

# CDM Federal Programs Corporation

May 31, 1989

Rose Harvell  
Project Officer  
U.S. Environmental Protection Agency  
401 M Street, Room 2834  
Washington, D.C. 20460

Superfund Records Center

DATE: Mottolo

BREAK: 3.2

OTHER: 468844



SDMS DocID 468844

PROJECT: EPA CONTRACT NO.: 68-01-7331  
DOCUMENT NO.: T993-C01-EP-DFLU-1  
SUBJECT: Data Validation Letter Report for  
Work Assignment C01011  
Mottolo, Raymond, New Hampshire  
Document No.: T993-C01-LR-DECU-3

Dear Ms. Harvell:

Please find enclosed the Data Validation Letter Report for Organics SAS Case #4372A for the Mottolo site, Raymond, New Hampshire as partial fulfillment of the reporting requirements for this work assignment.

If you have any comments regarding this submittal, please contact James Hewitt of Geoscience Consultants, Ltd. (GCL) at (301) 587-2088 within two weeks of receipt of this letter.

Sincerely,

CDM Federal Programs Corporation

*Karen Stone*

Karen Stone  
TES III Regional Manager

AR:rf

Enclosure

cc: Roger Duwart, EPA Primary Contact, CERCLA Region I  
Richard Leighton, EPA Regional Project Officer, CERCLA Region I  
Jack Shad, EPA HQ Coordinator, CERCLA Region I  
Harry Butler, CDM Federal Programs Corporation Program Manager  
Michael P. Riley, EPA Contracting Officer (letter only)  
Deborah Szaro, Deputy Project Officer, EPA Region I  
James Hewitt, GCL Program Manager, (letter only)

**LETTER REPORT  
ORGANIC SAS CASE #4372A  
MOTTOLO, RAYMOND, NEW HAMPSHIRE  
DATA VALIDATION**

**Prepared for**

**U.S. ENVIRONMENTAL PROTECTION AGENCY  
Office of Waste Programs Enforcement  
Washington, D.C. 20460**

Work Assignment No. : C01011  
EPA Region : I  
Facility I.D. No. :  
Contract No. : 68-01-7331  
CDM Federal Programs  
Corporation Document No. : T993-C01-LR-DECU-3  
Prepared By : GCL  
Work Assignment Project  
Manager : James Hewitt  
Telephone No. : (301) 587-2088  
Primary Contact : Roger Duwart  
Telephone Number : (617) 573-9628  
Date Prepared : May 31, 1989

ORGANIC REGIONAL DATA ASSESSMENT

CASE NO. SAS 4372 A SITE Motto 10  
 LABORATORY EC Jordan Co. NO. OF SAMPLES/  
 MATRIX 7/soil  
 SDG # AM690 REVIEWER (IF NOT ESD) GCL/CDM FPC  
 SOW# \_\_\_\_\_ REVIEWER'S NAME J. Hewitt/A. Rigassio  
 DPO: ACTION \_\_\_\_\_ FYI X COMPLETION DATE 5-5-89/5-8-89

DATA ASSESSMENT SUMMARY

	VOA	BNA	PEST	OTHER
1. HOLDING TIMES	<u>0<sup>1</sup></u>	<u>X<sup>2</sup></u>	<u>X<sup>2</sup></u>	_____
2. GC/MS TUNE/INSTR. PERFORM.	<u>0</u>	<u>0</u>	<u>X<sup>4</sup></u>	_____
3. CALIBRATIONS	<u>0<sup>3</sup></u>	<u>0<sup>3</sup></u>	<u>0<sup>3</sup></u>	_____
4. BLANKS	<u>X<sup>4</sup></u>	<u>X<sup>4</sup></u>	<u>X<sup>4</sup></u>	_____
5. SURROGATES	<u>0</u>	<u>0</u>	<u>0</u>	_____
6. MATRIX SPIKE/DUP	<u>0</u>	<u>0</u>	<u>M<sup>5</sup></u>	_____
7. OTHER QC	<u>M<sup>6</sup></u>	<u>0</u>	<u>0</u>	_____
8. INTERNAL STANDARDS	_____	_____	_____	_____
9. COMPOUND IDENTIFICATION	<u>0</u>	<u>0</u>	<u>0</u>	_____
10. SYSTEM PERFORMANCE	<u>0</u>	<u>0</u>	<u>M<sup>1</sup></u>	_____
11. OVERALL ASSESSMENT	_____	_____	_____	_____

O = Data had no problems/or qualified due to minor problems.  
 M = Data qualified due to major problems.  
 Z = Data unacceptable.  
 X = Problems, but do not affect data.

**CONCERN**  
 ACTION ITEMS: 0<sup>1</sup> - Estimate aromatic Compound results in Sample AM694. X<sup>2</sup> - HTs exceeded  
but no action for BNA + Pest/PCB frac. of AM690, AM691, AM692 + AM693. 0<sup>3</sup> - Analytes estimated  
in all samples due to calib. violations. X<sup>4</sup> - Methyl, di-n-butyl phthalate, bis(2-ch)phthal  
**AREAS OF CONCERN:** butyl benzyl phthal. + endrin found in blanks. Det. limits raised  
for all analytes except endrin. M<sup>5</sup> - Several pesticide cpds estimated due to imprecison  
X<sup>4</sup> - DBC90D > 2 for some samples, but no action. M<sup>6</sup> - field dup action on TECE

NOTABLE PERFORMANCE: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

~~FIGURE 6~~  
DOO SUMMARY FORM

1. SITE		EPA REGION <u>I</u>					
NAME <u>Mottolo</u>		PHASE _____					
LOCATION <u>Raymond, NH</u>		R11 R12 R13 ERA FS RD RA					
NUMBER <u>CD1011</u>		(CIRCLE ONE)					
2. MEDIA (CIRCLE ONE)	<input checked="" type="checkbox"/> SOIL	<input type="checkbox"/> GW	<input type="checkbox"/> SW/SED	<input type="checkbox"/> AIR	<input type="checkbox"/> BIO	<input type="checkbox"/> OTHER _____	
3. USE (CIRCLE ALL THAT APPLY)	SITE CHARAC. (H&S)	RISK ASSESS.	EVAL. ALTS.	ENGG. DESIGN	PRP DETER.	MONITORING REMEDIAL ACTION	<input checked="" type="checkbox"/> OTHER <u>oversight</u>
4. OBJECTIVE <u>To verify PRPs &amp; soil analysis results</u>							
5. SITE INFORMATION							
AREA <u>1 acre</u>		DEPTH TO GROUND WATER <u>7-8'</u>					
GROUND WATER USE <u>Potable Bedrock Wells</u>							
SOIL TYPES <u>Glacial Till / Outwash</u>							
SENSITIVE RECEPTORS <u>Brook A, Snake, River, Potable H<sub>2</sub>O Source</u>							
6. DATA TYPES (CIRCLE APPROPRIATE DATA TYPES)							
A. ANALYTICAL DATA				B. PHYSICAL DATA			
<input type="checkbox"/> pH	<input type="checkbox"/> CONDUCTIVITY	<input checked="" type="checkbox"/> PESTICIDES	<input type="checkbox"/> TOX	<input type="checkbox"/> PERMEABILITY	<input type="checkbox"/> HYDRAULIC HEAD		
<input type="checkbox"/> VOA	<input type="checkbox"/> ABN	<input checked="" type="checkbox"/> PCB	<input type="checkbox"/> TOC	<input type="checkbox"/> POROSITY	<input type="checkbox"/> PENETRATION TEST		
<input type="checkbox"/> TSP	<input type="checkbox"/> METALS	<input type="checkbox"/> BTX	<input type="checkbox"/> COC	<input type="checkbox"/> GRAIN SIZE	<input type="checkbox"/> HARDNESS		
<input checked="" type="checkbox"/> Tetrahydrofuran	<input checked="" type="checkbox"/> Methyl-t-butyl ether	<input type="checkbox"/> CYANIDE	<input type="checkbox"/> COC	<input type="checkbox"/> BULK DENSITY	_____		
7. SAMPLING METHOD (CIRCLE METHOD(S) TO BE USED)							
<input type="checkbox"/> ENVIRONMENTAL	<input type="checkbox"/> BIASED	<input checked="" type="checkbox"/> GRAB	<input type="checkbox"/> NON-INTRUSIVE	<input type="checkbox"/> PHASED			
<input type="checkbox"/> SOURCE	<input type="checkbox"/> GRID	<input type="checkbox"/> COMPOSITE	<input type="checkbox"/> INTRUSIVE	_____			
8. ANALYTICAL LEVELS (INDICATE LEVEL(S) AND EQUIPMENT & METHODS)							
LEVEL 1 FIELD SCREENING - EQUIPMENT _____							
LEVEL 2 FIELD ANALYSIS - EQUIPMENT _____							
LEVEL 3 NON-CLP LABORATORY - METHODS _____							
<input checked="" type="checkbox"/> LEVEL 4 CLP/RAS - METHODS <u>ABN + pest / PCB</u>							
<input checked="" type="checkbox"/> LEVEL 5 NON STANDARD <u>VOA + tetrahydrofuran + methyl-t-butyl ether</u>							
9. SAMPLING PROCEDURES							
BACKGROUND - 2 PER EVENT OR _____							
CRITICAL (LIST) _____							
PROCEDURES _____							
10. QUALITY CONTROL SAMPLES (CONFIRM OR SET STANDARD)							
A. FIELD				B. LABORATORY			
COLLOCATED - 5% OR _____				REAGENT BLANK - 1 PER ANALYSIS BATCH OR _____			
REPLICATE - 5% OR <u>two separate dup parts</u>				REPLICATE - 1 PER ANALYSIS BATCH OR _____			
FIELD BLANK - 5% OR _____				MATRIX SPIKE - 1 PER ANALYSIS BATCH OR _____			
TRIP BLANK - 1 PER DAY OR _____				OTHER <u>RAS CLP Requirements</u>			
11. BUDGET REQUIREMENTS							
BUDGET _____				SCHEDULE <u>Week of December 26<sup>th</sup>, 1988</u>			
STAFF <u>Mike Kulbersh / Joe Newton</u>				_____			
CONTRACTOR _____				PRIME CONTRACTOR <u>ODM FPC</u>			
SITE MANAGER <u>Michael Kulbersh</u>				DATE _____			

FOR DETAILS SEE SAMPLING & ANALYSIS PLAN

CDM SF DOO 1.002

May 17, 1989

Mr. Dennis Gagne  
U.S. Environmental Protection Agency  
JFK Federal Building  
Boston, Massachusetts 02203

Re: Case SAS4372  
E.C. Jordan Co. Laboratory  
Mottolo Superfund Site; Raymond, NH  
Volatiles: Five/Soil/AM690, AM691, AM694, AM695, AM959  
Semi-volatiles: Five/Soil/AM690, AM692, AM693, AM694, AM695  
Pesticide/PCB: Five/Soil/AM690, AM692, AM693, AM694, AM695

Dear Mr. Gagne:

A validation was performed on the organic analytical data from Case SAS4372 low level soil samples collected at the Mottolo site. The samples were collected by CDM Federal Programs Corporation. The samples were analyzed as a Special Analytical Service because analyses of two additional compounds, methyl-t-butyl ether and tetrahydrofuran was requested. The data was evaluated based on the following parameters:

- \* o Data completeness
- o Holding times
- \* o GC/MS tuning
- o Calibration
- o Blanks
- \* o Surrogate recoveries
- o Matrix spike/matrix spike duplicate
- o Field duplicates
- o Instrument Performance
- \* o Internal Standard Performance
- \* o Compound identification
- \* o Compound quantitation

\* All criteria were met for this parameter.

Table I summarizes the validation recommendations which were based on the following information:

#### Holding Times

Sample AM694 was analyzed for volatile organic compounds nine days after collection, two days in excess of the required holding time. Samples AM690, AM691 and AM693 were extracted for semi-volatile compounds eight days after collection, a day in excess of the required holding time. Samples AM690, AM692 and AM693 were extracted for pesticide/PCB compounds eight days after collection, a day in excess of holding time.

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Because of the high sensitivity of the volatile fraction, it is recommended to qualify positive and non-detected results for the aromatic compounds as estimated "J" and "UJ", respectively, in the volatile fraction of sample AM694. No action is required to qualify the semi-volatile or PCB fractions because exceeding the holding time by one day for these fractions is unlikely to affect the results substantially.

**Calibrations**

Initial and continuing calibration results are summarized as follows:

Volatile

<u>Compound</u>	<u>IC</u> 1/3/89	<u>CC</u> 1/9/89	<u>CC</u> 1/13/89
Tetrahydrofuran	x	x	x
Chloromethane		x	x
2-Butanone		x	
Methyl-t-butyl ether		x	x
Acetone	x	x	x
Bromoform			x
Associated samples	ALL	AM695	AM694

Semivolatile

<u>Compound</u>	<u>IC-A</u> 1/3/89	<u>CC</u> 1/9/89	<u>CC</u> 1/12/89
Isophorone	x	x	x
3,3-Dichloro-benzidine		x	
Indenopyrene		x	
Benzoic Acid			xx
Associated samples	ALL	AM690,AM692,AM693	NONE

Semivolatile cont.

<u>Compound</u>	<u>IC-B</u> 1/12/89	<u>CC</u> 1/16/89	<u>CC</u> 1/18/89
Isophorone			x
Hexachlorocyclo- pentadine			x
Benzoic Acid		x	xx
3,3-Dichlorobenzidine		x	x
Associated Samples		AM694,AM695	NONE
x	%RSD/%D criteria exceeded		
xx	%RSD/%D criteria exceeded and RF criteria exceeded		

Pesticide/PCB Calibration Results

<u>Compound</u>	<u>% D</u>	<u>Samples Affected</u>
Aldrin	16.3	A11

Blanks

Results of the Laboratory blank analysis are summarized in the following table:

<u>Compound</u>	<u>Maximum Concentration</u> <u>(ug/Kg)</u>	<u>Action Level</u> <u>(ug/Kg)</u>
Methylene Chloride	16	160
Di-n-butylphthalate	680	6800
Bis(2-ethylhexyl)phthalate	52	520
Butylbenzylphthalate	36	360
Endrin	4.3	21.5

Blank Actions:

- o value < CRQL; report CRQL followed by a U
- o value > CRQL and < action level; report value followed by U
- o value > CRQL and > action level; report value unqualified

The action level values were compared to sample values after application of sample dilution factors. From this comparison, it is recommended to

qualify positive results for di-n-butylphthalate in samples AM694 and AM695 as undetected "U", positive results for methylene chloride in all samples as undetected "U", and accept without qualification the positive result for endrin in sample AM695.

### Matrix Spike/Matrix Spike Duplicate Recoveries

The Matrix Spike/Matrix Spike Duplicate recoveries are summarized below:

#### Percent Recoveries

<u>TR#</u>	<u>Compound</u>	<u>%Recovery</u>
AM690 MS	Gamma-BHC	42
AM690 MS	Heptachlor	153
AM694 MS	Heptachlor	205
AM694 MSD	Heptachlor	203

#### Relative Percent Differences

<u>TR#</u>	<u>Compound</u>	<u>RPD</u>
AM690 MS	Gamma-BHC	53
AM690 MSD	Heptachlor	47
AM690 MS	Aldrin	53
AM690 MSD	Dieldrin	42

Based on the matrix spike/matrix spike duplicate results, it is recommended to qualify any positive results for the compounds gamma-BHC, heptachlor, aldrin or dieldrin as estimated "J" in the unspiked sample AM690 (nondetected results are acceptable).

### Field Duplicates

Sample AM959 is the duplicate of sample AM691, and sample AM692 is the duplicate of sample AM693. Field duplicate results are summarized below:

<u>Compound</u>	<u>AM959 conc.</u>	<u>AM691 conc.</u>	<u>RPD</u>
Tetrachloroethene	U	8	---

U - Undetected



May 17, 1989

In instances where one of the duplicates exhibits a positive result while the other does not it is recommended to qualify the positive result as estimated "J" and the non-detected result as estimated "UJ".

**Pesticide/PCB Instrument Performance**

Results of the DBC retention time check are summarized below:

<u>Sample</u>	<u>DBC % Difference</u>
AM694	3.2
AM695	2.6
AM694 MS	2.9

Although the results are above the two percent difference limit for packed columns, they do not warrant rejection of the data; therefore, no action is required to qualify the case.

Should you have any questions concerning this matter, do not hesitate to contact this office at (301) 587-2088.

Very truly yours,  
GEOSCIENCE CONSULTANTS, LTD.



James L. Hewitt, Jr., P.E.  
Program Manager

JLH\kt\tes3\CRON\WA99303.1tr

Attachment

Concurrence:

CDM Federal Programs Corporation



Anita Rigassio  
Environmental Scientist

MOTTOLO SITE  
CASE SAS4372  
TABLE I - RECOMMENDATIONS SUMMARY

TR#	VOLATILE	SEMI-VOLATILE	PESTICIDE/PCB
AM690	A <sup>5,8</sup>	A <sup>5,7</sup>	A <sup>3,5</sup>
AM691	A <sup>2,5,8</sup>		
AM692		A <sup>5,7</sup>	A <sup>5</sup>
AM693		A <sup>5,7</sup>	A <sup>5</sup>
AM694	A <sup>1,5,6,8</sup>	A <sup>4,5,6</sup>	A <sup>5</sup>
AM695	A <sup>5,6,8</sup>	A <sup>4,5,6</sup>	A <sup>5</sup>
AM959	A <sup>2,5,6,8</sup>		

- A - Accept all data.
- A<sup>1</sup> - Accept all data with the exception of the aromatic compounds which are qualified as estimated "UJ" due to exceedance of the holding time.
- A<sup>2</sup> - Accept all data with the exception of the compound tetrachloroethene which is qualified as estimated "UJ", if undetected, and "J", if detected, due to poor field precision results.
- A<sup>3</sup> - Accept all data with the exception of the compound gamma-BHC which is qualified as estimated "J" due to poor matrix spike results.
- A<sup>4</sup> - Accept all data with the exception of the compound di-n-butylphthalate which is qualified as undetected "U" due to blank contamination.
- A<sup>5</sup> - Accept all data with the exception of the following compounds, due to poor calibration results: acetone, tetrahydrofuran, isophorone, 3,3-dichlorobenzidine and aldrin. Qualify positive results "J" and undetected results "UJ".
- A<sup>6</sup> - Accept all data with the exception of the following compounds, due to poor calibration results: chloromethane, methyl-t-butyl ether, benzoic acid, 2-butanone (sample AM695 only) and bromoform (sample AM694 only). Qualify positive results "J" and non-detects "UJ".
- A<sup>7</sup> - Accept all data except for indenopyrene which is qualified as estimated "UJ", if undetected, and "J", if detected, due to poor calibration results.
- A<sup>8</sup> - Accept all data except for methylene chloride which is qualified as estimated "U" due to blank contamination.

MOTTOLO SITE  
CASE SAS4372

TABLE 11 - TENTATIVE IDENTIFIED COMPOUNDS (TICS) SUMMARY

TR#	AM959	AM690	AM691	AM692	AM693	AM694	AM695
COMPOUND							
3-penten-2-one,4-methyl		x		x			
phenol 2,6-bis(1,1-dimethyl)		x		x			
1,2-benzenedicarboxylic acid		x					
2-hexene 2,5-dimethyl		x					
2-penten-3,4-dimethyl		x					
3-pentanone 2,2,4,4-tetrame						x	x

wa993.tic

### DATA SUMMARY KEY

- A - Acceptable data.
- J - The associated numerical value is an estimated quantity.
- R - Reject data due to quality control criteria. The data are unusable (compound may or may not be present). Resampling and reanalysis is necessary for verification.
- U - The compound was analyzed for but was not detected. The associated numerical value is the sample quantitation limit.
- UJ- The compound was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.
- The compound was analyzed for, but was not detected. The sample quantitation limit is the same as the CRQL presented.

## Volatile

SITE: MOTTOLO

CASE: 4372

SAMPLE LOCATION:

SAMPLE NUMBER:

AM690 AM691 AM694 AM695 AM959

COMPOUND

CRQL

Chloromethane	10	11U	13U	11UJ	11UJ	12U
Bromomethane	10	11U	13U	11U	11U	12U
Vinyl Chloride	10	11U	13U	11U	11U	12U
Chloroethane	10	11U	13U	11U	11U	12U
Methylene Chloride	5	15U	15U	18U	14U	15U
Acetone	10	11UJ	5UJ	11UJ	49J	12UJ
Carbon Disulfide	5	5U	6U	6U	6U	6U
1,1-Dichloroethene	5	5U	6U	6U	6U	6U
1,1-Dichloroethane	5	5U	6U	6U	6U	6U
1,2-Dichloroethene(total)	5	5U	6U	6U	6U	6U
Chloroform	5	5U	6U	6U	6U	6U
1,2-Dichloroethane	5	5U	6U	6U	6U	6U
2-Butanone	10	11U	13U	11U	74J	5U
1,1,1-Trichloroethane	5	5U	6U	6U	6U	6U
Carbon Tetrachloride	5	5U	6U	6U	6U	6U
Vinyl Acetate	10	11U	13U	11U	11U	12U
Bromodichloromethane	5	5U	6U	6U	6U	6U
1,2-Dichloropropane	5	5U	6U	6U	6U	6U
cis-1,3-Dichloropropene	5	5U	6U	6U	6U	6U
Trichloroethene	5	5U	6U	6U	6U	6U
Dibromochloromethane	5	5U	6U	6U	6U	6U
1,1,2-Trichloroethane	5	5U	6U	6U	6U	6U
Benzene	5	5U	6U	6UJ	6U	6U
trans-1,3-Dichloropropene	5	5U	6U	6U	6U	6U
Bromoform	5	5U	6U	6UJ	6U	6U
4-Methyl-2-pentanone	10	11U	13U	11U	28	12U
2-Hexanone	10	11U	13U	11U	11U	12U
Tetrachloroethene	5	5U	8J	6U	6U	6UJ
1,1,2,2-Tetrachloroethane	5	5U	6U	6U	6U	6U
Toluene	5	5U	16	6UJ	160	10
Chlorobenzene	5	5U	6U	5UJ	6U	6U
Ethylbenzene	5	5U	6U	5UJ	330E	6U
Styrene	5	5U	6U	5UJ	6U	6U
Total Xylenes	5	5U	6U	5UJ	810E	6U
Tetrahydrofuran	5	5UJ	6UJ	5UJ	6UJ	6UJ
Methyl-t-butyl ether	5	5U	6U	5UJ	6UJ	6U

=====

DILUTION FACTOR: 1.0 1.0 1.0 1.0 1.0

DATE SAMPLED:

12/28/88 12/28/88 1/4/89 1/4/89 12/28/88

DATE ANALYZED:

1/3/89 1/3/89 1/13/89 1/9/89 1/3/89

% SOLIDS FACTOR:

0.91 0.78 0.87 0.87 0.84

## Semivolatile Soil Analysis

ug/kg

SITE: MOTTLO

CASE: 4372

SAMPLE NUMBER: AM690 AM692 AM693 AM694 AM695  
 SAMPLE LOCATION:  
 LABORATORY NUMBER:

COMPOUND	CRQL	AM690	AM692	AM693	AM694	AM695
Phenol	330	360U	380U	380U	380U	380U
bis(2-Chloroethyl) ether	330	360U	380U	380U	380U	380U
2-Chlorophenol	330	360U	380U	380U	380U	380U
1,3-Dichlorobenzene	330	360U	380U	380U	380U	380U
1,4-Dichlorobenzene	330	360U	380U	380U	380U	380U
Benzyl Alcohol	330	360U	380U	380U	380U	380U
1,2-Dichlorobenzene	330	360U	380U	380U	380U	380U
2-Methylphenol	330	360U	380U	380U	380U	380U
bis(2-Chloroisopropyl) ether	330	360U	380U	380U	380U	380U
4-Methylphenol	330	360U	380U	380U	380U	380U
N-Nitroso-di-n-propylamine	330	360U	380U	380U	380U	380U
Hexachloroethane	330	360U	380U	380U	380U	380U
Nitrobenzene	330	360U	380U	380U	380U	380U
Isophorone	330	360UJ	380UJ	380UJ	380U	380U
2-Nitrophenol	330	360U	380U	380U	380U	380U
2,4-Dimethylphenol	330	360U	380U	380U	380U	380U
Benzoic Acid	1600	1800	1900	1900	1900J	1900J
bis(2-Chloroethoxy)methane	330	360U	380U	380U	380U	380U
2,4-Dichlorophenol	330	360U	380U	380U	380U	380U
1,2,4-Trichlorobenzene	330	360U	380U	380U	380U	380U
Naphthalene	330	360U	380U	380U	380U	380U
4-Chloroaniline	330	360U	380U	380U	380U	380U
Hexachlorobutadiene	330	360U	380U	380U	380U	380U
4-Chloro-3-Methylphenol	330	360U	380U	380U	380U	380U
2-Methylnaphthalene	330	360U	380U	380U	380U	380U
Hexachlorocyclopentadiene	330	360U	380U	380U	380U	380U
2,4,6-Trichlorophenol	330	360U	380U	380U	380U	380U
2,4,5-Trichlorophenol	1600	1800U	1900U	1900U	1900U	1900U
2-Chloronaphthalene	330	360U	380U	380U	380U	380U
2-Nitroaniline	1600	1800U	1900U	1900U	1900U	1900U
Dimethylphthalate	330	360U	380U	380U	380U	380U
Acenaphthylene	330	360U	380U	380U	380U	380U
2,6-Dinitrotoluene	330	360U	380U	380U	380U	380U
3-Nitroaniline	1600	1800U	1900U	1900U	1900U	1900U
Acenaphthene	330	360U	380U	380U	380U	380U
2,4-Dinitrophenol	1600	1800U	1900U	1900U	1900U	1900U
4-Nitrophenol	1600	1800U	1900U	1900U	1900U	1900U
Dibenzofuran	330	360U	380U	380U	380U	380U
2,4-Dinitrotoluene	330	360U	380U	380U	380U	380U
Diethylphthalate	330	360U	380U	380U	380U	380U
4-Chlorophenyl-Phenylether	330	360U	380U	380U	380U	380U
Fluorene	330	360U	380U	380U	380U	380U
4-Nitroaniline	1600	1800U	1900U	1900U	1900U	1900U
4,6-Dinitro-2-Methylphenol	1600	1800U	1900U	1900U	1900U	1900U
N-nitrosodiphenylamine(1)	330	360U	380U	380U	380U	380U
4-Bromophenyl-Phenylether	330	360U	380U	380U	380U	380U

Semivolatile Results Cont.

SITE: MOTTOLO

CASE: 4372

SAMPLE NUMBER: AM690 AM692 AM693 AM694 AM695  
 SAMPLE LOCATION:  
 LABORATORY NUMBER:

COMPOUND	CRQL	AM690	AM692	AM693	AM694	AM695
Hexachlorobenzene	330	360U	380U	380U	380U	380U
Pentachlorophenol	1600	1800U	1900U	1900U	1900U	1900U
Phenanthrene	330	360U	380U	380U	380U	380U
Anthracene	330	360U	380U	380U	380U	380U
Di-n-butylphthalate	330	360U	430U	430U	1200U	1200U
Fluoranthene	330	360U	380U	380U	380U	380U
Pyrene	330	360U	380U	380U	380U	380U
Butylbenzylphthalate	330	360U	380U	380U	380U	380U
3,3'-Dichlorobenzidine	660	730UJ	770UJ	770UJ	760UJ	760UJ
Benzo(a)anthracene	330	360U	380U	380U	380U	380U
Chrysene	330	360U	380U	380U	380U	380U
Bis(2-ethylhexyl)phthalate	330	360U	380U	380U	380U	380U
Di-n-octylphthalate	330	360U	380U	380U	380U	380U
Benzo(b)fluoranthene	330	360U	380U	380U	380U	380U
Benzo(k)fluoranthene	330	360U	380U	380U	380U	380U
Benzo(a)pyrene	330	360U	380U	380U	380U	380U
Indeno(1,2,3-cd)pyrene	330	360UJ	380UJ	380UJ	380U	380U
Dibenz(a,h)anthracene	330	360U	380U	380U	380U	380U
Benzo(g,h,i)perylene	330	360U	380U	380U	380U	380U

=====

DILUTION FACTOR:	1.00	1.00	1.00	1.00	1.00
DATE SAMPLED:	12/28/88	12/28/88	12/28/88	01/04/89	01/04/89
DATE EXTRACTED:	01/05/89	01/05/89	01/05/89	01/11/89	01/11/89
DATE ANALYZED:	01/09/89	01/09/89	01/09/89	01/16/89	01/16/89
% SOLIDS FACTOR:	0.81	0.86	0.88	0.87	0.87

Pesticide/PCB Soil Analysis  
ug/kg

SITE: MOTTOLO

CASE: 4372

SAMPLE NUMBER: AM690 AM692 AM693 AM694 AM695  
SAMPLE LOCATION:

COMPOUND	CRQL	AM690	AM692	AM693	AM694	AM695
alpha-BHC	8.0	8.8U	9.3U	9.1U	9.2U	9.2U
beta-BHC	8.0	8.8U	9.3U	9.1U	9.2U	9.2U
delta-BHC	8.0	8.8U	9.3U	9.1U	9.2U	9.2U
gamma-BHC(Lindane)	8.0	.036J	7.7U	9.1U	9.2U	9.2U
Heptachlor	8.0	8.8U	9.3U	.13J	9.2U	9.2U
Aldrin	8.0	8.8UJ	9.3UJ	9.1UJ	9.2UJ	9.2UJ
Heptachlor Epoxide	8.0	8.8U	9.3U	9.1U	9.2U	9.2U
Endosulfan I	8.0	8.8U	9.3U	9.1U	9.2U	9.2U
Dieldrin	16.0	18U	19U	18U	18U	18U
4,4'-DDE	16.0	18U	19U	18U	18U	18U
Endrin	16.0	18U	19U	18U	18U	31
Endosulfan II	16.0	18U	19U	18U	18U	18U
4,4'-DDD	16.0	18U	19U	18U	18U	18U
Endosulfan Sulfate	16.0	18U	19U	18U	18U	18U
4,4'-DDT	16.0	18U	19U	18U	18U	18U
Methoxychlor	80.0	88U	93U	91U	92U	92U
Endrin Ketone	16.0	18U	19U	18U	18U	18U
alpha-Chlordane	80.0	88U	93U	91U	92U	92U
gamma-Chlordane	80.0	88U	93U	91U	92U	92U
Toxaphene	160.0	180U	190U	180U	180U	180U
Aroclor-1016	80.0	88U	93U	91U	92U	92U
Aroclor-1221	80.0	88U	93U	91U	92U	92U
Aroclor-1232	80.0	88U	93U	91U	92U	92U
Aroclor-1242	80.0	88U	93U	91U	92U	92U
Aroclor-1248	80.0	88U	93U	91U	92U	92U
Aroclor-1254	160.0	180U	190U	180U	180U	180U
Aroclor-1260	160.0	180U	190U	180U	180U	180U

=====

DILUTION FACTOR:	1.00	1.00	1.00	1.00	1.00
DATE SAMPLED:	12/28/88	12/28/88	12/28/88	1/4/89	1/4/89
DATE EXTRACTED:	1/5/89	1/5/89	1/5/89	1/11/89	1/11/89
DATE ANALYZED:	2/2/89	2/2/89	2/2/89	2/2/89	2/2/89
% SOLIDS FACTOR:	0.89	0.86	0.88	0.87	0.87



REGION I  
Data Review Worksheets

Site Name MOTTOLO  
Reference Number \_\_\_\_\_

REGION I REVIEW OF ORGANIC  
CONTRACT LABORATORY DATA PACKAGE

The hardcopied (laboratory name) E. C. JORDAN CO. data package received at Region I has been reviewed and the quality assurance and performance data summarized. The data review included:

Case No. \_\_\_\_\_ SAS No. 4372 Sampling Date(s) 12/28/88 and 1/4/89  
SDG No. AM690 Matrix SOIL Shipping Date(s) 12/29/88 1/5/89  
No. of Samples 7 Date Rec'd by Lab 2/31/88 1/6/89

Traffic Report Nos: AM690 AM959 AM691 AM692 AM693  
AM694 AM695  
Trip Blank No.: NONE  
Equipment Blank No.: NONE  
Field Dup Nos: AM959/AM691 AM692/AM693

SOW No. \_\_\_\_\_ requires that specific analytical work be done and that associated reports be provided by the laboratory to the Regions, EMSL-LV, and SMO. The general criteria used to determine the performance were based on an examination of:

- |                       |                                |
|-----------------------|--------------------------------|
| -Data Completeness    | -Matrix Spike/Matrix Spike Dup |
| -Holding Times        | -Field Duplicates              |
| -GC/MS Tuning         | -Internal Standard Performance |
| -Calibrations         | -Pesticide Inst. Performance   |
| -Blanks               | -Compound Identification       |
| -Surrogate Recoveries | -Compound Quantitation         |

Overall comments The samples were received by the lab in good condition.

Definitions and Qualifiers:

- A - Acceptable data.  
J - Approximate data due to quality control criteria.  
R - Reject data due to quality control criteria.  
U - Compound not detected.

Reviewer: Keith Trehy Date: 3/27/89



REGION I  
Data Review Worksheets

II. HOLDING TIMES

Complete table for all samples and circle the fractions which are not within criteria.

SAMPLE ID	DATE SAMPLED	VOA	BNA		PEST	
		DATE ANAL	DATE EXTR	DATE ANAL	DATE EXTR	DATE ANAL
AM690	12/28/88	1/3/89	1/5/89	1/9/89	1/5/89	2/2/89
AM691	12/28/88	1/3/89	1/5/89	1/9/89		
AM694	1/4/89	1/13/89	1/11/89	1/16/89	1/11/89	2/2/89
AM695	1/4/89	1/9/89	1/11/89	1/16/89	1/11/89	2/2/89
AM959	12/28/88	1/3/89				
AM693	12/28/88		1/5/89	1/9/89	1/5/89	2/2/89
AM692	12/28/88				1/5/89	2/2/89
A.....						

VOA - Unpreserved: Aromatic within 7 days, non-aromatic within 14 days of sample collection.  
 Preserved : Both within 14 days of sample collection.  
 Soils : Both within 14 days of sample collection.

BNA & PEST - Extracted within 7 days, analyzed within 40 days, soils and water.

ACTION:

1. If holding times are exceeded all positive results are estimate (J) and non-detects are estimated (UJ).
2. If holding times are grossly exceeded, the reviewer may determine that non-detects are unusable(\*)

REGION I  
Data Review Worksheets

III. GC/MS TUNING

The DFTPP performance results were reviewed and found to be within the specified criteria.

If no,  
Samples affected: \_\_\_\_\_

The BFB performance results were reviewed and found to be within the specified criteria.

If no,  
Samples affected: \_\_\_\_\_

If mass calibration is in error refer to the Region guidelines for expanded criteria. If necessary, all associated data as unusable (R).

REGION I  
Data Review Worksheets

IV A. VOLATILE CALIBRATION VERIFICATION

Date of Initial Calibration : 1/3/89  
 Dates of Continuing Calibrations: 1/9, 12/89  
 Instrument ID : 960202  
 Matrix/Level : SOIL/LOW

DATE	CRITERIA OUT RF, %RSD, RF, %D % RSD	COMPOUND (VALUE)
<u>1/3/89</u>	Samples Affected: <u>ALL</u>	<u>Acetone (82), Tetrahydrofuran (74.3)</u> <u>Methyl-t-butyl ether (46.6)</u>
<u>1/9/89</u>	Samples Affected: <u>AM695</u>	<u>Chloromethane (51.7), 2-Butanone (30.5), Tetrahydrofuran (55.2)</u>
<u>1/13/89</u>	Samples Affected: <u>AM694</u>	<u>Chloromethane (48.9), Acetone (37.9), Bromoform (30.5)</u>
<u>1/13/89</u>	Samples Affected: <u>AM694</u>	<u>Tetrahydrofuran (75.7), Methyl-t-butyl ether (36.4)</u>
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____

- All RF's, and RF's must be >0.05
- All %RSD's must be <30%
- All %D's must be <25%

ACTION:

- If any compound has an initial RF or a continuing RF of <0.05:
  - Flag positive results for that compound as estimated (J).
  - Flag non-detects for that compound as unusable (R).
- If any compound has a %RSD >30% or a %D >25%:
  - Flag positive results for that compound as estimated (J).
  - Flag non-detects for that compound as estimated (UJ) if the %RSD or %D is >50%.

A separate worksheet should be filled out for each initial curve.

REGION I  
Data Review Worksheets

IV B. SEMIVOLATILE CALIBRATION VERIFICATION

Date of Initial Calibration : 1/3/89  
Dates of Continuing Calibrations: 1/9, 12/89  
Instrument ID : 870101

<u>DATE</u>	<u>CRITERIA OUT</u> RF, %RSD, RF, %D <u>% RSD</u>	<u>COMPOUND</u>
<u>1/3/89</u>	<u>0% RSD</u> Samples Affected: <u>ALL</u>	<u>Isophorone (33.6)</u>
<u>1/12/89</u>	<u>% D</u> Samples Affected: <u>NONE</u>	<u>Isophorone (36.0) Benzoic Acid (74.8)</u>
<u>1/12/89</u>	<u>RF</u> Samples Affected: <u>NONE</u>	<u>Benzoic Acid (0.033)</u>
<u>1/9/89</u>	<u>% D</u> Samples Affected: <u>AM690, AM692, AM693</u>	<u>3,3-Dichlorobenzidine (54.8), Indenopyrene (34)</u>
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____
_____	Samples Affected: _____	_____

See worksheet IV-A for criteria and actions.  
A new worksheet should be filled out for each initial curve.

REGION I  
Data Review Worksheets

IV B. SEMIVOLATILE CALIBRATION VERIFICATION

Date of Initial Calibration : 1/12/89  
Dates of Continuing Calibrations: 1/16, 18/89  
Instrument ID : 870101

DATE	CRITERIA OUT RF, %RSD, RF, %D	COMPOUND
<u>1/16/89</u>	<u>% D</u> Samples Affected: <u>AM694, AM695</u>	<u>Benzoic acid (60.2), 3,3-Dichlorobenzidine (30.4)</u>
<u>1/18/89</u>	<u>% D</u> Samples Affected: <u>NONE</u>	<u>Isophorone (35.4), Hexachlorocyclopentadiene (25.4)</u>
<u>1/18/89</u>	<u>RF</u> Samples Affected: <u>NONE</u>	<u>Benzoic acid (.038)</u>
<u>1/18/89</u>	<u>% D</u> Samples Affected: <u>NONE</u>	<u>3,3-Dichlorobenzidine (38.5)</u>
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____
_____	_____ Samples Affected: _____	_____

See worksheet IV-A for criteria and actions.  
A new worksheet should be filled out for each initial curve.

REGION I  
Data Review Worksheet

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

List the contamination in the blanks below.

1. Laboratory Blanks

Level: LOW

<u>DATE</u>	<u>LAB ID</u>	<u>FRACTION/ MATRIX</u>	<u>COMPOUND</u>	<u>CONCENTRATION/ UNITS</u>
1/3/89	VBLK01	VOA/SOIL	Methylene Chloride	11 ug/kg
1/9/89	VBLK02	VOA/SOIL	Methylene Chloride	13 ug/kg
1/13/89	VBLK03	VOA/SOIL	Methylene Chloride	16 ug/kg
1/9/89	SBLK01	SVOA/SOIL	Di-n-Butylphthalate	450 ug/kg
1/9/89	SBLK01	SVOA/SOIL	Di-(2-Ethylhexyl) Phthalate	52 ug/kg
1/16/89	SBLK02	SVOA/SOIL	Di-n-Butylphthalate	680 ug/kg
1/16/89	SBLK02	SVOA/SOIL	Butylbenzylphthalate	36 ug/kg
2/2/89	PBLKAA	PCB/SOIL	Endrin	4.3 ug/kg
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

2. Equipment and Trip Blanks NONE

<u>DATE</u>	<u>TR #</u>	<u>FRACTION/ MATRIX</u>	<u>COMPOUND</u>	<u>CONCENTRATION/ UNITS</u>
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

A separate worksheet should be used for low and medium level blanks.



REGION I  
Data Review Worksheets

V B. BLANK ANALYSIS RESULTS (Section 3)

3. Blank Actions

Action levels should be based upon the highest concentration of contaminant determined in any blank. The action level for samples which have been concentrated or diluted should be multiplied by the concentration/dilution factor. No positive sample result should be reported unless the concentration of the compound in the sample exceeds the action level of 10 x's the amount in the blank for the common contaminants, or 5 x's the amount for any other compound. Specific actions are as follows:

1. The concentration is less than the CRQL, report the CRQL.
2. The concentration is greater than the CRQL, but less than the action level, report the concentration found U.
3. The concentration is greater than the action level, report the concentration unqualified.

For examples refer to the Regional Guidelines.

Common contaminants = methylene chloride, acetone, 2-butanone, toluene, and phthalates.

LEVEL: Low

<u>COMPOUND</u>	<u>MAX. CONC./</u> <u>UNITS</u>	<u>ACTION LEVEL/</u> <u>UNITS</u>	<u>CRQL</u>
Methylene Chloride	16 ug/kg	160 ug/kg	10
Di-n-butylphthalate	680 ug/kg	6800 ug/kg	330
Bis(2-ethylhexyl)phthalate	52 ug/kg	520 ug/kg	330
Butylbenzylphthalate	36 ug/kg	360 ug/kg	330
Endrin	4.3 ug/kg	21.5 ug/kg	16

A separate worksheet should be used for low and medium level blanks.

REGION I  
Data Review Worksheets

VI. SURROGATE SPIKE RECOVERIES

List the percent recoveries which do not meet the criteria for surrogate recovery.

*All surrogate recovery results were within criteria*

Matrix: Soil

TR #'S	TOL	<i>OKAY</i>	DCF	NBZ	<i>OK</i>	TPH	PHL	<i>OKAY</i>	TBP	DBC
		VOA			B/N			A		
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
_____	_____		_____	_____		_____	_____		_____	_____
QC Limits										
	to	to	to	to	to	to	to	to	to	to

Surrogate Actions: \*-Advisory only

	<u>PERCENT RECOVERY</u>		
	<u>&lt;10%</u>	<u>10%-CRR</u>	<u>&gt;CRR</u>
Positive sample results	J	J	J
Non-detected results	R	UJ	A

- CRR = Contract Required Recovery Range.  
 Surrogate action should be applied:
- If at least two surrogates in a B/N or A fraction or one surrogate in the VOA fraction are out of specification, but have recoveries of >10%.
  - If any one surrogate in a fraction shows <10% recovery.

REGION I  
Data Review Worksheets

VII A. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

1. Matrix Spike/Matrix Spike Duplicate Recoveries and Precision

TR Nos. AM690 , \_\_\_\_\_ Level: LOW Matrix: SOIL

List the percent recoveries and RPD's of compounds which do not meet the criteria stated on Form 3.

<u>FRACTION/ MS OR MSD</u>	<u>COMPOUND</u>	<u>%REC/ RPD</u>	<u>QC LIMITS</u>
<u>Pest/MS</u>	<u>Gamma-BHC</u>	<u>42/53</u>	<u>46-127/50</u>
<u>Pest/MSD</u>	<u>Heptachlor</u>	<u>153/47</u>	<u>35-130/31</u>
<u>↓</u>	<u>Aldrin</u>	<u>53</u>	<u>43</u>
<u>↓</u>	<u>Dieldrin</u>	<u>42</u>	<u>38</u>

QUALIFICATION IS LIMITED TO THE UNSPIKED SAMPLE ONLY.

1. If any compound does not meet the Contract Required Recovery range (CRR) follow the actions stated below:

	<u>PERCENT RECOVERY</u>		
	<u>&lt;10%</u>	<u>10% - CRR</u>	<u>&gt;CRR</u>
Positive Sample Results	J	J	J
Non-detected Results	R	A	A

2. If any compound does not meet the RPD criteria, flag positive results for that compound as estimated (J).

A separate worksheet should be used for each MS/MSD pair.

REGION I  
Data Review Worksheets

VII A. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

1. Matrix Spike/Matrix Spike Duplicate Recoveries and Precision

TR NOS. AM694 , \_\_\_\_\_ Level: LOW Matrix: SOIL

List the percent recoveries and RPD's of compounds which do not meet the criteria stated on Form 3.

<u>FRACTION/ MS OR MSD</u>	<u>COMPOUND</u>	<u>%REC/ RPD</u>	<u>QC LIMITS</u>
<u>Pest/MS</u>	<u>Heptachlor</u>	<u>205</u>	<u>35-130</u>
<u>Pest/MSD</u>	<u>Heptachlor</u>	<u>203</u>	<u>35-130</u>

QUALIFICATION IS LIMITED TO THE UNSPIKED SAMPLE ONLY.

1. If any compound does not meet the Contract Required Recovery range (CRR) follow the actions stated below:

	<u>PERCENT RECOVERY</u>		
	<u>&lt;10%</u>	<u>10% - CRR</u>	<u>&gt;CRR</u>
Positive Sample Results	J	J	J
Non-detected Results	R	A	A

2. If any compound does not meet the RPD criteria, flag positive results for that compound as estimated (J).

A separate worksheet should be used for each MS/MSD pair.

REGION I  
Data Review Worksheets

VII B. MATRIX SPIKE/MATRIX SPIKE DUPLICATE (Section 2)

3. Matrix Spike Duplicate - Unspiked Compounds

TR Nos. AM690 , AM694

List the concentrations of the unspiked compounds and determine the percent RSD's of the unspiked sample, matrix spike, and matrix spike duplicate. No limits have been developed for the RSD values of the unspiked compounds. .

<u>FRACTION</u>	<u>COMPOUND</u>	<u>SAMPLE, MS, MSD CONC</u>	<u>%RSD</u>
AM690 VOA	methyrene chloride	15 , 16 , 15	3.7
AM694 SVOA	Di-n-butyl phthalate	1200 , 290 , 300	87.5
AM690 Pest	Endrin	160 , 230 , 12	83

No action is required based on these results.

The reviewer must use professional judgement to determine if there is a need to qualify any of the unspiked compounds in the sample.

REGION I  
Data Review Worksheets

VIII. FIELD DATA POINT PRECISION

TR NO: AM959 , AM691

MATRIX: SOIL

List the concentrations of the compounds which do not meet the following RPD criteria:

1. An RPD of <30% for water duplicates.
2. An RPD of <50% for soil duplicates.

<u>FRACTION</u>	<u>COMPOUND</u>	<u>SAMPLE CONC</u>	<u>DUP SAMPLE CONC</u>	<u>RPD</u>
<u>VOA</u>	<u>tetrachloroethene</u>	<u>U</u>	<u>6</u>	<u>-</u>

ACTIONS: Flag AM959 compound (US) , AM691 compound (J)

1. If the results for any compounds do not meet the RPD criteria, flag the positive results for that compound as estimated.
2. If one value is non-detected, and one is above the CRQL:
  - a. Flag the positive result as estimated (J).
  - b. Flag the non-detected result as estimated (UJ).

NOTE: Professional judgement may be utilized to apply duplicate action to all samples of a similar matrix.

A separate worksheet should be filled out for each field duplicate pair.

REGION I  
Data Review Worksheets

IX. INTERNAL STANDARD PERFORMANCE

All samples are within *Internal standard criteria.*

List the internal standard areas of samples which do not meet the criteria of +100% or -50% of the internal standard area in the associated continuing calibration standard.

<u>SAMPLE ID</u>	<u>DATE</u>	<u>IS OUT</u>	<u>IS AREA/ RT</u>	<u>ACCEPTABLE RANGE</u>	<u>ACTION</u>
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____
_____	_____	_____	_____	_____	_____

**ACTION:**

1. If an IS area count is outside the criteria -50% or +100% of the associated standard:
  - a. Positive results for compounds quantitated using that IS are flagged as estimated (J) for that sample fraction.
  - b. Non-detects for compounds quantitated using that IS are flagged as estimated (UJ) for that sample fraction.
  - c. If extremely low area counts are reported, or if performance exhibits a major drop-off, then a severe loss of sensitivity is indicated. Non-detects should then be flagged as unusable (R).
2. If an IS retention time varies more than 30 seconds, the chromatographic profile for that sample must be examined to determine if any false positives or negatives exist. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for that sample fraction.

REGION I  
Data Review Worksheets

X A. PESTICIDE INSTRUMENT PERFORMANCE (Section 1)

*pesticide results met criteria*

1. DDT Retention Time *okay*

List the DDT standards which have a retention time (RT) of less than 12 minutes on the packed column (except OV-1 or OV-101).

<u>STANDARD ID</u>	<u>DATE/ TIME</u>	<u>RT</u>	<u>SAMPLES AFFECTED</u>	<u>ACTIONS</u>
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

ACTION:

If the RT is less than 12 minutes, examine the chromatography to evaluate the separation. If adequate separation is not achieved, flag all affected compound data as unusable (R).



REGION 1  
Data Review Worksheets

X B. PESTICIDE INSTRUMENT PERFORMANCE (Section 2)

2. Retention Time Windows *will in numbers*  
*Retention times were all within criteria*

List the compounds which are not within the established windows.

<u>COMPOUND</u>	<u>DATE</u> <u>(TIME)</u>	<u>RT</u>	<u>RT WINDOW</u>	<u>SAMPLES AFFECTED</u>
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

Check the sample chromatograms of the samples analyzed after the last in control standard for peaks within an expanded window. If no peaks are present, there is usually no effect on the data. Refer to Regional guidelines for information on qualifying data if peaks are present. If peaks are present, discuss actions below:

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

REGION I  
Data Review Worksheets

X C. PESTICIDE INSTRUMENT PERFORMANCE (Section 3)

Results are within criteria

3. DDT and Endrin Degradation *only*

List the standards which have a DDT or Endrin breakdown of greater than 20%.

<u>STANDARD ID</u>	<u>DDT OR ENDRIN</u>	<u>PERCENT BREAKDOWN</u>	<u>SAMPLES AFFECTED</u>	<u>DDD, DDE OR ENDRIN KETONE PRESENT</u>
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

If the percent breakdown for DDT is greater than 20%:

1. Flag all positive results for DDT as estimated (J) for all samples following the last in control standard. If no DDT was present, but DDD and/or DDE are positive, then flag the quantitation limit for DDT as unusable (R).
2. Flag all positive results for DDD +/- or DDE as estimated (J).

If the percent breakdown for Endrin is greater than 20%:

1. Flag all positive results for endrin as estimated (J) for all samples following the last in-control standard. If no endrin was detected, but endrin aldehyde and/or endrin ketone are positive, flag the quantitation limit for endrin as unusable (R).
2. Flag all positive results for endrin ketone as estimated (J).

REGION I  
Data Review Worksheets

X D. PESTICIDE INSTRUMENT PERFORMANCE (Section 4)

4. DBC Retention Time Check

List the percent difference for the DBC shift greater than 2% for packed columns, greater than 1.5% for wide-bore capillary columns, or greater than 0.3% for narrow-bore capillary columns.

<u>TR #'S</u>	<u>DBC %DIFFERENCE</u>	<u>ACTIONS</u>
<u>AM694</u>	<u>3.2</u>	<u></u>
<u>AM695</u>	<u>2.6</u>	<u></u>
<u>AM694MS</u>	<u>2.9</u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>
<u></u>	<u></u>	<u></u>

If the DBC does not meet the retention time criteria, the analysis may be flagged as unusable (R) for the affected samples, but qualification of the data is left up to the professional judgement of the reviewer. Discuss any qualification of the data below:

These few results slightly above the criteria do not justify  
rejection of the data.

REGION I  
Data Review worksheets

XI A. PESTICIDE CALIBRATION (Sections 1 and 2)

1. Initial Calibration

*Initial calibration results are within criteria.*

List the compounds which did not meet the Relative Standard Deviation (RSD) criteria of less than 10% for the initial calibration on the quantitation column.

<u>DATE</u>	<u>COMPOUND</u>	<u>%RSD</u>	<u>COLUMN</u>	<u>SAMPLES AFFECTED</u>
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____
_____	_____	_____	_____	_____

Flag all associated positive results as estimated (J) for samples which did not meet the %RSD criteria.

2. Analytical Sequence

Did the laboratory follow the correct 72 hour sequence described in the SOW? Yes or No *Yes*

If no,

The data may be affected. The data reviewer must use professional judgement to determine the severity of the effect and qualify the data accordingly. Discuss any actions below:

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

REGION I  
Data Review Worksheets

XI B. PESTICIDE CALIBRATION (Section 3)

3. Continuing Calibration

List the compounds which did not meet the percent difference (%D) criteria of <15% on the quantitation column or <20% on the confirmation for the continuing calibration.

<u>DATE</u>	<u>COMPOUND</u>	<u>%D</u>	<u>COLUMN</u>	<u>SAMPLES AFFECTED</u>
2/2/89	Aldrin	16.3	1	NONE

If the %D criteria is not met, flag all associated positive results as estimated (J).

REGION I  
Data Review Worksheets

XII. SAMPLE QUANTITATION

In the space below, please show a minimum of one sample calculation fraction:

VOA: AM691 Toluene soil

$$\frac{4267 \quad 50}{20596 (1.78) (0.829)} = 16.0$$

BNA: AM692 bis(2-ethylhexylphthalate)

$$\frac{6826 \quad 40}{44105 (1.599) (0.025) (0.86)} = 180$$

PEST/PCB:

AM695 Endrin

$$\frac{47274 \times 1 \mu\text{g} \times 20,000 \mu\text{l}}{233654 \times 5 \mu\text{l} \times 30 \text{g} \times 0.87} = 31 \frac{\mu\text{g}}{\text{kg}}$$