

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/25/2002 Time: 11:06:18

Sample: AE01222
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/25/02 11:05 Ended: 2/25/02 11:05 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		104		5.0

Comments for Sample AE01222 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-25-2002 Time: 11:06:29

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 14 of the 43 compounds in the LCS Standard were below their acceptance ranges.

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1502\1222.D

Vial: 28

Acq On : 16 Feb 2002 7:34 pm

Operator: 209 GALOS

Sample : gc/ms scan 1/17

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 20 14:06 2002

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Wed Feb 13 11:43:23 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.34	152	254085	20.00	ug/l	-0.02
10) Naphthalene-d8	15.97	136	943415	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.29	164	527854	20.00	ug/l	-0.03
29) Phenanthrene-d10	25.74	188	758915	20.00	ug/l	-0.03
38) Chrysene-d12	33.80	240	487886	20.00	ug/l	-0.05
44) Perylene-d12	38.09	264	325911	20.00	ug/l	-0.04

System Monitoring Compounds

37) 2,4-DDD	30.82	235	39630	5.20	ug/mL	0.32
Spiked Amount	5.000		Recovery	=	104.00%	

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-Dichlorobenzene	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-Chloropropan	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)methane	0.00	93	0	N.D.
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.
15) Naphthalene	16.03	128	868	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.
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	22.75	149	226	N.D.
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.
27) Fluorene	0.00	166	0	N.D.
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.

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

1222.D GCMS0208.M Wed Feb 20 14:07:49 2002

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Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1502\1222.D

Vial: 28

Acq On : 16 Feb 2002 7:34 pm

Operator: 209 GALOS

Sample : gc/ms scan 1/17

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 20 14:06 2002

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Wed Feb 13 11:43:23 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0208

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	25.79	178	189	N.D.		
34) Anthracene	25.79	178	190	N.D.		
35) Di-n-butylphthalate	27.70	149	5905	N.D.		
36) Fluoranthene	0.00	202	0	N.D.		
39) Pyrene	0.00	202	0	N.D.		
40) Butylbenzylphthalate	0.00	149	0	N.D.		
41) Benzo(a)anthracene	33.81	228	1274	N.D.		
42) Bis(2-ethylhexyl)phthalate	34.01	149	4444	N.D.		
43) Chrysene	33.81	228	1274	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	0.00	252	0	N.D.		
47) Benzo(k)fluoranthene	0.00	252	0	N.D.		
48) Benzo(a)pyrene	38.08	252	1121	N.D.		
49) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.		
50) Dibenz(a,h)anthracene	0.00	278	0	N.D.		
51) Benzo(g,h,i)perylene	0.00	276	0	N.D.		

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

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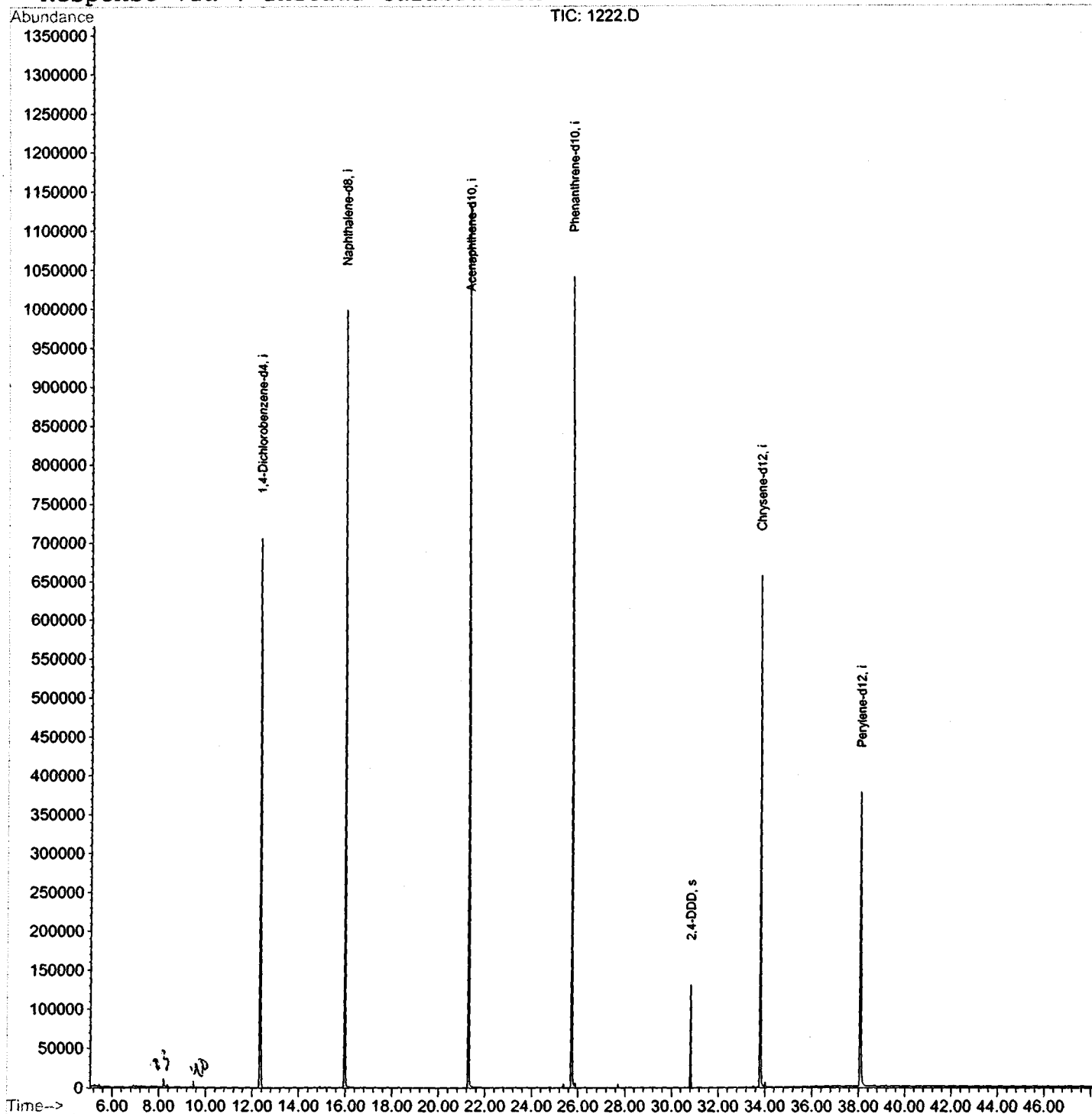
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Data File : C:\HPCHEM\1\DATA\FEB1502\1222.D
 Acq On : 16 Feb 2002 7:34 pm
 Sample : gc/ms scan 1/17
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Feb 20 14:06 2002

Vial: 28
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Feb 13 11:43:23 2002
 Response via : Initial Calibration



1222.D GCMS0208.M

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LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB1502\1222.D

Vial: 28

Acq On : 16 Feb 2002 7:34 pm

Operator: 209 GALOS

Sample : gc/ms scan 1/17

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

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	peak area	peak % max.	% of total
1	12.335	788	794	809	rVB	705599	1640740	74.30%	15.198%
2	15.971	1184	1190	1202	rBV	998287	1973588	89.38%	18.281%
3	21.286	1763	1769	1784	rBV	1135514	2208169	100.00%	20.454%
4	25.738	2247	2254	2265	rBV	1040818	2110158	95.56%	19.546%
5	30.815	2802	2807	2813	rVB	131710	245099	11.10%	2.270%
6	33.799	3126	3132	3145	rBV	657593	1488059	67.39%	13.784%
7	38.086	3591	3599	3626	rBV2	377635	1130043	51.18%	10.467%

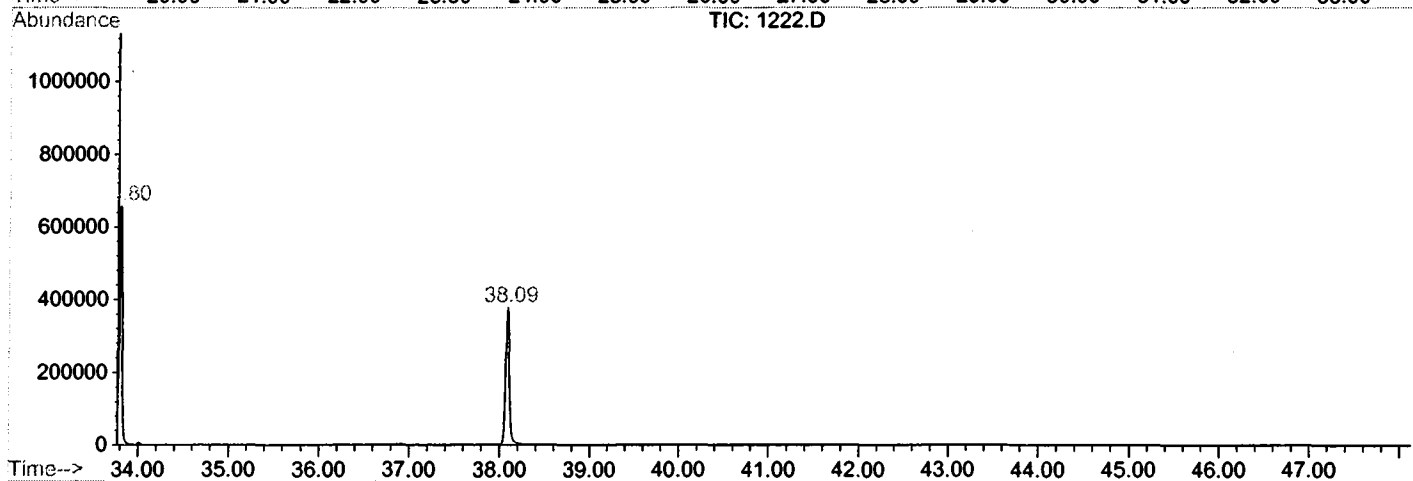
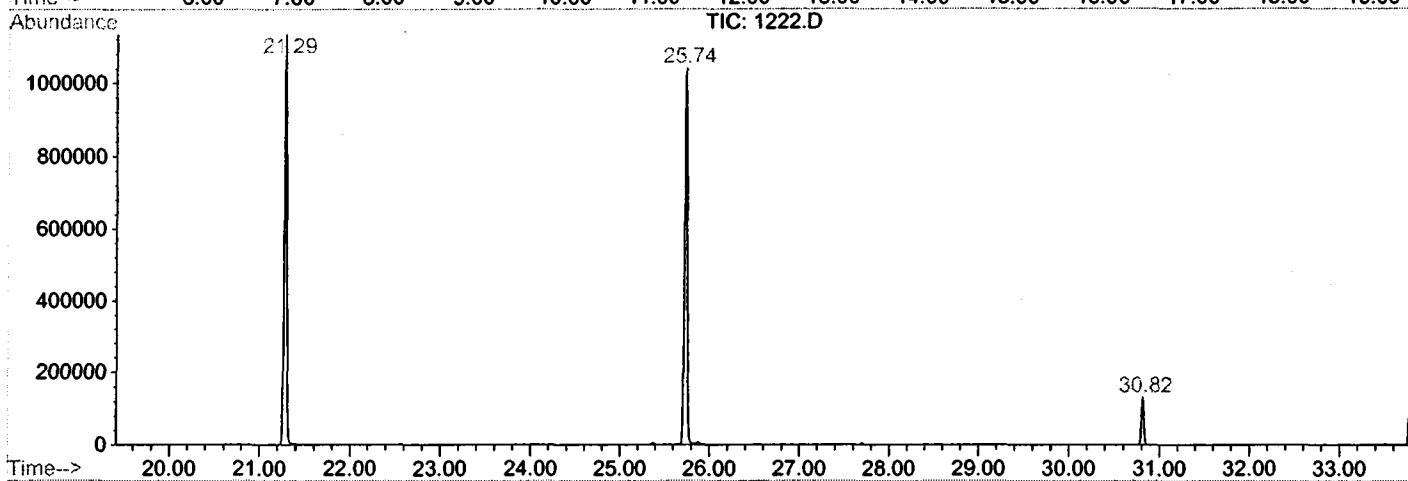
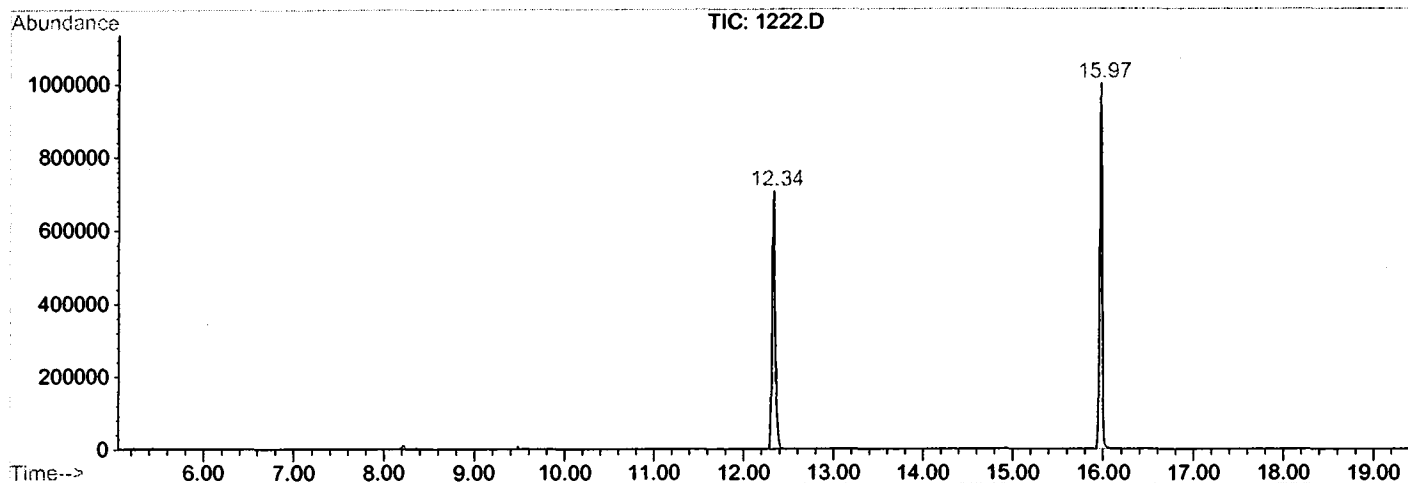
Sum of corrected areas: 10795856

1222.D GCMS0208.M

Wed Feb 20 14:30:03 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB1502\1222.D
 Operator : 209 GALOS
 Acquired : 16 Feb 2002 7:34 pm using AcqMethod GCMS0208
 Instrument : GC/MS 209
 Sample Name: gc/ms scan 1/17
 Misc Info :
 Vial Number: 28
 Quant File :GCMS0208.RES (RTE Integrator)



1222.D GCMS0208.M

Wed Feb 20 14:30:09 2002

Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 16 Feb 2002 7:34 pm
 Data File: C:\HPCHEM\1\DATA\FEB1502\1222.D
 Name: gc/ms scan 1/17
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
 Title: 209 625/TCLP calibration table
 Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc

1222.D GCMS0208.M			Wed Feb 20 14:30:13 2002					