User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/25/2002 Time: 12:19:19

Sample: AE07178
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/25/02 12:18 Ended: 4/25/02 12:18 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		114		10.0

Comments for Sample AE07178 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-25-2002 Time: 12:19:49

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\041802\7178.D Vial: 15
Acq On : 18 Apr 2002 8:18 pm Operator: Bohlk
Sample : FB, SAMPLE 7178 3/27/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 24 08:45:31 2002 Quant Results File: 5080402BBC.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 08:08:47 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.41	152	420562	40.00	ug/L	0.00
10) Naphthalene-d8	12.89	136	2030913	40.00	ug/L	0.00
17) Acenaphthene-d10	17.99	164	1148421	40.00	ug/L	0.00
30) Phenanthrene-d10	22.34	188	1635393	40.00	ug/L	0.00
39) Chrysene-d12	30.14	240	1358514	40.00	ug/L	-0.01
45) Perylene-d12	34.01	264	877849	40.00	ug/L	0.00
System Monitoring Compounds						
28) TCMX (SURR)	19.95	244	20880	4.56	ug/L	0.00
Spiked Amount 4.000			Recovery = 114.00%			
38) 2,4-DDD (SURR)	0.00	235	0	0.00	ug/L	
Spiked Amount 10.000			Recovery = 0.00%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D.		
3) bis(2-chloroethyl) ether	0.00	93	0	N.D.		
4) 1,3 - Dichlorobenze	0.00	146	0	N.D.		
5) 1,4 - Dichlorobenzene	0.00	146	0	N.D.		
6) 1,2 - Dichlorobenzene	0.00	146	0	N.D.		
7) 2,2 ' - oxybis (1-Chloropr	0.00	45	0	N.D.		
8) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.		
9) Hexachloroethane	0.00	117	0	N.D.		
11) Nitrobenzene	0.00	77	0	N.D.		
12) Isophorone	0.00	82	0	N.D.		
13) bis(2-Chloroethoxy) methan	0.00	93	0	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Naphthalene	0.00	128	0	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
19) 2-chloronaphthalene	0.00	162	0	N.D.		
20) Dimethylphthalate	0.00	163	0	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
25) Diethylphthalate	0.00	149	0	N.D.		
26) 4-Chlorophenyl-phenyl ethe	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	0.00	169	0	N.D.		
32) 4-Bromophenyl-phenyl ether	0.00	248	0	N.D.		
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	0.00	178	0	N.D.		
35) Anthracene	0.00	178	0	N.D.		
36) Di-n-butylphthalate	0.00	149	0	N.D.		
37) Fluoranthene	0.00	202	0	N.D.		
40) Pyrene	0.00	202	0	N.D.		
41) Butylbenzylphthalate	0.00	149	0	N.D.		
42) Benzo (a) antracene	0.00	228	0	N.D.	d	
43) Bis (2-ethylhexyl) phthala	30.76	149	64185	1.91	ug/L	94
44) Chrysene	0.00	228	0	N.D.	d	

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\041802\7178.D Vial: 15
Acq On : 18 Apr 2002 8:18 pm Operator: Bohlk
Sample : FB, SAMPLE 7178 3/27/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 24 08:45:31 2002 Quant Results File: 5080402BBC.RES

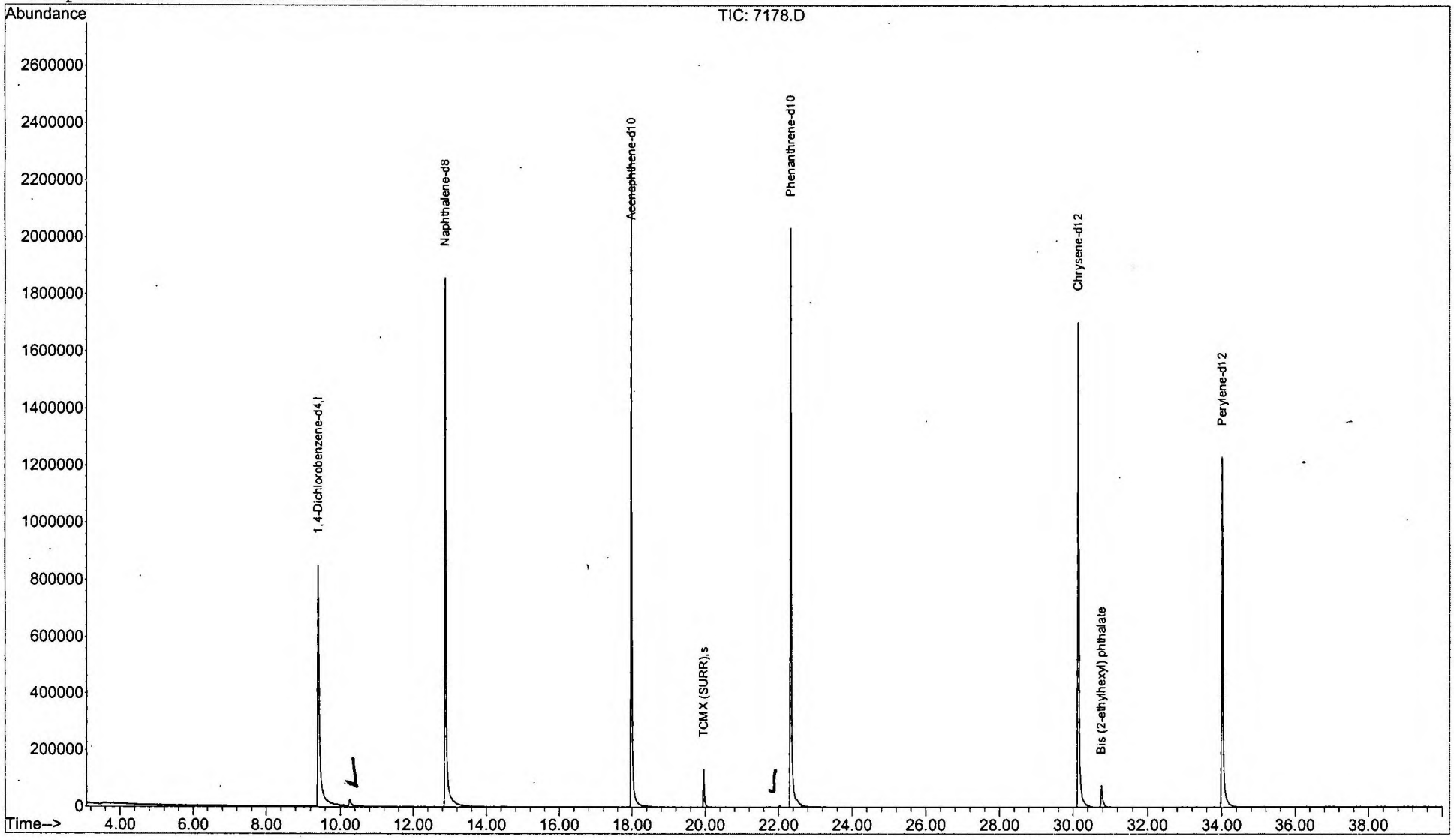
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 08:08:47 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
46)	Di-n-Octylphthalate	0.00	149	0	N.D.	
47)	Benzo (b) fluoranthene	0.00	252	0	N.D.	
48)	Benzo (k) fluoranthene	0.00	252	0	N.D.	
49)	Benzo (a) pyrene	0.00	252	0	N.D.	d
50)	Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.	
51)	Dibenz (a,h) anthracene	0.00	278	0	N.D.	
52)	Benzo (g,h,i) perylene	0.00	276	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Data File : C:\MSDCHEM\1\DATA\041802\7178.D Vial: 15
Acq On : 18 Apr 2002 8:18 pm Operator: Bohlk
Sample : FB, SAMPLE 7178 3/27/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 24 8:46 2002 Quant Results File: 5080402BBC.RES

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 08:08:47 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\041802\7178.D Vial: 15
Acq On : 18 Apr 2002 8:18 pm Operator: Bohlk
Sample : FB, SAMPLE 7178 3/27/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

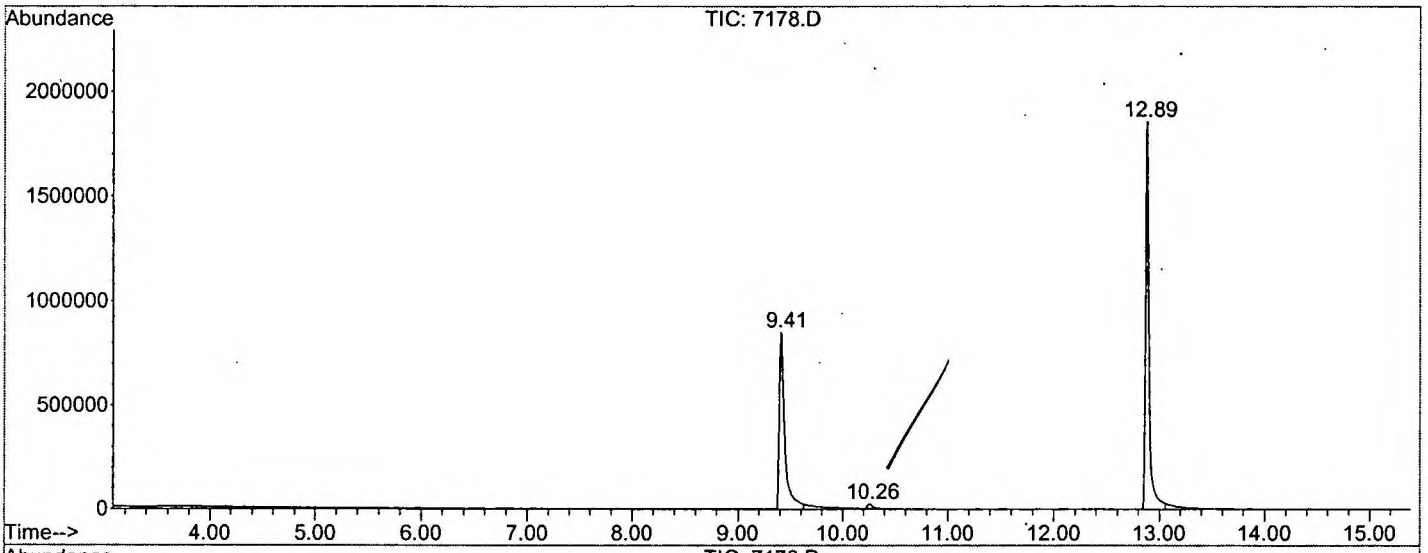
Signal : TIC

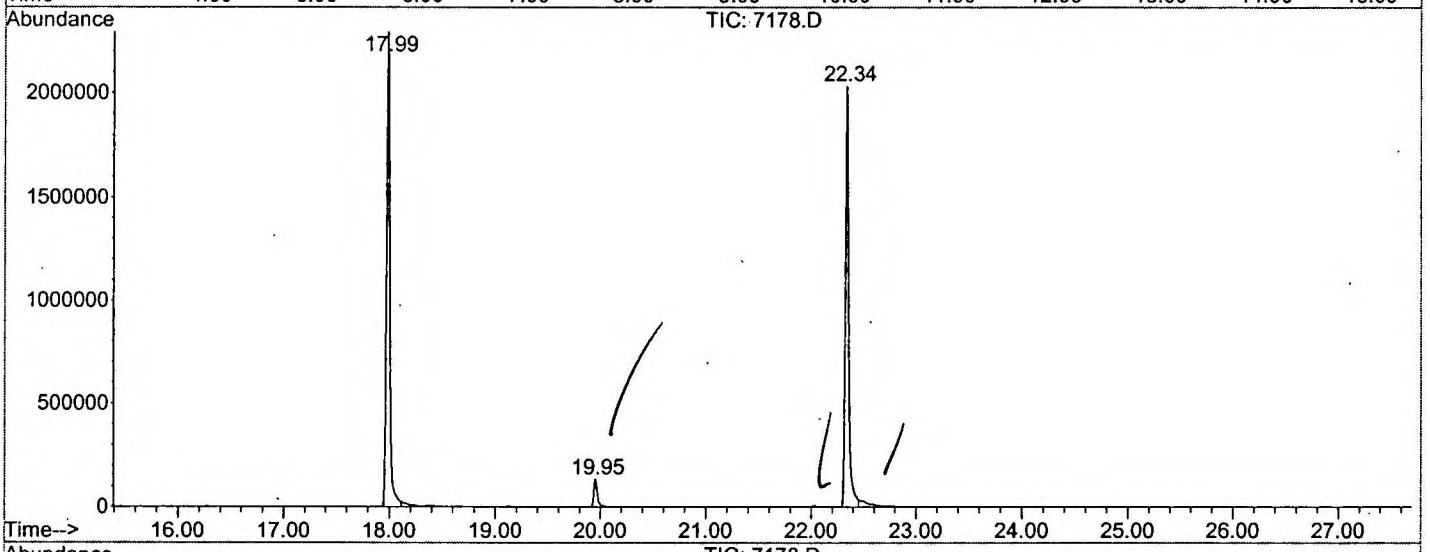
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.407	1071	1077	1114	rBV	849089	2976480	60.82%	11.937%
2	10.259	1215	1222	1236	rBV3	20869	63358	1.29%	0.254%
3	12.885	1661	1669	1698	rBV	1857421	4413112	90.18%	17.699%
4	17.985	2529	2537	2558	rBV	2291031	4893519	100.00%	19.625%
5	19.948	2864	2871	2889	rBV3	134286	312153	6.38%	1.252%
6	22.339	3269	3278	3297	rBV2	2033528	4660313	95.23%	18.690%
7	30.142	4597	4606	4625	rBV2	1702476	4135853	84.52%	16.587%
8	30.765	4705	4712	4731	rBV2	76821	242403	4.95%	0.972%
9	34.014	5256	5265	5293	rBV2	1230011	3237699	66.16%	12.985%

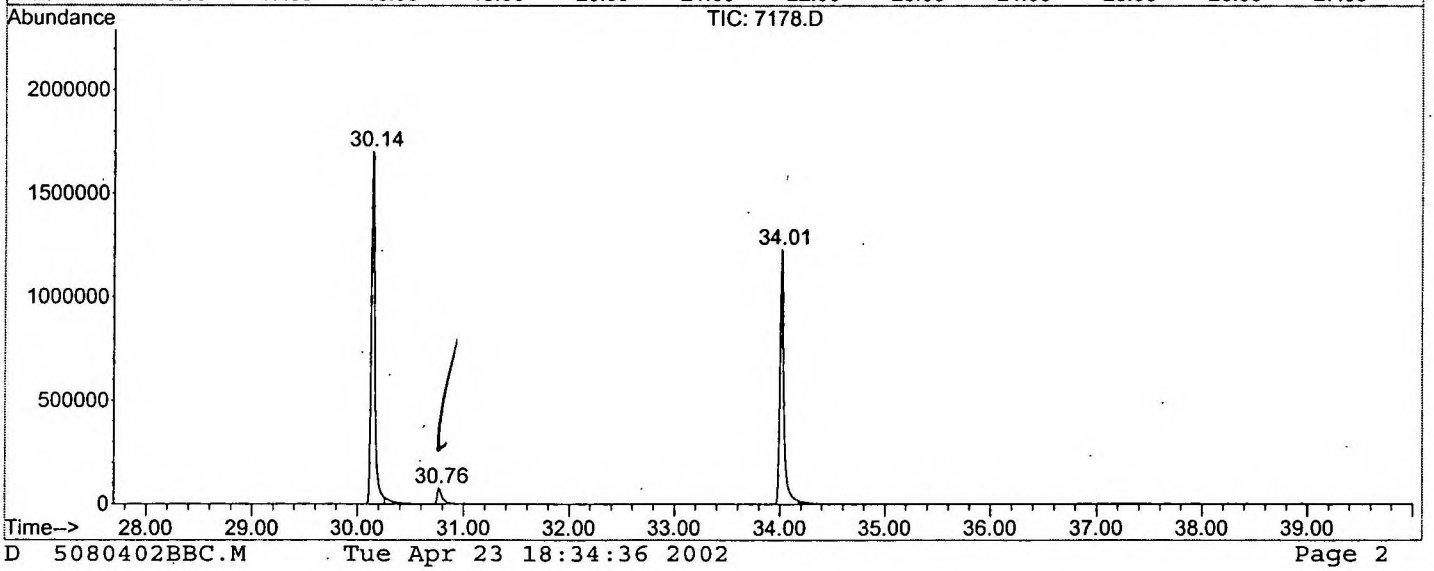
Sum of corrected areas: 24934890

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\041802\7178.D
Operator : Bohlk
Acquired : 18 Apr 2002 8:18 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: FB, SAMPLE 7178 3/27/02
Misc Info :
Vial Number: 15
Quant File :5080402BBC.RES (RTE Integrator)







Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\041802\7178.D Vial: 15
Acq On : 18 Apr 2002 8:18 pm Operator: Bohlk
Sample : FB, SAMPLE 7178 3/27/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

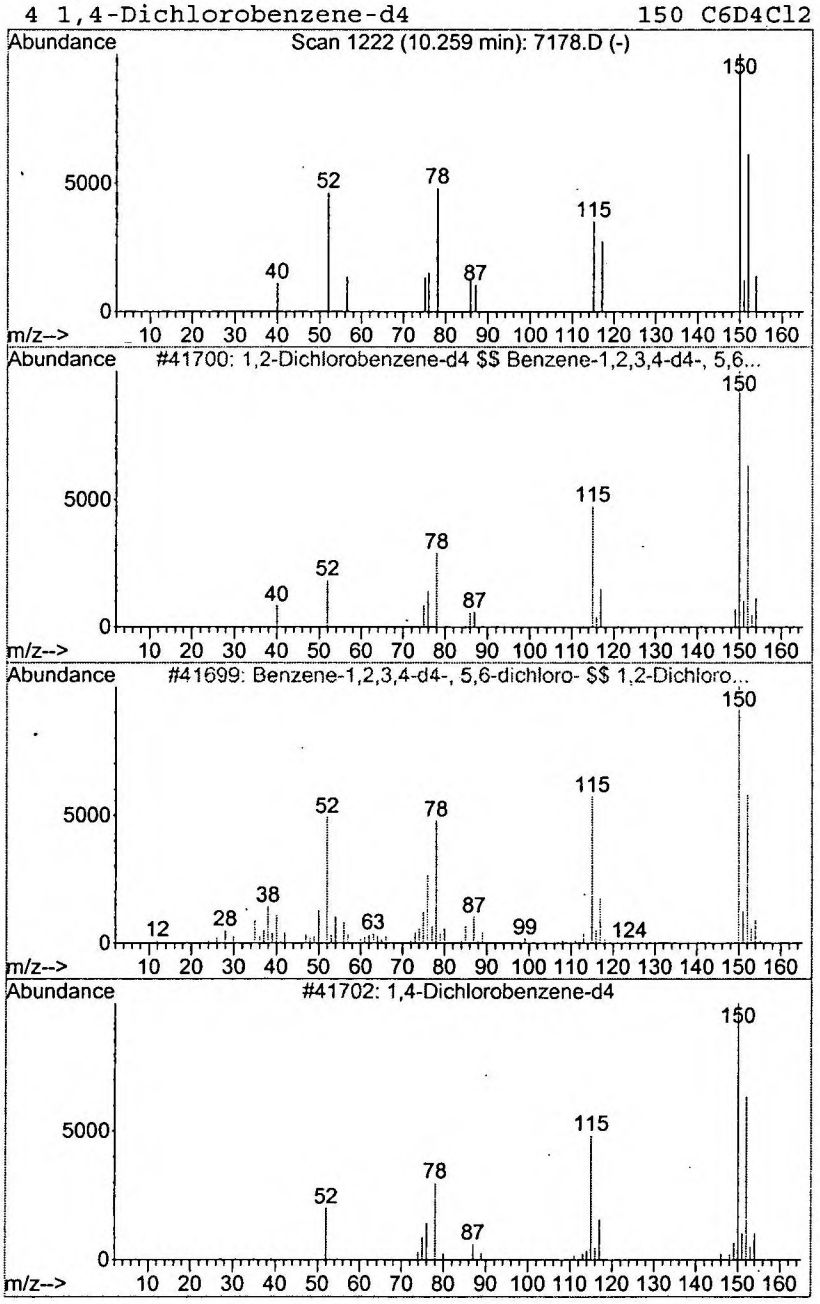
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

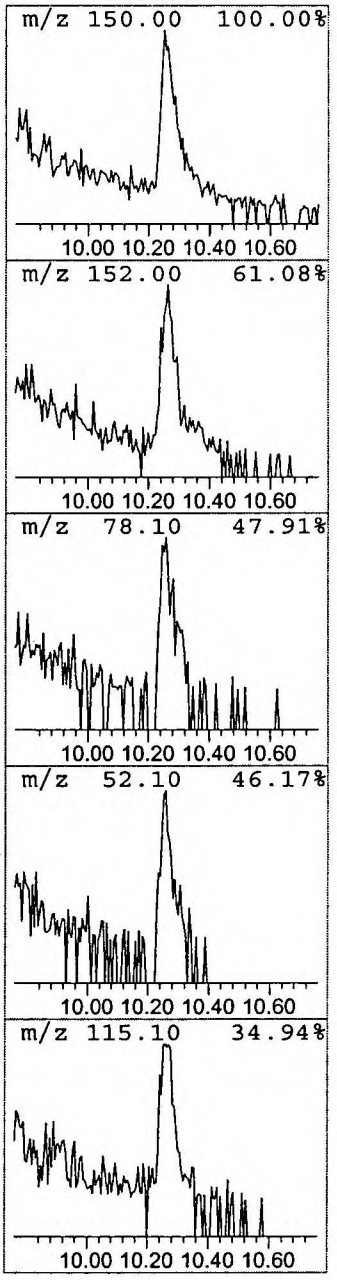
****

Peak Number 1 1,2-Dichlorobenzene-d4 $$ B... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	0.85 ug/L	63358	1,4-Dichlorobenzene-d4	9.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Dichlorobenzene-d4 $$ Benzen...	150	C6D4Cl2	002199-69-1	91
2	Benzene-1,2,3,4-d4-, 5,6-dichlor...	150	C6D4Cl2	002199-69-1	86
3	1,4-Dichlorobenzene-d4	150	C6D4Cl2	003855-82-1	86
4	1,4-Dichlorobenzene-d4	150	C6D4Cl2	003855-82-1	58





Tentatively Identified Compound (LSC) summary

Operator ID: Bohlk Date Acquired: 18 Apr 2002 8:18 pm
Data File: C:\MSDCHEM\1\DATA\041802\7178.D
Name: FB, SAMPLE 7178 3/27/02
Misc:
Method: C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title: Quantitation For Method 508gcscan
Library Searched: C:\DATABASE\WILEY7N.L

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1,2-Dichlorobenze...	10.26	0.9	ug/L	63358	1	9.41	2976480	40.0