

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 01/18/2002 Time: 17:17:58

Sample: AD30487
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 1/18/02 17:17 Ended: 1/18/02 17:17 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		93		5.0

Comments for Sample AD30487 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-18-2002 Time: 17:18:08

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\DEC3101\30487.D

Vial: 22

Acq On : 1 Jan 2002 7:21 am

Operator: 209 GALOS

Sample : gcms scan 12/14a

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Jan 1 8:10 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 : GCMS1126

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	11.68	152	780730	20.00	ug/l	-0.02
10) Naphthalene-d8	15.28	136	3020681	20.00	ug/l	-0.02
17) Acenaphthene-d10	20.56	164	1574464	20.00	ug/l	-0.01
29) Phenanthrene-d10	24.99	188	2189274	20.00	ug/l	0.00
38) Chrysene-d12	33.03	240	1261715	20.00	ug/l	0.00
44) Perylene-d12	37.12	264	680064	20.00	ug/l	0.00

System Monitoring Compounds

37) 2,4-DDD	30.07	235	111440	4.63	ug/mL	0.00
Spiked Amount	5.000		Recovery	=	92.60%	

Target Compounds

					Qvalue
2) N-nitroso-dimethyl amine	5.18	74	1040	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	655	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	21.24	165	176	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) Azobenzene/ 1,2-Diphenylhyd	0.00	77	0	N.D.	

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

30487.D GCMS0103.M Tue Jan 15 14:27:22 2002

Page 1

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC3101\30487.D

Vial: 22

Acq On : 1 Jan 2002 7:21 am

Operator: 209 GALOS

Sample : gcms scan 12/14a

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Jan 1 8:10 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 : GCMS1126

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	25.05	178	231	N.D.		
34) Anthracene	25.05	178	231	N.D.		
35) Di-n-butylphthalate	27.05	149	4336	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	3173	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.30	149	3304	N.D.		
43) Chrysene	33.02	228	3173	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	37.14	252	2449	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

30487.D GCMS0103.M

Tue Jan 15 14:27:25 2002

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30487.D

Acq On : 1 Jan 2002 7:21 am

Sample : gcms scan 12/14a

Misc :

MS Integration Params: GOOFY.P

Quant Time: Jan 1 8:10 2002

Vial: 22

Operator: 209 GALOS

Inst : GC/MS 209

Multiplr: 1.00

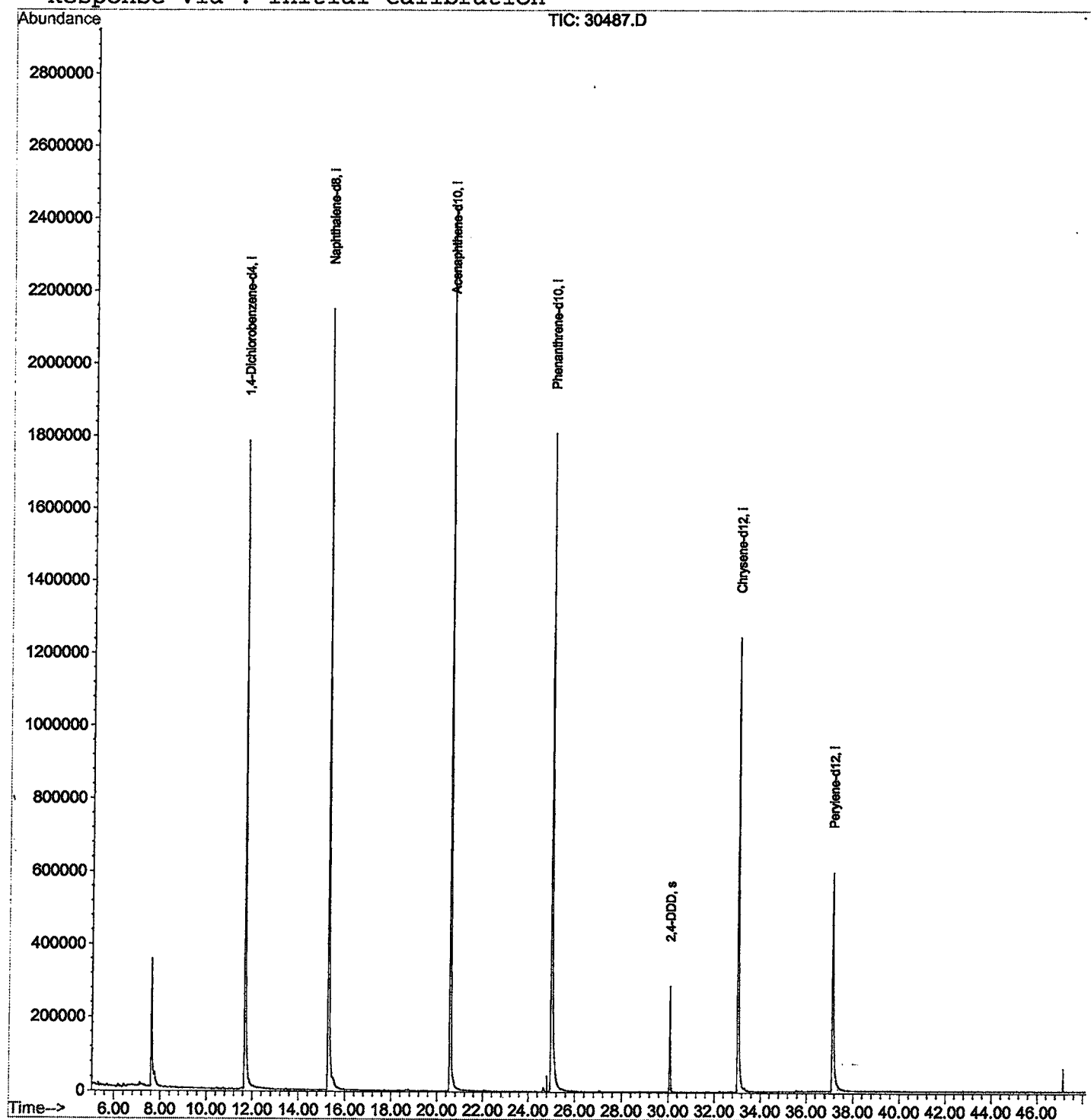
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Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Mon Jan 07 10:31:02 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30487.D

Vial: 22

Acq On : 1 Jan 2002 7:21 am

Operator: 209 GALOS

Sample : gcms scan 12/14a

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

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	7.617	276	280	292	rBV	347750	983820	14.96%	3.095%
2	7.746	292	294	316	rVB	39357	177578	2.70%	0.559%
3	11.675	717	722	753	rBV	1781226	4864291	73.95%	15.301%
4	15.283	1109	1115	1135	rBV	2147712	5983447	90.96%	18.821%
5	20.562	1684	1690	1718	rBV	2433552	6577814	100.00%	20.691%
6	24.987	2165	2172	2201	rBV2	1806995	6043842	91.88%	19.011%
7	30.063	2720	2725	2737	rBV	286361	686855	10.44%	2.161%
8	33.029	3041	3048	3072	rBV2	1243886	4008780	60.94%	12.610%
9	37.123	3486	3494	3523	rBV2	595139	2464508	37.47%	7.752%

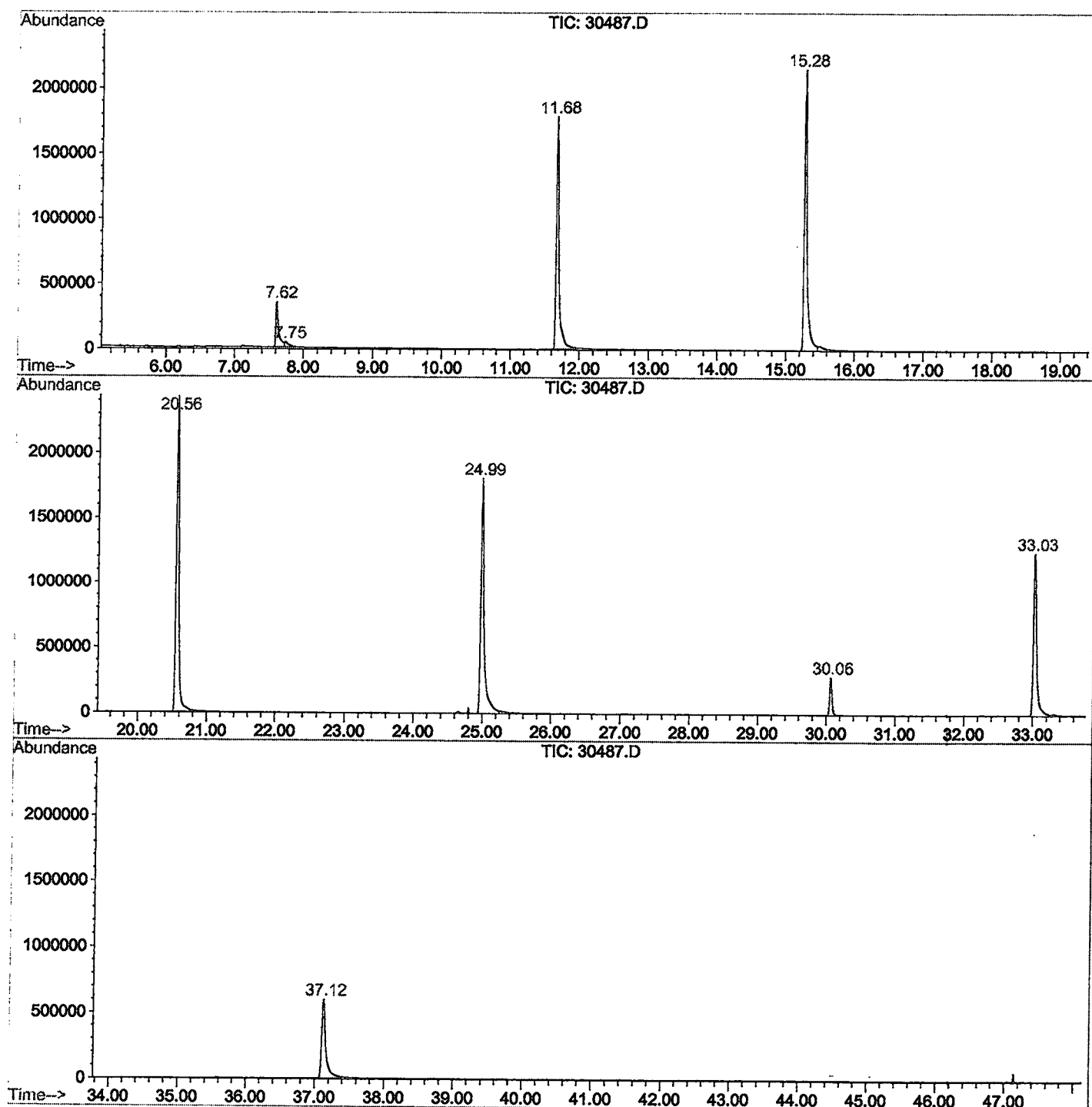
Sum of corrected areas: 31790935

30487.D GCMS1126.M

Tue Jan 15 14:51:54 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC3101\30487.D
 Operator : 209 GALOS
 Acquired : 1 Jan 2002 7:21 am using AcqMethod GCMS1126
 Instrument : GC/MS 209
 Sample Name: gcms scan 12/14a
 Misc Info :
 Vial Number: 22
 Quant File : GCMS1126.RES (RTE Integrator)



30487.D GCMS1126.M

Tue Jan 15 14:52:00 2002

Library Search Compound Report

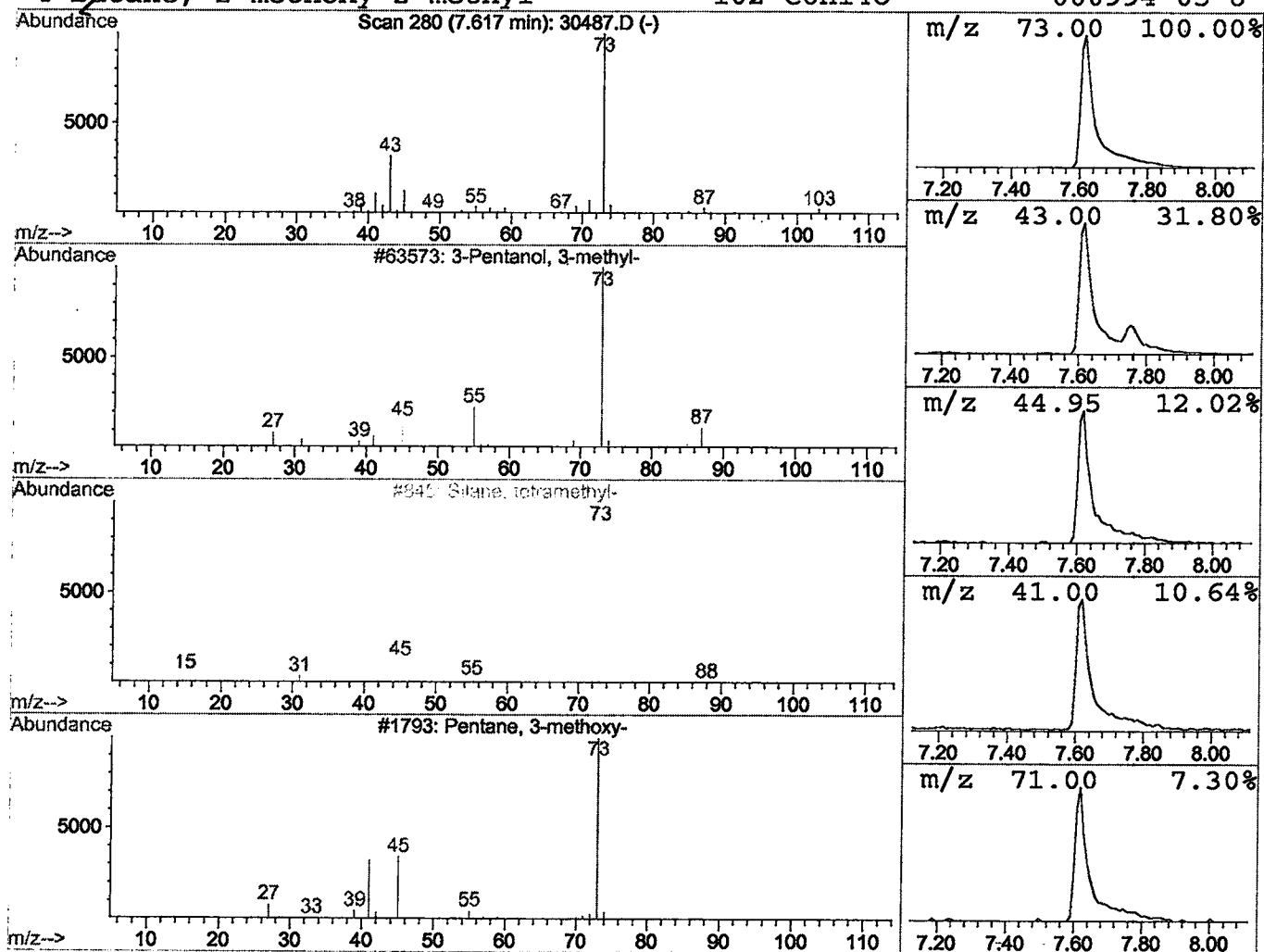
Data File : C:\HPCHEM\1\DATA\DEC3101\30487.D
 Acq On : 1 Jan 2002 7:21 am
 Sample : gcms scan 12/14a
 Misc :
 MS Integration Params: LSCINT.P

Vial: 22
 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 3-Pentanol, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
7.62	4.05 ug/l	983820	1,4-Dichlorobenzene-d4	11.68		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 3-methyl-		102	C6H14O	000077-74-7	36
2	Silane, tetramethyl-		88	C4H12Si	000075-76-3	9
3	Pentane, 3-methoxy-		102	C6H14O	036839-67-5	9
4	Butane, 2-methoxy-2-methyl-		102	C6H14O	000994-05-8	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30487.D
 Acq On : 1 Jan 2002 7:21 am
 Sample : gcms scan 12/14a
 Misc :
 MS Integration Params: LSCINT.P

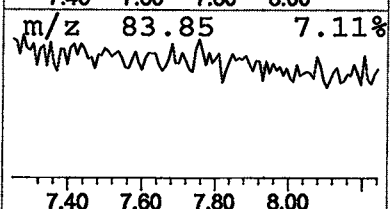
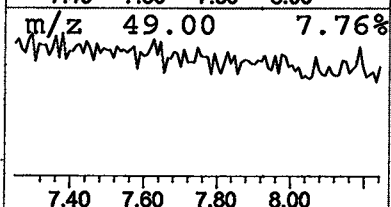
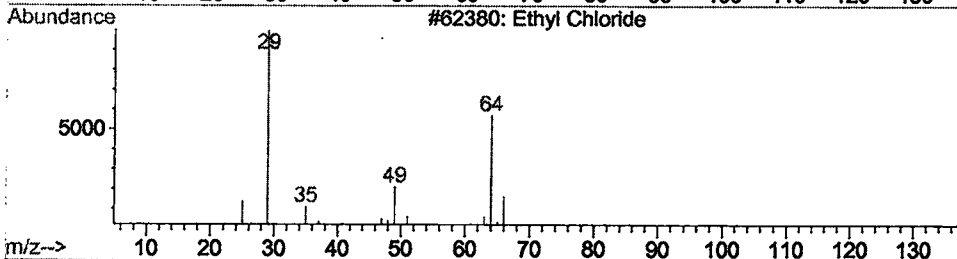
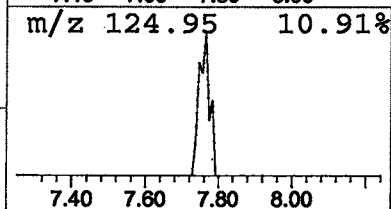
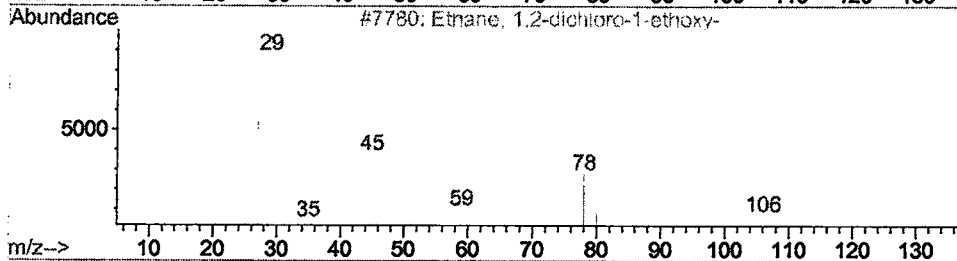
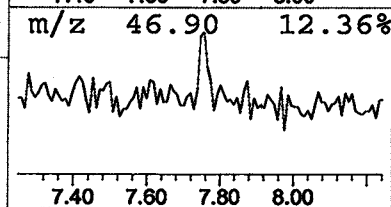
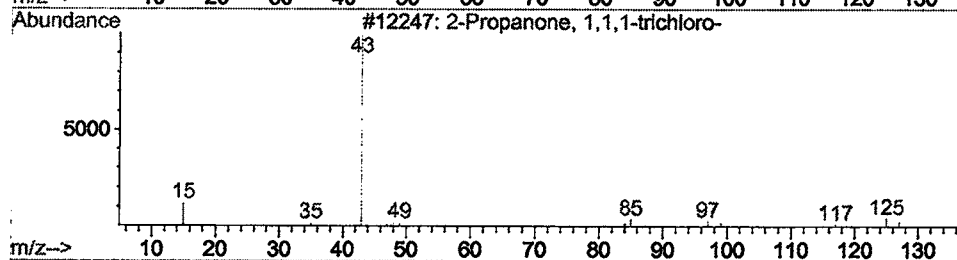
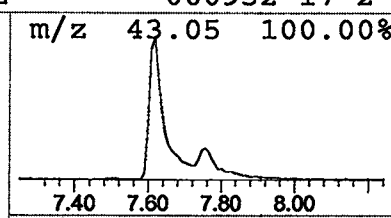
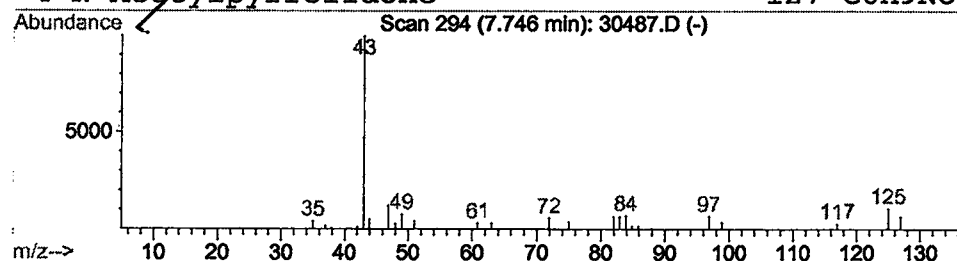
Vial: 22
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

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.
7.75	0.73 ug/l	177578	1,4-Dichlorobenzene-d4	11.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanone, 1,1,1-trichloro-	160	C3H3Cl3O	000918-00-3	50
2			Ethane, 1,2-dichloro-1-ethoxy-	142	C4H8Cl2O	000623-46-1	9
3			Ethyl Chloride	64	C2H5Cl	000075-00-3	7
4			N-Acetylpyrrolidone	127	C6H9NO2	000932-17-2	5



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 1 Jan 2002 7:21 am
 Data File: C:\HPCHEM\1\DATA\DEC3101\30487.D
 Name: gcms scan 12/14a
 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
3-Pentanol, 3-methyl	7.62	4.0	ug/l	983820	ISTD01	11.68	4864290	20.0
2-Propanone, 1,1,1-t	7.75	0.7	ug/l	177578	ISTD01	11.68	4864290	20.0

30487.D GCMS1126.M Tue Jan 15 14:52:13 2002