

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/22/2002 Time: 11:41:52

Sample: AE01967
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/22/02 11:18 Ended: 3/22/02 11:18 Analyst: DLV
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2	Surr. 2,4-DDD	USPEC	156		5.0

Comments for Sample AE01967 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L.

Date: 03-22-2002 Time: 11:41:54

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

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR1902\1967.D

Vial: 12

Acq On : 20 Mar 2002 3:47 pm

Operator:

Sample : gc/ms scan 3/5 fb

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Mar 21 9:06 2002

Quant Results File: GCMS0308.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0308.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Mar 08 08:54:27 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0308

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.27	152	560449	20.00	ug/l	-0.09
10) Naphthalene-d8	15.90	136	2186467m	20.00	ug/l	-0.10
17) Acenaphthene-d10	21.22	164	1181509	20.00	ug/l	-0.10
29) Phenanthrene-d10	25.66	188	1754673	20.00	ug/l	-0.10
38) Chrysene-d12	33.72	240	1116907	20.00	ug/l	-0.12
44) Perylene-d12	37.97	264	699756	20.00	ug/l	-0.15

System Monitoring Compounds

37) 2,4-DDD	30.74	235	99726	7.82	ug/mL	0.24
Spiked Amount	5.000		Recovery	= 156.40%		

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	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.	
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-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

1967.D GCMS0308.M Thu Mar 21 09:07:05 2002

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR1902\1967.D

Vial: 12

Acq On : 20 Mar 2002 3:47 pm

Operator:

Sample : gc/ms scan 3/5 fb

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Mar 21 9:06 2002

Quant Results File: GCMS0308.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0308.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Mar 08 08:54:27 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0308

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	25.67	178	454	N.D.		
34) Anthracene	25.67	178	454	N.D.		
35) Di-n-butylphthalate	27.63	149	1732	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	33.72	228	2301	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.93	149	2471	N.D.		
43) Chrysene	33.72	228	2301	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	36.78	252	481	N.D.		
47) Benzo(k)fluoranthene	36.84	252	686	N.D.		
48) Benzo(a)pyrene	37.77	252	473	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

1967.D GCMS0308.M

Thu Mar 21 09:07:09 2002

Page 2

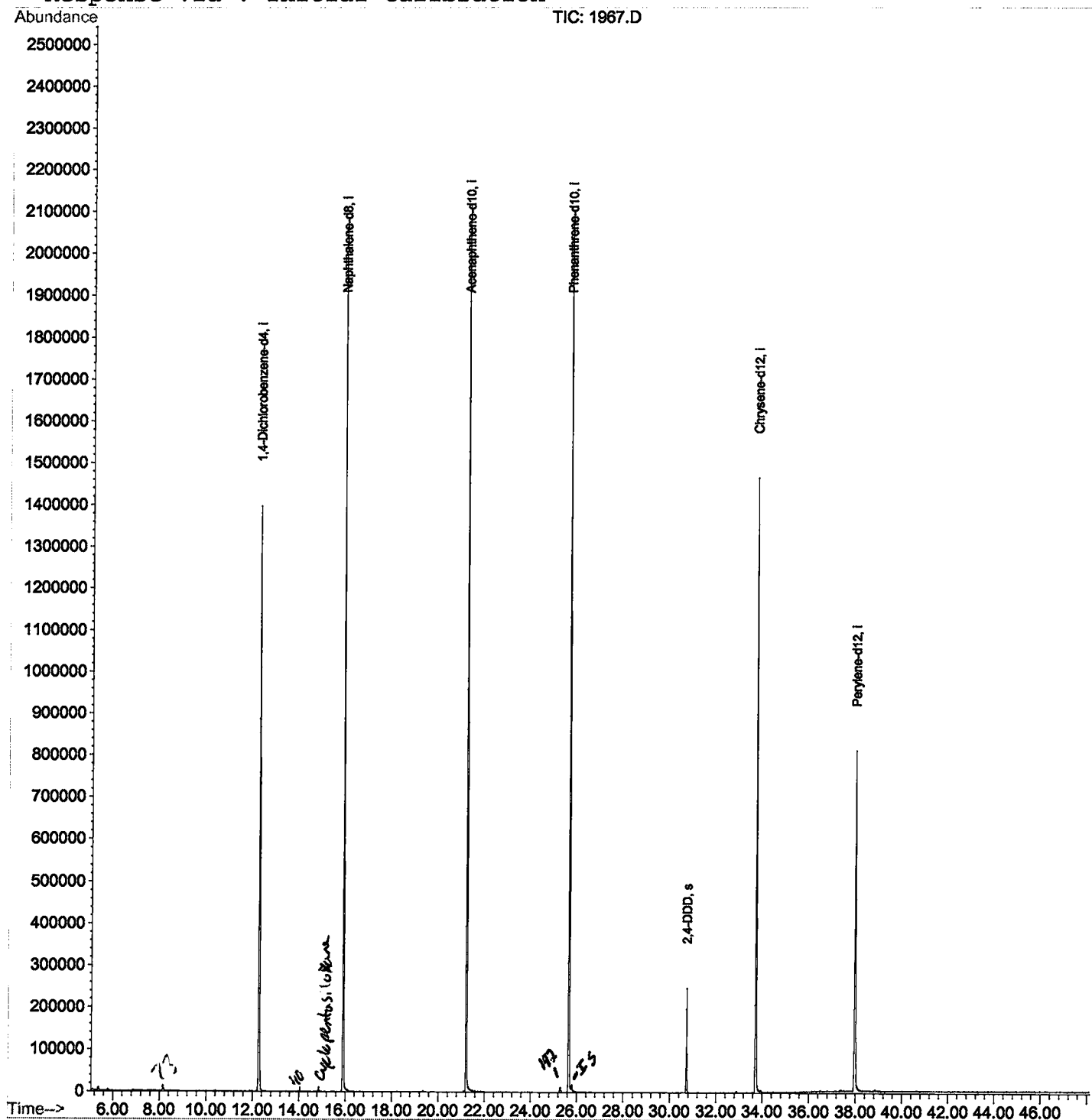
Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAR1902\1967.D
Acq On : 20 Mar 2002 3:47 pm
Sample : gc/ms scan 3/5 fb
Misc :
MS Integration Params: GOOFY.P
Quant Time: Mar 21 9:06 2002

Vial: 12
Operator:
Inst : GC/MS 209
Multiplr: 1.00

Quant Results File: GCMS0308.RES

Method : C:\HPCHEM\1\METHODS\GCMS0308.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Mar 08 08:54:27 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAR1902\1967.D
 Acq On : 20 Mar 2002 3:47 pm
 Sample : gc/ms scan 3/5 fb
 Misc :
 MS Integration Params: LSCINT.P

Vial: 12
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0308.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

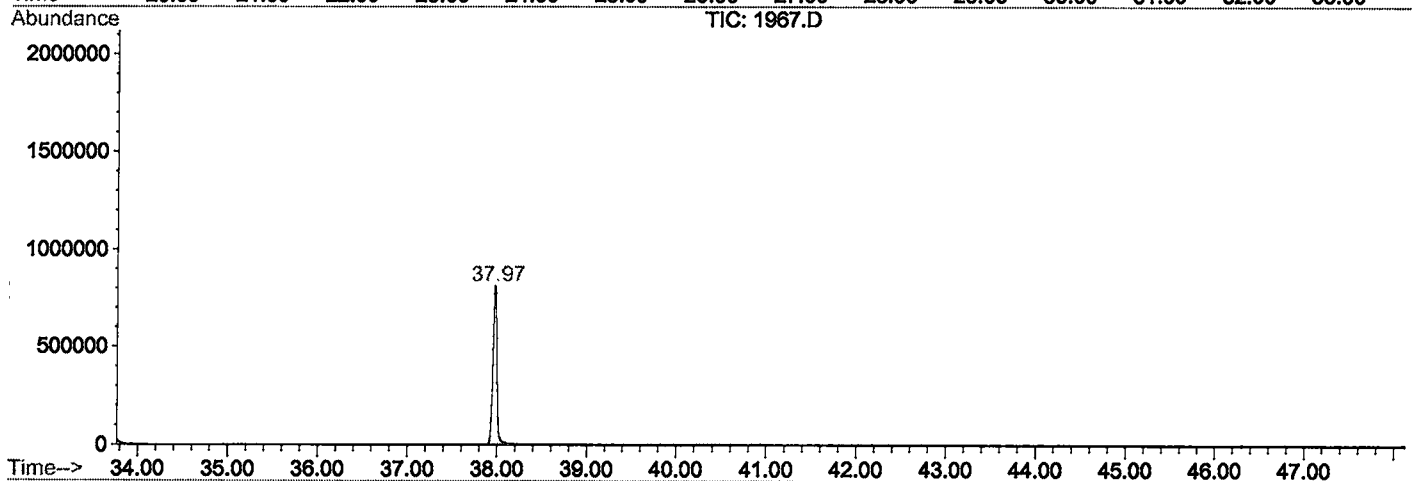
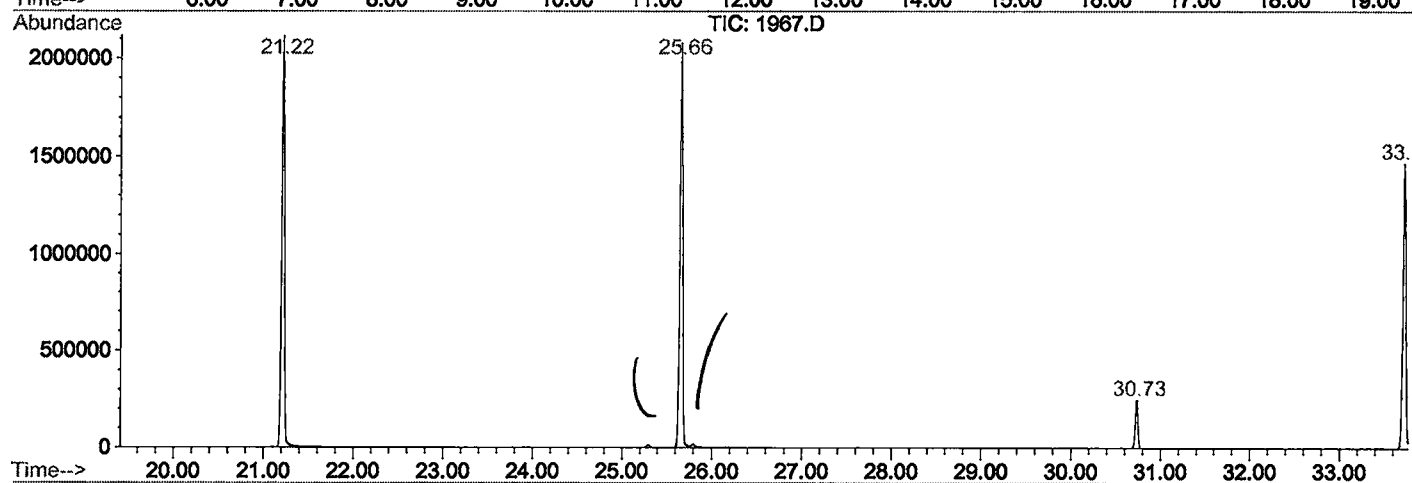
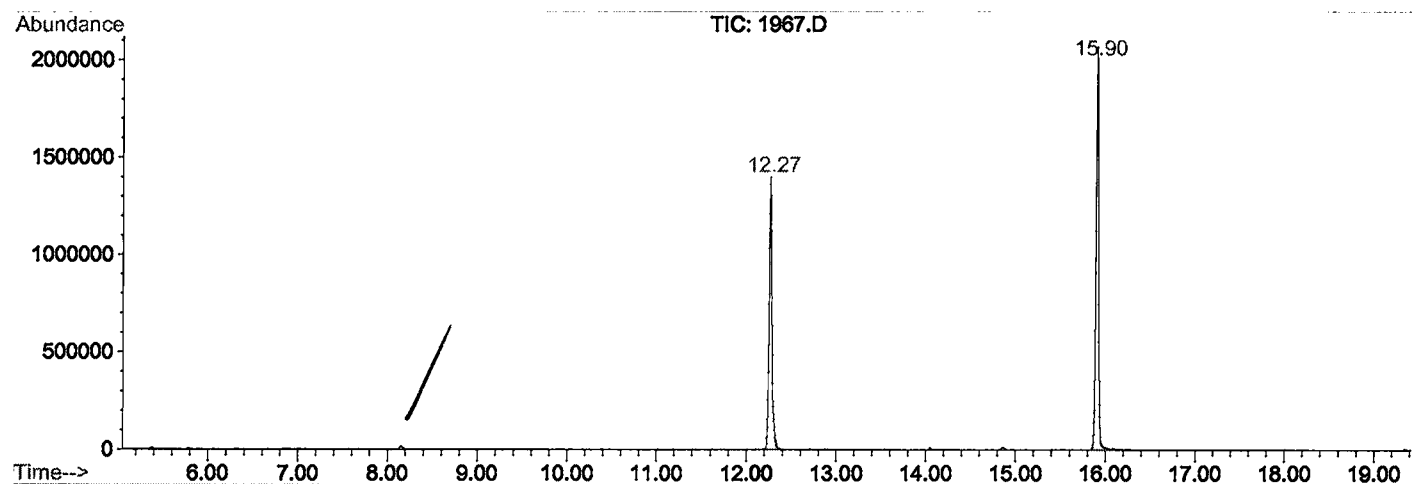
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	12.267	781	787	801	rVB	1395225	3101503	67.88%	13.731%
2	15.902	1176	1183	1201	rBV	2062645	4204379	92.02%	18.613%
3	21.218	1754	1762	1778	rBV	2117839	4569123	100.00%	20.228%
4	25.661	2239	2246	2256	rBV2	2078991	4431386	96.99%	19.618%
5	30.729	2793	2798	2805	rBV	246268	501466	10.98%	2.220%
6	33.721	3117	3124	3139	rBV2	1464525	3315822	72.57%	14.680%
7	37.972	3578	3587	3602	rBV2	807612	2464419	53.94%	10.910%

Sum of corrected areas: 22588098

1967.D GCMS0308.M Thu Mar 21 09:17:35 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAR1902\1967.D
 Operator :
 Acquired : 20 Mar 2002 3:47 pm using AcqMethod GCMS0308
 Instrument : GC/MS 209
 Sample Name: gc/ms scan 3/5 fb
 Misc Info :
 Vial Number: 12
 Quant File :GCMS0308.RES (RTE Integrator)



1967.D GCMS0308.M

Thu Mar 21 09:17:40 2002

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 20 Mar 2002 3:47 pm
 Data File: C:\HPCHEM\1\DATA\MAR1902\1967.D
 Name: gc/ms scan 3/5 fb
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0308.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

1967.D GCMS0308.M			Thu Mar 21 09:17:44 2002					