

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/29/2002 Time: 08:50:18

Sample: AE03507
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/29/02 08:49 Ended: 3/29/02 08:49 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2	Surr_2,4-DDD		103		5.0

Comments for Sample AE03507 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-29-2002 Time: 08:50:21

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 9 of the 43 compounds in the LCS Standard were high. New limits will be calculated.

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR2102\3507.D

Vial: 22

Acq On : 22 Mar 2002 7:59 am

Operator:

Sample : re-extract

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Mar 25 10:38 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	12.26	152	801206	20.00	ug/l	-0.09
10) Naphthalene-d8	15.91	136	3176925	20.00	ug/l	-0.09
17) Acenaphthene-d10	21.22	164	1721272	20.00	ug/l	-0.10
29) Phenanthrene-d10	25.67	188	2601822	20.00	ug/l	-0.10
38) Chrysene-d12	33.73	240	1582972	20.00	ug/l	-0.12
44) Perylene-d12	37.99	264	903912	20.00	ug/l	-0.13

System Monitoring Compounds

37) 2,4-DDD	30.74	235	97022	5.13	ug/mL	0.24
Spiked Amount	5.000		Recovery	=	102.60%	

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	15.96	128	216	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	21.64	165	105	N.D.
25) Diethylphthalate	22.69	149	679	N.D.
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.
27) Fluorene	0.00	166	0	N.D.
28) Azobenzene/ 1,2-Diphenylhyd	0.00	77	0	N.D.

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

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Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR2102\3507.D

Vial: 22

Acq On : 22 Mar 2002 7:59 am

Operator:

Sample : re-extract

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Mar 25 10:38 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	828	N.D.		
34) Anthracene	25.67	178	828	N.D.		
35) Di-n-butylphthalate	27.62	149	2029	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	33.73	228	3771	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.93	149	2631	N.D.		
43) Chrysene	33.73	228	3771	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	36.78	252	185	N.D.		
47) Benzo(k)fluoranthene	36.84	252	172	N.D.		
48) Benzo(a)pyrene	37.99	252	3402	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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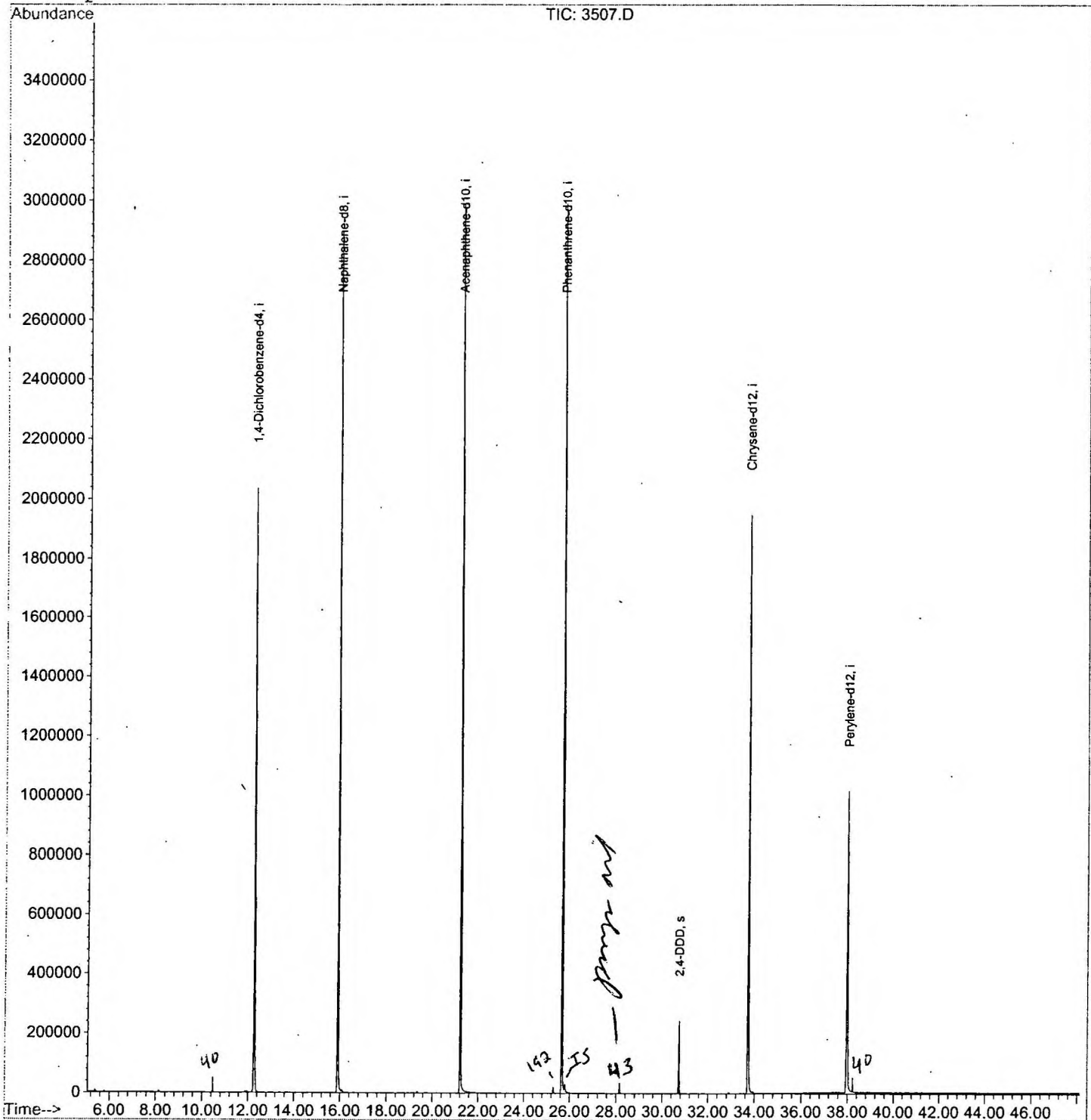
Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAR2102\3507.D
Acq On : 22 Mar 2002 7:59 am
Sample : re-extract
Misc :
MS Integration Params: GOOFY.P
Quant Time: Mar 25 10:38 2002

Vial: 22
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\MAR2102\3507.D

Vial: 22

Acq On : 22 Mar 2002 7:59 am

Operator:

Sample : re-extract

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

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	12.265	780	786	805	rVB	2033393	4415321	66.66%	13.807%
2	15.909	1176	1183	1199	rBV	2894730	6014432	90.80%	18.807%
3	21.215	1754	1761	1780	rBV	2984989	6623489	100.00%	20.711%
4	25.668	2238	2246	2256	rBV2	2990125	6619979	99.95%	20.700%
5	30.735	2792	2798	2804	rBV	242783	493096	7.44%	1.542%
6	33.728	3116	3124	3141	rBV2	1947683	4643835	70.11%	14.521%
7	37.988	3576	3588	3610	rBV2	1012868	3169734	47.86%	9.912%

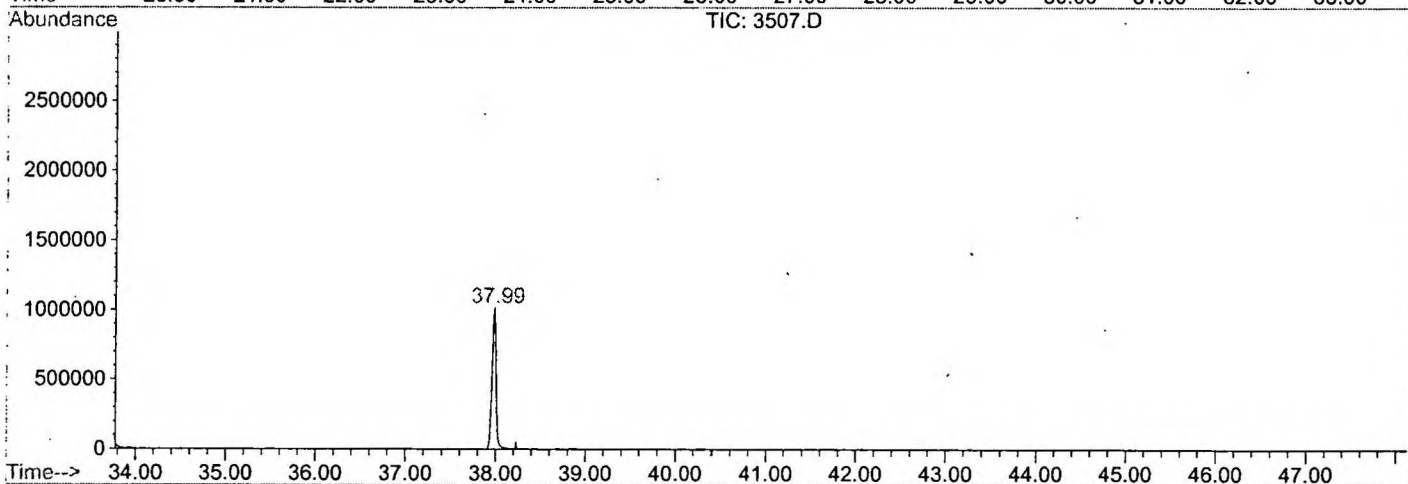
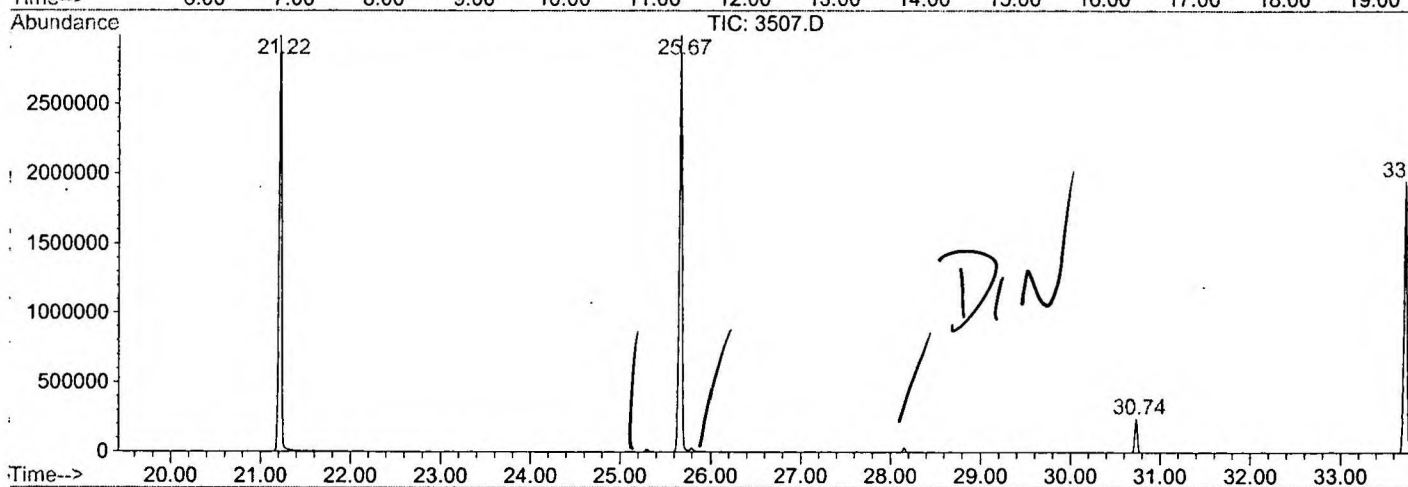
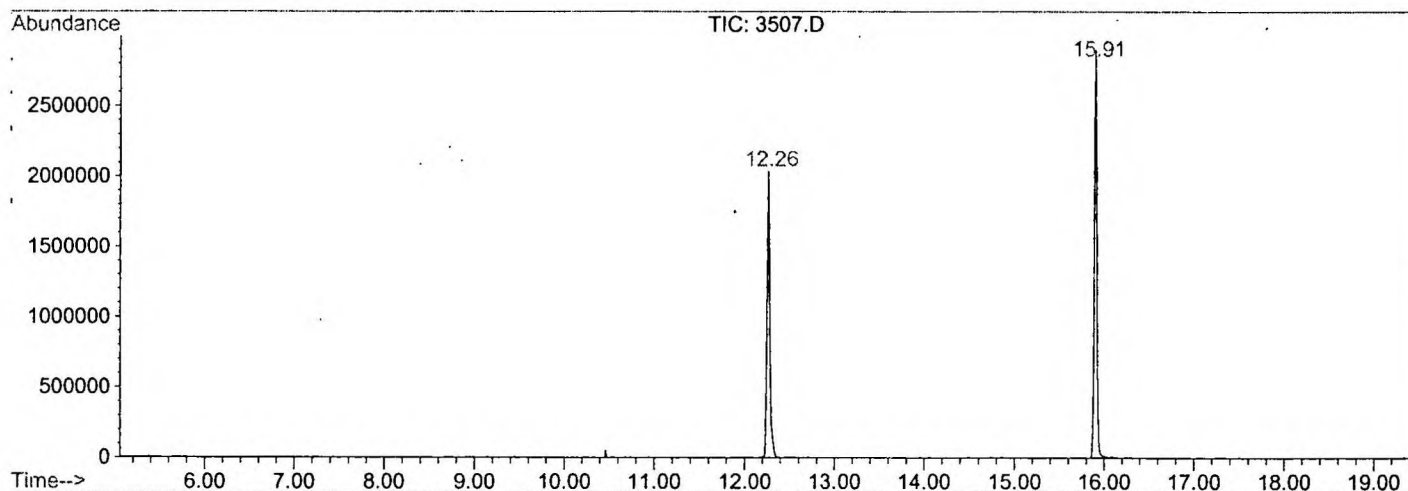
Sum of corrected areas: 31979886

3507.D GCMS0308.M

Mon Mar 25 14:23:39 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAR2102\3507.D
 Operator :
 Acquired : 22 Mar 2002 7:59 am using AcqMethod GCMS0308
 Instrument : GC/MS 209
 Sample Name: re-extract
 Misc Info :
 Vial Number: 22
 Quant File :GCMS0308.RES (RTE Integrator)



3507.D GCMS0308.M

Mon Mar 25 14:23:45 2002

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 22 Mar 2002 7:59 am
 Data File: C:\HPCHEM\1\DATA\MAR2102\3507.D
 Name: re-extract
 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

3507.D GCMS0308.M	Mon Mar 25	14:23:49	2002					