

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 04/04/2002 Time: 10:42:29

Sample: AE04735
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 4/4/02 10:41 Ended: 4/4/02 10:41 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2	Surr_2,4-DDD		115		5.0

Comments for Sample AE04735 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-04-2002 Time: 10:42:41

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 12 of the 43 compounds in the LCS Standard were high.

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR2802\4735.D

Vial: 8

Acq On : 28 Mar 2002 3:28 pm

Operator:

Sample : gc/ms scan 3/4

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Apr 1 9:53 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 29 10:31:28 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0308

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.25	152	400858	20.00	ug/l	-0.10
10) Naphthalene-d8	15.89	136	1460798	20.00	ug/l	-0.11
17) Acenaphthene-d10	21.19	164	830367	20.00	ug/l	-0.12
29) Phenanthrene-d10	25.63	188	1176823	20.00	ug/l	-0.13
38) Chrysene-d12	33.69	240	576127	20.00	ug/l	-0.15
44) Perylene-d12	37.94	264	322597	20.00	ug/l	-0.19

System Monitoring Compounds

37) 2,4-DDD	30.71	235	67700	5.73	ug/mL	0.21
Spiked Amount	5.000		Recovery	=	114.60%	

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	14.60	82	341	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	15.94	128	770	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	20.54	163	1615	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	22.67	149	1892	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

4735.D GCMS0308.M Mon Apr 01 09:54:45 2002

Page 1

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR2802\4735.D

Vial: 8

Acq On : 28 Mar 2002 3:28 pm

Operator:

Sample : gc/ms scan 3/4

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Apr 1 9:53 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 29 10:31:28 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	25.70	178	1005	N.D.		
34) Anthracene	25.70	178	1005	N.D.		
35) Di-n-butylphthalate	27.61	149	2517	N.D.		
36) Fluoranthene	29.33	202	954	N.D.		
39) Pyrene	30.01	202	750	N.D.		
40) Butylbenzylphthalate	0.00	149	0	N.D.		
41) Benzo(a)anthracene	33.69	228	1779	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.92	149	1340	N.D.		
43) Chrysene	33.77	228	883	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	36.83	252	1386	N.D.		
47) Benzo(k)fluoranthene	36.83	252	738	N.D.		
48) Benzo(a)pyrene	37.75	252	100	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

4735.D GCMS0308.M Mon Apr 01 09:54:48 2002

Page 2

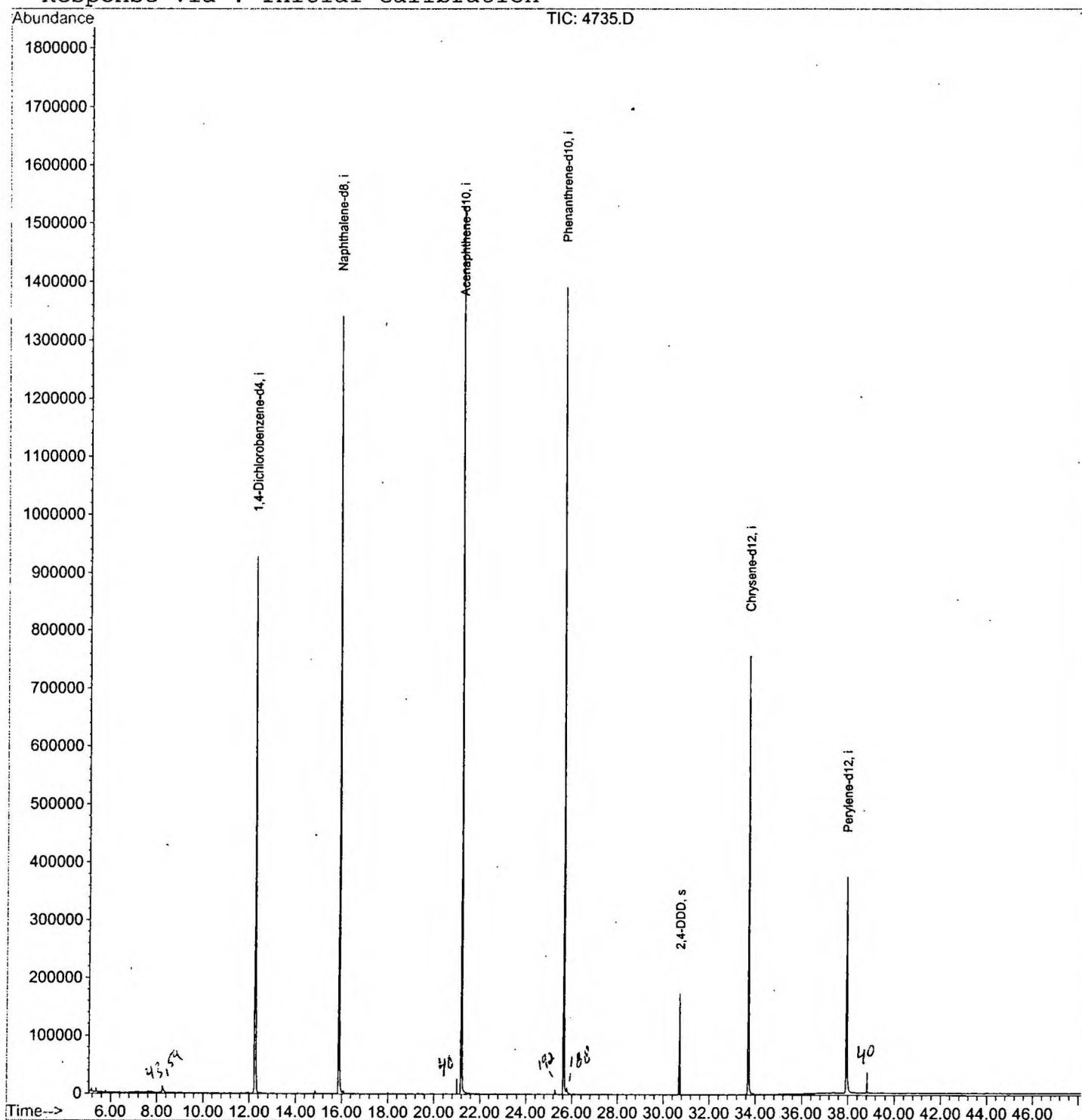
Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAR2802\4735.D
Acq On : 28 Mar 2002 3:28 pm
Sample : gc/ms scan 3/4
Misc :
MS Integration Params: GOOFY.P
Quant Time: Apr 1 9:53 2002

Vial: 8
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 29 10:31:28 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAR2802\4735.D
 Acq On : 28 Mar 2002 3:28 pm
 Sample : gc/ms scan 3/4
 Misc :
 MS Integration Params: LSCINT.P

Vial: 8
 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
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Signal : TIC

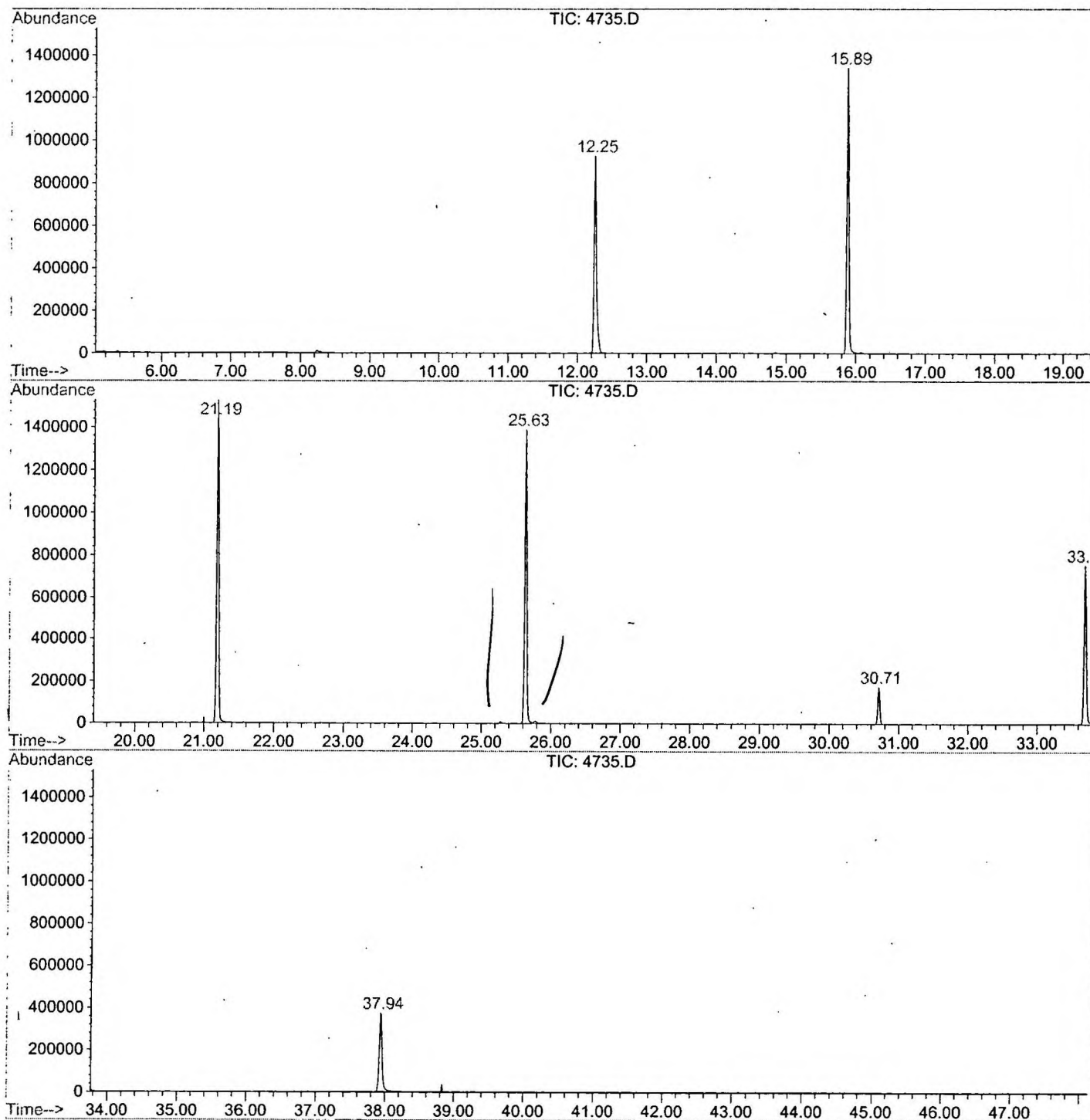
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	12.250	780	785	799	rVB	925514	2166405	68.91%	15.379%
2	15.885	1175	1181	1199	rBV	1341324	2764148	87.92%	19.623%
3	21.191	1753	1759	1769	rBV	1528087	3143851	100.00%	22.318%
4	25.635	2237	2243	2255	rBV2	1390555	2943894	93.64%	20.899%
5	30.711	2791	2796	2804	rBV	174224	332994	10.59%	2.364%
6	33.695	3115	3121	3137	rBV2	755944	1655236	52.65%	11.751%
7	37.936	3576	3583	3598	rBV2	372687	1079809	34.35%	7.666%

Sum of corrected areas: 14086337

4735.D GCMS0308.M Mon Apr 01 10:06:40 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAR2802\4735.D
 Operator :
 Acquired : 28 Mar 2002 3:28 pm using AcqMethod GCMS0308
 Instrument : GC/MS 209
 Sample Name: gc/ms scan 3/4
 Misc Info :
 Vial Number: 8
 Quant File :GCMS0308.RES (RTE Integrator)



4735.D GCMS0308.M

Mon Apr 01 10:06:45 2002

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 28 Mar 2002 3:28 pm
Data File: C:\HPCHEM\1\DATA\MAR2802\4735.D
Name: gc/ms scan 3/4
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

4735.D GCMS0308.M	Mon Apr 01 10:06:49 2002							