

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/29/2002 Time: 15:13:10

Sample: AE07107
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/29/02 15:12 Ended: 4/29/02 15:12 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		73		10.0

Comments for Sample AE07107 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-29-2002 Time: 15:13:52

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

The recoveries for 5 of the 43 compounds in the LCS Standard were low.

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 25 8:30 2002

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Apr 17 11:48:30 2002
 Response via : Initial Calibration
 DataAcq Meth : GCMS0412

Handwritten signature

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.21	152	560709	20.00	ug/l	-0.03
10) Naphthalene-d8	15.85	136	2055286	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.14	164	1151769	20.00	ug/l	-0.03
30) Phenanthrene-d10	25.59	188	1683081	20.00	ug/l	-0.03
39) Chrysene-d12	33.65	240	1072449	20.00	ug/l	-0.03
45) Perylene-d12	37.87	264	675318	20.00	ug/l	-0.05

System Monitoring Compounds

28) TCMX	23.28	209	42140	7.30	ug/L	-0.05
Spiked Amount	10.000		Recovery	=	73.00%	
38) 2,4-DDD	0.00	235	0	0.00	ug/mL	
Spiked Amount	5.000		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-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	22.63	149	684	N.D.	
26) 4-Chlorophenyl-phenylether	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	23.28	168	169	N.D.	
32) 4-Bromophenyl-phenylether	0.00	248	0	N.D.	

(#) = qualifier out of range (m) = manual integration
 7107.D GCMS0412.M Thu Apr 25 11:08:58 2002

Quantitation Report

(QT Reviewed)

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 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 25 8:30 2002

Vial: 13
 Operator:
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Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Apr 17 11:48:30 2002
 Response via : Initial Calibration
 DataAcq Meth : GCMS0412

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.59	178	390	N.D.		
35) Anthracene	25.59	178	390	N.D.		
36) Di-n-butylphthalate	27.56	149	3597	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)anthracene	33.64	228	2579	N.D.		
43) Bis(2-ethylhexyl)phthalate	33.85	149	805	N.D.		
44) Chrysene	33.64	228	2579	N.D.		
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	37.87	252	2507	N.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
 7107.D GCMS0412.M Thu Apr 25 11:08:59 2002

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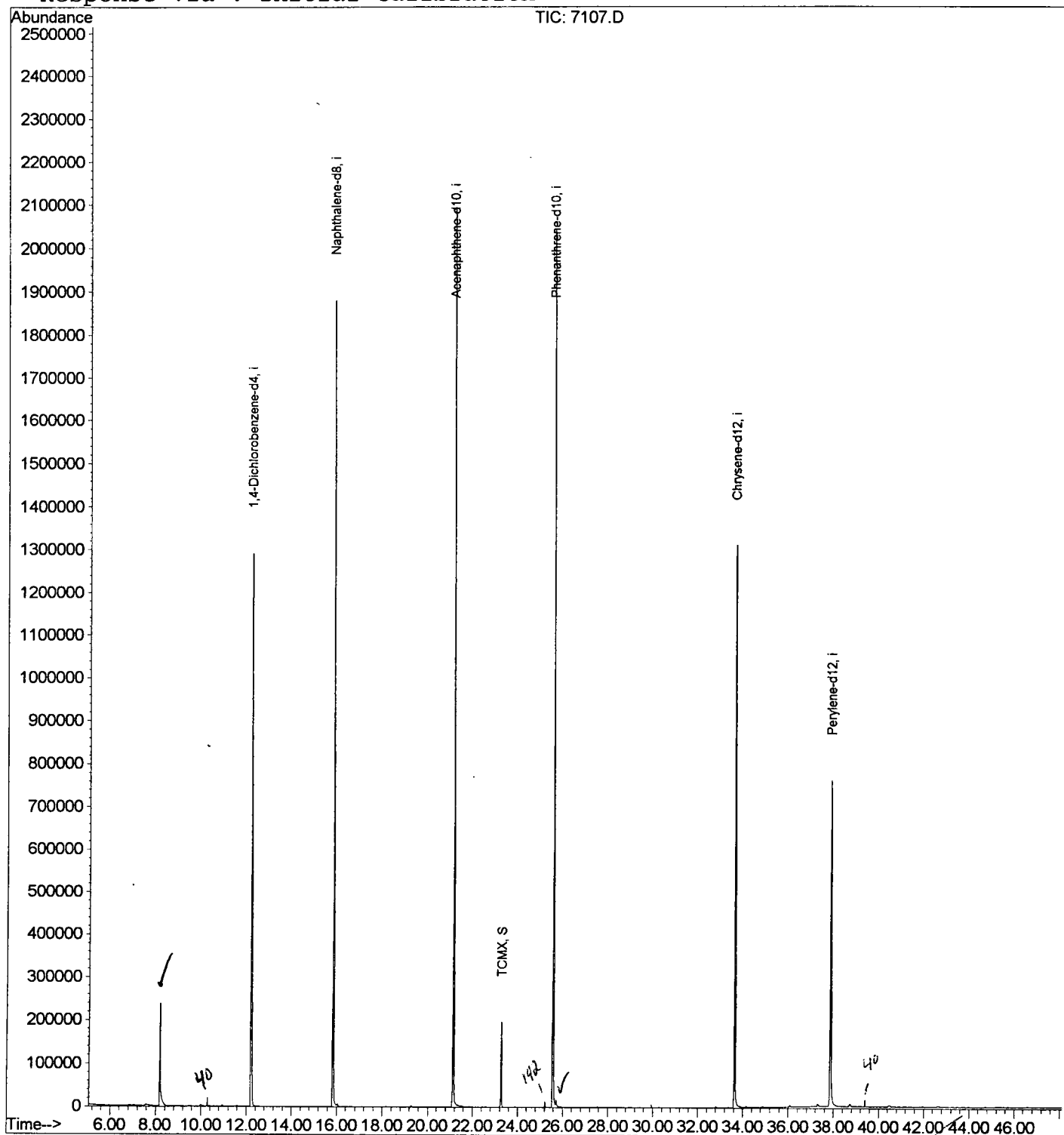
Quantitation Report

Data File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 25 8:30 2002

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0412.RES

Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Apr 17 11:48:30 2002
 Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: LSCINT.P

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0412.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

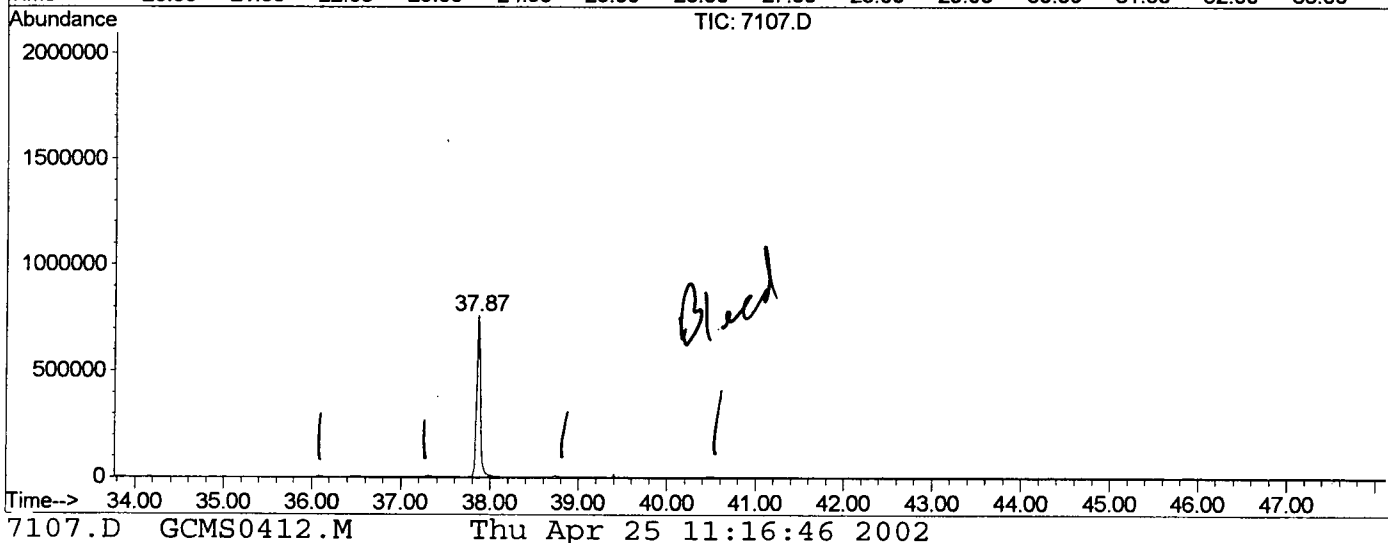
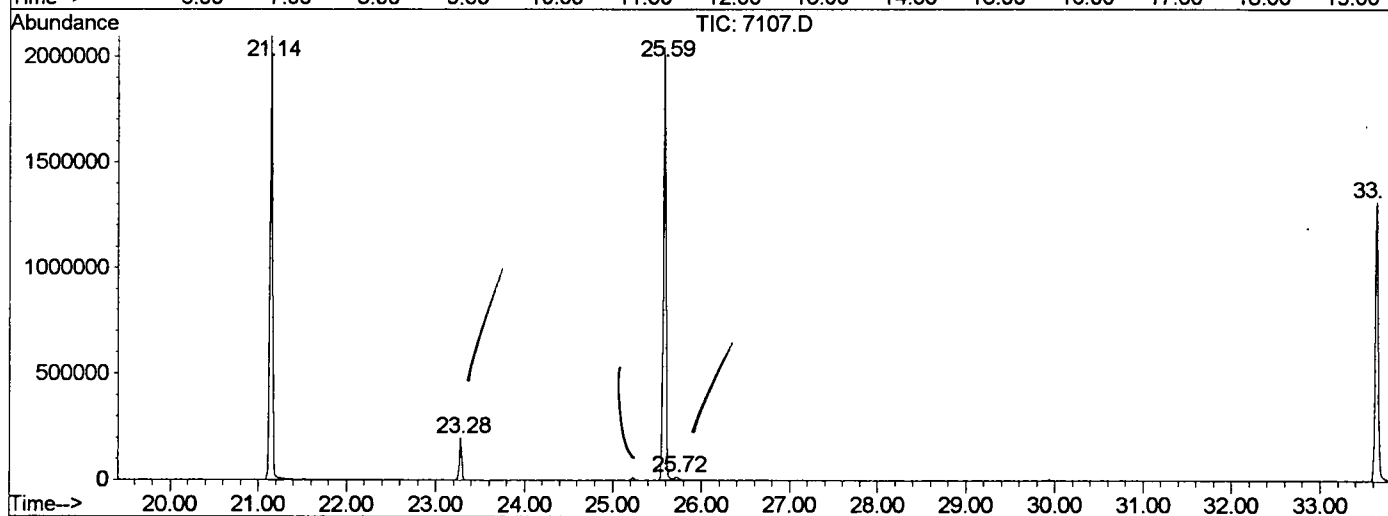
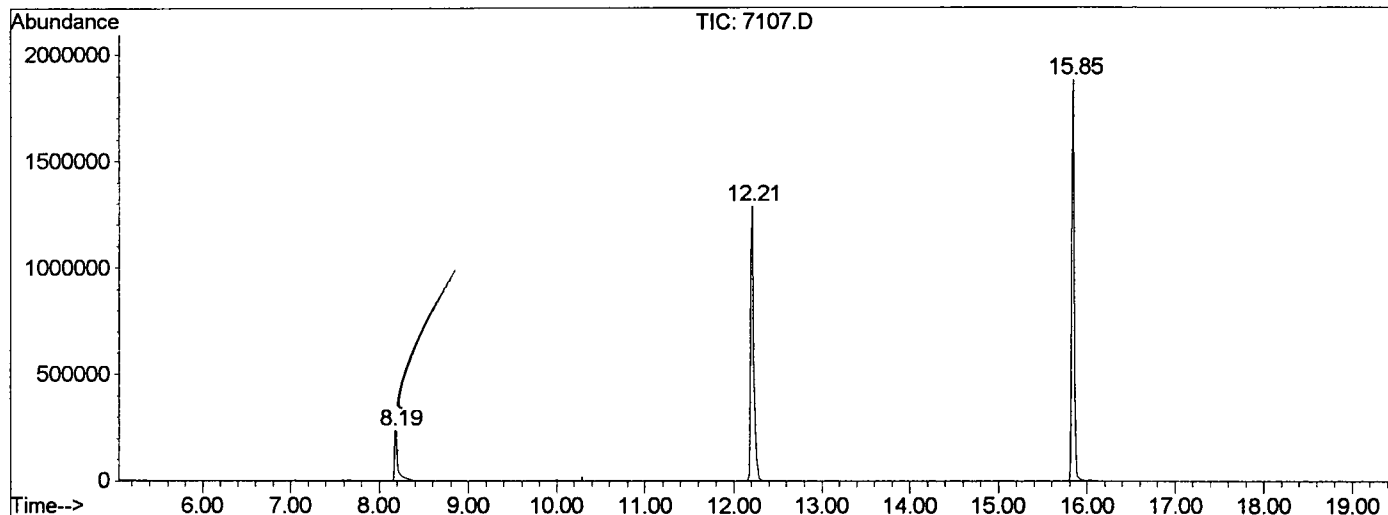
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.189	338	342	369	rBV	237273	607683	13.85%	2.774%
2	12.210	774	780	795	rBV	1291371	3024001	68.93%	13.803%
3	15.845	1170	1176	1193	rBV	1880451	3870514	88.23%	17.666%
4	21.142	1747	1753	1774	rBV	2095750	4386787	100.00%	20.023%
5	23.281	1978	1986	1997	rVB	197835	382555	8.72%	1.746%
6	25.585	2230	2237	2247	rBV2	2042910	4216386	96.12%	19.245%
7	25.723	2248	2252	2263	rVB	15945	44165	1.01%	0.202%
8	33.646	3108	3115	3129	rBV2	1311101	3090802	70.46%	14.107%
9	37.869	3566	3575	3594	rBV2	759088	2286180	52.12%	10.435%

Sum of corrected areas: 21909073

7107.D GCMS0412.M Thu Apr 25 11:16:45 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Operator :
 Acquired : 25 Apr 2002 3:27 am using AcqMethod GCMS0412
 Instrument : 209
 Sample Name: gc/ms scan 4/23
 Misc Info :
 Vial Number: 13
 Quant File :GCMS0412.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: LSCINT.P

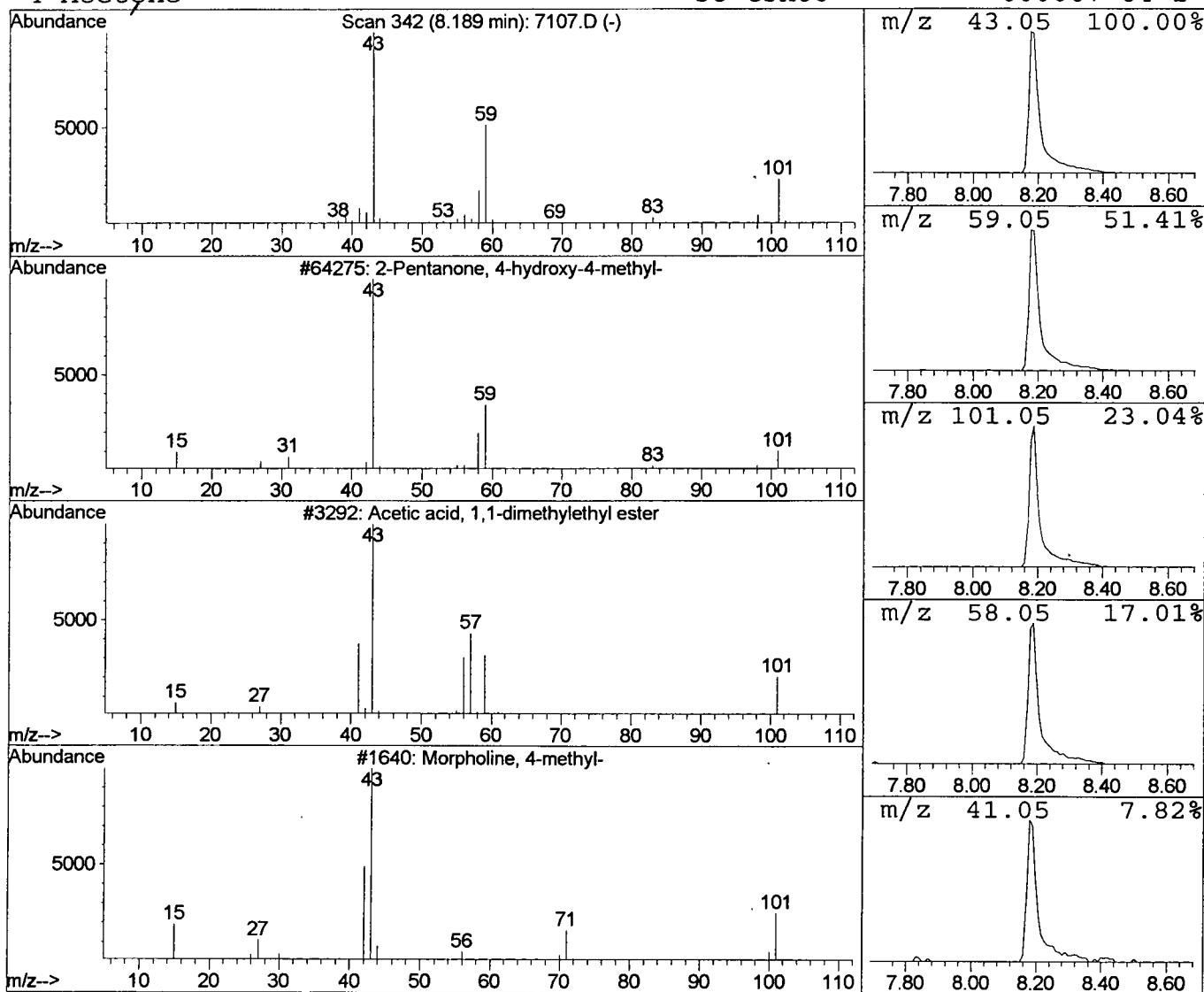
Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 1 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.19	4.02 ug/l	607683	1,4-Dichlorobenzene-d4	12.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	33
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Acetone	58	C3H6O	000067-64-1	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR2402\7107.D
 Acq On : 25 Apr 2002 3:27 am
 Sample : gc/ms scan 4/23
 Misc :
 MS Integration Params: LSCINT.P

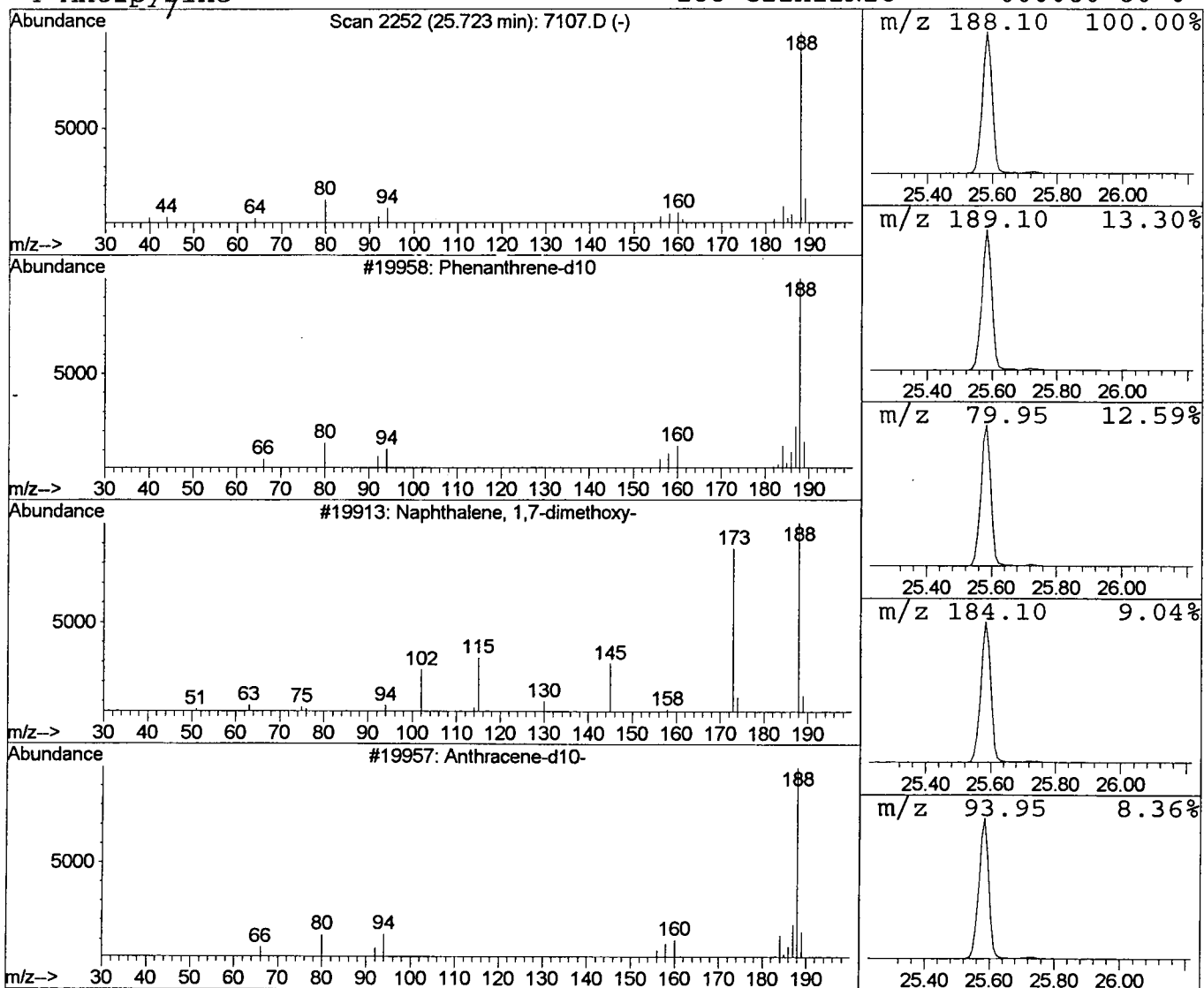
Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 2 Phenanthrene-d10 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.72	0.21 ug/l	44165	Phenanthrene-d10	25.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene-d10	188	C14D10	001517-22-2	72
2		Naphthalene, 1,7-dimethoxy-	188	C12H12O2	005309-18-2	42
3		Anthracene-d10-	188	C14D10	001719-06-8	42
4		Antipyrine	188	C11H12N2O	000060-80-0	40



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 25 Apr 2002 3:27 am
 Data File: C:\HPCHEM\1\DATA\APR2402\7107.D
 Name: gc/ms scan 4/23
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0412.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
-----	-----	-----	-----	-----	-----	-----	-----	-----
2-Pentanone, 4-hydro	8.19	4.0	ug/l	607683	ISTD01	12.21	3024000	20.0
Phenanthrene-d10	25.72	0.2	ug/l	44165	ISTD04	25.59	4216390	20.0

7107.D GCMS0412.M Thu Apr 25 11:16:48 2002