

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/27/2002 Time: 15:54:05

Sample: AD31499
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/27/02 15:53 Ended: 2/27/02 15:53 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		105		5.0

Comments for Sample AD31499 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-27-2002 Time: 15:54:08

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\FEB2102\31499.D
 Acq On : 23 Feb 2002 12:39 am
 Sample : gc/ms scan 2/21
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Feb 27 11:34 2002

Vial: 35
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

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 27 11:30:15 2002
 Response via : Initial Calibration
 DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	12.33	152	279689	20.00	ug/l	-0.02
10) Naphthalene-d8	15.97	136	1070686	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.29	164	558063	20.00	ug/l	-0.03
29) Phenanthrene-d10	25.74	188	796290	20.00	ug/l	-0.03
38) Chrysene-d12	33.81	240	542913	20.00	ug/l	-0.04
44) Perylene-d12	38.09	264	371329	20.00	ug/l	-0.04

System Monitoring Compounds

37) 2,4-DDD	30.81	235	43692	5.25	ug/mL	0.32
Spiked Amount	5.000		Recovery	=	105.00%	

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	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

31499.D GCMS0208.M Wed Feb 27 11:41:09 2002

Page 1

Quantitation Report

(QT Reviewed)

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 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	0.00	178	0	N.D.		
34) Anthracene	0.00	178	0	N.D.		
35) Di-n-butylphthalate	27.70	149	2537	N.D.		
36) Fluoranthene	29.44	202	197	N.D.		
39) Pyrene	30.11	202	284	N.D.		
40) Butylbenzylphthalate	0.00	149	0	N.D.		
41) Benzo(a)anthracene	33.80	228	1358	N.D.		
42) Bis(2-ethylhexyl)phthalate	34.01	149	11454	0.32	ug/l	95
43) Chrysene	33.80	228	1358	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	38.09	252	1448	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

31499.D GCMS0208.M

Wed Feb 27 11:41:13 2002

Page 2

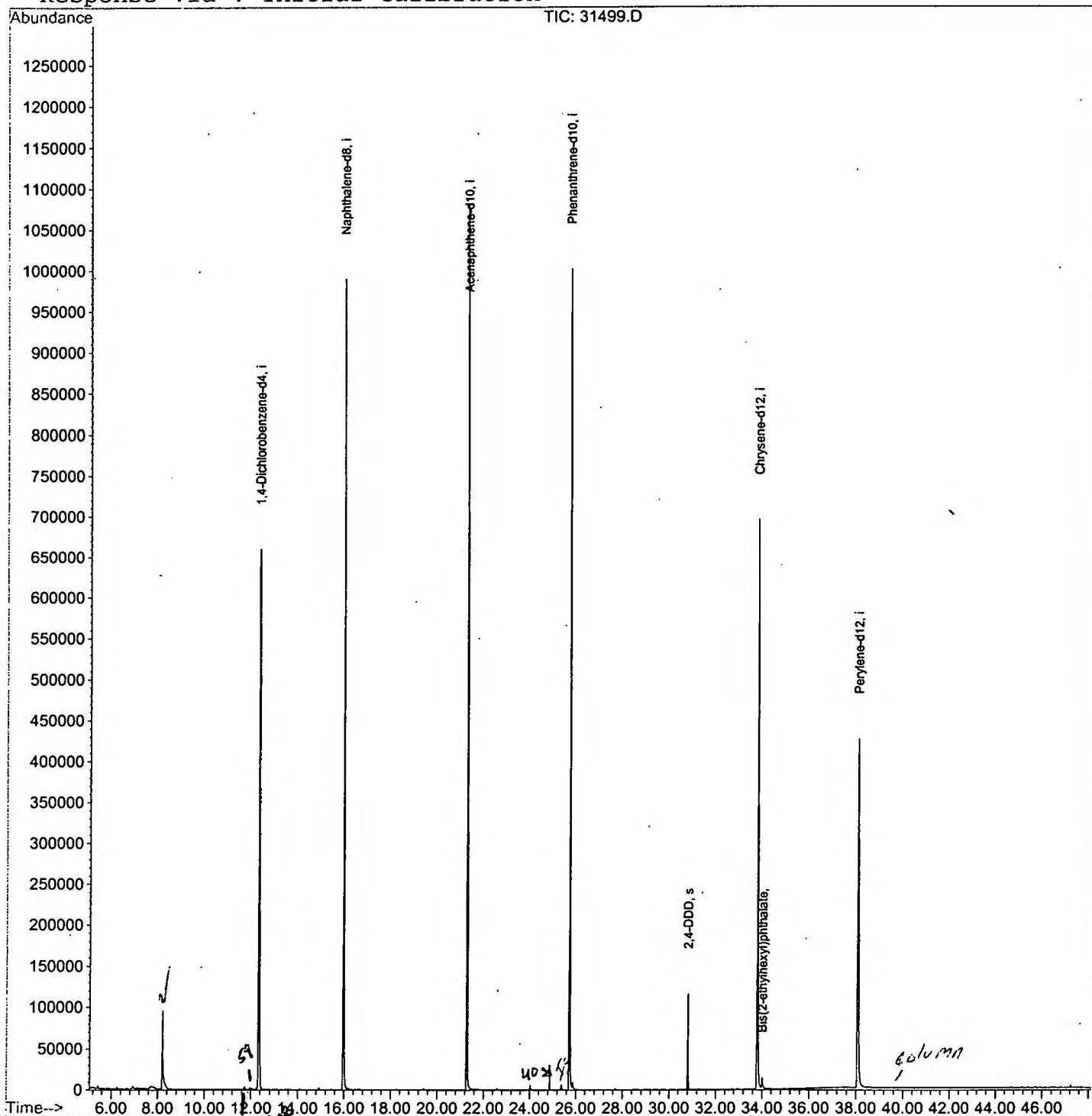
Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB2102\31499.D
Acq On : 23 Feb 2002 12:39 am
Sample : gc/ms scan 2/21
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 27 11:34 2002

Vial: 35
Operator:
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 27 11:30:15 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB2102\31499.D
 Acq On : 23 Feb 2002 12:39 am
 Sample : gc/ms scan 2/21
 Misc :
 MS Integration Params: LSCINT.P

Vial: 35
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0208.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

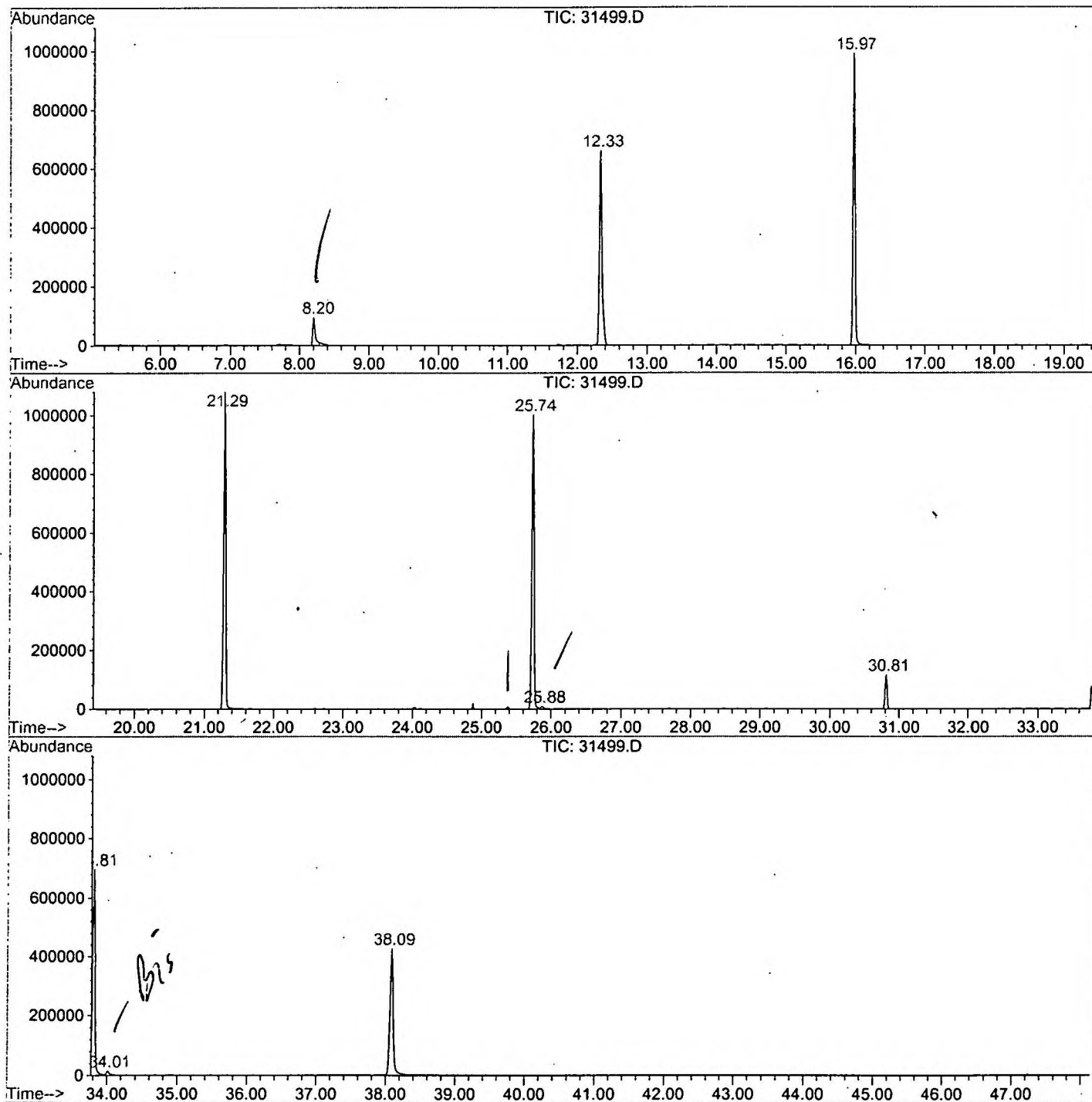
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.203	340	344	353	rBV	94828	226596	10.07%	1.948%
2	12.334	788	794	808	rVB	659181	1616261	71.84%	13.898%
3	15.970	1184	1190	1209	rBV	989776	2113636	93.94%	18.175%
4	21.285	1763	1769	1786	rBV	1081682	2249958	100.00%	19.347%
5	25.738	2247	2254	2265	rBV2	1002582	2131299	94.73%	18.327%
6	25.875	2265	2269	2282	rVB	8866	25261	1.12%	0.217%
7	30.814	2802	2807	2814	rVB	116290	233553	10.38%	2.008%
8	33.807	3125	3133	3148	rBV2	696305	1672988	74.36%	14.386%
9	34.009	3150	3155	3168	rVB	13905	44620	1.98%	0.384%
10	38.085	3590	3599	3616	rBV	425160	1315310	58.46%	11.310%

Sum of corrected areas: 11629482

31499.D GCMS0208.M Wed Feb 27 12:00:12 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB2102\31499.D
Operator :
Acquired : 23 Feb 2002 12:39 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: gc/ms scan 2/21
Misc Info :
Vial Number: 35
Quant File: GCMS0208.RES (RTE Integrator)



31499.D GCMS0208.M

Wed Feb 27 12:00:18 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\FEB2102\31499.D
Acq On : 23 Feb 2002 12:39 am
Sample : gc/ms scan 2/21
Misc :
MS Integration Params: LSCINT.P

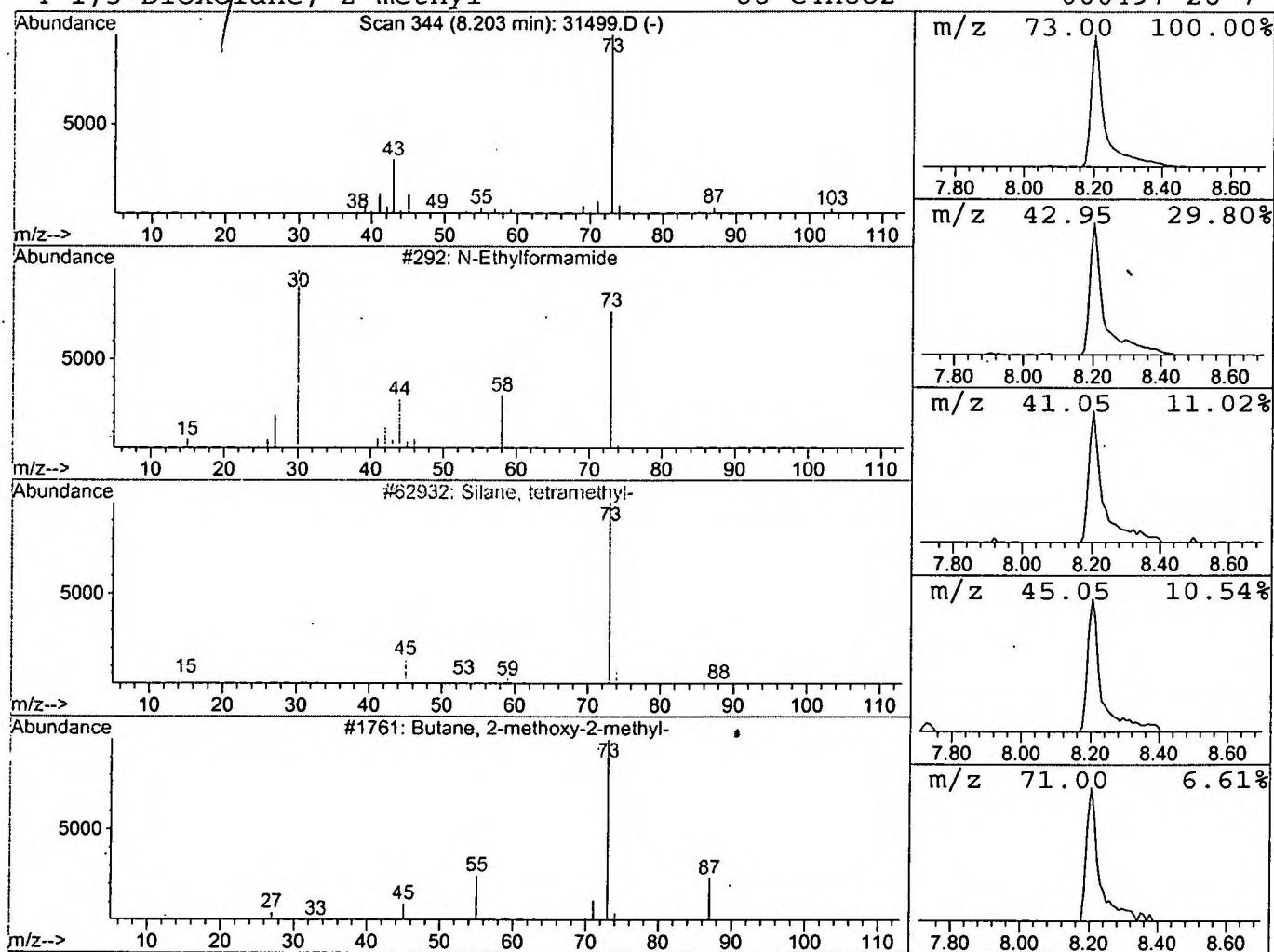
Vial: 35
Operator:
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

Peak Number 1 N-Ethylformamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	2.80 ug/l	226596	1,4-Dichlorobenzene-d4	12.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Ethylformamide	73	C3H7NO	000627-45-2	9
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	7



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\FEB2102\31499.D
Acq On : 23 Feb 2002 12:39 am
Sample : gc/ms scan 2/21
Misc :
MS Integration Params: LSCINT.P

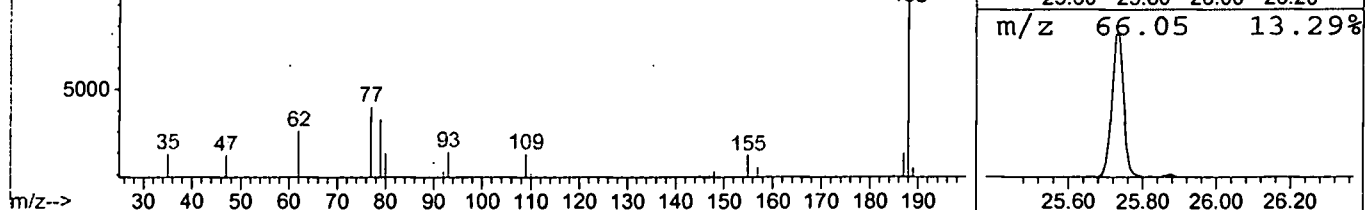
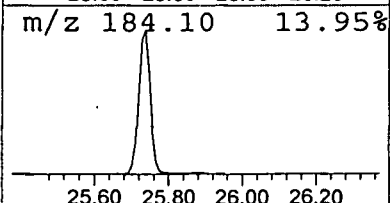
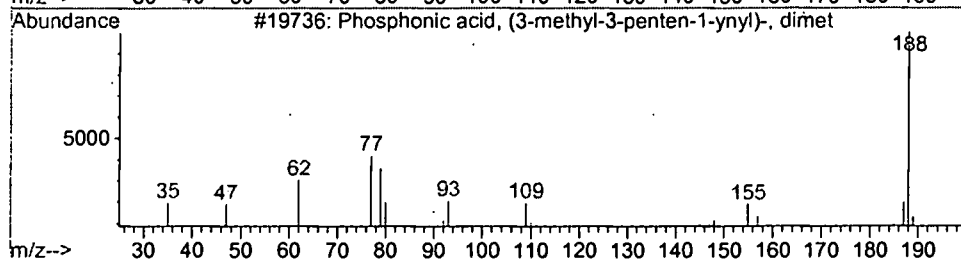
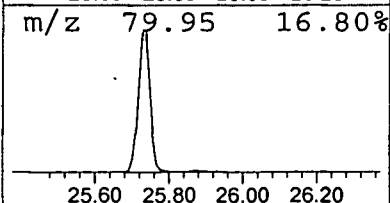
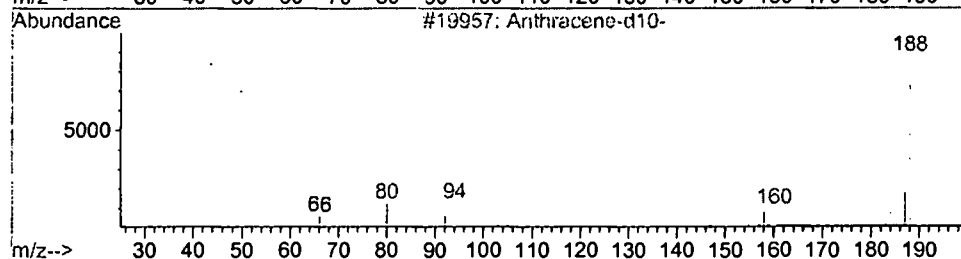
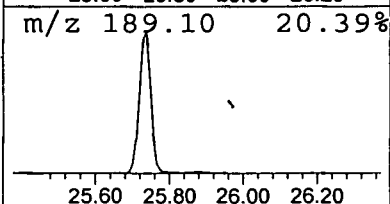
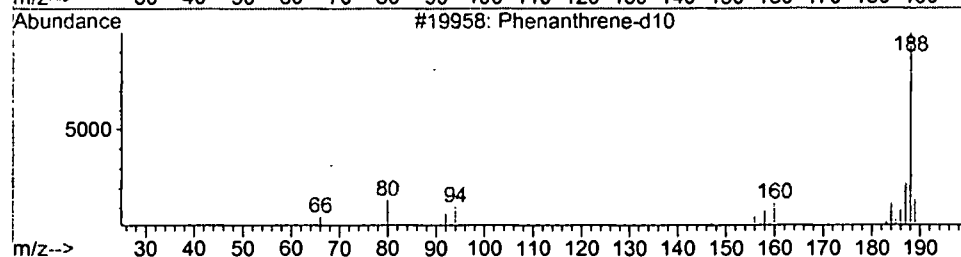
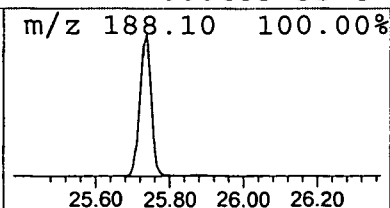
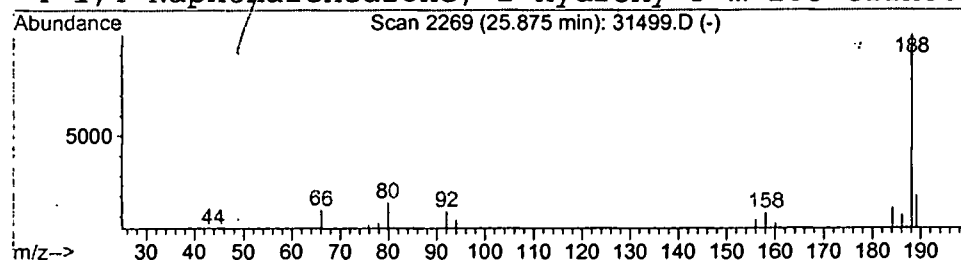
Vial: 35
Operator:
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.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.
25.88	0.24 ug/l	25261	Phenanthrene-d10	25.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene-d10	188	C14D10	001517-22-2	91
2			Anthracene-d10-	188	C14D10	001719-06-8	86
3			Phosphonic acid, (3-methyl-3-penten	188	C8H13O3P	022152-34-7	53
4			1,4-Naphthalenedione, 2-hydroxy-3-m	188	C11H8O3	000483-55-6	9



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 23 Feb 2002 12:39 am
 Data File: C:\HPCHEM\1\DATA\FEB2102\31499.D
 Name: gc/ms scan 2/21
 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
N-Ethylformamide	8.20	2.8 ug/l	226596	ISTD01	12.33	1616260	20.0
Phenanthrene-d10	25.88	0.2 ug/l	25261	ISTD04	25.74	2131300	20.0

31499.D GCMS0208.M Wed Feb 27 12:00:31 2002