User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 12/28/2001 Time: 14:53:21

Sample: AD29235
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/28/01 14:52 Ended: 12/28/01 14:52 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		127		5.0

mments for Sample AD29235 - Analysis $GCMS

estchester Dept. of Labs & Research
ser: Vinci, David L
ate: 12-28-2001 Time: 14:53:33

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D Vial: 8
Acq On : 12 Dec 2001 6:34 am Operator: 209 GALOS
Sample : gc/ms scan 12/4 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 19 12:46 2001 Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Nov 28 15:06:24 2001
Response via : Initial Calibration
DataAcq Meth : GCMS1126

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.67	152	875590	20.00	ug/l	-0.03
10) Naphthalene-d8	15.28	136	3340047	20.00	ug/l	-0.03
17) Acenaphthene-d10	20.55	164	1712414	20.00	ug/l	-0.03
29) Phenanthrene-d10	24.97	188	2346540	20.00	ug/l	-0.02
38) Chrysene-d12	33.00	240	1485832	20.00	ug/l	-0.03
44) Perylene-d12	37.10	264	1051283	20.00	ug/l	-0.02
System Monitoring Compounds						
37) 2,4-DDD	30.05	235	177162	6.34	ug/mL	-0.02
Spiked Amount 5.000			Recovery = 126.80%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.30	74	210	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	15.33	128	1655	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.	d	
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	21.02	165	216	N.D.		
25) Diethylphthalate	22.09	149	660	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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D Vial: 8
Acq On : 12 Dec 2001 6:34 am Operator: 209 GALOS
Sample : gc/ms scan 12/4 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 19 12:46 2001 Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Nov 28 15:06:24 2001
Response via : Initial Calibration
DataAcq Meth : GCMS1126

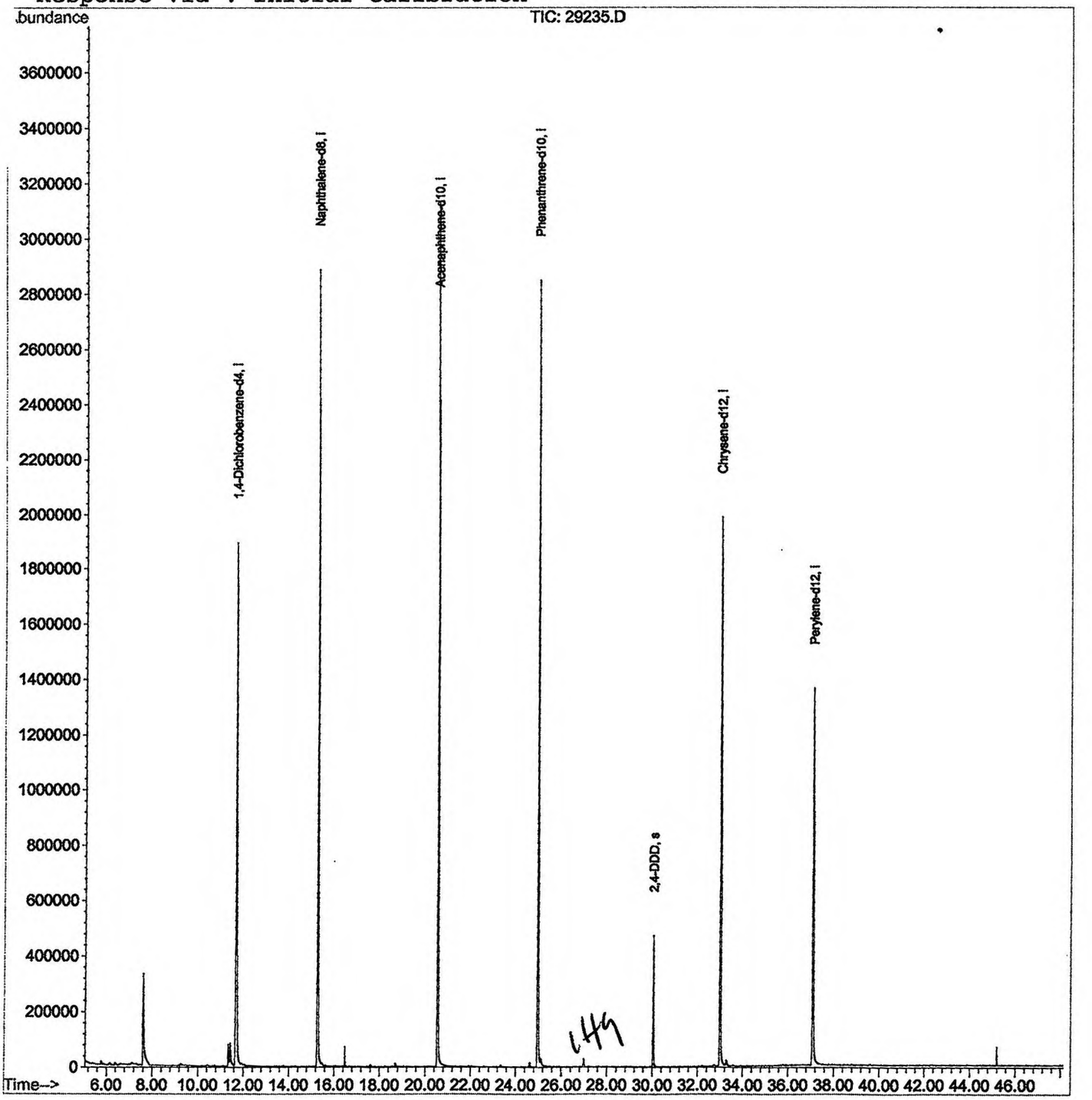
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32)	Hexachlorobenzene	0.00	284	0	N.D.	
33)	Phenanthrene	25.04	178	237	N.D.	
34)	Anthracene	25.04	178	237	N.D.	
35)	Di-n-butylphthalate	26.98	149	38831	N.D.	
36)	Fluoranthene	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	33.01	228	3186	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.29	149	18341	N.D.	
43)	Chrysene	33.01	228	3186	N.D.	
45)	Di-n-Octylphthalate	35.02	149	793	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	4301	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

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D Vial: 8
Acq On : 12 Dec 2001 6:34 am Operator: 209 GALOS
Sample : gc/ms scan 12/4 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 19 12:46 2001 Quant Results File: GCMS1126.RES

Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Dec 14 12:19:30 2001
Response via : Initial Calibration



LSC Area Percent Report

ata File : C:\HPCHEM\1\DATA\DEC1001\29235.D Vial: 8
eq On : 12 Dec 2001 6:34 am Operator: 209 GALOS
ample : gc/ms scan 12/4 Inst : GC/MS 209
isc : Multiplr: 1.00
S Integration Params: LSCINT.P

ethod : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
itle : 209 625/TCLP calibration table
moothing : OFF Filtering: 5
ampling : 1 Min Area: 1 % of largest Peak
tart Thrs: 0.2 Max Peaks: 100
top Thrs : 0 Peak Location: TOP

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

Signal : TIC

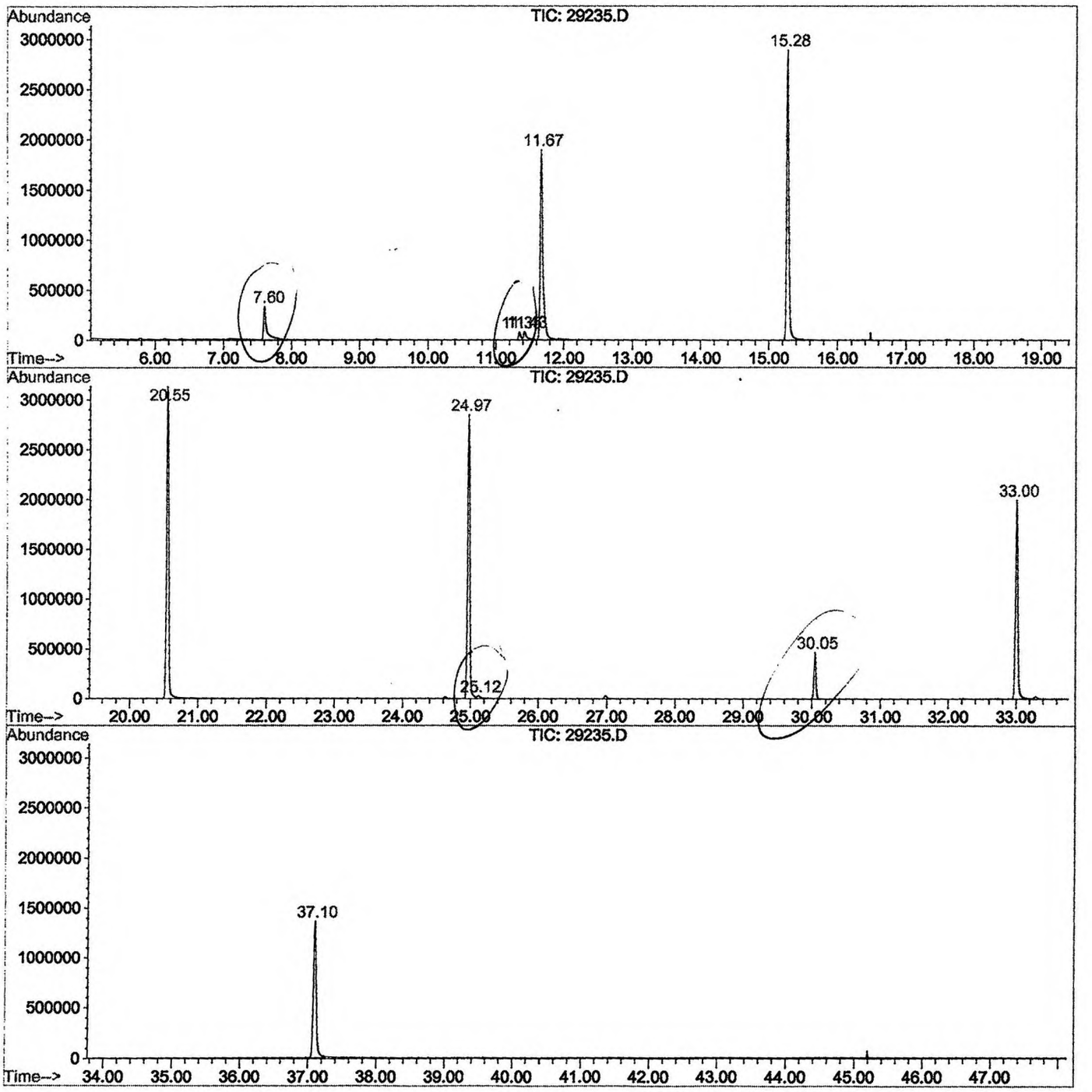
eak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.603	275	279	309	rBV	329226	1030334	14.72%	2.882%
2	11.348	683	687	692	rBV	78314	183997	2.63%	0.515%
3	11.431	692	696	715	rVB	81854	269635	3.85%	0.754%
4	11.670	715	722	745	rBV	1894119	5114626	73.06%	14.306%
5	15.277	1109	1115	1133	rBV	2886304	6331790	90.45%	17.710%
6	20.547	1683	1689	1711	rBV	3136165	7000548	100.00%	19.581%
7	24.972	2164	2171	2183	rBV2	2849001	6256330	89.37%	17.499%
8	25.119	2184	2187	2198	rVB2	24542	79646	1.14%	0.223%
9	30.049	2718	2724	2731	rBV	472046	954783	13.64%	2.671%
10	33.005	3039	3046	3064	rBV2	1988880	4658009	66.54%	13.029%
11	37.099	3484	3492	3509	rBV	1362569	3872131	55.31%	10.831%

Sum of corrected areas: 35751829

29235.D GCMS1126.M Wed Dec 19 12:52:37 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC1001\29235.D
Operator : 209 GALOS
Acquired : 12 Dec 2001 6:34 am using AcqMethod GCMS1126
Instrument : GC/MS 209
Sample Name: gc/ms scan 12/4
Misc Info :
Vial Number: 8
Quant File :GCMS1126.RES (RTE Integrator)



29235.D GCMS1126.M Wed Dec 19 12:52:44 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D
Acq On : 12 Dec 2001 6:34 am
Sample : gc/ms scan 12/4
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

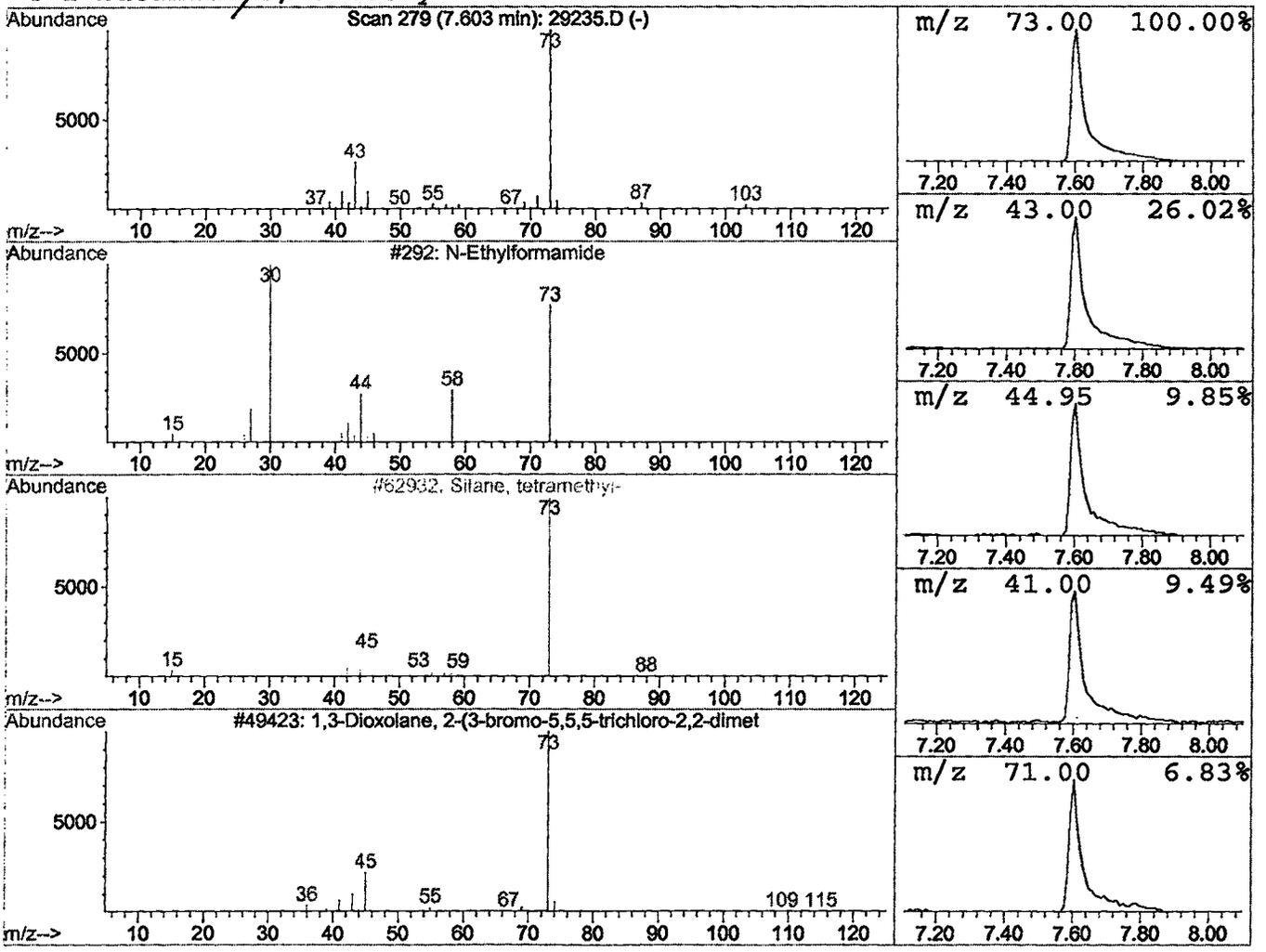
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 N-Ethylformamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	4.03 ug/l	1030330	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	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3	1,3-Dioxolane, 2-(3-bromo-5,5,5-tri	352	C10H16BrCl3O2	000000-00-0	9
4	1-Butanamine, 3-methyl-	87	C5H13N	000107-85-7	7



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D
Acq On : 12 Dec 2001 6:34 am
Sample : gc/ms scan 12/4
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

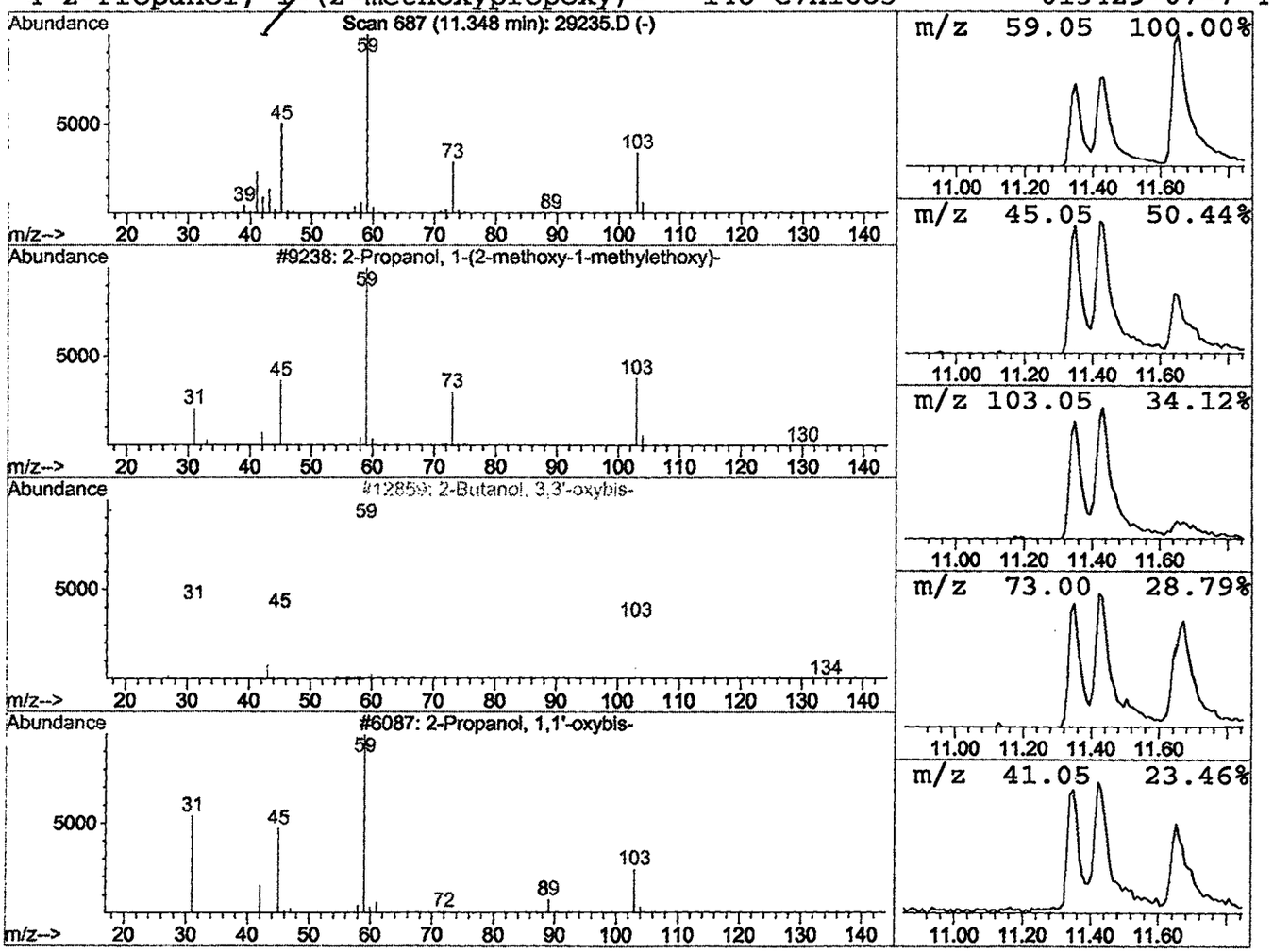
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	0.72 ug/l	183997	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-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	64
3	2-Propanol, 1,1'-oxybis-	134	C6H14O3	000110-98-5	56
4	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	47



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1001\29235.D Vial: 8
Acq On : 12 Dec 2001 6:34 am Operator: 209 GALOS
Sample : gc/ms scan 12/4 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

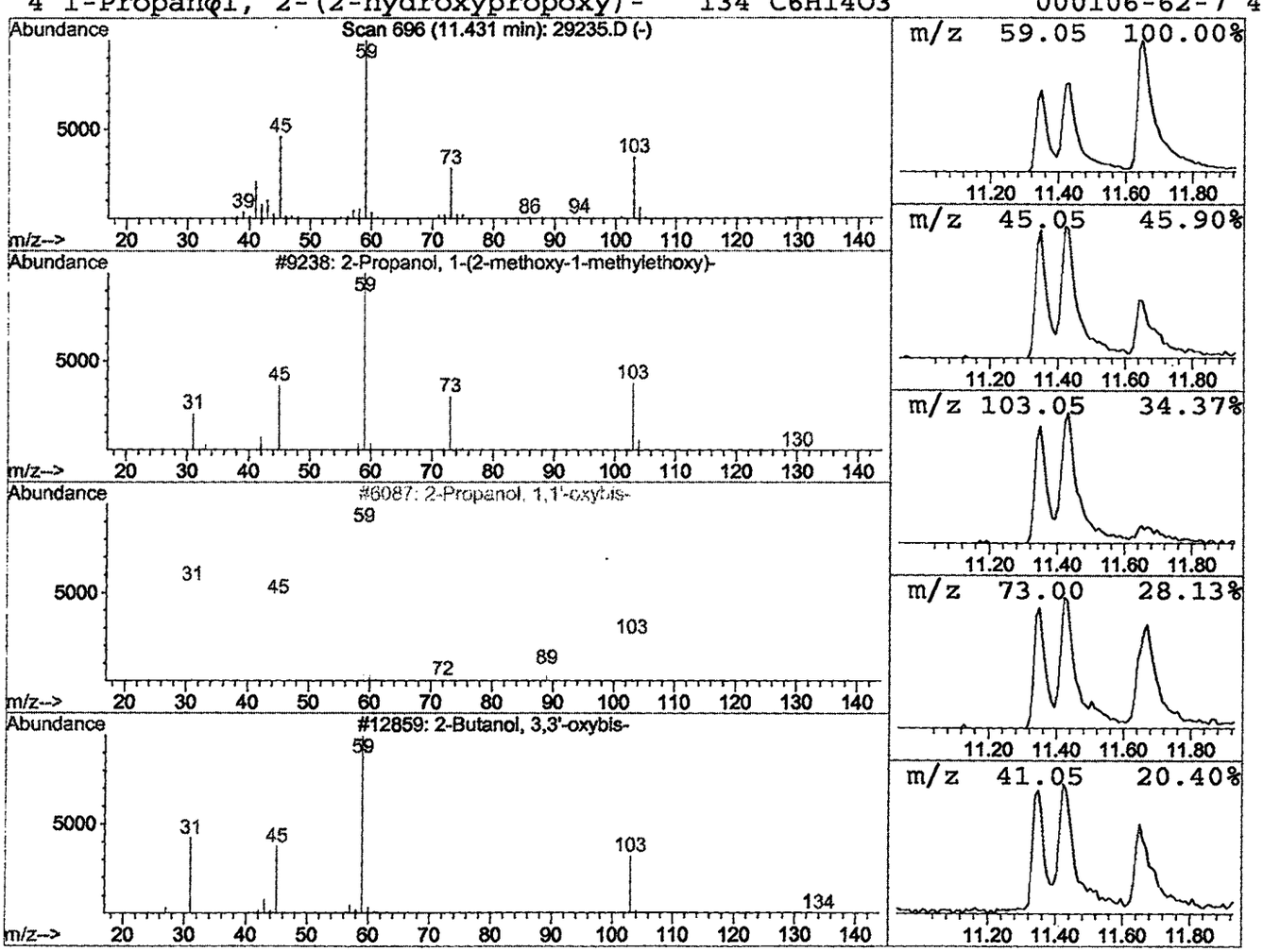
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	1.05 ug/l	269635	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	90
2	2-Propanol, 1,1'-oxybis-	134	C6H14O3	000110-98-5	56
3	2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	50
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	42



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 12 Dec 2001 6:34 am
Data File: C:\HPCHEM\1\DATA\DEC1001\29235.D
Name: gc/ms scan 12/4
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	4.0	ug/l	1030330	ISTD01	11.67	5114630	20.0
2-Propanol, 1-(2-met	11.35	0.7	ug/l	183997	ISTD01	11.67	5114630	20.0
2-Propanol, 1-(2-met	11.43	1.1	ug/l	269635	ISTD01	11.67	5114630	20.0

29235.D GCMS1126.M Wed Dec 19 12:53:14 2001