND	5.0	"
Silver	ND	2.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	ND	1.0	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041623 - EPA 3050B

LCS (8041623-BS1)

Prepared & Analyzed: 04/16/18

Arsenic	92.9	5.0	mg/kg	100		92.9	75-125			
Arsenic	92.9	5.0	"	100		92.9	75-125			
Barium	91.6	1.0	"	100		91.6	75-125			
Barium	91.6	1.0	"	100		91.6	75-125			
Cadmium	88.7	2.0	"	100		88.7	75-125			
Cadmium	88.7	2.0	"	100		88.7	75-125			
Chromium	90.3	2.0	"	100		90.3	75-125			
Chromium	90.3	2.0	"	100		90.3	75-125			
Lead	96.9	3.0	"	100		96.9	75-125			
Lead	96.9	3.0	"	100		96.9	75-125			

Matrix Spike (8041623-MS1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Arsenic	88.5	5.0	mg/kg	100	1.81	86.7	75-125			
Arsenic	88.5	5.0	"	100	1.81	86.7	75-125			
Barium	119	1.0	"	100	37.7	81.1	75-125			
Barium	119	1.0	"	100	37.7	81.1	75-125			
Cadmium	83.2	2.0	"	100	0.425	82.8	75-125			
Cadmium	83.2	2.0	"	100	0.425	82.8	75-125			
Chromium	89.3	2.0	"	100	5.37	84.0	75-125			
Chromium	89.3	2.0	"	100	5.37	84.0	75-125			
Lead	93.3	3.0	"	100	ND	93.3	75-125			
Lead	93.3	3.0	"	100	6.15	87.1	75-125			

Matrix Spike Dup (8041623-MSD1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Arsenic	95.2	5.0	mg/kg	100	1.81	93.4	75-125	7.37	20	
Arsenic	95.2	5.0	"	100	1.81	93.4	75-125	7.37	20	
Barium	123	1.0	"	100	37.7	85.1	75-125	3.25	20	
Barium	123	1.0	"	100	37.7	85.1	75-125	3.25	20	
Cadmium	85.4	2.0	"	100	0.425	85.0	75-125	2.67	20	
Cadmium	85.4	2.0	"	100	0.425	85.0	75-125	2.67	20	
Chromium	91.3	2.0	"	100	5.37	86.0	75-125	2.21	20	
Chromium	91.3	2.0	"	100	5.37	86.0	75-125	2.21	20	
Lead	99.2	3.0	"	100	ND	99.2	75-125	6.14	20	
Lead	99.2	3.0	"	100	6.15	93.0	75-125	6.14	20	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041624 - EPA 7471A Soil

Blank (8041624-BLK1)

Prepared & Analyzed: 04/16/18

Mercury	ND	0.10	mg/kg

LCS (8041624-BS1)

Prepared & Analyzed: 04/16/18

Mercury	0.405	0.10	mg/kg	0.417		97.2	80-120

Matrix Spike (8041624-MS1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Mercury	0.408	0.10	mg/kg	0.417	ND	98.0	75-125

Matrix Spike Dup (8041624-MSD1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Mercury	0.428	0.10	mg/kg	0.417	ND	103	75-125	4.64	20

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041633 - EPA 3550 ECD/GCMS

Blank (8041633-BLK1)

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	8.65		"	9.90		87.3	35-140			
Surrogate: Decachlorobiphenyl	10.1		"	9.90		102	35-140			

LCS (8041633-BS1)

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	70.1	10	ug/kg	99.0		70.8	40-130			
PCB-1260	66.8	10	"	99.0		67.5	40-130			
Surrogate: Tetrachloro-meta-xylene	9.51		"	9.90		96.0	35-140			
Surrogate: Decachlorobiphenyl	11.0		"	9.90		112	35-140			

Matrix Spike (8041633-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	60.2	10	ug/kg	101	ND	59.6	40-130			
PCB-1260	60.1	10	"	101	ND	59.5	40-130			
Surrogate: Tetrachloro-meta-xylene	8.99		"	10.1		89.0	35-140			
Surrogate: Decachlorobiphenyl	9.25		"	10.1		91.5	35-140			

Matrix Spike Dup (8041633-MSD1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	64.2	10	ug/kg	101	ND	63.6	40-130	6.49	30	
PCB-1260	62.7	10	"	101	ND	62.1	40-130	4.28	30	
Surrogate: Tetrachloro-meta-xylene	9.14		"	10.1		90.5	35-140			
Surrogate: Decachlorobiphenyl	9.02		"	10.1		89.3	35-140			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

Carbazole	ND	300	ug/kg
Aniline	ND	300	"
Phenol	ND	1000	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
2-Methylnaphthalene	ND	300	"
1-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
Pyridine	ND	300	"
2,4,6-Trichlorophenol	ND	1000	"

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Reported:
04/23/18 10:43

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

Surrogate: 2-Fluorophenol	1790		ug/kg	3380		53.0	15-121			
Surrogate: Phenol-d6	1920		"	3380		56.9	24-113			
Surrogate: Nitrobenzene-d5	2150		"	3380		63.5	21.3-119			
Surrogate: 2-Fluorobiphenyl	2510		"	3380		74.3	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3160		"	3380		93.4	18.1-105			
Surrogate: Terphenyl-d14	2960		"	3380		87.6	29.1-130			

LCS (8041634-BS1)

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1490	1000	ug/kg	3280		45.6	34-114			
2-Chlorophenol	1620	1000	"	3280		49.4	34-114			
1,4-Dichlorobenzene	1680	300	"	3280		51.4	34-114			
N-Nitrosodi-n-propylamine	1700	300	"	3280		51.9	30-110			
1,2,4-Trichlorobenzene	1900	300	"	3280		58.0	39-119			
4-Chloro-3-methylphenol	2150	1000	"	3280		65.6	50-130			
Acenaphthene	1720	300	"	3280		52.6	34-114			
Pentachlorophenol	2280	1000	"	3280		69.7	50-130			
Pyrene	1530	300	"	3280		46.8	30-110			
Surrogate: 2-Fluorophenol	1450		"	3280		44.3	15-121			
Surrogate: Phenol-d6	1650		"	3280		50.4	24-113			
Surrogate: Nitrobenzene-d5	1790		"	3280		54.7	21.3-119			
Surrogate: 2-Fluorobiphenyl	2130		"	3280		65.1	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2810		"	3280		85.6	18.1-105			
Surrogate: Terphenyl-d14	2600		"	3280		79.4	29.1-130			

Matrix Spike (8041634-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1750	1000	ug/kg	3390	ND	51.7	34-114			
2-Chlorophenol	1860	1000	"	3390	ND	54.9	34-114			
1,4-Dichlorobenzene	1830	300	"	3390	ND	54.0	34-114			
N-Nitrosodi-n-propylamine	1880	300	"	3390	ND	55.4	30-110			
1,2,4-Trichlorobenzene	2190	300	"	3390	ND	64.5	39-119			
4-Chloro-3-methylphenol	2550	1000	"	3390	ND	75.3	50-130			
Acenaphthene	1960	300	"	3390	ND	57.8	34-114			
Pentachlorophenol	2620	1000	"	3390	ND	77.3	50-130			
Pyrene	1980	300	"	3390	160	53.7	30-110			
Surrogate: 2-Fluorophenol	1600		"	3390		47.3	15-121			
Surrogate: Phenol-d6	1890		"	3390		55.8	24-113			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Matrix Spike (8041634-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Surrogate: Nitrobenzene-d5	2040		ug/kg	3390		60.1	21.3-119			
Surrogate: 2-Fluorobiphenyl	2310		"	3390		68.1	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3090		"	3390		91.2	18.1-105			
Surrogate: Terphenyl-d14	2670		"	3390		78.7	29.1-130			

Matrix Spike Dup (8041634-MSD1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1720	1000	ug/kg	3410	ND	50.4	34-114	2.00	42	
2-Chlorophenol	1840	1000	"	3410	ND	53.9	34-114	1.32	40	
1,4-Dichlorobenzene	1760	300	"	3410	ND	51.7	34-114	3.82	28	
N-Nitrosodi-n-propylamine	1840	300	"	3410	ND	54.0	30-110	1.82	38	
1,2,4-Trichlorobenzene	2090	300	"	3410	ND	61.4	39-119	4.29	28	
4-Chloro-3-methylphenol	2560	1000	"	3410	ND	74.9	50-130	0.108	42	
Acenaphthene	1960	300	"	3410	ND	57.3	34-114	0.119	31	
Pentachlorophenol	2860	1000	"	3410	ND	83.9	50-130	8.97	50	
Pyrene	1940	300	"	3410	160	52.1	30-110	2.05	31	
Surrogate: 2-Fluorophenol	1630		"	3410		47.8	15-121			
Surrogate: Phenol-d6	1870		"	3410		54.8	24-113			
Surrogate: Nitrobenzene-d5	1870		"	3410		54.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2210		"	3410		64.7	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3080		"	3410		90.1	18.1-105			
Surrogate: Terphenyl-d14	2940		"	3410		86.0	29.1-130			

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:43

Notes and Definitions

DET Analyte DETECTED
ND Analyte NOT DETECTED at or above the reporting limit
NR Not Reported
dry Sample results reported on a dry weight basis
RPD Relative Percent Difference

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Mike Jaroudi, Project Manager



Sunstar Laboratories

Chain of Custody Record

25712 Commerce Drive, Lake Forest, CA 92630
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Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332 ROSELAND, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/13/18 Page: 1 of 2
Project Name: LATHROP 40-Ac PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 718266 EDF #:

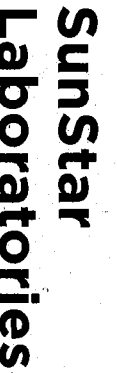
Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 - SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 - PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD + ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers	
V552	4/13/18	10:43	SOL	502	X			X				X	X				01	HOLD	1
V553		10:44						X				X	X				02	HOLD	1
V554		10:44						X				X	X				03	HOLD	1
V555		10:44						X				X	X				04	HOLD	1
V556		10:51						X				X	X				05	HOLD	1
V557		10:54						X				X	X				06	HOLD	1
V558		10:55						X				X	X				07	HOLD	1
V559		10:57						X				X	X				08	HOLD	1
V560		10:59						X				X	X				09	HOLD	1
V561		11:14															10	HOLD	1
V562		11:16															11	HOLD	1
V563		11:19															12	HOLD	1
V564		11:21															13	HOLD	1
V565		11:23															14	HOLD	1
V566		11:27															15	HOLD	1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 1:30PM</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 1:30</u>					Total # of containers <u>15</u>					Notes				
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 8:10</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 8:10</u>					Chain of Custody seals Y/N/A <u>Seals intact? Y/N/A</u>					Received good condition/cold <u>5-5</u>				
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 8:10</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4/13/18 8:10</u>					Turn around time: <u>Normal</u>									

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

COC 170174



25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSA ASSOCIATES, INC.

Date: 4/13/18

Page: 2 of 2

Address: PO Box 332, Roseville, CA 95661
Phone: (916) 677-470 Fax: (916) 677-471

Project Name: LATHROP 48-A PROPERTY
Collector: BRUCA Client Project #: 137-002

Project Manager: Joe Brusca

Batch #: 7181266 EDF #: 1

[illegible]

Sample disposal Instructions: Disposal @ \$2.00 each _____

Return to client _____

Pickup _____

GOC 170163



April 17, 2018

FAL Project ID: 11414

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11414**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Four solid samples were received on 4/2/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of ten business days for project **11414**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and a sample photo. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11414**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink that reads "Thomas C. Crabtree".

Thomas C. Crabtree
Director

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: **11414**

Received on: **04/02/2018**

Project Due: **04/17/2018** Storage: **R-4**

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11414-001-SA	0	137-002	VS1	EPA 1613 D/F	Soil	04/02/2018	11:45 am	04/02/2019
11414-002-SA	0	137-002	VS3	EPA 1613 D/F	Soil	04/02/2018	11:52 am	04/02/2019
11414-003-SA	0	137-002	VS5	EPA 1613 D/F	Soil	04/02/2018	11:57 am	04/02/2019
11414-004-SA	0	137-002	VS7	EPA 1613 D/F	Soil	04/02/2018	12:03 pm	04/02/2019

EPA Method 1613 PCDD/F



FAL ID: 11414-001-MB
Client ID: Method Blank
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-16-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.111		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.210		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.390		-	0.0503				
1,2,3,6,7,8-HxCDD	ND	0.391		-	0.0490	Total TCDD	ND	0.111	
1,2,3,7,8,9-HxCDD	ND	0.378		-	0.0488	Total PeCDD	ND	0.210	
1,2,3,4,6,7,8-HpCDD	ND	0.379		-	0.0541	Total HxCDD	ND	0.391	
OCDD	ND	0.662		-	0.0888	Total HpCDD	ND	0.379	
2,3,7,8-TCDF	ND	0.0839		-	0.0243				
1,2,3,7,8-PeCDF	ND	0.107		-	0.0285				
2,3,4,7,8-PeCDF	ND	0.105		-	0.0298				
1,2,3,4,7,8-HxCDF	ND	0.172		-	0.0255				
1,2,3,6,7,8-HxCDF	ND	0.176		-	0.0253				
2,3,4,6,7,8-HxCDF	ND	0.194		-	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.231		-	0.0367	Total TCDF	ND	0.0839	
1,2,3,4,6,7,8-HpCDF	ND	0.178		-	0.0321	Total PeCDF	ND	0.107	
1,2,3,4,7,8,9-HpCDF	ND	0.248		-	0.0396	Total HxCDF	ND	0.231	
OCDF	ND	0.277		-	0.0843	Total HpCDF	ND	0.248	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	96.3	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	89.5	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.1	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	97.8	23.0 - 140	
13C-OCDD	81.9	17.0 - 157	
13C-2,3,7,8-TCDF	100	24.0 - 169	
13C-1,2,3,7,8-PeCDF	79.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	82.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	106	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	109	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	106	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	108	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	92.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	98.1	26.0 - 138	
13C-OCDF	86.3	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 92.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/17/2018

Reviewed By: 

Date: 4/17/2018

000003 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11414-001-OPR
Client ID: OPR
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-16-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	11.2	6.70 - 15.8	
1,2,3,7,8-PeCDD	51.1	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	57.6	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	57.9	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	57.8	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	55.3	35.0 - 70.0	
OCDD	111	78.0 - 144	
2,3,7,8-TCDF	10.1	7.50 - 15.8	
1,2,3,7,8-PeCDF	47.1	40.0 - 67.0	
2,3,4,7,8-PeCDF	47.5	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	48.6	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	49.1	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	49.9	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	50.8	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	51.0	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	51.1	39.0 - 69.0	
OCDF	102	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	99.5	20.0 - 175	
13C-1,2,3,7,8-PeCDD	87.1	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	95.5	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	103	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	101	26.0 - 166	
13C-OCDD	87.7	13.0 - 198	
13C-2,3,7,8-TCDF	103	22.0 - 152	
13C-1,2,3,7,8-PeCDF	78.2	21.0 - 192	
13C-2,3,4,7,8-PeCDF	81.0	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	110	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	112	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	107	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	106	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	92.3	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	104	20.0 - 186	
13C-OCDF	87.3	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	96.0	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/17/2018

Reviewed By: 
Date: 4/17/2018

000004 of 000011

EPA Method 1613
PCDD/F



FAL ID: 11414-001-SA
Client ID: VS1
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: 04-02-2018
Amount: 5.03 g
% Solids: 96.30

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-17-2018
2005 WHO TEQ: 0.00672
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.149		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.322		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.388		-	0.0503				
1,2,3,6,7,8-HxCDD	ND	0.389		-	0.0490	Total TCDD	ND	0.149	
1,2,3,7,8,9-HxCDD	ND	0.376		-	0.0488	Total PeCDD	ND	0.322	
1,2,3,4,6,7,8-HpCDD	0.603	-	J	0.00603	0.0541	Total HxCDD	ND	0.389	
OCDD	2.29	-	J	0.000687	0.0888	Total HpCDD	1.16	-	J
2,3,7,8-TCDF	ND	0.112		-	0.0243				
1,2,3,7,8-PeCDF	ND	0.192		-	0.0285				
2,3,4,7,8-PeCDF	ND	0.184		-	0.0298				
1,2,3,4,7,8-HxCDF	ND	0.183		-	0.0255				
1,2,3,6,7,8-HxCDF	ND	0.183		-	0.0253				
2,3,4,6,7,8-HxCDF	ND	0.202		-	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.258		-	0.0367	Total TCDF	ND	0.112	
1,2,3,4,6,7,8-HpCDF	ND	0.251		-	0.0321	Total PeCDF	ND	0.192	
1,2,3,4,7,8,9-HpCDF	ND	0.325		-	0.0396	Total HxCDF	ND	0.258	
OCDF	ND	0.368		-	0.0843	Total HpCDF	ND	0.325	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	101	25.0 - 164	
13C-1,2,3,7,8-PeCDD	87.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	87.7	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	92.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	96.3	23.0 - 140	
13C-OCDD	79.9	17.0 - 157	
13C-2,3,7,8-TCDF	107	24.0 - 169	
13C-1,2,3,7,8-PeCDF	79.4	24.0 - 185	
13C-2,3,4,7,8-PeCDF	82.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	101	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	107	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	103	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	101	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	92.4	26.0 - 138	
13C-OCDF	81.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 98.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst:

Date: 4/17/2018

Reviewed By:

Date: 4/17/2018

000005 of 000011

EPA Method 1613
PCDD/F



FAL ID: 11414-002-SA
Client ID: VS3
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: 04-02-2018
Amount: 5.07 g
% Solids: 96.41

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-17-2018
2005 WHO TEQ: 0.00529
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.0914		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.201		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.343		-	0.0503				
1,2,3,6,7,8-HxCDD	ND	0.371		-	0.0490	Total TCDD	ND	0.0914	
1,2,3,7,8,9-HxCDD	ND	0.347		-	0.0488	Total PeCDD	ND	0.201	
1,2,3,4,6,7,8-HpCDD	0.460	-	J	0.00460	0.0541	Total HxCDD	ND	0.371	
OCDD	2.30	-	J	0.000690	0.0888	Total HpCDD	1.12	-	J
2,3,7,8-TCDF	ND	0.104		-	0.0243				
1,2,3,7,8-PeCDF	ND	0.150		-	0.0285				
2,3,4,7,8-PeCDF	ND	0.144		-	0.0298				
1,2,3,4,7,8-HxCDF	ND	0.103		-	0.0255				
1,2,3,6,7,8-HxCDF	ND	0.101		-	0.0253				
2,3,4,6,7,8-HxCDF	ND	0.115		-	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.142		-	0.0367	Total TCDF	ND	0.104	
1,2,3,4,6,7,8-HpCDF	ND	0.159		-	0.0321	Total PeCDF	ND	0.150	
1,2,3,4,7,8,9-HpCDF	ND	0.201		-	0.0396	Total HxCDF	ND	0.142	
OCDF	ND	0.355		-	0.0843	Total HpCDF	ND	0.201	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	95.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	83.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	84.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	89.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	86.9	23.0 - 140	
13C-OCDD	64.6	17.0 - 157	
13C-2,3,7,8-TCDF	102	24.0 - 169	
13C-1,2,3,7,8-PeCDF	74.9	24.0 - 185	
13C-2,3,4,7,8-PeCDF	78.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	99.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	103	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	95.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	94.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	83.0	26.0 - 138	
13C-OCDF	68.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 97.1 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/17/2018

Reviewed By: 

Date: 4/17/2018

000006 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11414-003-SA
Client ID: VS5
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: 04-02-2018
Amount: 5.00 g
% Solids: 95.28

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-17-2018
2005 WHO TEQ: 0.00850
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.0985		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.236		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.372		-	0.0503				
1,2,3,6,7,8-HxCDD	ND	0.359		-	0.0490	Total TCDD	ND	0.0985	
1,2,3,7,8,9-HxCDD	ND	0.353		-	0.0488	Total PeCDD	ND	0.236	
1,2,3,4,6,7,8-HpCDD	0.703	-	J	0.00703	0.0541	Total HxCDD	ND	0.372	
OCDD	4.89	-	J	0.00147	0.0888	Total HpCDD	1.46	-	J
2,3,7,8-TCDF	ND	0.0911		-	0.0243				
1,2,3,7,8-PeCDF	ND	0.185		-	0.0285				
2,3,4,7,8-PeCDF	ND	0.173		-	0.0298				
1,2,3,4,7,8-HxCDF	ND	0.210		-	0.0255				
1,2,3,6,7,8-HxCDF	ND	0.212		-	0.0253				
2,3,4,6,7,8-HxCDF	ND	0.230		-	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.282		-	0.0367	Total TCDF	ND	0.0911	
1,2,3,4,6,7,8-HpCDF	ND	0.186		-	0.0321	Total PeCDF	ND	0.185	
1,2,3,4,7,8,9-HpCDF	ND	0.247		-	0.0396	Total HxCDF	ND	0.282	
OCDF	ND	0.328		-	0.0843	Total HpCDF	ND	0.247	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	105	25.0 - 164	
13C-1,2,3,7,8-PeCDD	93.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	87.6	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	95.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	105	23.0 - 140	
13C-OCDD	87.5	17.0 - 157	
13C-2,3,7,8-TCDF	104	24.0 - 169	
13C-1,2,3,7,8-PeCDF	82.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	85.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	110	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	111	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	108	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	106	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	92.5	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	98.1	26.0 - 138	
13C-OCDF	91.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 97.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/17/2018

Reviewed By: 

Date: 4/17/2018

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EPA Method 1613 PCDD/F



FAL ID: 11414-004-SA
Client ID: VS7
Matrix: Soil
Batch No: X4472

Date Extracted: 04-13-2018
Date Received: 04-02-2018
Amount: 5.01 g
% Solids: 96.52

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-17-2018
2005 WHO TEQ: 0.665
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.138		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.331		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.290		-	0.0503				
1,2,3,6,7,8-HxCDD	0.714	-	J	0.0714	0.0490	Total TCDD	ND	0.138	
1,2,3,7,8,9-HxCDD	0.526	-	J	0.0526	0.0488	Total PeCDD	ND	0.331	
1,2,3,4,6,7,8-HpCDD	10.4	-		0.104	0.0541	Total HxCDD	5.34	-	
OCDD	67.9	-		0.0204	0.0888	Total HpCDD	19.1	-	
2,3,7,8-TCDF	0.523	-	J	0.0523	0.0243				
1,2,3,7,8-PeCDF	ND	0.204		-	0.0285				
2,3,4,7,8-PeCDF	0.637	-	J	0.191	0.0298				
1,2,3,4,7,8-HxCDF	0.571	-	J	0.0571	0.0255				
1,2,3,6,7,8-HxCDF	0.407	-	J	0.0407	0.0253				
2,3,4,6,7,8-HxCDF	0.366	-	J	0.0366	0.0279				
1,2,3,7,8,9-HpCDF	ND	0.218		-	0.0367	Total TCDF	3.35	-	
1,2,3,4,6,7,8-HpCDF	3.64	-	J	0.0364	0.0321	Total PeCDF	7.46	-	
1,2,3,4,7,8,9-HpCDF	ND	0.377		-	0.0396	Total HxCDF	6.54	-	
OCDF	7.64	-	J	0.00229	0.0843	Total HpCDF	8.50	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	104	25.0 - 164	
13C-1,2,3,7,8-PeCDD	89.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	85.1	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.7	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	101	23.0 - 140	
13C-OCDD	86.2	17.0 - 157	
13C-2,3,7,8-TCDF	104	24.0 - 169	
13C-1,2,3,7,8-PeCDF	80.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	82.7	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	103	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	107	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	105	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	105	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	86.9	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	97.9	26.0 - 138	
13C-OCDF	85.4	17.0 - 157	

Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	98.5	35.0 - 197	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/17/2018

Reviewed By: 

Date: 4/17/2018

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Frontier Analytical Laboratory
5172 Hillsdale Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
Fax: 916-934-0999

FAL USE ONLY

Laboratory Project No.: 11414
Temperature: 0 °C

Chain of Custody

www.frontieranalytical.com

Please Print in Pen Page 1 of 1

CLIENT INFORMATION				INVOICE INFORMATION (if different from client info)				PROJECT INFORMATION											
Company Name: Brusca Associates, Inc. Contact Name: Joe Brusca Address: 7633 Stonewood Court, Granite Bay, CA 95746 Phone: 916-677-1470 Fax: _____ Email: jbrusca@bruscaassociates.com				Company Name: Same Contact Name: _____ Address: _____ Phone: _____ Fax: _____ Email: _____				FAL Quote #: 3310 P.O. #: _____ Project #: <u>137-002</u> Project Name: <u>LATHROP 48-AC PROPERTY</u> TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (✓ one) * FAL must agree with price and RUSH TAT in writing.											
REPORT INFORMATION				REPORT DISTRIBUTION (email only is preferred)				ADDITIONAL INSTRUCTIONS											
Report Level: ☐ I/II ☐ III ☐ IV ☐ EDD: ☐ FAL Basic ☐ Geotracker ☐ Other: _____ ☐ Custom: Contact FAL ☐ California State Drinking Water Form System #: _____ Source #: _____ Sampler: _____ Employer: _____				☐ Hardcopy ☐ CD (.pdf including EDDs if requested) ☒ Email (.pdf including EDDs if requested)															
				# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/9A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS	**TEQ			
Sample ID				Date	Time	Matrix										☐ 2,3,7,8-TCDD only ☐ 2,3,7,8-TCDD/F only ☐ PCDD/F (Cl ₄ -Cl ₈)		☐ 1998 WHO ☒ 2005 WHO ☐ Other	
Collected				Remarks															
1	V51	4/2/18	11:45	SOIL		☒													
2	V53		11:52			☒													
3	V55		11:57			☒													
4	V57		12:03			☒													
5																			
6																			
7																			
8																			
9																			
10																			
11																			
12																			
13																			
14																			
15																			
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																			
Relinquished by: (Signature and Printed Name)				Date	Time	Received by: (Signature and Printed Name)				Date	Time								
				4/2/18	15:25					4/2/2018	1525								

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

White Copy - Report

Yellow Copy - Laboratory

Pink Copy - Originator

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11414**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/02/2018
Time Received:	03:25 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	4
Duplicates:	0
Storage Location:	R-4

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Blue Ice/Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



April 30, 2018

FAL Project ID: 11441

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11441**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Eight solid samples were received on 4/11/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of ten business days for project **11441**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and the sample photos. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11441**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink, appearing to read "Bradley B. Silverbush".

Bradley B. Silverbush
Director of Operations

FRONTIER ANALYTICAL LABORATORY

5172 Hillsdale Circle * El Dorado Hills, CA 95762
Tel (916) 934-0900 * Fax (916) 934-0999
www.frontieranalytical.com

000001 of 000016

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: 11441

Received on: 04/11/2018

Project Due: 04/26/2018

Storage: R-4

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11441-001-SA	0	137-002	VS18	EPA 1613 D/F	Soil	04/11/2018	11:22 am	04/11/2019
11441-002-SA	0	137-002	VS20	EPA 1613 D/F	Soil	04/11/2018	11:29 am	04/11/2019
11441-003-SA	0	137-002	VS22	EPA 1613 D/F	Soil	04/11/2018	11:34 am	04/11/2019
11441-004-SA	0	137-002	VS24	EPA 1613 D/F	Soil	04/11/2018	11:37 am	04/11/2019
11441-005-SA	0	137-002	VS35	EPA 1613 D/F	Soil	04/11/2018	12:43 pm	04/11/2019
11441-006-SA	0	137-002	VS37	EPA 1613 D/F	Soil	04/11/2018	12:48 pm	04/11/2019
11441-007-SA	0	137-002	VS39	EPA 1613 D/F	Soil	04/11/2018	12:52 pm	04/11/2019
11441-008-SA	0	137-002	VS41	EPA 1613 D/F	Soil	04/11/2018	12:57 pm	04/11/2019

EPA Method 1613 PCDD/F



FAL ID: 11441-001-MB
Client ID: Method Blank
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-27-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.226		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.321		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.379		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.367		-	0.0940	Total TCDD	ND	0.226	
1,2,3,7,8,9-HxCDD	ND	0.342		-	0.0823	Total PeCDD	ND	0.321	
1,2,3,4,6,7,8-HpCDD	ND	0.421		-	0.0842	Total HxCDD	ND	0.379	
OCDD	ND	1.22		-	0.172	Total HpCDD	ND	0.421	
2,3,7,8-TCDF	ND	0.187		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.222		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.239		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.186		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.196		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.200		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.228		-	0.0657	Total TCDF	ND	0.187	
1,2,3,4,6,7,8-HpCDF	ND	0.229		-	0.0747	Total PeCDF	ND	0.239	
1,2,3,4,7,8,9-HpCDF	ND	0.256		-	0.0883	Total HxCDF	ND	0.228	
OCDF	ND	0.545		-	0.170	Total HpCDF	ND	0.256	
Internal Standards	% Rec	QC Limits	Qual						
13C-2,3,7,8-TCDD	66.1	25.0 - 164							
13C-1,2,3,7,8-PeCDD	68.9	25.0 - 181							
13C-1,2,3,4,7,8-HxCDD	67.5	32.0 - 141							
13C-1,2,3,6,7,8-HxCDD	67.2	28.0 - 130							
13C-1,2,3,4,6,7,8-HpCDD	64.5	23.0 - 140							
13C-OCDD	56.3	17.0 - 157							
13C-2,3,7,8-TCDF	67.6	24.0 - 169							
13C-1,2,3,7,8-PeCDF	66.6	24.0 - 185							
13C-2,3,4,7,8-PeCDF	66.2	21.0 - 178							
13C-1,2,3,4,7,8-HxCDF	64.8	26.0 - 152							
13C-1,2,3,6,7,8-HxCDF	63.2	26.0 - 123							
13C-2,3,4,6,7,8-HxCDF	62.6	28.0 - 136							
13C-1,2,3,7,8,9-HxCDF	62.4	29.0 - 147							
13C-1,2,3,4,6,7,8-HpCDF	62.2	28.0 - 143							
13C-1,2,3,4,7,8,9-HpCDF	64.4	26.0 - 138							
13C-OCDF	59.4	17.0 - 157							
Cleanup Surrogate									
37Cl-2,3,7,8-TCDD	63.0	35.0 - 197							

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/30/2018

Reviewed By: 
Date: 4/30/2018

000003 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-001-OPR
Client ID: OPR
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-27-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.5	6.70 - 15.8	
1,2,3,7,8-PeCDD	51.4	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	47.6	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	49.1	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	47.5	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	49.4	35.0 - 70.0	
OCDD	104	78.0 - 144	
2,3,7,8-TCDF	9.90	7.50 - 15.8	
1,2,3,7,8-PeCDF	49.9	40.0 - 67.0	
2,3,4,7,8-PeCDF	49.4	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	49.3	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	49.4	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	49.4	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	49.9	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	49.7	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	50.3	39.0 - 69.0	
OCDF	96.8	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	85.2	20.0 - 175	
13C-1,2,3,7,8-PeCDD	86.3	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	92.2	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	88.6	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	86.5	26.0 - 166	
13C-OCDD	72.1	13.0 - 198	
13C-2,3,7,8-TCDF	90.5	22.0 - 152	
13C-1,2,3,7,8-PeCDF	82.7	21.0 - 192	
13C-2,3,4,7,8-PeCDF	85.3	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	85.0	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	83.7	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	83.3	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	83.5	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	83.3	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	86.4	20.0 - 186	
13C-OCDF	76.5	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	78.7	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000004 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-001-SA
Client ID: VS18
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.01 g
% Solids: 92.06

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.0453
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.166		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.185		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.201		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.215		-	0.0940	Total TCDD	ND	0.166	
1,2,3,7,8,9-HxCDD	0.351	-	J	0.0351	0.0823	Total PeCDD	ND	0.185	
1,2,3,4,6,7,8-HpCDD	0.868	-	J	0.00868	0.0842	Total HxCDD	0.817	-	J
OCDD	5.17	-	J	0.00155	0.172	Total HpCDD	1.99	-	J
2,3,7,8-TCDF	ND	0.129		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.143		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.153		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.182		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.189		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.197		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.207		-	0.0657	Total TCDF	0.314	-	J
1,2,3,4,6,7,8-HpCDF	ND	0.131		-	0.0747	Total PeCDF	ND	0.156	
1,2,3,4,7,8,9-HpCDF	ND	0.177		-	0.0883	Total HxCDF	ND	0.207	
OCDF	ND	0.373		-	0.170	Total HpCDF	ND	0.177	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	91.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	95.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	99.6	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	95.1	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.8	23.0 - 140	
13C-OCDD	82.1	17.0 - 157	
13C-2,3,7,8-TCDF	94.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	91.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	94.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	89.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	88.6	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	90.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	92.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	90.9	26.0 - 138	
13C-OCDF	82.6	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 83.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000005 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-002-SA
Client ID: VS20
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.01 g
% Solids: 90.84

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.252
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.216		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.269		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.226		-	0.0793				
1,2,3,6,7,8-HxCDD	0.844	-	J	0.0844	0.0940	Total TCDD	0.855	-	J
1,2,3,7,8,9-HxCDD	ND	0.216		-	0.0823	Total PeCDD	ND	0.269	
1,2,3,4,6,7,8-HpCDD	4.16	-	J	0.0416	0.0842	Total HxCDD	3.76	-	J
OCDD	19.3	-		0.00579	0.172	Total HpCDD	8.35	-	
2,3,7,8-TCDF	ND	0.170		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.184		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.196		-	0.0468				
1,2,3,4,7,8-HxCDF	0.339	-	J	0.0339	0.0437				
1,2,3,6,7,8-HxCDF	0.595	-	J	0.0595	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.148		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.156		-	0.0657	Total TCDF	1.39	-	
1,2,3,4,6,7,8-HpCDF	2.52	-	J	0.0252	0.0747	Total PeCDF	3.58	-	D,J,M
1,2,3,4,7,8,9-HpCDF	ND	0.247		-	0.0883	Total HxCDF	10.3	-	D,M
OCDF	4.00	-	J	0.00120	0.170	Total HpCDF	9.04	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	81.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	87.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	87.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	80.3	23.0 - 140	
13C-OCDD	68.9	17.0 - 157	
13C-2,3,7,8-TCDF	84.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	82.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	83.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	84.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	81.1	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	78.6	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	78.9	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	79.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	79.6	26.0 - 138	
13C-OCDF	69.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 78.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000006 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-003-SA
Client ID: VS22
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.01 g
% Solids: 92.99

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.0430
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.184		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.237		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.283		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.287		-	0.0940	Total TCDD	ND	0.184	
1,2,3,7,8,9-HxCDD	ND	0.261		-	0.0823	Total PeCDD	ND	0.237	
1,2,3,4,6,7,8-HpCDD	2.64	-	J	0.0264	0.0842	Total HxCDD	ND	0.287	
OCDD	53.2	-		0.0160	0.172	Total HpCDD	7.57	-	
2,3,7,8-TCDF	ND	0.132		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.191		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.191		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.220		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.204		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.225		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.259		-	0.0657	Total TCDF	ND	0.132	
1,2,3,4,6,7,8-HpCDF	ND	0.206		-	0.0747	Total PeCDF	ND	0.191	
1,2,3,4,7,8,9-HpCDF	ND	0.224		-	0.0883	Total HxCDF	ND	0.259	
OCDF	2.20	-	J	0.000660	0.170	Total HpCDF	0.950	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	88.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	89.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	97.5	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	93.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.0	23.0 - 140	
13C-OCDD	82.0	17.0 - 157	
13C-2,3,7,8-TCDF	92.4	24.0 - 169	
13C-1,2,3,7,8-PeCDF	87.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	90.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	89.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	87.8	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	88.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	93.2	26.0 - 138	
13C-OCDF	83.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.2 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000007 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-004-SA
Client ID: VS24
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.02 g
% Solids: 93.09

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 1.05
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.165		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.247		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.243		-	0.0793				
1,2,3,6,7,8-HxCDD	0.743	-	J	0.0743	0.0940	Total TCDD	ND	0.165	
1,2,3,7,8,9-HxCDD	0.510	-	J	0.0510	0.0823	Total PeCDD	ND	0.247	
1,2,3,4,6,7,8-HpCDD	12.8	-		0.128	0.0842	Total HxCDD	4.78	-	J
OCDD	112	-		0.0336	0.172	Total HpCDD	23.3	-	
2,3,7,8-TCDF	1.53	-		0.153	0.0269				
1,2,3,7,8-PeCDF	ND	0.228		-	0.0449				
2,3,4,7,8-PeCDF	0.995	-	J	0.299	0.0468				
1,2,3,4,7,8-HxCDF	1.02	-	J	0.102	0.0437				
1,2,3,6,7,8-HxCDF	0.528	-	J	0.0528	0.0417				
2,3,4,6,7,8-HxCDF	0.633	-	J	0.0633	0.0574				
1,2,3,7,8,9-HpCDF	0.486	-	J	0.0486	0.0657	Total TCDF	7.52	-	
1,2,3,4,6,7,8-HpCDF	4.12	-	J	0.0412	0.0747	Total PeCDF	6.47	-	
1,2,3,4,7,8,9-HpCDF	0.426	-	J	0.00426	0.0883	Total HxCDF	7.76	-	
OCDF	6.53	-	J	0.00196	0.170	Total HpCDF	9.52	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	89.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	92.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	95.1	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	92.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.4	23.0 - 140	
13C-OCDD	76.9	17.0 - 157	
13C-2,3,7,8-TCDF	93.7	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	87.5	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	86.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.1	26.0 - 138	
13C-OCDF	78.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 81.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000008 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-005-SA
Client ID: VS35
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.02 g
% Solids: 88.92

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.158
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.229		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.269		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.204		-	0.0793				
1,2,3,6,7,8-HxCDD	0.501	-	J	0.0501	0.0940	Total TCDD	ND	0.229	
1,2,3,7,8,9-HxCDD	0.433	-	J	0.0433	0.0823	Total PeCDD	ND	0.269	
1,2,3,4,6,7,8-HpCDD	1.86	-	J	0.0186	0.0842	Total HxCDD	0.935	-	J
OCDD	14.9	-		0.00447	0.172	Total HpCDD	3.51	-	J
2,3,7,8-TCDF	ND	0.194		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.202		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.223		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.172		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.188		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.202		-	0.0574				
1,2,3,7,8,9-HxCDF	0.368	-	J	0.0368	0.0657	Total TCDF	0.935	-	J
1,2,3,4,6,7,8-HpCDF	0.478	-	J	0.00478	0.0747	Total PeCDF	0.530	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.189		-	0.0883	Total HxCDF	0.871	-	J,M
OCDF	1.44	-	J	0.000432	0.170	Total HpCDF	1.81	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	85.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	82.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	84.7	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	80.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	73.8	23.0 - 140	
13C-OCDD	59.6	17.0 - 157	
13C-2,3,7,8-TCDF	84.6	24.0 - 169	
13C-1,2,3,7,8-PeCDF	80.1	24.0 - 185	
13C-2,3,4,7,8-PeCDF	78.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	79.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	76.4	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	74.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	74.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	74.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	73.9	26.0 - 138	
13C-OCDF	63.3	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 79.2 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000009 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-006-SA
Client ID: VS37
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.02 g
% Solids: 87.73

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.0529
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.147		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.207		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.187		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.213		-	0.0940	Total TCDD	ND	0.147	
1,2,3,7,8,9-HxCDD	ND	0.183		-	0.0823	Total PeCDD	ND	0.207	
1,2,3,4,6,7,8-HpCDD	1.75	-	J	0.0175	0.0842	Total HxCDD	0.648	-	J
OCDD	14.7	-		0.00441	0.172	Total HpCDD	3.88	-	J
2,3,7,8-TCDF	ND	0.142		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.154		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.160		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.133		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.145		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.146		-	0.0574				
1,2,3,7,8,9-HxCDF	0.310	-	J	0.0310	0.0657	Total TCDF	0.275	-	J
1,2,3,4,6,7,8-HpCDF	ND	0.190		-	0.0747	Total PeCDF	ND	0.160	
1,2,3,4,7,8,9-HpCDF	ND	0.222		-	0.0883	Total HxCDF	0.310	-	J
OCDF	ND	0.451		-	0.170	Total HpCDF	ND	0.222	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	98.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	105	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	99.7	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	93.5	23.0 - 140	
13C-OCDD	80.5	17.0 - 157	
13C-2,3,7,8-TCDF	98.5	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	95.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	97.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	92.7	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	93.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	91.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	94.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	95.0	26.0 - 138	
13C-OCDF	84.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000010 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-007-SA
Client ID: VS39
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.01 g
% Solids: 87.32

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.117
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.200		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.230		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.249		-	0.0793				
1,2,3,6,7,8-HxCDD	0.428	-	J	0.0428	0.0940	Total TCDD	ND	0.200	
1,2,3,7,8,9-HxCDD	0.492	-	J	0.0492	0.0823	Total PeCDD	ND	0.230	
1,2,3,4,6,7,8-HpCDD	1.99	-	J	0.0199	0.0842	Total HxCDD	1.92	-	J
OCDD	17.3	-		0.00519	0.172	Total HpCDD	4.52	-	J
2,3,7,8-TCDF	ND	0.151		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.147		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.156		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.188		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.207		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.212		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.246		-	0.0657	Total TCDF	ND	0.151	
1,2,3,4,6,7,8-HpCDF	ND	0.264		-	0.0747	Total PeCDF	ND	0.156	
1,2,3,4,7,8,9-HpCDF	ND	0.295		-	0.0883	Total HxCDF	ND	0.246	
OCDF	1.20	-	J	0.000360	0.170	Total HpCDF	0.655	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	89.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	94.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	96.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	92.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.4	23.0 - 140	
13C-OCDD	71.3	17.0 - 157	
13C-2,3,7,8-TCDF	90.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	90.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	89.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	86.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	87.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	85.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	86.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.0	26.0 - 138	
13C-OCDF	78.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 83.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000011 of 000016

EPA Method 1613 PCDD/F



FAL ID: 11441-008-SA
Client ID: VS41
Matrix: Soil
Batch No: X4491

Date Extracted: 04-26-2018
Date Received: 04-11-2018
Amount: 5.02 g
% Solids: 92.82

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-28-2018
2005 WHO TEQ: 0.215
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.180		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.226		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.236		-	0.0793				
1,2,3,6,7,8-HxCDD	0.529	-	J	0.0529	0.0940	Total TCDD	ND	0.180	
1,2,3,7,8,9-HxCDD	0.369	-	J	0.0369	0.0823	Total PeCDD	ND	0.226	
1,2,3,4,6,7,8-HpCDD	8.24	-		0.0824	0.0842	Total HxCDD	2.96	-	J
OCDD	67.3	-		0.0202	0.172	Total HpCDD	15.0	-	
2,3,7,8-TCDF	ND	0.154		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.202		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.226		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.177		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.181		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.175		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.200		-	0.0657	Total TCDF	ND	0.154	
1,2,3,4,6,7,8-HpCDF	2.01	-	J	0.0201	0.0747	Total PeCDF	ND	0.226	
1,2,3,4,7,8,9-HpCDF	ND	0.171		-	0.0883	Total HxCDF	1.98	-	J
OCDF	7.03	-	J	0.00211	0.170	Total HpCDF	5.80	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	89.9	25.0 - 164	
13C-1,2,3,7,8-PeCDD	91.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	96.8	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.6	23.0 - 140	
13C-OCDD	81.3	17.0 - 157	
13C-2,3,7,8-TCDF	95.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	88.1	24.0 - 185	
13C-2,3,4,7,8-PeCDF	91.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	86.3	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	85.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	87.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.9	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	88.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.6	26.0 - 138	
13C-OCDF	80.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/30/2018

Reviewed By: 

Date: 4/30/2018

000012 of 000016



Frontier Analytical Laboratory
5172 Hillsdale Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
Fax: 916-934-0999

FAL USE ONLY

Laboratory Project No.: 11441

Temperature: 0 °C

Chain of Custody

www.frontieranalytical.com

Please Print in Pen Page 1 of 1

CLIENT INFORMATION					INVOICE INFORMATION (if different from client info)										PROJECT INFORMATION						
Company Name: <u>Brusca Associates, Inc.</u>					Company Name: <u>Same</u>					FAL Quote #: <u>3310</u>											
Contact Name: <u>Joe Brusca</u>					Contact Name: _____					P.O. #: _____											
Address: <u>PO Box 332, Roseville, CA 95661</u>					Address: _____					Project #: <u>137-002</u>											
Phone: <u>916-677-1470</u> Fax: _____					Phone: _____ Fax: _____					Project Name: <u>LATHROP 48-Ac PROPERTY</u>											
Email: <u>jbrusca@bruscaassociates.com</u>					Email: _____					TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (✓ one)											
										* FAL must agree with price and RUSH TAT in writing.											
REPORT INFORMATION					REPORT DISTRIBUTION (email only is preferred)										ADDITIONAL INSTRUCTIONS						
Report Level: ☐ I/II ☐ III ☐ IV					☐ Hardcopy																
☐ EDD: ☐ FAL Basic ☐ Geotracker					☐ CD (.pdf including EDDs if requested)																
☐ Other: _____ ☐ Custom: Contact FAL					☐ Email (.pdf including EDDs if requested)																
☐ California State Drinking Water Form																					
System #: _____ Source #: _____																					
Sampler: _____ Employer: _____																					
Sample ID					Date	Time	Matrix	# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/9A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS	**TEQ	Remarks
					Collected														☐ 2,3,7,8-TCDD only	☐ 1998 WHO	
																			☐ 2,3,7,8-TCDD/F only	☒ 2005 WHO	
																			☒ PCDD/F (Cl ₄ -Cl ₈)	☐ Other	
1	VS18	4/11/18	11:22	SOIL		X															
2	VS20		11:29			X															
3	VS22		11:34			X															
4	VS24		11:37			X															
5	VS35		12:43			X															
6	VS37		12:48			X															
7	VS39		12:52			X															
8	VS41		12:57			X															
9																					
10																					
11																					
12																					
13																					
14																					
15																					
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																					
Relinquished by: (Signature and Printed Name)					Date	Time	Received by: (Signature and Printed Name)					Date	Time								
					4/11/18	3:45 pm	Kathy Zep					4/11/18	3:45								

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

White Copy - Report
Yellow Copy - Laboratory
Pink Copy - Originator

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: 11441

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/11/2018
Time Received:	03:45 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	8
Duplicates:	0
Storage Location:	R-4

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



May 2, 2018

FAL Project ID: 11471

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11471**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Five soil samples were received on 4/13/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of fifteen business days for project **11471**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and a sample photo. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11471**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink, appearing to read "Bradley B. Silverbush".

Bradley B. Silverbush
Director of Operations

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: 11471

Received on: 04/13/2018

Project Due: 05/07/2018 Storage: R-3

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11471-001-SA	0	137-002	VS53	EPA 1613 D/F	Soil	04/13/2018	10:44 am	04/15/2019
11471-002-SA	0	137-002	VS55	EPA 1613 D/F	Soil	04/13/2018	10:49 am	04/15/2019
11471-003-SA	0	137-002	VS57	EPA 1613 D/F	Soil	04/13/2018	10:54 am	04/15/2019
11471-004-SA	0	137-002	VS59	EPA 1613 D/F	Soil	04/13/2018	10:57 am	04/15/2019
11471-005-SA	0	137-002	VS65	EPA 1613 D/F	Soil	04/13/2018	11:23 am	04/15/2019

EPA Method 1613 PCDD/F



FAL ID: 11471-001-MB
Client ID: Method Blank
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.151		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.182		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.197		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.221		-	0.0940	Total TCDD	ND	0.151	
1,2,3,7,8,9-HxCDD	ND	0.191		-	0.0823	Total PeCDD	ND	0.182	
1,2,3,4,6,7,8-HpCDD	ND	0.180		-	0.0842	Total HxCDD	ND	0.221	
OCDD	ND	0.314		-	0.172	Total HpCDD	ND	0.180	
2,3,7,8-TCDF	ND	0.116		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.140		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.141		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.126		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.137		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.132		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.160		-	0.0657	Total TCDF	ND	0.116	
1,2,3,4,6,7,8-HpCDF	ND	0.122		-	0.0747	Total PeCDF	ND	0.141	
1,2,3,4,7,8,9-HpCDF	ND	0.211		-	0.0883	Total HxCDF	ND	0.160	
OCDF	ND	0.225		-	0.170	Total HpCDF	ND	0.211	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	104	25.0 - 164	
13C-1,2,3,7,8-PeCDD	107	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	110	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	107	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	105	23.0 - 140	
13C-OCDD	101	17.0 - 157	
13C-2,3,7,8-TCDF	111	24.0 - 169	
13C-1,2,3,7,8-PeCDF	104	24.0 - 185	
13C-2,3,4,7,8-PeCDF	108	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	99.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	99.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	97.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	98.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	75.9	26.0 - 138	
13C-OCDF	97.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 94.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000003 of 000012

EPA Method 1613 PCDD/F



FAL ID: 11471-001-OPR
Client ID: OPR
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 05-01-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.5	6.70 - 15.8	
1,2,3,7,8-PeCDD	50.2	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	50.0	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	49.6	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	47.4	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	51.5	35.0 - 70.0	
OCDD	99.1	78.0 - 144	
2,3,7,8-TCDF	11.3	7.50 - 15.8	
1,2,3,7,8-PeCDF	50.8	40.0 - 67.0	
2,3,4,7,8-PeCDF	50.8	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	51.2	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	52.2	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	51.5	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	51.8	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	51.6	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	52.6	39.0 - 69.0	
OCDF	101	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	76.9	20.0 - 175	
13C-1,2,3,7,8-PeCDD	84.4	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	97.8	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	96.9	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	87.3	26.0 - 166	
13C-OCDD	84.1	13.0 - 198	
13C-2,3,7,8-TCDF	80.3	22.0 - 152	
13C-1,2,3,7,8-PeCDF	88.2	21.0 - 192	
13C-2,3,4,7,8-PeCDF	90.0	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	88.7	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	86.6	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	86.1	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	74.7	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	81.5	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	78.0	20.0 - 186	
13C-OCDF	79.1	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	72.0	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 5/2/2018

Reviewed By: 
Date: 5/2/2018

000004 of 000012

EPA Method 1613
PCDD/F



FAL ID: 11471-001-SA
Client ID: VS53
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.04 g
% Solids: 92.01

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.0123
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.172		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.233		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.234		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.235		-	0.0940	Total TCDD	ND	0.172	
1,2,3,7,8,9-HxCDD	ND	0.215		-	0.0823	Total PeCDD	ND	0.233	
1,2,3,4,6,7,8-HpCDD	0.986	-	J	0.00986	0.0842	Total HxCDD	ND	0.235	
OCDD	8.05	-	J	0.00242	0.172	Total HpCDD	2.18	-	J
2,3,7,8-TCDF	ND	0.130		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.145		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.156		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.148		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.151		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.160		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.186		-	0.0657	Total TCDF	ND	0.130	
1,2,3,4,6,7,8-HpCDF	ND	0.246		-	0.0747	Total PeCDF	ND	0.156	
1,2,3,4,7,8,9-HpCDF	ND	0.285		-	0.0883	Total HxCDF	ND	0.186	
OCDF	ND	0.648		-	0.170	Total HpCDF	ND	0.285	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	81.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.8	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	105	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	102	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.7	23.0 - 140	
13C-OCDD	88.4	17.0 - 157	
13C-2,3,7,8-TCDF	76.6	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	93.7	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	93.9	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	92.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	90.9	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	88.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	91.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	93.4	26.0 - 138	
13C-OCDF	83.4	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 71.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000005 of 000012

EPA Method 1613 PCDD/F



FAL ID: 11471-002-SA
Client ID: VS55
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.05 g
% Solids: 91.40

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.0501
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.192		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.208		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.352		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.339		-	0.0940	Total TCDD	ND	0.192	
1,2,3,7,8,9-HxCDD	ND	0.317		-	0.0823	Total PeCDD	ND	0.208	
1,2,3,4,6,7,8-HpCDD	3.14	-	J	0.0314	0.0842	Total HxCDD	0.728	-	J
OCDD	32.0	-		0.00960	0.172	Total HpCDD	6.23	-	
2,3,7,8-TCDF	ND	0.191		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.209		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.219		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.167		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.156		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.177		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.212		-	0.0657	Total TCDF	ND	0.191	
1,2,3,4,6,7,8-HpCDF	0.850	-	J	0.00850	0.0747	Total PeCDF	0.874	-	J,M
1,2,3,4,7,8,9-HpCDF	ND	0.241		-	0.0883	Total HxCDF	1.53	-	J
OCDF	1.92	-	J	0.000576	0.170	Total HpCDF	2.98	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	79.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	107	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	105	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.1	23.0 - 140	
13C-OCDD	88.3	17.0 - 157	
13C-2,3,7,8-TCDF	80.7	24.0 - 169	
13C-1,2,3,7,8-PeCDF	92.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	93.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	96.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	96.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	92.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	82.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	79.6	26.0 - 138	
13C-OCDF	84.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 73.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000006 of 000012

EPA Method 1613
PCDD/F



FAL ID: 11471-003-SA
Client ID: VS57
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.09 g
% Solids: 91.21

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.0622
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.205		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.255		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.200		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.220		-	0.0940	Total TCDD	ND	0.205	
1,2,3,7,8,9-HxCDD	ND	0.192		-	0.0823	Total PeCDD	ND	0.255	
1,2,3,4,6,7,8-HpCDD	4.06	-	J	0.0406	0.0842	Total HxCDD	1.01	-	J
OCDD	31.8	-		0.00954	0.172	Total HpCDD	7.32	-	
2,3,7,8-TCDF	ND	0.153		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.190		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.206		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.165		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.182		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.182		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.229		-	0.0657	Total TCDF	ND	0.153	
1,2,3,4,6,7,8-HpCDF	1.11	-	J	0.0111	0.0747	Total PeCDF	ND	0.206	
1,2,3,4,7,8,9-HpCDF	ND	0.181		-	0.0883	Total HxCDF	0.433	-	J
OCDF	3.14	-	J	0.000942	0.170	Total HpCDF	2.56	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	80.3	25.0 - 164	
13C-1,2,3,7,8-PeCDD	89.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	107	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	107	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	91.7	23.0 - 140	
13C-OCDD	90.5	17.0 - 157	
13C-2,3,7,8-TCDF	81.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	95.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	94.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	95.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	93.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	93.6	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	84.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	80.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.6	26.0 - 138	
13C-OCDF	85.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 71.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000007 of 000012

EPA Method 1613 PCDD/F



FAL ID: 11471-004-SA
Client ID: VS59
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.02 g
% Solids: 93.03

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.332
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.154		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.260		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.323		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.351		-	0.0940	Total TCDD	0.433	-	J
1,2,3,7,8,9-HxCDD	ND	0.308		-	0.0823	Total PeCDD	ND	0.260	
1,2,3,4,6,7,8-HpCDD	3.03	-	J	0.0303	0.0842	Total HxCDD	ND	0.351	
OCDD	31.5	-		0.00945	0.172	Total HpCDD	5.92	-	
2,3,7,8-TCDF	ND	0.197		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.167		-	0.0449				
2,3,4,7,8-PeCDF	0.952	-	J	0.286	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.157		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.157		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.163		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.205		-	0.0657	Total TCDF	1.37	-	M
1,2,3,4,6,7,8-HpCDF	0.586	-	J	0.00586	0.0747	Total PeCDF	5.22	-	
1,2,3,4,7,8,9-HpCDF	ND	0.139		-	0.0883	Total HxCDF	1.02	-	J
OCDF	3.44	-	J	0.00103	0.170	Total HpCDF	2.28	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	80.5	25.0 - 164	
13C-1,2,3,7,8-PeCDD	88.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	106	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	102	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	87.8	23.0 - 140	
13C-OCDD	82.7	17.0 - 157	
13C-2,3,7,8-TCDF	80.7	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.0	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	96.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	93.0	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	90.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	83.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	81.6	26.0 - 138	
13C-OCDF	79.4	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 74.8 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000008 of 000012

EPA Method 1613 PCDD/F



FAL ID: 11471-005-SA
Client ID: VS65
Matrix: Soil
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.01 g
% Solids: 95.58

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.139
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.184		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.232		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.222		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.239		-	0.0940	Total TCDD	ND	0.184	
1,2,3,7,8,9-HxCDD	ND	0.211		-	0.0823	Total PeCDD	ND	0.232	
1,2,3,4,6,7,8-HpCDD	5.37	-		0.0537	0.0842	Total HxCDD	2.20	-	J
OCDD	52.8	-		0.0158	0.172	Total HpCDD	10.2	-	
2,3,7,8-TCDF	0.561	-	J	0.0561	0.0269				
1,2,3,7,8-PeCDF	ND	0.131		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.140		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.131		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.139		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.148		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.194		-	0.0657	Total TCDF	1.91	-	
1,2,3,4,6,7,8-HpCDF	1.18	-	J	0.0118	0.0747	Total PeCDF	0.776	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.143		-	0.0883	Total HxCDF	1.20	-	J
OCDF	4.63	-	J	0.00139	0.170	Total HpCDF	4.03	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	78.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.3	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	107	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	100	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	88.8	23.0 - 140	
13C-OCDD	86.0	17.0 - 157	
13C-2,3,7,8-TCDF	77.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	91.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	95.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	92.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	78.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	74.4	26.0 - 138	
13C-OCDF	82.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 69.2 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000009 of 000012



Frontier Analytical Laboratory
5172 Hillside Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
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FAL USE ONLY

Laboratory Project No.: 11471

Temperature: 0 °C

Chain of Custody

www.frontieranalytical.com

Please Print in Pen Page 1 of 1

CLIENT INFORMATION				INVOICE INFORMATION (if different from client info)				PROJECT INFORMATION								
Company Name: Brusca Associates, Inc.				Company Name: Same				FAL Quote #: 3310								
Contact Name: Joe Brusca				Contact Name:				P.O. #:								
Address: PO Box 332, Roseville, CA 95661				Address:				Project #: <u>137-002</u>								
Phone: 916-677-1470 Fax:				Phone: Fax:				Project Name: <u>LATHROP 48-AC PROPERTY</u>								
Email: jbrusca@bruscaassociates.com				Email:				TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (✓ one)								
								* FAL must agree with price and RUSH TAT in writing.								
REPORT INFORMATION				REPORT DISTRIBUTION (email only is preferred)				ADDITIONAL INSTRUCTIONS								
Report Level: ☐ I/II ☐ III ☐ IV				☐ Hardcopy				Kathy to Joe 15-TAT								
☐ EDD: ☐ FAL Basic ☐ Geotracker				☐ CD (.pdf including EDDs if requested)												
☐ Other: ☐ Custom: Contact FAL				☐ Email (.pdf including EDDs if requested)												
☐ California State Drinking Water Form																
System #: Source #:																
Sampler: Employer:																
Sample ID	Date	Time	Matrix	# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/9A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS	**TEQ
	Collected														☐ 2,3,7,8-TCDD only	☐ 1998 WHO
															☐ 2,3,7,8-TCDD/F only	☒ 2005 WHO
															☒ PCDD/F (Cl ₄ -Cl ₈)	☐ Other
1	VS53	4/13/18	10:44	SOIL												Remarks
2	VS55		10:49													
3	VS57		10:54													
4	VS59		10:57													
5	VS65		11:23													
6																
7																
8																
9																
10																
11																
12																
13																
14																
15																
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																
Relinquished by: (Signature and Printed Name)				Date	Time	Received by: (Signature and Printed Name)				Date	Time					
				4/13/18	2:12 PM	Kathy Zapp KZipp				4/13/18	1412					

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

White Copy - Report

Yellow Copy - Laboratory

Pink Copy - Originator

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11471**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/13/2018
Time Received:	02:12 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	5
Duplicates:	0
Storage Location:	R-3

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



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17 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48 Acre Property

Enclosed are the results of analyses for samples received by the laboratory on 04/05/18 09:55. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP1-1	T181153-01	Soil	04/04/18 11:59	04/05/18 09:55
SP1-5	T181153-05	Soil	04/04/18 12:12	04/05/18 09:55
SP1-6	T181153-06	Soil	04/04/18 12:13	04/05/18 09:55
SP1-8	T181153-08	Soil	04/04/18 12:22	04/05/18 09:55

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

DETECTIONS SUMMARY

Sample ID: SP1-1

Laboratory ID: T181153-01

Reporting

Analyte	Result	Limit	Units	Method	Notes
Antimony	1.5	1.0	ug/l	6020 ICP-MS	AO-1
Arsenic	30	1.0	ug/l	6020 ICP-MS	AO-1
Barium	480	1.0	ug/l	6020 ICP-MS	AO-1
Beryllium	1.2	1.0	ug/l	6020 ICP-MS	AO-1
Cadmium	2.1	1.0	ug/l	6020 ICP-MS	AO-1
Chromium	69	1.0	ug/l	6020 ICP-MS	AO-1
Cobalt	45	1.0	ug/l	6020 ICP-MS	AO-1
Copper	120	1.0	ug/l	6020 ICP-MS	AO-1
Lead	300	2.0	ug/l	6020 ICP-MS	AO-1
Mercury	0.18	0.10	ug/l	6020 ICP-MS	AO-1
Molybdenum	21	1.0	ug/l	6020 ICP-MS	AO-1
Nickel	60	1.0	ug/l	6020 ICP-MS	AO-1
Silver	5.3	1.0	ug/l	6020 ICP-MS	AO-1
Vanadium	260	1.0	ug/l	6020 ICP-MS	AO-1
Zinc	440	1.0	ug/l	6020 ICP-MS	AO-1

Sample ID: SP1-5

Laboratory ID: T181153-05

Reporting

Analyte	Result	Limit	Units	Method	Notes
Antimony	1.5	1.0	ug/l	6020 ICP-MS	AO-1
Arsenic	20	1.0	ug/l	6020 ICP-MS	AO-1
Barium	340	1.0	ug/l	6020 ICP-MS	AO-1
Cadmium	1.6	1.0	ug/l	6020 ICP-MS	AO-1
Chromium	60	1.0	ug/l	6020 ICP-MS	AO-1
Cobalt	39	1.0	ug/l	6020 ICP-MS	AO-1
Copper	200	1.0	ug/l	6020 ICP-MS	AO-1
Lead	360	2.0	ug/l	6020 ICP-MS	AO-1
Mercury	0.28	0.10	ug/l	6020 ICP-MS	AO-1
Molybdenum	16	1.0	ug/l	6020 ICP-MS	AO-1
Nickel	45	1.0	ug/l	6020 ICP-MS	AO-1
Silver	9.0	1.0	ug/l	6020 ICP-MS	AO-1
Vanadium	160	1.0	ug/l	6020 ICP-MS	AO-1

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

Sample ID: SP1-5

Laboratory ID: T181153-05

Analyte	Result	Reporting	Units	Method	Notes
		Limit			
Zinc	350	1.0	ug/l	6020 ICP-MS	AO-1

Sample ID: SP1-6

Laboratory ID: T181153-06

Analyte	Result	Reporting	Units	Method	Notes
		Limit			
Antimony	1.5	1.0	ug/l	6020 ICP-MS	AO-1
Arsenic	24	1.0	ug/l	6020 ICP-MS	AO-1
Barium	510	1.0	ug/l	6020 ICP-MS	AO-1
Beryllium	1.1	1.0	ug/l	6020 ICP-MS	AO-1
Cadmium	2.0	1.0	ug/l	6020 ICP-MS	AO-1
Chromium	56	1.0	ug/l	6020 ICP-MS	AO-1
Cobalt	39	1.0	ug/l	6020 ICP-MS	AO-1
Copper	110	1.0	ug/l	6020 ICP-MS	AO-1
Lead	390	2.0	ug/l	6020 ICP-MS	AO-1
Mercury	0.22	0.10	ug/l	6020 ICP-MS	AO-1
Molybdenum	16	1.0	ug/l	6020 ICP-MS	AO-1
Nickel	54	1.0	ug/l	6020 ICP-MS	AO-1
Silver	5.0	1.0	ug/l	6020 ICP-MS	AO-1
Vanadium	190	1.0	ug/l	6020 ICP-MS	AO-1
Zinc	380	1.0	ug/l	6020 ICP-MS	AO-1

Sample ID: SP1-8

Laboratory ID: T181153-08

Analyte	Result	Reporting	Units	Method	Notes
		Limit			
Antimony	1.0	1.0	ug/l	6020 ICP-MS	AO-1
Arsenic	27	1.0	ug/l	6020 ICP-MS	AO-1
Barium	520	1.0	ug/l	6020 ICP-MS	AO-1
Beryllium	1.1	1.0	ug/l	6020 ICP-MS	AO-1
Cadmium	1.1	1.0	ug/l	6020 ICP-MS	AO-1
Chromium	57	1.0	ug/l	6020 ICP-MS	AO-1
Cobalt	37	1.0	ug/l	6020 ICP-MS	AO-1
Copper	85	1.0	ug/l	6020 ICP-MS	AO-1
Lead	140	2.0	ug/l	6020 ICP-MS	AO-1
Mercury	0.14	0.10	ug/l	6020 ICP-MS	AO-1
Molybdenum	24	1.0	ug/l	6020 ICP-MS	AO-1

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

Sample ID: SP1-8

Laboratory ID: T181153-08

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Nickel	53	1.0		ug/l	6020 ICP-MS	AO-1
Silver	3.5	1.0		ug/l	6020 ICP-MS	AO-1
Vanadium	240	1.0		ug/l	6020 ICP-MS	AO-1
Zinc	370	1.0		ug/l	6020 ICP-MS	AO-1

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

SP1-1
T181153-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	1.5	1.0	ug/l	1	8041135	04/12/18	04/17/18	6020 ICP-MS	AO-1
Arsenic	30	1.0	"	"	"	"	"	"	AO-1
Barium	480	1.0	"	"	"	"	"	"	AO-1
Beryllium	1.2	1.0	"	"	"	"	"	"	AO-1
Cadmium	2.1	1.0	"	"	"	"	"	"	AO-1
Chromium	69	1.0	"	"	"	"	"	"	AO-1
Cobalt	45	1.0	"	"	"	"	"	"	AO-1
Copper	120	1.0	"	"	"	"	"	"	AO-1
Lead	300	2.0	"	"	"	"	"	"	AO-1
Mercury	0.18	0.10	"	"	"	"	"	"	AO-1
Molybdenum	21	1.0	"	"	"	"	"	"	AO-1
Nickel	60	1.0	"	"	"	"	"	"	AO-1
Selenium	ND	5.0	"	"	"	"	"	"	AO-1
Silver	5.3	1.0	"	"	"	"	"	"	AO-1
Thallium	ND	1.0	"	"	"	"	"	"	AO-1
Vanadium	260	1.0	"	"	"	"	"	"	AO-1
Zinc	440	1.0	"	"	"	"	"	"	AO-1

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

SP1-5
T181153-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	1.5	1.0	ug/l	1	8041135	04/11/18	04/17/18	6020 ICP-MS	AO-1
Arsenic	20	1.0	"	"	"	"	"	"	AO-1
Barium	340	1.0	"	"	"	"	"	"	AO-1
Beryllium	ND	1.0	"	"	"	"	"	"	AO-1
Cadmium	1.6	1.0	"	"	"	"	"	"	AO-1
Chromium	60	1.0	"	"	"	"	"	"	AO-1
Cobalt	39	1.0	"	"	"	"	"	"	AO-1
Copper	200	1.0	"	"	"	"	"	"	AO-1
Lead	360	2.0	"	"	"	"	"	"	AO-1
Mercury	0.28	0.10	"	"	"	"	"	"	AO-1
Molybdenum	16	1.0	"	"	"	"	"	"	AO-1
Nickel	45	1.0	"	"	"	"	"	"	AO-1
Selenium	ND	5.0	"	"	"	"	"	"	AO-1
Silver	9.0	1.0	"	"	"	"	"	"	AO-1
Thallium	ND	1.0	"	"	"	"	"	"	AO-1
Vanadium	160	1.0	"	"	"	"	"	"	AO-1
Zinc	350	1.0	"	"	"	"	"	"	AO-1

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

SP1-6
T181153-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	1.5	1.0	ug/l	1	8041135	04/12/18	04/17/18	6020 ICP-MS	AO-1
Arsenic	24	1.0	"	"	"	"	"	"	AO-1
Barium	510	1.0	"	"	"	"	"	"	AO-1
Beryllium	1.1	1.0	"	"	"	"	"	"	AO-1
Cadmium	2.0	1.0	"	"	"	"	"	"	AO-1
Chromium	56	1.0	"	"	"	"	"	"	AO-1
Cobalt	39	1.0	"	"	"	"	"	"	AO-1
Copper	110	1.0	"	"	"	"	"	"	AO-1
Lead	390	2.0	"	"	"	"	"	"	AO-1
Mercury	0.22	0.10	"	"	"	"	"	"	AO-1
Molybdenum	16	1.0	"	"	"	"	"	"	AO-1
Nickel	54	1.0	"	"	"	"	"	"	AO-1
Selenium	ND	5.0	"	"	"	"	"	"	AO-1
Silver	5.0	1.0	"	"	"	"	"	"	AO-1
Thallium	ND	1.0	"	"	"	"	"	"	AO-1
Vanadium	190	1.0	"	"	"	"	"	"	AO-1
Zinc	380	1.0	"	"	"	"	"	"	AO-1

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

SP1-8
T181153-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	1.0	1.0	ug/l	1	8041135	04/12/18	04/17/18	6020 ICP-MS	AO-1
Arsenic	27	1.0	"	"	"	"	"	"	AO-1
Barium	520	1.0	"	"	"	"	"	"	AO-1
Beryllium	1.1	1.0	"	"	"	"	"	"	AO-1
Cadmium	1.1	1.0	"	"	"	"	"	"	AO-1
Chromium	57	1.0	"	"	"	"	"	"	AO-1
Cobalt	37	1.0	"	"	"	"	"	"	AO-1
Copper	85	1.0	"	"	"	"	"	"	AO-1
Lead	140	2.0	"	"	"	"	"	"	AO-1
Mercury	0.14	0.10	"	"	"	"	"	"	AO-1
Molybdenum	24	1.0	"	"	"	"	"	"	AO-1
Nickel	53	1.0	"	"	"	"	"	"	AO-1
Selenium	ND	5.0	"	"	"	"	"	"	AO-1
Silver	3.5	1.0	"	"	"	"	"	"	AO-1
Thallium	ND	1.0	"	"	"	"	"	"	AO-1
Vanadium	240	1.0	"	"	"	"	"	"	AO-1
Zinc	370	1.0	"	"	"	"	"	"	AO-1

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041135 - EPA 3010A

Blank (8041135-BLK1)

Prepared: 04/11/18 Analyzed: 04/17/18

Antimony	ND	1.0	ug/l							AO-1
Arsenic	ND	1.0	"							AO-1
Barium	16.1	1.0	"							AO-1, QB-01
Beryllium	ND	1.0	"							AO-1
Cadmium	ND	1.0	"							AO-1
Chromium	1.93	1.0	"							AO-1, QB-01
Cobalt	ND	1.0	"							AO-1
Copper	1.50	1.0	"							AO-1, QB-01
Lead	ND	2.0	"							AO-1
Mercury	ND	0.10	"							AO-1
Molybdenum	ND	1.0	"							AO-1
Nickel	ND	1.0	"							AO-1
Selenium	ND	5.0	"							AO-1
Silver	ND	1.0	"							AO-1
Thallium	ND	1.0	"							AO-1
Vanadium	3.01	1.0	"							AO-1, QB-01
Zinc	11.1	1.0	"							AO-1, QB-01

LCS (8041135-BS1)

Prepared: 04/11/18 Analyzed: 04/17/18

Arsenic	48.2	1.0	ug/l	50.0		96.4	80-120			AO-1
Barium	48.8	1.0	"	50.0		97.5	80-120			AO-1
Cadmium	47.7	1.0	"	50.0		95.4	80-120			AO-1
Chromium	52.4	1.0	"	50.0		105	80-120			AO-1
Lead	51.3	2.0	"	50.0		103	80-120			AO-1

Duplicate (8041135-DUP1)

Source: T181153-05

Prepared: 04/11/18 Analyzed: 04/17/18

Barium	432	5.0	ug/l		341		23.6	200		AO-1
Lead	445	10	"		361		20.8	200		AO-1

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041135 - EPA 3010A

Matrix Spike (8041135-MS1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Arsenic	62.4	1.0	ug/l	50.0	19.8	85.2	75-125			AO-1
Barium	363	1.0	"	50.0	341	44.9	75-125			AO-1, QM-05
Cadmium	45.9	1.0	"	50.0	1.59	88.7	75-125			AO-1
Chromium	107	1.0	"	50.0	59.6	94.3	75-125			AO-1
Lead	368	2.0	"	50.0	361	12.3	75-125			AO-1, QM-05

Matrix Spike Dup (8041135-MSD1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Arsenic	61.3	1.0	ug/l	50.0	19.8	82.9	75-125	1.81	20	AO-1
Barium	372	1.0	"	50.0	341	62.8	75-125	2.43	20	AO-1, QM-05
Cadmium	45.7	1.0	"	50.0	1.59	88.2	75-125	0.541	20	AO-1
Chromium	107	1.0	"	50.0	59.6	94.9	75-125	0.296	20	AO-1
Lead	389	2.0	"	50.0	361	55.4	75-125	5.69	20	AO-1, QM-05

Post Spike (8041135-PS1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Barium	188		ug/l	2.50	0.341	NR	80-120			AO-1, QM-05
Lead	199		"	2.50	0.361	NR	80-120			AO-1, QM-05

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/17/18 15:58

Notes and Definitions

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

QB-01 The method blank contains analyte at a concentration above the MRL; however, concentration is less than 10% of the sample result, which is negligible according to method criteria.

AO-1 STLC Leach Ran 6020 Water

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commercentre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO BOX 332, ROSEMEAD, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/4/18 Page: 1 of 2
Project Name: LATHROP 48-Ac PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: T181153 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 PCBs 8092	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD AND ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers
SP1-1	4/4/18	11:59	SOLID	JAR	X	X	X	X	X	X	X	X	X	X	X	01		1
SP1-2	-	12:01	-	-	X	X	X	X	X	X	X	X	X	X	X	02		1
SP1-3	-	12:03	-	-	X	X	X	X	X	X	X	X	X	X	X	03		1
SP1-4	-	12:05	-	-	X	X	X	X	X	X	X	X	X	X	X	04		1
SP1-5	-	12:12	-	-	X	X	X	X	X	X	X	X	X	X	X	05		1
SP1-6	-	12:13	-	-	X	X	X	X	X	X	X	X	X	X	X	06		1
SP1-7	-	12:17	-	-	X	X	X	X	X	X	X	X	X	X	X	07		1
SP1-8	-	12:22	-	-	X	X	X	X	X	X	X	X	X	X	X	08		1
SP1-9	-	12:25	-	-	X	X	X	X	X	X	X	X	X	X	X	09		1
SP1-10	-	12:31	-	-	X	X	X	X	X	X	X	X	X	X	X	10		1
SP1-11	-	12:35	-	-	X	X	X	X	X	X	X	X	X	X	X	11		1
SP1-12	-	12:37	-	-	X	X	X	X	X	X	X	X	X	X	X	12		1
SP1-13	-	12:42	-	-	X	X	X	X	X	X	X	X	X	X	X	13		1
SP1-14	-	12:45	-	-	X	X	X	X	X	X	X	X	X	X	X	14		1
SP1-15	-	12:47	-	-	X	X	X	X	X	X	X	X	X	X	X	15		1
Relinquished by: (signature)	Date / Time	4/4/18 14:38	Received by: (signature)	Date / Time	4/4/18 14:38	Total # of containers												
Relinquished by: (signature)	Date / Time	4-5-18 955	Received by: (signature)	Date / Time	4-5-18 955	Chain of Custody seals DN/A Seals intact? DN/A Received good condition/cold 4.1												
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time		Notes * 3 DAY TAT ON METALS												

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

Turn around time: NORMAL
EXCEPT METALS -
3 DAY TAT ON METALS **COC 160906**



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Centre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, ROSEVILLE, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/4/18 Page: 2 of 2
Project Name: LATHROP 48-Ac PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: T181153 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260	8260 + OXY	8260 BTEX, OXY only	8270	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP1-16	4/4/18	12:51	SOLID	JAR	X	X	X	X	X	X	X	X	X	X	14		1
SP1-17		12:53			X	X	X	X	X	X	X	X	X	X	17		1
SP1-18		12:56			X	X	X	X	X	X	X	X	X	X	18		1
SP1-19		13:05			X	X	X	X	X	X	X	X	X	X	19		1
SP1-20		13:08			X	X	X	X	X	X	X	X	X	X	20		1
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Total # of containers												
<u>[Signature]</u>	4/4/18 14:38		<u>[Signature]</u>	4/4/18 14:38													
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Chain of Custody seals 9/N/A												
<u>[Signature]</u>	4-5-18 9:55		<u>[Signature]</u>	4-5-18 9:55	Seals intact? 0/N/A												
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Received good condition/cold												
<u>[Signature]</u>			<u>[Signature]</u>		4.1												

Notes
* 3 DAY TAT
ON METALS

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____
Turn around time: Normal
EXCEPT METALS -
3 DAY TAT ON METALS
COC 160912



April 19, 2018

FAL Project ID: 11423

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11423**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Twenty solid samples were received on 4/4/2018 in good condition. All twenty samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of ten business days for project **11423**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and a sample photo. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11423**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink, appearing to read "Bradley B. Silverbush".

Bradley B. Silverbush
Director of Operations

FRONTIER ANALYTICAL LABORATORY

5172 Hillsdale Circle * El Dorado Hills, CA 95762
Tel (916) 934-0900 * Fax (916) 934-0999
www.frontieranalytical.com

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Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: **11423**

Received on: **04/04/2018**

Project Due: **04/19/2018**

Storage: **R-4**

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11423-001-SA	0	137-002	SP1-1	EPA 1613 D/F	Solid	04/04/2018	11:59 am	04/04/2019
11423-002-SA	0	137-002	SP1-2	EPA 1613 D/F	Solid	04/04/2018	12:01 pm	04/04/2019
11423-003-SA	0	137-002	SP1-3	EPA 1613 D/F	Solid	04/04/2018	12:03 pm	04/04/2019
11423-004-SA	0	137-002	SP1-4	EPA 1613 D/F	Solid	04/04/2018	12:05 pm	04/04/2019
11423-005-SA	0	137-002	SP1-5	EPA 1613 D/F	Solid	04/04/2018	12:12 pm	04/04/2019
11423-006-SA	0	137-002	SP1-6	EPA 1613 D/F	Solid	04/04/2018	12:13 pm	04/04/2019
11423-007-SA	0	137-002	SP1-7	EPA 1613 D/F	Solid	04/04/2018	12:17 pm	04/04/2019
11423-008-SA	0	137-002	SP1-8	EPA 1613 D/F	Solid	04/04/2018	12:22 pm	04/04/2019
11423-009-SA	0	137-002	SP1-9	EPA 1613 D/F	Solid	04/04/2018	12:25 pm	04/04/2019
11423-010-SA	0	137-002	SP1-10	EPA 1613 D/F	Solid	04/04/2018	12:31 pm	04/04/2019
11423-011-SA	0	137-002	SP1-11	EPA 1613 D/F	Solid	04/04/2018	12:35 pm	04/04/2019
11423-012-SA	0	137-002	SP1-12	EPA 1613 D/F	Solid	04/04/2018	12:37 pm	04/04/2019
11423-013-SA	0	137-002	SP1-13	EPA 1613 D/F	Solid	04/04/2018	12:42 pm	04/04/2019
11423-014-SA	0	137-002	SP1-14	EPA 1613 D/F	Solid	04/04/2018	12:45 pm	04/04/2019
11423-015-SA	0	137-002	SP1-15	EPA 1613 D/F	Solid	04/04/2018	12:47 pm	04/04/2019
11423-016-SA	0	137-002	SP1-16	EPA 1613 D/F	Solid	04/04/2018	12:51 pm	04/04/2019
11423-017-SA	0	137-002	SP1-17	EPA 1613 D/F	Solid	04/04/2018	12:53 pm	04/04/2019
11423-018-SA	0	137-002	SP1-18	EPA 1613 D/F	Solid	04/04/2018	12:56 pm	04/04/2019
11423-019-SA	0	137-002	SP1-19	EPA 1613 D/F	Solid	04/04/2018	01:05 pm	04/04/2019
11423-020-SA	0	137-002	SP1-20	EPA 1613 D/F	Solid	04/04/2018	01:08 pm	04/04/2019

EPA Method 1613 PCDD/F



FAL ID: 11423-001-MB
Client ID: Method Blank
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.271		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.342		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.395		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.463		-	0.0940	Total TCDD	ND	0.271	
1,2,3,7,8,9-HxCDD	ND	0.393		-	0.0823	Total PeCDD	ND	0.342	
1,2,3,4,6,7,8-HpCDD	ND	0.544		-	0.0842	Total HxCDD	ND	0.463	
OCDD	ND	0.711		-	0.172	Total HpCDD	ND	0.544	
2,3,7,8-TCDF	ND	0.200		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.222		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.233		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.251		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.253		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.281		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.320		-	0.0657	Total TCDF	ND	0.200	
1,2,3,4,6,7,8-HpCDF	ND	0.339		-	0.0747	Total PeCDF	ND	0.233	
1,2,3,4,7,8,9-HpCDF	ND	0.430		-	0.0883	Total HxCDF	ND	0.320	
OCDF	ND	0.606		-	0.170	Total HpCDF	ND	0.430	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	80.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	82.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	82.5	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	84.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	78.2	23.0 - 140	
13C-OCDD	65.4	17.0 - 157	
13C-2,3,7,8-TCDF	83.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	78.9	24.0 - 185	
13C-2,3,4,7,8-PeCDF	78.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	80.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	79.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	75.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	73.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	75.9	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	77.0	26.0 - 138	
13C-OCDF	67.4	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 76.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000003 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-001-OPR
Client ID: OPR
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-18-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.6	6.70 - 15.8	
1,2,3,7,8-PeCDD	52.3	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	49.6	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	50.5	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	49.5	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	52.1	35.0 - 70.0	
OCDD	103	78.0 - 144	
2,3,7,8-TCDF	10.6	7.50 - 15.8	
1,2,3,7,8-PeCDF	52.4	40.0 - 67.0	
2,3,4,7,8-PeCDF	52.9	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	51.9	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	52.2	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	52.4	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	53.1	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	51.3	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	54.1	39.0 - 69.0	
OCDF	105	63.0 - 170	
Internal Standards			
13C-2,3,7,8-TCDD	78.8	20.0 - 175	
13C-1,2,3,7,8-PeCDD	85.3	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	82.7	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	80.1	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	78.0	26.0 - 166	
13C-OCDD	67.4	13.0 - 198	
13C-2,3,7,8-TCDF	81.0	22.0 - 152	
13C-1,2,3,7,8-PeCDF	80.3	21.0 - 192	
13C-2,3,4,7,8-PeCDF	79.5	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	77.7	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	74.5	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	74.6	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	74.5	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	77.4	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	77.0	20.0 - 186	
13C-OCDF	67.9	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	76.4	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/18/2018

Reviewed By: 
Date: 4/18/2018

000004 of 000033

EPA Method 1613
PCDD/F



FAL ID: 11423-001-MB
Client ID: Method Blank
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.162		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.199		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.283		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.287		-	0.0940	Total TCDD	ND	0.162	
1,2,3,7,8,9-HxCDD	ND	0.261		-	0.0823	Total PeCDD	ND	0.199	
1,2,3,4,6,7,8-HpCDD	ND	0.328		-	0.0842	Total HxCDD	ND	0.287	
OCDD	ND	0.708		-	0.172	Total HpCDD	ND	0.328	
2,3,7,8-TCDF	ND	0.110		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.171		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.172		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.261		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.256		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.276		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.309		-	0.0657	Total TCDF	ND	0.110	
1,2,3,4,6,7,8-HpCDF	ND	0.314		-	0.0747	Total PeCDF	ND	0.172	
1,2,3,4,7,8,9-HpCDF	ND	0.358		-	0.0883	Total HxCDF	ND	0.309	
OCDF	ND	0.688		-	0.170	Total HpCDF	ND	0.358	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	96.4	25.0 - 164	
13C-1,2,3,7,8-PeCDD	97.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	102	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	102	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.7	23.0 - 140	
13C-OCDD	72.0	17.0 - 157	
13C-2,3,7,8-TCDF	95.4	24.0 - 169	
13C-1,2,3,7,8-PeCDF	93.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	93.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	91.1	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	90.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	88.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	90.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	90.5	26.0 - 138	
13C-OCDF	75.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 89.1 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000005 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-001-OPR
Client ID: OPR
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-18-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.6	6.70 - 15.8	
1,2,3,7,8-PeCDD	52.1	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	48.7	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	49.2	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	46.8	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	50.3	35.0 - 70.0	
OCDD	103	78.0 - 144	
2,3,7,8-TCDF	10.4	7.50 - 15.8	
1,2,3,7,8-PeCDF	51.3	40.0 - 67.0	
2,3,4,7,8-PeCDF	52.4	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	50.7	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	51.2	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	51.3	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	52.7	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	52.9	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	52.7	39.0 - 69.0	
OCDF	103	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.0	20.0 - 175	
13C-1,2,3,7,8-PeCDD	97.1	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	100	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	102	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	91.1	26.0 - 166	
13C-OCDD	76.3	13.0 - 198	
13C-2,3,7,8-TCDF	96.9	22.0 - 152	
13C-1,2,3,7,8-PeCDF	93.1	21.0 - 192	
13C-2,3,4,7,8-PeCDF	92.9	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	94.3	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	91.2	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	89.1	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	86.5	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	87.3	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	88.1	20.0 - 186	
13C-OCDF	77.7	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	85.8	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/19/2018

Reviewed By: 
Date: 4/19/2018

000006 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-001-SA
Client ID: SP1-1
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.04 g
% Solids: 93.84

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 6.49
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	0.375	-	J	0.375	0.0273				
1,2,3,7,8-PeCDD	1.48	-	J	1.48	0.0570				
1,2,3,4,7,8-HxCDD	2.29	-	J	0.229	0.0793				
1,2,3,6,7,8-HxCDD	6.04	-		0.604	0.0940	Total TCDD	2.38	-	
1,2,3,7,8,9-HxCDD	3.47	-	J	0.347	0.0823	Total PeCDD	11.2	-	
1,2,3,4,6,7,8-HpCDD	171	-		1.71	0.0842	Total HxCDD	47.6	-	
OCDD	1270	-		0.381	0.172	Total HpCDD	294	-	
2,3,7,8-TCDF	2.26	-	F	0.226	0.0269				
1,2,3,7,8-PeCDF	0.749	-	J	0.0225	0.0449				
2,3,4,7,8-PeCDF	1.29	-	J	0.387	0.0468				
1,2,3,4,7,8-HxCDF	1.36	-	J	0.136	0.0437				
1,2,3,6,7,8-HxCDF	1.49	-	J	0.149	0.0417				
2,3,4,6,7,8-HxCDF	1.66	-	J	0.166	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.296		-	0.0657	Total TCDF	17.6	-	
1,2,3,4,6,7,8-HpCDF	23.4	-		0.234	0.0747	Total PeCDF	17.2	-	
1,2,3,4,7,8,9-HpCDF	1.88	-	J	0.0188	0.0883	Total HxCDF	32.0	-	
OCDF	91.2	-		0.0274	0.170	Total HpCDF	80.0	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	87.9	25.0 - 164	
13C-1,2,3,7,8-PeCDD	91.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	95.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	89.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	85.1	23.0 - 140	
13C-OCDD	70.8	17.0 - 157	
13C-2,3,7,8-TCDF	92.2	24.0 - 169	
13C-1,2,3,7,8-PeCDF	87.4	24.0 - 185	
13C-2,3,4,7,8-PeCDF	87.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	87.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	83.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	81.2	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	80.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	85.1	26.0 - 138	
13C-OCDF	73.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000007 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-002-SA
Client ID: SP1-2
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.03 g
% Solids: 94.12

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.668
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.209		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.390		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.353		-	0.0793				
1,2,3,6,7,8-HxCDD	1.12	-	J	0.112	0.0940	Total TCDD	0.542	-	J
1,2,3,7,8,9-HxCDD	0.700	-	J	0.0700	0.0823	Total PeCDD	ND	0.390	
1,2,3,4,6,7,8-HpCDD	16.1	-		0.161	0.0842	Total HxCDD	6.96	-	
OCDD	126	-		0.0378	0.172	Total HpCDD	30.6	-	
2,3,7,8-TCDF	0.628	-	J	0.0628	0.0269				
1,2,3,7,8-PeCDF	ND	0.182		-	0.0449				
2,3,4,7,8-PeCDF	0.464	-	J	0.139	0.0468				
1,2,3,4,7,8-HxCDF	0.443	-	J	0.0443	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.193		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.222		-	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.238		-	0.0657	Total TCDF	2.57	-	
1,2,3,4,6,7,8-HpCDF	3.80	-	J	0.0380	0.0747	Total PeCDF	3.62	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.255		-	0.0883	Total HxCDF	4.88	-	J
OCDF	9.14	-	J	0.00274	0.170	Total HpCDF	9.45	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	92.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	95.3	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	98.1	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.1	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	93.3	23.0 - 140	
13C-OCDD	83.7	17.0 - 157	
13C-2,3,7,8-TCDF	94.5	24.0 - 169	
13C-1,2,3,7,8-PeCDF	88.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	89.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	89.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.3	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	88.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	90.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	91.9	26.0 - 138	
13C-OCDF	82.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 82.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

EPA Method 1613 PCDD/F



FAL ID: 11423-003-SA
Client ID: SP1-3
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.02 g
% Solids: 94.41

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 1.10
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.194		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.281		-	0.0570				
1,2,3,4,7,8-HxCDD	0.702	-	J	0.0702	0.0793				
1,2,3,6,7,8-HxCDD	1.47	-	J	0.147	0.0940	Total TCDD	0.443	-	J
1,2,3,7,8,9-HxCDD	1.04	-	J	0.104	0.0823	Total PeCDD	0.610	-	J
1,2,3,4,6,7,8-HpCDD	24.3	-		0.243	0.0842	Total HxCDD	9.86	-	
OCDD	194	-		0.0582	0.172	Total HpCDD	45.0	-	
2,3,7,8-TCDF	1.01	-		0.101	0.0269				
1,2,3,7,8-PeCDF	ND	0.241		-	0.0449				
2,3,4,7,8-PeCDF	0.515	-	J	0.155	0.0468				
1,2,3,4,7,8-HxCDF	0.630	-	J	0.0630	0.0437				
1,2,3,6,7,8-HxCDF	0.422	-	J	0.0422	0.0417				
2,3,4,6,7,8-HxCDF	0.601	-	J	0.0601	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.248		-	0.0657	Total TCDF	3.15	-	
1,2,3,4,6,7,8-HpCDF	5.11	-		0.0511	0.0747	Total PeCDF	4.92	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.262		-	0.0883	Total HxCDF	7.74	-	
OCDF	14.0	-		0.00420	0.170	Total HpCDF	13.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	84.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	89.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	93.0	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	90.3	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	83.4	23.0 - 140	
13C-OCDD	66.9	17.0 - 157	
13C-2,3,7,8-TCDF	89.1	24.0 - 169	
13C-1,2,3,7,8-PeCDF	87.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	86.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	86.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	83.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	78.8	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	80.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	83.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	81.9	26.0 - 138	
13C-OCDF	71.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 81.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

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EPA Method 1613 PCDD/F



FAL ID: 11423-004-SA
Client ID: SP1-4
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.02 g
% Solids: 95.60

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.510
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.212		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.388		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.340		-	0.0793				
1,2,3,6,7,8-HxCDD	0.999	-	J	0.0999	0.0940	Total TCDD	ND	0.212	
1,2,3,7,8,9-HxCDD	0.825	-	J	0.0825	0.0823	Total PeCDD	ND	0.388	
1,2,3,4,6,7,8-HpCDD	18.1	-		0.181	0.0842	Total HxCDD	7.26	-	
OCDD	166	-		0.0498	0.172	Total HpCDD	33.2	-	
2,3,7,8-TCDF	0.653	-	J	0.0653	0.0269				
1,2,3,7,8-PeCDF	ND	0.216		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.237		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.439		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.443		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.469		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.517		-	0.0657	Total TCDF	3.18	-	
1,2,3,4,6,7,8-HpCDF	2.92	-	J	0.0292	0.0747	Total PeCDF	2.42	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.236		-	0.0883	Total HxCDF	3.80	-	J
OCDF	6.59	-	J	0.00198	0.170	Total HpCDF	7.14	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	84.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	92.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	90.2	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	87.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	84.0	23.0 - 140	
13C-OCDD	70.6	17.0 - 157	
13C-2,3,7,8-TCDF	88.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	85.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	87.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	83.6	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	81.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	80.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	80.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	82.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	83.4	26.0 - 138	
13C-OCDF	73.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 80.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000010 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-005-SA
Client ID: SP1-5
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.05 g
% Solids: 96.53

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 4.88
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.245		-	0.0273				
1,2,3,7,8-PeCDD	0.682	-	J	0.682	0.0570				
1,2,3,4,7,8-HxCDD	1.44	-	J	0.144	0.0793				
1,2,3,6,7,8-HxCDD	5.76	-		0.576	0.0940	Total TCDD	0.702	-	J
1,2,3,7,8,9-HxCDD	2.92	-	J	0.292	0.0823	Total PeCDD	7.92	-	M
1,2,3,4,6,7,8-HpCDD	152	-		1.52	0.0842	Total HxCDD	43.4	-	
OCDD	1540	-		0.462	0.172	Total HpCDD	280	-	
2,3,7,8-TCDF	2.49	-	F	0.249	0.0269				
1,2,3,7,8-PeCDF	0.683	-	J	0.0205	0.0449				
2,3,4,7,8-PeCDF	1.25	-	J	0.375	0.0468				
1,2,3,4,7,8-HxCDF	1.31	-	J	0.131	0.0437				
1,2,3,6,7,8-HxCDF	1.22	-	J	0.122	0.0417				
2,3,4,6,7,8-HxCDF	1.38	-	J	0.138	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.276		-	0.0657	Total TCDF	15.5	-	
1,2,3,4,6,7,8-HpCDF	13.8	-		0.138	0.0747	Total PeCDF	13.0	-	
1,2,3,4,7,8,9-HpCDF	1.51	-	J	0.0151	0.0883	Total HxCDF	18.9	-	
OCDF	36.7	-		0.0110	0.170	Total HpCDF	39.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	89.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	92.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	94.0	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	89.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	82.3	23.0 - 140	
13C-OCDD	65.1	17.0 - 157	
13C-2,3,7,8-TCDF	93.5	24.0 - 169	
13C-1,2,3,7,8-PeCDF	85.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	87.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	87.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	83.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	80.0	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	79.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	79.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	80.5	26.0 - 138	
13C-OCDF	68.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.1 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

EPA Method 1613 PCDD/F



FAL ID: 11423-006-SA
Client ID: SP1-6
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.02 g
% Solids: 96.16

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 5.00
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.232		-	0.0273				
1,2,3,7,8-PeCDD	1.06	-	J	1.06	0.0570				
1,2,3,4,7,8-HxCDD	1.72	-	J	0.172	0.0793				
1,2,3,6,7,8-HxCDD	4.93	-	J	0.493	0.0940	Total TCDD	ND	0.232	
1,2,3,7,8,9-HxCDD	3.17	-	J	0.317	0.0823	Total PeCDD	2.91	-	J
1,2,3,4,6,7,8-HpCDD	115	-		1.15	0.0842	Total HxCDD	31.8	-	
OCDD	1070	-		0.321	0.172	Total HpCDD	206	-	
2,3,7,8-TCDF	1.35	-		0.135	0.0269				
1,2,3,7,8-PeCDF	ND	0.269		-	0.0449				
2,3,4,7,8-PeCDF	1.38	-	J	0.414	0.0468				
1,2,3,4,7,8-HxCDF	2.27	-	J	0.227	0.0437				
1,2,3,6,7,8-HxCDF	1.47	-	J	0.147	0.0417				
2,3,4,6,7,8-HxCDF	1.81	-	J	0.181	0.0574				
1,2,3,7,8,9-HxCDF	0.677	-	J	0.0677	0.0657	Total TCDF	7.40	-	
1,2,3,4,6,7,8-HpCDF	27.9	-		0.279	0.0747	Total PeCDF	13.2	-	
1,2,3,4,7,8,9-HpCDF	1.81	-	J	0.0181	0.0883	Total HxCDF	39.9	-	
OCDF	65.4	-		0.0196	0.170	Total HpCDF	78.0	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	89.9	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	94.1	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	92.4	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.5	23.0 - 140	
13C-OCDD	78.6	17.0 - 157	
13C-2,3,7,8-TCDF	93.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	86.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	87.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	88.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	85.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	85.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	84.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.6	26.0 - 138	
13C-OCDF	77.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000012 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-007-SA
Client ID: SP1-7
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.05 g
% Solids: 95.81

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 1.65
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.251		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.356		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.431		-	0.0793				
1,2,3,6,7,8-HxCDD	2.12	-	J	0.212	0.0940	Total TCDD	ND	0.251	
1,2,3,7,8,9-HxCDD	1.44	-	J	0.144	0.0823	Total PeCDD	0.923	-	J
1,2,3,4,6,7,8-HpCDD	47.6	-		0.476	0.0842	Total HxCDD	14.3	-	
OCDD	398	-		0.119	0.172	Total HpCDD	84.9	-	
2,3,7,8-TCDF	1.43	-		0.143	0.0269				
1,2,3,7,8-PeCDF	ND	0.372		-	0.0449				
2,3,4,7,8-PeCDF	0.725	-	J	0.218	0.0468				
1,2,3,4,7,8-HxCDF	0.829	-	J	0.0829	0.0437				
1,2,3,6,7,8-HxCDF	0.626	-	J	0.0626	0.0417				
2,3,4,6,7,8-HxCDF	0.703	-	J	0.0703	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.344		-	0.0657	Total TCDF	6.16	-	
1,2,3,4,6,7,8-HpCDF	10.3	-		0.103	0.0747	Total PeCDF	5.88	-	
1,2,3,4,7,8,9-HpCDF	0.832	-	J	0.00832	0.0883	Total HxCDF	12.9	-	
OCDF	33.0	-		0.00990	0.170	Total HpCDF	30.7	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	64.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	67.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	70.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	69.0	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	66.8	23.0 - 140	
13C-OCDD	59.2	17.0 - 157	
13C-2,3,7,8-TCDF	68.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	66.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	63.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	67.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	64.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	63.4	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	64.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	64.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	65.7	26.0 - 138	
13C-OCDF	60.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 60.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000013 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-008-SA
Client ID: SP1-8
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.02 g
% Solids: 94.62

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 2.93
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.255		-	0.0273				
1,2,3,7,8-PeCDD	0.953	-	J	0.953	0.0570				
1,2,3,4,7,8-HxCDD	0.934	-	J	0.0934	0.0793				
1,2,3,6,7,8-HxCDD	2.85	-	J	0.285	0.0940	Total TCDD	ND	0.255	
1,2,3,7,8,9-HxCDD	1.73	-	J	0.173	0.0823	Total PeCDD	4.11	-	J
1,2,3,4,6,7,8-HpCDD	54.9	-		0.549	0.0842	Total HxCDD	24.4	-	
OCDD	453	-		0.136	0.172	Total HpCDD	105	-	
2,3,7,8-TCDF	1.73	-		0.173	0.0269				
1,2,3,7,8-PeCDF	ND	0.387		-	0.0449				
2,3,4,7,8-PeCDF	0.736	-	J	0.221	0.0468				
1,2,3,4,7,8-HxCDF	0.886	-	J	0.0886	0.0437				
1,2,3,6,7,8-HxCDF	0.699	-	J	0.0699	0.0417				
2,3,4,6,7,8-HxCDF	0.803	-	J	0.0803	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.281		-	0.0657	Total TCDF	9.10	-	
1,2,3,4,6,7,8-HpCDF	9.38	-		0.0938	0.0747	Total PeCDF	9.08	-	
1,2,3,4,7,8,9-HpCDF	0.802	-	J	0.00802	0.0883	Total HxCDF	13.2	-	
OCDF	22.1	-		0.00663	0.170	Total HpCDF	24.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	79.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	81.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	85.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	79.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	76.9	23.0 - 140	
13C-OCDD	60.1	17.0 - 157	
13C-2,3,7,8-TCDF	78.5	24.0 - 169	
13C-1,2,3,7,8-PeCDF	78.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	76.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	80.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	77.7	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	71.9	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	72.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	76.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	75.2	26.0 - 138	
13C-OCDF	63.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 75.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000014 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-009-SA
Client ID: SP1-9
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.03 g
% Solids: 93.92

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.139
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.223		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.339		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.420		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.431		-	0.0940	Total TCDD	ND	0.223	
1,2,3,7,8,9-HxCDD	ND	0.390		-	0.0823	Total PeCDD	ND	0.339	
1,2,3,4,6,7,8-HpCDD	5.47	-		0.0547	0.0842	Total HxCDD	1.72	-	J
OCDD	44.1	-		0.0132	0.172	Total HpCDD	10.7	-	
2,3,7,8-TCDF	0.560	-	J	0.0560	0.0269				
1,2,3,7,8-PeCDF	ND	0.326		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.319		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.349		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.357		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.400		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.430		-	0.0657	Total TCDF	1.15	-	
1,2,3,4,6,7,8-HpCDF	1.48	-	J	0.0148	0.0747	Total PeCDF	0.564	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.308		-	0.0883	Total HxCDF	0.710	-	J
OCDF	2.56	-	J	0.000768	0.170	Total HpCDF	3.02	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	88.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	93.2	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	92.2	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	90.3	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	86.3	23.0 - 140	
13C-OCDD	76.3	17.0 - 157	
13C-2,3,7,8-TCDF	93.2	24.0 - 169	
13C-1,2,3,7,8-PeCDF	89.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	91.8	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	86.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	82.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	81.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	83.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	82.9	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.8	26.0 - 138	
13C-OCDF	75.6	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

000015 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-010-SA
Client ID: SP1-10
Matrix: Solid
Batch No: X4473

Date Extracted: 04-16-2018
Date Received: 04-04-2018
Amount: 5.06 g
% Solids: 93.98

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 1.23
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.237		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.493		-	0.0570				
1,2,3,4,7,8-HxCDD	0.700	-	J	0.0700	0.0793				
1,2,3,6,7,8-HxCDD	1.71	-	J	0.171	0.0940	Total TCDD	ND	0.237	
1,2,3,7,8,9-HxCDD	1.34	-	J	0.134	0.0823	Total PeCDD	ND	0.493	
1,2,3,4,6,7,8-HpCDD	35.6	-		0.356	0.0842	Total HxCDD	13.4	-	
OCDD	285	-		0.0855	0.172	Total HpCDD	63.3	-	
2,3,7,8-TCDF	1.16	-		0.116	0.0269				
1,2,3,7,8-PeCDF	ND	0.336		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.363		-	0.0468				
1,2,3,4,7,8-HxCDF	0.791	-	J	0.0791	0.0437				
1,2,3,6,7,8-HxCDF	0.644	-	J	0.0644	0.0417				
2,3,4,6,7,8-HxCDF	0.736	-	J	0.0736	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.356		-	0.0657	Total TCDF	5.59	-	
1,2,3,4,6,7,8-HpCDF	7.49	-		0.0749	0.0747	Total PeCDF	4.71	-	J,M
1,2,3,4,7,8,9-HpCDF	ND	0.239		-	0.0883	Total HxCDF	10.9	-	
OCDF	17.5	-		0.00525	0.170	Total HpCDF	19.8	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	91.9	25.0 - 164	
13C-1,2,3,7,8-PeCDD	95.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	96.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.1	23.0 - 140	
13C-OCDD	80.8	17.0 - 157	
13C-2,3,7,8-TCDF	93.2	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	91.8	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	86.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.1	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	87.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	89.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	91.6	26.0 - 138	
13C-OCDF	81.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 89.8 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/18/2018

Reviewed By: 

Date: 4/18/2018

EPA Method 1613 PCDD/F



FAL ID: 11423-011-SA
Client ID: SP1-11
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.01 g
% Solids: 92.86

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 2.94
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.192		-	0.0273				
1,2,3,7,8-PeCDD	0.565	-	J	0.565	0.0570				
1,2,3,4,7,8-HxCDD	0.956	-	J	0.0956	0.0793				
1,2,3,6,7,8-HxCDD	2.57	-	J	0.257	0.0940	Total TCDD	ND	0.192	
1,2,3,7,8,9-HxCDD	1.68	-	J	0.168	0.0823	Total PeCDD	3.01	-	J
1,2,3,4,6,7,8-HpCDD	61.1	-		0.611	0.0842	Total HxCDD	19.0	-	
OCDD	556	-		0.167	0.172	Total HpCDD	114	-	
2,3,7,8-TCDF	1.58	-		0.158	0.0269				
1,2,3,7,8-PeCDF	0.472	-	J	0.0142	0.0449				
2,3,4,7,8-PeCDF	1.11	-	J	0.333	0.0468				
1,2,3,4,7,8-HxCDF	1.44	-	J	0.144	0.0437				
1,2,3,6,7,8-HxCDF	1.04	-	J	0.104	0.0417				
2,3,4,6,7,8-HxCDF	1.03	-	J	0.103	0.0574				
1,2,3,7,8,9-HxCDF	0.694	-	J	0.0694	0.0657	Total TCDF	9.88	-	
1,2,3,4,6,7,8-HpCDF	13.5	-		0.135	0.0747	Total PeCDF	11.5	-	
1,2,3,4,7,8,9-HpCDF	1.14	-	J	0.0114	0.0883	Total HxCDF	18.6	-	
OCDF	29.9	-		0.00897	0.170	Total HpCDF	35.7	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	96.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	100	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	105	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	99.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	93.6	23.0 - 140	
13C-OCDD	76.3	17.0 - 157	
13C-2,3,7,8-TCDF	99.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	98.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	99.3	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	95.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	93.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	91.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	91.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	91.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	92.2	26.0 - 138	
13C-OCDF	77.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 89.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

EPA Method 1613 PCDD/F



FAL ID: 11423-012-SA
Client ID: SP1-12
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.05 g
% Solids: 96.77

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 1.08
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.174		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.318		-	0.0570				
1,2,3,4,7,8-HxCDD	0.494	-	J	0.0494	0.0793				
1,2,3,6,7,8-HxCDD	1.73	-	J	0.173	0.0940	Total TCDD	ND	0.174	
1,2,3,7,8,9-HxCDD	1.06	-	J	0.106	0.0823	Total PeCDD	ND	0.318	
1,2,3,4,6,7,8-HpCDD	33.3	-		0.333	0.0842	Total HxCDD	11.1	-	
OCDD	302	-		0.0906	0.172	Total HpCDD	61.2	-	
2,3,7,8-TCDF	0.699	-	J	0.0699	0.0269				
1,2,3,7,8-PeCDF	ND	0.200		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.205		-	0.0468				
1,2,3,4,7,8-HxCDF	0.621	-	J	0.0621	0.0437				
1,2,3,6,7,8-HxCDF	0.522	-	J	0.0522	0.0417				
2,3,4,6,7,8-HxCDF	0.673	-	J	0.0673	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.144		-	0.0657	Total TCDF	3.68	-	
1,2,3,4,6,7,8-HpCDF	6.91	-		0.0691	0.0747	Total PeCDF	3.76	-	J
1,2,3,4,7,8,9-HpCDF	0.576	-	J	0.00576	0.0883	Total HxCDF	10.0	-	
OCDF	21.2	-		0.00636	0.170	Total HpCDF	19.4	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	95.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	102	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	99.8	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	98.0	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	90.0	23.0 - 140	
13C-OCDD	73.3	17.0 - 157	
13C-2,3,7,8-TCDF	96.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	97.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	97.8	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.9	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	86.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	84.6	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	86.9	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	90.6	26.0 - 138	
13C-OCDF	76.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000018 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-013-SA
Client ID: SP1-13
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.05 g
% Solids: 94.87

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 2.89
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.170		-	0.0273				
1,2,3,7,8-PeCDD	0.663	-	J	0.663	0.0570				
1,2,3,4,7,8-HxCDD	0.877	-	J	0.0877	0.0793				
1,2,3,6,7,8-HxCDD	2.60	-	J	0.260	0.0940	Total TCDD	ND	0.170	
1,2,3,7,8,9-HxCDD	1.45	-	J	0.145	0.0823	Total PeCDD	2.30	-	J
1,2,3,4,6,7,8-HpCDD	71.3	-		0.713	0.0842	Total HxCDD	18.8	-	
OCDD	629	-		0.189	0.172	Total HpCDD	128	-	
2,3,7,8-TCDF	1.22	-		0.122	0.0269				
1,2,3,7,8-PeCDF	0.397	-	J	0.0119	0.0449				
2,3,4,7,8-PeCDF	0.927	-	J	0.278	0.0468				
1,2,3,4,7,8-HxCDF	0.913	-	J	0.0913	0.0437				
1,2,3,6,7,8-HxCDF	0.845	-	J	0.0845	0.0417				
2,3,4,6,7,8-HxCDF	1.03	-	J	0.103	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.233		-	0.0657	Total TCDF	6.98	-	
1,2,3,4,6,7,8-HpCDF	12.5	-		0.125	0.0747	Total PeCDF	8.44	-	
1,2,3,4,7,8,9-HpCDF	0.901	-	J	0.00901	0.0883	Total HxCDF	16.6	-	
OCDF	37.1	-		0.0111	0.170	Total HpCDF	36.9	-	
Internal Standards	% Rec	QC Limits	Qual						
13C-2,3,7,8-TCDD	96.7	25.0 - 164							
13C-1,2,3,7,8-PeCDD	99.1	25.0 - 181							
13C-1,2,3,4,7,8-HxCDD	100	32.0 - 141							
13C-1,2,3,6,7,8-HxCDD	99.6	28.0 - 130							
13C-1,2,3,4,6,7,8-HpCDD	87.4	23.0 - 140							
13C-OCDD	75.2	17.0 - 157							
13C-2,3,7,8-TCDF	98.0	24.0 - 169							
13C-1,2,3,7,8-PeCDF	97.5	24.0 - 185							
13C-2,3,4,7,8-PeCDF	96.5	21.0 - 178							
13C-1,2,3,4,7,8-HxCDF	93.8	26.0 - 152							
13C-1,2,3,6,7,8-HxCDF	90.9	26.0 - 123							
13C-2,3,4,6,7,8-HxCDF	88.3	28.0 - 136							
13C-1,2,3,7,8,9-HxCDF	85.6	29.0 - 147							
13C-1,2,3,4,6,7,8-HpCDF	88.6	28.0 - 143							
13C-1,2,3,4,7,8,9-HpCDF	89.4	26.0 - 138							
13C-OCDF	76.8	17.0 - 157							
Cleanup Surrogate									
37Cl-2,3,7,8-TCDD	91.8	35.0 - 197							

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/19/2018

Reviewed By: 
Date: 4/19/2018

EPA Method 1613 PCDD/F



FAL ID: 11423-014-SA
Client ID: SP1-14
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.00 g
% Solids: 94.79

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.316
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.179		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.251		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.261		-	0.0793				
1,2,3,6,7,8-HxCDD	0.750	-	J	0.0750	0.0940	Total TCDD	ND	0.179	
1,2,3,7,8,9-HxCDD	ND	0.240		-	0.0823	Total PeCDD	ND	0.251	
1,2,3,4,6,7,8-HpCDD	14.3	-		0.143	0.0842	Total HxCDD	4.09	-	J
OCDD	147	-		0.0441	0.172	Total HpCDD	26.0	-	
2,3,7,8-TCDF	0.308	-	J	0.0308	0.0269				
1,2,3,7,8-PeCDF	ND	0.254		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.247		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.201		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.196		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.226		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.249		-	0.0657	Total TCDF	0.308	-	J
1,2,3,4,6,7,8-HpCDF	2.17	-	J	0.0217	0.0747	Total PeCDF	0.496	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.213		-	0.0883	Total HxCDF	1.97	-	J
OCDF	5.61	-	J	0.00168	0.170	Total HpCDF	5.67	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	92.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	97.8	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	96.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.0	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	85.6	23.0 - 140	
13C-OCDD	62.3	17.0 - 157	
13C-2,3,7,8-TCDF	95.2	24.0 - 169	
13C-1,2,3,7,8-PeCDF	96.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	97.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	87.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	83.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	81.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	81.5	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	83.5	26.0 - 138	
13C-OCDF	68.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 90.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000020 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-015-SA
Client ID: SP1-15
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.02 g
% Solids: 94.28

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 3.20
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.179		-	0.0273				
1,2,3,7,8-PeCDD	0.753	-	J	0.753	0.0570				
1,2,3,4,7,8-HxCDD	0.897	-	J	0.0897	0.0793				
1,2,3,6,7,8-HxCDD	2.91	-	J	0.291	0.0940	Total TCDD	0.467	-	J
1,2,3,7,8,9-HxCDD	1.93	-	J	0.193	0.0823	Total PeCDD	4.35	-	J
1,2,3,4,6,7,8-HpCDD	67.2	-		0.672	0.0842	Total HxCDD	22.6	-	
OCDD	546	-		0.164	0.172	Total HpCDD	122	-	
2,3,7,8-TCDF	1.78	-		0.178	0.0269				
1,2,3,7,8-PeCDF	0.368	-	J	0.0110	0.0449				
2,3,4,7,8-PeCDF	1.11	-	J	0.333	0.0468				
1,2,3,4,7,8-HxCDF	1.25	-	J	0.125	0.0437				
1,2,3,6,7,8-HxCDF	0.872	-	J	0.0872	0.0417				
2,3,4,6,7,8-HxCDF	1.02	-	J	0.102	0.0574				
1,2,3,7,8,9-HxCDF	0.450	-	J	0.0450	0.0657	Total TCDF	10.7	-	
1,2,3,4,6,7,8-HpCDF	14.0	-		0.140	0.0747	Total PeCDF	11.1	-	
1,2,3,4,7,8,9-HpCDF	0.992	-	J	0.00992	0.0883	Total HxCDF	19.9	-	
OCDF	33.5	-		0.0101	0.170	Total HpCDF	37.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	96.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	99.7	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	87.6	23.0 - 140	
13C-OCDD	75.9	17.0 - 157	
13C-2,3,7,8-TCDF	97.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	96.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.5	26.0 - 138	
13C-OCDF	76.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 92.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000021 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-016-SA
Client ID: SP1-16
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.00 g
% Solids: 94.97

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 1.97
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.180		-	0.0273				
1,2,3,7,8-PeCDD	0.556	-	J	0.556	0.0570				
1,2,3,4,7,8-HxCDD	0.701	-	J	0.0701	0.0793				
1,2,3,6,7,8-HxCDD	1.64	-	J	0.164	0.0940	Total TCDD	ND	0.180	
1,2,3,7,8,9-HxCDD	1.48	-	J	0.148	0.0823	Total PeCDD	1.78	-	J
1,2,3,4,6,7,8-HpCDD	34.8	-		0.348	0.0842	Total HxCDD	12.6	-	
OCDD	269	-		0.0807	0.172	Total HpCDD	62.9	-	
2,3,7,8-TCDF	0.863	-	J	0.0863	0.0269				
1,2,3,7,8-PeCDF	ND	0.209		-	0.0449				
2,3,4,7,8-PeCDF	0.519	-	J	0.156	0.0468				
1,2,3,4,7,8-HxCDF	0.552	-	J	0.0552	0.0437				
1,2,3,6,7,8-HxCDF	0.830	-	J	0.0830	0.0417				
2,3,4,6,7,8-HxCDF	1.06	-	J	0.106	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.219		-	0.0657	Total TCDF	4.02	-	
1,2,3,4,6,7,8-HpCDF	10.6	-		0.106	0.0747	Total PeCDF	5.39	-	
1,2,3,4,7,8,9-HpCDF	0.758	-	J	0.00758	0.0883	Total HxCDF	14.0	-	
OCDF	20.2	-		0.00606	0.170	Total HpCDF	23.7	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.3	25.0 - 164	
13C-1,2,3,7,8-PeCDD	94.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	101	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	98.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	88.4	23.0 - 140	
13C-OCDD	71.2	17.0 - 157	
13C-2,3,7,8-TCDF	95.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	93.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	94.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	93.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	85.4	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	86.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.8	26.0 - 138	
13C-OCDF	74.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.2 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000022 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-017-SA
Client ID: SP1-17
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.04 g
% Solids: 95.05

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.429
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.179		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.217		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.240		-	0.0793				
1,2,3,6,7,8-HxCDD	1.07	-	J	0.107	0.0940	Total TCDD	ND	0.179	
1,2,3,7,8,9-HxCDD	0.689	-	J	0.0689	0.0823	Total PeCDD	ND	0.217	
1,2,3,4,6,7,8-HpCDD	16.0	-		0.160	0.0842	Total HxCDD	6.81	-	
OCDD	131	-		0.0393	0.172	Total HpCDD	29.2	-	
2,3,7,8-TCDF	0.274	-	J	0.0274	0.0269				
1,2,3,7,8-PeCDF	ND	0.166		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.176		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.211		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.219		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.232		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.253		-	0.0657	Total TCDF	0.932	-	J,M
1,2,3,4,6,7,8-HpCDF	2.48	-	J	0.0248	0.0747	Total PeCDF	1.71	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.224		-	0.0883	Total HxCDF	3.08	-	J
OCDF	6.08	-	J	0.00182	0.170	Total HpCDF	6.12	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	97.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	98.8	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	97.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	88.8	23.0 - 140	
13C-OCDD	73.4	17.0 - 157	
13C-2,3,7,8-TCDF	95.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	93.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	94.7	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	92.3	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	87.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.7	26.0 - 138	
13C-OCDF	76.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000023 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-018-SA
Client ID: SP1-18
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.00 g
% Solids: 93.52

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-18-2018
2005 WHO TEQ: 0.891
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.171		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.292		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.272		-	0.0793				
1,2,3,6,7,8-HxCDD	0.832	-	J	0.0832	0.0940	Total TCDD	ND	0.171	
1,2,3,7,8,9-HxCDD	0.762	-	J	0.0762	0.0823	Total PeCDD	0.581	-	J
1,2,3,4,6,7,8-HpCDD	11.5	-		0.115	0.0842	Total HxCDD	6.39	-	
OCDD	89.3	-		0.0268	0.172	Total HpCDD	22.0	-	
2,3,7,8-TCDF	1.52	-		0.152	0.0269				
1,2,3,7,8-PeCDF	ND	0.203		-	0.0449				
2,3,4,7,8-PeCDF	0.821	-	J	0.246	0.0468				
1,2,3,4,7,8-HxCDF	0.671	-	J	0.0671	0.0437				
1,2,3,6,7,8-HxCDF	0.451	-	J	0.0451	0.0417				
2,3,4,6,7,8-HxCDF	0.489	-	J	0.0489	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.208		-	0.0657	Total TCDF	5.49	-	
1,2,3,4,6,7,8-HpCDF	2.91	-	J	0.0291	0.0747	Total PeCDF	5.82	-	
1,2,3,4,7,8,9-HpCDF	ND	0.321		-	0.0883	Total HxCDF	5.46	-	
OCDF	4.76	-	J	0.00143	0.170	Total HpCDF	6.72	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	98.4	25.0 - 164	
13C-1,2,3,7,8-PeCDD	101	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	101	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	101	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	91.5	23.0 - 140	
13C-OCDD	74.2	17.0 - 157	
13C-2,3,7,8-TCDF	100	24.0 - 169	
13C-1,2,3,7,8-PeCDF	97.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	97.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	95.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	91.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.9	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	89.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	91.3	26.0 - 138	
13C-OCDF	79.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 92.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000024 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-019-SA
Client ID: SP1-19
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.01 g
% Solids: 95.10

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-19-2018
2005 WHO TEQ: 2.84
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.197		-	0.0273				
1,2,3,7,8-PeCDD	0.774	-	J	0.774	0.0570				
1,2,3,4,7,8-HxCDD	0.932	-	J	0.0932	0.0793				
1,2,3,6,7,8-HxCDD	2.78	-	J	0.278	0.0940	Total TCDD	ND	0.197	
1,2,3,7,8,9-HxCDD	1.92	-	J	0.192	0.0823	Total PeCDD	3.60	-	J
1,2,3,4,6,7,8-HpCDD	58.5	-		0.585	0.0842	Total HxCDD	20.3	-	
OCDD	496	-		0.149	0.172	Total HpCDD	103	-	
2,3,7,8-TCDF	0.921	-	J	0.0921	0.0269				
1,2,3,7,8-PeCDF	0.364	-	J	0.0109	0.0449				
2,3,4,7,8-PeCDF	0.556	-	J	0.167	0.0468				
1,2,3,4,7,8-HxCDF	0.901	-	J	0.0901	0.0437				
1,2,3,6,7,8-HxCDF	1.06	-	J	0.106	0.0417				
2,3,4,6,7,8-HxCDF	1.31	-	J	0.131	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.248		-	0.0657	Total TCDF	4.61	-	
1,2,3,4,6,7,8-HpCDF	15.3	-		0.153	0.0747	Total PeCDF	8.55	-	
1,2,3,4,7,8,9-HpCDF	0.952	-	J	0.00952	0.0883	Total HxCDF	20.6	-	
OCDF	32.8	-		0.00984	0.170	Total HpCDF	37.7	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	95.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	93.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	98.7	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	97.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	88.3	23.0 - 140	
13C-OCDD	71.8	17.0 - 157	
13C-2,3,7,8-TCDF	98.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	95.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	87.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	85.8	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.3	26.0 - 138	
13C-OCDF	75.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000025 of 000033

EPA Method 1613 PCDD/F



FAL ID: 11423-020-SA
Client ID: SP1-20
Matrix: Solid
Batch No: X4475

Date Extracted: 04-17-2018
Date Received: 04-04-2018
Amount: 5.03 g
% Solids: 93.51

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-19-2018
2005 WHO TEQ: 3.62
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.210		-	0.0273				
1,2,3,7,8-PeCDD	0.856	-	J	0.856	0.0570				
1,2,3,4,7,8-HxCDD	1.19	-	J	0.119	0.0793				
1,2,3,6,7,8-HxCDD	3.04	-	J	0.304	0.0940	Total TCDD	0.546	-	J
1,2,3,7,8,9-HxCDD	2.00	-	J	0.200	0.0823	Total PeCDD	4.16	-	J
1,2,3,4,6,7,8-HpCDD	55.7	-		0.557	0.0842	Total HxCDD	20.1	-	
OCDD	455	-		0.137	0.172	Total HpCDD	100	-	
2,3,7,8-TCDF	2.58	-	F	0.258	0.0269				
1,2,3,7,8-PeCDF	0.736	-	J	0.0221	0.0449				
2,3,4,7,8-PeCDF	1.66	-	J	0.498	0.0468				
1,2,3,4,7,8-HxCDF	1.88	-	J	0.188	0.0437				
1,2,3,6,7,8-HxCDF	1.12	-	J	0.112	0.0417				
2,3,4,6,7,8-HxCDF	1.45	-	J	0.145	0.0574				
1,2,3,7,8,9-HxCDF	0.856	-	J	0.0856	0.0657	Total TCDF	14.6	-	
1,2,3,4,6,7,8-HpCDF	11.3	-		0.113	0.0747	Total PeCDF	13.2	-	
1,2,3,4,7,8,9-HpCDF	1.60	-	J	0.0160	0.0883	Total HxCDF	18.4	-	
OCDF	27.4	-		0.00822	0.170	Total HpCDF	32.2	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	97.5	25.0 - 164	
13C-1,2,3,7,8-PeCDD	101	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	104	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	99.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	93.2	23.0 - 140	
13C-OCDD	76.8	17.0 - 157	
13C-2,3,7,8-TCDF	102	24.0 - 169	
13C-1,2,3,7,8-PeCDF	96.9	24.0 - 185	
13C-2,3,4,7,8-PeCDF	98.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	93.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	91.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.2	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	89.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	91.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	93.5	26.0 - 138	
13C-OCDF	81.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 89.8 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/19/2018

Reviewed By: 

Date: 4/19/2018

000026 of 000033



Frontier Analytical Laboratory
5172 Hillsdale Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
Fax: 916-934-0999

FAL USE ONLY

Laboratory Project No.: 11423

Temperature: 0 °C

Chain of Custody

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Please Print in Pen Page 1 of 2

CLIENT INFORMATION					INVOICE INFORMATION (if different from client info)										PROJECT INFORMATION							
Company Name: <u>Brusca Associates, Inc.</u>					Company Name: <u>Same</u>					FAL Quote #: <u>3310</u>												
Contact Name: <u>Joe Brusca</u>					Contact Name: _____					P.O. #: _____												
Address: <u>PO Box 332, Roseville, CA 95661</u>					Address: _____					Project #: <u>137-002</u>												
Phone: <u>916-677-1470</u> Fax: _____					Phone: _____ Fax: _____					Project Name: <u>LATHROP 43-AC PROPERTY</u>												
Email: <u>jbrusca@bruscaassociates.com</u>					Email: _____					TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (✓ one)												
										* FAL must agree with price and RUSH TAT in writing.												
REPORT INFORMATION					REPORT DISTRIBUTION (email only is preferred)										ADDITIONAL INSTRUCTIONS							
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☐ California State Drinking Water Form																						
System #: _____ Source #: _____																						
Sampler: _____ Employer: _____																						
Sample ID					Date	Time	Matrix	# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS		**TEQ	
					Collected														☐ 2,3,7,8-TCDD only	☐ 1998 WHO		
																			☐ 2,3,7,8-TCDD/F only	☒ 2005 WHO		
																			☒ PCDD/F (Cl ₄ -Cl ₈)	☐ Other		
Remarks																						
1	SP1-1				4/4/13	11:59	SOLID		X													
2	SP1-2					12:01		X														
3	SP1-3					12:03		X														
4	SP1-4					12:05		X														
5	SP1-5					12:12		X														
6	SP1-6					12:13		X														
7	SP1-7					12:17		X														
8	SP1-8					12:22		X														
9	SP1-9					12:25		X														
10	SP1-10					12:31		X														
11	SP1-11					12:35		X														
12	SP1-12					12:37		X														
13	SP1-13					12:42		X														
14	SP1-14					12:45		X														
15	SP1-15					12:47		X														
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																						
Relinquished by: (Signature and Printed Name)					Date	Time	Received by: (Signature and Printed Name)					Date	Time									
					4/4/13	15:13						04/04/13	1513									

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

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Pink Copy - Originator



Frontier Analytical Laboratory
5172 Hillside Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
Fax: 916-934-0999

FAL USE ONLY

Laboratory Project No.: 11423
Temperature: 0 °C

Chain of Custody

www.frontieranalytical.com

Please Print in Pen Page 2 of 2

CLIENT INFORMATION					INVOICE INFORMATION (if different from client info)										PROJECT INFORMATION										
Company Name: <u>Brusca Associates, Inc.</u> Contact Name: <u>Joe Brusca</u> Address: <u>PO Box 332, Roseville, CA 95661</u> Phone: <u>916-677-1470</u> Fax: <u></u> Email: <u>jbrusca@bruscaassociates.com</u>					Company Name: <u>Same</u> Contact Name: <u></u> Address: <u></u> Phone: <u></u> Fax: <u></u> Email: <u></u>										FAL Quote #: <u>3310</u> P.O. #: <u></u> Project #: <u>137-002</u> Project Name: <u>LATHROP 48-AC PROPERTY</u> TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (√ one) * FAL must agree with price and RUSH TAT in writing.										
REPORT INFORMATION					REPORT DISTRIBUTION (email only is preferred)										ADDITIONAL INSTRUCTIONS										
Report Level: ☐ I/II ☐ III ☐ IV ☐ EDD: ☐ FAL Basic ☐ Geotracker ☐ Other: <u></u> ☐ Custom: Contact FAL ☐ California State Drinking Water Form System #: <u></u> Source #: <u></u> Sampler: <u></u> Employer: <u></u>					☐ Hardcopy ☐ CD (.pdf including EDDs if requested) ☐ Email (.pdf including EDDs if requested)																				
					# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/9A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS					**TEQ				
																	☐ 2,3,7,8-TCDD only	☐ 1998 WHO	☐ 2,3,7,8-TCDD/F only	☒ 2005 WHO	☐ Other				
																	Remarks								
1	SPL-16		4/4/13	12:51	SOLD		X																		
2	SPL-17			12:53		X																			
3	SPL-18			12:56		X																			
4	SPL-19			13:05		X																			
5	SPL-20			13:08		X																			
6																									
7																									
8																									
9																									
10																									
11																									
12																									
13																									
14																									
15																									
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																									
Relinquished by: (Signature and Printed Name)					Date	Time	Received by: (Signature and Printed Name)					Date	Time												
					4/4/13	15:13	Kathy Zipp K. Zipp					04/04/2013	1513												

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

White Copy - Report

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Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11423**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/04/2018
Time Received:	03:13 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	20
Duplicates:	0
Storage Location:	R-4

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

12 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48 Acre Property

Enclosed are the results of analyses for samples received by the laboratory on 04/05/18 09:55. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP1-1	T181153-01	Soil	04/04/18 11:59	04/05/18 09:55
SP1-2	T181153-02	Soil	04/04/18 12:01	04/05/18 09:55
SP1-3	T181153-03	Soil	04/04/18 12:03	04/05/18 09:55
SP1-4	T181153-04	Soil	04/04/18 12:05	04/05/18 09:55
SP1-5	T181153-05	Soil	04/04/18 12:12	04/05/18 09:55
SP1-6	T181153-06	Soil	04/04/18 12:13	04/05/18 09:55
SP1-7	T181153-07	Soil	04/04/18 12:17	04/05/18 09:55
SP1-8	T181153-08	Soil	04/04/18 12:22	04/05/18 09:55
SP1-9	T181153-09	Soil	04/04/18 12:25	04/05/18 09:55
SP1-10	T181153-10	Soil	04/04/18 12:31	04/05/18 09:55
SP1-11	T181153-11	Soil	04/04/18 12:35	04/05/18 09:55
SP1-12	T181153-12	Soil	04/04/18 12:37	04/05/18 09:55
SP1-13	T181153-13	Soil	04/04/18 12:42	04/05/18 09:55
SP1-14	T181153-14	Soil	04/04/18 12:45	04/05/18 09:55
SP1-15	T181153-15	Soil	04/04/18 12:47	04/05/18 09:55
SP1-16	T181153-16	Soil	04/04/18 12:51	04/05/18 09:55
SP1-17	T181153-17	Soil	04/04/18 12:53	04/05/18 09:55
SP1-18	T181153-18	Soil	04/04/18 12:56	04/05/18 09:55
SP1-19	T181153-19	Soil	04/04/18 13:05	04/05/18 09:55
SP1-20	T181153-20	Soil	04/04/18 13:08	04/05/18 09:55

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

DETECTIONS SUMMARY

Sample ID: SP1-1

Laboratory ID: T181153-01

Reporting

Analyte	Result	Limit	Units	Method	Notes
C13-C28 (DRO)	47	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	100	10	mg/kg	EPA 8015B	
Barium	54	1.0	mg/kg	EPA 6010B	
Chromium	5.4	2.0	mg/kg	EPA 6010B	
Cobalt	3.7	2.0	mg/kg	EPA 6010B	
Copper	7.8	1.0	mg/kg	EPA 6010B	
Lead	22	3.0	mg/kg	EPA 6010B	
Nickel	5.1	2.0	mg/kg	EPA 6010B	
Vanadium	12	5.0	mg/kg	EPA 6010B	
Zinc	36	1.0	mg/kg	EPA 6010B	
PCB-1254	35	10	ug/kg	EPA 8082	
Pyrene	2400	300	ug/kg	EPA 8270C	
Benzo (a) anthracene	900	300	ug/kg	EPA 8270C	
Benzo (b) fluoranthene	730	300	ug/kg	EPA 8270C	
Benzo (k) fluoranthene	410	300	ug/kg	EPA 8270C	
Benzo (a) pyrene	360	300	ug/kg	EPA 8270C	
Chrysene	1100	300	ug/kg	EPA 8270C	
Fluoranthene	2100	300	ug/kg	EPA 8270C	
Indeno (1,2,3-cd) pyrene	340	300	ug/kg	EPA 8270C	
Phenanthrene	1300	300	ug/kg	EPA 8270C	

Sample ID: SP1-2

Laboratory ID: T181153-02

Reporting

Analyte	Result	Limit	Units	Method	Notes
C13-C28 (DRO)	13	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	86	10	mg/kg	EPA 8015B	
Barium	65	1.0	mg/kg	EPA 6010B	
Chromium	5.4	2.0	mg/kg	EPA 6010B	
Cobalt	4.0	2.0	mg/kg	EPA 6010B	
Copper	3.7	1.0	mg/kg	EPA 6010B	
Nickel	5.4	2.0	mg/kg	EPA 6010B	
Vanadium	16	5.0	mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-2

Laboratory ID: T181153-02

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Zinc	29	1.0	mg/kg	EPA 6010B	

Sample ID: SP1-3

Laboratory ID: T181153-03

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	47	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	68	10	mg/kg	EPA 8015B	
Barium	61	1.0	mg/kg	EPA 6010B	
Chromium	5.3	2.0	mg/kg	EPA 6010B	
Cobalt	3.9	2.0	mg/kg	EPA 6010B	
Copper	4.0	1.0	mg/kg	EPA 6010B	
Nickel	5.1	2.0	mg/kg	EPA 6010B	
Vanadium	16	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	
PCB-1254	25	10	ug/kg	EPA 8082	
Pyrene	300	300	ug/kg	EPA 8270C	

Sample ID: SP1-4

Laboratory ID: T181153-04

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	24	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	75	10	mg/kg	EPA 8015B	
Barium	69	1.0	mg/kg	EPA 6010B	
Chromium	5.3	2.0	mg/kg	EPA 6010B	
Cobalt	4.6	2.0	mg/kg	EPA 6010B	
Copper	4.2	1.0	mg/kg	EPA 6010B	
Lead	8.4	3.0	mg/kg	EPA 6010B	
Nickel	5.9	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	32	1.0	mg/kg	EPA 6010B	
PCB-1254	13	10	ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-5

Laboratory ID: T181153-05

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	96	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	52	10	mg/kg	EPA 8015B	
Barium	41	1.0	mg/kg	EPA 6010B	
Chromium	4.6	2.0	mg/kg	EPA 6010B	
Cobalt	3.0	2.0	mg/kg	EPA 6010B	
Copper	11	1.0	mg/kg	EPA 6010B	
Lead	45	3.0	mg/kg	EPA 6010B	
Nickel	3.9	2.0	mg/kg	EPA 6010B	
Vanadium	8.0	5.0	mg/kg	EPA 6010B	
Zinc	30	1.0	mg/kg	EPA 6010B	
PCB-1254	47	10	ug/kg	EPA 8082	
Pyrene	420	300	ug/kg	EPA 8270C	
Fluoranthene	360	300	ug/kg	EPA 8270C	

Sample ID: SP1-6

Laboratory ID: T181153-06

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	44	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	78	10	mg/kg	EPA 8015B	
Barium	44	1.0	mg/kg	EPA 6010B	
Chromium	4.2	2.0	mg/kg	EPA 6010B	
Cobalt	3.0	2.0	mg/kg	EPA 6010B	
Copper	5.1	1.0	mg/kg	EPA 6010B	
Lead	24	3.0	mg/kg	EPA 6010B	
Nickel	4.2	2.0	mg/kg	EPA 6010B	
Vanadium	11	5.0	mg/kg	EPA 6010B	
Zinc	24	1.0	mg/kg	EPA 6010B	
PCB-1254	52	10	ug/kg	EPA 8082	

Sample ID: SP1-7

Laboratory ID: T181153-07

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	88	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	92	10	mg/kg	EPA 8015B	
Barium	60	1.0	mg/kg	EPA 6010B	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-7

Laboratory ID: T181153-07

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Chromium	5.3	2.0		mg/kg	EPA 6010B	
Cobalt	4.2	2.0		mg/kg	EPA 6010B	
Copper	5.1	1.0		mg/kg	EPA 6010B	
Lead	11	3.0		mg/kg	EPA 6010B	
Nickel	5.5	2.0		mg/kg	EPA 6010B	
Vanadium	14	5.0		mg/kg	EPA 6010B	
Zinc	29	1.0		mg/kg	EPA 6010B	
PCB-1254	43	10		ug/kg	EPA 8082	
Pyrene	490	300		ug/kg	EPA 8270C	
Fluoranthene	420	300		ug/kg	EPA 8270C	

Sample ID: SP1-8

Laboratory ID: T181153-08

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	28	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	64	10		mg/kg	EPA 8015B	
Barium	66	1.0		mg/kg	EPA 6010B	
Chromium	6.1	2.0		mg/kg	EPA 6010B	
Cobalt	3.6	2.0		mg/kg	EPA 6010B	
Copper	7.2	1.0		mg/kg	EPA 6010B	
Lead	13	3.0		mg/kg	EPA 6010B	
Nickel	5.1	2.0		mg/kg	EPA 6010B	
Vanadium	15	5.0		mg/kg	EPA 6010B	
Zinc	46	1.0		mg/kg	EPA 6010B	
PCB-1254	32	10		ug/kg	EPA 8082	
Pyrene	360	300		ug/kg	EPA 8270C	

Sample ID: SP1-9

Laboratory ID: T181153-09

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	44	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	28	10		mg/kg	EPA 8015B	
Barium	67	1.0		mg/kg	EPA 6010B	
Chromium	5.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.1	2.0		mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-9

Laboratory ID: T181153-09

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Copper	5.3	1.0		mg/kg	EPA 6010B	
Nickel	5.6	2.0		mg/kg	EPA 6010B	
Vanadium	17	5.0		mg/kg	EPA 6010B	
Zinc	35	1.0		mg/kg	EPA 6010B	
PCB-1254	11	10		ug/kg	EPA 8082	

Sample ID: SP1-10

Laboratory ID: T181153-10

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	18	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	58	10		mg/kg	EPA 8015B	
Barium	64	1.0		mg/kg	EPA 6010B	
Chromium	5.5	2.0		mg/kg	EPA 6010B	
Cobalt	3.9	2.0		mg/kg	EPA 6010B	
Copper	5.8	1.0		mg/kg	EPA 6010B	
Nickel	5.8	2.0		mg/kg	EPA 6010B	
Vanadium	16	5.0		mg/kg	EPA 6010B	
Zinc	28	1.0		mg/kg	EPA 6010B	
PCB-1254	13	10		ug/kg	EPA 8082	

Sample ID: SP1-11

Laboratory ID: T181153-11

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	75	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	87	10		mg/kg	EPA 8015B	
Barium	54	1.0		mg/kg	EPA 6010B	
Chromium	5.4	2.0		mg/kg	EPA 6010B	
Cobalt	3.7	2.0		mg/kg	EPA 6010B	
Copper	6.1	1.0		mg/kg	EPA 6010B	
Lead	13	3.0		mg/kg	EPA 6010B	
Nickel	4.9	2.0		mg/kg	EPA 6010B	
Vanadium	13	5.0		mg/kg	EPA 6010B	
Zinc	28	1.0		mg/kg	EPA 6010B	
PCB-1254	61	10		ug/kg	EPA 8082	

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-12

Laboratory ID: T181153-12

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	97	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	55	10		mg/kg	EPA 8015B	
Barium	55	1.0		mg/kg	EPA 6010B	
Chromium	4.5	2.0		mg/kg	EPA 6010B	
Cobalt	3.7	2.0		mg/kg	EPA 6010B	
Copper	4.6	1.0		mg/kg	EPA 6010B	
Lead	12	3.0		mg/kg	EPA 6010B	
Nickel	4.9	2.0		mg/kg	EPA 6010B	
Vanadium	14	5.0		mg/kg	EPA 6010B	
Zinc	26	1.0		mg/kg	EPA 6010B	
PCB-1254	24	10		ug/kg	EPA 8082	

Sample ID: SP1-13

Laboratory ID: T181153-13

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	71	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	74	10		mg/kg	EPA 8015B	
Barium	52	1.0		mg/kg	EPA 6010B	
Chromium	4.6	2.0		mg/kg	EPA 6010B	
Cobalt	3.6	2.0		mg/kg	EPA 6010B	
Copper	6.0	1.0		mg/kg	EPA 6010B	
Lead	22	3.0		mg/kg	EPA 6010B	
Nickel	4.5	2.0		mg/kg	EPA 6010B	
Vanadium	13	5.0		mg/kg	EPA 6010B	
Zinc	29	1.0		mg/kg	EPA 6010B	
PCB-1254	43	10		ug/kg	EPA 8082	

Sample ID: SP1-14

Laboratory ID: T181153-14

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	16	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	35	10		mg/kg	EPA 8015B	
Barium	52	1.0		mg/kg	EPA 6010B	
Chromium	4.2	2.0		mg/kg	EPA 6010B	
Cobalt	3.6	2.0		mg/kg	EPA 6010B	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-14

Laboratory ID: T181153-14

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Copper	3.2	1.0	mg/kg	EPA 6010B	
Lead	21	3.0	mg/kg	EPA 6010B	
Nickel	3.9	2.0	mg/kg	EPA 6010B	
Vanadium	12	5.0	mg/kg	EPA 6010B	
Zinc	21	1.0	mg/kg	EPA 6010B	
PCB-1254	19	10	ug/kg	EPA 8082	

Sample ID: SP1-15

Laboratory ID: T181153-15

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	170	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	240	10	mg/kg	EPA 8015B	
Barium	52	1.0	mg/kg	EPA 6010B	
Chromium	6.0	2.0	mg/kg	EPA 6010B	
Cobalt	3.9	2.0	mg/kg	EPA 6010B	
Copper	5.7	1.0	mg/kg	EPA 6010B	
Lead	18	3.0	mg/kg	EPA 6010B	
Nickel	5.0	2.0	mg/kg	EPA 6010B	
Vanadium	14	5.0	mg/kg	EPA 6010B	
Zinc	27	1.0	mg/kg	EPA 6010B	
PCB-1254	110	10	ug/kg	EPA 8082	

Sample ID: SP1-16

Laboratory ID: T181153-16

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	85	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	55	10	mg/kg	EPA 8015B	
Barium	64	1.0	mg/kg	EPA 6010B	
Chromium	5.6	2.0	mg/kg	EPA 6010B	
Cobalt	5.1	2.0	mg/kg	EPA 6010B	
Copper	5.2	1.0	mg/kg	EPA 6010B	
Nickel	6.5	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	31	1.0	mg/kg	EPA 6010B	
PCB-1254	25	10	ug/kg	EPA 8082	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-17

Laboratory ID: T181153-17

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	81	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	40	10		mg/kg	EPA 8015B	
Barium	49	1.0		mg/kg	EPA 6010B	
Chromium	5.0	2.0		mg/kg	EPA 6010B	
Cobalt	3.6	2.0		mg/kg	EPA 6010B	
Copper	3.9	1.0		mg/kg	EPA 6010B	
Lead	9.4	3.0		mg/kg	EPA 6010B	
Nickel	4.6	2.0		mg/kg	EPA 6010B	
Vanadium	13	5.0		mg/kg	EPA 6010B	
Zinc	23	1.0		mg/kg	EPA 6010B	
PCB-1254	18	10		ug/kg	EPA 8082	

Sample ID: SP1-18

Laboratory ID: T181153-18

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	26	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	82	10		mg/kg	EPA 8015B	
Barium	62	1.0		mg/kg	EPA 6010B	
Chromium	5.2	2.0		mg/kg	EPA 6010B	
Cobalt	3.9	2.0		mg/kg	EPA 6010B	
Copper	4.4	1.0		mg/kg	EPA 6010B	
Lead	13	3.0		mg/kg	EPA 6010B	
Nickel	5.7	2.0		mg/kg	EPA 6010B	
Vanadium	15	5.0		mg/kg	EPA 6010B	
Zinc	27	1.0		mg/kg	EPA 6010B	
PCB-1254	42	10		ug/kg	EPA 8082	

Sample ID: SP1-19

Laboratory ID: T181153-19

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	39	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	51	10		mg/kg	EPA 8015B	
Barium	58	1.0		mg/kg	EPA 6010B	
Chromium	5.4	2.0		mg/kg	EPA 6010B	
Cobalt	3.8	2.0		mg/kg	EPA 6010B	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Sample ID: SP1-19

Laboratory ID: T181153-19

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Copper	4.6	1.0	mg/kg	EPA 6010B	
Nickel	5.3	2.0	mg/kg	EPA 6010B	
Vanadium	14	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	
PCB-1254	21	10	ug/kg	EPA 8082	

Sample ID: SP1-20

Laboratory ID: T181153-20

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	76	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	150	10	mg/kg	EPA 8015B	
Barium	57	1.0	mg/kg	EPA 6010B	
Chromium	5.9	2.0	mg/kg	EPA 6010B	
Cobalt	3.9	2.0	mg/kg	EPA 6010B	
Copper	4.6	1.0	mg/kg	EPA 6010B	
Lead	10	3.0	mg/kg	EPA 6010B	
Nickel	5.3	2.0	mg/kg	EPA 6010B	
Vanadium	16	5.0	mg/kg	EPA 6010B	
Zinc	30	1.0	mg/kg	EPA 6010B	
PCB-1254	50	10	ug/kg	EPA 8082	
Pyrene	420	300	ug/kg	EPA 8270C	
Fluoranthene	340	300	ug/kg	EPA 8270C	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-1
T181153-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/09/18	EPA 8015B
C13-C28 (DRO)	47	10	"	"	"	"	"	"
C29-C40 (MORO)	100	10	"	"	"	"	"	"
Surrogate: <i>p</i> -Terphenyl		74.3 %	65-135		"	"	"	"

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B
Silver	ND	2.0	"	"	"	"	"	"
Arsenic	ND	5.0	"	"	"	"	"	"
Barium	54	1.0	"	"	"	"	"	"
Beryllium	ND	1.0	"	"	"	"	"	"
Cadmium	ND	2.0	"	"	"	"	"	"
Chromium	5.4	2.0	"	"	"	"	"	"
Cobalt	3.7	2.0	"	"	"	"	"	"
Copper	7.8	1.0	"	"	"	"	"	"
Lead	22	3.0	"	"	"	"	"	"
Molybdenum	ND	5.0	"	"	"	"	"	"
Nickel	5.1	2.0	"	"	"	"	"	"
Selenium	ND	5.0	"	"	"	"	"	"
Thallium	ND	2.0	"	"	"	"	"	"
Vanadium	12	5.0	"	"	"	"	"	"
Zinc	36	1.0	"	"	"	"	"	"

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-1

T181153-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	35	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		131 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		136 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	2400	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	900	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	730	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	410	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	
Benzo (a) pyrene	360	300	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-1

T181153-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzyl alcohol	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	1100	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	2100	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	340	300	"	"	"	"	"	"	
Isophorone	ND	300	"	"	"	"	"	"	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-1

T181153-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

2-Methylphenol	ND	1000	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	1300	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol		52.5 %	15-121		"	"	"	"
Surrogate: Phenol-d6		56.4 %	24-113		"	"	"	"
Surrogate: Nitrobenzene-d5		74.7 %	21.3-119		"	"	"	"
Surrogate: 2-Fluorobiphenyl		67.5 %	32.4-102		"	"	"	"
Surrogate: 2,4,6-Tribromophenol		74.9 %	18.1-105		"	"	"	"
Surrogate: Terphenyl-d14		118 %	29.1-130		"	"	"	"

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-2

T181153-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/09/18	EPA 8015B	
C13-C28 (DRO)	13	10	"	"	"	"	"	"	
C29-C40 (MORO)	86	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		75.8 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	65	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.4	2.0	"	"	"	"	"	"	
Cobalt	4.0	2.0	"	"	"	"	"	"	
Copper	3.7	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-2

T181153-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

119 % 35-140

" " " "

Surrogate: Decachlorobiphenyl

127 % 35-140

" " " "

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-2

T181153-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-2
T181153-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	58.2 %	15-121			"	"	"	"
Surrogate: Phenol-d6	58.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	79.6 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	70.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	70.2 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	116 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-3

T181153-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/09/18	EPA 8015B	
C13-C28 (DRO)	47	10	"	"	"	"	"	"	
C29-C40 (MORO)	68	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		77.9 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	61	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.3	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	4.0	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-3

T181153-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	25	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		122 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		132 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	300	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-3

T181153-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-3

T181153-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol		59.7 %	15-121		"	"	"	"
Surrogate: Phenol-d6		63.1 %	24-113		"	"	"	"
Surrogate: Nitrobenzene-d5		80.9 %	21.3-119		"	"	"	"
Surrogate: 2-Fluorobiphenyl		75.6 %	32.4-102		"	"	"	"
Surrogate: 2,4,6-Tribromophenol		73.9 %	18.1-105		"	"	"	"
Surrogate: Terphenyl-d14		121 %	29.1-130		"	"	"	"

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-4

T181153-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/09/18	EPA 8015B	
C13-C28 (DRO)	24	10	"	"	"	"	"	"	
C29-C40 (MORO)	75	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		76.2 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	69	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.3	2.0	"	"	"	"	"	"	
Cobalt	4.6	2.0	"	"	"	"	"	"	
Copper	4.2	1.0	"	"	"	"	"	"	
Lead	8.4	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	32	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-4

T181153-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	13	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		106 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		123 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-4

T181153-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-4
T181153-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	42.7 %	15-121			"	"	"	"
Surrogate: Phenol-d6	45.7 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	60.3 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	56.3 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	61.2 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	108 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-5
T181153-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	96	10	"	"	"	"	"	"	
C29-C40 (MORO)	52	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		74.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	41	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.6	2.0	"	"	"	"	"	"	
Cobalt	3.0	2.0	"	"	"	"	"	"	
Copper	11	1.0	"	"	"	"	"	"	
Lead	45	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	8.0	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-5
T181153-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	47	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		118 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		123 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	420	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-5
T181153-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	360	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-5
T181153-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	57.7 %	15-121			"	"	"	"
Surrogate: Phenol-d6	61.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	80.7 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	75.5 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	75.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	126 %	29.1-130			"	"	"	"

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-6
T181153-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	44	10	"	"	"	"	"	"	
C29-C40 (MORO)	78	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		78.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	44	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.2	2.0	"	"	"	"	"	"	
Cobalt	3.0	2.0	"	"	"	"	"	"	
Copper	5.1	1.0	"	"	"	"	"	"	
Lead	24	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	11	5.0	"	"	"	"	"	"	
Zinc	24	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-6
T181153-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	52	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		133 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		128 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-6
T181153-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-6
T181153-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	58.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	62.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	80.3 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	73.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	74.3 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	127 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-7

T181153-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	88	10	"	"	"	"	"	"	
C29-C40 (MORO)	92	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		81.8 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	60	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.3	2.0	"	"	"	"	"	"	
Cobalt	4.2	2.0	"	"	"	"	"	"	
Copper	5.1	1.0	"	"	"	"	"	"	
Lead	11	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	14	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Reported:
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SP1-7

T181153-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	43	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		128 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		140 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	490	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-7

T181153-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	420	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-7

T181153-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	63.6 %	15-121			"	"	"	"	
Surrogate: Phenol-d6	66.3 %	24-113			"	"	"	"	
Surrogate: Nitrobenzene-d5	86.5 %	21.3-119			"	"	"	"	
Surrogate: 2-Fluorobiphenyl	80.1 %	32.4-102			"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	80.0 %	18.1-105			"	"	"	"	
Surrogate: Terphenyl-d14	136 %	29.1-130			"	"	"	"	S-GC

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-8
T181153-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	28	10	"	"	"	"	"	"	
C29-C40 (MORO)	64	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		77.5 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	66	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.1	2.0	"	"	"	"	"	"	
Cobalt	3.6	2.0	"	"	"	"	"	"	
Copper	7.2	1.0	"	"	"	"	"	"	
Lead	13	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	46	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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SP1-8
T181153-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	32	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		118 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		140 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	360	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-8

T181153-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-8
T181153-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	57.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	60.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	77.3 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	72.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	71.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	118 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-9

T181153-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	44	10	"	"	"	"	"	"	
C29-C40 (MORO)	28	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		78.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	67	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.7	2.0	"	"	"	"	"	"	
Cobalt	4.1	2.0	"	"	"	"	"	"	
Copper	5.3	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	35	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-9

T181153-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	11	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		112 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		112 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-9

T181153-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-9
T181153-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	54.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	70.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	66.0 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	75.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	116 %	29.1-130			"	"	"	"

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-10
T181153-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	18	10	"	"	"	"	"	"	
C29-C40 (MORO)	58	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		80.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	64	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.5	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	5.8	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.8	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-10
T181153-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	13	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		116 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		133 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-10
T181153-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-10
T181153-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	53.1 %	15-121			"	"	"	"
Surrogate: Phenol-d6	56.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	73.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	66.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	69.5 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	119 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-11
T181153-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	75	10	"	"	"	"	"	"	
C29-C40 (MORO)	87	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		80.8 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	54	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.4	2.0	"	"	"	"	"	"	
Cobalt	3.7	2.0	"	"	"	"	"	"	
Copper	6.1	1.0	"	"	"	"	"	"	
Lead	13	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	13	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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SP1-11
T181153-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	61	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		115 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		129 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-11
T181153-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-11
T181153-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	57.8 %	15-121			"	"	"	"
Surrogate: Phenol-d6	61.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	77.9 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	72.1 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	72.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	124 %	29.1-130			"	"	"	"

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-12
T181153-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	97	10	"	"	"	"	"	"	
C29-C40 (MORO)	55	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		83.8 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	55	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.5	2.0	"	"	"	"	"	"	
Cobalt	3.7	2.0	"	"	"	"	"	"	
Copper	4.6	1.0	"	"	"	"	"	"	
Lead	12	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	14	5.0	"	"	"	"	"	"	
Zinc	26	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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04/12/18 15:44

SP1-12
T181153-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	24	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		122 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		134 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

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04/12/18 15:44

SP1-12
T181153-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-12
T181153-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.8 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.1 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	72.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	70.1 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	71.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	126 %	29.1-130			"	"	"	"

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-13
T181153-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	71	10	"	"	"	"	"	"	
C29-C40 (MORO)	74	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		75.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	52	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.6	2.0	"	"	"	"	"	"	
Cobalt	3.6	2.0	"	"	"	"	"	"	
Copper	6.0	1.0	"	"	"	"	"	"	
Lead	22	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	13	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-13
T181153-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	43	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		118 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		140 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-13
T181153-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-13
T181153-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	48.9 %	15-121			"	"	"	"
Surrogate: Phenol-d6	53.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	68.6 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	67.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	71.4 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	125 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-14
T181153-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	16	10	"	"	"	"	"	"	
C29-C40 (MORO)	35	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		72.6 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	52	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.2	2.0	"	"	"	"	"	"	
Cobalt	3.6	2.0	"	"	"	"	"	"	
Copper	3.2	1.0	"	"	"	"	"	"	
Lead	21	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	12	5.0	"	"	"	"	"	"	
Zinc	21	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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04/12/18 15:44

SP1-14
T181153-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	19	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		116 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		122 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-14
T181153-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-14
T181153-14 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	55.7 %	15-121	"	"	"	"	"	"
Surrogate: Phenol-d6	56.3 %	24-113	"	"	"	"	"	"
Surrogate: Nitrobenzene-d5	74.6 %	21.3-119	"	"	"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.0 %	32.4-102	"	"	"	"	"	"
Surrogate: 2,4,6-Tribromophenol	68.6 %	18.1-105	"	"	"	"	"	"
Surrogate: Terphenyl-d14	120 %	29.1-130	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-15
T181153-15 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	170	10	"	"	"	"	"	"	
C29-C40 (MORO)	240	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		73.1 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	52	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.0	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	5.7	1.0	"	"	"	"	"	"	
Lead	18	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.0	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	14	5.0	"	"	"	"	"	"	
Zinc	27	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-15
T181153-15 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	110	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		119 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		133 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-15
T181153-15 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-15
T181153-15 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	45.2 %	15-121			"	"	"	"
Surrogate: Phenol-d6	49.3 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	62.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	61.0 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	62.6 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	107 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-16
T181153-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	85	10	"	"	"	"	"	"	
C29-C40 (MORO)	55	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		75.3 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	64	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.6	2.0	"	"	"	"	"	"	
Cobalt	5.1	2.0	"	"	"	"	"	"	
Copper	5.2	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	31	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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04/12/18 15:44

SP1-16
T181153-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	25	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		109 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		111 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-16
T181153-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-16
T181153-16 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	49.5 %	15-121			"	"	"	"
Surrogate: Phenol-d6	52.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	67.4 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	66.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	68.2 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	120 %	29.1-130			"	"	"	"

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-17
T181153-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	81	10	"	"	"	"	"	"	
C29-C40 (MORO)	40	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		88.5 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	49	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.0	2.0	"	"	"	"	"	"	
Cobalt	3.6	2.0	"	"	"	"	"	"	
Copper	3.9	1.0	"	"	"	"	"	"	
Lead	9.4	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	13	5.0	"	"	"	"	"	"	
Zinc	23	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-17
T181153-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	18	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		125 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		131 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-17
T181153-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-17
T181153-17 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/09/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	56.8 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	76.6 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	73.5 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	70.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	111 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-18
T181153-18 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	26	10	"	"	"	"	"	"	
C29-C40 (MORO)	82	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		76.5 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	62	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.2	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	4.4	1.0	"	"	"	"	"	"	
Lead	13	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.7	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	27	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-18
T181153-18 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	42	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		120 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		136 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-18
T181153-18 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-18
T181153-18 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	61.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	64.3 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	80.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	73.7 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	72.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	115 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-19
T181153-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	39	10	"	"	"	"	"	"	
C29-C40 (MORO)	51	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		73.6 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	58	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.4	2.0	"	"	"	"	"	"	
Cobalt	3.8	2.0	"	"	"	"	"	"	
Copper	4.6	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.3	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	14	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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Reported:
04/12/18 15:44

SP1-19
T181153-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/06/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	21	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		101 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		116 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-19
T181153-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-19
T181153-19 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	54.3 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.3 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	72.8 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	72.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	120 %	29.1-130			"	"	"	"

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Project Number: 137-002
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Reported:
04/12/18 15:44

SP1-20
T181153-20 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040518	04/05/18	04/10/18	EPA 8015B	
C13-C28 (DRO)	76	10	"	"	"	"	"	"	
C29-C40 (MORO)	150	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		81.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040603	04/06/18	04/06/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	57	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.9	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	4.6	1.0	"	"	"	"	"	"	
Lead	10	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.3	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040605	04/06/18	04/06/18	EPA 7471A Soil	

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SP1-20
T181153-20 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040520	04/05/18	04/07/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	50	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		134 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		131 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	420	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-20
T181153-20 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	340	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

SP1-20
T181153-20 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040523	04/05/18	04/10/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.3 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.0 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	72.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	73.2 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	117 %	29.1-130			"	"	"	"

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Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040518 - EPA 3550B GC

Blank (8040518-BLK1)

Prepared: 04/05/18 Analyzed: 04/09/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	77.1		"	100		77.1	65-135			

LCS (8040518-BS1)

Prepared: 04/05/18 Analyzed: 04/09/18

C13-C28 (DRO)	500	10	mg/kg	500		101	75-125			
Surrogate: <i>p</i> -Terphenyl	81.2		"	100		81.2	65-135			

Matrix Spike (8040518-MS1)

Source: T181153-01

Prepared: 04/05/18 Analyzed: 04/09/18

C13-C28 (DRO)	580	10	mg/kg	500	47	107	75-125			
Surrogate: <i>p</i> -Terphenyl	84.4		"	100		84.4	65-135			

Matrix Spike Dup (8040518-MSD1)

Source: T181153-01

Prepared: 04/05/18 Analyzed: 04/09/18

C13-C28 (DRO)	520	10	mg/kg	495	47	95.8	75-125	10.8	20	
Surrogate: <i>p</i> -Terphenyl	74.9		"	99.0		75.6	65-135			

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Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040603 - EPA 3050B

Blank (8040603-BLK1)

Prepared & Analyzed: 04/06/18

Antimony	ND	3.0	mg/kg
Silver	ND	2.0	"
Arsenic	ND	5.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	2.0	"
Chromium	ND	2.0	"
Cobalt	ND	2.0	"
Copper	ND	1.0	"
Lead	ND	3.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Selenium	ND	5.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	1.60	1.0	"

QB-01

LCS (8040603-BS1)

Prepared & Analyzed: 04/06/18

Arsenic	105	5.0	mg/kg	100	105	75-125
Barium	105	1.0	"	100	105	75-125
Cadmium	104	2.0	"	100	104	75-125
Chromium	105	2.0	"	100	105	75-125
Lead	109	3.0	"	100	109	75-125

Matrix Spike (8040603-MS1)

Source: T181153-01

Prepared & Analyzed: 04/06/18

Arsenic	96.5	5.0	mg/kg	100	2.35	94.2	75-125
Barium	155	1.0	"	100	53.7	101	75-125
Cadmium	95.9	2.0	"	100	0.220	95.7	75-125
Chromium	103	2.0	"	100	5.36	97.6	75-125
Lead	117	3.0	"	100	21.6	95.8	75-125

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Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040603 - EPA 3050B

Matrix Spike Dup (8040603-MSD1)

Source: T181153-01

Prepared & Analyzed: 04/06/18

Arsenic	96.2	5.0	mg/kg	100	2.35	93.9	75-125	0.314	20	
Barium	158	1.0	"	100	53.7	104	75-125	1.77	20	
Cadmium	92.7	2.0	"	100	0.220	92.5	75-125	3.42	20	
Chromium	99.9	2.0	"	100	5.36	94.6	75-125	3.02	20	
Lead	118	3.0	"	100	21.6	96.2	75-125	0.371	20	

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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040605 - EPA 7471A Soil

Blank (8040605-BLK1)

Prepared & Analyzed: 04/06/18

Mercury	ND	0.10	mg/kg							

LCS (8040605-BS1)

Prepared & Analyzed: 04/06/18

Mercury	0.433	0.10	mg/kg	0.417		104	80-120			

Matrix Spike (8040605-MS1)

Source: T181153-01

Prepared & Analyzed: 04/06/18

Mercury	0.446	0.10	mg/kg	0.417	ND	107	75-125			

Matrix Spike Dup (8040605-MSD1)

Source: T181153-01

Prepared & Analyzed: 04/06/18

Mercury	0.447	0.10	mg/kg	0.417	ND	107	75-125	0.150	20	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040520 - EPA 3550 ECD/GCMS

Blank (8040520-BLK1)

Prepared: 04/05/18 Analyzed: 04/06/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	13.4		"	10.0		134	35-140			
Surrogate: Decachlorobiphenyl	14.4		"	10.0		144	35-140			S-GC

LCS (8040520-BS1)

Prepared: 04/05/18 Analyzed: 04/06/18

PCB-1016	63.2	10	ug/kg	100		63.2	40-130			
PCB-1260	63.7	10	"	100		63.7	40-130			
Surrogate: Tetrachloro-meta-xylene	13.0		"	10.0		130	35-140			
Surrogate: Decachlorobiphenyl	14.6		"	10.0		146	35-140			S-GC

Matrix Spike (8040520-MS1)

Source: T181153-01

Prepared: 04/05/18 Analyzed: 04/06/18

PCB-1016	62.7	10	ug/kg	98.0	ND	64.0	40-130			
PCB-1260	56.1	10	"	98.0	ND	57.2	40-130			
Surrogate: Tetrachloro-meta-xylene	11.7		"	9.80		120	35-140			
Surrogate: Decachlorobiphenyl	12.9		"	9.80		132	35-140			

Matrix Spike Dup (8040520-MSD1)

Source: T181153-01

Prepared: 04/05/18 Analyzed: 04/06/18

PCB-1016	62.6	10	ug/kg	98.0	ND	63.9	40-130	0.156	30	
PCB-1260	52.3	10	"	98.0	ND	53.3	40-130	7.06	30	
Surrogate: Tetrachloro-meta-xylene	10.4		"	9.80		106	35-140			
Surrogate: Decachlorobiphenyl	11.7		"	9.80		120	35-140			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040523 - EPA 3550 ECD/GCMS

Blank (8040523-BLK1)

Prepared: 04/05/18 Analyzed: 04/09/18

Carbazole	ND	300	ug/kg
Phenol	ND	1000	"
Aniline	ND	300	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
1-Methylnaphthalene	ND	300	"
2-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

SunStar Laboratories, Inc.

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Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040523 - EPA 3550 ECD/GCMS

Blank (8040523-BLK1)

Prepared: 04/05/18 Analyzed: 04/09/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
Pyridine	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
2,4,6-Trichlorophenol	ND	1000	"

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Brusca Associates Inc.
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Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040523 - EPA 3550 ECD/GCMS

Blank (8040523-BLK1)

Prepared: 04/05/18 Analyzed: 04/09/18

Surrogate: 2-Fluorophenol	1800		ug/kg	3320		54.2	15-121			
Surrogate: Phenol-d6	1820		"	3320		54.8	24-113			
Surrogate: Nitrobenzene-d5	2540		"	3320		76.3	21.3-119			
Surrogate: 2-Fluorobiphenyl	2260		"	3320		67.9	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2120		"	3320		63.7	18.1-105			
Surrogate: Terphenyl-d14	4060		"	3320		122	29.1-130			

LCS (8040523-BS1)

Prepared: 04/05/18 Analyzed: 04/09/18

Phenol	1960	1000	ug/kg	3320		59.0	34-114			
2-Chlorophenol	1990	1000	"	3320		59.8	34-114			
1,4-Dichlorobenzene	2100	300	"	3320		63.1	34-114			
N-Nitrosodi-n-propylamine	2580	300	"	3320		77.6	30-110			
1,2,4-Trichlorobenzene	2300	300	"	3320		69.2	39-119			
4-Chloro-3-methylphenol	2780	1000	"	3320		83.7	50-130			
Acenaphthene	2510	300	"	3320		75.7	34-114			
Pentachlorophenol	1730	1000	"	3320		52.1	50-130			
Pyrene	2550	300	"	3320		76.8	30-110			
Surrogate: 2-Fluorophenol	1760		"	3320		53.1	15-121			
Surrogate: Phenol-d6	1780		"	3320		53.7	24-113			
Surrogate: Nitrobenzene-d5	2520		"	3320		75.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2250		"	3320		67.7	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2360		"	3320		71.2	18.1-105			
Surrogate: Terphenyl-d14	4160		"	3320		125	29.1-130			

Matrix Spike (8040523-MS1)

Source: T181153-01

Prepared: 04/05/18 Analyzed: 04/09/18

Phenol	2040	1000	ug/kg	3300	ND	61.8	34-114			
2-Chlorophenol	1920	1000	"	3300	ND	58.2	34-114			
1,4-Dichlorobenzene	1940	300	"	3300	ND	58.8	34-114			
N-Nitrosodi-n-propylamine	2240	300	"	3300	ND	68.0	30-110			
1,2,4-Trichlorobenzene	2310	300	"	3300	ND	70.1	39-119			
4-Chloro-3-methylphenol	2920	1000	"	3300	ND	88.5	50-130			
Acenaphthene	2580	300	"	3300	98.0	75.1	34-114			
Pentachlorophenol	1860	1000	"	3300	ND	56.5	50-130			
Pyrene	5710	300	"	3300	2380	101	30-110			
Surrogate: 2-Fluorophenol	1670		"	3300		50.7	15-121			
Surrogate: Phenol-d6	1910		"	3300		58.0	24-113			

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Mike Jaroudi, Project Manager



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Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040523 - EPA 3550 ECD/GCMS

Matrix Spike (8040523-MS1)		Source: T181153-01		Prepared: 04/05/18 Analyzed: 04/09/18						
Surrogate: Nitrobenzene-d5	2430		ug/kg	3300		73.6	21.3-119			
Surrogate: 2-Fluorobiphenyl	2210		"	3300		66.9	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2470		"	3300		74.9	18.1-105			
Surrogate: Terphenyl-d14	3700		"	3300		112	29.1-130			
Matrix Spike Dup (8040523-MSD1)		Source: T181153-01		Prepared: 04/05/18 Analyzed: 04/09/18						
Phenol	1920	1000	ug/kg	3290	ND	58.3	34-114	6.11	42	
2-Chlorophenol	1790	1000	"	3290	ND	54.5	34-114	6.95	40	
1,4-Dichlorobenzene	1850	300	"	3290	ND	56.1	34-114	4.99	28	
N-Nitrosodi-n-propylamine	2130	300	"	3290	ND	64.8	30-110	5.10	38	
1,2,4-Trichlorobenzene	2150	300	"	3290	ND	65.5	39-119	7.17	28	
4-Chloro-3-methylphenol	3050	1000	"	3290	ND	92.8	50-130	4.37	42	
Acenaphthene	2450	300	"	3290	98.0	71.5	34-114	4.93	31	
Pentachlorophenol	1820	1000	"	3290	ND	55.3	50-130	2.53	50	
Pyrene	4600	300	"	3290	2380	67.6	30-110	21.5	31	
Surrogate: 2-Fluorophenol	1600		"	3290		48.7	15-121			
Surrogate: Phenol-d6	1680		"	3290		51.1	24-113			
Surrogate: Nitrobenzene-d5	2250		"	3290		68.4	21.3-119			
Surrogate: 2-Fluorobiphenyl	2070		"	3290		62.9	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2380		"	3290		72.3	18.1-105			
Surrogate: Terphenyl-d14	3650		"	3290		111	29.1-130			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48 Acre Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/12/18 15:44

Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

QB-01 The method blank contains analyte at a concentration above the MRL; however, concentration is less than 10% of the sample result, which is negligible according to method criteria.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO BOX 332, ROSEMEAD, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/4/18 Page: 1 of 2
Project Name: LATHROP 48-Ac PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: T181153 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 PCBs 8092	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	LEAD AND ARSENIC	Laboratory ID #	Comments/Preservative	Total # of containers
SP1-1	4/4/18	11:59	SOLID	JAR	X	X	X	X	X	X	X	X	X	X	X	01		1
SP1-2	-	12:01	-	-	X	X	X	X	X	X	X	X	X	X	X	02		1
SP1-3	-	12:03	-	-	X	X	X	X	X	X	X	X	X	X	X	03		1
SP1-4	-	12:05	-	-	X	X	X	X	X	X	X	X	X	X	X	04		1
SP1-5	-	12:12	-	-	X	X	X	X	X	X	X	X	X	X	X	05		1
SP1-6	-	12:13	-	-	X	X	X	X	X	X	X	X	X	X	X	06		1
SP1-7	-	12:17	-	-	X	X	X	X	X	X	X	X	X	X	X	07		1
SP1-8	-	12:22	-	-	X	X	X	X	X	X	X	X	X	X	X	08		1
SP1-9	-	12:25	-	-	X	X	X	X	X	X	X	X	X	X	X	09		1
SP1-10	-	12:31	-	-	X	X	X	X	X	X	X	X	X	X	X	10		1
SP1-11	-	12:35	-	-	X	X	X	X	X	X	X	X	X	X	X	11		1
SP1-12	-	12:37	-	-	X	X	X	X	X	X	X	X	X	X	X	12		1
SP1-13	-	12:42	-	-	X	X	X	X	X	X	X	X	X	X	X	13		1
SP1-14	-	12:45	-	-	X	X	X	X	X	X	X	X	X	X	X	14		1
SP1-15	-	12:47	-	-	X	X	X	X	X	X	X	X	X	X	X	15		1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/4/18 14:38</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4/4/18 14:38</u>					Total # of containers					Notes			
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4-5-18 955</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4-5-18 955</u>					Chain of Custody seals DN/A					* 3 DAY TAT ON METALS			
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4-5-18 955</u>					Received by: (signature) <u>[Signature]</u> Date / Time <u>4-5-18 955</u>					Received good condition/cold					4.1			

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

Turn around time: NORMAL
EXCEPT METALS -
3 DAY TAT ON METALS **COC 160906**



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Centre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, ROSEVILLE, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/4/18 Page: 2 of 2
Project Name: LATHROP 48-Ac PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: T181153 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260	8260 + OXY	8260 BTEX, OXY only	8270	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP1-16	4/4/18	12:51	Solid	JAR	X	X	X	X	X	X	X	X	X	X	14		1
SP1-17		12:53			X	X	X	X	X	X	X	X	X	X	17		1
SP1-18		12:56			X	X	X	X	X	X	X	X	X	X	18		1
SP1-19		13:05			X	X	X	X	X	X	X	X	X	X	19		1
SP1-20		13:08			X	X	X	X	X	X	X	X	X	X	20		1
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Total # of containers												Notes
<u>[Signature]</u>	4/4/18 14:38		<u>[Signature]</u>	4/4/18 14:38													* 3 DAY TAT ON METALS
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Chain of Custody seals 9/N/A												
<u>[Signature]</u>	4-5-18 9:55		<u>[Signature]</u>	4-5-18 9:55	Seals intact? 0/N/A												
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Received good condition/cold												
<u>[Signature]</u>			<u>[Signature]</u>		4.1												

Sample disposal instructions: Disposal @ \$2.00 each _____

Return to client _____ Pickup _____

Turn around time: Normal
EXCEPT METALS -
3 DAY TAT ON METALS
COC 160912



25712 Commercentre Drive
Lake Forest, California 92630
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18 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/06/18 10:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP2-1	T181179-01	Soil	04/05/18 08:54	04/06/18 10:00

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

DETECTIONS SUMMARY

Sample ID: SP2-1

Laboratory ID: T181179-01

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Arsenic	17	1.0	ug/l	6020 ICP-MS	AO-1
Barium	450	1.0	ug/l	6020 ICP-MS	AO-1
Beryllium	1.3	1.0	ug/l	6020 ICP-MS	AO-1
Chromium	65	1.0	ug/l	6020 ICP-MS	AO-1
Cobalt	49	1.0	ug/l	6020 ICP-MS	AO-1
Copper	96	1.0	ug/l	6020 ICP-MS	AO-1
Lead	170	2.0	ug/l	6020 ICP-MS	AO-1
Mercury	0.33	0.10	ug/l	6020 ICP-MS	AO-1
Molybdenum	9.1	1.0	ug/l	6020 ICP-MS	AO-1
Nickel	61	1.0	ug/l	6020 ICP-MS	AO-1
Silver	5.0	1.0	ug/l	6020 ICP-MS	AO-1
Vanadium	220	1.0	ug/l	6020 ICP-MS	AO-1
Zinc	340	1.0	ug/l	6020 ICP-MS	AO-1

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

SP2-1

T181179-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8041135	04/12/18	04/17/18	6020 ICP-MS	AO-1
Arsenic	17	1.0	"	"	"	"	"	"	AO-1
Barium	450	1.0	"	"	"	"	"	"	AO-1
Beryllium	1.3	1.0	"	"	"	"	"	"	AO-1
Cadmium	ND	1.0	"	"	"	"	"	"	AO-1
Chromium	65	1.0	"	"	"	"	"	"	AO-1
Cobalt	49	1.0	"	"	"	"	"	"	AO-1
Copper	96	1.0	"	"	"	"	"	"	AO-1
Lead	170	2.0	"	"	"	"	"	"	AO-1
Mercury	0.33	0.10	"	"	"	"	"	"	AO-1
Molybdenum	9.1	1.0	"	"	"	"	"	"	AO-1
Nickel	61	1.0	"	"	"	"	"	"	AO-1
Selenium	ND	5.0	"	"	"	"	"	"	AO-1
Silver	5.0	1.0	"	"	"	"	"	"	AO-1
Thallium	ND	1.0	"	"	"	"	"	"	AO-1
Vanadium	220	1.0	"	"	"	"	"	"	AO-1
Zinc	340	1.0	"	"	"	"	"	"	AO-1

SunStar Laboratories, Inc.

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PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041135 - EPA 3010A

Blank (8041135-BLK1)

Prepared: 04/11/18 Analyzed: 04/17/18

Antimony	ND	1.0	ug/l							AO-1
Arsenic	ND	1.0	"							AO-1
Barium	16.1	1.0	"							AO-1, QB-01
Beryllium	ND	1.0	"							AO-1
Cadmium	ND	1.0	"							AO-1
Chromium	1.93	1.0	"							AO-1, QB-01
Cobalt	ND	1.0	"							AO-1
Copper	1.50	1.0	"							AO-1, QB-01
Lead	ND	2.0	"							AO-1
Mercury	ND	0.10	"							AO-1
Molybdenum	ND	1.0	"							AO-1
Nickel	ND	1.0	"							AO-1
Selenium	ND	5.0	"							AO-1
Silver	ND	1.0	"							AO-1
Thallium	ND	1.0	"							AO-1
Vanadium	3.01	1.0	"							AO-1, QB-01
Zinc	11.1	1.0	"							AO-1, QB-01

LCS (8041135-BS1)

Prepared: 04/11/18 Analyzed: 04/17/18

Arsenic	48.2	1.0	ug/l	50.0		96.4	80-120			AO-1
Barium	48.8	1.0	"	50.0		97.5	80-120			AO-1
Cadmium	47.7	1.0	"	50.0		95.4	80-120			AO-1
Chromium	52.4	1.0	"	50.0		105	80-120			AO-1
Lead	51.3	2.0	"	50.0		103	80-120			AO-1

Duplicate (8041135-DUP1)

Source: T181153-05

Prepared: 04/11/18 Analyzed: 04/17/18

Barium	432	5.0	ug/l		341		23.6	200		AO-1
Lead	445	10	"		361		20.8	200		AO-1

SunStar Laboratories, Inc.

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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041135 - EPA 3010A

Matrix Spike (8041135-MS1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Arsenic	62.4	1.0	ug/l	50.0	19.8	85.2	75-125			AO-1
Barium	363	1.0	"	50.0	341	44.9	75-125			AO-1, QM-05
Cadmium	45.9	1.0	"	50.0	1.59	88.7	75-125			AO-1
Chromium	107	1.0	"	50.0	59.6	94.3	75-125			AO-1
Lead	368	2.0	"	50.0	361	12.3	75-125			AO-1, QM-05

Matrix Spike Dup (8041135-MSD1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Arsenic	61.3	1.0	ug/l	50.0	19.8	82.9	75-125	1.81	20	AO-1
Barium	372	1.0	"	50.0	341	62.8	75-125	2.43	20	AO-1, QM-05
Cadmium	45.7	1.0	"	50.0	1.59	88.2	75-125	0.541	20	AO-1
Chromium	107	1.0	"	50.0	59.6	94.9	75-125	0.296	20	AO-1
Lead	389	2.0	"	50.0	361	55.4	75-125	5.69	20	AO-1, QM-05

Post Spike (8041135-PS1)		Source: T181153-05			Prepared: 04/11/18 Analyzed: 04/17/18					
Barium	188		ug/l	2.50	0.341	NR	80-120			AO-1, QM-05
Lead	199		"	2.50	0.361	NR	80-120			AO-1, QM-05

SunStar Laboratories, Inc.

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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/18/18 09:13

Notes and Definitions

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

QB-01 The method blank contains analyte at a concentration above the MRL; however, concentration is less than 10% of the sample result, which is negligible according to method criteria.

AO-1 STLC Leach Ran 6020 Water

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

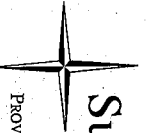
dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commercentre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO BOX 332, ROSEVILLE, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/5/18 Page: 1 Of 1
Project Name: LATHROP 48-A PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 7181178 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 - SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 - PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP2-1	4/5/18	8:54	Solid	JAR	X			X				X	X	X	01		1
SP2-2		8:57			X			X				X	X	X	02		1
SP2-3		8:59			X			X				X	X	X	03		1
SP2-4		9:02			X			X				X	X	X	04		1
SP2-5		9:04			X			X				X	X	X	05		1
SP2-6		9:08			X			X				X	X	X	06		1
SP2-7		9:14			X			X				X	X	X	07		1
SP2-8		9:18			X			X				X	X	X	08		1
SP2-9		9:21			X			X				X	X	X	09		1
SP2-10		9:24			X			X				X	X	X	10		1
SP2-11		9:26			X			X				X	X	X	11		1
SP2-12		9:29			X			X				X	X	X	12		1
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Total # of containers												Notes
<u>[Signature]</u>	4/5/18-13:10		<u>[Signature]</u>	4/5/18 13:10													* 3 DAY TAT ON METALS
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													
<u>[Signature]</u>	4/5/18 10:00		<u>[Signature]</u>	4/5/18 10:00													
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													
<u>[Signature]</u>	4/5/18 10:00		<u>[Signature]</u>	4/5/18 10:00													

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

Turn around time: NORMAL
EXCEPT 3 DAY
TAT ON METALS

COC 160913



April 26, 2018

FAL Project ID: 11426

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11426**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Twelve solid samples were received on 4/5/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of ten business days for project **11426**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and the sample photos. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11426**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink that reads "Thomas C. Crabtree".

Thomas C. Crabtree
Director

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: 11426

Received on: 04/05/2018

Project Due: 04/20/2018

Storage: R-4

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11426-001-SA	0	137-002	SP2-1	EPA 1613 D/F	Solid	04/05/2018	08:54 am	04/05/2019
11426-002-SA	0	137-002	SP2-2	EPA 1613 D/F	Solid	04/05/2018	08:57 am	04/05/2019
11426-003-SA	0	137-002	SP2-3	EPA 1613 D/F	Solid	04/05/2018	08:59 am	04/05/2019
11426-004-SA	0	137-002	SP2-4	EPA 1613 D/F	Solid	04/05/2018	09:02 am	04/05/2019
11426-005-SA	0	137-002	SP2-5	EPA 1613 D/F	Solid	04/05/2018	09:04 am	04/05/2019
11426-006-SA	0	137-002	SP2-6	EPA 1613 D/F	Solid	04/05/2018	09:08 am	04/05/2019
11426-007-SA	0	137-002	SP2-7	EPA 1613 D/F	Solid	04/05/2018	09:14 am	04/05/2019
11426-008-SA	0	137-002	SP2-8	EPA 1613 D/F	Solid	04/05/2018	09:18 am	04/05/2019
11426-009-SA	0	137-002	SP2-9	EPA 1613 D/F	Solid	04/05/2018	09:21 am	04/05/2019
11426-010-SA	0	137-002	SP2-10	EPA 1613 D/F	Solid	04/05/2018	09:24 am	04/05/2019
11426-011-SA	0	137-002	SP2-11	EPA 1613 D/F	Solid	04/05/2018	09:26 am	04/05/2019
11426-012-SA	0	137-002	SP2-12	EPA 1613 D/F	Solid	04/05/2018	09:29 am	04/05/2019

EPA Method 1613 PCDD/F



FAL ID: 11426-001-MB
Client ID: Method Blank
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-20-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.123		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.223		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.289		-	0.0503				
1,2,3,6,7,8-HxCDD	ND	0.294		-	0.0490	Total TCDD	ND	0.123	
1,2,3,7,8,9-HxCDD	ND	0.282		-	0.0488	Total PeCDD	ND	0.223	
1,2,3,4,6,7,8-HpCDD	ND	0.426		-	0.0541	Total HxCDD	ND	0.294	
OCDD	ND	0.812		-	0.0888	Total HpCDD	ND	0.426	
2,3,7,8-TCDF	ND	0.109		-	0.0243				
1,2,3,7,8-PeCDF	ND	0.179		-	0.0285				
2,3,4,7,8-PeCDF	ND	0.192		-	0.0298				
1,2,3,4,7,8-HxCDF	ND	0.144		-	0.0255				
1,2,3,6,7,8-HxCDF	ND	0.146		-	0.0253				
2,3,4,6,7,8-HxCDF	ND	0.160		-	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.202		-	0.0367	Total TCDF	ND	0.109	
1,2,3,4,6,7,8-HpCDF	ND	0.207		-	0.0321	Total PeCDF	ND	0.192	
1,2,3,4,7,8,9-HpCDF	ND	0.270		-	0.0396	Total HxCDF	ND	0.202	
OCDF	ND	0.418		-	0.0843	Total HpCDF	ND	0.270	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	98.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	85.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	98.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	82.1	23.0 - 140	
13C-OCDD	70.4	17.0 - 157	
13C-2,3,7,8-TCDF	98.2	24.0 - 169	
13C-1,2,3,7,8-PeCDF	80.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	79.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.0	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	92.2	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	96.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	83.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	94.6	26.0 - 138	
13C-OCDF	75.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 92.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000003 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-001-OPR
Client ID: OPR
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-20-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	11.7	6.70 - 15.8	
1,2,3,7,8-PeCDD	53.0	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	53.3	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	54.1	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	56.4	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	55.6	35.0 - 70.0	
OCDD	110	78.0 - 144	
2,3,7,8-TCDF	11.9	7.50 - 15.8	
1,2,3,7,8-PeCDF	52.4	40.0 - 67.0	
2,3,4,7,8-PeCDF	54.1	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	50.7	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	51.2	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	53.1	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	50.1	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	56.1	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	55.0	39.0 - 69.0	
OCDF	105	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.6	20.0 - 175	
13C-1,2,3,7,8-PeCDD	89.0	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	86.6	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	97.1	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	79.5	26.0 - 166	
13C-OCDD	63.1	13.0 - 198	
13C-2,3,7,8-TCDF	96.6	22.0 - 152	
13C-1,2,3,7,8-PeCDF	81.9	21.0 - 192	
13C-2,3,4,7,8-PeCDF	79.1	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	94.5	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	99.5	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	91.2	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	96.8	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	81.2	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	91.3	20.0 - 186	
13C-OCDF	70.2	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	90.8	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 4/25/2018

Reviewed By: 
Date: 4/26/2018

000004 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-001-SA
Client ID: SP2-1
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.06 g
% Solids: 93.19

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 1.77
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.153		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.326		-	0.0468				
1,2,3,4,7,8-HxCDD	0.545	-	J	0.0545	0.0503				
1,2,3,6,7,8-HxCDD	1.66	-	J	0.166	0.0490	Total TCDD	0.439	-	J
1,2,3,7,8,9-HxCDD	1.01	-	J	0.101	0.0488	Total PeCDD	1.26	-	J,M
1,2,3,4,6,7,8-HpCDD	40.1	-		0.401	0.0541	Total HxCDD	14.4	-	
OCDD	334	-		0.100	0.0888	Total HpCDD	80.5	-	
2,3,7,8-TCDF	1.59	-		0.159	0.0243				
1,2,3,7,8-PeCDF	0.430	-	J	0.0129	0.0285				
2,3,4,7,8-PeCDF	0.859	-	J	0.258	0.0298				
1,2,3,4,7,8-HxCDF	1.18	-	J	0.118	0.0255				
1,2,3,6,7,8-HxCDF	1.50	-	J	0.150	0.0253				
2,3,4,6,7,8-HxCDF	0.823	-	J	0.0823	0.0279				
1,2,3,7,8,9-HxCDF	0.471	-	J	0.0471	0.0367	Total TCDF	10.4	-	
1,2,3,4,6,7,8-HpCDF	9.74	-		0.0974	0.0321	Total PeCDF	7.53	-	
1,2,3,4,7,8,9-HpCDF	1.54	-	J	0.0154	0.0396	Total HxCDF	10.8	-	
OCDF	20.3	-		0.00609	0.0843	Total HpCDF	21.5	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	81.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	83.5	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	84.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	61.8	23.0 - 140	
13C-OCDD	45.5	17.0 - 157	
13C-2,3,7,8-TCDF	88.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	75.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	73.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	83.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	78.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	80.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	70.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	75.2	26.0 - 138	
13C-OCDF	48.6	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 88.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000005 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-002-SA
Client ID: SP2-2
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.00 g
% Solids: 95.91

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 4.65
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.176		-	0.0315				
1,2,3,7,8-PeCDD	0.494	-	J	0.494	0.0468				
1,2,3,4,7,8-HxCDD	1.03	-	J	0.103	0.0503				
1,2,3,6,7,8-HxCDD	4.52	-	J	0.452	0.0490	Total TCDD	0.836	-	J
1,2,3,7,8,9-HxCDD	2.42	-	J	0.242	0.0488	Total PeCDD	5.77	-	M
1,2,3,4,6,7,8-HpCDD	123	-		1.23	0.0541	Total HxCDD	35.3	-	
OCDD	1110	-		0.333	0.0888	Total HpCDD	236	-	
2,3,7,8-TCDF	2.17	-	F	0.217	0.0243				
1,2,3,7,8-PeCDF	0.818	-	J	0.0245	0.0285				
2,3,4,7,8-PeCDF	1.38	-	J	0.414	0.0298				
1,2,3,4,7,8-HxCDF	2.55	-	J	0.255	0.0255				
1,2,3,6,7,8-HxCDF	3.14	-	J	0.314	0.0253				
2,3,4,6,7,8-HxCDF	1.76	-	J	0.176	0.0279				
1,2,3,7,8,9-HxCDF	1.01	-	J	0.101	0.0367	Total TCDF	17.3	-	
1,2,3,4,6,7,8-HpCDF	24.3	-		0.243	0.0321	Total PeCDF	12.6	-	
1,2,3,4,7,8,9-HpCDF	4.09	-	J	0.0409	0.0396	Total HxCDF	22.7	-	
OCDF	51.3	-		0.0154	0.0843	Total HpCDF	54.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	102	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	95.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	81.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	73.0	23.0 - 140	
13C-OCDD	60.3	17.0 - 157	
13C-2,3,7,8-TCDF	102	24.0 - 169	
13C-1,2,3,7,8-PeCDF	83.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	81.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	82.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	83.4	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	84.9	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	83.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	72.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	82.3	26.0 - 138	
13C-OCDF	60.6	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 94.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000006 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-003-SA
Client ID: SP2-3
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.03 g
% Solids: 91.60

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.73
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.154		-	0.0315				
1,2,3,7,8-PeCDD	0.568	-	J	0.568	0.0468				
1,2,3,4,7,8-HxCDD	0.857	-	J	0.0857	0.0503				
1,2,3,6,7,8-HxCDD	1.98	-	J	0.198	0.0490	Total TCDD	0.652	-	J
1,2,3,7,8,9-HxCDD	1.51	-	J	0.151	0.0488	Total PeCDD	3.75	-	J,M
1,2,3,4,6,7,8-HpCDD	41.2	-		0.412	0.0541	Total HxCDD	18.4	-	
OCDD	301	-		0.0903	0.0888	Total HpCDD	79.8	-	
2,3,7,8-TCDF	2.72	-	F	0.272	0.0243				
1,2,3,7,8-PeCDF	0.515	-	J	0.0155	0.0285				
2,3,4,7,8-PeCDF	1.57	-	J	0.471	0.0298				
1,2,3,4,7,8-HxCDF	1.42	-	J	0.142	0.0255				
1,2,3,6,7,8-HxCDF	0.828	-	J	0.0828	0.0253				
2,3,4,6,7,8-HxCDF	0.938	-	J	0.0938	0.0279				
1,2,3,7,8,9-HxCDF	0.478	-	J	0.0478	0.0367	Total TCDF	15.1	-	M
1,2,3,4,6,7,8-HpCDF	9.19	-		0.0919	0.0321	Total PeCDF	12.8	-	
1,2,3,4,7,8,9-HpCDF	0.629	-	J	0.00629	0.0396	Total HxCDF	16.2	-	
OCDF	18.8	-		0.00564	0.0843	Total HpCDF	24.5	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	101	25.0 - 164	
13C-1,2,3,7,8-PeCDD	87.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	89.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.2	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	80.5	23.0 - 140	
13C-OCDD	71.2	17.0 - 157	
13C-2,3,7,8-TCDF	98.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	78.4	24.0 - 185	
13C-2,3,4,7,8-PeCDF	77.8	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	98.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	100	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	94.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	98.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	83.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	94.7	26.0 - 138	
13C-OCDF	73.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 93.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000007 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-004-SA
Client ID: SP2-4
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.04 g
% Solids: 92.56

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.96
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.191		-	0.0315				
1,2,3,7,8-PeCDD	0.588	-	J	0.588	0.0468				
1,2,3,4,7,8-HxCDD	0.810	-	J	0.0810	0.0503				
1,2,3,6,7,8-HxCDD	2.24	-	J	0.224	0.0490	Total TCDD	0.388	-	J
1,2,3,7,8,9-HxCDD	1.42	-	J	0.142	0.0488	Total PeCDD	3.70	-	J,M
1,2,3,4,6,7,8-HpCDD	56.2	-		0.562	0.0541	Total HxCDD	18.2	-	
OCDD	495	-		0.149	0.0888	Total HpCDD	105	-	
2,3,7,8-TCDF	1.93	-		0.193	0.0243				
1,2,3,7,8-PeCDF	0.497	-	J	0.0149	0.0285				
2,3,4,7,8-PeCDF	1.33	-	J	0.399	0.0298				
1,2,3,4,7,8-HxCDF	1.65	-	J	0.165	0.0255				
1,2,3,6,7,8-HxCDF	1.20	-	J	0.120	0.0253				
2,3,4,6,7,8-HxCDF	0.978	-	J	0.0978	0.0279				
1,2,3,7,8,9-HxCDF	0.563	-	J	0.0563	0.0367	Total TCDF	12.4	-	M
1,2,3,4,6,7,8-HpCDF	13.6	-		0.136	0.0321	Total PeCDF	10.8	-	
1,2,3,4,7,8,9-HpCDF	1.58	-	J	0.0158	0.0396	Total HxCDF	16.5	-	
OCDF	40.7	-		0.0122	0.0843	Total HpCDF	38.4	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	101	25.0 - 164	
13C-1,2,3,7,8-PeCDD	96.8	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	91.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	92.4	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	79.7	23.0 - 140	
13C-OCDD	67.8	17.0 - 157	
13C-2,3,7,8-TCDF	101	24.0 - 169	
13C-1,2,3,7,8-PeCDF	89.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	87.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	95.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	93.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	94.1	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	80.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.1	26.0 - 138	
13C-OCDF	69.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 94.1 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000008 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-005-SA
Client ID: SP2-5
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.04 g
% Solids: 94.36

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 1.64
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.172		-	0.0315				
1,2,3,7,8-PeCDD	0.341	-	J	0.341	0.0468				
1,2,3,4,7,8-HxCDD	0.445	-	J	0.0445	0.0503				
1,2,3,6,7,8-HxCDD	1.30	-	J	0.130	0.0490	Total TCDD	0.322	-	J
1,2,3,7,8,9-HxCDD	0.941	-	J	0.0941	0.0488	Total PeCDD	2.50	-	J,M
1,2,3,4,6,7,8-HpCDD	33.3	-		0.333	0.0541	Total HxCDD	12.5	-	
OCDD	280	-		0.0840	0.0888	Total HpCDD	64.2	-	
2,3,7,8-TCDF	1.38	-		0.138	0.0243				
1,2,3,7,8-PeCDF	0.322	-	J	0.00966	0.0285				
2,3,4,7,8-PeCDF	0.716	-	J	0.215	0.0298				
1,2,3,4,7,8-HxCDF	0.572	-	J	0.0572	0.0255				
1,2,3,6,7,8-HxCDF	0.462	-	J	0.0462	0.0253				
2,3,4,6,7,8-HxCDF	0.530	-	J	0.0530	0.0279				
1,2,3,7,8,9-HxCDF	0.287	-	J	0.0287	0.0367	Total TCDF	8.35	-	M
1,2,3,4,6,7,8-HpCDF	5.46	-		0.0546	0.0321	Total PeCDF	6.62	-	
1,2,3,4,7,8,9-HpCDF	0.530	-	J	0.00530	0.0396	Total HxCDF	8.96	-	
OCDF	12.0	-		0.00360	0.0843	Total HpCDF	14.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	99.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	98.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	85.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	91.1	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	74.9	23.0 - 140	
13C-OCDD	65.3	17.0 - 157	
13C-2,3,7,8-TCDF	97.7	24.0 - 169	
13C-1,2,3,7,8-PeCDF	90.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	84.3	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.3	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	97.0	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.8	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	91.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	81.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	85.7	26.0 - 138	
13C-OCDF	65.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 95.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000009 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-006-SA
Client ID: SP2-6
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.01 g
% Solids: 92.42

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.71
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.116		-	0.0315				
1,2,3,7,8-PeCDD	0.596	-	J	0.596	0.0468				
1,2,3,4,7,8-HxCDD	0.431	-	J	0.0431	0.0503				
1,2,3,6,7,8-HxCDD	1.62	-	J	0.162	0.0490	Total TCDD	2.23	-	
1,2,3,7,8,9-HxCDD	1.08	-	J	0.108	0.0488	Total PeCDD	3.86		J,M
1,2,3,4,6,7,8-HpCDD	32.3	-		0.323	0.0541	Total HxCDD	14.8	-	
OCDD	257	-		0.0771	0.0888	Total HpCDD	63.0	-	
2,3,7,8-TCDF	1.95	-		0.195	0.0243				
1,2,3,7,8-PeCDF	0.532	-	J	0.0160	0.0285				
2,3,4,7,8-PeCDF	1.29	-	J	0.387	0.0298				
1,2,3,4,7,8-HxCDF	2.13	-	J	0.213	0.0255				
1,2,3,6,7,8-HxCDF	2.19	-	J	0.219	0.0253				
2,3,4,6,7,8-HxCDF	1.23	-	J	0.123	0.0279				
1,2,3,7,8,9-HpCDF	0.780	-	J	0.0780	0.0367	Total TCDF	12.7	-	
1,2,3,4,6,7,8-HpCDF	13.5	-		0.135	0.0321	Total PeCDF	10.5	-	
1,2,3,4,7,8,9-HpCDF	2.59	-	J	0.0259	0.0396	Total HxCDF	14.2	-	
OCDF	20.3	-		0.00609	0.0843	Total HpCDF	25.4	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	100	25.0 - 164	
13C-1,2,3,7,8-PeCDD	85.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	93.0	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	81.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	68.3	23.0 - 140	
13C-OCDD	56.9	17.0 - 157	
13C-2,3,7,8-TCDF	106	24.0 - 169	
13C-1,2,3,7,8-PeCDF	81.9	24.0 - 185	
13C-2,3,4,7,8-PeCDF	80.3	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	84.6	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	86.4	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.1	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	83.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	71.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	78.7	26.0 - 138	
13C-OCDF	57.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 97.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000010 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-007-SA
Client ID: SP2-7
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.04 g
% Solids: 94.30

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.45
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.186		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.233		-	0.0468				
1,2,3,4,7,8-HxCDD	0.351	-	J	0.0351	0.0503				
1,2,3,6,7,8-HxCDD	0.685	-	J	0.0685	0.0490	Total TCDD	0.371	-	J
1,2,3,7,8,9-HxCDD	0.618	-	J	0.0618	0.0488	Total PeCDD	0.969	-	J,M
1,2,3,4,6,7,8-HpCDD	15.4	-		0.154	0.0541	Total HxCDD	6.62	-	
OCDD	126	-		0.0378	0.0888	Total HpCDD	30.5	-	
2,3,7,8-TCDF	1.66	-		0.166	0.0243				
1,2,3,7,8-PeCDF	0.618	-	J	0.0185	0.0285				
2,3,4,7,8-PeCDF	2.70	-	J	0.810	0.0298				
1,2,3,4,7,8-HxCDF	5.26	-		0.526	0.0255				
1,2,3,6,7,8-HxCDF	1.83	-	J	0.183	0.0253				
2,3,4,6,7,8-HxCDF	1.55	-	J	0.155	0.0279				
1,2,3,7,8,9-HxCDF	1.56	-	J	0.156	0.0367	Total TCDF	12.8	-	
1,2,3,4,6,7,8-HpCDF	5.68	-		0.0568	0.0321	Total PeCDF	15.3	-	
1,2,3,4,7,8,9-HpCDF	1.45	-	J	0.0145	0.0396	Total HxCDF	19.6	-	
OCDF	11.6	-		0.00348	0.0843	Total HpCDF	15.2	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	99.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	91.2	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	80.9	23.0 - 140	
13C-OCDD	69.5	17.0 - 157	
13C-2,3,7,8-TCDF	103	24.0 - 169	
13C-1,2,3,7,8-PeCDF	85.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	82.8	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	96.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	104	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	97.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	99.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	85.3	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	93.8	26.0 - 138	
13C-OCDF	71.5	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 88.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000011 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-008-SA
Client ID: SP2-8
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.01 g
% Solids: 96.02

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.57
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.212		-	0.0315				
1,2,3,7,8-PeCDD	0.487	-	J	0.487	0.0468				
1,2,3,4,7,8-HxCDD	0.602	-	J	0.0602	0.0503				
1,2,3,6,7,8-HxCDD	1.78	-	J	0.178	0.0490	Total TCDD	0.491	-	J
1,2,3,7,8,9-HxCDD	1.22	-	J	0.122	0.0488	Total PeCDD	3.77	-	J,M
1,2,3,4,6,7,8-HpCDD	41.4	-		0.414	0.0541	Total HxCDD	15.4	-	
OCDD	358	-		0.107	0.0888	Total HpCDD	79.8	-	
2,3,7,8-TCDF	2.47	-	F	0.247	0.0243				
1,2,3,7,8-PeCDF	0.486	-	J	0.0146	0.0285				
2,3,4,7,8-PeCDF	1.40	-	J	0.420	0.0298				
1,2,3,4,7,8-HxCDF	1.67	-	J	0.167	0.0255				
1,2,3,6,7,8-HxCDF	1.17	-	J	0.117	0.0253				
2,3,4,6,7,8-HxCDF	0.815	-	J	0.0815	0.0279				
1,2,3,7,8,9-HxCDF	0.489	-	J	0.0489	0.0367	Total TCDF	13.9	-	
1,2,3,4,6,7,8-HpCDF	8.92	-		0.0892	0.0321	Total PeCDF	10.6	-	
1,2,3,4,7,8,9-HpCDF	1.21	-	J	0.0121	0.0396	Total HxCDF	13.3	-	
OCDF	24.3	-		0.00729	0.0843	Total HpCDF	24.3	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.2	25.0 - 164	
13C-1,2,3,7,8-PeCDD	79.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	85.6	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	87.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	71.6	23.0 - 140	
13C-OCDD	64.6	17.0 - 157	
13C-2,3,7,8-TCDF	99.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	76.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	72.3	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	94.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	91.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	88.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	79.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.5	26.0 - 138	
13C-OCDF	64.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 86.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000012 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-009-SA
Client ID: SP2-9
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.00 g
% Solids: 93.81

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 3.43
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.165		-	0.0315				
1,2,3,7,8-PeCDD	0.809	-	J	0.809	0.0468				
1,2,3,4,7,8-HxCDD	0.719	-	J	0.0719	0.0503				
1,2,3,6,7,8-HxCDD	2.06	-	J	0.206	0.0490	Total TCDD	0.523	-	J
1,2,3,7,8,9-HxCDD	1.67	-	J	0.167	0.0488	Total PeCDD	4.41	-	J,M
1,2,3,4,6,7,8-HpCDD	46.6	-		0.466	0.0541	Total HxCDD	21.4	-	
OCDD	309	-		0.0927	0.0888	Total HpCDD	90.0	-	
2,3,7,8-TCDF	4.32	-	F	0.432	0.0243				
1,2,3,7,8-PeCDF	0.699	-	J	0.0210	0.0285				
2,3,4,7,8-PeCDF	1.99	-	J	0.597	0.0298				
1,2,3,4,7,8-HxCDF	1.96	-	J	0.196	0.0255				
1,2,3,6,7,8-HxCDF	1.15	-	J	0.115	0.0253				
2,3,4,6,7,8-HxCDF	1.04	-	J	0.104	0.0279				
1,2,3,7,8,9-HxCDF	0.505	-	J	0.0505	0.0367	Total TCDF	21.8	-	M
1,2,3,4,6,7,8-HpCDF	9.07	-		0.0907	0.0321	Total PeCDF	15.5	-	
1,2,3,4,7,8,9-HpCDF	0.828	-	J	0.00828	0.0396	Total HxCDF	15.3	-	
OCDF	14.0	-		0.00420	0.0843	Total HpCDF	20.5	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	84.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	81.2	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	90.3	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	74.0	23.0 - 140	
13C-OCDD	62.8	17.0 - 157	
13C-2,3,7,8-TCDF	95.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	77.4	24.0 - 185	
13C-2,3,4,7,8-PeCDF	76.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	91.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	91.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	92.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	80.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	86.6	26.0 - 138	
13C-OCDF	67.6	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.7 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000013 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-010-SA
Client ID: SP2-10
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.01 g
% Solids: 92.69

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 1.08
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.170		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.244		-	0.0468				
1,2,3,4,7,8-HxCDD	ND	0.267		-	0.0503				
1,2,3,6,7,8-HxCDD	0.772	-	J	0.0772	0.0490	Total TCDD	ND	0.170	
1,2,3,7,8,9-HxCDD	0.515	-	J	0.0515	0.0488	Total PeCDD	0.350	-	J
1,2,3,4,6,7,8-HpCDD	15.3	-		0.153	0.0541	Total HxCDD	5.24	-	
OCDD	132	-		0.0396	0.0888	Total HpCDD	28.9	-	
2,3,7,8-TCDF	1.34	-		0.134	0.0243				
1,2,3,7,8-PeCDF	0.313	-	J	0.00939	0.0285				
2,3,4,7,8-PeCDF	0.758	-	J	0.227	0.0298				
1,2,3,4,7,8-HxCDF	1.14	-	J	0.114	0.0255				
1,2,3,6,7,8-HxCDF	0.933	-	J	0.0933	0.0253				
2,3,4,6,7,8-HxCDF	0.617	-	J	0.0617	0.0279				
1,2,3,7,8,9-HxCDF	0.453	-	J	0.0453	0.0367	Total TCDF	7.51	-	
1,2,3,4,6,7,8-HpCDF	6.46	-		0.0646	0.0321	Total PeCDF	5.33	-	
1,2,3,4,7,8,9-HpCDF	0.797	-	J	0.00797	0.0396	Total HxCDF	7.31	-	
OCDF	8.97	-	J	0.00269	0.0843	Total HpCDF	12.2	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	98.4	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	92.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	90.7	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	73.5	23.0 - 140	
13C-OCDD	60.7	17.0 - 157	
13C-2,3,7,8-TCDF	101	24.0 - 169	
13C-1,2,3,7,8-PeCDF	82.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	82.7	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	94.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	96.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	95.2	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	95.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	80.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.7	26.0 - 138	
13C-OCDF	66.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 97.8 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000014 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-011-SA
Client ID: SP2-11
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.01 g
% Solids: 91.92

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 1.75
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.189		-	0.0315				
1,2,3,7,8-PeCDD	0.401	-	J	0.401	0.0468				
1,2,3,4,7,8-HxCDD	0.496	-	J	0.0496	0.0503				
1,2,3,6,7,8-HxCDD	1.42	-	J	0.142	0.0490	Total TCDD	ND	0.189	
1,2,3,7,8,9-HxCDD	0.911	-	J	0.0911	0.0488	Total PeCDD	3.26	-	J,M
1,2,3,4,6,7,8-HpCDD	31.0	-		0.310	0.0541	Total HxCDD	13.6	-	
OCDD	262	-		0.0786	0.0888	Total HpCDD	61.4	-	
2,3,7,8-TCDF	1.66	-		0.166	0.0243				
1,2,3,7,8-PeCDF	0.294	-	J	0.00882	0.0285				
2,3,4,7,8-PeCDF	0.785	-	J	0.236	0.0298				
1,2,3,4,7,8-HxCDF	0.711	-	J	0.0711	0.0255				
1,2,3,6,7,8-HxCDF	0.617	-	J	0.0617	0.0253				
2,3,4,6,7,8-HxCDF	0.612	-	J	0.0612	0.0279				
1,2,3,7,8,9-HxCDF	ND	0.347		-	0.0367	Total TCDF	9.31	-	
1,2,3,4,6,7,8-HpCDF	6.04	-		0.0604	0.0321	Total PeCDF	7.08	-	
1,2,3,4,7,8,9-HpCDF	0.628	-	J	0.00628	0.0396	Total HxCDF	8.60	-	
OCDF	14.0	-		0.00420	0.0843	Total HpCDF	15.9	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	99.4	25.0 - 164	
13C-1,2,3,7,8-PeCDD	102	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	91.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	97.4	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	81.3	23.0 - 140	
13C-OCDD	72.6	17.0 - 157	
13C-2,3,7,8-TCDF	105	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.5	24.0 - 185	
13C-2,3,4,7,8-PeCDF	90.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	101	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	110	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	103	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	103	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	93.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	101	26.0 - 138	
13C-OCDF	77.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 93.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000015 of 000020

EPA Method 1613 PCDD/F



FAL ID: 11426-012-SA
Client ID: SP2-12
Matrix: Solid
Batch No: X4479

Date Extracted: 04-19-2018
Date Received: 04-05-2018
Amount: 5.01 g
% Solids: 93.26

ICal: PCDDFAL3-12-22-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-21-2018
2005 WHO TEQ: 2.21
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.199		-	0.0315				
1,2,3,7,8-PeCDD	ND	0.308		-	0.0468				
1,2,3,4,7,8-HxCDD	0.444	-	J	0.0444	0.0503				
1,2,3,6,7,8-HxCDD	1.31	-	J	0.131	0.0490	Total TCDD	0.457	-	J
1,2,3,7,8,9-HxCDD	0.957	-	J	0.0957	0.0488	Total PeCDD	1.40	-	J,M
1,2,3,4,6,7,8-HpCDD	25.9	-		0.259	0.0541	Total HxCDD	11.1	-	
OCDD	225	-		0.0675	0.0888	Total HpCDD	50.4	-	
2,3,7,8-TCDF	4.13	-	F	0.413	0.0243				
1,2,3,7,8-PeCDF	0.542	-	J	0.0163	0.0285				
2,3,4,7,8-PeCDF	2.04	-	J	0.612	0.0298				
1,2,3,4,7,8-HxCDF	2.32	-	J	0.232	0.0255				
1,2,3,6,7,8-HxCDF	0.924	-	J	0.0924	0.0253				
2,3,4,6,7,8-HxCDF	0.986	-	J	0.0986	0.0279				
1,2,3,7,8,9-HxCDF	0.638	-	J	0.0638	0.0367	Total TCDF	19.5	-	
1,2,3,4,6,7,8-HpCDF	7.25	-		0.0725	0.0321	Total PeCDF	12.5	-	
1,2,3,4,7,8,9-HpCDF	0.813	-	J	0.00813	0.0396	Total HxCDF	14.0	-	
OCDF	15.4	-		0.00462	0.0843	Total HpCDF	18.2	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	87.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	83.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	81.8	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	86.3	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	69.5	23.0 - 140	
13C-OCDD	53.7	17.0 - 157	
13C-2,3,7,8-TCDF	89.5	24.0 - 169	
13C-1,2,3,7,8-PeCDF	73.1	24.0 - 185	
13C-2,3,4,7,8-PeCDF	71.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.9	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.0	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	90.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	81.5	26.0 - 138	
13C-OCDF	59.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 4/25/2018

Reviewed By: 

Date: 4/26/2018

000016 of 000020

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed. FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11426**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/05/2018
Time Received:	01:46 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	12
Duplicates:	0
Storage Location:	R-4

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	N/A
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



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13 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/06/18 10:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP2-1	T181179-01	Soil	04/05/18 08:54	04/06/18 10:00
SP2-2	T181179-02	Soil	04/05/18 08:57	04/06/18 10:00
SP2-3	T181179-03	Soil	04/05/18 08:59	04/06/18 10:00
SP2-4	T181179-04	Soil	04/05/18 09:02	04/06/18 10:00
SP2-5	T181179-05	Soil	04/05/18 09:04	04/06/18 10:00
SP2-6	T181179-06	Soil	04/05/18 09:08	04/06/18 10:00
SP2-7	T181179-07	Soil	04/05/18 09:14	04/06/18 10:00
SP2-8	T181179-08	Soil	04/05/18 09:18	04/06/18 10:00
SP2-9	T181179-09	Soil	04/05/18 09:21	04/06/18 10:00
SP2-10	T181179-10	Soil	04/05/18 09:24	04/06/18 10:00
SP2-11	T181179-11	Soil	04/05/18 09:26	04/06/18 10:00
SP2-12	T181179-12	Soil	04/05/18 09:29	04/06/18 10:00

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

DETECTIONS SUMMARY

Sample ID: SP2-1

Laboratory ID: T181179-01

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	62	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	59	10	mg/kg	EPA 8015B	
Barium	58	1.0	mg/kg	EPA 6010B	
Chromium	6.2	2.0	mg/kg	EPA 6010B	
Cobalt	4.9	2.0	mg/kg	EPA 6010B	
Copper	5.2	1.0	mg/kg	EPA 6010B	
Lead	13	3.0	mg/kg	EPA 6010B	
Nickel	5.9	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	33	1.0	mg/kg	EPA 6010B	
Mercury	0.52	0.10	mg/kg	EPA 7471A Soil	
PCB-1260	29	10	ug/kg	EPA 8082	

Sample ID: SP2-2

Laboratory ID: T181179-02

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	24	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	42	10	mg/kg	EPA 8015B	
Barium	55	1.0	mg/kg	EPA 6010B	
Chromium	5.5	2.0	mg/kg	EPA 6010B	
Cobalt	4.7	2.0	mg/kg	EPA 6010B	
Copper	3.5	1.0	mg/kg	EPA 6010B	
Lead	11	3.0	mg/kg	EPA 6010B	
Nickel	5.3	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	29	1.0	mg/kg	EPA 6010B	
PCB-1016	66	10	ug/kg	EPA 8082	
PCB-1260	72	10	ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Sample ID: SP2-3

Laboratory ID: T181179-03

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	18	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	55	10		mg/kg	EPA 8015B	
Barium	63	1.0		mg/kg	EPA 6010B	
Chromium	6.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.2	2.0		mg/kg	EPA 6010B	
Copper	5.5	1.0		mg/kg	EPA 6010B	
Nickel	6.2	2.0		mg/kg	EPA 6010B	
Vanadium	22	5.0		mg/kg	EPA 6010B	
Zinc	33	1.0		mg/kg	EPA 6010B	
PCB-1254	36	10		ug/kg	EPA 8082	

Sample ID: SP2-4

Laboratory ID: T181179-04

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	94	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	69	10		mg/kg	EPA 8015B	
Barium	62	1.0		mg/kg	EPA 6010B	
Chromium	6.5	2.0		mg/kg	EPA 6010B	
Cobalt	4.9	2.0		mg/kg	EPA 6010B	
Copper	5.5	1.0		mg/kg	EPA 6010B	
Lead	9.8	3.0		mg/kg	EPA 6010B	
Nickel	6.6	2.0		mg/kg	EPA 6010B	
Vanadium	22	5.0		mg/kg	EPA 6010B	
Zinc	30	1.0		mg/kg	EPA 6010B	
PCB-1260	77	10		ug/kg	EPA 8082	

Sample ID: SP2-5

Laboratory ID: T181179-05

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	51	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	46	10		mg/kg	EPA 8015B	
Barium	59	1.0		mg/kg	EPA 6010B	
Chromium	6.2	2.0		mg/kg	EPA 6010B	
Cobalt	4.4	2.0		mg/kg	EPA 6010B	
Copper	5.6	1.0		mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Sample ID: SP2-5

Laboratory ID: T181179-05

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Lead	17	3.0		mg/kg	EPA 6010B	
Nickel	5.9	2.0		mg/kg	EPA 6010B	
Vanadium	18	5.0		mg/kg	EPA 6010B	
Zinc	31	1.0		mg/kg	EPA 6010B	
PCB-1260	32	10		ug/kg	EPA 8082	
Pyrene	950	300		ug/kg	EPA 8270C	
Benzo (a) anthracene	340	300		ug/kg	EPA 8270C	
Chrysene	440	300		ug/kg	EPA 8270C	
Fluoranthene	880	300		ug/kg	EPA 8270C	
Phenanthrene	530	300		ug/kg	EPA 8270C	

Sample ID: SP2-6

Laboratory ID: T181179-06

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	90	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	49	10		mg/kg	EPA 8015B	
Barium	68	1.0		mg/kg	EPA 6010B	
Chromium	7.1	2.0		mg/kg	EPA 6010B	
Cobalt	5.0	2.0		mg/kg	EPA 6010B	
Copper	5.7	1.0		mg/kg	EPA 6010B	
Lead	13	3.0		mg/kg	EPA 6010B	
Nickel	7.5	2.0		mg/kg	EPA 6010B	
Vanadium	24	5.0		mg/kg	EPA 6010B	
Zinc	33	1.0		mg/kg	EPA 6010B	
PCB-1260	58	10		ug/kg	EPA 8082	
Pyrene	1200	300		ug/kg	EPA 8270C	
Benzo (a) anthracene	490	300		ug/kg	EPA 8270C	
Benzo (b) fluoranthene	410	300		ug/kg	EPA 8270C	
Chrysene	570	300		ug/kg	EPA 8270C	
Fluoranthene	1100	300		ug/kg	EPA 8270C	
Phenanthrene	750	300		ug/kg	EPA 8270C	

Sample ID: SP2-7

Laboratory ID: T181179-07

Analyte	Result	Reporting		Units	Method	Notes
		Limit				

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Sample ID: SP2-7

Laboratory ID: T181179-07

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	60	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	57	10		mg/kg	EPA 8015B	
Barium	58	1.0		mg/kg	EPA 6010B	
Chromium	6.0	2.0		mg/kg	EPA 6010B	
Cobalt	4.6	2.0		mg/kg	EPA 6010B	
Copper	4.1	1.0		mg/kg	EPA 6010B	
Nickel	6.4	2.0		mg/kg	EPA 6010B	
Vanadium	22	5.0		mg/kg	EPA 6010B	
Zinc	24	1.0		mg/kg	EPA 6010B	
PCB-1260	86	10		ug/kg	EPA 8082	

Sample ID: SP2-8

Laboratory ID: T181179-08

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	170	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	120	10		mg/kg	EPA 8015B	
Barium	61	1.0		mg/kg	EPA 6010B	
Chromium	5.5	2.0		mg/kg	EPA 6010B	
Cobalt	4.4	2.0		mg/kg	EPA 6010B	
Copper	6.5	1.0		mg/kg	EPA 6010B	
Lead	11	3.0		mg/kg	EPA 6010B	
Nickel	5.1	2.0		mg/kg	EPA 6010B	
Vanadium	16	5.0		mg/kg	EPA 6010B	
Zinc	27	1.0		mg/kg	EPA 6010B	
PCB-1260	58	10		ug/kg	EPA 8082	

Sample ID: SP2-9

Laboratory ID: T181179-09

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	370	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	380	10		mg/kg	EPA 8015B	
Barium	52	1.0		mg/kg	EPA 6010B	
Chromium	5.5	2.0		mg/kg	EPA 6010B	
Cobalt	5.6	2.0		mg/kg	EPA 6010B	
Copper	8.5	1.0		mg/kg	EPA 6010B	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Sample ID: SP2-9

Laboratory ID: T181179-09

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Lead	32	3.0	mg/kg	EPA 6010B	
Nickel	9.5	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	
PCB-1260	58	10	ug/kg	EPA 8082	

Sample ID: SP2-10

Laboratory ID: T181179-10

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	150	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	76	10	mg/kg	EPA 8015B	
Barium	61	1.0	mg/kg	EPA 6010B	
Chromium	6.1	2.0	mg/kg	EPA 6010B	
Cobalt	4.6	2.0	mg/kg	EPA 6010B	
Copper	4.1	1.0	mg/kg	EPA 6010B	
Nickel	6.2	2.0	mg/kg	EPA 6010B	
Vanadium	21	5.0	mg/kg	EPA 6010B	
Zinc	26	1.0	mg/kg	EPA 6010B	
PCB-1016	72	10	ug/kg	EPA 8082	
PCB-1260	45	10	ug/kg	EPA 8082	

Sample ID: SP2-11

Laboratory ID: T181179-11

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	110	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	68	10	mg/kg	EPA 8015B	
Barium	58	1.0	mg/kg	EPA 6010B	
Chromium	6.7	2.0	mg/kg	EPA 6010B	
Cobalt	4.4	2.0	mg/kg	EPA 6010B	
Copper	6.5	1.0	mg/kg	EPA 6010B	
Lead	17	3.0	mg/kg	EPA 6010B	
Nickel	6.4	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	31	1.0	mg/kg	EPA 6010B	
PCB-1260	26	10	ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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PO Box 332
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Sample ID: SP2-12

Laboratory ID: T181179-12

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	140	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	150	10		mg/kg	EPA 8015B	
Barium	58	1.0		mg/kg	EPA 6010B	
Chromium	5.8	2.0		mg/kg	EPA 6010B	
Cobalt	4.5	2.0		mg/kg	EPA 6010B	
Copper	5.8	1.0		mg/kg	EPA 6010B	
Lead	11	3.0		mg/kg	EPA 6010B	
Nickel	5.4	2.0		mg/kg	EPA 6010B	
Vanadium	17	5.0		mg/kg	EPA 6010B	
Zinc	29	1.0		mg/kg	EPA 6010B	
PCB-1260	61	10		ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-1

T181179-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	62	10	"	"	"	"	"	"	
C29-C40 (MORO)	59	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		68.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	58	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.2	2.0	"	"	"	"	"	"	
Cobalt	4.9	2.0	"	"	"	"	"	"	
Copper	5.2	1.0	"	"	"	"	"	"	
Lead	13	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	33	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	0.52	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-1

T181179-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	29	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		134 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		132 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-1

T181179-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-1

T181179-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	46.4 %	15-121			"	"	"	"
Surrogate: Phenol-d6	52.9 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	57.7 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	63.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	86.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	90.3 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-2

T181179-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	24	10	"	"	"	"	"	"	
C29-C40 (MORO)	42	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		80.9 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	55	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.5	2.0	"	"	"	"	"	"	
Cobalt	4.7	2.0	"	"	"	"	"	"	
Copper	3.5	1.0	"	"	"	"	"	"	
Lead	11	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.3	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-2

T181179-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	66	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	72	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		140 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		135 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-2

T181179-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-2

T181179-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	47.9 %	15-121			"	"	"	"
Surrogate: Phenol-d6	52.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	60.8 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	66.9 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	83.2 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	83.5 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-3

T181179-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	18	10	"	"	"	"	"	"	
C29-C40 (MORO)	55	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		70.2 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	63	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.7	2.0	"	"	"	"	"	"	
Cobalt	4.2	2.0	"	"	"	"	"	"	
Copper	5.5	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	22	5.0	"	"	"	"	"	"	
Zinc	33	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-3

T181179-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	36	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		137 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		144 %	35-140		"	"	"	"	S-GC

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-3

T181179-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-3

T181179-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		54.9 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		60.2 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		65.4 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		66.0 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		86.1 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		85.7 %	29.1-130		"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-4

T181179-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	94	10	"	"	"	"	"	"	
C29-C40 (MORO)	69	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		69.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	62	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.5	2.0	"	"	"	"	"	"	
Cobalt	4.9	2.0	"	"	"	"	"	"	
Copper	5.5	1.0	"	"	"	"	"	"	
Lead	9.8	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	22	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-4

T181179-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	77	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		138 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		134 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-4

T181179-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-4

T181179-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		55.5 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		60.4 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		65.1 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		65.4 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		87.0 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		88.3 %	29.1-130		"	"	"	"	

SunStar Laboratories, Inc.

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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-5

T181179-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	51	10	"	"	"	"	"	"	
C29-C40 (MORO)	46	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		83.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	59	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.2	2.0	"	"	"	"	"	"	
Cobalt	4.4	2.0	"	"	"	"	"	"	
Copper	5.6	1.0	"	"	"	"	"	"	
Lead	17	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	18	5.0	"	"	"	"	"	"	
Zinc	31	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

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04/13/18 15:53

SP2-5

T181179-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	32	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		135 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		137 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	950	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	340	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	
Benzo (a) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-5

T181179-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzyl alcohol	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	440	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	880	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	
Isophorone	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-5

T181179-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

2-Methylphenol	ND	1000	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	530	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		54.2 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		59.1 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		62.7 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		66.3 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		90.3 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		89.8 %	29.1-130		"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-6
T181179-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	90	10	"	"	"	"	"	"	
C29-C40 (MORO)	49	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		70.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	68	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	7.1	2.0	"	"	"	"	"	"	
Cobalt	5.0	2.0	"	"	"	"	"	"	
Copper	5.7	1.0	"	"	"	"	"	"	
Lead	13	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	24	5.0	"	"	"	"	"	"	
Zinc	33	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project Number: 137-002
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Reported:
04/13/18 15:53

SP2-6
T181179-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	58	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		137 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		127 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Aniline	ND	300	"	"	"	"	"	"	
Phenol	ND	1000	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	1200	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	490	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	410	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	
Benzo (a) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-6

T181179-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzyl alcohol	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"
4-Chloroaniline	ND	300	"	"	"	"	"	"
2-Chloronaphthalene	ND	300	"	"	"	"	"	"
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"
Chrysene	570	300	"	"	"	"	"	"
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"
Dibenzofuran	ND	300	"	"	"	"	"	"
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"
Diethyl phthalate	ND	300	"	"	"	"	"	"
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"
Dimethyl phthalate	ND	300	"	"	"	"	"	"
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"
Fluoranthene	1100	300	"	"	"	"	"	"
Fluorene	ND	300	"	"	"	"	"	"
Hexachlorobenzene	ND	1500	"	"	"	"	"	"
Hexachlorobutadiene	ND	300	"	"	"	"	"	"
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"
Hexachloroethane	ND	300	"	"	"	"	"	"
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"
Isophorone	ND	300	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-6

T181179-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

2-Methylphenol	ND	1000	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	750	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol		55.3 %	15-121		"	"	"	"
Surrogate: Phenol-d6		60.1 %	24-113		"	"	"	"
Surrogate: Nitrobenzene-d5		65.6 %	21.3-119		"	"	"	"
Surrogate: 2-Fluorobiphenyl		67.2 %	32.4-102		"	"	"	"
Surrogate: 2,4,6-Tribromophenol		92.0 %	18.1-105		"	"	"	"
Surrogate: Terphenyl-d14		88.9 %	29.1-130		"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-7

T181179-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	60	10	"	"	"	"	"	"	
C29-C40 (MORO)	57	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		72.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	58	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.0	2.0	"	"	"	"	"	"	
Cobalt	4.6	2.0	"	"	"	"	"	"	
Copper	4.1	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	22	5.0	"	"	"	"	"	"	
Zinc	24	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-7

T181179-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	86	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		133 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		125 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-7

T181179-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-7

T181179-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	56.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	63.3 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	65.7 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.3 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	90.1 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	94.1 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-8

T181179-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	170	10	"	"	"	"	"	"	
C29-C40 (MORO)	120	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		77.2 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	61	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.5	2.0	"	"	"	"	"	"	
Cobalt	4.4	2.0	"	"	"	"	"	"	
Copper	6.5	1.0	"	"	"	"	"	"	
Lead	11	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	16	5.0	"	"	"	"	"	"	
Zinc	27	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-8

T181179-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	58	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		139 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		134 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-8

T181179-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-8

T181179-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	54.7 %	15-121			"	"	"	"
Surrogate: Phenol-d6	62.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	64.4 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	68.7 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	90.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	96.9 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-9

T181179-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	370	10	"	"	"	"	"	"	
C29-C40 (MORO)	380	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		79.2 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	52	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.5	2.0	"	"	"	"	"	"	
Cobalt	5.6	2.0	"	"	"	"	"	"	
Copper	8.5	1.0	"	"	"	"	"	"	
Lead	32	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	9.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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04/13/18 15:53

SP2-9

T181179-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	58	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		132 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		119 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-9

T181179-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-9

T181179-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol		50.7 %	15-121		"	"	"	"
Surrogate: Phenol-d6		58.9 %	24-113		"	"	"	"
Surrogate: Nitrobenzene-d5		62.2 %	21.3-119		"	"	"	"
Surrogate: 2-Fluorobiphenyl		71.7 %	32.4-102		"	"	"	"
Surrogate: 2,4,6-Tribromophenol		88.9 %	18.1-105		"	"	"	"
Surrogate: Terphenyl-d14		86.1 %	29.1-130		"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-10
T181179-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	150	10	"	"	"	"	"	"	
C29-C40 (MORO)	76	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		70.1 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	61	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.1	2.0	"	"	"	"	"	"	
Cobalt	4.6	2.0	"	"	"	"	"	"	
Copper	4.1	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	21	5.0	"	"	"	"	"	"	
Zinc	26	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Project Number: 137-002
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Reported:
04/13/18 15:53

SP2-10
T181179-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	72	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	45	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		142 %	35-140		"	"	"	"	S-GC
<i>Surrogate: Decachlorobiphenyl</i>		120 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-10
T181179-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-10
T181179-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	51.5 %	15-121			"	"	"	"
Surrogate: Phenol-d6	56.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	60.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	67.2 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	91.0 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	95.8 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-11
T181179-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	110	10	"	"	"	"	"	"	
C29-C40 (MORO)	68	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		82.1 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	58	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.7	2.0	"	"	"	"	"	"	
Cobalt	4.4	2.0	"	"	"	"	"	"	
Copper	6.5	1.0	"	"	"	"	"	"	
Lead	17	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	31	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Reported:
04/13/18 15:53

SP2-11
T181179-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	26	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		134 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		129 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-11

T181179-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-11

T181179-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	54.1 %	15-121	"	"	"	"	"	"
Surrogate: Phenol-d6	60.5 %	24-113	"	"	"	"	"	"
Surrogate: Nitrobenzene-d5	62.2 %	21.3-119	"	"	"	"	"	"
Surrogate: 2-Fluorobiphenyl	67.8 %	32.4-102	"	"	"	"	"	"
Surrogate: 2,4,6-Tribromophenol	96.6 %	18.1-105	"	"	"	"	"	"
Surrogate: Terphenyl-d14	93.2 %	29.1-130	"	"	"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-12
T181179-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8040617	04/06/18	04/06/18	EPA 8015B	
C13-C28 (DRO)	140	10	"	"	"	"	"	"	
C29-C40 (MORO)	150	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		77.4 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8040616	04/06/18	04/09/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	58	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.8	2.0	"	"	"	"	"	"	
Cobalt	4.5	2.0	"	"	"	"	"	"	
Copper	5.8	1.0	"	"	"	"	"	"	
Lead	11	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8040615	04/06/18	04/09/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-12
T181179-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8040620	04/06/18	04/10/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	61	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		137 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		127 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-12
T181179-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

SP2-12
T181179-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8040619	04/06/18	04/11/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	50.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	58.5 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	59.6 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	67.5 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	93.0 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	94.0 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040617 - EPA 3550B GC

Blank (8040617-BLK1)

Prepared & Analyzed: 04/06/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	81.1		"	100		81.1	65-135			

LCS (8040617-BS1)

Prepared & Analyzed: 04/06/18

C13-C28 (DRO)	450	10	mg/kg	500		90.2	75-125			
Surrogate: <i>p</i> -Terphenyl	77.4		"	100		77.4	65-135			

LCS Dup (8040617-BSD1)

Prepared & Analyzed: 04/06/18

C13-C28 (DRO)	490	10	mg/kg	500		98.3	75-125	8.56	20	
Surrogate: <i>p</i> -Terphenyl	82.8		"	100		82.8	65-135			

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Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040616 - EPA 3050B

Blank (8040616-BLK1)

Prepared: 04/06/18 Analyzed: 04/09/18

Antimony	ND	3.0	mg/kg
Silver	ND	2.0	"
Arsenic	ND	5.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	2.0	"
Chromium	ND	2.0	"
Cobalt	ND	2.0	"
Copper	ND	1.0	"
Lead	ND	3.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Selenium	ND	5.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	ND	1.0	"

LCS (8040616-BS1)

Prepared: 04/06/18 Analyzed: 04/09/18

Arsenic	95.3	5.0	mg/kg	100	95.3	75-125
Barium	95.6	1.0	"	100	95.6	75-125
Cadmium	95.8	2.0	"	100	95.8	75-125
Chromium	95.8	2.0	"	100	95.8	75-125
Lead	98.3	3.0	"	100	98.3	75-125

Matrix Spike (8040616-MS1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

Arsenic	99.8	5.0	mg/kg	99.0	4.03	96.7	75-125
Barium	142	1.0	"	99.0	45.4	97.3	75-125
Cadmium	97.8	2.0	"	99.0	ND	98.7	75-125
Chromium	102	2.0	"	99.0	4.73	98.5	75-125
Lead	99.4	3.0	"	99.0	1.84	98.5	75-125

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04/13/18 15:53

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040616 - EPA 3050B

Matrix Spike Dup (8040616-MSD1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

Arsenic	83.9	4.5	mg/kg	90.9	4.03	87.8	75-125	17.3	20	
Barium	130	0.91	"	90.9	45.4	93.2	75-125	8.51	20	
Cadmium	82.3	1.8	"	90.9	ND	90.6	75-125	17.1	20	
Chromium	89.3	1.8	"	90.9	4.73	93.0	75-125	13.5	20	
Lead	84.0	2.7	"	90.9	1.84	90.3	75-125	16.8	20	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040615 - EPA 7471A Soil

Blank (8040615-BLK1)

Prepared: 04/06/18 Analyzed: 04/09/18

Mercury	ND	0.10	mg/kg

LCS (8040615-BS1)

Prepared: 04/06/18 Analyzed: 04/09/18

Mercury	0.438	0.10	mg/kg	0.410	107	80-120

Matrix Spike (8040615-MS1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

Mercury	0.453	0.10	mg/kg	0.417	ND	109	75-125

Matrix Spike Dup (8040615-MSD1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

Mercury	0.414	0.10	mg/kg	0.391	ND	106	75-125	8.91	20

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040620 - EPA 3550 ECD/GCMS

Blank (8040620-BLK1)

Prepared: 04/06/18 Analyzed: 04/09/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	13.3		"	9.90		134	35-140			
Surrogate: Decachlorobiphenyl	15.9		"	9.90		160	35-140			S-GC

LCS (8040620-BS1)

Prepared: 04/06/18 Analyzed: 04/09/18

PCB-1016	84.5	10	ug/kg	99.0		85.3	40-130			
PCB-1260	68.5	10	"	99.0		69.2	40-130			
Surrogate: Tetrachloro-meta-xylene	13.6		"	9.90		138	35-140			
Surrogate: Decachlorobiphenyl	16.0		"	9.90		162	35-140			S-GC

Matrix Spike (8040620-MS1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

PCB-1016	68.6	10	ug/kg	98.0	ND	70.0	40-130			
PCB-1260	62.9	10	"	98.0	ND	64.2	40-130			
Surrogate: Tetrachloro-meta-xylene	14.8		"	9.80		151	35-140			QM-05
Surrogate: Decachlorobiphenyl	15.9		"	9.80		162	35-140			QM-05

Matrix Spike Dup (8040620-MSD1)

Source: T181174-01

Prepared: 04/06/18 Analyzed: 04/09/18

PCB-1016	74.2	10	ug/kg	99.0	ND	74.9	40-130	7.75	30	
PCB-1260	68.0	10	"	99.0	ND	68.7	40-130	7.76	30	
Surrogate: Tetrachloro-meta-xylene	15.3		"	9.90		155	35-140			QM-05
Surrogate: Decachlorobiphenyl	16.4		"	9.90		165	35-140			QM-05

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040619 - EPA 3550 ECD/GCMS

Blank (8040619-BLK1)

Prepared: 04/06/18 Analyzed: 04/11/18

Carbazole	ND	300	ug/kg
Aniline	ND	300	"
Phenol	ND	1000	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
1-Methylnaphthalene	ND	300	"
2-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040619 - EPA 3550 ECD/GCMS

Blank (8040619-BLK1)

Prepared: 04/06/18 Analyzed: 04/11/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
Pyridine	ND	300	"
2,4,6-Trichlorophenol	ND	1000	"

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Reported:
04/13/18 15:53

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040619 - EPA 3550 ECD/GCMS

Blank (8040619-BLK1)

Prepared: 04/06/18 Analyzed: 04/11/18

Surrogate: 2-Fluorophenol	1920		ug/kg	3380		56.7	15-121			
Surrogate: Phenol-d6	1890		"	3380		55.9	24-113			
Surrogate: Nitrobenzene-d5	2290		"	3380		67.7	21.3-119			
Surrogate: 2-Fluorobiphenyl	2380		"	3380		70.3	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2600		"	3380		77.0	18.1-105			
Surrogate: Terphenyl-d14	2820		"	3380		83.6	29.1-130			

LCS (8040619-BS1)

Prepared: 04/06/18 Analyzed: 04/11/18

Phenol	1910	1000	ug/kg	3270		58.5	34-114			
2-Chlorophenol	2220	1000	"	3270		68.0	34-114			
1,4-Dichlorobenzene	2250	300	"	3270		68.9	34-114			
N-Nitrosodi-n-propylamine	2620	300	"	3270		80.1	30-110			
1,2,4-Trichlorobenzene	2420	300	"	3270		73.9	39-119			
4-Chloro-3-methylphenol	2270	1000	"	3270		69.4	50-130			
Acenaphthene	2480	300	"	3270		76.0	34-114			
Pentachlorophenol	3980	1000	"	3270		122	50-130			
Pyrene	2150	300	"	3270		65.8	30-110			
Surrogate: 2-Fluorophenol	1900		"	3270		58.1	15-121			
Surrogate: Phenol-d6	1970		"	3270		60.4	24-113			
Surrogate: Nitrobenzene-d5	2250		"	3270		68.8	21.3-119			
Surrogate: 2-Fluorobiphenyl	2190		"	3270		66.9	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2840		"	3270		87.0	18.1-105			
Surrogate: Terphenyl-d14	2800		"	3270		85.7	29.1-130			

LCS Dup (8040619-BSD1)

Prepared: 04/06/18 Analyzed: 04/11/18

Phenol	1820	1000	ug/kg	3320		54.8	34-114	4.88	42	
2-Chlorophenol	2140	1000	"	3320		64.3	34-114	3.95	40	
1,4-Dichlorobenzene	2180	300	"	3320		65.7	34-114	3.11	28	
N-Nitrosodi-n-propylamine	2660	300	"	3320		80.1	30-110	1.61	38	
1,2,4-Trichlorobenzene	2380	300	"	3320		71.5	39-119	1.72	28	
4-Chloro-3-methylphenol	2310	1000	"	3320		69.5	50-130	1.76	42	
Acenaphthene	2500	300	"	3320		75.3	34-114	0.788	31	
Pentachlorophenol	4080	1000	"	3320		123	50-130	2.43	50	
Pyrene	2200	300	"	3320		66.2	30-110	2.31	31	
Surrogate: 2-Fluorophenol	1800		"	3320		54.0	15-121			
Surrogate: Phenol-d6	1870		"	3320		56.4	24-113			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8040619 - EPA 3550 ECD/GCMS

LCS Dup (8040619-BSD1)

Prepared: 04/06/18 Analyzed: 04/11/18

Surrogate: Nitrobenzene-d5	2210		ug/kg	3320		66.4	21.3-119			
Surrogate: 2-Fluorobiphenyl	2180		"	3320		65.6	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2870		"	3320		86.3	18.1-105			
Surrogate: Terphenyl-d14	2910		"	3320		87.7	29.1-130			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/13/18 15:53

Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

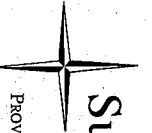
dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commercentre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, ROSEVILLE, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/5/18 Page: 1 Of 1
Project Name: LATHROP 48-A PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 7181178 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 - SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 - PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	8010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP2-1	4/5/18	8:54	Solid	JAR	X			X				X	X	X	01		1
SP2-2		8:57			X			X				X	X	X	02		1
SP2-3		8:59			X			X				X	X	X	03		1
SP2-4		9:02			X			X				X	X	X	04		1
SP2-5		9:04			X			X				X	X	X	05		1
SP2-6		9:08			X			X				X	X	X	06		1
SP2-7		9:14			X			X				X	X	X	07		1
SP2-8		9:18			X			X				X	X	X	08		1
SP2-9		9:21			X			X				X	X	X	09		1
SP2-10		9:24			X			X				X	X	X	10		1
SP2-11		9:26			X			X				X	X	X	11		1
SP2-12		9:29			X			X				X	X	X	12		1
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Total # of containers												Notes
<u>[Signature]</u>	4/5/18-13:10		<u>[Signature]</u>	4/5/18 13:10													* 3 DAY TAT ON METALS
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													
<u>[Signature]</u>	4/5/18 10:00		<u>[Signature]</u>	4/5/18 10:00													
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													
<u>[Signature]</u>	4/5/18 10:00		<u>[Signature]</u>	4/5/18 10:00													

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

Turn around time: NORMAL
EXCEPT 3 DAY
TAT ON METALS

COC 160913



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

30 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/06/18 10:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP2-6	T181179-06	Soil	04/05/18 09:08	04/06/18 10:00
SP2-9	T181179-09	Soil	04/05/18 09:21	04/06/18 10:00

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

DETECTIONS SUMMARY

Sample ID: SP2-6

Laboratory ID: T181179-06

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Arsenic	28	1.0	ug/l	6020 ICP-MS	STLC
Barium	520	1.0	ug/l	6020 ICP-MS	STLC
Beryllium	1.9	1.0	ug/l	6020 ICP-MS	STLC
Chromium	54	1.0	ug/l	6020 ICP-MS	STLC
Cobalt	44	1.0	ug/l	6020 ICP-MS	STLC
Copper	99	1.0	ug/l	6020 ICP-MS	STLC
Lead	110	2.0	ug/l	6020 ICP-MS	STLC
Mercury	0.21	0.10	ug/l	6020 ICP-MS	STLC
Molybdenum	7.2	1.0	ug/l	6020 ICP-MS	STLC
Nickel	73	1.0	ug/l	6020 ICP-MS	STLC
Silver	3.0	1.0	ug/l	6020 ICP-MS	STLC
Vanadium	250	1.0	ug/l	6020 ICP-MS	STLC
Zinc	320	1.0	ug/l	6020 ICP-MS	STLC

Sample ID: SP2-9

Laboratory ID: T181179-09

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Antimony	0.97	0.25	ug/l	6020 ICP-MS	STLC
Arsenic	8.9	0.25	ug/l	6020 ICP-MS	STLC
Barium	120	0.25	ug/l	6020 ICP-MS	STLC
Beryllium	0.36	0.25	ug/l	6020 ICP-MS	STLC
Cadmium	0.25	0.25	ug/l	6020 ICP-MS	STLC
Chromium	16	0.25	ug/l	6020 ICP-MS	STLC
Cobalt	13	0.25	ug/l	6020 ICP-MS	STLC
Copper	34	0.25	ug/l	6020 ICP-MS	STLC
Lead	51	0.25	ug/l	6020 ICP-MS	STLC
Mercury	0.091	0.025	ug/l	6020 ICP-MS	STLC
Molybdenum	5.7	0.25	ug/l	6020 ICP-MS	STLC
Nickel	15	0.25	ug/l	6020 ICP-MS	STLC
Selenium	1.6	1.2	ug/l	6020 ICP-MS	STLC
Silver	0.73	0.25	ug/l	6020 ICP-MS	STLC
Vanadium	80	0.25	ug/l	6020 ICP-MS	STLC

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

Sample ID: SP2-9

Laboratory ID: T181179-09

Analyte	Result	Reporting		Units	Method	Notes
			Limit			
Zinc	68		0.25	ug/l	6020 ICP-MS	STLC

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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

SP2-6
T181179-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8042318	04/23/18	04/26/18	6020 ICP-MS	STLC
Arsenic	28	1.0	"	"	"	"	"	"	STLC
Barium	520	1.0	"	"	"	"	"	"	STLC
Beryllium	1.9	1.0	"	"	"	"	04/27/18	"	STLC
Cadmium	ND	1.0	"	"	"	"	04/26/18	"	STLC
Chromium	54	1.0	"	"	"	"	"	"	STLC
Cobalt	44	1.0	"	"	"	"	"	"	STLC
Copper	99	1.0	"	"	"	"	"	"	STLC
Lead	110	2.0	"	"	"	"	"	"	STLC
Mercury	0.21	0.10	"	"	"	"	"	"	STLC
Molybdenum	7.2	1.0	"	"	"	"	"	"	STLC
Nickel	73	1.0	"	"	"	"	"	"	STLC
Selenium	ND	5.0	"	"	"	"	"	"	STLC
Silver	3.0	1.0	"	"	"	"	"	"	STLC
Thallium	ND	1.0	"	"	"	"	"	"	STLC
Vanadium	250	1.0	"	"	"	"	"	"	STLC
Zinc	320	1.0	"	"	"	"	"	"	STLC

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

SP2-9

T181179-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	0.97	0.25	ug/l	1	8042319	04/23/18	04/30/18	6020 ICP-MS	STLC
Arsenic	8.9	0.25	"	"	"	"	"	"	STLC
Barium	120	0.25	"	"	"	"	"	"	STLC
Beryllium	0.36	0.25	"	"	"	"	"	"	STLC
Cadmium	0.25	0.25	"	"	"	"	"	"	STLC
Chromium	16	0.25	"	"	"	"	"	"	STLC
Cobalt	13	0.25	"	"	"	"	"	"	STLC
Copper	34	0.25	"	"	"	"	"	"	STLC
Lead	51	0.25	"	"	"	"	"	"	STLC
Mercury	0.091	0.025	"	"	"	"	"	"	STLC
Molybdenum	5.7	0.25	"	"	"	"	"	"	STLC
Nickel	15	0.25	"	"	"	"	"	"	STLC
Selenium	1.6	1.2	"	"	"	"	"	"	STLC
Silver	0.73	0.25	"	"	"	"	"	"	STLC
Thallium	ND	0.25	"	"	"	"	"	"	STLC
Vanadium	80	0.25	"	"	"	"	"	"	STLC
Zinc	68	0.25	"	"	"	"	"	"	STLC

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Project Manager: Joe Brusca

Reported:
04/30/18 17:08

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Blank (8042318-BLK1)

Prepared: 04/23/18 Analyzed: 04/26/18

Antimony	ND	1.0	ug/l
Arsenic	ND	1.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	1.0	"
Chromium	ND	1.0	"
Cobalt	ND	1.0	"
Copper	ND	1.0	"
Lead	ND	2.0	"
Mercury	ND	0.10	"
Molybdenum	ND	1.0	"
Nickel	ND	1.0	"
Selenium	ND	5.0	"
Silver	ND	1.0	"
Thallium	ND	1.0	"
Vanadium	ND	1.0	"
Zinc	ND	1.0	"

LCS (8042318-BS1)

Prepared: 04/23/18 Analyzed: 04/26/18

Arsenic	57.4	1.0	ug/l	50.0	115	80-120
Barium	56.4	1.0	"	50.0	113	80-120
Cadmium	56.5	1.0	"	50.0	113	80-120
Chromium	55.6	1.0	"	50.0	111	80-120
Lead	56.2	2.0	"	50.0	112	80-120

Duplicate (8042318-DUP1)

Source: T181179-06

Prepared: 04/23/18 Analyzed: 04/26/18

Barium	548	5.0	ug/l	521	4.95	200
Chromium	54.5	5.0	"	53.8	1.28	200

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Matrix Spike (8042318-MS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	72.5	1.0	ug/l	50.0	27.8	89.2	75-125			
Barium	611	1.0	"	50.0	521	178	75-125			QM-PS
Cadmium	46.0	1.0	"	50.0	0.966	90.1	75-125			
Chromium	121	1.0	"	50.0	53.8	134	75-125			QM-PS
Lead	168	2.0	"	50.0	112	112	75-125			

Matrix Spike Dup (8042318-MSD1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	73.4	1.0	ug/l	50.0	27.8	91.2	75-125	1.36	20	
Barium	595	1.0	"	50.0	521	147	75-125	2.61	20	QM-PS
Cadmium	47.4	1.0	"	50.0	0.966	92.9	75-125	2.97	20	
Chromium	124	1.0	"	50.0	53.8	141	75-125	2.85	20	QM-PS
Lead	164	2.0	"	50.0	112	104	75-125	2.22	20	

Post Spike (8042318-PS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Barium	287		ug/l	25.0	261	104	80-120			
Chromium	56.2		"	25.0	26.9	117	80-120			

Batch 8042319 - STLC Metals

Blank (8042319-BLK1)		Prepared: 04/23/18 Analyzed: 04/30/18								
Antimony	ND	0.25	ug/l							STLC
Arsenic	0.366	0.25	"							STLC
Barium	4.73	0.25	"							QB-01, STLC
Beryllium	ND	0.25	"							STLC
Cadmium	ND	0.25	"							STLC
Chromium	0.518	0.25	"							QB-01, STLC
Cobalt	ND	0.25	"							STLC
Copper	0.663	0.25	"							QB-01, STLC
Lead	0.269	0.25	"							STLC
Mercury	ND	0.025	"							STLC
Molybdenum	0.282	0.25	"							STLC
Nickel	ND	0.25	"							STLC
Selenium	ND	1.2	"							STLC
Silver	ND	0.25	"							STLC
Thallium	ND	0.25	"							STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042319 - STLC Metals

Blank (8042319-BLK1)

Prepared: 04/23/18 Analyzed: 04/30/18

Vanadium	1.01	0.25	ug/l							QB-01, STLC
Zinc	0.690	0.25	"							QB-01, STLC

LCS (8042319-BS1)

Prepared: 04/23/18 Analyzed: 04/30/18

Arsenic	45.9	0.25	ug/l	50.0		91.7	80-120			STLC
Barium	51.1	0.25	"	50.0		102	80-120			STLC
Cadmium	45.4	0.25	"	50.0		90.8	80-120			STLC
Chromium	45.9	0.25	"	50.0		91.8	80-120			STLC
Lead	44.8	0.25	"	50.0		89.5	80-120			STLC

Matrix Spike (8042319-MS1)

Source: T181179-09

Prepared: 04/23/18 Analyzed: 04/30/18

Arsenic	55.4	0.25	ug/l	50.0	8.92	93.0	75-125			STLC
Barium	174	0.25	"	50.0	115	117	75-125			STLC
Cadmium	46.3	0.25	"	50.0	0.251	92.0	75-125			STLC
Chromium	67.0	0.25	"	50.0	15.5	103	75-125			STLC
Lead	102	0.25	"	50.0	50.7	103	75-125			STLC

Matrix Spike Dup (8042319-MSD1)

Source: T181179-09

Prepared: 04/23/18 Analyzed: 04/30/18

Arsenic	70.7	0.28	ug/l	56.8	8.92	109	75-125	24.3	20	QR-04, STLC
Barium	211	0.28	"	56.8	115	169	75-125	19.4	20	QR-04, STLC
Cadmium	60.2	0.28	"	56.8	0.251	105	75-125	26.2	20	QR-04, STLC
Chromium	82.6	0.28	"	56.8	15.5	118	75-125	20.9	20	QR-04, STLC
Lead	123	0.28	"	56.8	50.7	127	75-125	17.9	20	QR-04, STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/30/18 17:08

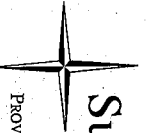
Notes and Definitions

STLC	Sample prepared in accordance with CCR Chapter 11, Article 5, Appendix II, Waste Extraction Test (WET).
QR-04	The percent recovery and/or RPD was outside acceptance criteria. Results accepted based upon percent recovery results in duplicate QC sample and the CCV and CCB results.
QM-PS	The percent recovery and/or RPD are outside acceptance criteria. Results accepted based upon percent recovery results in the post spike and/or serial dilution.
QB-01	The method blank contains analyte at a concentration above the MRL; however, concentration is less than 10% of the sample result, which is negligible according to method criteria.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commercentre Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO BOX 332, ROSEVILLE, CA 95661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: JOE BRUSCA

Date: 4/5/18 Page: 1 Of 1
Project Name: LATHROP 48-A PROPERTY
Collector: BRUSCA Client Project #: 137-002
Batch #: 7181178 EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260 - SVOCs 8270	8260 + OXY	8260 BTEX, OXY only	8270 - PCBs 8082	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	8010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP2-1	4/5/18	8:54	Solid	JAR	X			X				X	X	X	01		1
SP2-2		8:57			X			X				X	X	X	02		1
SP2-3		8:59			X			X				X	X	X	03		1
SP2-4		9:02			X			X				X	X	X	04		1
SP2-5		9:04			X			X				X	X	X	05		1
SP2-6		9:08			X			X				X	X	X	06		1
SP2-7		9:14			X			X				X	X	X	07		1
SP2-8		9:18			X			X				X	X	X	08		1
SP2-9		9:21			X			X				X	X	X	09		1
SP2-10		9:24			X			X				X	X	X	10		1
SP2-11		9:26			X			X				X	X	X	11		1
SP2-12		9:29			X			X				X	X	X	12		1
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time	Total # of containers												Notes
<u>[Signature]</u>	4/5/18-13:10		<u>[Signature]</u>	4/5/18 13:10													* 3 DAY TAT ON METALS
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													
<u>[Signature]</u>	4/5/18 10:00		<u>[Signature]</u>	4/5/18 10:00													
Relinquished by: (signature)	Date / Time		Received by: (signature)	Date / Time													

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

Turn around time: NORMAL

Chain of Custody seals Y/N/NA

Seals intact? Y/N/NA

Received good condition/cold Y/N

EXCEPT 3 DAY TAT ON METALS

COC 160913



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27 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/13/18 09:30. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP3-3	T181249-03	Soil	04/12/18 11:00	04/13/18 09:30
SP3-12	T181249-12	Soil	04/12/18 11:26	04/13/18 09:30
SP4-1	T181249-13	Soil	04/12/18 11:35	04/13/18 09:30

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

DETECTIONS SUMMARY

Sample ID: SP3-3

Laboratory ID: T181249-03

Reporting

Analyte	Result	Limit	Units	Method	Notes
Arsenic	13	1.0	ug/l	6020 ICP-MS	STLC
Barium	340	1.0	ug/l	6020 ICP-MS	STLC
Cadmium	1.5	1.0	ug/l	6020 ICP-MS	STLC
Chromium	40	1.0	ug/l	6020 ICP-MS	STLC
Cobalt	33	1.0	ug/l	6020 ICP-MS	STLC
Copper	120	1.0	ug/l	6020 ICP-MS	STLC
Lead	170	2.0	ug/l	6020 ICP-MS	STLC
Mercury	0.28	0.10	ug/l	6020 ICP-MS	STLC
Molybdenum	7.1	1.0	ug/l	6020 ICP-MS	STLC
Nickel	41	1.0	ug/l	6020 ICP-MS	STLC
Silver	3.1	1.0	ug/l	6020 ICP-MS	STLC
Vanadium	99	1.0	ug/l	6020 ICP-MS	STLC
Zinc	210	1.0	ug/l	6020 ICP-MS	STLC

Sample ID: SP3-12

Laboratory ID: T181249-12

Reporting

Analyte	Result	Limit	Units	Method	Notes
Arsenic	9.0	1.0	ug/l	6020 ICP-MS	STLC
Barium	360	1.0	ug/l	6020 ICP-MS	STLC
Beryllium	1.3	1.0	ug/l	6020 ICP-MS	STLC
Cadmium	1.6	1.0	ug/l	6020 ICP-MS	STLC
Chromium	40	1.0	ug/l	6020 ICP-MS	STLC
Cobalt	49	1.0	ug/l	6020 ICP-MS	STLC
Copper	110	1.0	ug/l	6020 ICP-MS	STLC
Lead	200	2.0	ug/l	6020 ICP-MS	STLC
Mercury	0.52	0.10	ug/l	6020 ICP-MS	STLC
Molybdenum	4.8	1.0	ug/l	6020 ICP-MS	STLC
Nickel	54	1.0	ug/l	6020 ICP-MS	STLC
Silver	7.0	1.0	ug/l	6020 ICP-MS	STLC
Vanadium	110	1.0	ug/l	6020 ICP-MS	STLC
Zinc	230	1.0	ug/l	6020 ICP-MS	STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

Sample ID: SP4-1

Laboratory ID: T181249-13

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Arsenic	11	1.0		ug/l	6020 ICP-MS	STLC
Barium	230	1.0		ug/l	6020 ICP-MS	STLC
Cadmium	1.9	1.0		ug/l	6020 ICP-MS	STLC
Chromium	33	1.0		ug/l	6020 ICP-MS	STLC
Cobalt	31	1.0		ug/l	6020 ICP-MS	STLC
Copper	110	1.0		ug/l	6020 ICP-MS	STLC
Lead	170	2.0		ug/l	6020 ICP-MS	STLC
Mercury	0.32	0.10		ug/l	6020 ICP-MS	STLC
Molybdenum	7.4	1.0		ug/l	6020 ICP-MS	STLC
Nickel	28	1.0		ug/l	6020 ICP-MS	STLC
Silver	3.0	1.0		ug/l	6020 ICP-MS	STLC
Vanadium	75	1.0		ug/l	6020 ICP-MS	STLC
Zinc	250	1.0		ug/l	6020 ICP-MS	STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager

Page 3 of 9



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

SP3-3
T181249-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8042318	04/23/18	04/26/18	6020 ICP-MS	STLC
Arsenic	13	1.0	"	"	"	"	"	"	STLC
Barium	340	1.0	"	"	"	"	"	"	STLC
Beryllium	ND	1.0	"	"	"	"	04/27/18	"	STLC
Cadmium	1.5	1.0	"	"	"	"	04/26/18	"	STLC
Chromium	40	1.0	"	"	"	"	"	"	STLC
Cobalt	33	1.0	"	"	"	"	"	"	STLC
Copper	120	1.0	"	"	"	"	"	"	STLC
Lead	170	2.0	"	"	"	"	"	"	STLC
Mercury	0.28	0.10	"	"	"	"	"	"	STLC
Molybdenum	7.1	1.0	"	"	"	"	"	"	STLC
Nickel	41	1.0	"	"	"	"	"	"	STLC
Selenium	ND	5.0	"	"	"	"	"	"	STLC
Silver	3.1	1.0	"	"	"	"	"	"	STLC
Thallium	ND	1.0	"	"	"	"	"	"	STLC
Vanadium	99	1.0	"	"	"	"	"	"	STLC
Zinc	210	1.0	"	"	"	"	"	"	STLC

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

SP3-12
T181249-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8042318	04/23/18	04/26/18	6020 ICP-MS	STLC
Arsenic	9.0	1.0	"	"	"	"	"	"	STLC
Barium	360	1.0	"	"	"	"	"	"	STLC
Beryllium	1.3	1.0	"	"	"	"	04/27/18	"	STLC
Cadmium	1.6	1.0	"	"	"	"	04/26/18	"	STLC
Chromium	40	1.0	"	"	"	"	"	"	STLC
Cobalt	49	1.0	"	"	"	"	"	"	STLC
Copper	110	1.0	"	"	"	"	"	"	STLC
Lead	200	2.0	"	"	"	"	"	"	STLC
Mercury	0.52	0.10	"	"	"	"	"	"	STLC
Molybdenum	4.8	1.0	"	"	"	"	"	"	STLC
Nickel	54	1.0	"	"	"	"	"	"	STLC
Selenium	ND	5.0	"	"	"	"	"	"	STLC
Silver	7.0	1.0	"	"	"	"	"	"	STLC
Thallium	ND	1.0	"	"	"	"	"	"	STLC
Vanadium	110	1.0	"	"	"	"	"	"	STLC
Zinc	230	1.0	"	"	"	"	"	"	STLC

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

SP4-1

T181249-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8042318	04/23/18	04/26/18	6020 ICP-MS	STLC
Arsenic	11	1.0	"	"	"	"	"	"	STLC
Barium	230	1.0	"	"	"	"	"	"	STLC
Beryllium	ND	1.0	"	"	"	"	04/27/18	"	STLC
Cadmium	1.9	1.0	"	"	"	"	04/26/18	"	STLC
Chromium	33	1.0	"	"	"	"	"	"	STLC
Cobalt	31	1.0	"	"	"	"	"	"	STLC
Copper	110	1.0	"	"	"	"	"	"	STLC
Lead	170	2.0	"	"	"	"	"	"	STLC
Mercury	0.32	0.10	"	"	"	"	"	"	STLC
Molybdenum	7.4	1.0	"	"	"	"	"	"	STLC
Nickel	28	1.0	"	"	"	"	"	"	STLC
Selenium	ND	5.0	"	"	"	"	"	"	STLC
Silver	3.0	1.0	"	"	"	"	"	"	STLC
Thallium	ND	1.0	"	"	"	"	"	"	STLC
Vanadium	75	1.0	"	"	"	"	"	"	STLC
Zinc	250	1.0	"	"	"	"	"	"	STLC

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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Blank (8042318-BLK1)

Prepared: 04/23/18 Analyzed: 04/26/18

Antimony	ND	1.0	ug/l
Arsenic	ND	1.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	1.0	"
Chromium	ND	1.0	"
Cobalt	ND	1.0	"
Copper	ND	1.0	"
Lead	ND	2.0	"
Mercury	ND	0.10	"
Molybdenum	ND	1.0	"
Nickel	ND	1.0	"
Selenium	ND	5.0	"
Silver	ND	1.0	"
Thallium	ND	1.0	"
Vanadium	ND	1.0	"
Zinc	ND	1.0	"

LCS (8042318-BS1)

Prepared: 04/23/18 Analyzed: 04/26/18

Arsenic	57.4	1.0	ug/l	50.0	115	80-120
Barium	56.4	1.0	"	50.0	113	80-120
Cadmium	56.5	1.0	"	50.0	113	80-120
Chromium	55.6	1.0	"	50.0	111	80-120
Lead	56.2	2.0	"	50.0	112	80-120

Duplicate (8042318-DUP1)

Source: T181179-06

Prepared: 04/23/18 Analyzed: 04/26/18

Barium	548	5.0	ug/l	521	4.95	200
Chromium	54.5	5.0	"	53.8	1.28	200

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Matrix Spike (8042318-MS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	72.5	1.0	ug/l	50.0	27.8	89.2	75-125			
Barium	611	1.0	"	50.0	521	178	75-125			QM-PS
Cadmium	46.0	1.0	"	50.0	0.966	90.1	75-125			
Chromium	121	1.0	"	50.0	53.8	134	75-125			QM-PS
Lead	168	2.0	"	50.0	112	112	75-125			

Matrix Spike Dup (8042318-MSD1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	73.4	1.0	ug/l	50.0	27.8	91.2	75-125	1.36	20	
Barium	595	1.0	"	50.0	521	147	75-125	2.61	20	QM-PS
Cadmium	47.4	1.0	"	50.0	0.966	92.9	75-125	2.97	20	
Chromium	124	1.0	"	50.0	53.8	141	75-125	2.85	20	QM-PS
Lead	164	2.0	"	50.0	112	104	75-125	2.22	20	

Post Spike (8042318-PS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Barium	287		ug/l	25.0	261	104	80-120			
Chromium	56.2		"	25.0	26.9	117	80-120			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:10

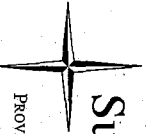
Notes and Definitions

STLC	Sample prepared in accordance with CCR Chapter 11, Article 5, Appendix II, Waste Extraction Test (WET).
QM-PS	The percent recovery and/or RPD are outside acceptance criteria. Results accepted based upon percent recovery results in the post spike and/or serial dilution.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, Rosemead, CA 91661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: Joe Brusca

Date: 4/12/18 Page: 1 of 1
Project Name: LATHROP 43-Ac Property
Collector: BRUSCA Client Project #: 137-002
Batch #: 7/8/247 EDF #:

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260	8260 + OXY	8260 BTEX, OXY only	8270	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP3-1	4/12/18	10:51	SOL	JAR	X	X	X	X	X	X	X	X	X	X	01		1
SP3-2		10:55			X	X	X	X	X	X	X	X	X	X	02		1
SP3-3		11:00			X	X	X	X	X	X	X	X	X	X	03		1
SP3-4		11:02			X	X	X	X	X	X	X	X	X	X	04		1
SP3-5		11:05			X	X	X	X	X	X	X	X	X	X	05		1
SP3-6		11:09			X	X	X	X	X	X	X	X	X	X	06		1
SP3-7		11:14			X	X	X	X	X	X	X	X	X	X	07		1
SP3-8		11:16			X	X	X	X	X	X	X	X	X	X	08		1
SP3-9		11:19			X	X	X	X	X	X	X	X	X	X	09		1
SP3-10		11:21			X	X	X	X	X	X	X	X	X	X	10		1
SP3-11		11:23			X	X	X	X	X	X	X	X	X	X	11		1
SP3-12		11:26			X	X	X	X	X	X	X	X	X	X	12		1
SP4-1		11:35			X	X	X	X	X	X	X	X	X	X	13		1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18-15:45</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 15:45</u>																	
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:50</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u>																	
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u>																	

Turn around time: NORMAL

Chain of Custody seals Y/N/A
Seals intact? Y/N/A
Received good condition/cold 36

Notes

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

COC 160914



May 1, 2018

FAL Project ID: 11455

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11455**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Thirteen solid samples were received on 4/12/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of fifteen business days for project **11455**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and the sample photos. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11455**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink, appearing to read "Bradley B. Silverbush".

Bradley B. Silverbush
Director of Operations

FRONTIER ANALYTICAL LABORATORY

5172 Hillsdale Circle * El Dorado Hills, CA 95762
Tel (916) 934-0900 * Fax (916) 934-0999
www.frontieranalytical.com

000001 of 000022

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: **11455**

Received on: **04/12/2018**

Project Due: **05/04/2018**

Storage: **R-4**

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11455-001-SA	0	137-002	SP3-1	EPA 1613 D/F	Solid	04/12/2018	10:51 am	04/12/2019
11455-002-SA	0	137-002	SP3-2	EPA 1613 D/F	Solid	04/12/2018	10:55 am	04/12/2019
11455-003-SA	0	137-002	SP3-3	EPA 1613 D/F	Solid	04/12/2018	11:00 am	04/12/2019
11455-004-SA	0	137-002	SP3-4	EPA 1613 D/F	Solid	04/12/2018	11:02 am	04/12/2019
11455-005-SA	0	137-002	SP3-5	EPA 1613 D/F	Solid	04/12/2018	11:05 am	04/12/2019
11455-006-SA	0	137-002	SP3-6	EPA 1613 D/F	Solid	04/12/2018	11:09 am	04/12/2019
11455-007-SA	0	137-002	SP3-7	EPA 1613 D/F	Solid	04/12/2018	11:14 am	04/12/2019
11455-008-SA	0	137-002	SP3-8	EPA 1613 D/F	Solid	04/12/2018	11:16 am	04/12/2019
11455-009-SA	0	137-002	SP3-9	EPA 1613 D/F	Solid	04/12/2018	11:19 am	04/12/2019
11455-010-SA	0	137-002	SP3-10	EPA 1613 D/F	Solid	04/12/2018	11:21 am	04/12/2019
11455-011-SA	0	137-002	SP3-11	EPA 1613 D/F	Solid	04/12/2018	11:23 am	04/12/2019
11455-012-SA	0	137-002	SP3-12	EPA 1613 D/F	Solid	04/12/2018	11:26 am	04/12/2019
11455-013-SA	0	137-002	SP4-1	EPA 1613 D/F	Solid	04/12/2018	11:35 am	04/12/2019

EPA Method 1613
PCDD/F



FAL ID: 11455-001-MB
Client ID: Method Blank
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.137		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.215		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.222		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.236		-	0.0940	Total TCDD	ND	0.137	
1,2,3,7,8,9-HxCDD	ND	0.210		-	0.0823	Total PeCDD	ND	0.215	
1,2,3,4,6,7,8-HpCDD	ND	0.275		-	0.0842	Total HxCDD	ND	0.236	
OCDD	ND	0.430		-	0.172	Total HpCDD	ND	0.275	
2,3,7,8-TCDF	ND	0.107		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.148		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.159		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.155		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.151		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.161		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.183		-	0.0657	Total TCDF	ND	0.107	
1,2,3,4,6,7,8-HpCDF	ND	0.180		-	0.0747	Total PeCDF	ND	0.159	
1,2,3,4,7,8,9-HpCDF	ND	0.195		-	0.0883	Total HxCDF	ND	0.183	
OCDF	ND	0.422		-	0.170	Total HpCDF	ND	0.195	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	96.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	94.6	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.4	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	83.4	23.0 - 140	
13C-OCDD	62.6	17.0 - 157	
13C-2,3,7,8-TCDF	98.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	90.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.6	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	85.6	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	83.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	82.7	26.0 - 138	
13C-OCDF	66.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 87.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000003 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-001-OPR
Client ID: OPR
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 04-30-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.8	6.70 - 15.8	
1,2,3,7,8-PeCDD	53.0	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	47.7	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	50.6	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	47.5	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	51.0	35.0 - 70.0	
OCDD	101	78.0 - 144	
2,3,7,8-TCDF	10.7	7.50 - 15.8	
1,2,3,7,8-PeCDF	50.3	40.0 - 67.0	
2,3,4,7,8-PeCDF	51.6	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	51.1	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	50.1	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	51.1	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	51.1	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	51.1	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	52.3	39.0 - 69.0	
OCDF	104	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	92.4	20.0 - 175	
13C-1,2,3,7,8-PeCDD	96.5	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	97.1	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	100	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	88.8	26.0 - 166	
13C-OCDD	81.3	13.0 - 198	
13C-2,3,7,8-TCDF	97.6	22.0 - 152	
13C-1,2,3,7,8-PeCDF	91.3	21.0 - 192	
13C-2,3,4,7,8-PeCDF	92.6	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	89.0	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	90.0	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	88.6	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	87.6	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	90.3	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	91.3	20.0 - 186	
13C-OCDF	80.4	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	89.4	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000004 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-001-SA
Client ID: SP3-1
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.01 g
% Solids: 92.26

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 1.79
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.206		-	0.0273				
1,2,3,7,8-PeCDD	0.545	-	J	0.545	0.0570				
1,2,3,4,7,8-HxCDD	0.893	-	J	0.0893	0.0793				
1,2,3,6,7,8-HxCDD	1.80	-	J	0.180	0.0940	Total TCDD	ND	0.206	
1,2,3,7,8,9-HxCDD	1.33	-	J	0.133	0.0823	Total PeCDD	4.20	-	J
1,2,3,4,6,7,8-HpCDD	27.0	-		0.270	0.0842	Total HxCDD	17.8	-	
OCDD	194	-		0.0582	0.172	Total HpCDD	55.7	-	
2,3,7,8-TCDF	0.767	-	J	0.0767	0.0269				
1,2,3,7,8-PeCDF	ND	0.224		-	0.0449				
2,3,4,7,8-PeCDF	0.672	-	J	0.202	0.0468				
1,2,3,4,7,8-HxCDF	0.658	-	J	0.0658	0.0437				
1,2,3,6,7,8-HxCDF	0.555	-	J	0.0555	0.0417				
2,3,4,6,7,8-HxCDF	0.473	-	J	0.0473	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.202		-	0.0657	Total TCDF	5.58	-	
1,2,3,4,6,7,8-HpCDF	5.84	-		0.0584	0.0747	Total PeCDF	3.97	-	J
1,2,3,4,7,8,9-HpCDF	0.735	-	J	0.00735	0.0883	Total HxCDF	8.08	-	
OCDF	15.8	-		0.00474	0.170	Total HpCDF	16.7	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	90.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	95.1	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	93.1	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	87.0	23.0 - 140	
13C-OCDD	76.8	17.0 - 157	
13C-2,3,7,8-TCDF	92.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	88.2	24.0 - 185	
13C-2,3,4,7,8-PeCDF	88.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	87.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	86.7	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	84.5	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	82.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	84.9	26.0 - 138	
13C-OCDF	76.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 88.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000005 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-002-SA
Client ID: SP3-2
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.01 g
% Solids: 94.99

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 1.62
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.202		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.456		-	0.0570				
1,2,3,4,7,8-HxCDD	0.707	-	J	0.0707	0.0793				
1,2,3,6,7,8-HxCDD	1.56	-	J	0.156	0.0940	Total TCDD	ND	0.202	
1,2,3,7,8,9-HxCDD	0.878	-	J	0.0878	0.0823	Total PeCDD	ND	0.456	
1,2,3,4,6,7,8-HpCDD	24.6	-		0.246	0.0842	Total HxCDD	11.8	-	
OCDD	173	-		0.0519	0.172	Total HpCDD	47.0	-	
2,3,7,8-TCDF	1.80	-		0.180	0.0269				
1,2,3,7,8-PeCDF	ND	0.264		-	0.0449				
2,3,4,7,8-PeCDF	1.33	-	J	0.399	0.0468				
1,2,3,4,7,8-HxCDF	1.32	-	J	0.132	0.0437				
1,2,3,6,7,8-HxCDF	0.729	-	J	0.0729	0.0417				
2,3,4,6,7,8-HxCDF	0.855	-	J	0.0855	0.0574				
1,2,3,7,8,9-HxCDF	0.513	-	J	0.0513	0.0657	Total TCDF	10.2	-	
1,2,3,4,6,7,8-HpCDF	7.87	-		0.0787	0.0747	Total PeCDF	8.81	-	
1,2,3,4,7,8,9-HpCDF	0.678	-	J	0.00678	0.0883	Total HxCDF	11.6	-	
OCDF	15.2	-		0.00456	0.170	Total HpCDF	18.8	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	82.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	84.0	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	88.2	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	85.4	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	80.3	23.0 - 140	
13C-OCDD	71.8	17.0 - 157	
13C-2,3,7,8-TCDF	85.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	78.1	24.0 - 185	
13C-2,3,4,7,8-PeCDF	79.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	80.8	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	80.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	78.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	77.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	79.1	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	79.5	26.0 - 138	
13C-OCDF	71.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 79.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000006 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-003-SA
Client ID: SP3-3
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.02 g
% Solids: 92.65

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 0.933
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.149		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.331		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.230		-	0.0793				
1,2,3,6,7,8-HxCDD	1.90	-	J	0.190	0.0940	Total TCDD	0.697	-	J
1,2,3,7,8,9-HxCDD	0.732	-	J	0.0732	0.0823	Total PeCDD	0.448	-	J
1,2,3,4,6,7,8-HpCDD	34.3	-		0.343	0.0842	Total HxCDD	12.3	-	
OCDD	302	-		0.0906	0.172	Total HpCDD	60.9	-	
2,3,7,8-TCDF	0.431	-	J	0.0431	0.0269				
1,2,3,7,8-PeCDF	ND	0.197		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.204		-	0.0468				
1,2,3,4,7,8-HxCDF	0.531	-	J	0.0531	0.0437				
1,2,3,6,7,8-HxCDF	0.440	-	J	0.0440	0.0417				
2,3,4,6,7,8-HxCDF	0.389	-	J	0.0389	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.150		-	0.0657	Total TCDF	2.97	-	
1,2,3,4,6,7,8-HpCDF	4.89	-	J	0.0489	0.0747	Total PeCDF	2.01	-	J
1,2,3,4,7,8,9-HpCDF	0.534	-	J	0.00534	0.0883	Total HxCDF	7.73	-	
OCDF	10.5	-		0.00315	0.170	Total HpCDF	13.0	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	95.1	25.0 - 164	
13C-1,2,3,7,8-PeCDD	95.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	98.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	94.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	88.6	23.0 - 140	
13C-OCDD	79.4	17.0 - 157	
13C-2,3,7,8-TCDF	101	24.0 - 169	
13C-1,2,3,7,8-PeCDF	93.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	95.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.6	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.0	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	85.9	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	84.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.5	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	88.5	26.0 - 138	
13C-OCDF	79.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 92.8 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000007 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-004-SA
Client ID: SP3-4
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.02 g
% Solids: 93.88

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 0.522
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.189		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.314		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.222		-	0.0793				
1,2,3,6,7,8-HxCDD	0.820	-	J	0.0820	0.0940	Total TCDD	ND	0.189	
1,2,3,7,8,9-HxCDD	0.707	-	J	0.0707	0.0823	Total PeCDD	ND	0.314	
1,2,3,4,6,7,8-HpCDD	13.9	-		0.139	0.0842	Total HxCDD	6.13	-	
OCDD	104	-		0.0312	0.172	Total HpCDD	26.4	-	
2,3,7,8-TCDF	0.634	-	J	0.0634	0.0269				
1,2,3,7,8-PeCDF	ND	0.230		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.239		-	0.0468				
1,2,3,4,7,8-HxCDF	0.397	-	J	0.0397	0.0437				
1,2,3,6,7,8-HxCDF	0.301	-	J	0.0301	0.0417				
2,3,4,6,7,8-HxCDF	0.308	-	J	0.0308	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.128		-	0.0657	Total TCDF	3.03	-	
1,2,3,4,6,7,8-HpCDF	2.93	-	J	0.0293	0.0747	Total PeCDF	1.77	-	J
1,2,3,4,7,8,9-HpCDF	0.326	-	J	0.00326	0.0883	Total HxCDF	4.20	-	J
OCDF	8.17	-	J	0.00245	0.170	Total HpCDF	8.09	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	95.0	25.0 - 164	
13C-1,2,3,7,8-PeCDD	97.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	99.5	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.3	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	91.3	23.0 - 140	
13C-OCDD	82.7	17.0 - 157	
13C-2,3,7,8-TCDF	95.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.4	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	88.7	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	87.4	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	86.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	88.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.7	26.0 - 138	
13C-OCDF	81.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 88.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000008 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-005-SA
Client ID: SP3-5
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.04 g
% Solids: 90.61

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 2.76
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.204		-	0.0273				
1,2,3,7,8-PeCDD	0.870	-	J	0.870	0.0570				
1,2,3,4,7,8-HxCDD	0.714	-	J	0.0714	0.0793				
1,2,3,6,7,8-HxCDD	2.01	-	J	0.201	0.0940	Total TCDD	2.30	-	
1,2,3,7,8,9-HxCDD	1.56	-	J	0.156	0.0823	Total PeCDD	7.61	-	
1,2,3,4,6,7,8-HpCDD	29.9	-		0.299	0.0842	Total HxCDD	22.3	-	
OCDD	195	-		0.0585	0.172	Total HpCDD	65.0	-	
2,3,7,8-TCDF	2.41	-	F	0.241	0.0269				
1,2,3,7,8-PeCDF	0.523	-	J	0.0157	0.0449				
2,3,4,7,8-PeCDF	1.00	-	J	0.300	0.0468				
1,2,3,4,7,8-HxCDF	1.17	-	J	0.117	0.0437				
1,2,3,6,7,8-HxCDF	1.47	-	J	0.147	0.0417				
2,3,4,6,7,8-HxCDF	0.968	-	J	0.0968	0.0574				
1,2,3,7,8,9-HxCDF	0.495	-	J	0.0495	0.0657	Total TCDF	11.3	-	
1,2,3,4,6,7,8-HpCDF	11.1	-		0.111	0.0747	Total PeCDF	8.69	-	
1,2,3,4,7,8,9-HpCDF	1.63	-	J	0.0163	0.0883	Total HxCDF	12.1	-	
OCDF	16.6	-		0.00498	0.170	Total HpCDF	20.9	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	90.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	98.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	96.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.6	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	90.3	23.0 - 140	
13C-OCDD	80.2	17.0 - 157	
13C-2,3,7,8-TCDF	95.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.0	24.0 - 185	
13C-2,3,4,7,8-PeCDF	94.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.5	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.0	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	86.7	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	87.4	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	87.6	26.0 - 138	
13C-OCDF	78.4	17.0 - 157	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	88.5	35.0 - 197	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 5/1/2018

Reviewed By: 
Date: 5/1/2018

EPA Method 1613 PCDD/F



FAL ID: 11455-006-SA
Client ID: SP3-6
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.01 g
% Solids: 94.32

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 0.995
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.187		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.334		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.245		-	0.0793				
1,2,3,6,7,8-HxCDD	1.51	-	J	0.151	0.0940	Total TCDD	ND	0.187	
1,2,3,7,8,9-HxCDD	0.597	-	J	0.0597	0.0823	Total PeCDD	ND	0.334	
1,2,3,4,6,7,8-HpCDD	30.5	-		0.305	0.0842	Total HxCDD	11.1	-	
OCDD	265	-		0.0795	0.172	Total HpCDD	57.0	-	
2,3,7,8-TCDF	0.552	-	J	0.0552	0.0269				
1,2,3,7,8-PeCDF	ND	0.177		-	0.0449				
2,3,4,7,8-PeCDF	0.411	-	J	0.123	0.0468				
1,2,3,4,7,8-HxCDF	0.572	-	J	0.0572	0.0437				
1,2,3,6,7,8-HxCDF	0.567	-	J	0.0567	0.0417				
2,3,4,6,7,8-HxCDF	0.476	-	J	0.0476	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.133		-	0.0657	Total TCDF	3.25	-	
1,2,3,4,6,7,8-HpCDF	5.04	-		0.0504	0.0747	Total PeCDF	2.10	-	J
1,2,3,4,7,8,9-HpCDF	0.646	-	J	0.00646	0.0883	Total HxCDF	5.48	-	
OCDF	11.1	-		0.00333	0.170	Total HpCDF	12.0	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	92.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	93.9	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	99.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.0	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	90.2	23.0 - 140	
13C-OCDD	79.5	17.0 - 157	
13C-2,3,7,8-TCDF	96.3	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	90.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	87.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.0	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	87.2	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.2	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	90.7	26.0 - 138	
13C-OCDF	83.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 89.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000010 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-007-SA
Client ID: SP3-7
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.03 g
% Solids: 93.06

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 0.823
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.180		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.299		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.210		-	0.0793				
1,2,3,6,7,8-HxCDD	1.64	-	J	0.164	0.0940	Total TCDD	0.811	-	J
1,2,3,7,8,9-HxCDD	0.759	-	J	0.0759	0.0823	Total PeCDD	ND	0.299	
1,2,3,4,6,7,8-HpCDD	28.8	-		0.288	0.0842	Total HxCDD	11.1	-	
OCDD	232	-		0.0696	0.172	Total HpCDD	51.5	-	
2,3,7,8-TCDF	0.428	-	J	0.0428	0.0269				
1,2,3,7,8-PeCDF	ND	0.246		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.252		-	0.0468				
1,2,3,4,7,8-HxCDF	0.458	-	J	0.0458	0.0437				
1,2,3,6,7,8-HxCDF	0.541	-	J	0.0541	0.0417				
2,3,4,6,7,8-HxCDF	0.454	-	J	0.0454	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.141		-	0.0657	Total TCDF	3.89	-	M
1,2,3,4,6,7,8-HpCDF	3.06	-	J	0.0306	0.0747	Total PeCDF	1.90	-	J
1,2,3,4,7,8,9-HpCDF	0.506	-	J	0.00506	0.0883	Total HxCDF	4.59	-	J
OCDF	5.74	-	J	0.00172	0.170	Total HpCDF	7.36	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	94.7	25.0 - 164	
13C-1,2,3,7,8-PeCDD	96.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	97.4	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	91.2	23.0 - 140	
13C-OCDD	82.6	17.0 - 157	
13C-2,3,7,8-TCDF	97.1	24.0 - 169	
13C-1,2,3,7,8-PeCDF	91.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.9	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	89.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.6	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	87.8	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	86.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	87.7	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.3	26.0 - 138	
13C-OCDF	81.9	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 88.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000011 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-008-SA
Client ID: SP3-8
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.01 g
% Solids: 93.69

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 04-30-2018
2005 WHO TEQ: 1.34
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.243		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.352		-	0.0570				
1,2,3,4,7,8-HxCDD	0.422	-	J	0.0422	0.0793				
1,2,3,6,7,8-HxCDD	1.29	-	J	0.129	0.0940	Total TCDD	ND	0.243	
1,2,3,7,8,9-HxCDD	1.00	-	J	0.100	0.0823	Total PeCDD	ND	0.352	
1,2,3,4,6,7,8-HpCDD	32.3	-		0.323	0.0842	Total HxCDD	9.74	-	
OCDD	267	-		0.0801	0.172	Total HpCDD	56.5	-	
2,3,7,8-TCDF	1.04	-		0.104	0.0269				
1,2,3,7,8-PeCDF	ND	0.232		-	0.0449				
2,3,4,7,8-PeCDF	0.654	-	J	0.196	0.0468				
1,2,3,4,7,8-HxCDF	1.10	-	J	0.110	0.0437				
1,2,3,6,7,8-HxCDF	0.615	-	J	0.0615	0.0417				
2,3,4,6,7,8-HxCDF	0.724	-	J	0.0724	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.154		-	0.0657	Total TCDF	5.68	-	M
1,2,3,4,6,7,8-HpCDF	10.2	-		0.102	0.0747	Total PeCDF	6.52	-	
1,2,3,4,7,8,9-HpCDF	0.794	-	J	0.00794	0.0883	Total HxCDF	11.3	-	
OCDF	32.1	-		0.00963	0.170	Total HpCDF	29.1	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	87.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	88.4	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	92.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	88.5	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	84.8	23.0 - 140	
13C-OCDD	74.9	17.0 - 157	
13C-2,3,7,8-TCDF	90.6	24.0 - 169	
13C-1,2,3,7,8-PeCDF	83.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	84.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	83.3	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	81.9	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	82.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	82.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	82.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	84.6	26.0 - 138	
13C-OCDF	78.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 84.5 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000012 of 000022

EPA Method 1613
PCDD/F



FAL ID: 11455-009-SA
Client ID: SP3-9
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.03 g
% Solids: 92.21

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.731
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	0.470	-	J	0.470	0.0273				
1,2,3,7,8-PeCDD	ND	0.223		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.170		-	0.0793				
1,2,3,6,7,8-HxCDD	0.526	-	J	0.0526	0.0940	Total TCDD	0.470	-	J
1,2,3,7,8,9-HxCDD	0.354	-	J	0.0354	0.0823	Total PeCDD	1.69	-	J
1,2,3,4,6,7,8-HpCDD	8.05	-		0.0805	0.0842	Total HxCDD	3.49	-	J
OCDD	57.1	-		0.0171	0.172	Total HpCDD	14.6	-	
2,3,7,8-TCDF	0.546	-	J	0.0546	0.0269				
1,2,3,7,8-PeCDF	ND	0.179		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.180		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.203		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.212		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.220		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.235		-	0.0657	Total TCDF	1.29	-	
1,2,3,4,6,7,8-HpCDF	1.82	-	J	0.0182	0.0747	Total PeCDF	1.00	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.173		-	0.0883	Total HxCDF	1.77	-	J
OCDF	7.04	-	J	0.00211	0.170	Total HpCDF	5.65	-	
Internal Standards	% Rec	QC Limits	Qual						
13C-2,3,7,8-TCDD	91.9	25.0 - 164							
13C-1,2,3,7,8-PeCDD	94.3	25.0 - 181							
13C-1,2,3,4,7,8-HxCDD	97.8	32.0 - 141							
13C-1,2,3,6,7,8-HxCDD	95.7	28.0 - 130							
13C-1,2,3,4,6,7,8-HpCDD	91.1	23.0 - 140							
13C-OCDD	81.5	17.0 - 157							
13C-2,3,7,8-TCDF	95.6	24.0 - 169							
13C-1,2,3,7,8-PeCDF	91.5	24.0 - 185							
13C-2,3,4,7,8-PeCDF	91.8	21.0 - 178							
13C-1,2,3,4,7,8-HxCDF	88.0	26.0 - 152							
13C-1,2,3,6,7,8-HxCDF	85.7	26.0 - 123							
13C-2,3,4,6,7,8-HxCDF	87.3	28.0 - 136							
13C-1,2,3,7,8,9-HxCDF	86.4	29.0 - 147							
13C-1,2,3,4,6,7,8-HpCDF	88.1	28.0 - 143							
13C-1,2,3,4,7,8,9-HpCDF	90.2	26.0 - 138							
13C-OCDF	82.4	17.0 - 157							
Cleanup Surrogate									
37Cl-2,3,7,8-TCDD	85.7	35.0 - 197							

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 5/1/2018

Reviewed By: 
Date: 5/1/2018

000013 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-010-SA
Client ID: SP3-10
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.01 g
% Solids: 90.08

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.492
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.159		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.282		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.237		-	0.0793				
1,2,3,6,7,8-HxCDD	0.838	-	J	0.0838	0.0940	Total TCDD	0.436	-	J
1,2,3,7,8,9-HxCDD	0.589	-	J	0.0589	0.0823	Total PeCDD	ND	0.282	
1,2,3,4,6,7,8-HpCDD	11.2	-		0.112	0.0842	Total HxCDD	6.24	-	
OCDD	72.6	-		0.0218	0.172	Total HpCDD	20.4	-	
2,3,7,8-TCDF	0.885	-	J	0.0885	0.0269				
1,2,3,7,8-PeCDF	ND	0.186		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.198		-	0.0468				
1,2,3,4,7,8-HxCDF	0.366	-	J	0.0366	0.0437				
1,2,3,6,7,8-HxCDF	0.381	-	J	0.0381	0.0417				
2,3,4,6,7,8-HxCDF	0.262	-	J	0.0262	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.109		-	0.0657	Total TCDF	3.32	-	
1,2,3,4,6,7,8-HpCDF	2.17	-	J	0.0217	0.0747	Total PeCDF	1.37	-	J,M
1,2,3,4,7,8,9-HpCDF	0.367	-	J	0.00367	0.0883	Total HxCDF	2.68	-	J
OCDF	3.33	-	J	0.000999	0.170	Total HpCDF	4.31	-	J

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	93.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	95.3	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	101	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	91.3	23.0 - 140	
13C-OCDD	81.7	17.0 - 157	
13C-2,3,7,8-TCDF	98.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	90.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.2	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.1	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	91.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.8	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	87.3	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	86.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.5	26.0 - 138	
13C-OCDF	79.8	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 91.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000014 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-011-SA
Client ID: SP3-11
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.02 g
% Solids: 92.68

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.616
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.221		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.283		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.254		-	0.0793				
1,2,3,6,7,8-HxCDD	0.963	-	J	0.0963	0.0940	Total TCDD	ND	0.221	
1,2,3,7,8,9-HxCDD	0.676	-	J	0.0676	0.0823	Total PeCDD	ND	0.283	
1,2,3,4,6,7,8-HpCDD	16.7	-		0.167	0.0842	Total HxCDD	5.96	-	
OCDD	139	-		0.0417	0.172	Total HpCDD	31.3	-	
2,3,7,8-TCDF	0.516	-	J	0.0516	0.0269				
1,2,3,7,8-PeCDF	ND	0.247		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.264		-	0.0468				
1,2,3,4,7,8-HxCDF	0.372	-	J	0.0372	0.0437				
1,2,3,6,7,8-HxCDF	0.374	-	J	0.0374	0.0417				
2,3,4,6,7,8-HxCDF	0.374	-	J	0.0374	0.0574				
1,2,3,7,8,9-HxCDF	0.337	-	J	0.0337	0.0657	Total TCDF	1.88	-	
1,2,3,4,6,7,8-HpCDF	3.84	-	J	0.0384	0.0747	Total PeCDF	1.73	-	J
1,2,3,4,7,8,9-HpCDF	0.460	-	J	0.00460	0.0883	Total HxCDF	5.04	-	
OCDF	11.7	-		0.00351	0.170	Total HpCDF	11.1	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	90.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	96.3	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	101	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	96.8	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	92.4	23.0 - 140	
13C-OCDD	85.1	17.0 - 157	
13C-2,3,7,8-TCDF	96.9	24.0 - 169	
13C-1,2,3,7,8-PeCDF	88.6	24.0 - 185	
13C-2,3,4,7,8-PeCDF	89.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	90.9	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	88.4	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	89.2	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	87.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	88.4	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	89.8	26.0 - 138	
13C-OCDF	83.0	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 87.9 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000015 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-012-SA
Client ID: SP3-12
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.03 g
% Solids: 94.10

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 1.79
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.229		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.384		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.267		-	0.0793				
1,2,3,6,7,8-HxCDD	2.30	-	J	0.230	0.0940	Total TCDD	0.432	-	J
1,2,3,7,8,9-HxCDD	1.11	-	J	0.111	0.0823	Total PeCDD	ND	0.384	
1,2,3,4,6,7,8-HpCDD	46.8	-		0.468	0.0842	Total HxCDD	16.3	-	
OCDD	395	-		0.119	0.172	Total HpCDD	82.1	-	
2,3,7,8-TCDF	1.12	-		0.112	0.0269				
1,2,3,7,8-PeCDF	ND	0.271		-	0.0449				
2,3,4,7,8-PeCDF	0.896	-	J	0.269	0.0468				
1,2,3,4,7,8-HxCDF	1.24	-	J	0.124	0.0437				
1,2,3,6,7,8-HxCDF	1.22	-	J	0.122	0.0417				
2,3,4,6,7,8-HxCDF	0.904	-	J	0.0904	0.0574				
1,2,3,7,8,9-HpCDF	0.422	-	J	0.0422	0.0657	Total TCDF	6.85	-	
1,2,3,4,6,7,8-HpCDF	8.09	-		0.0809	0.0747	Total PeCDF	4.91	-	J
1,2,3,4,7,8,9-HpCDF	1.54	-	J	0.0154	0.0883	Total HxCDF	9.17	-	
OCDF	14.6	-		0.00438	0.170	Total HpCDF	17.9	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	78.4	25.0 - 164	
13C-1,2,3,7,8-PeCDD	86.6	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	88.3	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	84.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	80.1	23.0 - 140	
13C-OCDD	73.7	17.0 - 157	
13C-2,3,7,8-TCDF	79.7	24.0 - 169	
13C-1,2,3,7,8-PeCDF	81.9	24.0 - 185	
13C-2,3,4,7,8-PeCDF	83.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	78.6	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	77.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	77.8	28.0 - 136	
13C-1,2,3,7,8,9-HpCDF	75.6	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	77.0	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	78.5	26.0 - 138	
13C-OCDF	71.7	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 74.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000016 of 000022

EPA Method 1613 PCDD/F



FAL ID: 11455-013-SA
Client ID: SP4-1
Matrix: Solid
Batch No: X4492

Date Extracted: 04-27-2018
Date Received: 04-12-2018
Amount: 5.04 g
% Solids: 96.08

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 2.91
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.184		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.309		-	0.0570				
1,2,3,4,7,8-HxCDD	0.471	-	J	0.0471	0.0793				
1,2,3,6,7,8-HxCDD	4.20	-	J	0.420	0.0940	Total TCDD	0.754	-	J
1,2,3,7,8,9-HxCDD	1.88	-	J	0.188	0.0823	Total PeCDD	1.40	-	J
1,2,3,4,6,7,8-HpCDD	79.8	-		0.798	0.0842	Total HxCDD	32.4	-	
OCDD	683	-		0.205	0.172	Total HpCDD	150	-	
2,3,7,8-TCDF	1.73	-		0.173	0.0269				
1,2,3,7,8-PeCDF	0.461	-	J	0.0138	0.0449				
2,3,4,7,8-PeCDF	1.02	-	J	0.306	0.0468				
1,2,3,4,7,8-HxCDF	1.81	-	J	0.181	0.0437				
1,2,3,6,7,8-HxCDF	1.98	-	J	0.198	0.0417				
2,3,4,6,7,8-HxCDF	1.53	-	J	0.153	0.0574				
1,2,3,7,8,9-HxCDF	0.615	-	J	0.0615	0.0657	Total TCDF	12.4	-	
1,2,3,4,6,7,8-HpCDF	13.8	-		0.138	0.0747	Total PeCDF	7.70	-	
1,2,3,4,7,8,9-HpCDF	2.31	-	J	0.0231	0.0883	Total HxCDF	16.5	-	
OCDF	24.5	-		0.00735	0.170	Total HpCDF	31.6	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	88.8	25.0 - 164	
13C-1,2,3,7,8-PeCDD	92.7	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	92.9	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	89.9	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	86.4	23.0 - 140	
13C-OCDD	78.9	17.0 - 157	
13C-2,3,7,8-TCDF	94.0	24.0 - 169	
13C-1,2,3,7,8-PeCDF	89.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	92.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	81.9	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	81.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	80.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	82.0	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	82.5	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	86.5	26.0 - 138	
13C-OCDF	78.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 85.3 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/1/2018

Reviewed By: 

Date: 5/1/2018

000017 of 000022



Frontier Analytical Laboratory
5172 Hillsdale Circle
El Dorado Hills, CA 95762
Tel: 916-934-0900
Fax: 916-934-0999

FAL USE ONLY

Laboratory Project No.: 11455
Temperature: 0 °C

Chain of Custody

www.frontieranalytical.com

Please Print in Pen Page 1 of 1

CLIENT INFORMATION				INVOICE INFORMATION (if different from client info)				PROJECT INFORMATION								
Company Name: Brusca Associates, Inc.				Company Name: Same				FAL Quote #: 3310								
Contact Name: Joe Brusca				Contact Name:				P.O. #:								
Address: PO Box 332, Roseville, CA 95661				Address:				Project #: 137-002								
Phone: 916-677-1470 Fax:				Phone: Fax:				Project Name: LATROP 48-Ac Property								
Email: jbrusca@bruscaassociates.com				Email:				TAT (business days): ☐ 15 ☒ 10 ☐ 5* ☐ 3* (✓ one)								
								* FAL must agree with price and RUSH TAT in writing.								
REPORT INFORMATION				REPORT DISTRIBUTION (email only is preferred)				ADDITIONAL INSTRUCTIONS								
Report Level: ☐ I/II ☐ III ☐ IV				☐ Hardcopy				Joe to Kathy 15-TAT 4/12/18								
☐ EDD: ☐ FAL Basic ☐ Geotracker				☐ CD (.pdf including EDDs if requested)												
☐ Other: ☐ Custom: Contact FAL				☐ Email (.pdf including EDDs if requested)												
☐ California State Drinking Water Form																
System #: Source #:																
Sampler: Employer:																
Sample ID	Date	Time	Matrix	# of containers	EPA 1613**	EPA 8290**	DLM 02.0	EPA 8280**	Appendix IX	EPA TO-9/A	EPA 23/23A	EPA 1668	FAL 15	Other	**CONGENERS	**TEQ
		Collected													☐ 2,3,7,8-TCDD only	☐ 1998 WHO
															☐ 2,3,7,8-TCDD/F only	☒ 2005 WHO
															☒ PCDD/F (Cl ₄ -Cl ₈)	☐ Other
															Remarks	
1	SP3-1	4/12/18 10:51	Solo		X											
2	SP3-2	10:55			X											
3	SP3-3	11:00			X											
4	SP3-4	11:02			X											
5	SP3-5	11:05			X											
6	SP3-6	11:09			X											
7	SP3-7	11:14			X											
8	SP3-8	11:16			X											
9	SP3-9	11:19			X											
10	SP3-10	11:21			X											
11	SP3-11	11:23			X											
12	SP3-12	11:26			X											
13	SP4-1	11:35			X											
14																
15																
Samples will be disposed of 90 days after sample receipt unless other arrangements have been made and agreed upon in writing.																
Relinquished by: (Signature and Printed Name)				Date	Time	Received by: (Signature and Printed Name)				Date	Time					
[Signature]				4/12/18	3:03P	Kathy [Signature]				4/12/18	1503					

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed.
FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

White Copy - Report

Yellow Copy - Laboratory

Pink Copy - Originator

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11455**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/12/2018
Time Received:	03:03 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	13
Duplicates:	0
Storage Location:	R-4

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

20 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/13/18 09:30. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP3-1	T181249-01	Soil	04/12/18 10:51	04/13/18 09:30
SP3-2	T181249-02	Soil	04/12/18 10:55	04/13/18 09:30
SP3-3	T181249-03	Soil	04/12/18 11:00	04/13/18 09:30
SP3-4	T181249-04	Soil	04/12/18 11:02	04/13/18 09:30
SP3-5	T181249-05	Soil	04/12/18 11:05	04/13/18 09:30
SP3-6	T181249-06	Soil	04/12/18 11:09	04/13/18 09:30
SP3-7	T181249-07	Soil	04/12/18 11:14	04/13/18 09:30
SP3-8	T181249-08	Soil	04/12/18 11:16	04/13/18 09:30
SP3-9	T181249-09	Soil	04/12/18 11:19	04/13/18 09:30
SP3-10	T181249-10	Soil	04/12/18 11:21	04/13/18 09:30
SP3-11	T181249-11	Soil	04/12/18 11:23	04/13/18 09:30
SP3-12	T181249-12	Soil	04/12/18 11:26	04/13/18 09:30
SP4-1	T181249-13	Soil	04/12/18 11:35	04/13/18 09:30

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



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Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

DETECTIONS SUMMARY

Sample ID: SP3-1

Laboratory ID: T181249-01

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	29	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	29	10	mg/kg	EPA 8015B	
Barium	61	1.0	mg/kg	EPA 6010B	
Chromium	5.8	2.0	mg/kg	EPA 6010B	
Cobalt	4.5	2.0	mg/kg	EPA 6010B	
Copper	4.5	1.0	mg/kg	EPA 6010B	
Nickel	6.0	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	30	1.0	mg/kg	EPA 6010B	
Pyrene	650	300	ug/kg	EPA 8270C	
Benzo (a) anthracene	300	300	ug/kg	EPA 8270C	
Benzo (b) fluoranthene	300	300	ug/kg	EPA 8270C	
Chrysene	340	300	ug/kg	EPA 8270C	
Fluoranthene	730	300	ug/kg	EPA 8270C	
Phenanthrene	460	300	ug/kg	EPA 8270C	

Sample ID: SP3-2

Laboratory ID: T181249-02

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	150	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	160	10	mg/kg	EPA 8015B	
Barium	72	1.0	mg/kg	EPA 6010B	
Chromium	5.6	2.0	mg/kg	EPA 6010B	
Cobalt	4.7	2.0	mg/kg	EPA 6010B	
Copper	8.0	1.0	mg/kg	EPA 6010B	
Nickel	6.6	2.0	mg/kg	EPA 6010B	
Vanadium	17	5.0	mg/kg	EPA 6010B	
Zinc	33	1.0	mg/kg	EPA 6010B	
PCB-1254	20	10	ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Sample ID: SP3-3

Laboratory ID: T181249-03

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	51	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	37	10		mg/kg	EPA 8015B	
Barium	41	1.0		mg/kg	EPA 6010B	
Chromium	4.4	2.0		mg/kg	EPA 6010B	
Cobalt	3.5	2.0		mg/kg	EPA 6010B	
Copper	5.0	1.0		mg/kg	EPA 6010B	
Lead	20	3.0		mg/kg	EPA 6010B	
Nickel	4.1	2.0		mg/kg	EPA 6010B	
Vanadium	10	5.0		mg/kg	EPA 6010B	
Zinc	24	1.0		mg/kg	EPA 6010B	
PCB-1254	17	10		ug/kg	EPA 8082	

Sample ID: SP3-4

Laboratory ID: T181249-04

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	100	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	96	10		mg/kg	EPA 8015B	
Barium	53	1.0		mg/kg	EPA 6010B	
Chromium	5.1	2.0		mg/kg	EPA 6010B	
Cobalt	4.4	2.0		mg/kg	EPA 6010B	
Copper	4.3	1.0		mg/kg	EPA 6010B	
Nickel	5.7	2.0		mg/kg	EPA 6010B	
Vanadium	15	5.0		mg/kg	EPA 6010B	
Zinc	26	1.0		mg/kg	EPA 6010B	
PCB-1254	15	10		ug/kg	EPA 8082	

Sample ID: SP3-5

Laboratory ID: T181249-05

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	32	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	15	10		mg/kg	EPA 8015B	
Barium	64	1.0		mg/kg	EPA 6010B	
Chromium	4.9	2.0		mg/kg	EPA 6010B	
Cobalt	4.2	2.0		mg/kg	EPA 6010B	
Copper	5.2	1.0		mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Sample ID: SP3-5

Laboratory ID: T181249-05

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Nickel	5.1	2.0	mg/kg	EPA 6010B	
Vanadium	15	5.0	mg/kg	EPA 6010B	
Zinc	29	1.0	mg/kg	EPA 6010B	

Sample ID: SP3-6

Laboratory ID: T181249-06

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	56	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	32	10	mg/kg	EPA 8015B	
Barium	43	1.0	mg/kg	EPA 6010B	
Chromium	5.1	2.0	mg/kg	EPA 6010B	
Cobalt	3.8	2.0	mg/kg	EPA 6010B	
Copper	4.4	1.0	mg/kg	EPA 6010B	
Lead	14	3.0	mg/kg	EPA 6010B	
Nickel	4.5	2.0	mg/kg	EPA 6010B	
Vanadium	11	5.0	mg/kg	EPA 6010B	
Zinc	23	1.0	mg/kg	EPA 6010B	
PCB-1254	12	10	ug/kg	EPA 8082	

Sample ID: SP3-7

Laboratory ID: T181249-07

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	39	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	19	10	mg/kg	EPA 8015B	
Barium	43	1.0	mg/kg	EPA 6010B	
Chromium	4.6	2.0	mg/kg	EPA 6010B	
Cobalt	3.8	2.0	mg/kg	EPA 6010B	
Copper	3.3	1.0	mg/kg	EPA 6010B	
Lead	10	3.0	mg/kg	EPA 6010B	
Nickel	4.1	2.0	mg/kg	EPA 6010B	
Vanadium	11	5.0	mg/kg	EPA 6010B	
Zinc	25	1.0	mg/kg	EPA 6010B	
PCB-1254	12	10	ug/kg	EPA 8082	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Sample ID: SP3-8

Laboratory ID: T181249-08

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	29	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	17	10		mg/kg	EPA 8015B	
Barium	67	1.0		mg/kg	EPA 6010B	
Chromium	5.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.9	2.0		mg/kg	EPA 6010B	
Copper	4.6	1.0		mg/kg	EPA 6010B	
Lead	17	3.0		mg/kg	EPA 6010B	
Nickel	6.6	2.0		mg/kg	EPA 6010B	
Vanadium	17	5.0		mg/kg	EPA 6010B	
Zinc	26	1.0		mg/kg	EPA 6010B	
PCB-1254	19	10		ug/kg	EPA 8082	

Sample ID: SP3-9

Laboratory ID: T181249-09

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	85	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	82	10		mg/kg	EPA 8015B	
Barium	59	1.0		mg/kg	EPA 6010B	
Chromium	5.9	2.0		mg/kg	EPA 6010B	
Cobalt	5.6	2.0		mg/kg	EPA 6010B	
Copper	4.8	1.0		mg/kg	EPA 6010B	
Nickel	6.6	2.0		mg/kg	EPA 6010B	
Vanadium	18	5.0		mg/kg	EPA 6010B	
Zinc	29	1.0		mg/kg	EPA 6010B	
PCB-1254	16	10		ug/kg	EPA 8082	

Sample ID: SP3-10

Laboratory ID: T181249-10

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	360	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	380	10		mg/kg	EPA 8015B	
Barium	57	1.0		mg/kg	EPA 6010B	
Chromium	5.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.9	2.0		mg/kg	EPA 6010B	
Copper	4.6	1.0		mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Sample ID: SP3-10

Laboratory ID: T181249-10

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Nickel	5.7	2.0	mg/kg	EPA 6010B	
Vanadium	14	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	
PCB-1254	18	10	ug/kg	EPA 8082	

Sample ID: SP3-11

Laboratory ID: T181249-11

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	26	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	33	10	mg/kg	EPA 8015B	
Barium	56	1.0	mg/kg	EPA 6010B	
Chromium	5.0	2.0	mg/kg	EPA 6010B	
Cobalt	4.4	2.0	mg/kg	EPA 6010B	
Copper	4.0	1.0	mg/kg	EPA 6010B	
Nickel	5.7	2.0	mg/kg	EPA 6010B	
Vanadium	15	5.0	mg/kg	EPA 6010B	
Zinc	28	1.0	mg/kg	EPA 6010B	

Sample ID: SP3-12

Laboratory ID: T181249-12

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	140	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	130	10	mg/kg	EPA 8015B	
Barium	38	1.0	mg/kg	EPA 6010B	
Chromium	3.8	2.0	mg/kg	EPA 6010B	
Cobalt	3.9	2.0	mg/kg	EPA 6010B	
Copper	7.6	1.0	mg/kg	EPA 6010B	
Lead	38	3.0	mg/kg	EPA 6010B	
Nickel	3.6	2.0	mg/kg	EPA 6010B	
Vanadium	8.3	5.0	mg/kg	EPA 6010B	
Zinc	22	1.0	mg/kg	EPA 6010B	
PCB-1254	34	10	ug/kg	EPA 8082	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Sample ID: SP4-1

Laboratory ID: T181249-13

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	170	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	220	10		mg/kg	EPA 8015B	
Barium	31	1.0		mg/kg	EPA 6010B	
Chromium	3.5	2.0		mg/kg	EPA 6010B	
Cobalt	3.2	2.0		mg/kg	EPA 6010B	
Copper	7.0	1.0		mg/kg	EPA 6010B	
Lead	25	3.0		mg/kg	EPA 6010B	
Nickel	2.9	2.0		mg/kg	EPA 6010B	
Vanadium	7.8	5.0		mg/kg	EPA 6010B	
Zinc	21	1.0		mg/kg	EPA 6010B	
PCB-1016	170	10		ug/kg	EPA 8082	
PCB-1254	110	10		ug/kg	EPA 8082	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-1
T181249-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	29	10	"	"	"	"	"	"	
C29-C40 (MORO)	29	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		88.9 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	61	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.8	2.0	"	"	"	"	"	"	
Cobalt	4.5	2.0	"	"	"	"	"	"	
Copper	4.5	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.0	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
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SP3-1

T181249-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		78.9 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		76.1 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	650	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	300	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	300	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	
Benzo (a) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-1

T181249-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzyl alcohol	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	340	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	730	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	
Isophorone	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-1

T181249-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

2-Methylphenol	ND	1000	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	460	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		44.2 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		53.9 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		54.1 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		68.9 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		88.8 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		73.1 %	29.1-130		"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-2

T181249-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	150	10	"	"	"	"	"	"	
C29-C40 (MORO)	160	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		79.5 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	72	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.6	2.0	"	"	"	"	"	"	
Cobalt	4.7	2.0	"	"	"	"	"	"	
Copper	8.0	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	33	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-2

T181249-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	20	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		91.4 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		90.7 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-2

T181249-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-2

T181249-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	45.7 %	15-121			"	"	"	"
Surrogate: Phenol-d6	58.5 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	57.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.0 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	83.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	75.7 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-3

T181249-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	51	10	"	"	"	"	"	"	
C29-C40 (MORO)	37	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		84.4 %		65-135	"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	41	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.4	2.0	"	"	"	"	"	"	
Cobalt	3.5	2.0	"	"	"	"	"	"	
Copper	5.0	1.0	"	"	"	"	"	"	
Lead	20	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	10	5.0	"	"	"	"	"	"	
Zinc	24	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-3

T181249-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	17	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		96.1 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		94.5 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-3

T181249-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-3

T181249-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		51.2 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		59.0 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		59.2 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		69.2 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		85.0 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		73.2 %	29.1-130		"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-4

T181249-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	100	10	"	"	"	"	"	"	
C29-C40 (MORO)	96	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		88.1 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	53	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.1	2.0	"	"	"	"	"	"	
Cobalt	4.4	2.0	"	"	"	"	"	"	
Copper	4.3	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.7	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	26	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

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04/20/18 11:17

SP3-4

T181249-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	15	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		81.2 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		67.1 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-4

T181249-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-4

T181249-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol		38.5 %	15-121		"	"	"	"
Surrogate: Phenol-d6		47.3 %	24-113		"	"	"	"
Surrogate: Nitrobenzene-d5		44.5 %	21.3-119		"	"	"	"
Surrogate: 2-Fluorobiphenyl		58.9 %	32.4-102		"	"	"	"
Surrogate: 2,4,6-Tribromophenol		81.1 %	18.1-105		"	"	"	"
Surrogate: Terphenyl-d14		75.8 %	29.1-130		"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-5
T181249-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	32	10	"	"	"	"	"	"	
C29-C40 (MORO)	15	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		91.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	64	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.9	2.0	"	"	"	"	"	"	
Cobalt	4.2	2.0	"	"	"	"	"	"	
Copper	5.2	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Reported:
04/20/18 11:17

SP3-5
T181249-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	

Surrogate: Tetrachloro-meta-xylene

92.6 % 35-140

" " " "

Surrogate: Decachlorobiphenyl

85.5 % 35-140

" " " "

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-5

T181249-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-5
T181249-05 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	43.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	51.4 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	50.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	57.7 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	78.7 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	64.3 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-6
T181249-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	56	10	"	"	"	"	"	"	
C29-C40 (MORO)	32	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		84.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	43	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.1	2.0	"	"	"	"	"	"	
Cobalt	3.8	2.0	"	"	"	"	"	"	
Copper	4.4	1.0	"	"	"	"	"	"	
Lead	14	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.5	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	11	5.0	"	"	"	"	"	"	
Zinc	23	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Reported:
04/20/18 11:17

SP3-6
T181249-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	12	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		81.4 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		92.4 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-6

T181249-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-6

T181249-06 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	40.9 %	15-121			"	"	"	"
Surrogate: Phenol-d6	48.5 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	48.9 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	61.2 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	86.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	76.3 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-7

T181249-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	39	10	"	"	"	"	"	"	
C29-C40 (MORO)	19	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		86.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	43	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	4.6	2.0	"	"	"	"	"	"	
Cobalt	3.8	2.0	"	"	"	"	"	"	
Copper	3.3	1.0	"	"	"	"	"	"	
Lead	10	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	4.1	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	11	5.0	"	"	"	"	"	"	
Zinc	25	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-7

T181249-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	12	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		88.8 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		91.8 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-7

T181249-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-7

T181249-07 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		41.1 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		47.3 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		49.2 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		60.5 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		74.5 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		70.2 %	29.1-130		"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-8

T181249-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	29	10	"	"	"	"	"	"	
C29-C40 (MORO)	17	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		88.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	67	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.7	2.0	"	"	"	"	"	"	
Cobalt	4.9	2.0	"	"	"	"	"	"	
Copper	4.6	1.0	"	"	"	"	"	"	
Lead	17	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	17	5.0	"	"	"	"	"	"	
Zinc	26	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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04/20/18 11:17

SP3-8

T181249-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	19	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		84.8 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		82.3 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-8

T181249-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-8
T181249-08 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	47.6 %	15-121			"	"	"	"
Surrogate: Phenol-d6	57.6 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	58.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	69.6 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	91.0 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	77.5 %	29.1-130			"	"	"	"

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-9

T181249-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	85	10	"	"	"	"	"	"	
C29-C40 (MORO)	82	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		80.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	59	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.9	2.0	"	"	"	"	"	"	
Cobalt	5.6	2.0	"	"	"	"	"	"	
Copper	4.8	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	18	5.0	"	"	"	"	"	"	
Zinc	29	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-9

T181249-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	16	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		89.7 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		69.3 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-9

T181249-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-9

T181249-09 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	52.2 %	15-121			"	"	"	"
Surrogate: Phenol-d6	61.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	60.1 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	73.5 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	90.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	75.3 %	29.1-130			"	"	"	"

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Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-10
T181249-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	360	10	"	"	"	"	"	"	
C29-C40 (MORO)	380	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		80.0 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	57	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.7	2.0	"	"	"	"	"	"	
Cobalt	4.9	2.0	"	"	"	"	"	"	
Copper	4.6	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.7	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	14	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Reported:
04/20/18 11:17

SP3-10
T181249-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	18	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		81.2 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		61.2 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-10

T181249-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-10
T181249-10 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	54.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	65.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	62.5 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	73.1 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	87.8 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	80.1 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-11
T181249-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	26	10	"	"	"	"	"	"	
C29-C40 (MORO)	33	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		79.6 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	56	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	5.0	2.0	"	"	"	"	"	"	
Cobalt	4.4	2.0	"	"	"	"	"	"	
Copper	4.0	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	5.7	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	15	5.0	"	"	"	"	"	"	
Zinc	28	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Reported:
04/20/18 11:17

SP3-11
T181249-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		87.7 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		83.4 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-11
T181249-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-11
T181249-11 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	48.4 %	15-121			"	"	"	"
Surrogate: Phenol-d6	59.2 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	55.9 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	68.4 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	95.9 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	69.3 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-12
T181249-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	140	10	"	"	"	"	"	"	
C29-C40 (MORO)	130	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		84.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	38	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	3.8	2.0	"	"	"	"	"	"	
Cobalt	3.9	2.0	"	"	"	"	"	"	
Copper	7.6	1.0	"	"	"	"	"	"	
Lead	38	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	3.6	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	8.3	5.0	"	"	"	"	"	"	
Zinc	22	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

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Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-12
T181249-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	34	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		94.7 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		74.2 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-12
T181249-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP3-12
T181249-12 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	45.2 %	15-121			"	"	"	"
Surrogate: Phenol-d6	53.8 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	53.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	63.8 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	82.6 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	68.3 %	29.1-130			"	"	"	"

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP4-1

T181249-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041323	04/13/18	04/14/18	EPA 8015B	
C13-C28 (DRO)	170	10	"	"	"	"	"	"	
C29-C40 (MORO)	220	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		84.7 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041326	04/13/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	31	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	3.5	2.0	"	"	"	"	"	"	
Cobalt	3.2	2.0	"	"	"	"	"	"	
Copper	7.0	1.0	"	"	"	"	"	"	
Lead	25	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	2.9	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	7.8	5.0	"	"	"	"	"	"	
Zinc	21	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041325	04/13/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP4-1

T181249-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	170	10	ug/kg	1	8041319	04/13/18	04/17/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	110	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		88.1 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		69.7 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP4-1

T181249-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

SP4-1

T181249-13 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041321	04/13/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	48.0 %	15-121			"	"	"	"
Surrogate: Phenol-d6	56.9 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	58.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	71.7 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	84.5 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	76.6 %	29.1-130			"	"	"	"

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Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041323 - EPA 3550B GC

Blank (8041323-BLK1)

Prepared: 04/13/18 Analyzed: 04/14/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							
Surrogate: <i>p</i> -Terphenyl	90.7		"	102		88.9	65-135			

LCS (8041323-BS1)

Prepared: 04/13/18 Analyzed: 04/14/18

C13-C28 (DRO)	410	10	mg/kg	490		83.7	75-125			
Surrogate: <i>p</i> -Terphenyl	79.2		"	98.0		80.8	65-135			

LCS Dup (8041323-BSD1)

Prepared: 04/13/18 Analyzed: 04/14/18

C13-C28 (DRO)	490	10	mg/kg	510		95.4	75-125	17.0	20	
Surrogate: <i>p</i> -Terphenyl	85.0		"	102		83.3	65-135			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041326 - EPA 3050B

Blank (8041326-BLK1)

Prepared: 04/13/18 Analyzed: 04/16/18

Antimony	ND	3.0	mg/kg							
Silver	ND	2.0	"							
Arsenic	ND	5.0	"							
Barium	ND	1.0	"							
Beryllium	ND	1.0	"							
Cadmium	ND	2.0	"							
Chromium	ND	2.0	"							
Cobalt	ND	2.0	"							
Copper	ND	1.0	"							
Lead	ND	3.0	"							
Molybdenum	ND	5.0	"							
Nickel	ND	2.0	"							
Selenium	ND	5.0	"							
Thallium	ND	2.0	"							
Vanadium	ND	5.0	"							
Zinc	ND	1.0	"							

LCS (8041326-BS1)

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	96.4	5.0	mg/kg	100		96.4	75-125			
Barium	97.6	1.0	"	100		97.6	75-125			
Cadmium	97.8	2.0	"	100		97.8	75-125			
Chromium	97.4	2.0	"	100		97.4	75-125			
Lead	105	3.0	"	100		105	75-125			

Matrix Spike (8041326-MS1)

Source: T181248-01

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	82.6	5.0	mg/kg	96.2	1.96	83.9	75-125			
Barium	195	1.0	"	96.2	125	72.2	75-125			QM-05
Cadmium	84.0	2.0	"	96.2	0.911	86.4	75-125			
Chromium	109	2.0	"	96.2	29.8	82.6	75-125			
Lead	96.6	3.0	"	96.2	ND	100	75-125			

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Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041326 - EPA 3050B

Matrix Spike Dup (8041326-MSD1)

Source: T181248-01

Prepared: 04/13/18 Analyzed: 04/16/18

Arsenic	86.4	5.0	mg/kg	94.3	1.96	89.6	75-125	4.51	20	
Barium	187	1.0	"	94.3	125	65.2	75-125	4.15	20	QM-05
Cadmium	82.3	2.0	"	94.3	0.911	86.3	75-125	2.00	20	
Chromium	111	2.0	"	94.3	29.8	86.3	75-125	1.80	20	
Lead	95.6	3.0	"	94.3	ND	101	75-125	1.02	20	

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Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041325 - EPA 7471A Soil

Blank (8041325-BLK1)

Prepared: 04/13/18 Analyzed: 04/16/18

Mercury	ND	0.10	mg/kg

LCS (8041325-BS1)

Prepared: 04/13/18 Analyzed: 04/16/18

Mercury	0.413	0.10	mg/kg	0.397	104	80-120

Matrix Spike (8041325-MS1)

Source: T181248-01

Prepared: 04/13/18 Analyzed: 04/16/18

Mercury	0.460	0.10	mg/kg	0.403	0.0498	102	75-125

Matrix Spike Dup (8041325-MSD1)

Source: T181248-01

Prepared: 04/13/18 Analyzed: 04/16/18

Mercury	0.470	0.10	mg/kg	0.410	0.0498	103	75-125	2.10	20

SunStar Laboratories, Inc.

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041319 - EPA 3550 ECD/GCMS

Blank (8041319-BLK1)

Prepared: 04/13/18 Analyzed: 04/17/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	9.06		"	10.2		88.8	35-140			
Surrogate: Decachlorobiphenyl	10.6		"	10.2		104	35-140			

LCS (8041319-BS1)

Prepared: 04/13/18 Analyzed: 04/17/18

PCB-1016	59.6	10	ug/kg	100		59.6	40-130			
PCB-1260	64.4	10	"	100		64.4	40-130			
Surrogate: Tetrachloro-meta-xylene	9.02		"	10.0		90.2	35-140			
Surrogate: Decachlorobiphenyl	10.7		"	10.0		107	35-140			

Matrix Spike (8041319-MS1)

Source: T181249-01

Prepared: 04/13/18 Analyzed: 04/16/18

PCB-1016	71.5	10	ug/kg	99.0	ND	72.2	40-130			
PCB-1260	62.3	10	"	99.0	ND	62.9	40-130			
Surrogate: Tetrachloro-meta-xylene	7.61		"	9.90		76.9	35-140			
Surrogate: Decachlorobiphenyl	8.74		"	9.90		88.2	35-140			

Matrix Spike Dup (8041319-MSD1)

Source: T181249-01

Prepared: 04/13/18 Analyzed: 04/16/18

PCB-1016	68.7	10	ug/kg	102	ND	67.3	40-130	4.01	30	
PCB-1260	56.7	10	"	102	ND	55.6	40-130	9.31	30	
Surrogate: Tetrachloro-meta-xylene	7.48		"	10.2		73.3	35-140			
Surrogate: Decachlorobiphenyl	8.06		"	10.2		79.0	35-140			

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04/20/18 11:17

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041321 - EPA 3550 ECD/GCMS

Blank (8041321-BLK1)

Prepared: 04/13/18 Analyzed: 04/17/18

Carbazole	ND	300	ug/kg
Phenol	ND	1000	"
Aniline	ND	300	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
2-Methylnaphthalene	ND	300	"
1-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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04/20/18 11:17

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041321 - EPA 3550 ECD/GCMS

Blank (8041321-BLK1)

Prepared: 04/13/18 Analyzed: 04/17/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
Pyridine	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
2,4,6-Trichlorophenol	ND	1000	"

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Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041321 - EPA 3550 ECD/GCMS

Blank (8041321-BLK1)

Prepared: 04/13/18 Analyzed: 04/17/18

Surrogate: 2-Fluorophenol	1750		ug/kg	3300		53.0	15-121			
Surrogate: Phenol-d6	1930		"	3300		58.6	24-113			
Surrogate: Nitrobenzene-d5	2070		"	3300		62.7	21.3-119			
Surrogate: 2-Fluorobiphenyl	2330		"	3300		70.5	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2680		"	3300		81.2	18.1-105			
Surrogate: Terphenyl-d14	2480		"	3300		75.1	29.1-130			

LCS (8041321-BS1)

Prepared: 04/13/18 Analyzed: 04/17/18

Phenol	1550	1000	ug/kg	3290		47.0	34-114			
2-Chlorophenol	1700	1000	"	3290		51.6	34-114			
1,4-Dichlorobenzene	1760	300	"	3290		53.5	34-114			
N-Nitrosodi-n-propylamine	1880	300	"	3290		57.2	30-110			
1,2,4-Trichlorobenzene	2080	300	"	3290		63.3	39-119			
4-Chloro-3-methylphenol	2210	1000	"	3290		67.2	50-130			
Acenaphthene	2210	300	"	3290		67.2	34-114			
Pentachlorophenol	2380	1000	"	3290		72.4	50-130			
Pyrene	1810	300	"	3290		55.0	30-110			
Surrogate: 2-Fluorophenol	1430		"	3290		43.4	15-121			
Surrogate: Phenol-d6	1580		"	3290		48.0	24-113			
Surrogate: Nitrobenzene-d5	1720		"	3290		52.2	21.3-119			
Surrogate: 2-Fluorobiphenyl	2080		"	3290		63.4	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2520		"	3290		76.6	18.1-105			
Surrogate: Terphenyl-d14	2200		"	3290		67.0	29.1-130			

Matrix Spike (8041321-MS1)

Source: T181249-01

Prepared: 04/13/18 Analyzed: 04/16/18

Phenol	1770	1000	ug/kg	3300	ND	53.8	34-114			
2-Chlorophenol	1820	1000	"	3300	ND	55.2	34-114			
1,4-Dichlorobenzene	1950	300	"	3300	ND	59.0	34-114			
N-Nitrosodi-n-propylamine	1860	300	"	3300	ND	56.3	30-110			
1,2,4-Trichlorobenzene	2380	300	"	3300	ND	72.1	39-119			
4-Chloro-3-methylphenol	2700	1000	"	3300	ND	81.7	50-130			
Acenaphthene	2440	300	"	3300	ND	73.8	34-114			
Pentachlorophenol	2510	1000	"	3300	ND	76.2	50-130			
Pyrene	2420	300	"	3300	654	53.5	30-110			
Surrogate: 2-Fluorophenol	1590		"	3300		48.3	15-121			
Surrogate: Phenol-d6	1870		"	3300		56.6	24-113			

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041321 - EPA 3550 ECD/GCMS

Matrix Spike (8041321-MS1)		Source: T181249-01		Prepared: 04/13/18 Analyzed: 04/16/18						
Surrogate: Nitrobenzene-d5	1850		ug/kg	3300		56.1	21.3-119			
Surrogate: 2-Fluorobiphenyl	2380		"	3300		72.0	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2780		"	3300		84.2	18.1-105			
Surrogate: Terphenyl-d14	2330		"	3300		70.7	29.1-130			
Matrix Spike Dup (8041321-MSD1)		Source: T181249-01		Prepared: 04/13/18 Analyzed: 04/16/18						
Phenol	1630	1000	ug/kg	3260	ND	49.9	34-114	8.72	42	
2-Chlorophenol	1730	1000	"	3260	ND	53.2	34-114	4.93	40	
1,4-Dichlorobenzene	1650	300	"	3260	ND	50.7	34-114	16.4	28	
N-Nitrosodi-n-propylamine	1830	300	"	3260	ND	56.3	30-110	1.26	38	
1,2,4-Trichlorobenzene	2110	300	"	3260	ND	64.8	39-119	12.0	28	
4-Chloro-3-methylphenol	2550	1000	"	3260	ND	78.2	50-130	5.71	42	
Acenaphthene	2430	300	"	3260	ND	74.7	34-114	0.100	31	
Pentachlorophenol	2500	1000	"	3260	ND	76.7	50-130	0.579	50	
Pyrene	3330	300	"	3260	654	82.1	30-110	31.7	31	QM-07
Surrogate: 2-Fluorophenol	1310		"	3260		40.3	15-121			
Surrogate: Phenol-d6	1680		"	3260		51.7	24-113			
Surrogate: Nitrobenzene-d5	1690		"	3260		51.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2170		"	3260		66.6	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2790		"	3260		85.5	18.1-105			
Surrogate: Terphenyl-d14	2440		"	3260		74.9	29.1-130			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/20/18 11:17

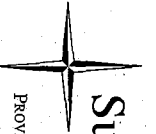
Notes and Definitions

QM-07	The spike recovery and or RPD was outside acceptance limits for the MS and/or MSD. The batch was accepted based on acceptable LCS recovery.
QM-05	The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



SunStar
Laboratories, Inc.

Chain of Custody Record

PROVIDING QUALITY ANALYTICAL SERVICES NATIONWIDE
25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Client: BRUSCA ASSOCIATES, INC.
Address: PO Box 332, Rosemead, CA 91661
Phone: (916) 677-1470 Fax: (916) 677-1471
Project Manager: Joe Brusca

Date: 4/12/18 Page: 1 of 1
Project Name: LATHROP 43-Ac Property
Collector: BRUSCA Client Project #: 137-002
Batch #: 7/8/247 EDF #:

Sample ID	Date Sampled	Time	Sample Type	Container Type	8260	8260 + OXY	8260 BTEX, OXY only	8270	8021 BTEX	8015M (gasoline)	8015M (diesel)	8015M Ext./Carbon Chain	6010/7000 Title 22 Metals	6020 ICP-MS Metals	Laboratory ID #	Comments/Preservative	Total # of containers
SP3-1	4/12/18	10:51	SOL	JAR	X	X	X	X	X	X	X	X	X	X	01		1
SP3-2		10:55			X	X	X	X	X	X	X	X	X	X	02		1
SP3-3		11:00			X	X	X	X	X	X	X	X	X	X	03		1
SP3-4		11:02			X	X	X	X	X	X	X	X	X	X	04		1
SP3-5		11:05			X	X	X	X	X	X	X	X	X	X	05		1
SP3-6		11:09			X	X	X	X	X	X	X	X	X	X	06		1
SP3-7		11:14			X	X	X	X	X	X	X	X	X	X	07		1
SP3-8		11:16			X	X	X	X	X	X	X	X	X	X	08		1
SP3-9		11:19			X	X	X	X	X	X	X	X	X	X	09		1
SP3-10		11:21			X	X	X	X	X	X	X	X	X	X	10		1
SP3-11		11:23			X	X	X	X	X	X	X	X	X	X	11		1
SP3-12		11:26			X	X	X	X	X	X	X	X	X	X	12		1
SP4-1		11:35			X	X	X	X	X	X	X	X	X	X	13		1
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18-15:45</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 15:45</u>																	
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:50</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u>																	
Relinquished by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u> Received by: (signature) <u>[Signature]</u> Date / Time <u>4/12/18 9:30</u>																	

Turn around time: NORMAL

Chain of Custody seals Y/N/A
Seals intact? Y/N/A
Received good condition/cold 3/6

Notes

Sample disposal instructions: Disposal @ \$2.00 each

Return to client

Pickup

COC 160914



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

27 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/14/18 08:10. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP5-1	T181265-01	Soil	04/13/18 10:00	04/14/18 08:10

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

DETECTIONS SUMMARY

Sample ID: SP5-1

Laboratory ID: T181265-01

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Arsenic	31	1.0		ug/l	6020 ICP-MS	STLC
Barium	400	1.0		ug/l	6020 ICP-MS	STLC
Beryllium	1.2	1.0		ug/l	6020 ICP-MS	STLC
Chromium	47	1.0		ug/l	6020 ICP-MS	STLC
Cobalt	32	1.0		ug/l	6020 ICP-MS	STLC
Copper	55	1.0		ug/l	6020 ICP-MS	STLC
Lead	46	2.0		ug/l	6020 ICP-MS	STLC
Molybdenum	7.3	1.0		ug/l	6020 ICP-MS	STLC
Nickel	47	1.0		ug/l	6020 ICP-MS	STLC
Vanadium	200	1.0		ug/l	6020 ICP-MS	STLC
Zinc	220	1.0		ug/l	6020 ICP-MS	STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commercentre Drive
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

SP5-1
T181265-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Metals by EPA 6020 Method

Antimony	ND	1.0	ug/l	1	8042318	04/23/18	04/26/18	6020 ICP-MS	STLC
Arsenic	31	1.0	"	"	"	"	"	"	STLC
Barium	400	1.0	"	"	"	"	"	"	STLC
Beryllium	1.2	1.0	"	"	"	"	04/27/18	"	STLC
Cadmium	ND	1.0	"	"	"	"	04/26/18	"	STLC
Chromium	47	1.0	"	"	"	"	"	"	STLC
Cobalt	32	1.0	"	"	"	"	"	"	STLC
Copper	55	1.0	"	"	"	"	"	"	STLC
Lead	46	2.0	"	"	"	"	"	"	STLC
Mercury	ND	0.10	"	"	"	"	"	"	STLC
Molybdenum	7.3	1.0	"	"	"	"	"	"	STLC
Nickel	47	1.0	"	"	"	"	"	"	STLC
Selenium	ND	5.0	"	"	"	"	"	"	STLC
Silver	ND	1.0	"	"	"	"	"	"	STLC
Thallium	ND	1.0	"	"	"	"	"	"	STLC
Vanadium	200	1.0	"	"	"	"	"	"	STLC
Zinc	220	1.0	"	"	"	"	"	"	STLC

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Blank (8042318-BLK1)

Prepared: 04/23/18 Analyzed: 04/26/18

Antimony	ND	1.0	ug/l
Arsenic	ND	1.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	1.0	"
Chromium	ND	1.0	"
Cobalt	ND	1.0	"
Copper	ND	1.0	"
Lead	ND	2.0	"
Mercury	ND	0.10	"
Molybdenum	ND	1.0	"
Nickel	ND	1.0	"
Selenium	ND	5.0	"
Silver	ND	1.0	"
Thallium	ND	1.0	"
Vanadium	ND	1.0	"
Zinc	ND	1.0	"

LCS (8042318-BS1)

Prepared: 04/23/18 Analyzed: 04/26/18

Arsenic	57.4	1.0	ug/l	50.0	115	80-120
Barium	56.4	1.0	"	50.0	113	80-120
Cadmium	56.5	1.0	"	50.0	113	80-120
Chromium	55.6	1.0	"	50.0	111	80-120
Lead	56.2	2.0	"	50.0	112	80-120

Duplicate (8042318-DUP1)

Source: T181179-06

Prepared: 04/23/18 Analyzed: 04/26/18

Barium	548	5.0	ug/l	521	4.95	200
Chromium	54.5	5.0	"	53.8	1.28	200

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Lake Forest, California 92630
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

Metals by EPA 6020 Method - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8042318 - STLC Metals

Matrix Spike (8042318-MS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	72.5	1.0	ug/l	50.0	27.8	89.2	75-125			
Barium	611	1.0	"	50.0	521	178	75-125			QM-PS
Cadmium	46.0	1.0	"	50.0	0.966	90.1	75-125			
Chromium	121	1.0	"	50.0	53.8	134	75-125			QM-PS
Lead	168	2.0	"	50.0	112	112	75-125			

Matrix Spike Dup (8042318-MSD1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Arsenic	73.4	1.0	ug/l	50.0	27.8	91.2	75-125	1.36	20	
Barium	595	1.0	"	50.0	521	147	75-125	2.61	20	QM-PS
Cadmium	47.4	1.0	"	50.0	0.966	92.9	75-125	2.97	20	
Chromium	124	1.0	"	50.0	53.8	141	75-125	2.85	20	QM-PS
Lead	164	2.0	"	50.0	112	104	75-125	2.22	20	

Post Spike (8042318-PS1)		Source: T181179-06			Prepared: 04/23/18 Analyzed: 04/26/18					
Barium	287		ug/l	25.0	261	104	80-120			
Chromium	56.2		"	25.0	26.9	117	80-120			

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/27/18 11:08

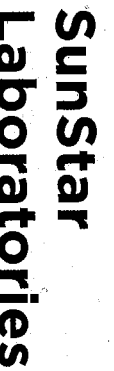
Notes and Definitions

STLC	Sample prepared in accordance with CCR Chapter 11, Article 5, Appendix II, Waste Extraction Test (WET).
QM-PS	The percent recovery and/or RPD are outside acceptance criteria. Results accepted based upon percent recovery results in the post spike and/or serial dilution.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

25712 Commerce Drive, Lake Forest, CA 92630
949-297-5020

Address: PO Box 332, Roseville, CA 95661

Phone: (916) 677-1470 Fax: (916) 677-1471

Project Manager: Joe Brusca

Page: 11 Of 11

Project Name: LAT1420P 4B-Ac 120P=RTT

Collector: Brusca Client Project #: 137-002

Batch #: 2 EDF #: 7181265

EDF #:

Sample disposal Instructions: Disposal @ \$2.00 each _____

Return to client _____

Pickup _____

Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>	4/13/18 1:30 PM	<i>[Signature]</i>	4/13/18 1330
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>		<i>[Signature]</i>	
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>		<i>[Signature]</i>	

Total # of containers	
Chain of Custody seals	Y/N/A
Seals intact?	Y/N/A
Received good condition/cold	
Turn around time:	5.3

Notes

Sample ID
SP5-1
SP5-2
SP5-3
SP5-4

Date Sampled 4/13/14

Time
10:00
10:04
10:06
10:09

Sample Type
Solid

Container Type
JAR

~~XXXX~~ 8260 SVOLs 8270

				8260 + OXY

				8260 BTEX, OXY only

~~XXXX~~ 0270 PCBs 8082

				8021 BTEX

				8015M (gasoline)

				8015M (diesel)

~~X~~	~~X~~	~~X~~	~~X~~	8015M Ext./Carbon Chain

~~X~~	~~X~~	~~X~~	~~X~~	6010/7000 Title 22 Metals

6020 ICP-MS Metals

					0020 101 -MS Metals
					~~LEAD AND ASSY~~

01	Laboratory ID #
02	
03	
04	

Comments/Preservative

— — — — Total # of containers

GOC 170169



May 2, 2018

FAL Project ID: 11472

Mr. Joe Brusca
Brusca Associates
7633 Stonewood Court
Granite Bay, CA 95746

Dear Mr. Brusca,

The following results are associated with Frontier Analytical Laboratory project **11472**. This corresponds to your **Lathrop 48-Ac Property** project under project number **137-002**. Four solid samples were received on 4/13/2018 in good condition. These samples were extracted and analyzed by EPA Method 1613 for tetra through octa chlorinated dibenzo dioxins and furans. The Toxic Equivalency (TEQ) for your samples has been calculated using the 2005 World Health Organization's (WHO's) toxic equivalency factors (TEFs). The total TEQ is reported on the upper right hand corner of each sample data sheet. Brusca Associates requested a turnaround time of fifteen business days for project **11472**.

The following report consists of an Analytical Data section and a Sample Receipt section. The Analytical Data section contains our sample tracking log and the analytical results. The Sample Receipt section contains your chain of custody, our sample login form and a sample photo. The attached results are specifically for the samples referenced in this report only. These results meet all NELAC requirements and shall not be reproduced except in full. Frontier Analytical Laboratory's State of Oregon NELAP certificate number is **4041**. Our State of California ELAP certificate number is **2934**. This report has been emailed to you as a portable document format (PDF) file. A hardcopy will not be sent to you unless specifically requested.

If you have any questions regarding project **11472**, please contact me at (916) 934-0900. Thank you for choosing Frontier Analytical Laboratory for your analytical testing needs.

Sincerely,

A handwritten signature in black ink, appearing to read "Bradley B. Silverbush".

Bradley B. Silverbush
Director of Operations

FRONTIER ANALYTICAL LABORATORY

5172 Hillsdale Circle * El Dorado Hills, CA 95762
Tel (916) 934-0900 * Fax (916) 934-0999
www.frontieranalytical.com

000001 of 000011

Frontier Analytical Laboratory

Sample Tracking Log

FAL Project ID: 11472

Received on: 04/13/2018

Project Due: 05/07/2018

Storage: R-3

FAL Sample ID	Dup	Client Project ID	Client Sample ID	Requested Method	Matrix	Sampling Date	Sampling Time	Hold Time Due Date
11472-001-SA	0	137-002	SP5-1	EPA 1613 D/F	Solid	04/13/2018	10:00 am	04/15/2019
11472-002-SA	0	137-002	SP5-2	EPA 1613 D/F	Solid	04/13/2018	10:04 am	04/15/2019
11472-003-SA	0	137-002	SP5-3	EPA 1613 D/F	Solid	04/13/2018	10:06 am	04/15/2019
11472-004-SA	0	137-002	SP5-4	EPA 1613 D/F	Solid	04/13/2018	10:09 am	04/15/2019

EPA Method 1613 PCDD/F



FAL ID: 11472-001-MB
Client ID: Method Blank
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.0
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.151		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.182		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.197		-	0.0793				
1,2,3,6,7,8-HxCDD	ND	0.221		-	0.0940	Total TCDD	ND	0.151	
1,2,3,7,8,9-HxCDD	ND	0.191		-	0.0823	Total PeCDD	ND	0.182	
1,2,3,4,6,7,8-HpCDD	ND	0.180		-	0.0842	Total HxCDD	ND	0.221	
OCDD	ND	0.314		-	0.172	Total HpCDD	ND	0.180	
2,3,7,8-TCDF	ND	0.116		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.140		-	0.0449				
2,3,4,7,8-PeCDF	ND	0.141		-	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.126		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.137		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.132		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.160		-	0.0657	Total TCDF	ND	0.116	
1,2,3,4,6,7,8-HpCDF	ND	0.122		-	0.0747	Total PeCDF	ND	0.141	
1,2,3,4,7,8,9-HpCDF	ND	0.211		-	0.0883	Total HxCDF	ND	0.160	
OCDF	ND	0.225		-	0.170	Total HpCDF	ND	0.211	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	104	25.0 - 164	
13C-1,2,3,7,8-PeCDD	107	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	110	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	107	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	105	23.0 - 140	
13C-OCDD	101	17.0 - 157	
13C-2,3,7,8-TCDF	111	24.0 - 169	
13C-1,2,3,7,8-PeCDF	104	24.0 - 185	
13C-2,3,4,7,8-PeCDF	108	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	99.4	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	98.5	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	99.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	97.8	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	98.8	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	75.9	26.0 - 138	
13C-OCDF	97.2	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 94.0 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000003 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11472-001-OPR
Client ID: OPR
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: NA
Amount: 5.00 g

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: ng/ml

Acquired: 05-01-2018
2005 WHO TEQ: NA

Compound	Conc	QC Limits	Qual
2,3,7,8-TCDD	10.5	6.70 - 15.8	
1,2,3,7,8-PeCDD	50.2	35.0 - 71.0	
1,2,3,4,7,8-HxCDD	50.0	35.0 - 82.0	
1,2,3,6,7,8-HxCDD	49.6	38.0 - 67.0	
1,2,3,7,8,9-HxCDD	47.4	32.0 - 81.0	
1,2,3,4,6,7,8-HpCDD	51.5	35.0 - 70.0	
OCDD	99.1	78.0 - 144	
2,3,7,8-TCDF	11.3	7.50 - 15.8	
1,2,3,7,8-PeCDF	50.8	40.0 - 67.0	
2,3,4,7,8-PeCDF	50.8	34.0 - 80.0	
1,2,3,4,7,8-HxCDF	51.2	36.0 - 67.0	
1,2,3,6,7,8-HxCDF	52.2	42.0 - 65.0	
2,3,4,6,7,8-HxCDF	51.5	35.0 - 78.0	
1,2,3,7,8,9-HxCDF	51.8	39.0 - 65.0	
1,2,3,4,6,7,8-HpCDF	51.6	41.0 - 61.0	
1,2,3,4,7,8,9-HpCDF	52.6	39.0 - 69.0	
OCDF	101	63.0 - 170	
Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	76.9	20.0 - 175	
13C-1,2,3,7,8-PeCDD	84.4	21.0 - 227	
13C-1,2,3,4,7,8-HxCDD	97.8	21.0 - 193	
13C-1,2,3,6,7,8-HxCDD	96.9	25.0 - 163	
13C-1,2,3,4,6,7,8-HpCDD	87.3	26.0 - 166	
13C-OCDD	84.1	13.0 - 198	
13C-2,3,7,8-TCDF	80.3	22.0 - 152	
13C-1,2,3,7,8-PeCDF	88.2	21.0 - 192	
13C-2,3,4,7,8-PeCDF	90.0	13.0 - 328	
13C-1,2,3,4,7,8-HxCDF	88.7	19.0 - 202	
13C-1,2,3,6,7,8-HxCDF	86.6	21.0 - 159	
13C-2,3,4,6,7,8-HxCDF	86.1	22.0 - 176	
13C-1,2,3,7,8,9-HxCDF	74.7	17.0 - 205	
13C-1,2,3,4,6,7,8-HpCDF	81.5	21.0 - 158	
13C-1,2,3,4,7,8,9-HpCDF	78.0	20.0 - 186	
13C-OCDF	79.1	13.0 - 198	
Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	72.0	31.0 - 191	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 5/2/2018

Reviewed By: 
Date: 5/2/2018

000004 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11472-001-SA
Client ID: SP5-1
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.01 g
% Solids: 88.35

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.759
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.226		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.227		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.213		-	0.0793				
1,2,3,6,7,8-HxCDD	0.953	-	J	0.0953	0.0940	Total TCDD	ND	0.226	
1,2,3,7,8,9-HxCDD	0.696	-	J	0.0696	0.0823	Total PeCDD	0.602	-	J
1,2,3,4,6,7,8-HpCDD	16.1	-		0.161	0.0842	Total HxCDD	7.07	-	
OCDD	104	-		0.0312	0.172	Total HpCDD	29.2	-	
2,3,7,8-TCDF	0.815	-	J	0.0815	0.0269				
1,2,3,7,8-PeCDF	ND	0.178		-	0.0449				
2,3,4,7,8-PeCDF	0.547	-	J	0.164	0.0468				
1,2,3,4,7,8-HxCDF	0.507	-	J	0.0507	0.0437				
1,2,3,6,7,8-HxCDF	0.379	-	J	0.0379	0.0417				
2,3,4,6,7,8-HxCDF	0.336	-	J	0.0336	0.0574				
1,2,3,7,8,9-HpCDF	ND	0.219		-	0.0657	Total TCDF	3.42	-	
1,2,3,4,6,7,8-HpCDF	3.24	-	J	0.0324	0.0747	Total PeCDF	3.71	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.262		-	0.0883	Total HxCDF	5.28	-	
OCDF	6.45	-	J	0.00194	0.170	Total HpCDF	8.39	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	78.5	25.0 - 164	
13C-1,2,3,7,8-PeCDD	88.5	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	109	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	105	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	89.2	23.0 - 140	
13C-OCDD	90.8	17.0 - 157	
13C-2,3,7,8-TCDF	79.1	24.0 - 169	
13C-1,2,3,7,8-PeCDF	92.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	95.4	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	99.2	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	95.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	93.3	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	82.9	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	78.5	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	82.2	26.0 - 138	
13C-OCDF	84.2	17.0 - 157	

Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	72.4	35.0 - 197	

A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1

B Analyte is present in Method Blank

C Chemical Interference

D Presence of Diphenyl Ethers

DNQ Analyte concentration is below calibration range

E Analyte concentration is above calibration range

F Analyte confirmation on secondary column

J Analyte concentration is below calibration range

M Maximum possible concentration

ND Analyte Not Detected at Detection Limit Level

NP Not Provided

P Pre-filtered through a Whatman 0.7um GF/F filter

S Sample acceptance criteria not met

X Matrix interferences

* Result taken from dilution or reinjection

Analyst:

Date: 5/2/2018

Reviewed By:

Date: 5/2/2018

000005 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11472-002-SA
Client ID: SP5-2
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.03 g
% Solids: 92.60

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 0.670
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.189		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.287		-	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.213		-	0.0793				
1,2,3,6,7,8-HxCDD	0.736	-	J	0.0736	0.0940	Total TCDD	ND	0.189	
1,2,3,7,8,9-HxCDD	0.589	-	J	0.0589	0.0823	Total PeCDD	ND	0.287	
1,2,3,4,6,7,8-HpCDD	10.9	-		0.109	0.0842	Total HxCDD	5.21	-	
OCDD	73.3	-		0.0220	0.172	Total HpCDD	20.4	-	
2,3,7,8-TCDF	0.829	-	J	0.0829	0.0269				
1,2,3,7,8-PeCDF	0.269	-	J	0.00807	0.0449				
2,3,4,7,8-PeCDF	0.478	-	J	0.143	0.0468				
1,2,3,4,7,8-HxCDF	0.572	-	J	0.0572	0.0437				
1,2,3,6,7,8-HxCDF	0.459	-	J	0.0459	0.0417				
2,3,4,6,7,8-HxCDF	0.407	-	J	0.0407	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.159		-	0.0657	Total TCDF	4.70	-	
1,2,3,4,6,7,8-HpCDF	2.66	-	J	0.0266	0.0747	Total PeCDF	4.34	-	J
1,2,3,4,7,8,9-HpCDF	ND	0.246		-	0.0883	Total HxCDF	4.67	-	J
OCDF	5.95	-	J	0.00179	0.170	Total HpCDF	6.39	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	85.9	25.0 - 164	
13C-1,2,3,7,8-PeCDD	93.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	112	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	108	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	98.8	23.0 - 140	
13C-OCDD	89.6	17.0 - 157	
13C-2,3,7,8-TCDF	82.8	24.0 - 169	
13C-1,2,3,7,8-PeCDF	94.7	24.0 - 185	
13C-2,3,4,7,8-PeCDF	94.6	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	103	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	99.3	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	98.4	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	88.7	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	80.9	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	85.6	26.0 - 138	
13C-OCDF	84.9	17.0 - 157	

Cleanup Surrogate			
37Cl-2,3,7,8-TCDD	76.2	35.0 - 197	

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 
Date: 5/2/2018

Reviewed By: 
Date: 5/2/2018

000006 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11472-003-SA
Client ID: SP5-3
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.03 g
% Solids: 90.60

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-01-2018
2005 WHO TEQ: 1.24
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.204		-	0.0273				
1,2,3,7,8-PeCDD	0.340	-	J	0.340	0.0570				
1,2,3,4,7,8-HxCDD	ND	0.279		-	0.0793				
1,2,3,6,7,8-HxCDD	0.910	-	J	0.0910	0.0940	Total TCDD	0.472	-	J
1,2,3,7,8,9-HxCDD	0.499	-	J	0.0499	0.0823	Total PeCDD	2.65	-	J
1,2,3,4,6,7,8-HpCDD	12.7	-		0.127	0.0842	Total HxCDD	6.81	-	
OCDD	82.4	-		0.0247	0.172	Total HpCDD	24.6	-	
2,3,7,8-TCDF	1.04	-		0.104	0.0269				
1,2,3,7,8-PeCDF	0.272	-	J	0.00816	0.0449				
2,3,4,7,8-PeCDF	0.747	-	J	0.224	0.0468				
1,2,3,4,7,8-HxCDF	0.920	-	J	0.0920	0.0437				
1,2,3,6,7,8-HxCDF	0.565	-	J	0.0565	0.0417				
2,3,4,6,7,8-HxCDF	0.484	-	J	0.0484	0.0574				
1,2,3,7,8,9-HxCDF	0.372	-	J	0.0372	0.0657	Total TCDF	5.58	-	
1,2,3,4,6,7,8-HpCDF	2.91	-	J	0.0291	0.0747	Total PeCDF	5.24	-	M
1,2,3,4,7,8,9-HpCDF	0.556	-	J	0.00556	0.0883	Total HxCDF	6.19	-	
OCDF	5.75	-	J	0.00173	0.170	Total HpCDF	7.89	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	81.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	90.1	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	112	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	110	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	97.8	23.0 - 140	
13C-OCDD	91.7	17.0 - 157	
13C-2,3,7,8-TCDF	80.4	24.0 - 169	
13C-1,2,3,7,8-PeCDF	95.3	24.0 - 185	
13C-2,3,4,7,8-PeCDF	98.1	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	102	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	99.8	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	96.1	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	88.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	84.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	83.3	26.0 - 138	
13C-OCDF	88.4	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 73.6 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000007 of 000011

EPA Method 1613 PCDD/F



FAL ID: 11472-004-SA
Client ID: SP5-4
Matrix: Solid
Batch No: X4493

Date Extracted: 04-30-2018
Date Received: 04-13-2018
Amount: 5.01 g
% Solids: 91.50

ICal: PCDDFAL4-12-20-17
GC Column: DB5MS
Units: pg/g

Acquired: 05-02-2018
2005 WHO TEQ: 0.539
Basis: Dry Weight

Compound	Conc	DL	Qual	2005 WHO Tox	MDL	Compound	Conc	DL	Qual
2,3,7,8-TCDD	ND	0.193		-	0.0273				
1,2,3,7,8-PeCDD	ND	0.231		-	0.0570				
1,2,3,4,7,8-HxCDD	0.441	-	J	0.0441	0.0793				
1,2,3,6,7,8-HxCDD	0.971	-	J	0.0971	0.0940	Total TCDD	ND	0.193	
1,2,3,7,8,9-HxCDD	0.806	-	J	0.0806	0.0823	Total PeCDD	2.34	-	J
1,2,3,4,6,7,8-HpCDD	15.6	-		0.156	0.0842	Total HxCDD	10.1	-	
OCDD	90.0	-		0.0270	0.172	Total HpCDD	30.6	-	
2,3,7,8-TCDF	ND	0.226		-	0.0269				
1,2,3,7,8-PeCDF	ND	0.191		-	0.0449				
2,3,4,7,8-PeCDF	0.324	-	J	0.0972	0.0468				
1,2,3,4,7,8-HxCDF	ND	0.289		-	0.0437				
1,2,3,6,7,8-HxCDF	ND	0.266		-	0.0417				
2,3,4,6,7,8-HxCDF	ND	0.283		-	0.0574				
1,2,3,7,8,9-HxCDF	ND	0.344		-	0.0657	Total TCDF	0.380	-	J
1,2,3,4,6,7,8-HpCDF	3.07	-	J	0.0307	0.0747	Total PeCDF	1.62	-	J
1,2,3,4,7,8,9-HpCDF	0.399	-	J	0.00399	0.0883	Total HxCDF	3.21	-	J
OCDF	6.79	-	J	0.00204	0.170	Total HpCDF	9.66	-	

Internal Standards	% Rec	QC Limits	Qual
13C-2,3,7,8-TCDD	81.6	25.0 - 164	
13C-1,2,3,7,8-PeCDD	89.8	25.0 - 181	
13C-1,2,3,4,7,8-HxCDD	110	32.0 - 141	
13C-1,2,3,6,7,8-HxCDD	108	28.0 - 130	
13C-1,2,3,4,6,7,8-HpCDD	95.1	23.0 - 140	
13C-OCDD	90.6	17.0 - 157	
13C-2,3,7,8-TCDF	79.4	24.0 - 169	
13C-1,2,3,7,8-PeCDF	95.8	24.0 - 185	
13C-2,3,4,7,8-PeCDF	95.5	21.0 - 178	
13C-1,2,3,4,7,8-HxCDF	97.0	26.0 - 152	
13C-1,2,3,6,7,8-HxCDF	94.2	26.0 - 123	
13C-2,3,4,6,7,8-HxCDF	94.8	28.0 - 136	
13C-1,2,3,7,8,9-HxCDF	85.5	29.0 - 147	
13C-1,2,3,4,6,7,8-HpCDF	79.6	28.0 - 143	
13C-1,2,3,4,7,8,9-HpCDF	80.0	26.0 - 138	
13C-OCDF	87.1	17.0 - 157	

Cleanup Surrogate

37Cl-2,3,7,8-TCDD 73.4 35.0 - 197

- A Isotopic Labeled Standard outside QC range but signal to noise ratio is >10:1
- B Analyte is present in Method Blank
- C Chemical Interference
- D Presence of Diphenyl Ethers
- DNQ Analyte concentration is below calibration range
- E Analyte concentration is above calibration range
- F Analyte confirmation on secondary column
- J Analyte concentration is below calibration range
- M Maximum possible concentration
- ND Analyte Not Detected at Detection Limit Level
- NP Not Provided
- P Pre-filtered through a Whatman 0.7um GF/F filter
- S Sample acceptance criteria not met
- X Matrix interferences
- * Result taken from dilution or reinjection

Analyst: 

Date: 5/2/2018

Reviewed By: 

Date: 5/2/2018

000008 of 000011

Client understands that all terms described in the proposals, quotations, and/or the general terms provided in the current FAL price schedules will be followed. FAL reserves the rights to terminate its service or withhold delivery of reports, if in FAL's sole discretion the terms of the project have been broken.

Frontier Analytical Laboratory

Sample Login Form

FAL Project ID: **11472**

Client:	Brusca Associates, Inc.
Client Project ID:	137-002
Date Received:	04/13/2018
Time Received:	02:12 pm
Received By:	KZ
Logged In By:	KZ
# of Samples Received:	4
Duplicates:	0
Storage Location:	R-3

Method of Delivery:	Courier
Tracking Number:	NA
Shipping Container Received Intact	Yes
Custody seals(s) present?	No
Custody seals(s) intact?	No
Sample Arrival Temperature (C)	0
Cooling Method	Ice
Chain Of Custody Present?	Yes
Return Shipping Container To Client	Yes
Test aqueous sample for residual Chlorine	No
Sodium Thiosulfate Added	No
Adequate Sample Volume	Yes
Appropriate Sample Container	Yes
pH Range of Aqueous Sample	N/A
Anomalies or additional comments:	



25712 Commercentre Drive
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949.297.5027 Fax

23 April 2018

Joe Brusca
Brusca Associates Inc.
PO Box 332
Roseville, CA 95661
RE: Lathrop 48-Ac Property

Enclosed are the results of analyses for samples received by the laboratory on 04/14/18 08:10. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Mike Jaroudi
Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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949.297.5027 Fax

Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
SP5-1	T181265-01	Soil	04/13/18 10:00	04/14/18 08:10
SP5-2	T181265-02	Soil	04/13/18 10:04	04/14/18 08:10
SP5-3	T181265-03	Soil	04/13/18 10:06	04/14/18 08:10
SP5-4	T181265-04	Soil	04/13/18 10:09	04/14/18 08:10

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

DETECTIONS SUMMARY

Sample ID: SP5-1

Laboratory ID: T181265-01

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	53	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	61	10	mg/kg	EPA 8015B	
Barium	59	1.0	mg/kg	EPA 6010B	
Chromium	8.6	2.0	mg/kg	EPA 6010B	
Cobalt	4.3	2.0	mg/kg	EPA 6010B	
Copper	4.4	1.0	mg/kg	EPA 6010B	
Nickel	6.2	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	33	1.0	mg/kg	EPA 6010B	
PCB-1254	17	10	ug/kg	EPA 8082	

Sample ID: SP5-2

Laboratory ID: T181265-02

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	70	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	58	10	mg/kg	EPA 8015B	
Barium	63	1.0	mg/kg	EPA 6010B	
Chromium	6.4	2.0	mg/kg	EPA 6010B	
Cobalt	5.3	2.0	mg/kg	EPA 6010B	
Copper	4.9	1.0	mg/kg	EPA 6010B	
Nickel	7.2	2.0	mg/kg	EPA 6010B	
Vanadium	20	5.0	mg/kg	EPA 6010B	
Zinc	30	1.0	mg/kg	EPA 6010B	
PCB-1254	14	10	ug/kg	EPA 8082	

Sample ID: SP5-3

Laboratory ID: T181265-03

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
C13-C28 (DRO)	35	10	mg/kg	EPA 8015B	
C29-C40 (MORO)	50	10	mg/kg	EPA 8015B	
Barium	61	1.0	mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

Sample ID: SP5-3

Laboratory ID: T181265-03

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Chromium	7.0	2.0		mg/kg	EPA 6010B	
Cobalt	4.5	2.0		mg/kg	EPA 6010B	
Copper	8.6	1.0		mg/kg	EPA 6010B	
Nickel	6.4	2.0		mg/kg	EPA 6010B	
Vanadium	18	5.0		mg/kg	EPA 6010B	
Zinc	30	1.0		mg/kg	EPA 6010B	
PCB-1254	35	10		ug/kg	EPA 8082	

Sample ID: SP5-4

Laboratory ID: T181265-04

Analyte	Result	Reporting		Units	Method	Notes
		Limit				
C13-C28 (DRO)	17	10		mg/kg	EPA 8015B	
C29-C40 (MORO)	56	10		mg/kg	EPA 8015B	
Barium	64	1.0		mg/kg	EPA 6010B	
Chromium	6.7	2.0		mg/kg	EPA 6010B	
Cobalt	4.9	2.0		mg/kg	EPA 6010B	
Copper	5.2	1.0		mg/kg	EPA 6010B	
Nickel	7.2	2.0		mg/kg	EPA 6010B	
Vanadium	20	5.0		mg/kg	EPA 6010B	
Zinc	31	1.0		mg/kg	EPA 6010B	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-1
T181265-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B
C13-C28 (DRO)	53	10	"	"	"	"	"	"
C29-C40 (MORO)	61	10	"	"	"	"	"	"
Surrogate: <i>p</i> -Terphenyl		107 %	65-135		"	"	"	"

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B
Silver	ND	2.0	"	"	"	"	"	"
Arsenic	ND	5.0	"	"	"	"	"	"
Barium	59	1.0	"	"	"	"	"	"
Beryllium	ND	1.0	"	"	"	"	"	"
Cadmium	ND	2.0	"	"	"	"	"	"
Chromium	8.6	2.0	"	"	"	"	"	"
Cobalt	4.3	2.0	"	"	"	"	"	"
Copper	4.4	1.0	"	"	"	"	"	"
Lead	ND	3.0	"	"	"	"	"	"
Molybdenum	ND	5.0	"	"	"	"	"	"
Nickel	6.2	2.0	"	"	"	"	"	"
Selenium	ND	5.0	"	"	"	"	"	"
Thallium	ND	2.0	"	"	"	"	"	"
Vanadium	20	5.0	"	"	"	"	"	"
Zinc	33	1.0	"	"	"	"	"	"

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-1

T181265-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	17	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		84.8 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		84.4 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-1

T181265-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-1

T181265-01 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		43.4 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		54.5 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		56.1 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		65.3 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		90.2 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		85.5 %	29.1-130		"	"	"	"	

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-2

T181265-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	70	10	"	"	"	"	"	"	
C29-C40 (MORO)	58	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	63	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.4	2.0	"	"	"	"	"	"	
Cobalt	5.3	2.0	"	"	"	"	"	"	
Copper	4.9	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-2

T181265-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	14	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		86.2 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		85.8 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Mike Jaroudi, Project Manager



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Brusca Associates Inc.
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Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-2

T181265-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-2

T181265-02 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
2-Methylphenol	ND	1000	"	"	"	"	"	"	
4-Methylphenol	ND	1000	"	"	"	"	"	"	
Naphthalene	ND	300	"	"	"	"	"	"	
2-Nitroaniline	ND	300	"	"	"	"	"	"	
3-Nitroaniline	ND	300	"	"	"	"	"	"	
4-Nitroaniline	ND	300	"	"	"	"	"	"	
Nitrobenzene	ND	1000	"	"	"	"	"	"	
2-Nitrophenol	ND	1000	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"	
Phenanthrene	ND	300	"	"	"	"	"	"	
Azobenzene	ND	300	"	"	"	"	"	"	
Pyridine	ND	300	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		49.5 %	15-121		"	"	"	"	
Surrogate: Phenol-d6		60.8 %	24-113		"	"	"	"	
Surrogate: Nitrobenzene-d5		60.2 %	21.3-119		"	"	"	"	
Surrogate: 2-Fluorobiphenyl		71.2 %	32.4-102		"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		94.0 %	18.1-105		"	"	"	"	
Surrogate: Terphenyl-d14		84.5 %	29.1-130		"	"	"	"	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-3

T181265-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	35	10	"	"	"	"	"	"	
C29-C40 (MORO)	50	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		106 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	61	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	7.0	2.0	"	"	"	"	"	"	
Cobalt	4.5	2.0	"	"	"	"	"	"	
Copper	8.6	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	6.4	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	18	5.0	"	"	"	"	"	"	
Zinc	30	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

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04/23/18 10:35

SP5-3

T181265-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	35	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
<i>Surrogate: Tetrachloro-meta-xylene</i>		83.3 %	35-140		"	"	"	"	
<i>Surrogate: Decachlorobiphenyl</i>		81.8 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-3

T181265-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-3

T181265-03 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	56.9 %	15-121			"	"	"	"
Surrogate: Phenol-d6	68.6 %	24-113			"	"	"	"
Surrogate: Nitrobenzene-d5	68.0 %	21.3-119			"	"	"	"
Surrogate: 2-Fluorobiphenyl	83.3 %	32.4-102			"	"	"	"
Surrogate: 2,4,6-Tribromophenol	99.1 %	18.1-105			"	"	"	"
Surrogate: Terphenyl-d14	88.9 %	29.1-130			"	"	"	"

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-4

T181265-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup

C6-C12 (GRO)	ND	10	mg/kg	1	8041214	04/18/18	04/18/18	EPA 8015B	
C13-C28 (DRO)	17	10	"	"	"	"	"	"	
C29-C40 (MORO)	56	10	"	"	"	"	"	"	
Surrogate: <i>p</i> -Terphenyl		105 %	65-135		"	"	"	"	

Metals by EPA 6010B

Antimony	ND	3.0	mg/kg	1	8041623	04/16/18	04/16/18	EPA 6010B	
Silver	ND	2.0	"	"	"	"	"	"	
Arsenic	ND	5.0	"	"	"	"	"	"	
Barium	64	1.0	"	"	"	"	"	"	
Beryllium	ND	1.0	"	"	"	"	"	"	
Cadmium	ND	2.0	"	"	"	"	"	"	
Chromium	6.7	2.0	"	"	"	"	"	"	
Cobalt	4.9	2.0	"	"	"	"	"	"	
Copper	5.2	1.0	"	"	"	"	"	"	
Lead	ND	3.0	"	"	"	"	"	"	
Molybdenum	ND	5.0	"	"	"	"	"	"	
Nickel	7.2	2.0	"	"	"	"	"	"	
Selenium	ND	5.0	"	"	"	"	"	"	
Thallium	ND	2.0	"	"	"	"	"	"	
Vanadium	20	5.0	"	"	"	"	"	"	
Zinc	31	1.0	"	"	"	"	"	"	

Cold Vapor Extraction EPA 7470/7471

Mercury	ND	0.10	mg/kg	1	8041624	04/16/18	04/16/18	EPA 7471A Soil	

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Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-4

T181265-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Polychlorinated Biphenyls by EPA Method 8082

PCB-1016	ND	10	ug/kg	1	8041633	04/16/18	04/18/18	EPA 8082	
PCB-1221	ND	10	"	"	"	"	"	"	
PCB-1232	ND	10	"	"	"	"	"	"	
PCB-1242	ND	10	"	"	"	"	"	"	
PCB-1248	ND	10	"	"	"	"	"	"	
PCB-1254	ND	10	"	"	"	"	"	"	
PCB-1260	ND	10	"	"	"	"	"	"	
Surrogate: Tetrachloro-meta-xylene		94.1 %	35-140		"	"	"	"	
Surrogate: Decachlorobiphenyl		95.9 %	35-140		"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Phenol	ND	1000	"	"	"	"	"	"	
Aniline	ND	300	"	"	"	"	"	"	
2-Chlorophenol	ND	1000	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	300	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	300	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	300	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	1000	"	"	"	"	"	"	
2-Methylnaphthalene	ND	300	"	"	"	"	"	"	
1-Methylnaphthalene	ND	300	"	"	"	"	"	"	
Acenaphthene	ND	300	"	"	"	"	"	"	
4-Nitrophenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	300	"	"	"	"	"	"	
Pentachlorophenol	ND	1000	"	"	"	"	"	"	
Pyrene	ND	300	"	"	"	"	"	"	
Acenaphthylene	ND	300	"	"	"	"	"	"	
Anthracene	ND	300	"	"	"	"	"	"	
Benzo (a) anthracene	ND	300	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	300	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	1000	"	"	"	"	"	"	

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Mike Jaroudi, Project Manager



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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-4

T181265-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzo (a) pyrene	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C	
Benzyl alcohol	ND	300	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	300	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	300	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	300	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	300	"	"	"	"	"	"	
4-Chloroaniline	ND	300	"	"	"	"	"	"	
2-Chloronaphthalene	ND	300	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	300	"	"	"	"	"	"	
Chrysene	ND	300	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	300	"	"	"	"	"	"	
Dibenzofuran	ND	300	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	300	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	300	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	300	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	1000	"	"	"	"	"	"	
Diethyl phthalate	ND	300	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	1000	"	"	"	"	"	"	
Dimethyl phthalate	ND	300	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	1000	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	1000	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	1000	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	300	"	"	"	"	"	"	
Fluoranthene	ND	300	"	"	"	"	"	"	
Fluorene	ND	300	"	"	"	"	"	"	
Hexachlorobenzene	ND	1500	"	"	"	"	"	"	
Hexachlorobutadiene	ND	300	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	1000	"	"	"	"	"	"	
Hexachloroethane	ND	300	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	300	"	"	"	"	"	"	

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

SP5-4

T181265-04 (Soil)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes

SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	300	ug/kg	1	8041634	04/16/18	04/17/18	EPA 8270C
2-Methylphenol	ND	1000	"	"	"	"	"	"
4-Methylphenol	ND	1000	"	"	"	"	"	"
Naphthalene	ND	300	"	"	"	"	"	"
2-Nitroaniline	ND	300	"	"	"	"	"	"
3-Nitroaniline	ND	300	"	"	"	"	"	"
4-Nitroaniline	ND	300	"	"	"	"	"	"
Nitrobenzene	ND	1000	"	"	"	"	"	"
2-Nitrophenol	ND	1000	"	"	"	"	"	"
N-Nitrosodimethylamine	ND	300	"	"	"	"	"	"
N-Nitrosodiphenylamine	ND	300	"	"	"	"	"	"
2,3,5,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
2,3,4,6-Tetrachlorophenol	ND	300	"	"	"	"	"	"
Phenanthrene	ND	300	"	"	"	"	"	"
Azobenzene	ND	300	"	"	"	"	"	"
Pyridine	ND	300	"	"	"	"	"	"
2,4,5-Trichlorophenol	ND	1000	"	"	"	"	"	"
2,4,6-Trichlorophenol	ND	1000	"	"	"	"	"	"
Surrogate: 2-Fluorophenol	57.5 %	15-121	"	"	"	"	"	"
Surrogate: Phenol-d6	65.7 %	24-113	"	"	"	"	"	"
Surrogate: Nitrobenzene-d5	65.1 %	21.3-119	"	"	"	"	"	"
Surrogate: 2-Fluorobiphenyl	75.3 %	32.4-102	"	"	"	"	"	"
Surrogate: 2,4,6-Tribromophenol	95.4 %	18.1-105	"	"	"	"	"	"
Surrogate: Terphenyl-d14	86.6 %	29.1-130	"	"	"	"	"	"

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Reported:
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Extractable Petroleum Hydrocarbons by 8015B with Silica Gel Cleanup - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041214 - EPA 3550B GC

Blank (8041214-BLK1)

Prepared: 04/12/18 Analyzed: 04/18/18

C6-C12 (GRO)	ND	10	mg/kg							
C13-C28 (DRO)	ND	10	"							
C29-C40 (MORO)	ND	10	"							

Surrogate: <i>p</i> -Terphenyl	97.8		"	99.0		98.8	65-135			

LCS (8041214-BS1)

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	500	10	mg/kg	495		101	75-125			
Surrogate: <i>p</i> -Terphenyl	101		"	99.0		102	65-135			

Matrix Spike (8041214-MS1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	560	10	mg/kg	505	53	100	75-125			
Surrogate: <i>p</i> -Terphenyl	106		"	101		105	65-135			

Matrix Spike Dup (8041214-MSD1)

Source: T181265-01

Prepared: 04/12/18 Analyzed: 04/18/18

C13-C28 (DRO)	590	10	mg/kg	500	53	108	75-125	5.92	20	
Surrogate: <i>p</i> -Terphenyl	104		"	100		104	65-135			

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Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041623 - EPA 3050B

Blank (8041623-BLK1)

Prepared & Analyzed: 04/16/18

Antimony	ND	3.0	mg/kg
Silver	ND	2.0	"
Arsenic	ND	5.0	"
Barium	ND	1.0	"
Beryllium	ND	1.0	"
Cadmium	ND	2.0	"
Chromium	ND	2.0	"
Cobalt	ND	2.0	"
Copper	ND	1.0	"
Lead	ND	3.0	"
Molybdenum	ND	5.0	"
Nickel	ND	2.0	"
Selenium	ND	5.0	"
Thallium	ND	2.0	"
Vanadium	ND	5.0	"
Zinc	ND	1.0	"

LCS (8041623-BS1)

Prepared & Analyzed: 04/16/18

Arsenic	92.9	5.0	mg/kg	100	92.9	75-125
Barium	91.6	1.0	"	100	91.6	75-125
Cadmium	88.7	2.0	"	100	88.7	75-125
Chromium	90.3	2.0	"	100	90.3	75-125
Lead	96.9	3.0	"	100	96.9	75-125

Matrix Spike (8041623-MS1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Arsenic	88.5	5.0	mg/kg	100	1.81	86.7	75-125
Barium	119	1.0	"	100	37.7	81.1	75-125
Cadmium	83.2	2.0	"	100	0.425	82.8	75-125
Chromium	89.3	2.0	"	100	5.37	84.0	75-125
Lead	93.3	3.0	"	100	ND	93.3	75-125

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Project Manager: Joe Brusca

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Metals by EPA 6010B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041623 - EPA 3050B

Matrix Spike Dup (8041623-MSD1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Arsenic	95.2	5.0	mg/kg	100	1.81	93.4	75-125	7.37	20	
Barium	123	1.0	"	100	37.7	85.1	75-125	3.25	20	
Cadmium	85.4	2.0	"	100	0.425	85.0	75-125	2.67	20	
Chromium	91.3	2.0	"	100	5.37	86.0	75-125	2.21	20	
Lead	99.2	3.0	"	100	ND	99.2	75-125	6.14	20	

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04/23/18 10:35

Cold Vapor Extraction EPA 7470/7471 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041624 - EPA 7471A Soil

Blank (8041624-BLK1)

Prepared & Analyzed: 04/16/18

Mercury	ND	0.10	mg/kg

LCS (8041624-BS1)

Prepared & Analyzed: 04/16/18

Mercury	0.405	0.10	mg/kg	0.417		97.2	80-120

Matrix Spike (8041624-MS1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Mercury	0.408	0.10	mg/kg	0.417	ND	98.0	75-125

Matrix Spike Dup (8041624-MSD1)

Source: T181256-01

Prepared & Analyzed: 04/16/18

Mercury	0.428	0.10	mg/kg	0.417	ND	103	75-125	4.64	20

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04/23/18 10:35

Polychlorinated Biphenyls by EPA Method 8082 - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041633 - EPA 3550 ECD/GCMS

Blank (8041633-BLK1)

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	ND	10	ug/kg							
PCB-1221	ND	10	"							
PCB-1232	ND	10	"							
PCB-1242	ND	10	"							
PCB-1248	ND	10	"							
PCB-1254	ND	10	"							
PCB-1260	ND	10	"							
Surrogate: Tetrachloro-meta-xylene	8.65		"	9.90		87.3	35-140			
Surrogate: Decachlorobiphenyl	10.1		"	9.90		102	35-140			

LCS (8041633-BS1)

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	70.1	10	ug/kg	99.0		70.8	40-130			
PCB-1260	66.8	10	"	99.0		67.5	40-130			
Surrogate: Tetrachloro-meta-xylene	9.51		"	9.90		96.0	35-140			
Surrogate: Decachlorobiphenyl	11.0		"	9.90		112	35-140			

Matrix Spike (8041633-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	60.2	10	ug/kg	101	ND	59.6	40-130			
PCB-1260	60.1	10	"	101	ND	59.5	40-130			
Surrogate: Tetrachloro-meta-xylene	8.99		"	10.1		89.0	35-140			
Surrogate: Decachlorobiphenyl	9.25		"	10.1		91.5	35-140			

Matrix Spike Dup (8041633-MSD1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/18/18

PCB-1016	64.2	10	ug/kg	101	ND	63.6	40-130	6.49	30	
PCB-1260	62.7	10	"	101	ND	62.1	40-130	4.28	30	
Surrogate: Tetrachloro-meta-xylene	9.14		"	10.1		90.5	35-140			
Surrogate: Decachlorobiphenyl	9.02		"	10.1		89.3	35-140			

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Project Manager: Joe Brusca

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Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

Carbazole	ND	300	ug/kg
Phenol	ND	1000	"
Aniline	ND	300	"
2-Chlorophenol	ND	1000	"
1,4-Dichlorobenzene	ND	300	"
N-Nitrosodi-n-propylamine	ND	300	"
1,2,4-Trichlorobenzene	ND	300	"
4-Chloro-3-methylphenol	ND	1000	"
1-Methylnaphthalene	ND	300	"
2-Methylnaphthalene	ND	300	"
Acenaphthene	ND	300	"
4-Nitrophenol	ND	1000	"
2,4-Dinitrotoluene	ND	300	"
Pentachlorophenol	ND	1000	"
Pyrene	ND	300	"
Acenaphthylene	ND	300	"
Anthracene	ND	300	"
Benzo (a) anthracene	ND	300	"
Benzo (b) fluoranthene	ND	300	"
Benzo (k) fluoranthene	ND	300	"
Benzo (g,h,i) perylene	ND	1000	"
Benzo (a) pyrene	ND	300	"
Benzyl alcohol	ND	300	"
Bis(2-chloroethoxy)methane	ND	300	"
Bis(2-chloroethyl)ether	ND	300	"
Bis(2-chloroisopropyl)ether	ND	300	"
Bis(2-ethylhexyl)phthalate	ND	300	"
4-Bromophenyl phenyl ether	ND	300	"
Butyl benzyl phthalate	ND	300	"
4-Chloroaniline	ND	300	"
2-Chloronaphthalene	ND	300	"
4-Chlorophenyl phenyl ether	ND	300	"
Chrysene	ND	300	"
Dibenz (a,h) anthracene	ND	300	"
Dibenzofuran	ND	300	"
Di-n-butyl phthalate	ND	300	"

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Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

1,2-Dichlorobenzene	ND	300	ug/kg
1,3-Dichlorobenzene	ND	300	"
2,4-Dichlorophenol	ND	1000	"
Diethyl phthalate	ND	300	"
2,4-Dimethylphenol	ND	1000	"
Dimethyl phthalate	ND	300	"
4,6-Dinitro-2-methylphenol	ND	1000	"
2,4-Dinitrophenol	ND	1000	"
2,6-Dinitrotoluene	ND	1000	"
Di-n-octyl phthalate	ND	300	"
Fluoranthene	ND	300	"
Fluorene	ND	300	"
Hexachlorobenzene	ND	1500	"
Hexachlorobutadiene	ND	300	"
Hexachlorocyclopentadiene	ND	1000	"
Hexachloroethane	ND	300	"
Indeno (1,2,3-cd) pyrene	ND	300	"
Isophorone	ND	300	"
2-Methylphenol	ND	1000	"
4-Methylphenol	ND	1000	"
Naphthalene	ND	300	"
2-Nitroaniline	ND	300	"
3-Nitroaniline	ND	300	"
4-Nitroaniline	ND	300	"
Nitrobenzene	ND	1000	"
2-Nitrophenol	ND	1000	"
N-Nitrosodimethylamine	ND	300	"
N-Nitrosodiphenylamine	ND	300	"
2,3,5,6-Tetrachlorophenol	ND	300	"
2,3,4,6-Tetrachlorophenol	ND	300	"
Phenanthrene	ND	300	"
Azobenzene	ND	300	"
2,4,5-Trichlorophenol	ND	1000	"
Pyridine	ND	300	"
2,4,6-Trichlorophenol	ND	1000	"

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Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Blank (8041634-BLK1)

Prepared: 04/16/18 Analyzed: 04/17/18

Surrogate: 2-Fluorophenol	1790		ug/kg	3380		53.0	15-121			
Surrogate: Phenol-d6	1920		"	3380		56.9	24-113			
Surrogate: Nitrobenzene-d5	2150		"	3380		63.5	21.3-119			
Surrogate: 2-Fluorobiphenyl	2510		"	3380		74.3	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3160		"	3380		93.4	18.1-105			
Surrogate: Terphenyl-d14	2960		"	3380		87.6	29.1-130			

LCS (8041634-BS1)

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1490	1000	ug/kg	3280		45.6	34-114			
2-Chlorophenol	1620	1000	"	3280		49.4	34-114			
1,4-Dichlorobenzene	1680	300	"	3280		51.4	34-114			
N-Nitrosodi-n-propylamine	1700	300	"	3280		51.9	30-110			
1,2,4-Trichlorobenzene	1900	300	"	3280		58.0	39-119			
4-Chloro-3-methylphenol	2150	1000	"	3280		65.6	50-130			
Acenaphthene	1720	300	"	3280		52.6	34-114			
Pentachlorophenol	2280	1000	"	3280		69.7	50-130			
Pyrene	1530	300	"	3280		46.8	30-110			
Surrogate: 2-Fluorophenol	1450		"	3280		44.3	15-121			
Surrogate: Phenol-d6	1650		"	3280		50.4	24-113			
Surrogate: Nitrobenzene-d5	1790		"	3280		54.7	21.3-119			
Surrogate: 2-Fluorobiphenyl	2130		"	3280		65.1	32.4-102			
Surrogate: 2,4,6-Tribromophenol	2810		"	3280		85.6	18.1-105			
Surrogate: Terphenyl-d14	2600		"	3280		79.4	29.1-130			

Matrix Spike (8041634-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1750	1000	ug/kg	3390	ND	51.7	34-114			
2-Chlorophenol	1860	1000	"	3390	ND	54.9	34-114			
1,4-Dichlorobenzene	1830	300	"	3390	ND	54.0	34-114			
N-Nitrosodi-n-propylamine	1880	300	"	3390	ND	55.4	30-110			
1,2,4-Trichlorobenzene	2190	300	"	3390	ND	64.5	39-119			
4-Chloro-3-methylphenol	2550	1000	"	3390	ND	75.3	50-130			
Acenaphthene	1960	300	"	3390	ND	57.8	34-114			
Pentachlorophenol	2620	1000	"	3390	ND	77.3	50-130			
Pyrene	1980	300	"	3390	160	53.7	30-110			
Surrogate: 2-Fluorophenol	1600		"	3390		47.3	15-121			
Surrogate: Phenol-d6	1890		"	3390		55.8	24-113			

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Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes

Batch 8041634 - EPA 3550 ECD/GCMS

Matrix Spike (8041634-MS1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Surrogate: Nitrobenzene-d5	2040		ug/kg	3390		60.1	21.3-119			
Surrogate: 2-Fluorobiphenyl	2310		"	3390		68.1	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3090		"	3390		91.2	18.1-105			
Surrogate: Terphenyl-d14	2670		"	3390		78.7	29.1-130			

Matrix Spike Dup (8041634-MSD1)

Source: T181265-01

Prepared: 04/16/18 Analyzed: 04/17/18

Phenol	1720	1000	ug/kg	3410	ND	50.4	34-114	2.00	42	
2-Chlorophenol	1840	1000	"	3410	ND	53.9	34-114	1.32	40	
1,4-Dichlorobenzene	1760	300	"	3410	ND	51.7	34-114	3.82	28	
N-Nitrosodi-n-propylamine	1840	300	"	3410	ND	54.0	30-110	1.82	38	
1,2,4-Trichlorobenzene	2090	300	"	3410	ND	61.4	39-119	4.29	28	
4-Chloro-3-methylphenol	2560	1000	"	3410	ND	74.9	50-130	0.108	42	
Acenaphthene	1960	300	"	3410	ND	57.3	34-114	0.119	31	
Pentachlorophenol	2860	1000	"	3410	ND	83.9	50-130	8.97	50	
Pyrene	1940	300	"	3410	160	52.1	30-110	2.05	31	
Surrogate: 2-Fluorophenol	1630		"	3410		47.8	15-121			
Surrogate: Phenol-d6	1870		"	3410		54.8	24-113			
Surrogate: Nitrobenzene-d5	1870		"	3410		54.9	21.3-119			
Surrogate: 2-Fluorobiphenyl	2210		"	3410		64.7	32.4-102			
Surrogate: 2,4,6-Tribromophenol	3080		"	3410		90.1	18.1-105			
Surrogate: Terphenyl-d14	2940		"	3410		86.0	29.1-130			

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
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Brusca Associates Inc.
PO Box 332
Roseville CA, 95661

Project: Lathrop 48-Ac Property
Project Number: 137-002
Project Manager: Joe Brusca

Reported:
04/23/18 10:35

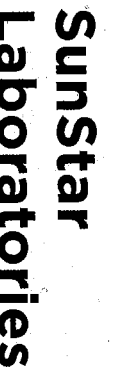
Notes and Definitions

DET Analyte DETECTED
ND Analyte NOT DETECTED at or above the reporting limit
NR Not Reported
dry Sample results reported on a dry weight basis
RPD Relative Percent Difference

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Mike Jaroudi, Project Manager



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Project Manager: Joe Brusca

Page: 11 Of 11

Project Name: LATIT4PROP 4B-Ac PROP=RTT

Collector: Brusca Client Project #: 137-002

Batch #: 2 EDF #: 7181265

EDF #:

Sample disposal Instructions: Disposal @ \$2.00 each _____

Return to client _____

Pickup _____

Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>	4/13/18 1:30 PM	<i>[Signature]</i>	4/13/18 1330
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>		<i>[Signature]</i>	
Relinquished by: (signature)	Date / Time	Received by: (signature)	Date / Time
<i>[Signature]</i>		<i>[Signature]</i>	

	Total # of containers
	Chain of Custody seals Y/N/A
	Seals intact? Y/N/A
	Received good condition/cold
Turn around time:	Normal
	5/3

Notes

Sample ID
SP5-1
SP5-2
SP5-3
SP5-4

Date Sampled 4/13/14

Time	10:00	10:04	10:06	10:09

Sample Type	Solid

Container Type
JAR

~~XXXX~~ 8260 SVOLs 8270

				8260 + OXY

				8260 BTEX, OXY only

~~XXXX~~ 0270 PCBs 3032

				8021 BTEX

				8015M (gasoline)

				8015M (diesel)

~~2~~	~~2~~	~~2~~	~~2~~	8015M Ext./Carbon Chain

~~X~~	~~X~~	~~X~~	~~X~~	6010/7000 Title 22 Metals

6020 ICP-MS Metals

					0020 101 -MS Metals
					~~LEAD AND ASSY~~

01	Laboratory ID #
02	
03	
04	

Comments/Preservative

— — — — Total # of containers

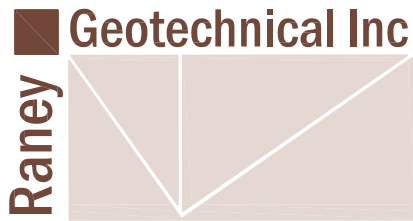
GOC 170169

APPENDIX B – Foundation Investigation

FOUNDATION INVESTIGATION
MURPHY PARKWAY WAREHOUSE
Murphy Parkway near D'Arcy Parkway
Lathrop, California

RANEY GEOTECHNICAL INC. JOB NO. 146-618





December 19, 2016

Buzz Oates Development LP
Attention: Troy Estacio
8615 Elder Creek Road
Sacramento, CA 95828

**FOUNDATION INVESTIGATION
MURPHY PARKWAY WAREHOUSE**

Murphy Parkway near D'Arcy Parkway
Lathrop, California
Job No. 146-618

INTRODUCTION

Our firm has completed a Foundation Investigation for a warehouse proposed for construction east of Murphy Parkway in Lathrop. This report presents the results of the investigation. The purpose of this report is to provide data pertinent to earthwork construction, foundation design, floor support, and pavement design.

Field exploration for this investigation has included the drilling of ten test borings to depths of ten to 20 feet below the existing site grade, and one boring to a depth of 50 feet. Undisturbed soil samples were obtained from the borings for classification and laboratory testing. Disturbed samples of the near surface soils also were obtained from proposed pavement areas. The test boring and sampling locations are shown on Plate 1, *Plot Plan*; logs of the test borings completed for this study are shown on Plates 2 through 16, *Log of Boring*. The nomenclature used to describe the soils on the logs is defined on Plate 17, *Unified Soil Classification System*. Moisture, density and unconfined compressive strength test data are presented on the logs at the depths of each sample tested. Two samples of the subsurface sandy soils were subjected to sieve analyses to provide information used in our liquefaction analyses; these test results are graphed on Plate 18, *Grain Size Analyses*. Near surface soil samples were subjected to direct shear tests to evaluate strength properties; these test results are tabulated on Plate 19, *Direct Shear Data*. Samples of anticipated pavement subgrade soils were subjected to resistance value tests to evaluate pavement subgrade properties; the test results are shown on Plates 20 and 21, *Resistance Value Data*. A sample of the subgrade soils was mixed with cement, remolded, and tested in unconfined compression to evaluate reactivity with the treatment chemical; these test results are tabulated on Plate 22, *Treated Soil Compressive Strength Data*. Two samples of the subsurface soils were submitted to an associate chemical laboratory for testing of corrosion properties; these test results are reproduced on Plates 23 and 24. The results of liquefaction analyses are depicted on Plates 25, *Liquefaction Analysis*.

PROPOSED CONSTRUCTION

We understand the property will be developed for a single distribution warehouse. The warehouse will have plan dimensions of about 510 feet by 1680 feet. The building construction is expected to be of concrete tilt-up-panel design with a concrete slab-on-grade floor. Depressed portland cement concrete paved loading docks would be provided along the north and south sides of the building. Asphalt concrete paved automobile parking and driveways would be placed along the west side of the building. Truck driveways would surround the building and connect the docks to Murphy Parkway.

Building foundation loads are expected to be moderate and typical to this type of construction.

For the purposes of this report we have assumed that the building pad level will be near the average existing site grade.

SITE CONDITIONS AND CONCLUSIONS

SURFACE

The subject site encompasses 48.6 acres and is situated between about 500 feet and 2800 feet east of the northerly end of Murphy Parkway. The property is bordered on the north by the shuttered Pilkington glass plant, and on the south by warehouse, trucking and distribution businesses. A pasture and lands of a Simplot fertilizer plant are to the east of the subject parcel. A shuttered cogeneration plant that apparently supplied the Pilkington plant, occupies the northerly portion of the parcel between Murphy Parkway and the subject site. An earth berm, power lines, and chain link fencing are along the northerly border at the glass plant. A drainage swale with trees is along the southerly edge of the site. Wire fencing is along the westerly border.

Most of the property is relatively flat and is currently used as sheep pasture. The easterly portion supports dry weeds and new grass growth. Patchy weeds are present on the westerly portion. An eight-to ten-foot deep depression of an apparent former pond is on the northwesterly portion of the site. It appears the pond may have been used for disposal by the adjacent glass or cogeneration plants. Aerial photos also suggest that glass had been stored or disposed of on the northwesterly corner of the property, and was later spread further to the south. We observed considerable amounts of broken glass on the surface and mixed into the surface soils of the west edge of the subject property as well as on the southerly portion of the cogeneration parcel to the west.

SUBSURFACE

Boring 2 was drilled on the northwesterly corner of the property and found about five feet of fill on the surface. The fill was observed to consist loose to medium dense silty fine sands and to contain glass fragments and shards. Boring 2 was located in an area that is shown on aerial photographs to have supported large piles of glass beads or fragments.

Except for the surface fills at Boring 2, the test borings revealed relatively uniform soil conditions across

the property. The test borings generally encountered surface soils consisting of loose to medium dense, brown silty fine sands with some fine sandy silts and extending to depths of four to six feet.

Beneath the surface sandy soils medium dense and sometimes variably cemented light brown fine sandy silts and silty fine sands often were encountered. These silty soils were observed to extend to depths varying from six to about nine feet. The silty layers often were underlain by medium dense to very dense white to light gray clean fine to medium sands. The clean sands were observed to extend to depths varying from 12 to 15 feet. Beneath these sands and extending to the 50 foot maximum depth drilled, interlayered light brown to gray stiff clayey fine sandy silts and very stiff silty fine sandy clays, as well as medium dense to very dense gray fine to medium sands were encountered.

Borings 14 and 15 were drilled within the former pond areas on the northwesterly portion of the site and were drilled at elevations that appear to be seven to eight feet below the site grades outside of the pond limits. These two borings encountered interlayered sands and silts that are consistent with the soil profile of at similar elevations of the other test borings.

BEARING CAPACITY

The soils within the upper six or nine feet include loose sand and silt layers that can be subject to some densification and settlement under heavy loads. The native undisturbed soils are considered capable of supporting the planned construction. The building may be supported on spread foundations, however foundation bearing pressures must be limited to control settlements of the loose near surface soils. With the recommended earthwork preparation procedures, new engineered fills placed and compacted in accordance with the recommendations of this report are expected to have sufficient strength for support of the planned construction.

EXPANSIVE SOILS

The near surface soils consist primarily of low plasticity sands and silt. These soils are considered to be of low expansion potential. Expansive soils are not expected to have a significant effect on the planned construction.

GROUNDWATER

Groundwater was encountered in the test borings at depths varying from 13 to 17 feet below the ground surface. After drilling, water levels in the borings appeared to stabilize at depths of about varying from about 11 to 14 feet below the site grade.

Although the groundwater level will fluctuate somewhat with rainfall and season, the level is expected to remain below final grade levels and not have a significant effect on completed surface improvements. Deep utility or other construction excavations may experience groundwater inflow. Foundation excavations at loading docks may encounter groundwater if constructed during a high groundwater period. Construction excavations extending more than about six feet below the original site grade may engage clean sands that can be subject to sloughing and caving regardless of the groundwater level. Below groundwater levels, the site silty soils also may be subject to sloughing and caving. Sloughing

and caving soils may require formwork to keep them open.

The surface soils include low permeability layers and have variable drainage characteristics. The surface soils have high silt contents that render them sensitive to moisture contents. High moisture content soils are likely to be unstable under construction equipment, and require considerable aeration in order to achieve a moisture content that will allow compaction. The surface soils are likely to have high moisture conditions during the wet season and for several weeks following rains or irrigation. Subsurface soils within a few feet of groundwater are likely to have high moisture contents year round. The potential for high soil moisture contents should be considered in construction scheduling.

SOIL LIQUEFACTION POTENTIAL

Soil liquefaction is the loss of strength of low- to no- cohesion soils (usually sands) that occurs when pore water pressure exceeds the confining stress (weight) of the soils. Liquefaction normally occurs only under saturated conditions and in soils with a low relative density. Liquefaction can occur during earthquakes as vibrations induce soils to readjust to a more compact state. Experience has shown that earthquake induced liquefaction normally occurs only within the upper 50 to 60 feet of the soil profile.

The test borings show loose to medium dense sands and sandy silts within the upper nine feet. Layers of medium dense to dense sandy soils are indicated between depths of six and 50 feet. The soils below depths of 13 to 14 feet below the building pad level are likely to be saturated year round. Slightly higher groundwater levels may be possible, but the duration of such high groundwater levels is expected to be limited.

We have used the methods of Seed and others as implemented in the LiquefyPro software by CivilTech Corporation to assess the liquefaction susceptibility of the sandy soils, as well as to evaluate potential settlements of the loose near surface soils during seismic shaking. The method correlates standard penetration resistance and liquefaction potential based on historical case studies. In determining liquefaction potential, groundwater depth, confining pressures, and, intensity and duration of potential ground shaking are considered.

Probabilistic seismic hazards mapping by the United States Geological Survey indicates that the peak ground acceleration produced by maximum credible earthquakes on nearby faults (10 percent probability of exceedance in 50 years) is likely to be on the order of 0.28 g. Interpretations according to the ASCE 7-10 standard indicates that the peak ground acceleration produced by maximum credible earthquakes with a two percent probability of exceedance in 50 years is likely to be on the order of 0.43g. The controlling earthquakes in developing these accelerations appear to mostly be events of magnitude 6.5 to 6.7 occurring within the Coast Range/Central Valley (CRCV) boundary zone about 21 kilometers to the southwest. Using these earthquake data, our analysis indicates that seismic induced liquefaction on this project is unlikely, although some sandy soils at depths of near 40 feet may approach a state of liquefaction under the highest levels of shaking. Settlements of the soils during maximum seismic shaking are estimated to be on the order of one quarter inch. The results of the analyses can be viewed on Plate 25, *Liquefaction Analysis*.

CORROSION POTENTIAL

Two samples of the native soils from anticipated pipe horizons were submitted to an associate laboratory for corrosion evaluation tests; the test data are presented on Plates 23 and 24.

We have used criteria established by the Ductile Iron Pipe Research Association to evaluate the corrosivity of the soils tested to iron pipe. The criteria assigns point values to the test results as well as a subjective evaluation of moisture conditions at pipe level; a total point value of ten or higher indicates the soils are potentially corrosive to iron pipe. Our evaluation indicates total point values of 2. We conclude that the soils are not corrosive to gray or ductile iron pipe.

Based on the Portland Cement Association (PCA) guidelines, the soluble sulfate content of the site soils indicates that the relative degree of sulfate attack on concrete is negligible, and therefore use of Type I, Type II, or Type 1-II portland cement is appropriate.

RECOMMENDATIONS

EARTHWORK

The building pad and pavement areas should be cleared of surface vegetation, trees and root systems, rubble fragments exceeding three inches in maximum dimension, rubbish, and any other debris. Any earth berms, stockpiled soils and any other existing fills should be removed. Fills were found in Boring 2 to a depth of about five feet on the northwesterly corner of the property. These fills should be overexcavated to allow replacement as engineered fill. Test excavations should be made on the northwesterly corner of the property and adjacent areas with our representative present to check for the presence of old fill materials. Unsuitable soils identified by our representative should be removed.

Any underground pipes within two feet of original or final grade (whichever is lower) should be removed. Abandoned pipes exceeding two inches in diameter should be removed from the building areas regardless of depth. Low areas should be cleaned out of any loose or saturated materials. Excavations required for the removal of the above items, as well as any other loose or unstable soil deposits identified by our representative, should be cleaned of loose, saturated or soft materials so that firm undisturbed soils are exposed. Deep excavations required for the removal of the above items should be sloped back to a dish shaped configuration allowing through passage of compaction equipment. Any deep excavations should be restored to grade with engineered fill placed and compacted in accordance with the recommendations of this report.

Areas designated to receive engineered fill as well as building pad and pavement areas left at existing grade should be scarified to a depth of eight inches, brought to a uniform moisture content of at least optimum, and compacted in place to at least 90 percent of the maximum dry density determined by ASTM D1557-02 test procedure. If recompacted subgrades are unstable or compaction cannot be achieved due to high soil moisture contents, our firm should be contacted for further recommendations.

Engineered fill should be placed in lifts not exceeding six inches in compacted thickness, brought to a

moisture content of at least optimum, and compacted. Fill placed in the building pad area at depths of more than two feet below the final soil building pad level should be compacted to at least 95 percent of the ASTM D1557 maximum dry density. All other fills should be compacted to at least 90 percent of the maximum dry density in accordance with the above standard. On-site soils are suitable for use as engineered fill provided the soils do not contain significant vegetable matter, rubble, rubbish, or other undesirable substances. Import materials, if any, should consist of sands, silts, or gravels with a plasticity index of ten or less and should be tested and approved by this firm prior to importation to the site.

To enhance slab performance and provide a more stable base for working during the rainy season, the upper portion of the building pad and any adjacent concrete flatwork areas should be treated with portland cement. The treatment should extend at least five feet outside of building wall lines, or to the outer edges of surrounding portland cement concrete walkways or aprons, whichever is greater in extent. The soils should be treated to a depth of at least 12 inches with at least five percent portland cement as measured by dry unit weight of the compacted soil. The treated soil should be brought to a uniform over optimum moisture content, thoroughly mixed with the cement, and compacted to at least 92 percent of the maximum dry density determined by the ASTM D1557-02 test procedure. Not more than three hours should elapse between the time the cement is mixed with the soil/water and final compaction is completed. The above concentration of treatment chemical are based on assumptions made regarding the soil composition of the upper 12 inches of the building pad. Some modification to this concentration may be required. The percentage and type of treatment chemical be used should be established by our engineers based on a review of the types of soils present on each area to be treated. Treatment should conform to applicable provisions of the Caltrans Standard Specifications, Sections 24 and 27. The treated pad should either be kept wet for a period of at least three days after compaction or seal coated.

Untreated pavement subgrades, if any, should be compacted to at least 90 percent of the ASTM D1557-02 maximum dry density regardless of whether the final grade is achieved by cutting, filling or left at final grade.

Permanent excavation and embankment slopes should not exceed an inclination of one vertical on two horizontal. A representative of this firm should be present during grading operations to test and observe earthwork construction.

FOUNDATIONS

The proposed building may be supported upon continuous and/or isolated spread foundations based in undisturbed or recompacted natural ground, engineered fill, or a combination of these materials. Building foundations should extend to a minimum depth of 18 inches below the building pad. Foundations in dock areas should extend at least 12 inches below the lowest surrounding soil subgrade level. Foundation excavations should be observed by an Engineer from this office to ensure that proper bearing soils are engaged. A minimum foundation width of 12 inches should be maintained. Foundations so established may be designed for maximum allowable soil bearing pressures of 2000 pounds per square foot (psf) for dead load, 2500 psf for dead plus live load, and 3300 psf for total load including the effects of wind and seismic forces.

Foundations for trash enclosure walls, signs, landscape retaining walls, and other appurtenant construction should extend a minimum of 12 inches below the building pad or lowest surrounding soil grade, and have a minimum width of 12 inches. Such foundations may be sized for allowable soil bearing pressures of 1500 psf for dead load, 2300 psf for dead plus live load, and 3000 psf for dead plus live plus seismic/wind loads.

The weight of foundation concrete below grade may be disregarded in sizing computations. Continuous foundations should be reinforced with at least four No. 4 bars--two each, top and bottom, as a minimum. The actual reinforcement should be determined by the Structural Engineer.

Foundation excavations should not be allowed to stand open for extended periods prior to concrete placement and should be clean and free of all loose and/or soft materials when concrete is placed. The bearing materials should be in a firm, moist condition. Some dock foundation excavations may engage clean or saturated sands and experience sloughing and caving; forming may be required.

Resistance to lateral forces may be computed using either friction or passive pressure, but not both, except as recommended below. A coefficient of friction of 0.30 may be utilized for design. Passive earth pressure may be considered equivalent to a fluid weighing 250 pounds per cubic foot. A combination of both friction and passive pressure may be utilized provided the larger mode of resistance is reduced by 50 percent. The recommended passive pressure and friction coefficient values have been modified by appropriate factors of safety and may be applied directly in design calculations.

SEISMIC DESIGN

In design using the lateral force provisions of the 2013 and 2016 California Building Codes, the parameters in Table 1 may be used.

TABLE 1

Period (seconds)	Mapped Spectral Response Accelerations (g)		Site Class	Site Coefficients		Maximum Considered Earthquake Spectral Response Accelerations (g)		Design Spectral Response Accelerations (g)	
0.2	S_s	1.065	D	F_a	1.074	S_{MS}	1.144	S_{DS}	0.762
1	S_1	0.372		F_v	1.655	S_{M1}	0.616	S_{D1}	0.411

LOADING DOCK WALLS

Restrained dock walls should be capable of resisting an at-rest soil pressure equivalent to that exerted by a fluid weighing 50 pounds per cubic foot. Wall design also should consider pressures associated with traffic or other adjacent surcharge. Care should be taken to avoid exertion of excessive compaction

pressures in backfilling of walls. The above loading is based on the assumption that hydrostatic pressures will not develop behind the walls.

Walls should be drained as needed to relieve potential hydrostatic pressure. Where floor slabs or pavements adjoin the tops of walls, this should suffice to exclude significant water and further drainage measures would not be required.

Only nonexpansive silts, sands or gravels should be used to backfill dock area walls. Clays should not be used in wall backfill. Use of 3/8-inch pea gravel or clean 3/4-inch crushed rock for wall backfill is acceptable and can facilitate placement and compaction procedures. Backfill should be placed in level lifts not exceeding 12 inches in compacted thickness. Each lift should be compacted at a uniform near optimum moisture content to at least 90 percent of the ASTM D1557-02 maximum dry density.

SLAB-ON-GRADE

Thickness Requirements

Because of stresses induced by truck cranes during panel erection a minimum six-inch thick floor slab is suggested for all floor areas. Our calculations indicate a six-inch thick slab would be capable of supporting typical 3000- to 5000-pound capacity forklifts carrying loads of less than 3000 pounds. If the floor will support more heavily loaded forklifts, or if high racks are used, then a six-inch slab may not be sufficient and our firm should be contacted for further recommendations. A modulus of subgrade reaction of 150 pounds per cubic inch may be used for the treated building pad in determining slab thickness for forklift loads.

Due to the industrial use, we suggest the floor slab be reinforced with No. 4 bars spaced on 24-inch centers in each direction as a minimum. The reinforcement should be chaired at the middepth of the slab. The above discussion of slab thickness is based on use of quality strength concrete (4000 pounds per square inch minimum 28-day compressive strength). We suggest slab concrete be placed at a slump of three to four inches. Fibermesh® may be used in concrete to increase toughness, if desired.

The floor slab should be thickened by at least 20 percent of the above recommended thickness at any drive-through doors. In addition, edges should be thickened wherever heavy materials will be stored within five feet of a free slab edge. A free slab edge is defined as any joint or edge where load transfer to adjacent areas is not provided, such as at building wall lines or undoweled expansion joints. The transition to the thickened edge may be achieved by tapering the slab thickness over a distance of five feet. Tapering may be accomplished by reducing the underlying aggregate base thickness at the edges.

Joints

Shrinkage crack control joints should consist of sawcut grooves penetrating at least one-fourth of the slab thickness. Control joints should not be spaced farther apart than about 30 times the slab thickness. Construction cold joints and expansion joints in the warehouse slab should include dowels to provide load transfer. Dowels should be three-quarters inch in diameter, 14 inches long, and spaced on 12 inch centers for six inch thick slabs. One end of dowels at expansion and crack control joints should be

greased and wrapped in plastic to allow horizontal movement. All reinforcement and dowels should be placed at the middepth of the slab.

Underlayment and Moisture Control

The floor slab may be supported on the chemically treated building pad prepared as recommended above. In warehouse areas where minor moisture penetration through the slabs can be tolerated, the slab should be underlain by a minimum four-inch thick layer of Caltrans Class 2 aggregate base to serve as a leveling course. The aggregate base should be wetted immediately prior to placement of slab concrete. Moisture sensitive areas such as office areas receiving impervious floor coverings should be underlain by a minimum four-inch thick layer of clean three-quarter inch crushed rock graded such that 100 percent passes a one-inch sieve and none passes a No. 4 sieve, to serve as a capillary moisture break. In such areas the drainage rock should be covered with a plastic membrane at least 10-mils thick as a moisture vapor retarder. One to two inches of clean sand may be spread over the membrane for protection and to aid concrete curing, if desired. Alternatively, the membrane may be placed directly on the building pad beneath the rock.

With the use of water-based floor adhesives, impervious floor coverings are extremely sensitive to slab moisture. Under some conditions, the small amount of moisture vapor which bypasses the vapor membrane, or even the excess water remaining in the slab from placement, can be sufficient to cause debonding and discoloration problems. To minimize moisture vapor problems, the capillary break gravel must be present to the minimum recommended thickness and the membrane must be continuous throughout the slab area. Any membrane seams should overlap by at least one foot. The membrane should be cut tight to penetrations. Tears and punctures should be sealed with membrane manufacturer-approved tape, or overlain by a second patching membrane. Slab concrete should be placed at as low a water-cement ratio as practical. The under-slab gravel layer should be protected from precipitation and other moisture; wetting of the sand over the membrane prior to concrete placement should be minimized. The edges of the slab at the building perimeter should be thickened to form a cutoff between the building exterior and under-slab gravel layers. If impervious floor coverings are planned and greater assurance against moisture problems is desired, consideration should be given to waterproofing of slabs with a quality commercial concrete sealant. A sealant or other waterproofing system may be necessary for the satisfactory performance of wood laminates, sheet vinyl, and other impervious flooring.

PAVEMENTS

Resistance (R) value tests are used to evaluate pavement subgrade properties. Our investigation indicates that the pavement subgrades are likely to consist of very silty sands of fair to good pavement support qualities. R values can range from five for the poorest quality clay subgrades to 70 or higher for high quality sand and gravel subgrades. Resistance value tests performed on samples of the site surface soils resulted in design R values of 64 and 65. We consider these R values to be somewhat high for the range of soils that are likely to be present on pavement subgrades, and therefore have used an R value of 45 in our calculations. The R value of 45, together with the results of unconfined compressive strength tests on soils treated with cement, have been used in the Caltrans Design Method for Flexible Pavements and in Portland Cement Association's design method to evaluate pavement sections.

The Caltrans design method uses traffic indices to account for vehicle loads, frequency, and design life. A design life of 20 years is commonly used for commercial pavements.

Based upon these data, the standard Oates structural pavement section consisting of three inches of asphalt concrete (AC) over three inches of aggregate base (AB) over 12 inches of chemically treated soil is considered to have a traffic capacity equivalent to TI 7.0. A TI of 7.0 is considered capable of supporting up to about 60 fully loaded, five-axle semi-trucks per week. Pavement sections designed for TI 8.0 are considered capable of supporting up to about 190 fully loaded, five axle semi-trucks per week. Our calculations indicate that pavement sections for TI 8.0 may consist of four inches of AC over three inches of AB over 12 inches of chemically treated subgrade, or three and one-half inches of AC over four inches of AB over 12 inches of chemically treated subgrade. These pavement sections, together with alternative asphalt concrete pavement sections without chemically treated subgrades, are summarized in Table 2. We can provide additional pavement section alternatives upon request.

TABLE 2
ASPHALT CONCRETE PAVEMENT SECTION
ALTERNATIVES

Design Traffic Index/ Use	Type B Asphalt Concrete (inches)	Class 2 Aggregate Base (inches)	Chemically Treated Soil (inches)
4.5	2.5	4	--
Auto Parking	2.5	--	10
7.0	3	8	--
Up to 60	3	3	12
Trucks per Week			
8.0	3.5	9	--
120 to 190	4	8	--
Trucks per Week	3.5	4	12
	4	3	12

Materials and construction within structural pavement sections should conform to the applicable provisions of the 2010 Caltrans Standard Specifications

We recommend chemically treated subgrades be treated with portland cement to provide reactivity with the expected sandy subgrade soils. The pavement areas should be treated after the subgrades have been cut or filled to rough final grade and utility lines have been installed. The subgrade soils should be treated with at least five percent portland cement, as measured by dry unit weight of the compacted soil. The cement and soil should be brought to a uniform over-optimum moisture content, thoroughly mixed and compacted to at least 92 percent of the maximum dry density determined by the ASTM D1557-02 test procedure. Not more than three hours should elapse between the time the cement is mixed with the soil/water and final compaction is completed. The above concentration of cement is based on assumptions made regarding the soil composition of the pavement subgrades. The actual percentage and

type of treatment chemicals to be used should be established by our engineers based on a review of the types of soils present on each area to be treated.

All treatment should conform to applicable provisions of the 2010 Caltrans Standard Specifications, Sections 24 and 27, except with regard to compaction as specified above. The treated subgrade should be either kept wet for a period of at least three days after final compaction, or seal coated as recommended in Section 24.

Treated soils tend to experience shrinkage and cracking during curing. Such cracking does not affect the strength of the treated materials, but can lead to cosmetic reflective cracking through asphalt concrete. Depending on the degree of cracking, crack sealing or similar maintenance may be required to ensure long term pavement performance. Use of the nominal layer of Caltrans Class 2 aggregate base between the treated soil layer and the asphalt concrete will decrease the incidence of the described shrinkage cracking.

Recommended portland cement concrete pavement sections for the traffic indices discussed above are presented in Table 3.

TABLE 3
PORTLAND CEMENT CONCRETE SECTION ALTERNATIVES

Design Traffic Index	Portland Cement Concrete (inches)	Class 2 Aggregate Base (inches)	Chemically Treated Soil (inches)
6.0	6	2	12
7.0	6.5	2	12
8.0	7	2	12

The aggregate base layer is considered a leveling course and its thickness may be varied in the field for leveling purposes. Chemically treated soil subgrades should be constructed as recommended in the Foundation investigation report.

Any one section of Portland cement concrete pavements at depressed loading docks would be expected to support only a few trucks per day, including parking tractor-trailers and inadvertent travel of maneuvering vehicles. Based on this, traffic index 6.0 is considered appropriate for dock pavements.

The above portland cement concrete pavement sections are based on use of concrete with a minimum modulus of rupture of 490 pounds per square inch. We anticipate this will require use of concrete with a minimum compressive strength at 28 days of 3000 pounds per square inch. Thickened slab edges should be provided at isolation joints between the concrete pavement and adjacent construction, at the transition from portland cement concrete pavement to asphalt concrete pavement, and adjacent to unpaved areas. The thickened edge should be at least one and one-half inches greater than the unreinforced thicknesses recommended above, and should be tapered back to the design slab thickness over a distance of five feet. Narrow concrete strips used for support of trailer parking dollies should have minimum thicknesses of 7.5 inches, 8 inches, and 8.5 inches in traffic index 6.0, 7.0, and 8.0 areas,

respectively.

Shrinkage crack control joints should consist of premolded inserts or sawcut grooves penetrating at least one-fourth of the slab thickness. Control joints in unreinforced pavements should not be spaced farther apart than about 25 times the pavement thickness. Construction cold joints should include dowels to provide load transfer. Table 4 shows the recommended dowel bar sizes and spacings.

TABLE 4
DOWEL SIZE AND SPACING

Pavement Thickness (inches)	Dowel Diameter (inches)	Dowel Length (inches)	Dowel Spacing center to center (inches)
6.0 - 6.5	3 / 4	16	12
7.0	1	18	12

Where construction cold joints also serve as shrinkage crack control joints, one end of the dowels should be greased and wrapped in plastic to allow horizontal movement. All reinforcement and dowels should be placed at the middepth of the slab.

Where the slab edges are not confined such as adjacent to lower unpaved areas, the outer slab sections of large pavement areas may tend to spread laterally towards the free edge. For unreinforced pavements we recommend non-crack control construction cold joints parallel to and within 25 feet of the free edge include tie bars to hold joints closed. Tie bars should be No. 4 bars at least 21 inches long and spaced on 20 inch centers. Any crack control joints, including sawcut types and construction joints, parallel to and within 25 feet of the free edge should include dowels across the joint. Dowel sizes and spacings should be in accordance with the above recommendations.

LIMITATIONS

This report necessarily assumes uniform variation of soils between borings. Our recommendations are based upon this assumed uniformity and the information provided regarding the proposed construction. If unusual conditions are encountered during construction, the contractor or his representative should notify this firm immediately so that alternate written recommendations can be made.

This report is applicable only to the proposed construction, as described herein, and should not be utilized for design or construction on any other site.

Murphy Parkway Warehouse
December 19, 2016
Job No. 146-618
Page 13

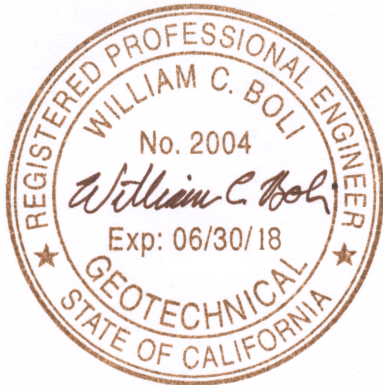
oOo

The following Plates are attached and complete this report:

Plate 1 - Plot Plan
Plates 2 through 16 - Log of Boring, Borings 1 through 15
Plate 17 - Unified Soil Classification System
Plate 18 - Grain Size Distribution
Plate 19 - Direct Shear Data
Plates 20 and 21 - Resistance Value Data
Plate 22 - Treated Soil Compressive Strength Data
Plates 23 and 24 - Corrosion Test Data
Plate 25 - Liquefaction Analysis

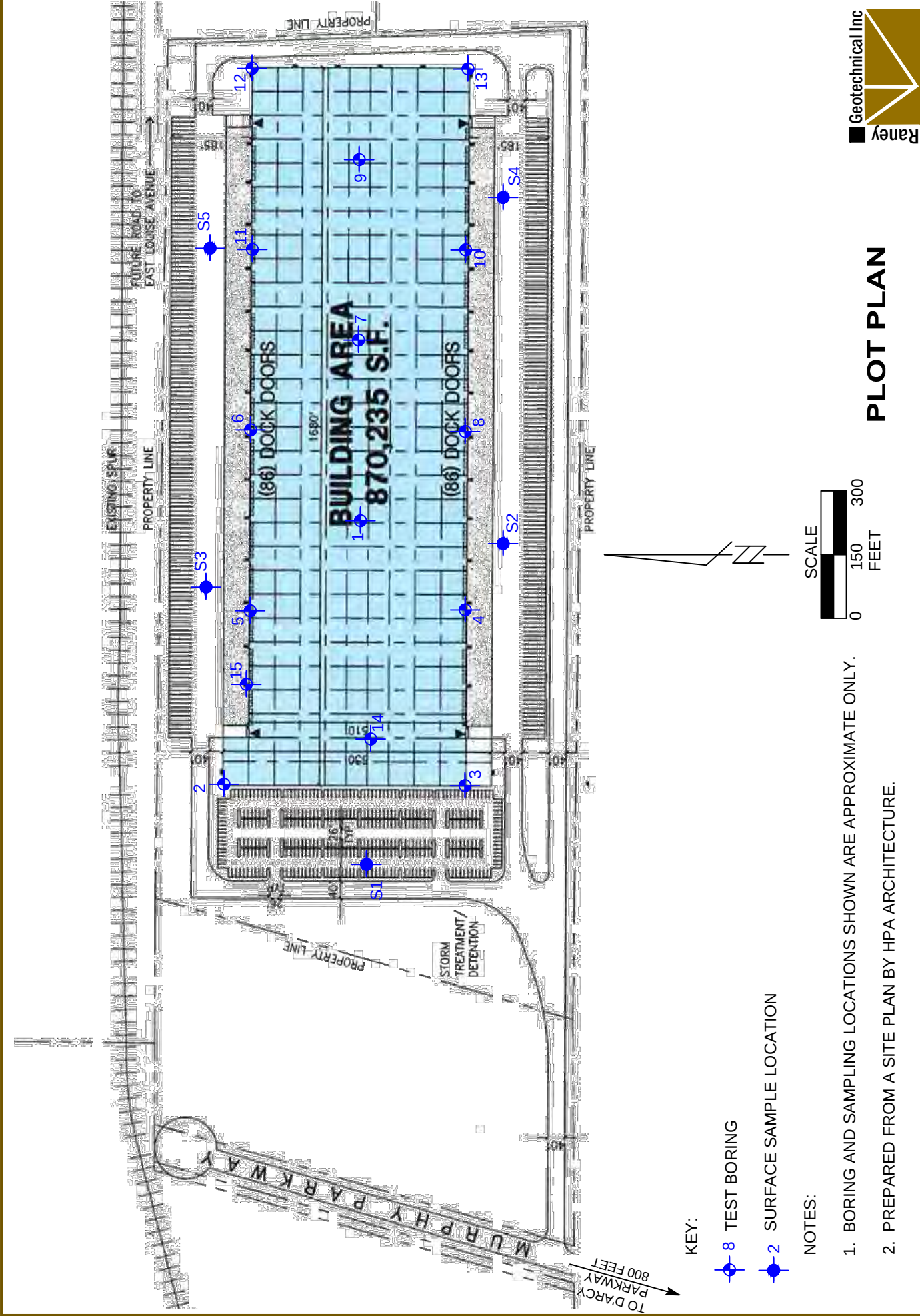
Sincerely,

RANEY GEOTECHNICAL, INC.



William C. Boli
Geotechnical Engineer No. 2004

(1) addressee



PROJECT NUMBER: 146-618
DRAWN BY: WCB
DATE: 12/8/16
PLATE NUMBER: 2

BORING 1

DRILLED: 11/17/16

DEPTH IN FEET			PCF	%	TSF
0					
107	13.5	3.0			
102	22.4	5.2			
103	3.5				
103	23.3				
20					
75	43.1	1.1			
30					
90	31.1	2.9			
40					
109	17.2				
50					

63/10"	SM	BROWN SILTY VERY FINE SAND--LOOSE, MOIST SLIGHTLY CEMENTED
90/10"	ML	GRAY-BROWN VERY FINE SANDY SILT--VERY DENSE, MOIST WITH RUST SPOTS LIGHT BROWN GRADING FINE SANDY
43	SP	LIGHT BROWN AND GRAY FINE TO MEDIUM SAND--DENSE, MOIST
		▼ GROUNDWATER LEVEL AFTER DRILLING 11/17/16
24	ML	LIGHT GRAY-BROWN CLAYEY FINE SANDY SILT--VERY STIFF, SATURATED
	SP	▼ GROUNDWATER FOUND DURING DRILLING 11/17/16
69		GRAY-BROWN FINE TO COARSE SAND--DENSE, SATURATED GRADING WITH LESS FINE SAND
13	CL	DARK GRAY SILTY CLAY--STIFF, SATURATED LIGHT BLUE-GRAY
26	SP	BROWN FINE SAND--MEDIUM DENSE, SATURATED
	CL	DARK GRAY SILTY CLAY--VERY STIFF, SATURATED
	ML	BROWN FINE SANDY CLAYEY SILT--VERY STIFF, SATURATED
	CL	BLUE-GRAY SLIGHTLY FINE SANDY VERY SILTY CLAY WITH ORANGE-BROWN STREAKS--VERY STIFF, SATURATED GRADING LESS FINE SANDY
28		
16	SM	DARK GRAY SILTY M. TO C. SAND-- MED. DENSE GRADING MORE SILTY AND WITH FINE SAND
	CL	DARK GRAY SLIGHTLY SILTY CLAY--VERY STIFF, SATURATED
	SP	DARK GRAY FINE TO MEDIUM SAND--DENSE, SATURATED GRADING WITH SOME COARSE SAND
43		
66		VERY DENSE

NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
3. UNDISTURBED SAMPLE OBTAINED WITH 2" I.D. MODIFIED CALIFORNIA SAMPLER.
4. SAMPLER PENETRATION RESISTANCE IN BLOWS PER FOOT OR FRACTION THEREOF; 140-POUND HAMMER, 30" DROP.

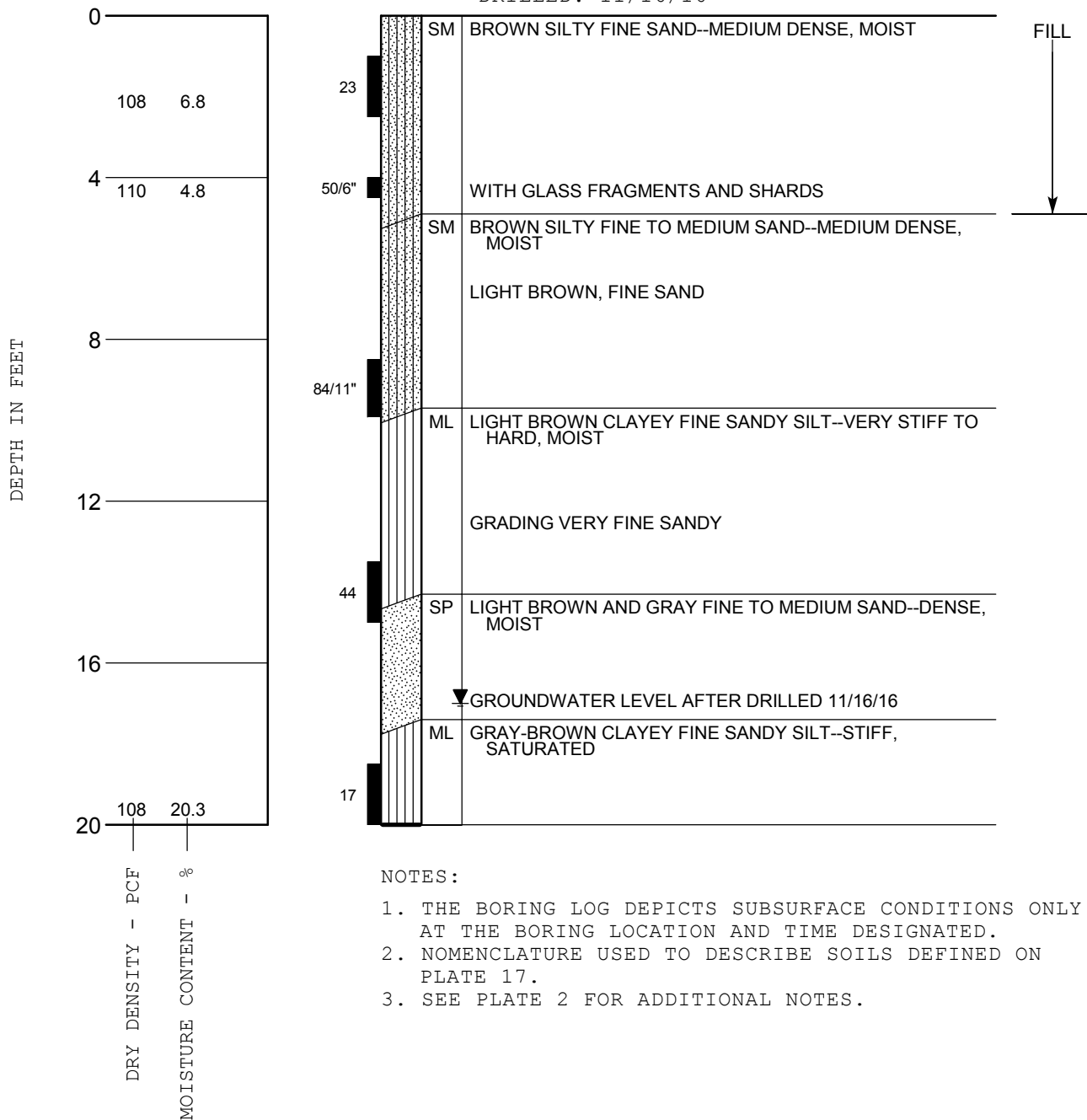
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 3

BORING 2

DRILLED: 11/16/16



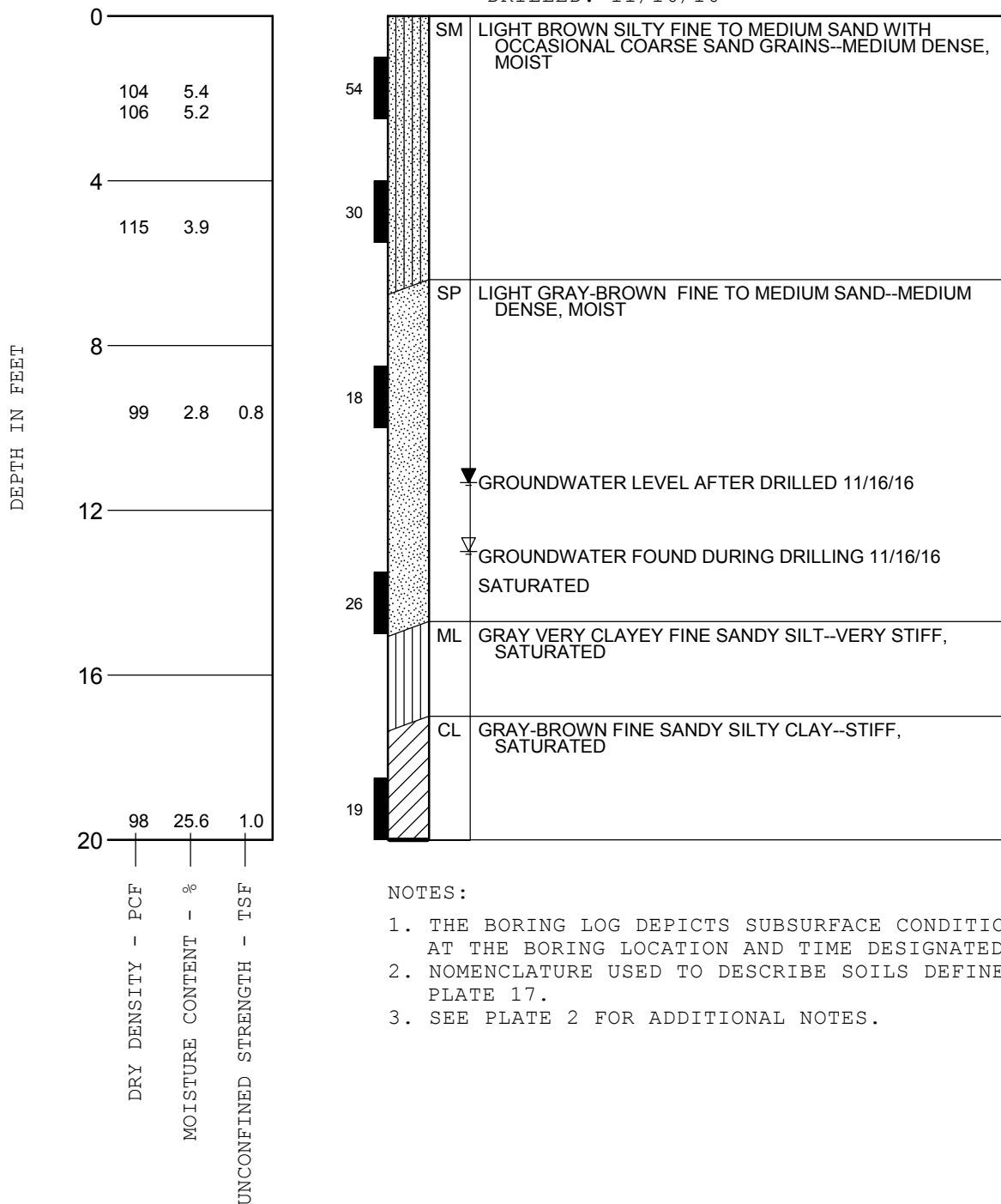
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 4

BORING 3

DRILLED: 11/16/16



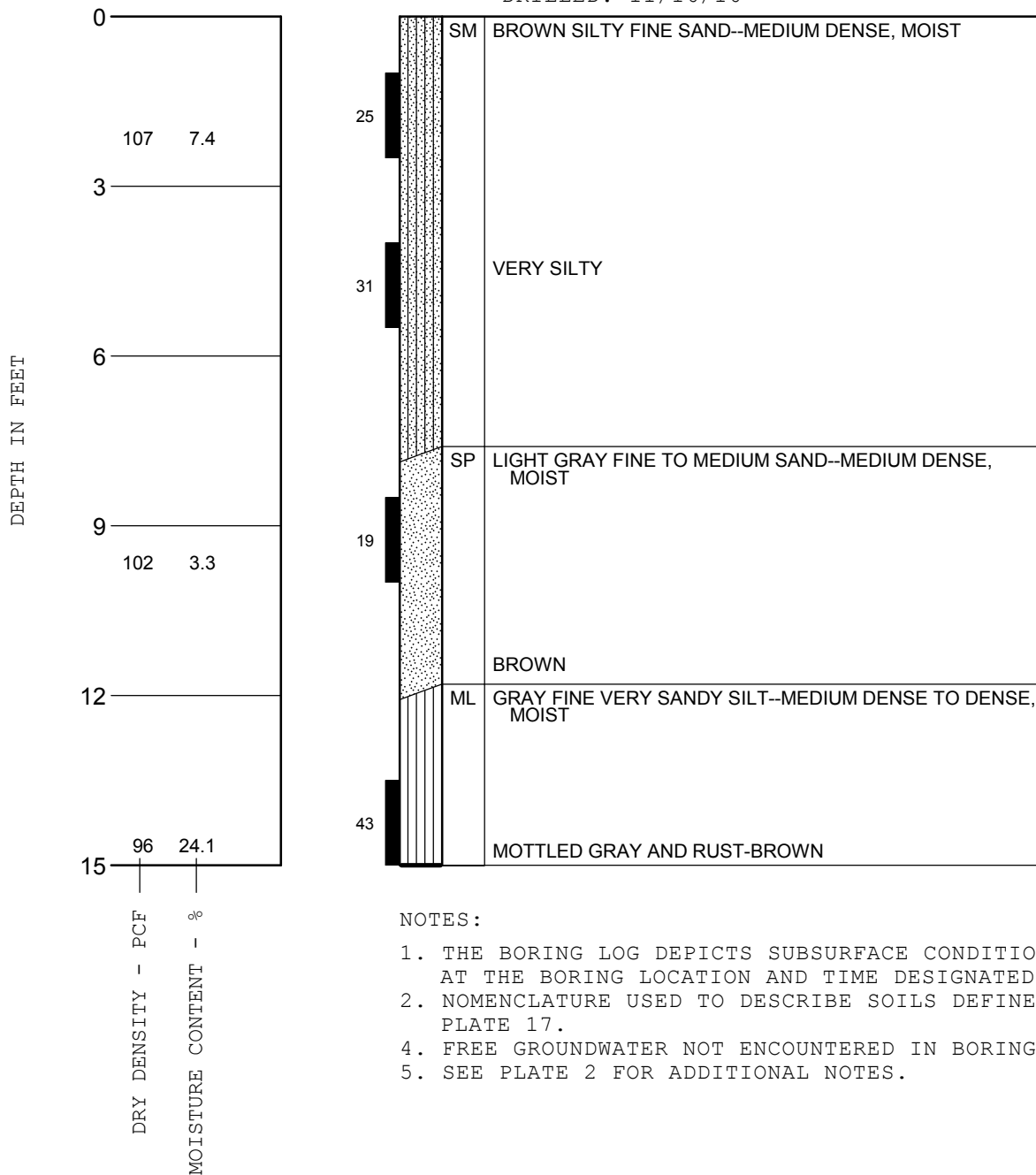
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 5

BORING 4

DRILLED: 11/16/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
4. FREE GROUNDWATER NOT ENCOUNTERED IN BORING 4.
5. SEE PLATE 2 FOR ADDITIONAL NOTES.

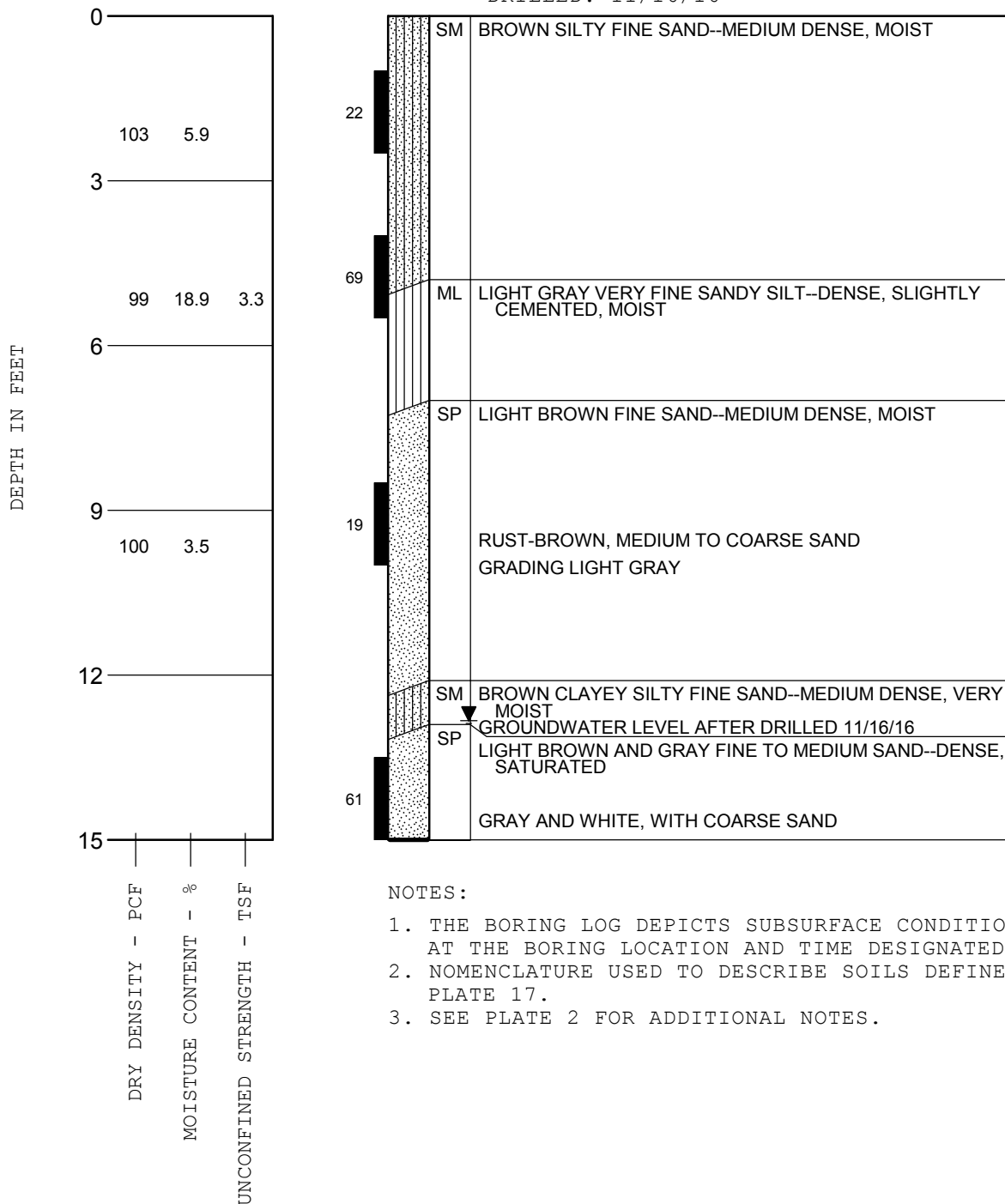
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 6

BORING 5

DRILLED: 11/16/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
3. SEE PLATE 2 FOR ADDITIONAL NOTES.

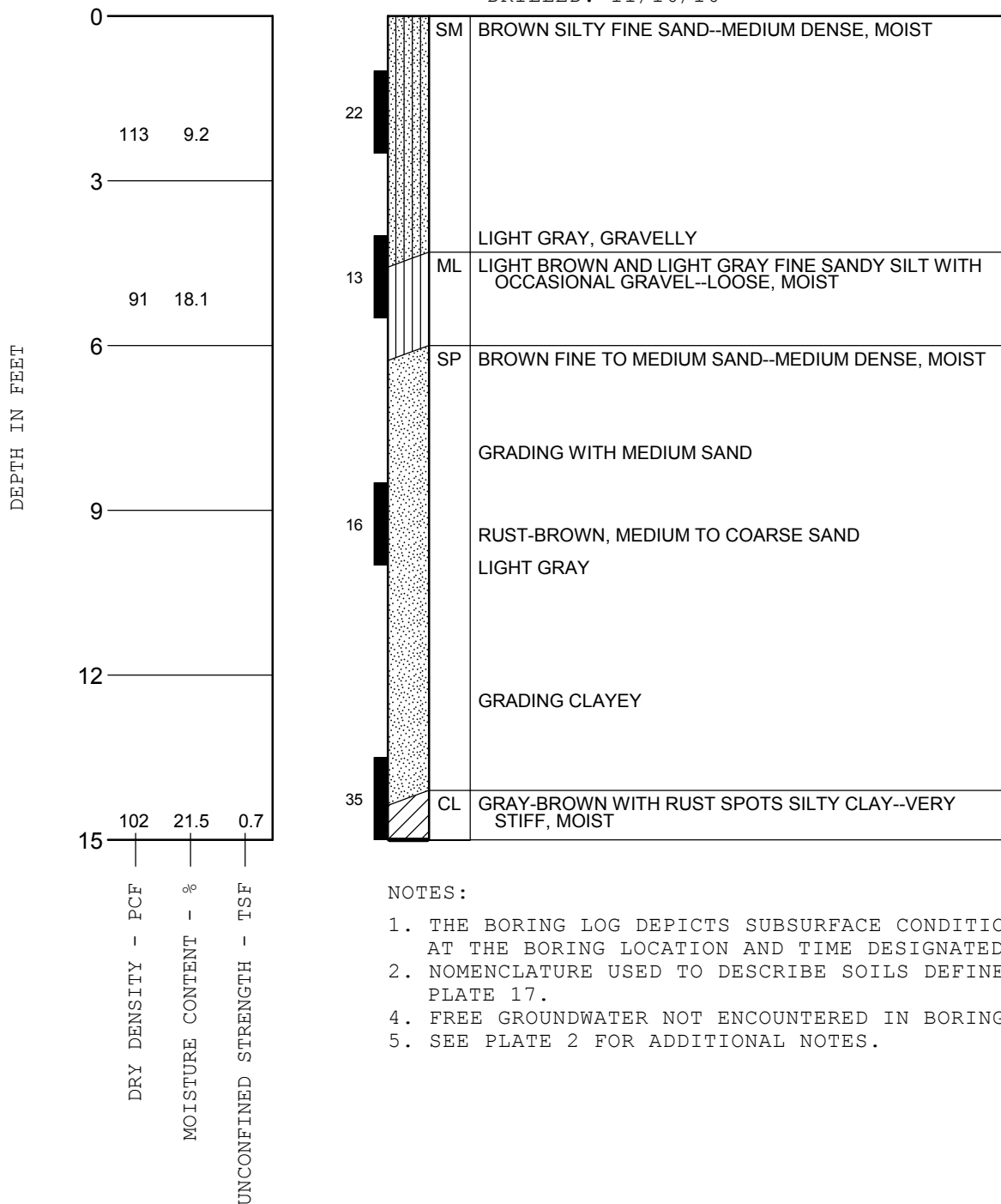
LOG OF BORING



PROJECT NUMBER: 146-618
DRAWN BY: WCB
DATE: 12/8/16
PLATE NUMBER: 7

BORING 6

DRILLED: 11/16/16



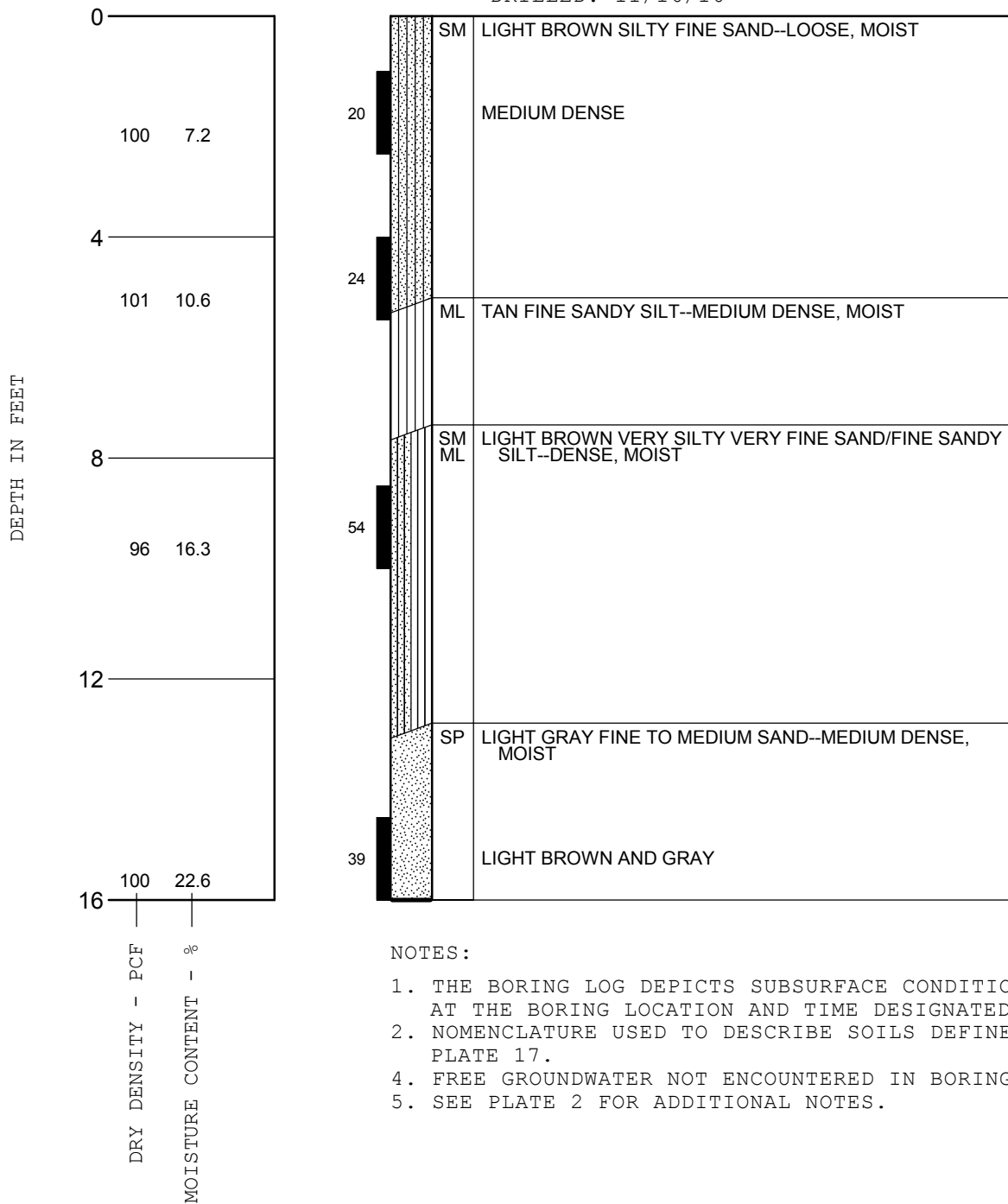
LOG OF BORING



PROJECT NUMBER: 146-618
DRAWN BY: WCB
DATE: 12/8/16
PLATE NUMBER: 8

BORING 7

DRILLED: 11/16/16



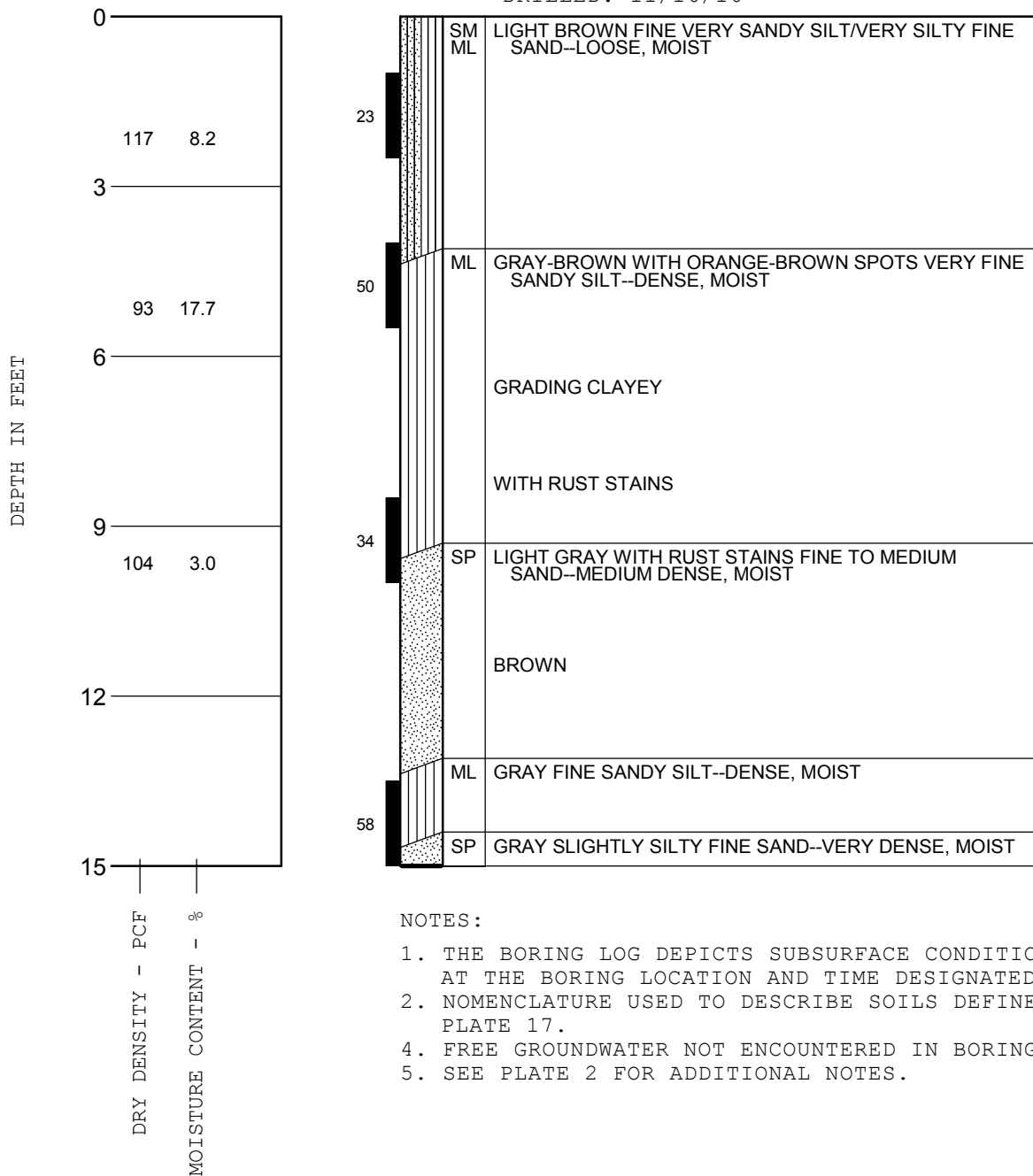
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 9

BORING 8

DRILLED: 11/16/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
4. FREE GROUNDWATER NOT ENCOUNTERED IN BORING 8.
5. SEE PLATE 2 FOR ADDITIONAL NOTES.

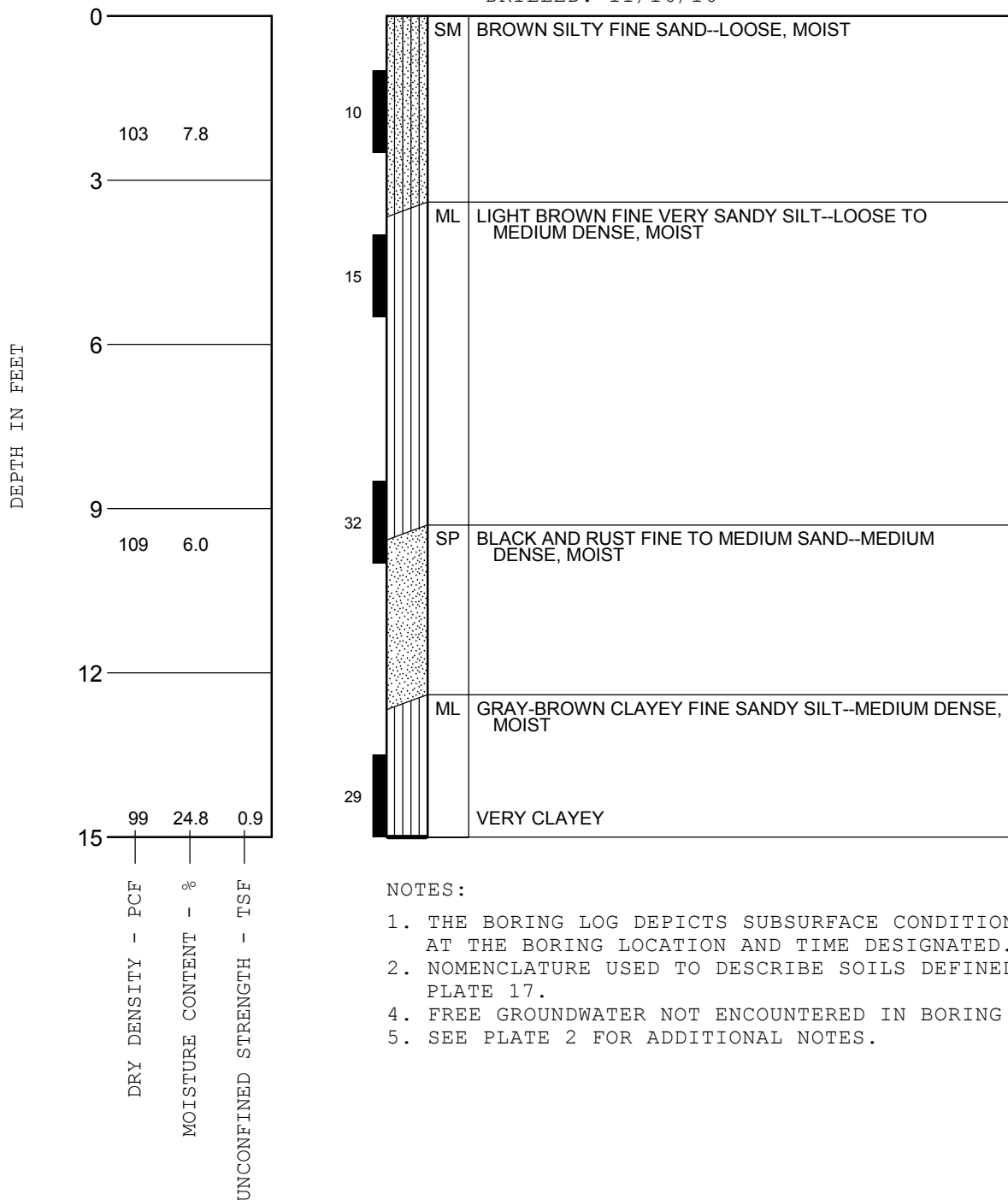
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 10

BORING 9

DRILLED: 11/16/16



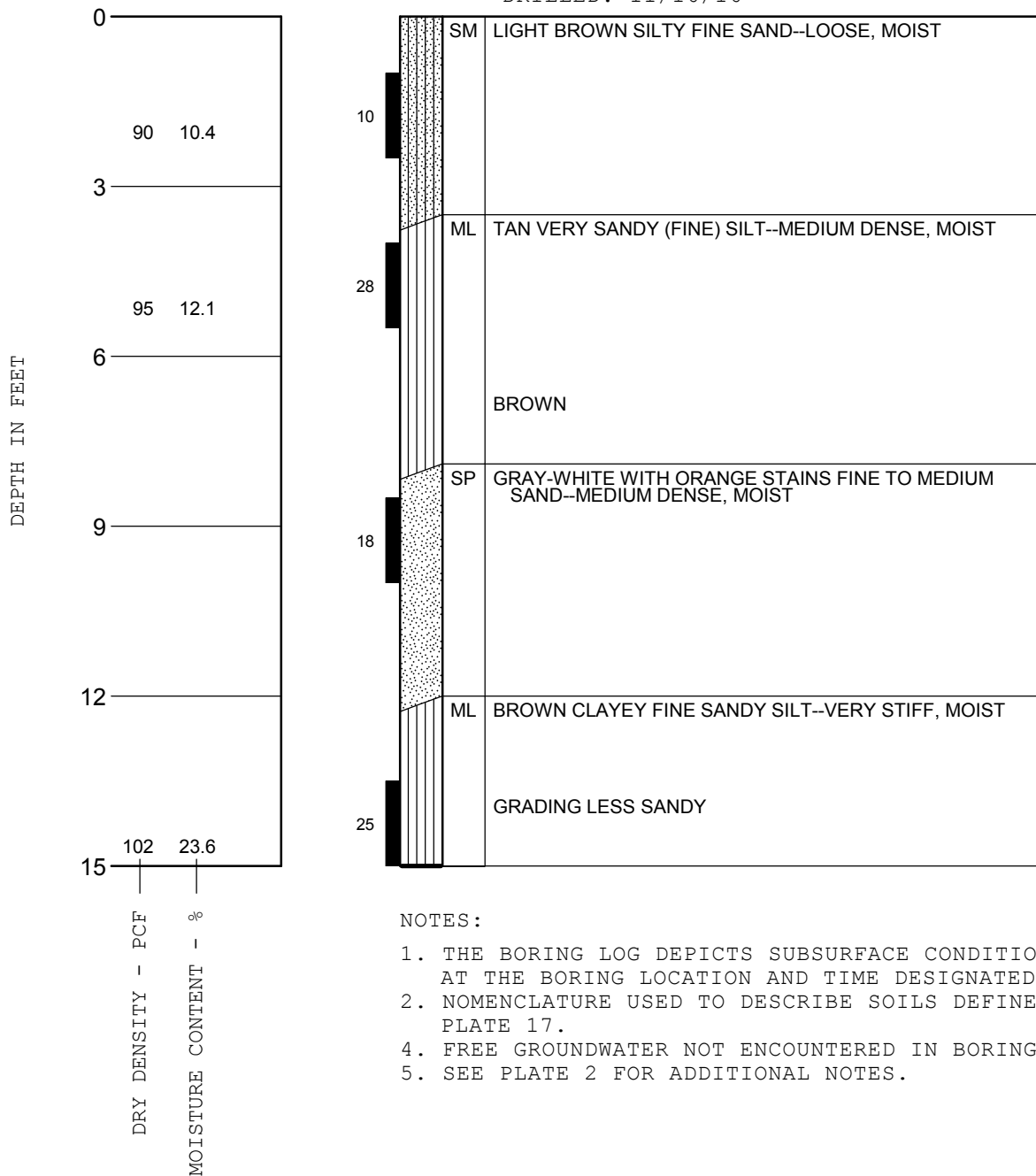
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 11

BORING 10

DRILLED: 11/16/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
4. FREE GROUNDWATER NOT ENCOUNTERED IN BORING 10.
5. SEE PLATE 2 FOR ADDITIONAL NOTES.

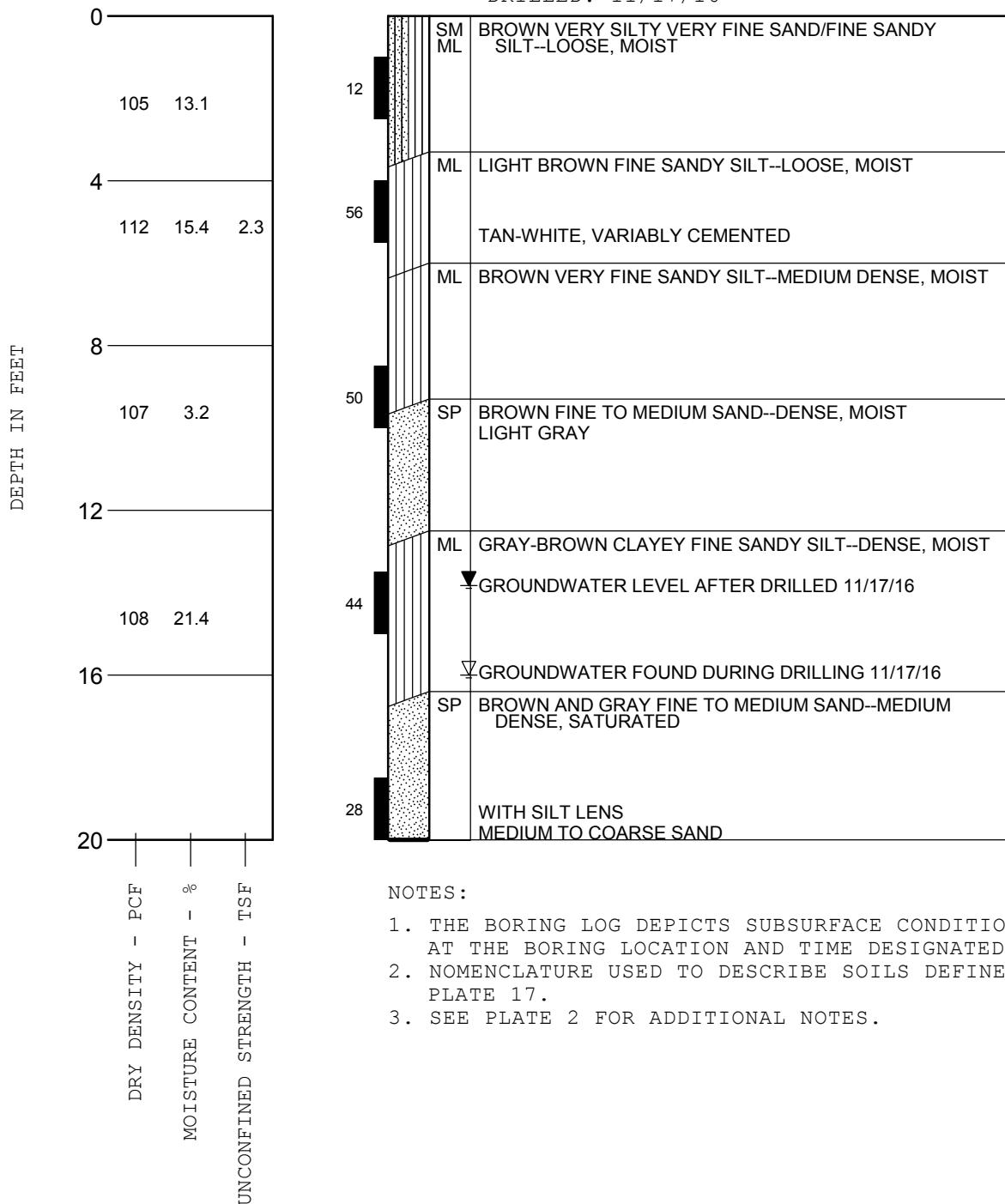
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 12

BORING 11

DRILLED: 11/17/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
3. SEE PLATE 2 FOR ADDITIONAL NOTES.

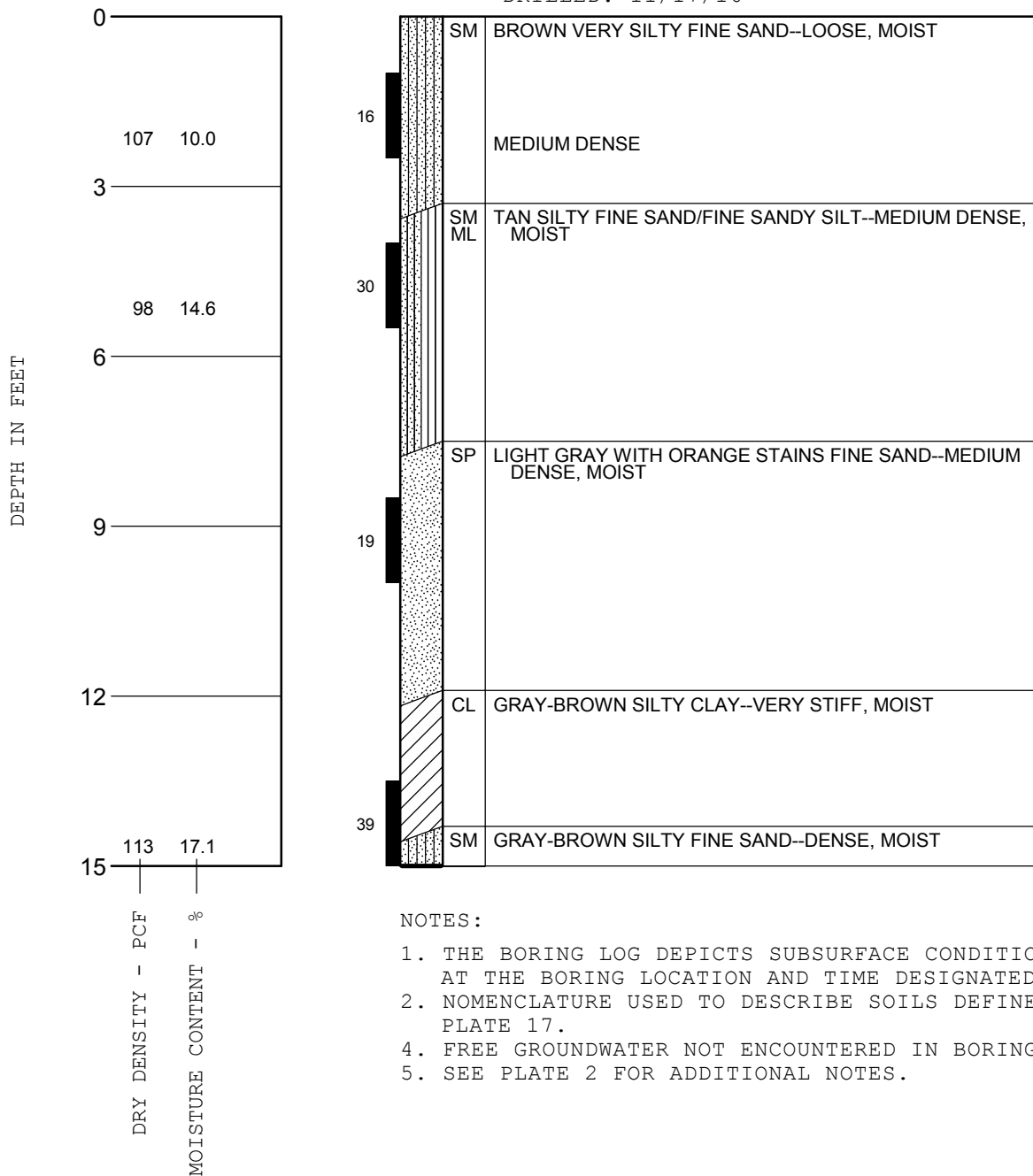
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 13

BORING 12

DRILLED: 11/17/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
4. FREE GROUNDWATER NOT ENCOUNTERED IN BORING 12.
5. SEE PLATE 2 FOR ADDITIONAL NOTES.

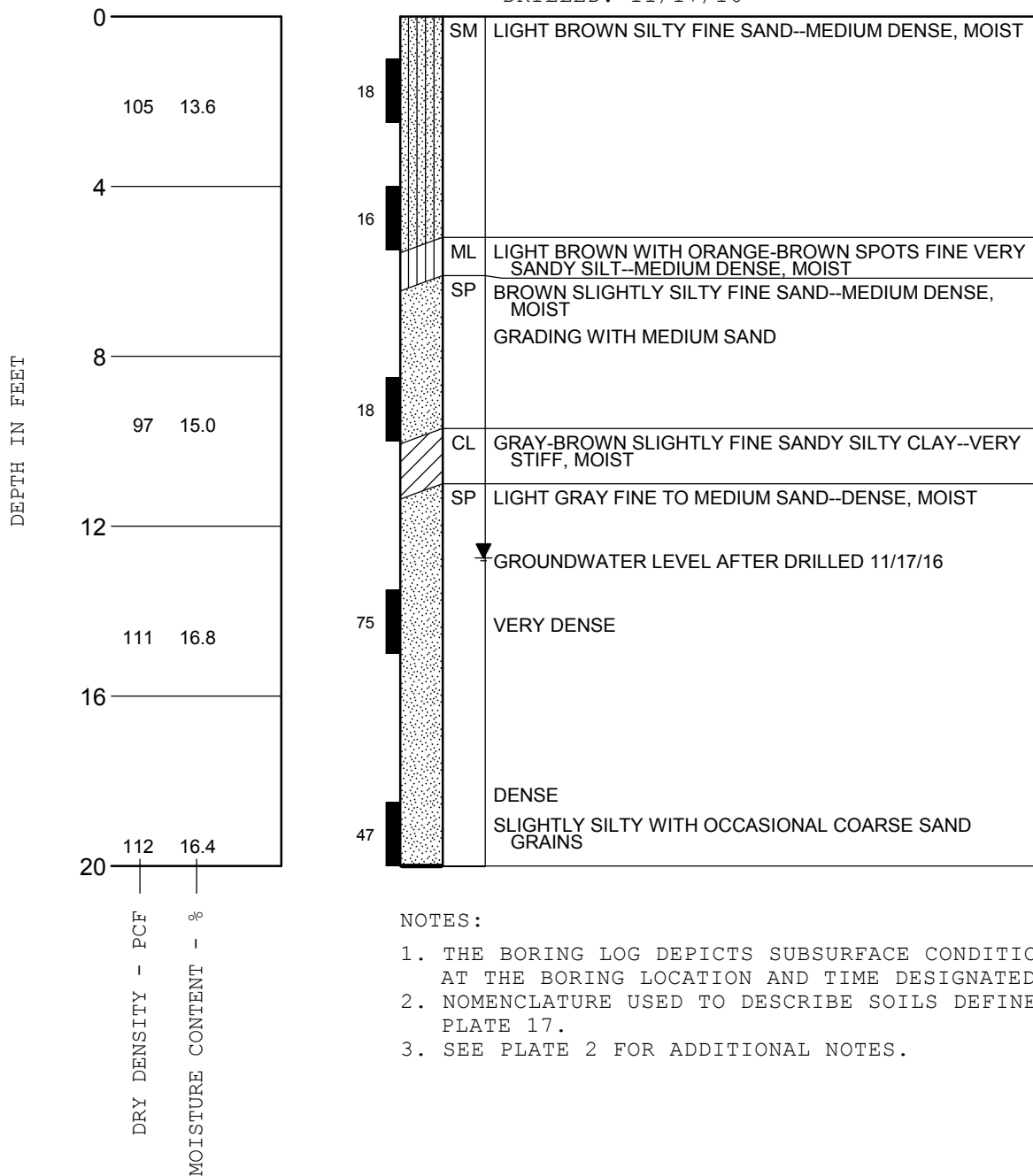
LOG OF BORING



PROJECT NUMBER: 146-618
 DRAWN BY: WCB
 DATE: 12/8/16
 PLATE NUMBER: 14

BORING 13

DRILLED: 11/17/16



NOTES:

1. THE BORING LOG DEPICTS SUBSURFACE CONDITIONS ONLY AT THE BORING LOCATION AND TIME DESIGNATED.
2. NOMENCLATURE USED TO DESCRIBE SOILS DEFINED ON PLATE 17.
3. SEE PLATE 2 FOR ADDITIONAL NOTES.

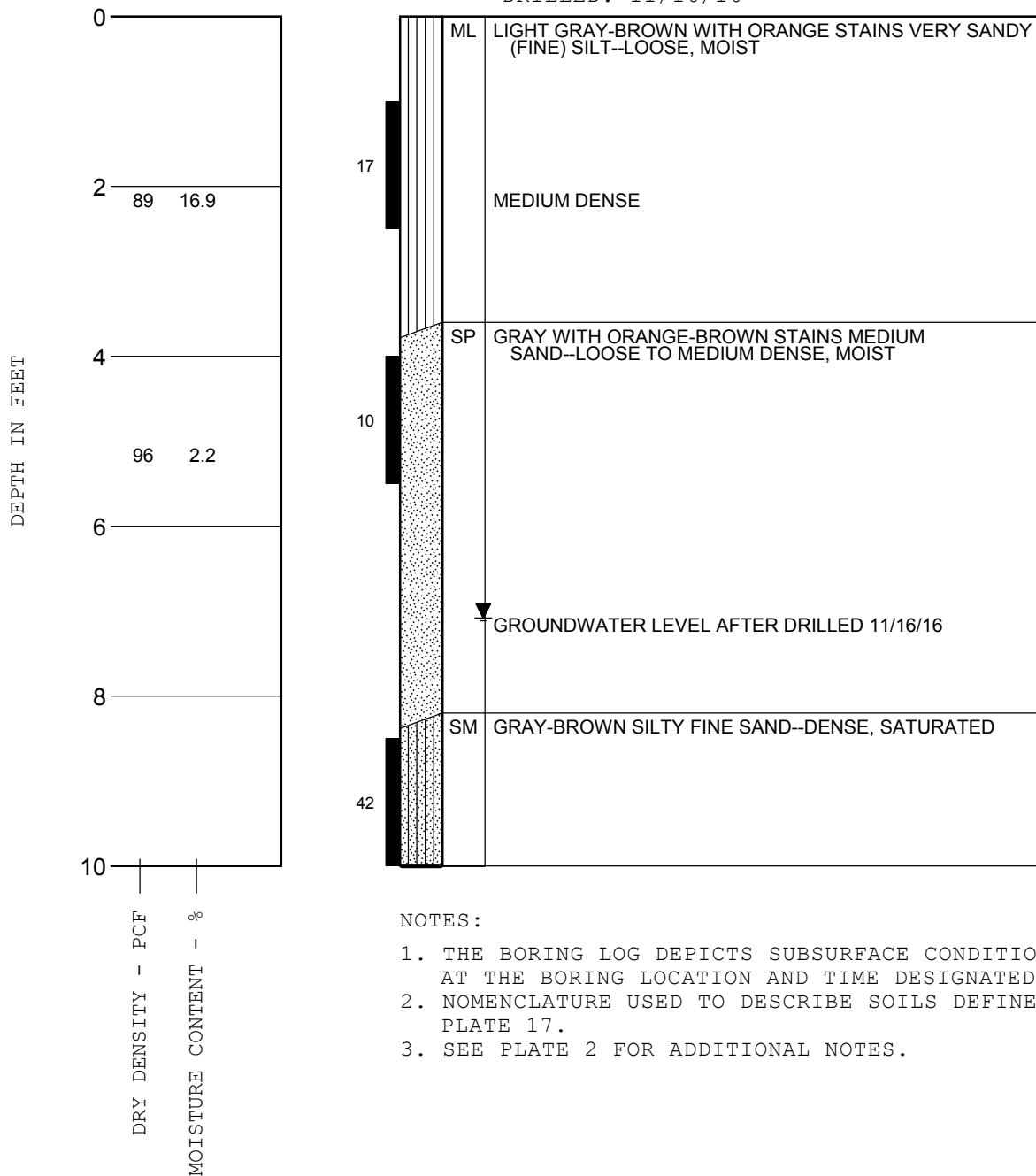
LOG OF BORING



PROJECT NUMBER: 146-618
DRAWN BY: WCB
DATE: 12/8/16
PLATE NUMBER: 15

BORING 14

DRILLED: 11/16/16



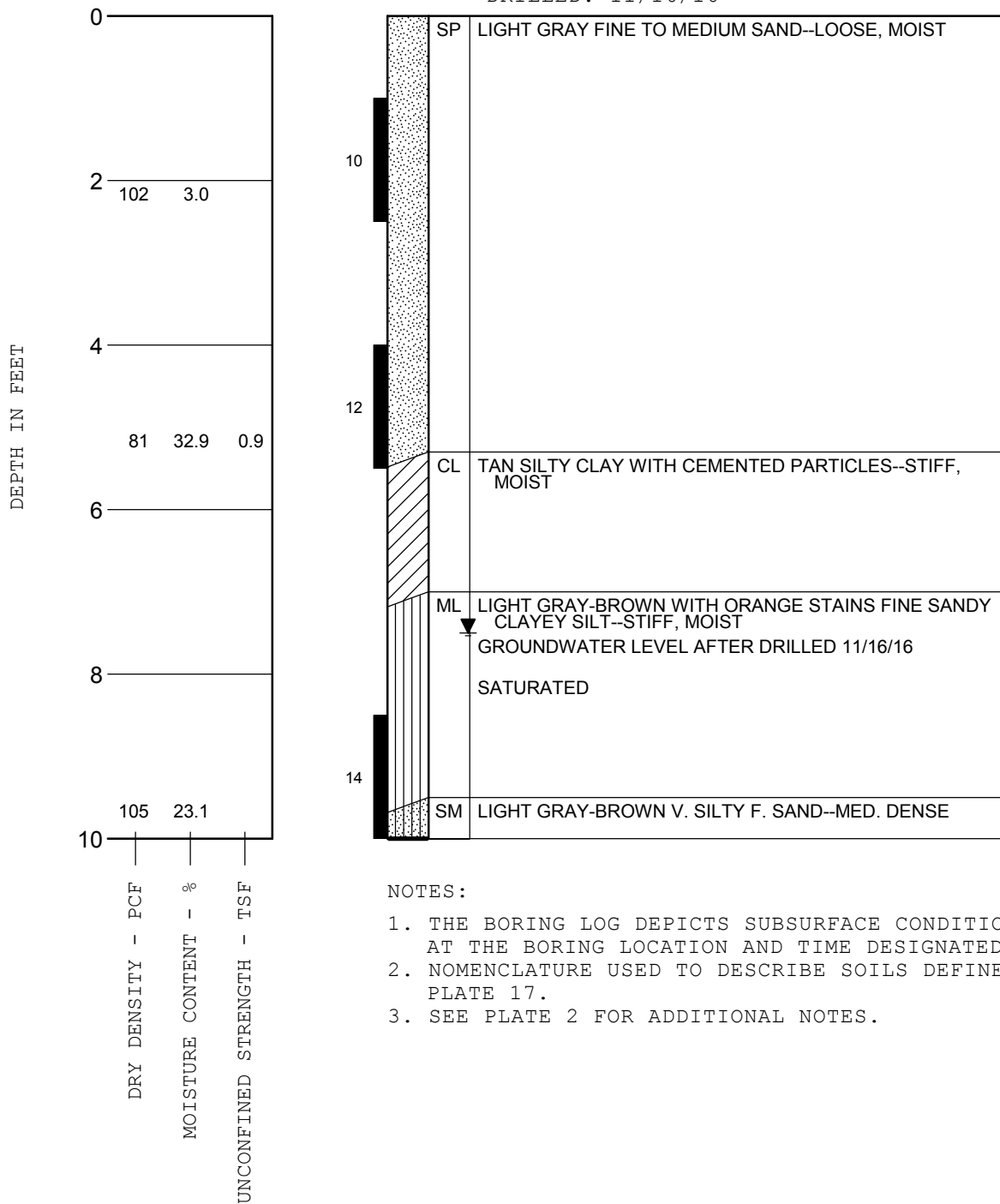
LOG OF BORING



PROJECT NUMBER: 146-618
DRAWN BY: WCB
DATE: 12/8/16
PLATE NUMBER: 16

BORING 15

DRILLED: 11/16/16



LOG OF BORING



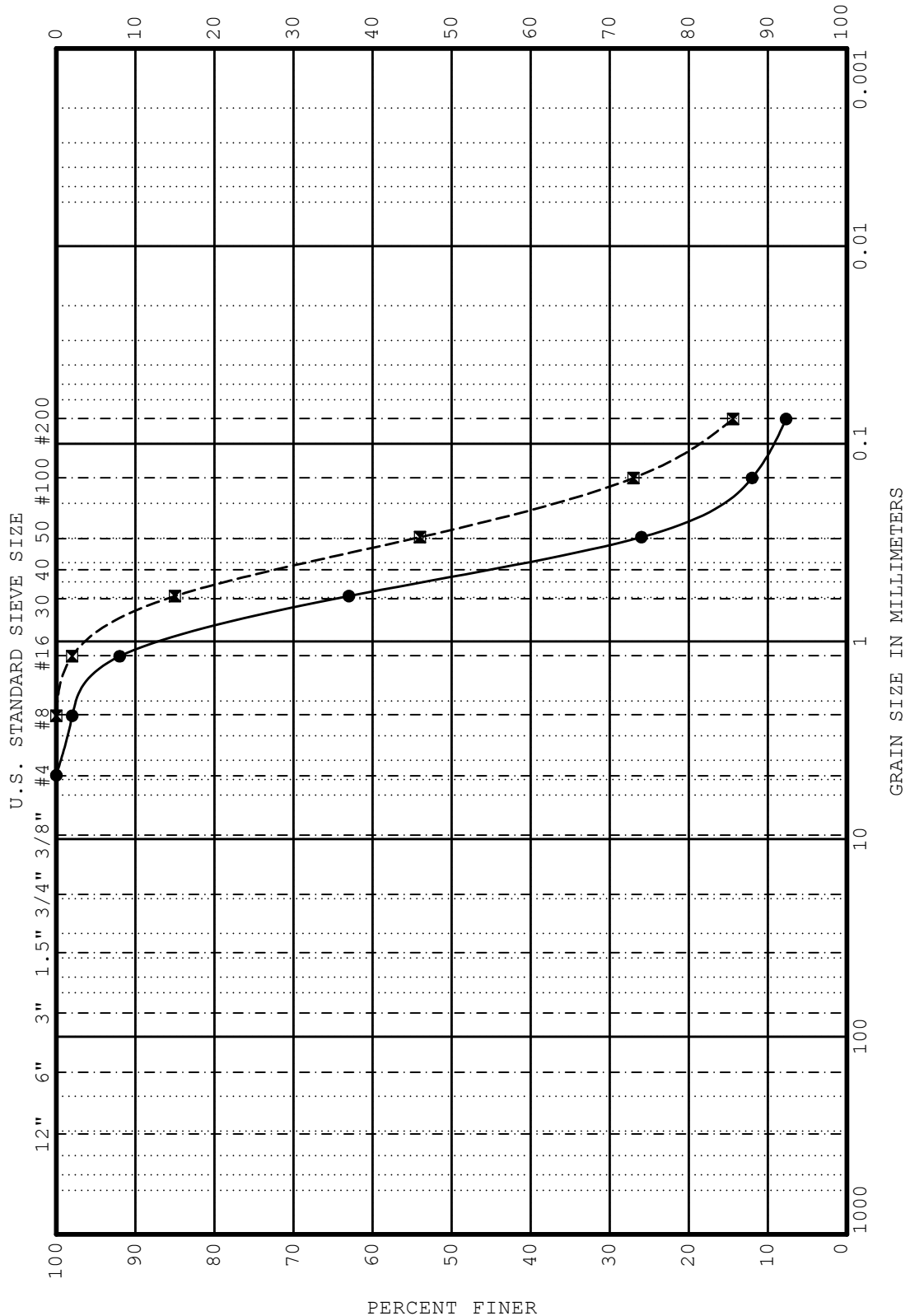
GRAPH	SYMBOL	DESCRIPTION	MAJOR DIVISIONS			
	GW	WELL GRADED GRAVELS, GRAVEL-SAND MIXTURES	CLEAN GRAVELS WITH LESS THAN 5% FINES	GRAVEL AND GRAVELLY SOILS	COARSE GRAINED SOILS MORE THAN 50% <u>LARGER</u> THAN NO. 200 SIEVE	
	GP	POORLY GRADED GRAVELS, GRAVEL-SAND MIXTURES				
	GM	SILTY GRAVELS, GRAVEL-SAND-SILT MIXTURES	GRAVELS WITH MORE THAN 12% FINES	MORE THAN 50% OF COARSE FRACTION <u>RETAINED</u> ON NO. 4 SIEVE		
	GC	CLAYEY GRAVELS, GRAVEL-SAND-CLAY MIXTURES				
	SW	WELL GRADED SANDS, GRAVELLY SANDS	CLEAN SANDS WITH LESS THAN 5% FINES	SANDS AND SANDY SOILS		
	SP	POORLY GRADED SANDS, GRAVELLY SANDS				
	SM	SILTY SANDS, SAND-SILT MIXTURES	SANDS WITH MORE THAN 12% FINES	MORE THAN 50% OF COARSE FRACTION <u>PASSING</u> NO. 4 SIEVE		
	SC	CLAYEY SANDS, SAND-CLAY MIXTURES				
	ML	INORGANIC SILTS, ROCK FLOUR, OR CLAYEY SILTS WITH SLIGHT PLASTICITY	LIQUID LIMIT <u>LESS</u> THAN 50		SILTS AND CLAYS	
	CL	INORGANIC CLAYS OF LOW TO MEDIUM PLASTICITY, GRAVELLY CLAYS, SANDY CLAYS, SILTY CLAYS, LEAN CLAYS				
	OL	ORGANIC SILTS AND ORGANIC SILTY CLAYS OF LOW PLASTICITY				
		MH	INORGANIC SILTS, MICACEOUS OR DIATOMACEOUS SILTS, ELASTIC SILTS	LIQUID LIMIT <u>GREATER</u> THAN 50		SILTS AND CLAYS
		CH	INORGANIC CLAYS OF HIGH PLASTICITY, FAT CLAYS			
		OH	ORGANIC CLAYS AND ORGANIC SILTS OF MEDIUM TO HIGH PLASTICITY			
	PT	PEAT, HUMUS, SWAMP SOILS WITH HIGH ORGANIC CONTENT	HIGHLY ORGANIC SOILS			

UNIFIED SOIL CLASSIFICATION SYSTEM

18

The graph plots 'PERCENT FINER' (Y-axis, 0 to 100) against 'GRAIN SIZE IN MILLIMETERS' (X-axis, logarithmic scale from 1000 to 0.001). Two curves are shown: a solid line with circular markers and a dashed line with square markers. The dashed line represents a finer material than the solid line.

Grain Size (mm)	Percent Finer (Solid Line)	Percent Finer (Dashed Line)
1000	100	100
600	100	100
300	100	100
150	100	100
75	100	100
37.5	100	100
19	100	100
10	100	100
4.75	95	95
2.5	85	90
1.18	75	85
0.6	65	75
0.3	55	65
0.15	45	55
0.075	35	45
0.0425	25	35
0.025	15	25
0.015	5	15
0.0075	0	5
0.00425	0	0



BOULDERS	COBBLES	GRAVEL			SAND		SILT	CLAY
		COARSE	FINE	COARSE	MEDIUM	FINE		

GRAIN SIZE DISTRIBUTION

DIRECT SHEAR DATA

TEST DESIGNATION: UNCONSOLIDATED, QUICK

STRAIN RATE: 0.030 inches/minute

SAMPLE CONDITION: Undisturbed

MATERIAL DESCRIPTION	SAMPLE I.D.	MOISTURE CONTENT (%)	DRY DENSITY (pcf)	NORMAL STRESS (psf)	SHEAR STRESS (psf)	
					3% Strain	10% Strain
Light brown silty fine to medium sand	Boring 3 1.6'	7.4	100	1000	720	960
	Boring 3 1.8'	5.4	104	2000	1200	1820
	Boring 3 2.3'	5.2	106	4000	2270	2960
Light brown silty fine sand	Boring 10 1.9'	11.2	95	1000	930	1100
	Boring 10 1.5'	12.7	84	2000	1510	1820
	Boring 10 1.7'	10.4	90	4000	3780	5910

DIRECT SHEAR DATA

**RESISTANCE VALUE TEST
CALIFORNIA TEST METHOD 301G**

SAMPLE LOCATION: SAMPLE S2
DEPTH: 12"-24"
MATERIAL DESCRIPTION: Dark gray fine very sandy silt

TEST NUMBER	DRY DENSITY (PCF)	MOISTURE CONTENT (%)	EXUDATION PRESSURE (PSI)	EXPANSION PRESSURE (PSF)	RESISTANCE VALUE
1	118	11.8	104	17	60
2	118	11.3	252	26	64
3	119	10.8	427	43	66

Resistance value at 300 psi exudation pressure = 65

RESISTANCE VALUE DATA

**RESISTANCE VALUE TEST
CALIFORNIA TEST METHOD 301G**

SAMPLE LOCATION: SAMPLE S3
DEPTH: 6"-18"
MATERIAL DESCRIPTION: Gray-brown silty fine sand

TEST NUMBER	DRY DENSITY (PCF)	MOISTURE CONTENT (%)	EXUDATION PRESSURE (PSI)	EXPANSION PRESSURE (PSF)	RESISTANCE VALUE
1	115	12.4	88	0	62
2	115	11.4	215	0	63
3	114	10.4	646	0	66

Resistance value at 300 psi exudation pressure = 64

RESISTANCE VALUE DATA

CALIFORNIA TEST METHOD NO. 373

Sample Location: S5

Depth: 6"-18"

Material Description: Brown silty fine sand

Additives: Portland Cement as shown

Sample No.	Additive Amounts	Dry Density (pcf)	Moisture Content (%)	Maximum Dry Density/ Optimum Moisture Content	Unconfined Compressive Strength
A	4% cement	115.0	8.3	118 pcf @ 10.4%	357 psi
B	4% cement	115.0	9.4		328 psi
C	4% cement	117.8	10.4		414 psi
D	4% cement	117.4	11.5		329 psi
E	4% cement	117.7	10.4		363 psi
F	6% cement	120.6	10.4		605 psi
G	6% cement	120.5	10.4		565 psi

TREATED SOIL COMPRESSIVE STRENGTH DATA



Sunland Analytical

11419 Sunrise Gold Circle, #10
Rancho Cordova, CA 95742
(916) 852-8557

Date Reported 12/02/2016
Date Submitted 11/28/2016

To: William Boli
Raney Geotechnical
3140 Beacon Blvd. Ste#100
West Sacramento, CA 95691

From: Gene Oliphant, Ph.D. \ Randy Horney
General Manager \ Lab Manager

The reported analysis was requested for the following location:
Location : 146-618 MURPHY PKWAY Site ID : B4 @ 4-5.5FT.
Thank you for your business.

* For future reference to this analysis please use SUN # 73282-152961.

EVALUATION FOR SOIL CORROSION

Soil pH	7.80		
Moisture	10.9	%	
Minimum Resistivity	1.53	ohm-cm (x1000)	
Chloride	47.7	ppm	00.00477 %
Sulfate	125.6	ppm	00.01256 %
Redox Potential	(+) 192	mv	
Sulfides		Présence -	NEGATIVE

METHODS

pH and Min.Resistivity CA DOT Test #643 Mod.(Sm.Cell)
Sulfate CA DOT Test #417, Chloride CA DOT Test #422
Redox Potential ASTM G-200, Sulfides AWWA C105/A25.5



Sunland Analytical

11419 Sunrise Gold Circle, #10
Rancho Cordova, CA 95742
(916) 852-8557

Date Reported 12/02/2016
Date Submitted 11/28/2016

To: William Boli
Raney Geotechnical
3140 Beacon Blvd. Ste#100
West Sacramento, CA 95691

From: Gene Oliphant, Ph.D. \ Randy Horney
General Manager \ Lab Manager

The reported analysis was requested for the following location:
Location : 146-618 MURPHY PKWAY Site ID : B9 @ 4-5.5FT.
Thank you for your business.

* For future reference to this analysis please use SUN # 73282-152962.

EVALUATION FOR SOIL CORROSION

Soil pH	7.95		
Moisture	14.2	%	
Minimum Resistivity	1.66	ohm-cm (x1000)	
Chloride	29.4 ppm	00.00294	%
Sulfate	78.4 ppm	00.00784	%
Redox Potential	(+) 210	mv	
Sulfides	Presence - NEGATIVE		

METHODS

pH and Min.Resistivity CA DOT Test #643 Mod.(Sm.Cell)
Sulfate CA DOT Test #417, Chloride CA DOT Test #422
Redox Potential ASTM G-200, Sulfides AWWA C105/A25.5

LIQUEFACTION ANALYSIS

Murphy Parkway Warehouse

Hole No.=B-1 Water Depth=13.8 ft
Ground Improvement of Fill=1 ft

Magnitude=6.6
Acceleration=0.43g

