



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 7
25 FUNSTON ROAD
KANSAS CITY, KANSAS 66115

Date: JAN 31 1992

MEMORANDUM

SUBJECT: Data Transmittal for Activity #: IS651
Site Description: Syntex Facility

FROM: Andrea Jirka *H*
Chief, Laboratory Branch, ENSV

TO: Robert Morby
Chief, Superfund Branch, WSTM

ATTN: Mark Bazina

Attached is the data transmittal for the above referenced site. This should be considered a _____ Partial or X Complete data transmittal (completes transmittal of _____). If you have have any questions or comments, please contact Dee Simmons at 551-5129.

Attachments

cc: Data Files

NOTE: Please see Mary Gerken, SPFD-WSTM, if you want an electronic copy of the data.



DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, MO 64115

EPA ACNO: 18651 SAMNO: 001 QCC: MEDIA: WATER PL: S P F D

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE:
LOCATION: VERONA MO PROJECT NUM: A41 PT: LONGITUDE:

SAMPLE DES: Groundwater well MW-1
LOCATION: VERONA MO REG: 10/31/91 TIME: 9:00 FROM REF PT:
CASE/BATCH/SNO: / / LAB: END: 10/31/91 9:15 EAST:
STORET/SARDAD NO: DOWN:

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MCP	NAME
2 VOA VIALS	ICED	UV	VOLATILES
CURT	5 ML HNO3	UM	METALS <i>+Hg(w/34)</i>
GLASS	ICED	US	SEMIVOLATILES
GLASS	ICED	UP	PESTICIDES

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: OPERABLE UNIT:

Sample Collected from well MW-1

- Temp. - 14.2°*
- Pin - 6.50'*
- Cond. - 1040*

SAMPLE COLLECTED BY: *M. J. Soydas*

DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 64115

EY: 92 ACTNO: 19451 SAMNO: 002 QCC: MEDIA: WATER PL: S P F D

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE:
LOCATION: VERONA NO PROJECT NUM: 641 PT: LONGITUDE:

SAMPLE DES: Groundwater well MW-5
LOCATION: VERONA NO: BEG: 10/31/91 TIME: 9:30 FROM REF PT:
CASE/BATCH/SNO: / / LAB: END: 10/31/91 9:50 EAST:
SIURET/SARUAD NO: DOWN:

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	QCP	NAME
2 VOA VIALS	ICED	UV	VOLATILES ✓
UBT	5 ML HNO3	UM	METALS <i>with (unseen)</i>
GLASS	ICED	US	SEMIVOLATILES
GLASS	ICED	UP	PESTICIDES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: OPERABLE UNIT:

Sample collected from Well MW-5

- Temp 16.8*
- Ph 6.88, 6.90, 6.88*
- Concl. 480, 460, 420*

SAMPLE COLLECTED BY: Michael J. Boydas

00001

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII

ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 64115

EY: 92 ACTNO: 18451 SAMNO: 003 QCD: MEDIA: WATER PI: S P F D

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: _____

LOCATION: VERONA MO PROJECT NUM: A41 PC: LONGITUDE: _____

SAMPLE DES: Groundwater Well MW-11 DATE TIME FROM REF PT

LOCATION: VERONA MO REF: 10/31/91 10:05 EAST: _____

CASE/BATCH/SMD: / / / LAB: END: 10/31/91 10:30 NORTH: _____

STORET/SAROAD NO: DOWN: _____

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
2 VOA VIALS	ICED	UV	VOLATILES ✓
1 CRT	5 MI HNO3	WM	METALS <u>Heavy metals</u> ✓
GLASS	ICED	WS	SEMIVOLATILES ✓
GLASS	ICED	WP	PESTICIDES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: _____ OPERABLE UNIT: _____

Sample Collected from Well MW-11

SAMPLE COLLECTED BY: Michael J Hayes

DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 64115

FY: 92 ACTNO: 19651 SAMNO: 004 UCC: MEDIA: WATER PL: S P F 0

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: _____

LOCATION: VERONA MO PROJECT NUM: A41 PT: LONGITUDE: _____

SAMPLE DES: Well No ~~WA~~ MW-16
LOCATION: VERONA MO
CASE/BATCH/SNO: _____ LAB: _____
STORET/SAROAD NO: _____
DATE: 10/31/91 TIME: 10:50 FROM REF PT: _____
BEG: 10/31/91 EAST: _____
END: 10/31/91 TIME: 11:00 NORTH: _____
DOWN: _____

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MVP	NAME
2 VOA VIALS	ICED	UV	VOLATILES ✓
CUBI	5 ML HNO3	UM	METALS <u>4/13/91 (2/13/91)</u>
GLASS	ICED	US	SEMI-VOLATILES ✓
GLASS	ICED	WP	PESTICIDES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: _____ OPERABLE UNIT: _____

Sample Collected from well ~~H6~~ 16

- pH 7.30, 7.40, 7.46, 7.46
- T. 15.9°C
- Cond. \ 290, 280, 260

SAMPLE COLLECTED BY: Michael J. Boydas

DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII

ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

EY: 92 ACINO: 18651 SAMNO: 005 QCC: MEDIA: WATER PL: S P F D

ACTIVITY DES: SYNTEX FACILITY

REF LATITUDE: ---

LOCATION: VERONA

MO: PROJECT NUM: A41

PT: LONGITUDE: ---

SAMPLE DES: Mw-16B

LOCATION: VERONA

MO

BEG: 10/21/91

TIME

FROM REF PT

EAST: ---

CASE/BATCH/SMD: / /

LAB: ---

END: 10/21/91

11:10

NORTH: ---

STORET/SAROAD NO: ---

DOWN: ---

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MCP	NAME
2 VOA VIALS	ICED	UV	VOLATILES ✓
CUBT	5 ML HNO3	WM	METALS ✓
GLASS	ICED	WS	SEMIVOLATILES ✓
GLASS	ICED	WP	PESTICIDES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: --- OPERABLE UNIT: ---

Sample collected from MW-16B

Ph 7.84, 7.88, 7.82, 7.78

T. 15.9°C

Cond. 290, 280

SAMPLE COLLECTED BY: Michael J. Goydas

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FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

EPA ID: 92 ACTNO: 18651 SAMNO: 906 QDC: MEDIA: WATER PL: S P F D

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: ---
LOCATION: VERONA MO PROJECT NUM: A41 PT: LONGITUDE: ---

SAMPLE DES: MW-9
LOCATION: VERONA MO REG: 10/31/91 TIME: 11:00 FROM REF PT: ---
CASE/BATCH/SNO: / / LAB: FND: 10/31/91 TIME: 11:10 NORTH: ---
STORET/SAROAD NO: DOWN: ---

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MCP	NAME
2 VOA VIALS	ICED	WP	VOLATILES
CUBT	5 ML HNO3	WM	METALS (MS/UM34)
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: --- OPERABLE UNIT: ---

Sample Collected from MW-9
Ph 6.88, 6.90, 6.92
T. 15.3 °C
Cond. 1780, 770, 760

SAMPLE COLLECTED BY: Michael J. Hayes

EY: 92 ACTNO: 18451 SAMNO: 007 QCC: _ MEDIA: WATER PL: S P F D
ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: _ _ _
LOCATION: VERONA MO PROJECT NUM: A41 PT: LONGITUDE: _ _ _
SAMPLE DES: Groundwater well MW 15 10/31/91 TIME FROM REF PT
LOCATION: VERONA MO REG: 10/31/91 11:35 EAST: _ _ _
CASE/BATCH/SMD: _ / _ / _ LAB: _ END: 10/31/91 12:05 NORTH: _ _ _
STORET/SARDAD NO: _ _ _ DOWN: _ _ _

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
2 VOA VIALS	ICED	UV	VOLATILES
CUBI	5 MI HNO3	UM	METALS <u>(Mn, Zn, Pb)</u>
GLASS	ICED	MS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: _ _ _ OPERABLE UNIT: _ _ _

Sample Collected from MW-15

Ph 7.41, 7.38
T. 13.9°C
Concl 1520, 540, 480

SAMPLE COLLECTED BY: Michael J. Boydas

DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

EY: 92 ACTNO: 05451 SAMNO: 008 QCC: _ MEDIA: WATER PL: S P F 0

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: _ _ _
LOCATION: VERONA MO PROJECT NUM: A41 PT: LONGITUDE: _ _ _

SAMPLE DES: Groundwater Well MW 15-B
LOCATION: VERONA MO DATE: 10/31/91 TIME: 11:35 FROM REF PT: _____
CASE/BATCH/SNO: _____/_____/_____ LAB: _____ BEG: 10/31/91 END: 10/31/91 NORTH: _____
STORET/SAROAD NO: _____ DOWN: _____

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
2 VOA VIALS	ICED	UV	VOLATILES ✓
CUBT	5 ML HNO3	NM	METALS <u>TAS (WASZ)</u>
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: _____ OPERABLE UNIT: _____

Sample Collected from MW-15B

pH 6.62, 6.69, 6.88, 6.69

T. 15.9°C

Cond. 1200, 1400, 1200

SAMPLE COLLECTED BY: Michael J. Soydas

DRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

ACTNO: 92 ACTNO: JS651 SAMNO: 009 QCC: F MEDIA: WATER PI: S P F D

ACTIVITY DES: SYNTEX FACILITY REF LATITUDE: _____
LOCATION: VERONA NO PROJECT NUM: A41 PT: LONGITUDE: _____

SAMPLE DES: TRIP BLANK
LOCATION: VERONA NO ^{10/31/91} TIME FROM REF PT
CASE/BATCH/SMD: _____ / _____ / _____ LAB: _____ BEG: ~~10/27/91~~ ^{10/31/91} 8:00 EAST: _____
STORET/SAROAD NO: _____ END: 10/31/91 12:05 NORTH: _____
DOWN: _____

ANALYSIS REQUESTED:
CONTAINER PRESERVATIVE MGP NAME -
2 VOA VIALS ICED UV VOLATILES ✓

COMMENTS: FOR SUPERFUND ONLY: SUBSITE IDENTIFIER: _____ OPERABLE UNIT: _____

Trip F

SAMPLE COLLECTED BY : Michael J. Boydas

**CHAIN OF CUSTODY RECORD
ENVIRONMENTAL PROTECTION AGENCY REGION VII**

11/14/91
1 of 1

ACTIVITY LEADER(Print) <u>M. Goydas (MEE)</u> <u>M. Bogina (EPA.) M. Dodan (BEE)</u>	NAME OF SURVEY OR ACTIVITY <u>IS 651</u> <u>SYNTEX - Verona, Ground water</u>	DATE OF COLLECTION <u>91</u> <u>31</u> DAY <u>10</u> MONTH <u>91</u> YEAR	SHEET <u>1</u> of <u>1</u>
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SAMPLE NUMBER	TYPE OF CONTAINERS				SAMPLED MEDIA					RECEIVING LABORATORY REMARKS/OTHER INFORMATION (condition of samples upon receipt, other sample numbers, etc.)	
	1-LITER CUBITAINER	800Z BOTTLE	BOTTLE	BOTTLE	VOA SET (2 VIALS EA)	water	soil	sediment	dust		other
	NUMBERS OF CONTAINERS PER SAMPLE NUMBER										
IS 651 001	1	2			1	X					
002	1	2			1						
003	1	2			1						
004	1	2			1						
005	1	2			1						
006	1	2			1						
007	1	2			1						
008	1	2			1						
V 009 F					1						
<i>No other samples</i>											

DESCRIPTION OF SHIPMENT <u>33</u> PIECE(S) CONSISTING OF _____ BOX(ES) <u>4</u> ICE CHEST(S); OTHER _____	MODE OF SHIPMENT ____ COMMERCIAL CARRIER: _____ ____ COURIER <input checked="" type="checkbox"/> SAMPLER CONVEYED _____ (SHIPPING DOCUMENT NUMBER)
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PERSONNEL CUSTODY RECORD			
RELINQUISHED BY (SAMPLER) <u>Michael J. Goydas</u>	DATE <u>8/05/91</u> <u>11/1/91</u>	TIME <u>8:05am</u>	RECEIVED BY <u>[Signature]</u>
<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED			<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED
REASON FOR CHANGE OF CUSTODY <u>Reed @ Lab</u>			
RELINQUISHED BY	DATE	TIME	RECEIVED BY
<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED			<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED
REASON FOR CHANGE OF CUSTODY			
RELINQUISHED BY	DATE	TIME	RECEIVED BY
<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED			<input type="checkbox"/> SEALED <input type="checkbox"/> UNSEALED
REASON FOR CHANGE OF CUSTODY			

ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

ICF Technology

Mantech Environmental Technology, Inc.

ESAT Region VII
Mantech Env. Tech., Inc.
25 Funston Road
Kansas City, KS 66115
(913) 551-5000

TO: Barry Evans, Data Review Task Monitor/ENSV
THRU: Harold Brown, Ph.D., ESAT Contract Manager/ENSV
FROM: Rebecca K. Estep, ESAT Data Reviewer/Mantech *Re Re*
THRU: Ronald Ross, Manager/ESAT/Mantech
DATE: December 20, 1991
SUBJECT: Review of inorganic data for Syntex Facility.

TID#: 07-9109-016
ASSIGNMENT#: 067
ICF ACCT#: 045-05-016-02
Mantech S.O.#: 1110-016
ESAT Document#: ESAT-VII-016-5067

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," July 1, 1988 revision.

The following comments and attached data sheets are a result of Mantech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE NO.:	17360	LABORATORY:	DATA C
CONTRACT NO.:	-----	METHOD NO:	CS0390I
SITE:	Syntex Facility	ACTIVITY:	IS651
REVIEWER:	Rebecca K. Estep	MATRIX:	Water

Samples for Total Metals and Mercury (Hg) Analysis

<u>SMO SAMPLE NO.</u>	<u>EPA SAMPLE NO.</u>	<u>SMO SAMPLE NO.</u>	<u>EPA SAMPLE NO.</u>
MGH914	IS651982P	MGH920	IS651001
MGH921	IS651002	MGH922	IS651003
MGH923	IS651004	MGH924	IS651005
MGH925	IS651006	MGH926	IS651007
MGH927	IS651008		

6. DUPLICATES

A. Aluminum (Al), chromium (Cr), and copper (Cu) were outside quality control limit requirements in the duplicate analysis. Aluminum (Al) in samples IS651982P, IS651003, IS651005, IS651007, and IS651008 and copper (Cu) in samples IS651001 and IS651002 would have been "J" coded, however, these qualifications were overridden due to the blank rules. Aluminum (Al) in samples IS651001, IS651002, IS651006, IS651002S, and IS651901C, chromium (Cr) in samples IS651982P, IS651001, IS651002, IS651005, IS651006, IS651007, IS651002S, and IS651901C, and copper (Cu) in samples IS651901C and IS651002S were "J" coded due to the duplicate rules.

7. MATRIX SPIKE

A. A matrix spike was analyzed with arsenic (As) and selenium (Se) outside quality control limit requirements for percent recovery. Arsenic (As) in samples IS651982P, IS651001, IS651006, and IS651901C and selenium (Se) in samples IS651982P and IS651901C were "J" coded due to the matrix spike rules. Arsenic (As) in samples IS651002, IS651003, IS651004, IS651005, IS651007, IS651008, IS651002L, and IS651900M was invalidated due to the matrix spike rules.

8. GRAPHITE FURNACE ATOMIC ABSORPTION (GFAA) SPECTROSCOPY

A. Selenium (Se) and thallium (Tl) were outside quality control limit requirements for post digestion spike recoveries in several samples, however, since these analytes were non-detect in all associated samples, no data were qualified.

B. The method of standard additions was performed for lead (Pb) in sample IS651001 and the correlation coefficient was within quality control limit requirements, thus, no data were qualified.

9. PERFORMANCE EVALUATION SAMPLE

A. One performance evaluation audit sample was submitted to the laboratory for analysis associated with this case and SDG number with all of the audit compounds detected. Aluminum (Al) was detected additionally.

10. ICP SERIAL DILUTION

A. All analytes were within quality control limit requirements according to the ICP serial dilution rules.

11. COMPOUND IDENTIFICATION and QUANTITATION

A. In level 2 review, no calculations from raw data (quantitation reports) are performed.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

ICF Technology

Mantech Environmental Technology, Inc.

ESAT Region VII
Mantech Env. Tech., Inc.
25 Funston Road
Kansas City, KS 66115
(913) 551-5000

TO: Barry Evans, Data Review Task Monitor/ENSV
THRU: Harold Brown, Ph.D., Regional Project Officer/ENSV

FROM: Carla S. Phelps, ESAT Data Reviewer/Mantech CP R
THRU: Ronald Ross, Manager/ESAT/Mantech

DATE: January 27, 1992

SUBJECT: Review of organic data for Syntex Facility.

TID#: 07-9109-016
ASSIGNMENT#: 099A
ICF ACCT#: 045-05-016-02
Mantech S.O.#: 1110-016
ESAT Document#: ESAT-VII-016-5099A

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses," February 1, 1988 revision.

The following comments and attached data sheets are a result of Mantech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE NO.:	17360	LABORATORY:	ENSECO
CONTRACT NO.:	68-D0-0151	METHOD NO:	CS03900
SITE:	Syntex Facility	EPA ACTIVITY:	IS651
REVIEWER:	Carla S. Phelps	MATRIX:	Water

<u>SMO SAMPLE NO.</u>	<u>EPA SAMPLE NO.</u>
GL678	IS651001
GL679	IS651002
GL680	IS651003
GL681	IS651004
GL682	IS651005
GL683	IS651006
GL684	IS651007
GL685	IS651008
* GL686	* IS651009F
* GL687	* IS651979P
** GL688	** IS651980P

* volatile analyses only

** base/neutral and acid analyses only

D. The base/neutral and acid initial calibrations were within quality control limit requirements for average relative response factor and percent relative standard deviation. The continuing calibrations were within quality control limit requirements for relative response factor for base/neutral and acid analyses.

E. Several compounds in the continuing calibrations were outside quality control limit requirements for percent difference (greater than 25%). Since all associated samples were non-detect, no data were qualified due to the continuing calibration rules.

4. INTERNAL STANDARD RESPONSE

A. Internal standard response areas were within quality control limit requirements for the volatile and base/neutral and acid fractions.

5. BLANKS

A. One trip blank and two method blanks were analyzed for the volatile fraction with methylene chloride and acetone detected. All compounds were detected below the contract required quantitation limit (CRQL) except for acetone in sample IS651009F. Acetone in samples IS651004, IS651001S and IS651001W were "U" coded due to the blank rules. Since methylene chloride was non-detect in all associated samples, no data were qualified.

B. Two method blanks were analyzed for the base/neutral and acid fractions with di-n-butylphthalate detected below the contract required quantitation limit (CRQL). Since di-n-butylphthalate in all associated samples were non-detect, no data were qualified due to the blank rules.

6. SURROGATE RECOVERY

A. All surrogates were within quality control limit requirements for percent recovery for the volatile fraction.

B. All surrogates were within quality control limit requirements for percent recovery for the base/neutral and acid fractions except for 2-chlorophenol-d4 in samples IS651004, IS651005, IS651006, IS651007, IS651008, IS651001W, and IS651904M; 2,4,6-tribromophenol in sample IS651008; nitrobenzene-d5 in sample IS651001S; 2-fluorobiphenyl in sample IS651001S; and terphenyl-d14 in sample IS651001S. Since 1,2,4-trichlorobenzene, acenaphthene, 1,4-dichlorobenzene, and n-nitroso-di-n-propylamine were also outside quality control limit requirements for percent recovery in the matrix spike, these compounds were "J" coded due to the poor surrogate recoveries.

D. Phenol, 2-chlorophenol, 4-nitrophenol, and pentachlorophenol, for the base/neutral and acid fractions in sample IS651001W, were "J" coded due to the compounds exceeding the calibration range.

E. Several compounds in several samples for volatile and base/neutral and acid analyses were detected above the instrument detection limit (IDL) but below the contract required quantitation limit (CRQL). These results were raised to the CRQL and "U" coded.

10. SUMMARY

A. Acetone in three samples were "U" coded due to the blank rule; 1,2,4-trichlorobenzene, acenaphthene, 1,4-dichlorobenzene, and n-nitroso-di-n-propylamine were "J" coded due to poor surrogate recoveries; acetone in one sample was "J" coded due to calibration outliers; phenol, 2-chlorophenol, 4-nitrophenol, and pentachlorophenol in one sample were "J" coded due to the compounds exceeding the calibration range. Acetone in one sample for the volatile fraction was "J" coded due to the compound both exceeding the calibration range and calibration outliers. Bis(2-ethylhexyl)phthalate for the base/neutral and acid fractions, was "J" coded due to holding times from extraction to analyses being exceeded.

B. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Organic Analysis dated March 1990, with the exceptions noted above.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

ICF Technology

ManTech Environmental Technology, Inc.

ESAT Region VII
ManTech Env. Tech., Inc.
25 Funston Road
Kansas City, KS 66115
(913) 551-5000

TO: Barry Evans, Data Review Task Monitor/ENSV
THRU: Harold Brown, Ph.D., ESAT Contract Manager/ENSV
FROM: Stephen R. Bales, ESAT Data Reviewer/ManTech
THRU: Ronald Ross, Manager/ESAT/ManTech *SB* *RE*
DATE: January 27, 1992
SUBJECT: Review of organic data for Syntex Facility.

TID#: 07-9109-016
ASSIGNMENT#: 099B
ICF ACCT#: 045-05-016-02
ManTech S.O.#: 1110-016
ESAT Document#: ESAT-VII-016-5099B

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Pesticides and PCBs".

The following comments and attached data sheets are a result of ManTech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE NO.:	17360	LABORATORY:	ENSECO
CONTRACT NO.:	68-D0-0151	METHOD NO:	CS03900
SITE:	Syntex Facility	EPA ACTIVITY:	IS651
REVIEWER:	Stephen R. Bales	MATRIX:	Water

Samples for Pesticide Analysis

<u>SMO SAMPLE NO.</u>	<u>EPA SAMPLE NO.</u>	<u>SMO SAMPLE NO.</u>	<u>EPA SAMPLE NO.</u>
GL678	IS651-001	GL679	IS651-002
GL680	IS651-003	GL681	IS651-004
GL682	IS651-005	GL683	IS651-006
GL684	IS651-007	GL685	IS651-008
GL689	IS651-981P		

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

A. A matrix spike/matrix spike duplicate was analyzed for pesticides/PCB's with the relative percent difference and percent recovery within quality control limits.

8. PESTICIDE CLEANUP CHECKS

A. All the percent recoveries for the florisil cartridge were outside quality control limits. No data were qualified due to the florisil cartridge.

9. PERFORMANCE EVALUATION SAMPLE

A. A performance evaluation sample was submitted to the laboratory for pesticide/PCB's analysis associated with this case with all audit compounds detected. The analysis detected 4,4'-DDE greater than the instrument detection limit but less than the contract required quantitation limit. Thus, 4,4'-DDE was raised to the CRQL and "U" coded.

10. COMPOUND IDENTIFICATION and QUANTITATION

A. In a level 2 review, no calculations from raw data (quantitation reports) are performed.

11. SUMMARY

A. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Organic Analysis dated March 1990.

ANALYSIS REQUEST REPORT

VALIDATED DATA

FOR ACTIVITY: IS651

S P F D

02/03/92 13:10:58

ALL REAL SAMPLES AND FIELD Q.C.

* FINAL REPORT

FY: 92 ACTIVITY: IS651 DESCRIPTION: SYNTEX FACILITY LOCATION: VERONA MISSOURI
 STATUS: ACTIVE TYPE: SAMPLING - CONTRACT LAB ANALYSIS PROJECT: A41
 LABO DUE DATE IS 12/31/91. REPORT DUE DATE IS 4/28/92.
 INSPECTION DATE: 10/31/91 ALL SAMPLES RECEIVED DATE: 11/01/91
 ALL DATA APPROVED BY LABO DATE: 01/31/92 FINAL REPORT TRANSMITTED DATE: 01/31/92
 EXPECTED LABO TURNAROUND TIME IS 60 DAYS EXPECTED REPORT TURNAROUND TIME IS 180 DAYS
 ACTUAL LABO TURNAROUND TIME IS 91 DAYS ACTUAL REPORT TURNAROUND TIME IS 92 DAYS
 SITE CODE: SITE:

SAMP. NO.	QCC	M	DESCRIPTION	SAMPLE STATUS	#	CITY	STATE	AIRS/ STORET LOC NO	LAY- SECT	ER	BEG. DATE	BEG. TIME	END. DATE	END. TIME
001	W		SYNTEX FACILITY-MW-1 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	09:00	10/31/91	09:15
002	W		SYNTEX FACILITY-MW-5 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	09:30	10/31/91	09:50
003	W		SYNTEX FACILITY-MW-11 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	10:05	10/31/91	10:30
004	W		SYNTEX FACILITY-MW-16 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	10:50	10/31/91	11:00
005	W		SYNTEX FACILITY-MW-16B SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	11:10	10/31/91	11:30
006	W		SYNTEX FACILITY-MW-9 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	11:00	10/31/91	11:10
007	W		SYNTEX FACILITY-MW-15 SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	11:35	10/31/91	12:05
008	W		SYNTEX FACILITY-MW-15B SAMPLE LOCATION	1		VERONA	MISSOURI				10/31/91	11:35	10/31/91	12:05
009	F	W	SYNTEX FACILITY-VOA TRIP BLANK	1		VERONA	MISSOURI				10/31/91	08:00	10/31/91	12:05

EXPLANATION OF CODES AND INFORMATION ON ANALYSIS REQUEST DETAIL REPORT

SAMPLE INFORMATION:

SAMP. NO. = SAMPLE IDENTIFICATION NUMBER (A 3-DIGIT NUMBER WHICH IN COMBINATION WITH THE ACTIVITY NUMBER AND QCC, PROVIDES AN UNIQUE NUMBER FOR EACH SAMPLE FOR IDENTIFICATION PURPOSES)

QCC = QUALITY CONTROL CODE (A ONE-LETTER CODE USED TO DESIGNATE SPECIFIC QC SAMPLES. THIS FIELD WILL BE BLANK FOR ALL NON-QC OR ACTUAL SAMPLES):

A = TRUE VALUE FOR CALIBRATION STANDARD
 B = CONCENTRATION RESULTING FROM DUPLICATE LAB SPIKE
 C = MEASURED VALUE FOR CALIBRATION STANDARD
 D = MEASURED VALUE FOR FILED DUPLICATE
 F = MEASURED VALUE FOR FIELD BLANK
 G = MEASURED VALUE FOR METHOD STANDARD
 H = TRUE VALUE FOR METHOD STANDARD
 K = CONCENTRATION RESULTING FROM DUPLICATE FIELD SPIKE
 L = MEASURED VALUE FOR LAB DUPLICATE
 M = MEASURED VALUE FOR LAB BLANK
 N = MEASURED VALUE FOR DUPLICATE FIELD SPIKE
 P = MEASURED VALUE FOR PERFORMANCE STANDARD
 R = CONCENTRATION RESULTING FROM LAB SPIKE
 S = MEASURED VALUE FOR LAB SPIKE
 T = TRUE VALUE OF PERFORMANCE STANDARD
 W = MEASURED VALUE FOR DUPLICATE LAB SPIKE
 Y = MEASURED VALUE FOR FIELD SPIKE
 Z = CONCENTRATION RESULTING FROM FIELD SPIKE

= MEDIA CODE (A ONE-LETTER CODE DESIGNATING THE MEDIA OF THE SAMPLE):

A = AIR
 H = OTHER (DOES NOT FIT ANY OTHER CATEGORY)
 S = SOLID (SOIL, SEDIMENT, SLUDGE)
 T = TISSUE (PLANT & ANIMAL)
 W = WATER (GROUND WATER, SURFACE WATER, WASTE WATER, DRINKING WATER)

DESCRIPTION = A SHORT DESCRIPTION OF THE LOCATION WHERE SAMPLE WAS COLLECTED

AIRS/STORET LOC. NO. = THE SPECIFIC LOCATION IDENTIFICATION NUMBER FOR EITHER OF THESE NATIONAL DATABASE SYSTEMS, AS APPROPRIATE

DATE/TIME INFORMATION = SPECIFIC INFORMATION REGARDING WHEN THE SAMPLE WAS COLLECTED

BEG. DATE = DATE SAMPLING WAS STARTED
 BEG. TIME = TIME SAMPLING WAS STARTED
 END DATE = DATE SAMPLING WAS COMPLETED
 END TIME = TIME SAMPLING WAS COMPLETED

NOTE: A GRAB SAMPLE WILL CONTAIN ONLY BEG. DATE/TIME
 A TIMED COMPOSITE SAMPLE WILL CONTAIN BOTH BEG AND END DATE/TIME TO DESIGNATE DURATION OF SAMPLE COLLECTION

OTHER CODES: V = VALIDATED

ANALYTICAL RESULTS/MEASUREMENTS INFORMATION:

COMPOUND = MGP (MEDIA-GROUP-PARAMETER) CODE AND NAME OF THE MEASURED CONSTITUENT OR CHARACTERISTIC OF EACH SAMPLE

UNITS = SPECIFIC UNITS IN WHICH RESULTS ARE REPORTED:

C = CENTIGRADE (CELSIUS) DEGREES
 CFS = CUBIC FEET PER SECOND
 GPM = GALLONS PER MINUTE
 IN = INCHES
 I.D. = SPECIES IDENTIFICATION
 KG = KILOGRAM
 L = LITER
 LB = POUNDS
 MG = MILLIGRAMS (1 X 10⁻³ GRAMS)
 MGD = MILLION GALLONS PER DAY
 MPH = MILES PER HOUR
 MV = MILLIVOLT
 M/F = MALE/FEMALE
 M2 = SQUARE METER
 M3 = CUBIC METER
 NA = NOT APPLICABLE
 NG = NANOGRAMS (1 X 10⁻⁹ GRAMS)
 NTU = NEPHELOMETRIC TURBIDITY UNITS
 PC/L = PICO (1 X 10⁻¹²) CURRIES PER LITER
 PG = PICOGRAMS (1 X 10⁻¹² GRAMS)
 P/CM2 = PICOGRAMS PER SQUARE CENTIMETER
 SCM = STANDARD CUBIC METER (1 ATM, 25 C)
 SQ FT = SQUARE FEET
 SU = STANDARD UNITS (PH)
 UG = MICROGRAMS (1 X 10⁻⁶ GRAMS)
 UMHOS = MICROMHOS/CM (CONDUCTIVITY UNITS)
 U/CC2 = MICROGRAMS PER 100 SQUARE CENTIMETERS
 U/CM2 = MICROGRAMS PER SQUARE CENTIMETER
 1000G = 1000 GALLONS
 +/- = POSITIVE/NEGATIVE
 # = NUMBER

DATA QUALIFIERS = SPECIFIC CODES USED IN CONJUNCTION WITH DATA VALUES TO PROVIDE ADDITIONAL INFORMATION ON THE REPORTED RESULTS, OR USED TO EXPLAIN THE ABSENCE OF A SPECIFIC VALUE:

BLANK = IF FIELD IS BLANK, NO REMARKS OR QUALIFIERS ARE PERTINENT. FOR FINAL REPORTED DATA, THIS MEANS THAT THE VALUES HAVE BEEN REVIEWED AND FOUND TO BE ACCEPTABLE FOR USE.

I = INVALID SAMPLE/DATA - VALUE NOT REPORTED
 J = DATA REPORTED BUT NOT VALID BY APPROVED QC PROCEDURES
 K = ACTUAL VALUE OF SAMPLE IS < VALUE REPORTED
 L = ACTUAL VALUE OF SAMPLE IS > VALUE REPORTED
 M = DETECTED BUT BELOW THE LEVEL OF REPORTED VALUE FOR ACCURATE QUANTIFICATION
 O = PARAMETER NOT ANALYZED
 U = ACTUAL VALUE OF SAMPLE IS < THE MEASUREMENT DETECTION LIMIT (REPORTED VALUE)

ANALYSIS REQUEST DETAIL REPORT

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VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
WFO1 TEMPERATURE, WATER	'C	14.2	16.8		15.9	15.9
WFO5 PH, FIELD	SU	6.50	6.88		7.46	7.88
WF10 CONDUCTIVITY (FIELD)	UMHOS	1040	480		290	290
WM01 SILVER, TOTAL, BY ICAP	UG/L	10 U	10 U	10 U	10 U	10 U
WM02 ALUMINUM, TOTAL, BY ICAP	UG/L	28000 J	61000 J	560 U	200 U	510 U
WM03 ARSENIC, TOTAL, BY ICAP	UG/L	14 J	N/A I	N/A I	N/A I	N/A I
WM04 BARIUM, TOTAL, BY ICAP	UG/L	2100	2000	210	200 U	200 U
WM05 BERYLLIUM, TOTAL, BY ICAP	UG/L	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U
WM06 CADMIUM, TOTAL, BY ICAP	UG/L	7.3 U	5.0 U	5.0 U	5.0 U	5.0 U
WM07 COBALT, TOTAL, BY ICAP	UG/L	120	50 U	50 U	50 U	50 U
WM08 CHROMIUM, TOTAL, BY ICAP	UG/L	340 J	97 J	10 U	10 U	17 J
WM09 COPPER, TOTAL, BY ICAP	UG/L	36 U	66 U	25 U	25 U	25 U
WM10 IRON, TOTAL, BY ICAP	UG/L	56000	63000	720	190 U	460
WM11 MANGANESE, TOTAL, BY ICAP	UG/L	42000	4400	180	230	1100
WM12 MOLYBDENUM, TOTAL, BY ICAP	UG/L	N/A 0	N/A 0	N/A 0	N/A 0	N/A 0
WM13 NICKEL, TOTAL, BY ICAP	UG/L	390	67 U	40 U	40 U	40 U
WM14 LEAD, TOTAL, BY ICAP	UG/L	32	30	3.0 U	3.0 U	3.0 U
WM15 ANTIMONY, TOTAL, BY ICAP	UG/L	60 U	60 U	60 U	60 U	60 U
WM16 SELENIUM, TOTAL, BY ICAP	UG/L	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U
WM17 TITANIUM, TOTAL, BY ICAP	UG/L	N/A 0	N/A 0	N/A 0	N/A 0	N/A 0
WM18 THALLIUM, TOTAL, BY ICAP	UG/L	10 U	10 U	10 U	10 U	10 U
WM19 VANADIUM, TOTAL, BY ICAP	UG/L	61	100	50 U	50 U	50 U
WM20 ZINC, TOTAL, BY ICAP	UG/L	170	220	29 U	20 U	52 U
WM21 CALCIUM, TOTAL, BY ICAP	MG/L	190	56	60	34	56
WM22 MAGNESIUM, TOTAL, BY ICAP	MG/L	14	9.6	11	18	5.0 U
WM23 SODIUM, TOTAL, BY ICAP	MG/L	120	43	5.0 U	9.0	12

ANALYSIS REQUEST DETAIL REPORT

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VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
WM24 POTASSIUM, TOTAL, BY ICAP	MG/L	7.6	10	5.0 U	5.0 U	5.0 U
WM34 MERCURY, TOTAL, BY COLD VAPOR AA	UG/L	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
WPO1 BHC, ALPHA, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO2 BHC, BETA, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO3 BHC, DELTA	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO4 BHC, GAMMA-(LINDANE), BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO5 ALDRIN, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO6 DIELDRIN, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WPO7 ENDOSULFAN I, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WPO8 ENDOSULFAN II, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WPO9 ENDOSULFAN SULFATE, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP10 ENDRIN, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP11 ENDRIN ALDEHYDE, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP13 DDE-4,4'-	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP14 DDD-4,4'-	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP15 DDT-4,4'-	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP16 TOXAPHENE, BY GC/EC	UG/L	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U
WP17 PCB-AROCLOR 1016	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP18 PCB-AROCLOR 1221	UG/L	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U
WP19 PCB-AROCLOR 1232	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP20 PCB-AROCLOR 1242	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP21 PCB-AROCLOR 1248	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP22 PCB-AROCLOR 1254	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP23 PCB-AROCLOR 1260	UG/L	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
WP24 CHLORDANE, TECHNICAL, BY GC/EC	UG/L	N/A 0	N/A 0	N/A 0	N/A 0	N/A 0
WP25 HEPTACHLOR, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U

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VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
WP26 HEPTACHLOR EPOXIDE, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WP83 ENDRIN KETONE, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP84 METHOXYCHLOR, BY GC/EC	UG/L	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U
WP85 CHLORDANE, ALPHA	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WP86 CHLORDANE, GAMMA	UG/L	0.050 U	0.050 U	0.050 U	0.050 U	0.050 U
WS01 PHENOL, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS03 ETHER, BIS(2-CHLOROETHYL), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS04 2-CHLOROPHENOL	UG/L	10 U	10 U	10 U	10 U	10 U
WS05 DICHLOROBENZENE, 1,3-, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS06 DICHLOROBENZENE, 1,4-	UG/L	10 U	10 U	10 U	10 U	10 U
WS07 BENZYL ALCOHOL	UG/L	N/A 0	N/A 0	N/A 0	N/A 0	N/A 0
WS08 DICHLOROBENZENE, 1,2-, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS09 CRESOL, ORTHO(2-METHYLPHENOL)	UG/L	10 U	10 U	10 U	10 U	10 U
WS10 ETHER, BIS(2-CHLOROISOPROPYL), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS11 CRESOL, PARA-(4-METHYLPHENOL)	UG/L	10 U	10 U	10 U	10 U	10 U
WS12 N-NITROSODIPROPYLAMINE	UG/L	10 U	10 U	10 U	10 U	10 U
WS13 HEXACHLOROETHANE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS14 NITROBENZENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS15 ISOPHORONE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS16 NITROPHENOL, 2-	UG/L	10 U	10 U	10 U	10 U	10 U
WS17 DIMETHYLPHENOL, 2,4, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS18 BENZOIC ACID, BY GC/MS	UG/L	N/A 0	N/A 0	N/A 0	N/A 0	N/A 0
WS19 METHANE, BIS(2-CHLOROETHOXY), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS20 2,4-DICHLOROPHENOL	UG/L	10 U	10 U	10 U	10 U	10 U
WS21 TRICHLOROBENZENE, 1,2,4, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS22 NAPHTHALENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U

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VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
WS23 CHLOROANILINE, 4-	UG/L	10 U	10 U	10 U	10 U	10 U
WS24 HEXACHLOROBUTADIENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS25 PHENOL, 4-CHLORO-3-METHYL	UG/L	10 U	10 U	10 U	10 U	10 U
WS26 2-METHYLNAPHTHALENE	UG/L	10 U	10 U	10 U	10 U	10 U
WS27 HEXACHLOROCYCLOPENTADIENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS28 TRICHLOROPHENOL, 2, 4, 6	UG/L	10 U	10 U	10 U	10 U	10 U
WS29 TRICHLOROPHENOL, 2, 4, 5	UG/L	25 U	25 U	25 U	25 U	25 U
WS30 2-CHLORONAPHTHALENE	UG/L	10 U	10 U	10 U	10 U	10 U
WS31 NITROANILINE, 2-(ORTHO)	UG/L	25 U	25 U	25 U	25 U	25 U
WS32 PHTHALATE, DIMETHYL, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS33 ACENAPHTHYLENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS34 NITROANILINE, 3-	UG/L	25 U	25 U	25 U	25 U	25 U
WS35 ACENAPHTHENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS36 DINITROPHENOL, 2, 4, BY GC/MS	UG/L	25 U	25 U	25 U	25 U	25 U
WS37 NITROPHENOL, 4-	UG/L	25 U	25 U	25 U	25 U	25 U
WS38 DIBENZOFURAN	UG/L	10 U	10 U	10 U	10 U	10 U
WS39 DINITROTOLUENE, 2, 4, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS40 DINITROTOLUENE, 2, 6	UG/L	10 U	10 U	10 U	10 U	10 U
WS41 PHTHALATE, DIETHYL, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS42 ETHER, 4-CHLOROPHENYL PHENYL	UG/L	10 U	10 U	10 U	10 U	10 U
WS43 FLUORENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS44 NITROANILINE, 4-	UG/L	25 U	25 U	25 U	25 U	25 U
WS45 PHENOL, 4, 6-DINITRO-2-METHYL	UG/L	25 U	25 U	25 U	25 U	25 U
WS46 N-NITROSODIPHENYLAMINE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS47 ETHER, 4-BROMOPHENYL PHENYL	UG/L	10 U	10 U	10 U	10 U	10 U
WS48 HEXACHLOROENZENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U

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VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
WS49 PENTACHLOROPHENOL, BY GC/MS	UG/L	25 U	25 U	25 U	25 U	25 U
WS50 PHENANTHRENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS51 ANTHRACENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS52 PHTHALATE, DI-N-BUTYL-, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS53 FLUORANTHENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS54 PYRENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS55 PHTHALATE, BUTYL BENZYL	UG/L	10 U	10 U	10 U	10 U	10 U
WS56 DICHLOROBENZIDINE, 3,3'	UG/L	10 U	10 U	10 U	10 U	10 U
WS57 ANTHRACENE, BENZO(A), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS58 PHTHALATE, BIS(2-ETHYLHEXYL), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS59 CHRYSENE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS60 PHTHALATE, DI-N-OCTYL-, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS61 FLUORANTHENE, BENZO(B), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS62 FLUORANTHENE, BENZO(K), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS63 PYRENE, BENZO(A), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS64 PYRENE, INDENO(1,2,3-CD)	UG/L	10 U	10 U	10 U	10 U	10 U
WS65 ANTHRACENE, DIBENZO(A,H), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS66 PERYLENE, BENZO(G,H,I), BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WS67 CARBAZOLE	UG/L	10 U	10 U	10 U	10 U	10 U
WV03 CHLOROMETHANE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WV04 BROMOMETHANE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WV05 VINYL CHLORIDE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WV06 CHLOROETHANE, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U
WV07 METHYLENE CHLORIDE (DICHLOROMETHANE)	UG/L	10 U	10 U	10 U	10 U	10 U
WV08 DICHLOROETHYLENE, 1,1	UG/L	10 U	10 U	10 U	10 U	10 U
WV09 DICHLOROETHANE, 1,1, BY GC/MS	UG/L	10 U	10 U	10 U	10 U	10 U

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VALIDATED DATA

COMPOUND	UNITS	001		002		003		004		005	
WV10 DICHLOROETHYLENE, TRANS,1,2-	UG/L	10	U	10	U	10	U	10	U	10	U
WV11 CHLOROFORM, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV12 DICHLOROETHANE,1,2, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV13 TRICHLOROETHANE,1,1,1-, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV14 CARBON TETRACHLORIDE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV15 BROMODICHLOROMETHANE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV16 DICHLOROPROPANE,1,2, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV17 BENZENE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV19 TRICHLOROETHYLENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV20 DICHLOROPROPYLENE,CIS-1,3, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV21 DIBROMOCHLOROMETHANE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV22 TRICHLOROETHANE,1,1,2-, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV24 BROMOFORM, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV25 TETRACHLOROETHYLENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV26 TOLUENE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV27 TETRACHLOROETHANE,1,1,2,2, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV28 CHLOROBENZENE, BY GC/MS	UG/L	10	U	13		10	U	10	U	10	U
WV29 ETHYL BENZENE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV30 ACETONE, BY GC/MS	UG/L	10	U	10	U	10	U	26	U	10	U
WV31 CARBON DISULFIDE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV32 METHYL ETHYL KETONE (2-BUTANONE)	UG/L	10	U	10	U	10	U	10	U	10	U
WV34 2-HEXANONE	UG/L	10	U	10	U	10	U	10	U	10	U
WV35 4-METHYL-2-PENTANONE	UG/L	10	U	10	U	10	U	10	U	10	U
WV36 STYRENE, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV37 XYLENES, TOTAL, BY GC/MS	UG/L	10	U	10	U	10	U	10	U	10	U
WV40 DICHLOROPROPYLENE,TRANS-1,3	UG/L	10	U	10	U	10	U	10	U	10	U

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	001	002	003	004	005
ZZ01 SAMPLE NUMBER	NA	001	002	003	004	005
ZZ02 ACTIVITY CODE	NA	IS651	IS651	IS651	IS651	IS651

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
WFO1 TEMPERATURE, WATER	'C	15.3	13.9	15.9	
WFO5 PH, FIELD	SU	6.92	7.41	6.69	
WF10 CONDUCTIVITY (FIELD)	UMHOS	780	520	1400	
WM01 SILVER, TOTAL, BY ICAP	UG/L	10 U	10 U	10 U	
WM02 ALUMINUM, TOTAL, BY ICAP	UG/L	6400 J	220 U	260 U	
WM03 ARSENIC, TOTAL, BY ICAP	UG/L	22 J	N/A I	N/A I	
WM04 BARIUM, TOTAL, BY ICAP	UG/L	650	200 U	1200	
WM05 BERYLLIUM, TOTAL, BY ICAP	UG/L	5.0 U	5.0 U	5.0 U	
WM06 CADMIUM, TOTAL, BY ICAP	UG/L	5.0 U	5.0 U	5.0 U	
WM07 COBALT, TOTAL, BY ICAP	UG/L	50 U	50 U	50 U	
WM08 CHROMIUM, TOTAL, BY ICAP	UG/L	15 J	11 J	10 U	
WM09 COPPER, TOTAL, BY ICAP	UG/L	25 U	25 U	25 U	
WM10 IRON, TOTAL, BY ICAP	UG/L	30000	180 U	14000	
WM11 MANGANESE, TOTAL, BY ICAP	UG/L	4400	79	4300	
WM12 MOLYBDENUM, TOTAL, BY ICAP	UG/L	N/A O	N/A O	N/A O	
WM13 NICKEL, TOTAL, BY ICAP	UG/L	40 U	40 U	40 U	
WM14 LEAD, TOTAL, BY ICAP	UG/L	4.6	3.0 U	3.0 U	
WM15 ANTIMONY, TOTAL, BY ICAP	UG/L	60 U	62	60 U	
WM16 SELENIUM, TOTAL, BY ICAP	UG/L	5.0 U	5.0 U	5.0 U	
WM17 TITANIUM, TOTAL, BY ICAP	UG/L	N/A O	N/A O	N/A O	
WM18 THALLIUM, TOTAL, BY ICAP	UG/L	10 U	10 U	10 U	
WM19 VANADIUM, TOTAL, BY ICAP	UG/L	50 U	50 U	50 U	
WM20 ZINC, TOTAL, BY ICAP	UG/L	42 U	53 U	76	
WM21 CALCIUM, TOTAL, BY ICAP	MG/L	66	93	88	
WM22 MAGNESIUM, TOTAL, BY ICAP	MG/L	5.0 U	7.4	8.1	
WM23 SODIUM, TOTAL, BY ICAP	MG/L	63	5.3	89	

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F	
WM24 POTASSIUM, TOTAL, BY ICAP	MG/L	5.0	U	5.0	U	6.9
WM34 MERCURY, TOTAL, BY COLD VAPOR AA	UG/L	0.20	U	0.20	U	0.20
WPO1 BHC, ALPHA, BY GC/EC	UG/L	0.050	U	0.050	U	0.050
WPO2 BHC, BETA, BY GC/EC	UG/L	0.050	U	0.050	U	0.050
WPO3 BHC, DELTA	UG/L	0.050	U	0.050	U	0.050
WPO4 BHC, GAMMA-(LINDANE), BY GC/EC	UG/L	0.050	U	0.050	U	0.050
WPO5 ALDRIN, BY GC/EC	UG/L	0.050	U	0.050	U	0.050
WPO6 DIELDRIN, BY GC/EC	UG/L	0.10	U	0.10	U	0.10
WPO7 ENDOSULFAN I, BY GC/EC	UG/L	0.050	U	0.050	U	0.050
WPO8 ENDOSULFAN II, BY GC/EC	UG/L	0.10	U	0.10	U	0.10
WPO9 ENDOSULFAN SULFATE, BY GC/EC	UG/L	0.10	U	0.10	U	0.10
WP10 ENDRIN, BY GC/EC	UG/L	0.10	U	0.10	U	0.10
WP11 ENDRIN ALDEHYDE, BY GC/EC	UG/L	0.10	U	0.10	U	0.10
WP13 DDE-4,4'-	UG/L	0.10	U	0.10	U	0.10
WP14 DDD-4,4'-	UG/L	0.10	U	0.10	U	0.10
WP15 DDT-4,4'-	UG/L	0.10	U	0.10	U	0.10
WP16 TOXAPHENE, BY GC/EC	UG/L	5.0	U	5.0	U	5.0
WP17 PCB-AROCLOR 1016	UG/L	1.0	U	1.0	U	1.0
WP18 PCB-AROCLOR 1221	UG/L	2.0	U	2.0	U	2.0
WP19 PCB-AROCLOR 1232	UG/L	1.0	U	1.0	U	1.0
WP20 PCB-AROCLOR 1242	UG/L	1.0	U	1.0	U	1.0
WP21 PCB-AROCLOR 1248	UG/L	1.0	U	1.0	U	1.0
WP22 PCB-AROCLOR 1254	UG/L	1.0	U	1.0	U	1.0
WP23 PCB-AROCLOR 1260	UG/L	1.0	U	1.0	U	1.0
WP24 CHLORDANE, TECHNICAL, BY GC/EC	UG/L	N/A	0	N/A	0	N/A
WP25 HEPTACHLOR, BY GC/EC	UG/L	0.050	U	0.050	U	0.050

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
WP26 HEPTACHLOR EPOXIDE, BY GC/EC	UG/L	0.050 U	0.050 U	0.050 U	
WP83 ENDRIN KETONE, BY GC/EC	UG/L	0.10 U	0.10 U	0.10 U	
WP84 METHOXYCHLOR, BY GC/EC	UG/L	0.50 U	0.50 U	0.50 U	
WP85 CHLORDANE, ALPHA	UG/L	0.050 U	0.050 U	0.050 U	
WP86 CHLORDANE, GAMMA	UG/L	0.050 U	0.050 U	0.050 U	
WS01 PHENOL, BY GC/MS	UG/L	10 U	10 U	10 U	
WS03 ETHER, BIS(2-CHLOROETHYL), BY GC/MS	UG/L	10 U	10 U	10 U	
WS04 2-CHLOROPHENOL	UG/L	10 U	10 U	10 U	
WS05 DICHLOROBENZENE, 1,3-, BY GC/MS	UG/L	10 U	10 U	10 U	
WS06 DICHLOROBENZENE, 1,4-	UG/L	25 U	10 U	65 U	
WS07 BENZYL ALCOHOL	UG/L	N/A 0	N/A 0	N/A 0	
WS08 DICHLOROBENZENE, 1,2-, BY GC/MS	UG/L	10 U	10 U	10 U	
WS09 CRESOL, ORTHO(2-METHYLPHENOL)	UG/L	10 U	10 U	10 U	
WS10 ETHER, BIS(2-CHLOROISOPROPYL), BY GC/MS	UG/L	10 U	10 U	10 U	
WS11 CRESOL, PARA-(4-METHYLPHENOL)	UG/L	10 U	10 U	10 U	
WS12 N-NITROSODIPROPYLAMINE	UG/L	10 U	10 U	10 U	
WS13 HEXACHLOROETHANE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS14 NITROBENZENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS15 ISOPHORONE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS16 NITROPHENOL, 2-	UG/L	10 U	10 U	10 U	
WS17 DIMETHYLPHENOL, 2,4, BY GC/MS	UG/L	10 U	10 U	10 U	
WS18 BENZOIC ACID, BY GC/MS	UG/L	N/A 0	N/A 0	N/A 0	
WS19 METHANE, BIS(2-CHLOROETHOXY), BY GC/MS	UG/L	10 U	10 U	10 U	
WS20 2,4-DICHLOROPHENOL	UG/L	10 U	10 U	10 U	
WS21 TRICHLOROBENZENE, 1,2,4, BY GC/MS	UG/L	10 U	10 U	10 U	
WS22 NAPHTHALENE, BY GC/MS	UG/L	10 U	10 U	10 U	

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
WS23 CHLOROANILINE, 4-	UG/L	10 U	10 U	10 U	
WS24 HEXACHLOROBUTADIENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS25 PHENOL, 4-CHLORO-3-METHYL	UG/L	10 U	10 U	10 U	
WS26 2-METHYLNAPHTHALENE	UG/L	10 U	10 U	10 U	
WS27 HEXACHLOROCYCLOPENTADIENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS28 TRICHLOROPHENOL, 2,4,6	UG/L	10 U	10 U	10 U	
WS29 TRICHLOROPHENOL, 2,4,5	UG/L	25 U	25 U	25 U	
WS30 2-CHLORONAPHTHALENE	UG/L	10 U	10 U	10 U	
WS31 NITROANILINE, 2-(ORTHO)	UG/L	25 U	25 U	25 U	
WS32 PHTHALATE, DIMETHYL, BY GC/MS	UG/L	10 U	10 U	10 U	
WS33 ACENAPHTHYLENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS34 NITROANILINE, 3-	UG/L	25 U	25 U	25 U	
WS35 ACENAPHTHENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS36 DINITROPHENOL, 2,4, BY GC/MS	UG/L	25 U	25 U	25 U	
WS37 NITROPHENOL, 4-	UG/L	25 U	25 U	25 U	
WS38 DIBENZOFURAN	UG/L	10 U	10 U	10 U	
WS39 DINITROTOLUENE, 2,4, BY GC/MS	UG/L	10 U	10 U	10 U	
WS40 DINITROTOLUENE, 2,6	UG/L	10 U	10 U	10 U	
WS41 PHTHALATE, DIETHYL, BY GC/MS	UG/L	10 U	10 U	10 U	
WS42 ETHER, 4-CHLOROPHENYL PHENYL	UG/L	10 U	10 U	10 U	
WS43 FLUORENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS44 NITROANILINE, 4-	UG/L	25 U	25 U	25 U	
WS45 PHENOL, 4,6-DINITRO-2-METHYL	UG/L	25 U	25 U	25 U	
WS46 N-NITROSODIPHENYLAMINE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS47 ETHER, 4-BROMOPHENYL PHENYL	UG/L	10 U	10 U	10 U	
WS48 HEXACHLOROBENZENE, BY GC/MS	UG/L	10 U	10 U	10 U	

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ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
WS49 PENTACHLOROPHENOL, BY GC/MS	UG/L	25 U	25 U	25 U	
WS50 PHENANTHRENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS51 ANTHRACENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS52 PHTHALATE, DI-N-BUTYL-, BY GC/MS	UG/L	10 U	10 U	10 U	
WS53 FLUORANTHENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS54 PYRENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS55 PHTHALATE, BUTYL BENZYL	UG/L	10 U	10 U	10 U	
WS56 DICHLOROBENZIDINE, 3,3'	UG/L	10 U	10 U	10 U	
WS57 ANTHRACENE, BENZO(A), BY GC/MS	UG/L	10 U	10 U	10 U	
WS58 PHTHALATE, BIS(2-ETHYLHEXYL), BY GC/MS	UG/L	10 U	10 U	10 U	
WS59 CHRYSENE, BY GC/MS	UG/L	10 U	10 U	10 U	
WS60 PHTHALATE, DI-N-OCTYL-, BY GC/MS	UG/L	10 U	10 U	10 U	
WS61 FLUORANTHENE, BENZO(B), BY GC/MS	UG/L	10 U	10 U	10 U	
WS62 FLUORANTHENE, BENZO(K), BY GC/MS	UG/L	10 U	10 U	10 U	
WS63 PYRENE, BENZO(A), BY GC/MS	UG/L	10 U	10 U	10 U	
WS64 PYRENE, INDENO(1,2,3-CD)	UG/L	10 U	10 U	10 U	
WS65 ANTHRACENE, DIBENZO(A,H), BY GC/MS	UG/L	10 U	10 U	10 U	
WS66 PERYLENE, BENZO(G,H,I), BY GC/MS	UG/L	10 U	10 U	10 U	
WS67 CARBAZOLE	UG/L	10 U	10 U	10 U	
WV03 CHLOROMETHANE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV04 BROMOMETHANE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV05 VINYL CHLORIDE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV06 CHLOROETHANE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV07 METHYLENE CHLORIDE (DICHLOROMETHANE)	UG/L	25 U	10 U	20 U	10 U
WV08 DICHLOROETHYLENE, 1,1	UG/L	25 U	10 U	20 U	10 U
WV09 DICHLOROETHANE, 1,1, BY GC/MS	UG/L	25 U	10 U	20 U	10 U

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
WV10 DICHLOROETHYLENE, TRANS,1,2-	UG/L	25 U	10 U	20 U	10 U
WV11 CHLOROFORM, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV12 DICHLOROETHANE,1,2, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV13 TRICHLOROETHANE,1,1,1-, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV14 CARBON TETRACHLORIDE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV15 BROMODICHLOROMETHANE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV16 DICHLOROPROPANE,1,2, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV17 BENZENE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV19 TRICHLOROETHYLENE	UG/L	25 U	10 U	20 U	10 U
WV20 DICHLOROPROPYLENE,CIS-1,3, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV21 DIBROMOCHLOROMETHANE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV22 TRICHLOROETHANE,1,1,2-, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV24 BROMOFORM, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV25 TETRACHLOROETHYLENE	UG/L	25 U	10 U	20 U	10 U
WV26 TOLUENE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV27 TETRACHLOROETHANE,1,1,2,2, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV28 CHLOROBENZENE, BY GC/MS	UG/L	110	10 U	140	10 U
WV29 ETHYL BENZENE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV30 ACETONE, BY GC/MS	UG/L	25 U	10 U	5100 J	45 J
WV31 CARBON DISULFIDE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV32 METHYL ETHYL KETONE (2-BUTANONE)	UG/L	25 U	10 U	20 U	10 U
WV34 2-HEXANONE	UG/L	25 U	10 U	20 U	10 U
WV35 4-METHYL-2-PENTANONE	UG/L	25 U	10 U	20 U	10 U
WV36 STYRENE, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV37 XYLENES, TOTAL, BY GC/MS	UG/L	25 U	10 U	20 U	10 U
WV40 DICHLOROPROPYLENE,TRANS-1,3	UG/L	10 U	10 U	10 U	10 U

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 2-IS651

VALIDATED DATA

COMPOUND	UNITS	006	007	008	009F
ZZ01 SAMPLE NUMBER	NA	006	007	008	009
ZZ02 ACTIVITY CODE	NA	IS651	IS651	IS651	IS651

ACTIVITY IS651 SYNTEX FACILITY

THE PROJECT LEADER SHOULD CIRCLE ONE - STORET, AIRS, OR ARCHIVE.

CIRCLE ONE: STORET AIRS ARCHIVE

DATA APPROVED BY LABO FOR TRANSMISSION TO PROJECT LEADER ON 01/31/92 09:48:05 BY _____

