

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/28/2002 Time: 12:30:55

Sample: AD30011
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/28/02 12:30 Ended: 1/28/02 12:30 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 AD30011 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-28-2002 Time: 12:31:12

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\DEC1801\30011.D

Vial: 14

Acq On : 19 Dec 2001 4:47 am

Operator: 209 GALOS

Sample : gc/ms scan 12/11

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Jan 15 11:04 2002

Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Dec 28 15:11:00 2001

Response via : Initial Calibration

DataAcq Meth : NP091101

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.67	152	751595	20.00	ug/l	-0.02
10) Naphthalene-d8	15.27	136	2878085	20.00	ug/l	-0.03
17) Acenaphthene-d10	20.55	164	1457171	20.00	ug/l	-0.02
29) Phenanthrene-d10	24.98	188	2073307	20.00	ug/l	-0.02
38) Chrysene-d12	33.01	240	1274608	20.00	ug/l	-0.02
44) Perylene-d12	37.10	264	845640	20.00	ug/l	-0.02

System Monitoring Compounds

37) 2,4-DDD	30.05	235	117730	5.16	ug/mL	-0.02
Spiked Amount	5.000		Recovery	= 103.20%		

Target Compounds

					Qvalue
2) N-nitroso-dimethyl amine	5.31	74	168	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	15.34	128	1202	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	21.10	165	203	N.D.	
25) Diethylphthalate	22.10	149	994	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

30011.D GCMS1126.M

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Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D

Vial: 14

Acq On : 19 Dec 2001 4:47 am

Operator: 209 GALOS

Sample : gc/ms scan 12/11

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Jan 15 11:04 2002

Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Dec 28 15:11:00 2001

Response via : Initial Calibration

DataAcq Meth : NP091101

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	25.03	178	355	N.D.		
34) Anthracene	25.03	178	355	N.D.		
35) Di-n-butylphthalate	27.01	149	6016	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	33.02	228	3108	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.29	149	4787	N.D.		
43) Chrysene	33.02	228	3108	N.D.		
45) Di-n-Octylphthalate	34.86	149	105	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	37.10	252	3085	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

30011.D GCMS1126.M

Tue Jan 15 11:05:44 2002

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Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D

Vial: 14

Acq On : 19 Dec 2001 4:47 am

Operator: 209 GALOS

Sample : gc/ms scan 12/11

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Jan 15 11:04 2002

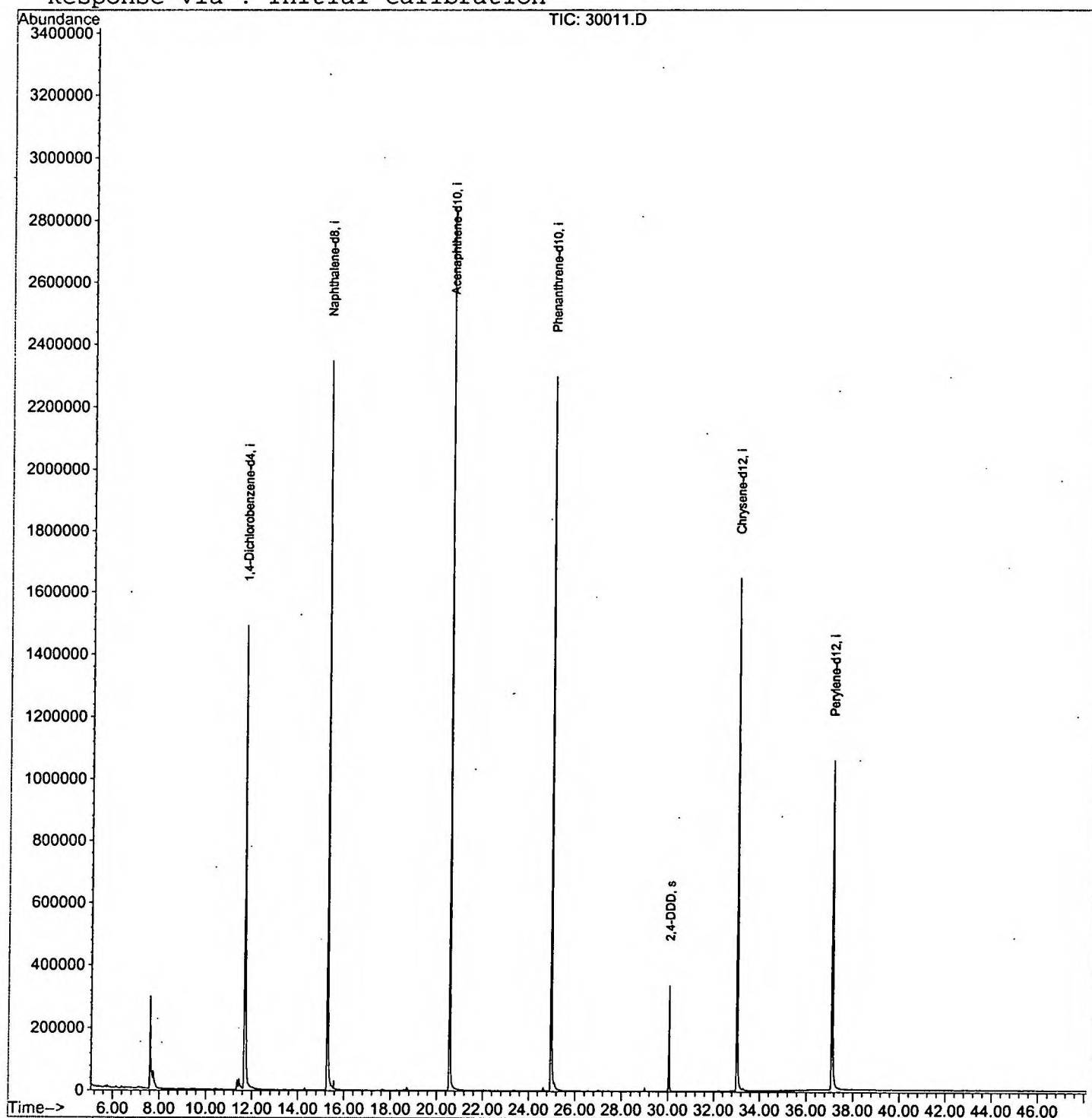
Quant Results File: GCMS1126.RES

Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Dec 28 15:11:00 2001

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D

Acq On : 19 Dec 2001 4:47 am

Sample : gc/ms scan 12/11

Misc :

MS Integration Params: LSCINT.P

Vial: 14

Operator: 209 GALOS

Inst : GC/MS 209

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS1126.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	7.598	274	278	290	rBV	296293	853569	13.87%	2.781%
2	7.736	290	293	315	rVB	55831	248006	4.03%	0.808%
3	11.362	683	688	692	rBV	31007	77848	1.26%	0.254%
4	11.436	692	696	703	rVV2	27882	83075	1.35%	0.271%
5	11.674	716	722	750	rBV	1487377	4488616	72.93%	14.624%
6	15.273	1109	1114	1132	rBV	2346072	5560963	90.36%	18.118%
7	20.552	1683	1689	1712	rBV	2843952	6154356	100.00%	20.052%
8	24.977	2164	2171	2183	rBV2	2297135	5436362	88.33%	17.712%
9	30.053	2718	2724	2736	rBV	336160	678341	11.02%	2.210%
10	33.009	3039	3046	3065	rBV2	1647814	4030180	65.48%	13.131%
11	37.104	3485	3492	3512	rBV	1055697	3081374	50.07%	10.039%

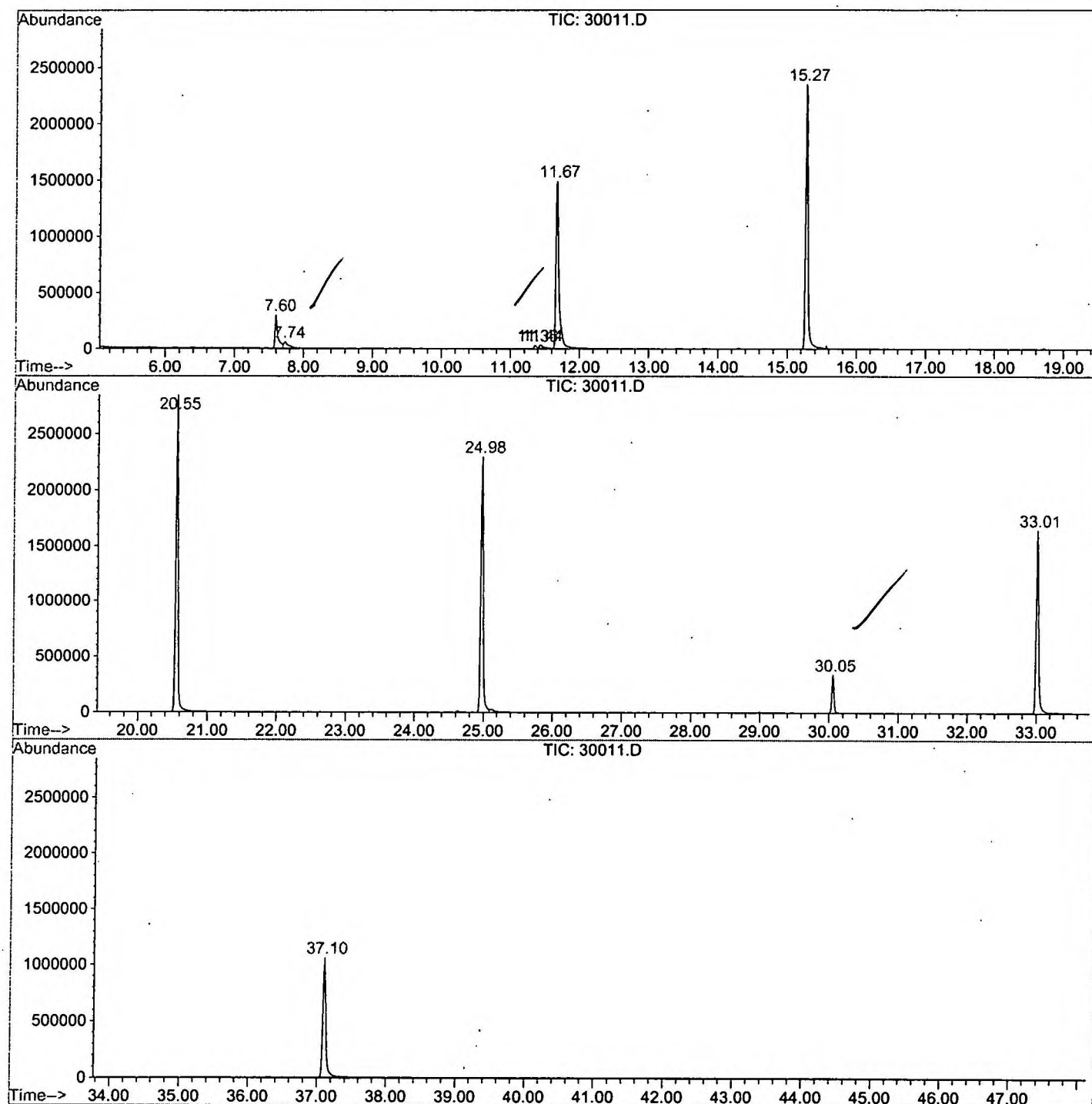
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30011.D GCMS1126.M

Tue Jan 15 15:30:40 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC1801\30011.D
Operator : 209 GALOS
Acquired : 19 Dec 2001 4:47 am using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms scan 12/11
Misc Info :
Vial Number: 14
Quant File : GCMS1126.RES (RTE Integrator)



30011.D GCMS1126.M

Tue Jan 15 15:30:45 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D
 Acq On : 19 Dec 2001 4:47 am
 Sample : gc/ms scan 12/11
 Misc :
 MS Integration Params: LSCINT.P

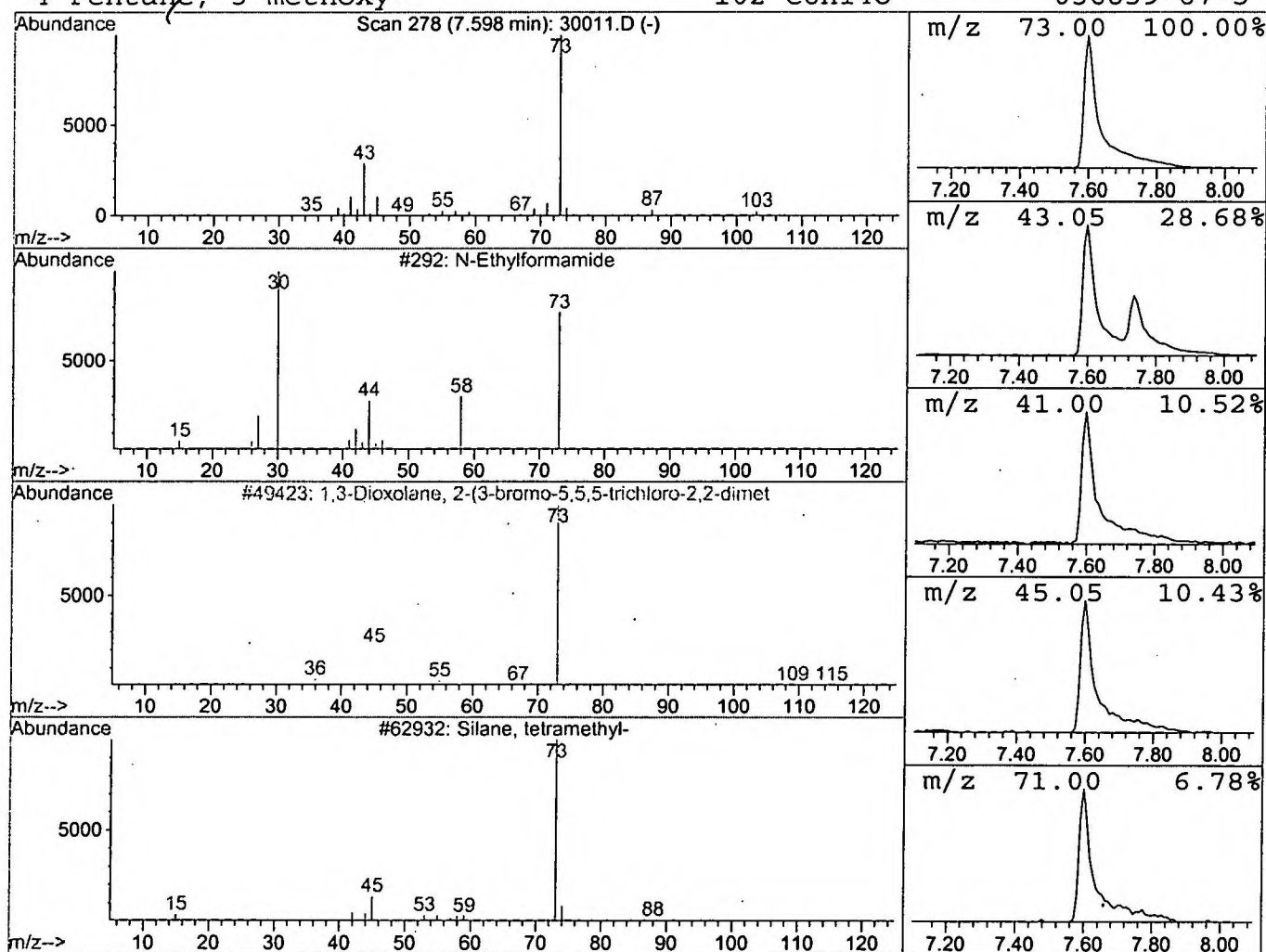
Vial: 14
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.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.
7.60	3.80 ug/l	853569	1,4-Dichlorobenzene-d4	11.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Ethylformamide	73	C3H7NO	000627-45-2	9
2			1,3-Dioxolane, 2-(3-bromo-5,5,5-tri	352	C10H16BrCl3O2	000000-00-0	9
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D
Acq On : 19 Dec 2001 4:47 am
Sample : gc/ms scan 12/11
Misc :
MS Integration Params: LSCINT.P

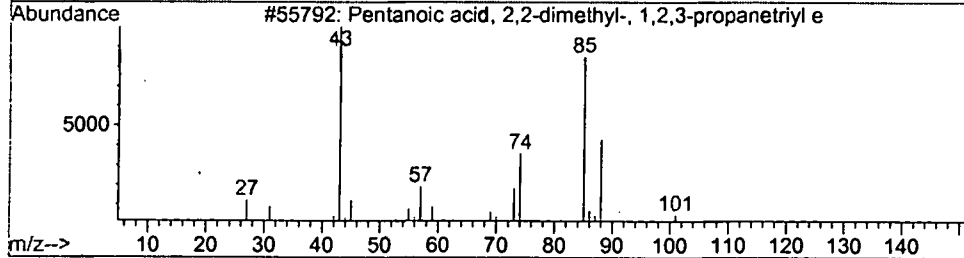
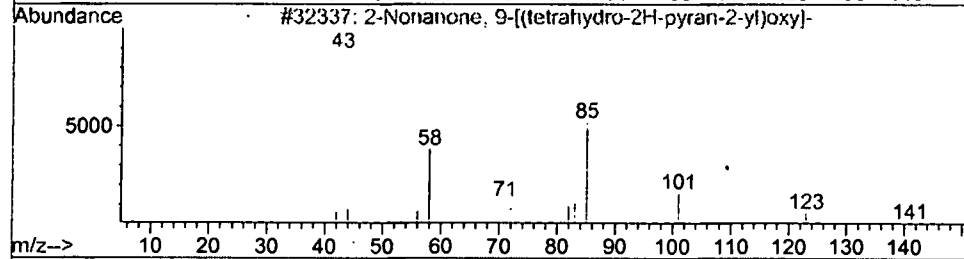
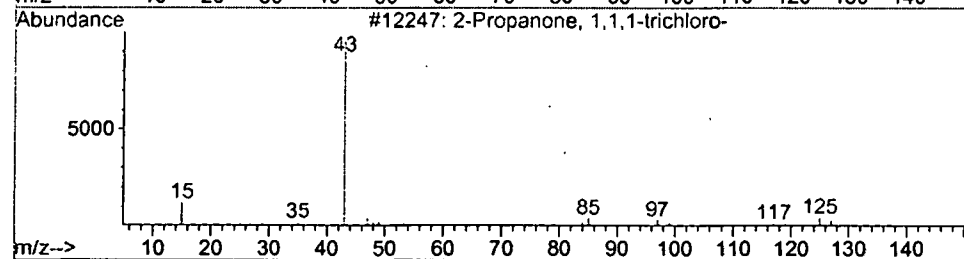
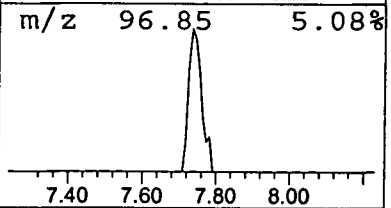
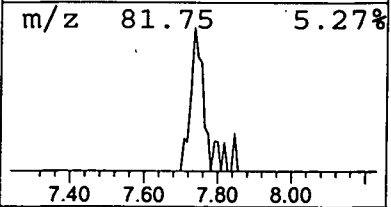
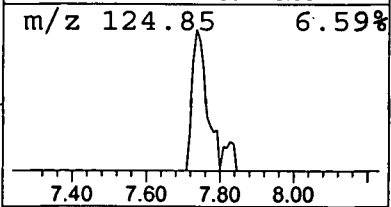
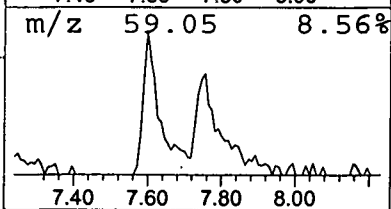
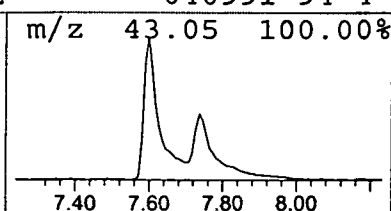
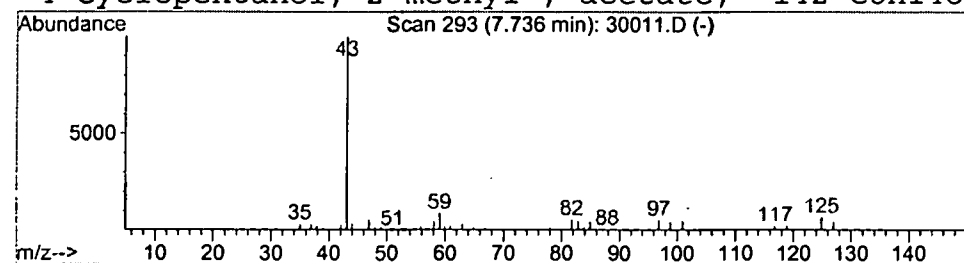
Vial: 14
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

Peak Number 2 2-Propanone, 1,1,1-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.74	1.11 ug/l	248006	1,4-Dichlorobenzene-d4	11.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanone, 1,1,1-trichloro-	160	C3H3Cl3O	000918-00-3	38
2			2-Nonanone, 9-[(tetrahydro-2H-pyran	242	C14H26O3	054699-41-1	9
3			Pentanoic acid, 2,2-dimethyl-, 1,2,	428	C24H44O6	057346-62-0	7
4			Cyclopentanol, 2-methyl-, acetate,	142	C8H14O2	040991-94-4	7



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1801\30011.D
 Acq On : 19 Dec 2001 4:47 am
 Sample : gc/ms scan 12/11
 Misc :
 MS Integration Params: LSCINT.P

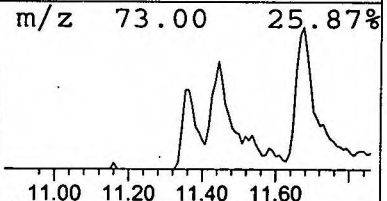
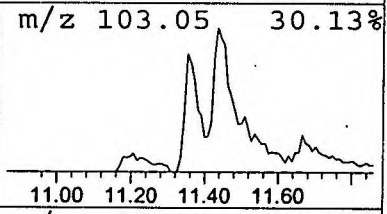
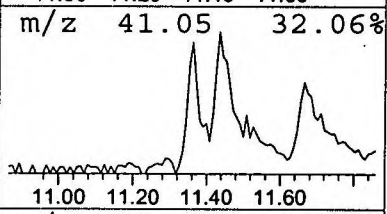
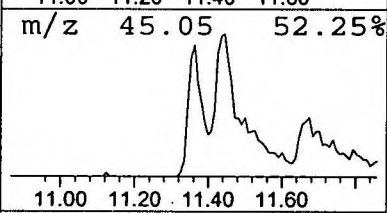
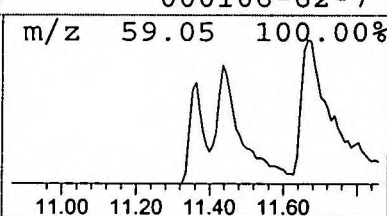
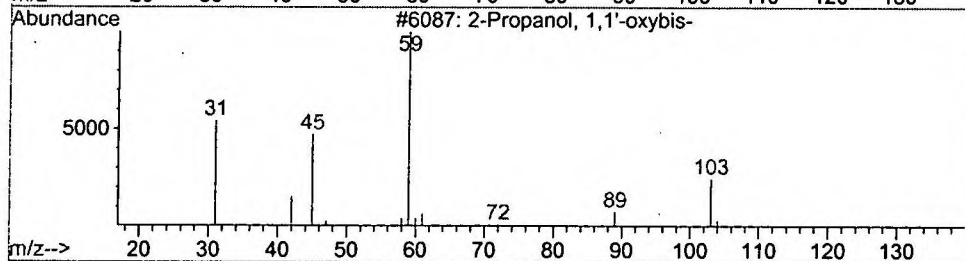
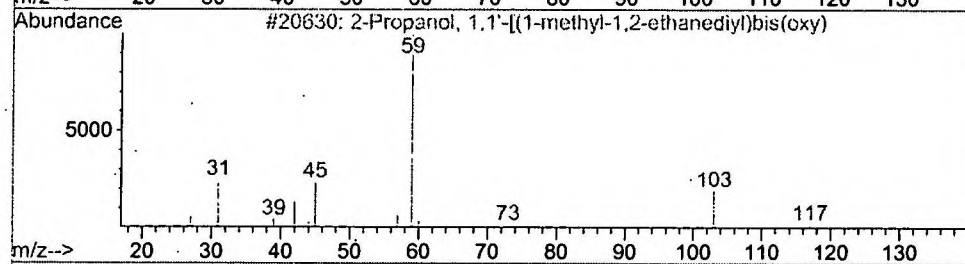
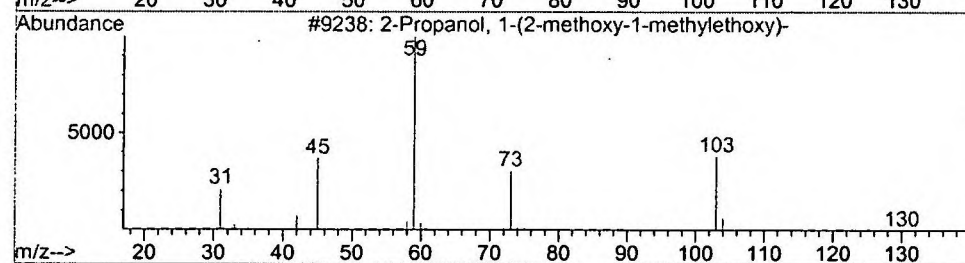
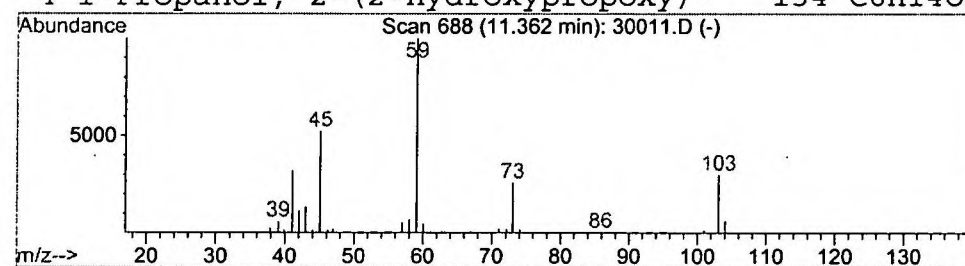
Vial: 14
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 3 2-Propanol, 1-(2-methoxy-1-met Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	0.35 ug/l	77848	1,4-Dichlorobenzene-d4	11.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methylet	148	C7H16O3	020324-32-7	72	
2	2-Propanol, 1,1'-[(1-methyl-1,2-eth	192	C9H20O4	001638-16-0	53	
3	2-Propanol, 1,1'-oxybis-	134	C6H14O3	000110-98-5	50	
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	45	



Library Search Compound Report

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 Misc :
 MS Integration Params: LSCINT.P

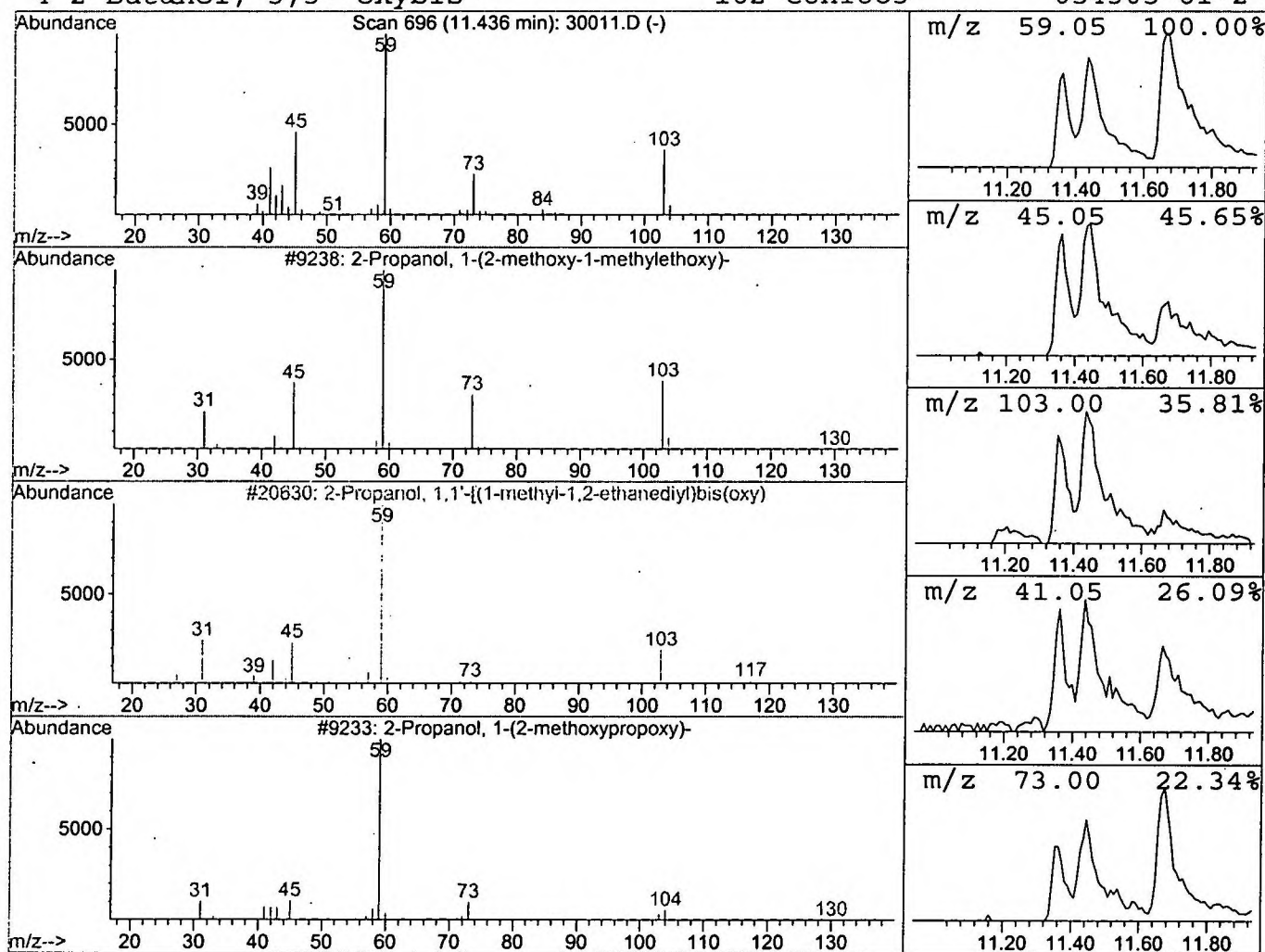
Vial: 14
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 4 2-Propanol, 1-(2-methoxy-1-met Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	0.37 ug/l	83075	1,4-Dichlorobenzene-d4	11.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methylet	148	C7H16O3	020324-32-7	72
2			2-Propanol, 1,1'-[(1-methyl-1,2-eth	192	C9H20O4	001638-16-0	53
3			2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	40
4			2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	39



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 19 Dec 2001 4:47 am
Data File: C:\HPCHEM\1\DATA\DEC1801\30011.D
Name: gc/ms scan 12/11
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS1126.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	7.60	3.8	ug/l	853569	ISTD01	11.67	4488620	20.0
2-Propanone, 1,1,1-t	7.74	1.1	ug/l	248006	ISTD01	11.67	4488620	20.0
2-Propanol, 1-(2-met	11.36	0.3	ug/l	77848	ISTD01	11.67	4488620	20.0
2-Propanol, 1-(2-met	11.44	0.4	ug/l	83075	ISTD01	11.67	4488620	20.0

30011.D GCMS1126.M Tue Jan 15 15:31:08 2002