User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/11/2001 Time: 08:18:14

Sample: AD22582
Analysis: $GCMS GC/MS Scan For Unknowns
Started: 10/11/01 08:17 Ended: 10/11/01 08:17 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D Vial: 10
Acq On : 28 Sep 2001 9:15 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Sep 28 22:04 2001 Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.79	152	487790	20.00	ug/l	-0.12
19) Naphthalene-d8	15.41	136	1909592	20.00	ug/l	-0.12
31) Acenaphthene-d10	20.69	164	897022	20.00	ug/l	-0.13
49) Phenanthrene-d10	25.12	188	1265974	20.00	ug/l	-0.12
59) Chrysene-d12	33.16	240	879862	20.00	ug/l	-0.12
68) Perylene-d12	37.28	264	653708	20.00	ug/l	-0.14
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
5) Phenol-d5	0.00	99	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
6) 2-Chlorophenol-d4	11.79	132	178	0.00	ug/l	0.38
Spiked Amount 75.000			Recovery =	0.00%		
7) 1,2-Dichlorobenzene-d4	12.30	152	5230	0.22	ug/l	-0.13
Spiked Amount 50.000			Recovery =	0.44%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.83	172	1862	0.03	ug/l	0.02
Spiked Amount 50.000			Recovery =	0.06%		
33) 2,4,6-Tribromophenol	0.00	330	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
62) Terphenyl-d14	0.00	244	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
Target Compounds						Qvalue
2) Pyridine	5.20	79	1045	N.D.		
3) N-nitroso-dimethyl amine	0.00	74	0	N.D.		
8) Phenol	0.00	94	0	N.D.		
9) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
10) 2-Chlorophenol	0.00	128	0	N.D.		
11) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
12) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
13) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
14) 2-Methylphenol	0.00	108	0	N.D.		
15) 2,2'-oxybis(1-Chloropropan	0.00	45	0	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		
17) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.		
18) Hexachloroethane	0.00	117	0	N.D.		
21) Nitrobenzene	0.00	77	0	N.D.		
22) Isophorone	0.00	82	0	N.D.		

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

Quantitation Report (QT Reviewed)

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Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
23) 2-Nitrophenol	0.00	139	0	N.D.	
24) 2,4-Dimethylphenol	0.00	107	0	N.D.	
25) bis(2-Chloroethoxy)methane	0.00	93	0	N.D.	
26) 2,4-Dichlorophenol	0.00	162	0	N.D.	
27) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
28) Naphthalene	0.00	128	0	N.D.	
29) Hexachlorobutadiene	0.00	225	0	N.D.	
30) 4-Chloro-3-methylphenol	0.00	107	0	N.D.	
34) Hexachlorocyclopentadiene	0.00	237	0	N.D.	
35) 2,4,6-Trichlorophenol	0.00	196	0	N.D.	
36) 2,4,5-Trichlorophenol	0.00	196	0	N.D.	
37) 2-Chloronaphthalene	0.00	162	0	N.D.	
38) Dimethylphthalate	0.00	163	0	N.D.	
39) 2,6-Dinitrotoluene	0.00	165	0	N.D.	
40) Acenaphthylene	0.00	152	0	N.D.	
41) Acenaphthene	0.00	153	0	N.D.	
42) 2,4-Dinitrophenol	0.00	184	0	N.D.	
43) 4-Nitrophenol	21.52	65	270	N.D.	
44) 2,4-Dinitrotoluene	20.94	165	105	N.D.	
45) Diethylphthalate	22.19	149	183	N.D.	
46) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.	
47) Fluorene	0.00	166	0	N.D.	
48) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.	
50) 4,6-Dinitro-2-methylphenol	0.00	198	0	N.D.	
51) N-Nitrosodiphenylamine	0.00	168	0	N.D.	
52) 4-Bromophenyl-phenylether	0.00	248	0	N.D.	
53) Hexachlorobenzene	0.00	284	0	N.D.	
54) Pentachlorophenol	0.00	266	0	N.D.	
55) Phenanthrene	25.12	178	312	N.D.	
56) Anthracene	25.12	178	312	N.D.	
57) Di-n-butylphthalate	27.11	149	2538	N.D.	
58) Fluoranthene	0.00	202	0	N.D.	
60) 3,3'-Dichlorobenzidine	0.00	252	0	N.D.	
61) Benzidine	0.00	184	0	N.D.	
63) Pyrene	0.00	202	0	N.D.	
64) Butylbenzylphthalate	0.00	149	0	N.D.	
65) Benzo(a)anthracene	33.17	228	2208	N.D.	
66) Bis(2-ethylhexyl)phthalate	33.42	149	4088	N.D.	
67) Chrysene	33.17	228	2208	N.D.	
69) Di-n-Octylphthalate	35.16	149	1434	N.D.	
70) Benzo(b)fluoranthene	0.00	252	0	N.D.	

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

Quantitation Report (QT Reviewed)

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Acq On : 28 Sep 2001 9:15 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Sep 28 22:04 2001 Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
72)	Benzo(a)pyrene	37.10	252	202	N.D.	
73)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
74)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
75)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

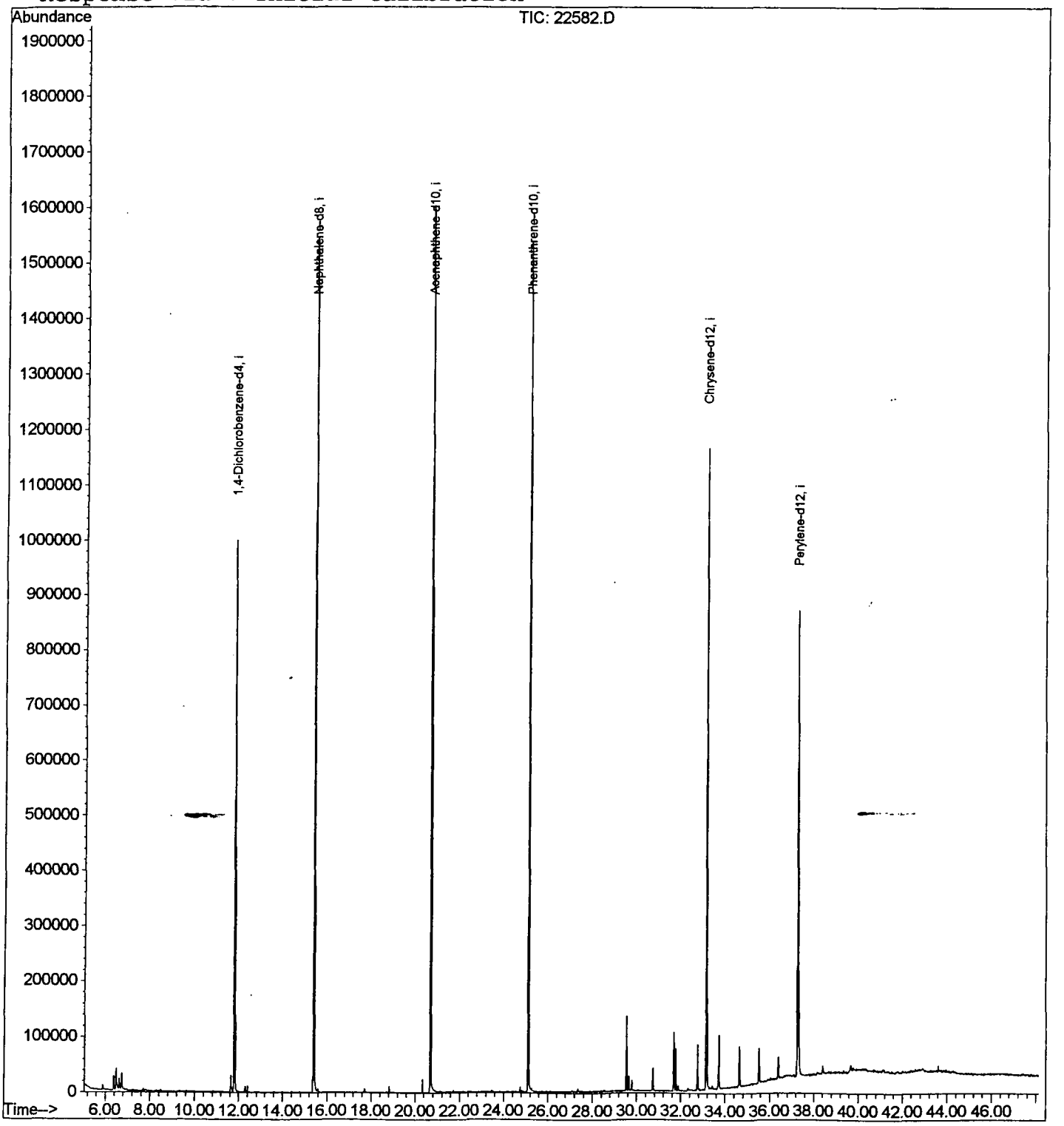
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
Quant Time: Sep 28 22:04 2001

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: NP091101.RES

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



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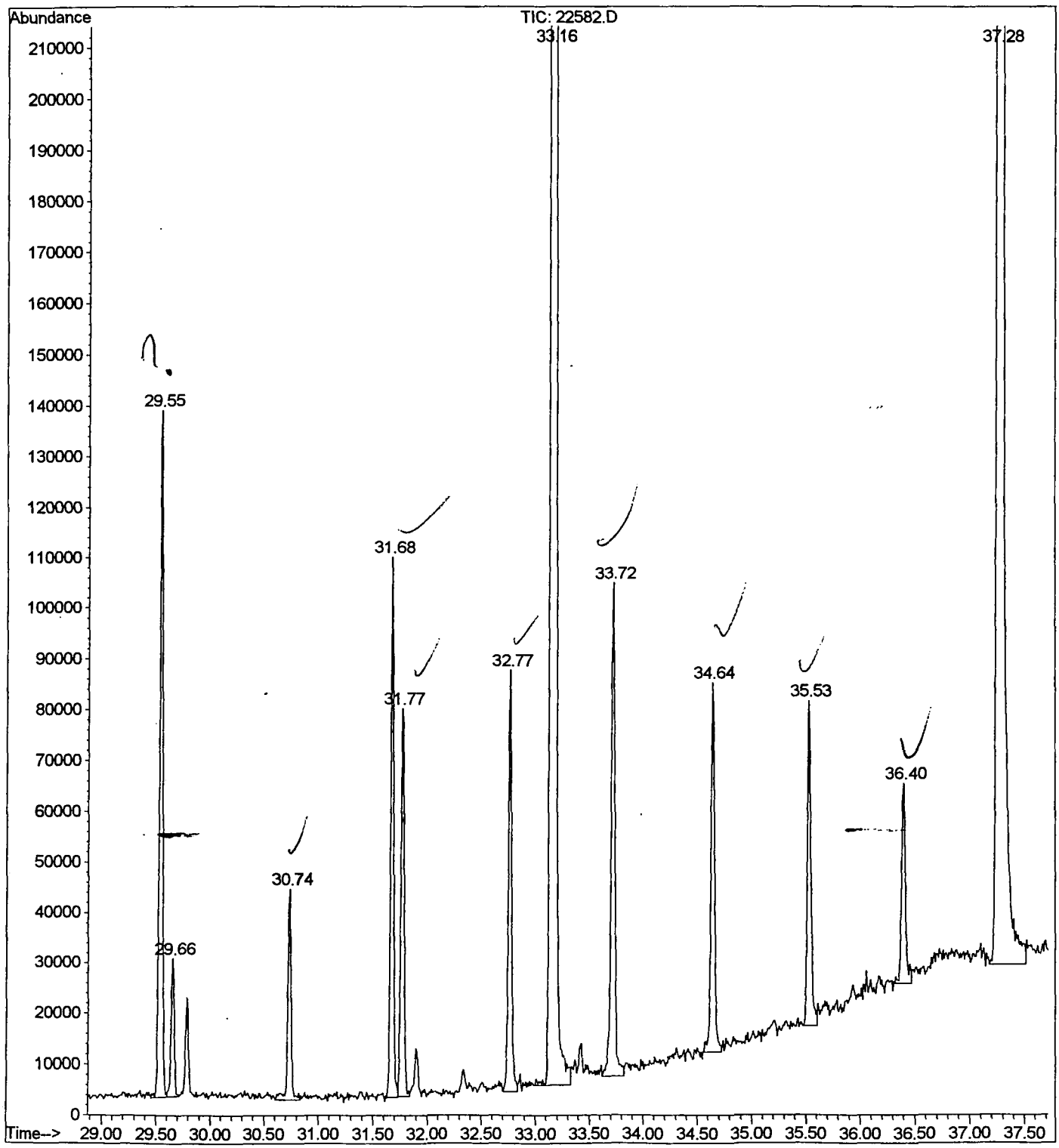
Operator : 209 GALOS
Acquired : 28 Sep 2001 9:15 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 10



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D Vial: 10
Acq On : 28 Sep 2001 9:15 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
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

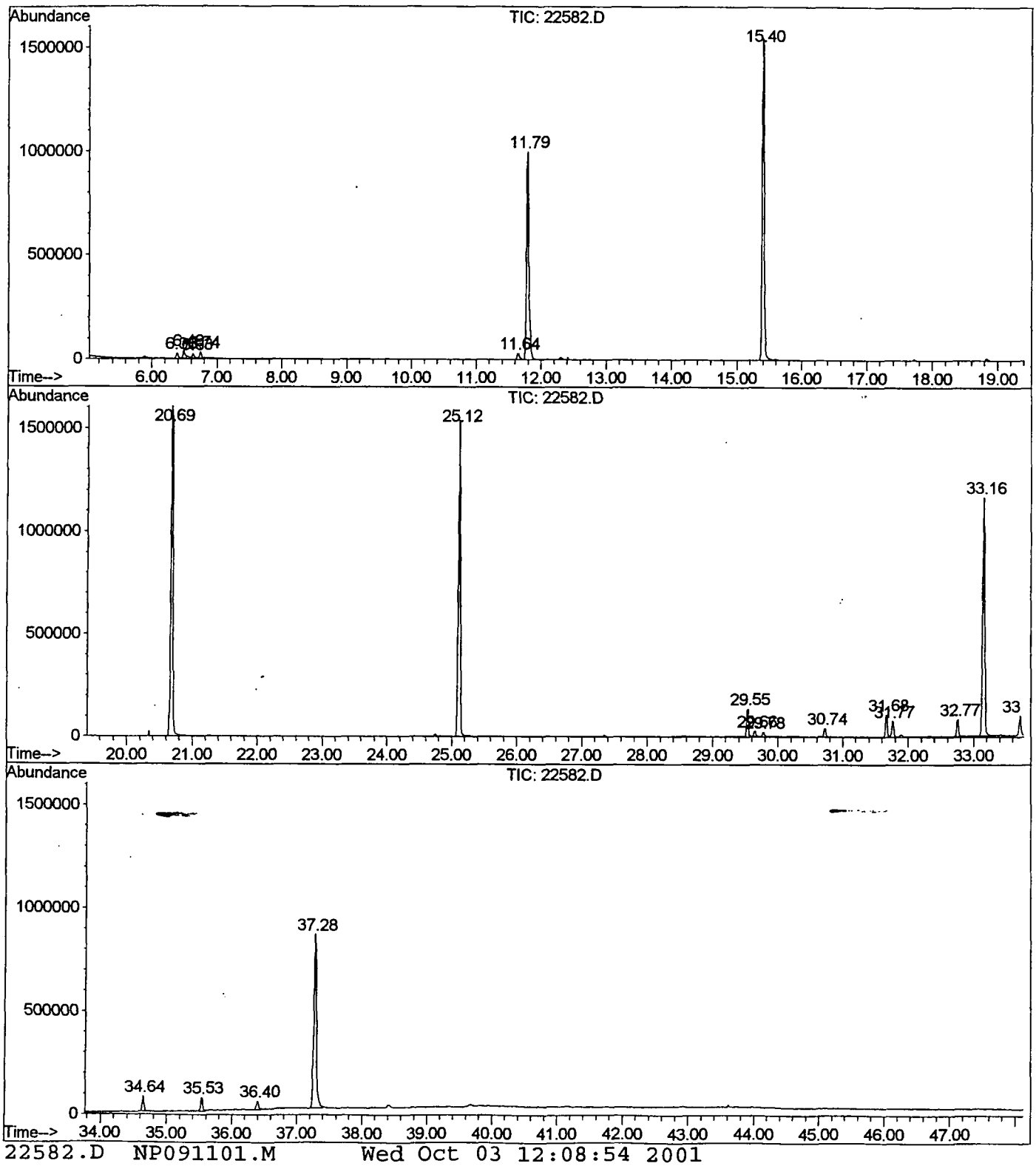
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.384	141	146	152	rBV	26124	53564	1.50%	0.269%
2	6.494	154	158	162	rBV	38504	82023	2.30%	0.412%
3	6.631	169	173	179	rVB	18917	38620	1.08%	0.194%
4	6.742	181	185	192	rVB	28198	55422	1.55%	0.278%
5	11.644	713	719	729	rBV	29467	70536	1.98%	0.354%
6	11.791	729	735	753	rBV	999333	2529919	70.86%	12.707%
7	15.399	1122	1128	1144	rBV	1547735	3483189	97.56%	17.495%
8	20.686	1697	1704	1719	rBV	1604417	3570238	100.00%	17.932%
9	25.121	2180	2187	2200	rBV2	1539286	3267136	91.51%	16.409%
10	29.545	2663	2669	2674	rBV	135360	268733	7.53%	1.350%
11	29.656	2677	2681	2686	rVB	26901	53528	1.50%	0.269%
12	29.784	2691	2695	2703	rVB	19392	39557	1.11%	0.199%
13	30.739	2794	2799	2804	rVB	40886	82051	2.30%	0.412%
14	31.675	2896	2901	2907	rBV	106704	208581	5.84%	1.048%
15	31.767	2907	2911	2917	rVV	76479	162025	4.54%	0.814%
16	32.768	3013	3020	3027	rBV	82989	174249	4.88%	0.875%
17	33.162	3056	3063	3076	rBV2	1163038	2752481	77.10%	13.824%
18	33.722	3113	3124	3130	rBV	97255	215720	6.04%	1.083%
19	34.641	3220	3224	3234	rVB	71822	147673	4.14%	0.742%
20	35.531	3316	3321	3326	rBV	63927	138906	3.89%	0.698%
21	36.403	3411	3416	3421	rBV	38985	89450	2.51%	0.449%
22	37.284	3504	3512	3524	rBV2	842515	2426587	67.97%	12.188%

Sum of corrected areas: 19910188

22582.D NP091101.M Wed Oct 03 12:08:52 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Operator : 209 GALOS
Acquired : 28 Sep 2001 9:15 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 10
Quant File :NP091101.RES (RTE Integrator)



22582.D NP091101.M Wed Oct 03 12:08:54 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

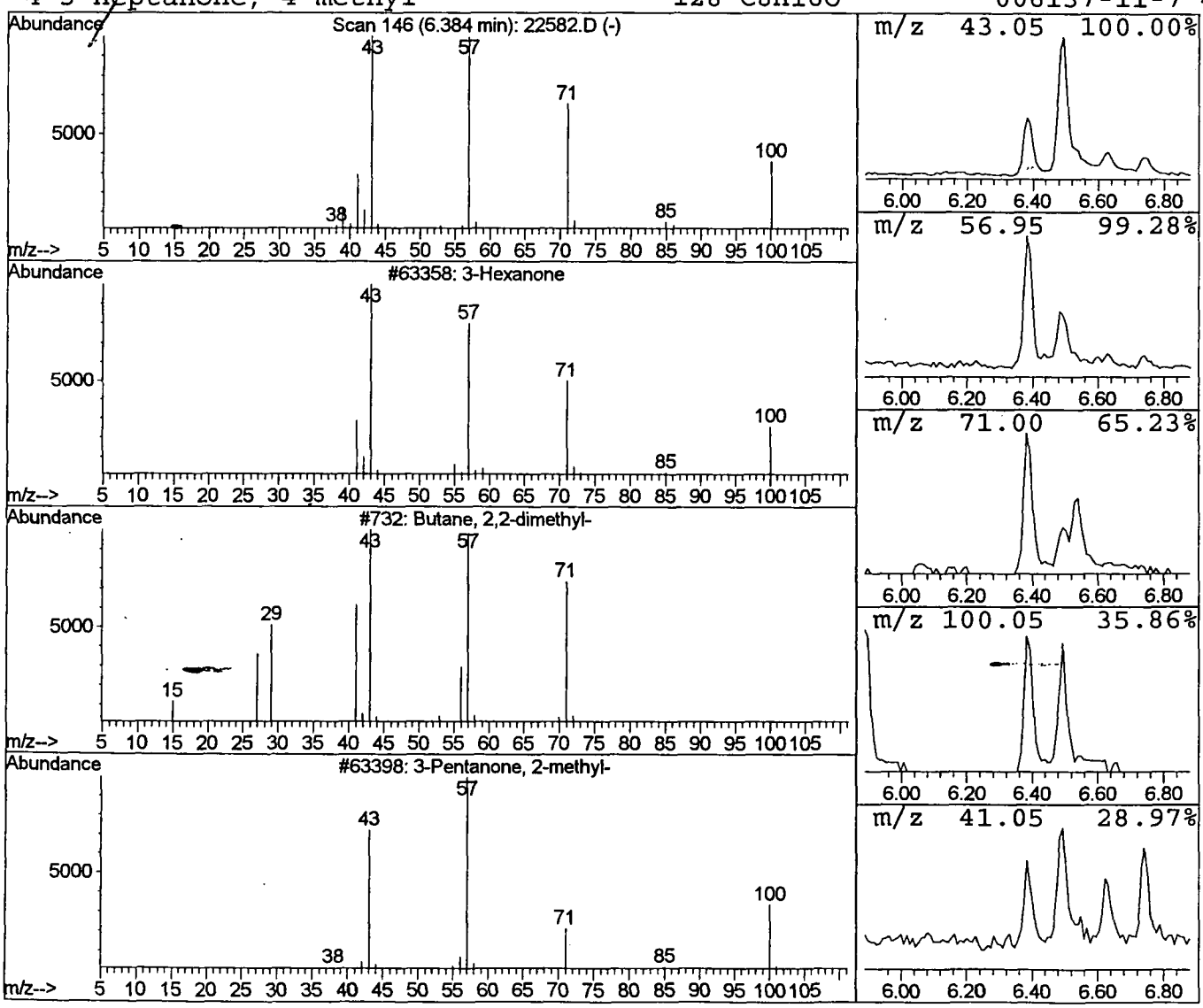
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexanone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	0.42 ug/l	53564	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	53
3	3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	53
4	3-Heptanone, 4-methyl-	128	C8H16O	006137-11-7	45



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

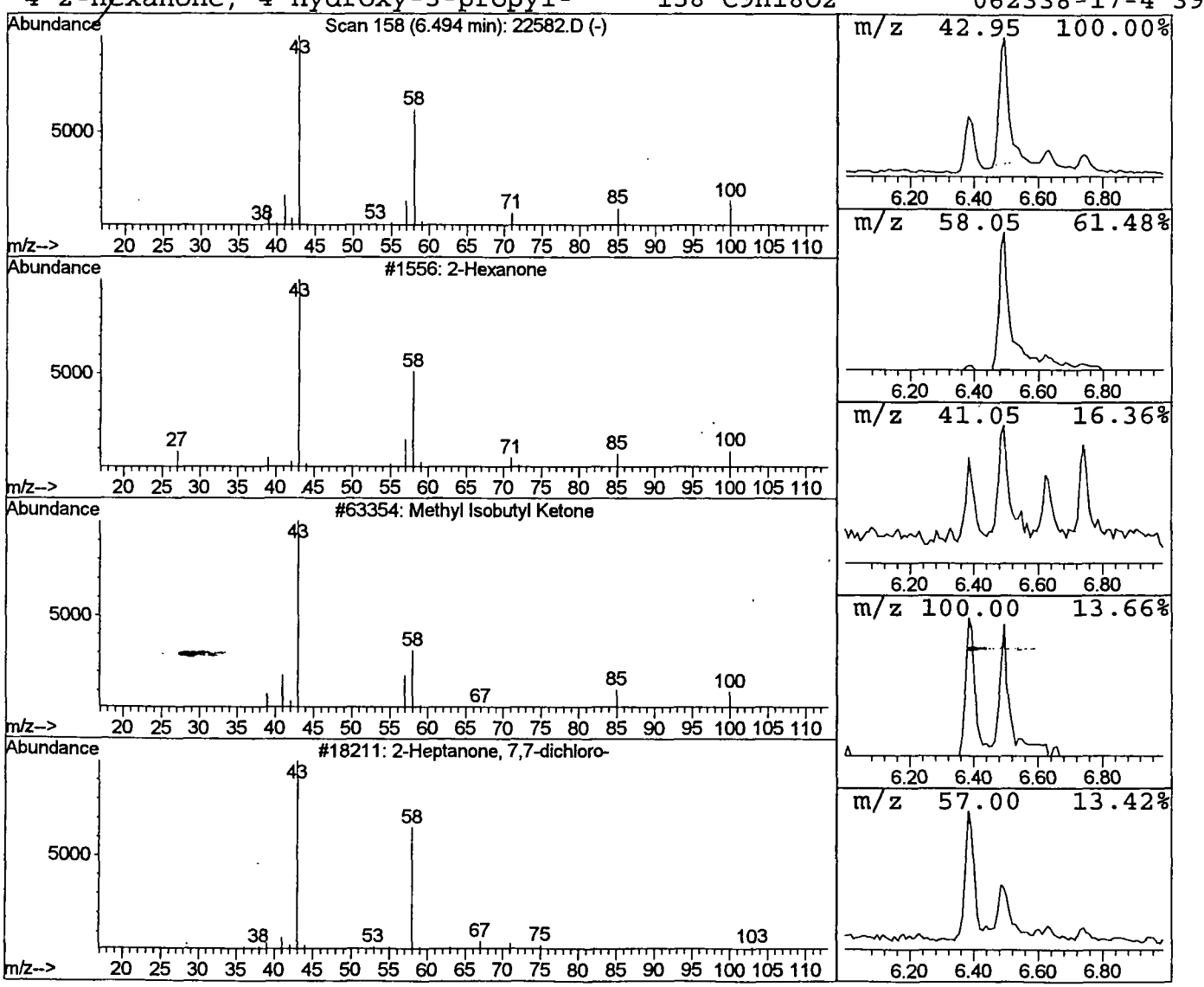
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	0.65 ug/l	82023	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58
3	2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	000000-00-0	40
4	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	39



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

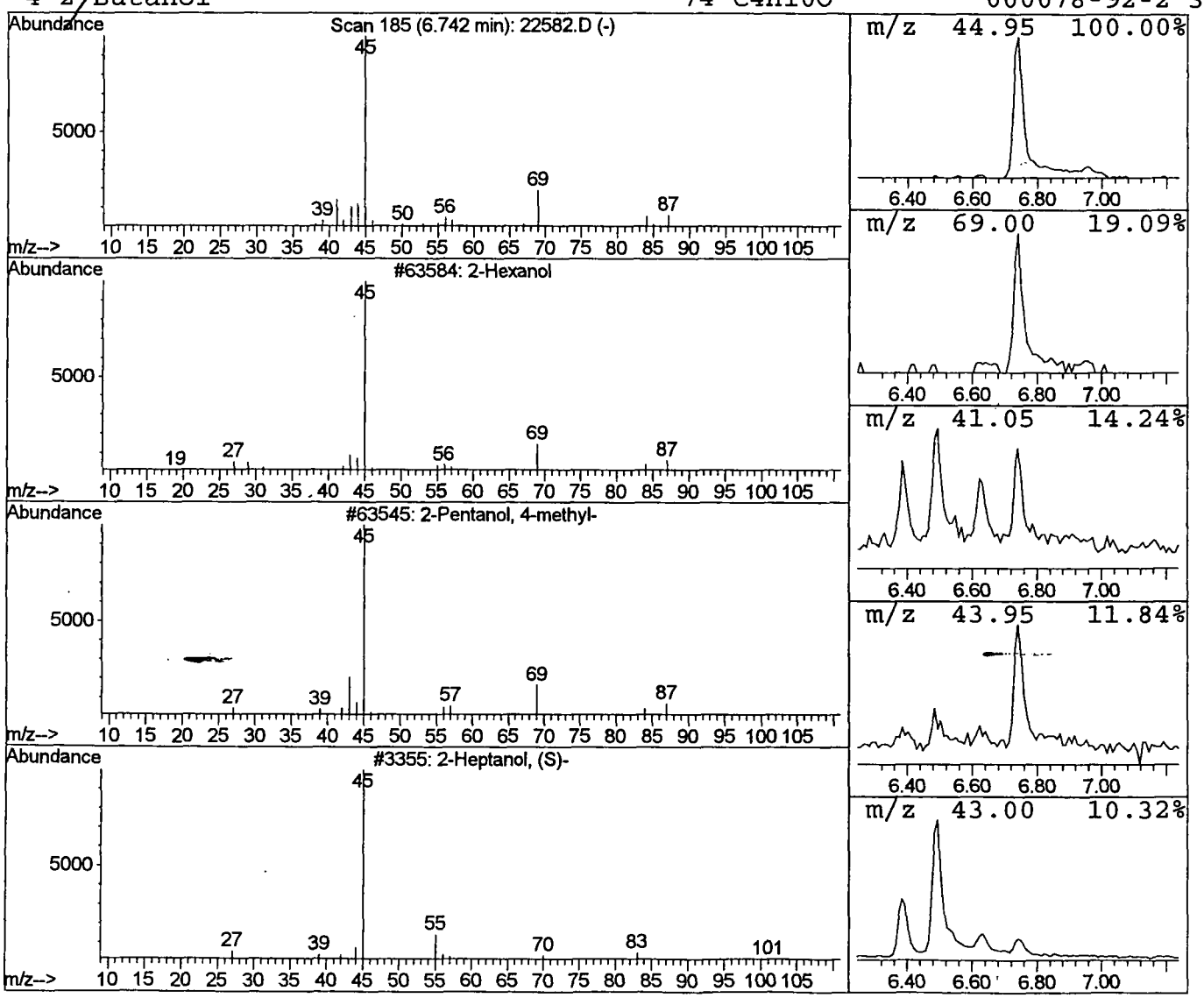
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 2-Hexanol — Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	0.44 ug/l	55422	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	74
3	2-Heptanol, (S)-	116	C7H16O	006033-23-4	40
4	2-Butanol	74	C4H10O	000078-92-2	39



Library Search Compound Report

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Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

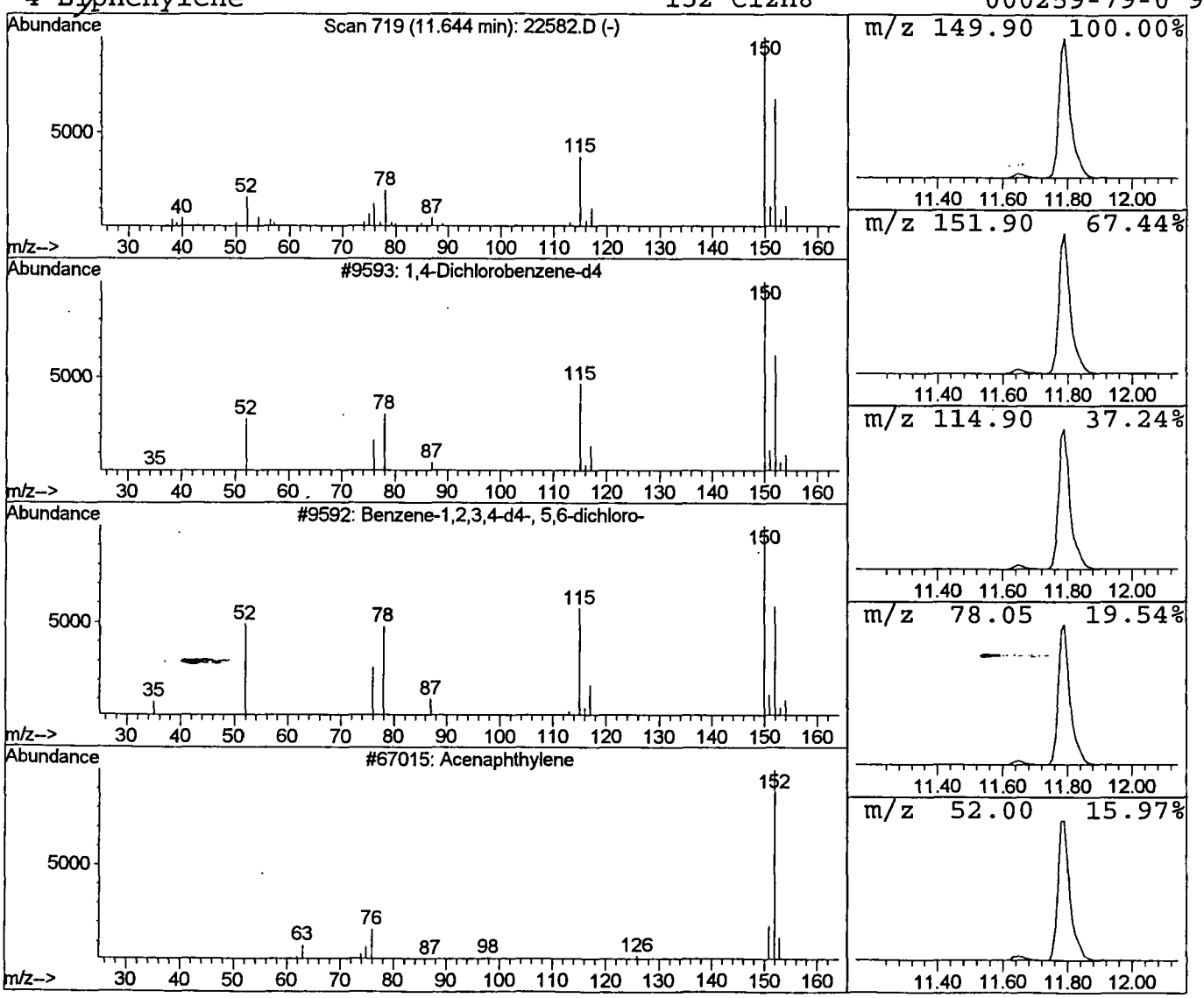
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 1,4-Dichlorobenzene-d4 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	0.56 ug/l	70536	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dichlorobenzene-d4	150	C6Cl2D4	003855-82-1	53
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	25
3	Acenaphthylene	152	C12H8	000208-96-8	10
4	Biphenylene	152	C12H8	000259-79-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

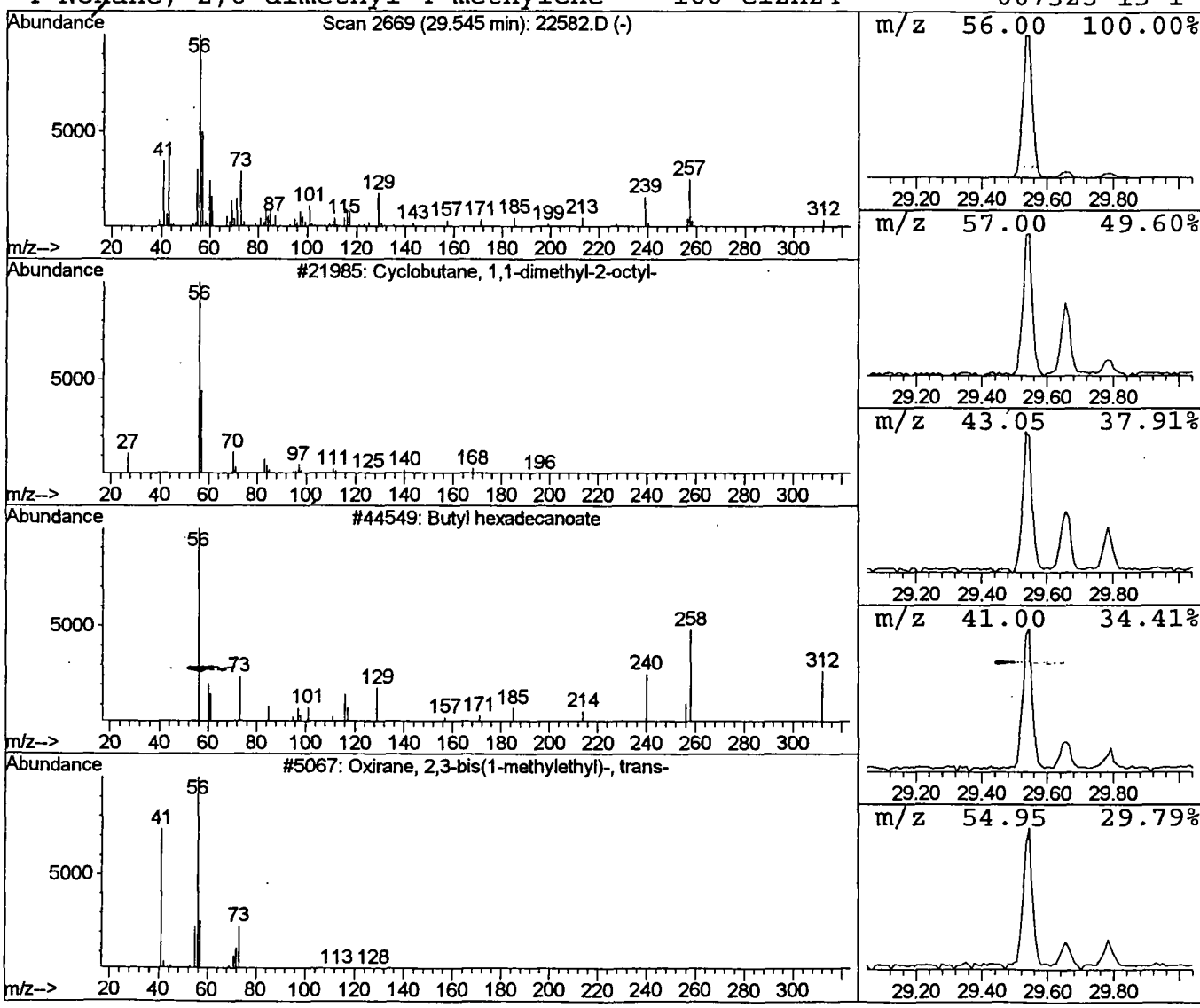
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Cyclobutane, 1,1-dimethyl-2-oc Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.55	1.95 ug/l	268733	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	27
2	Butyl hexadecanoate	312	C20H40O2	000000-00-0	25
3	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	23
4	Nonane, 2,8-dimethyl-4-methylene-	168	C12H24	007323-15-1	16



Library Search Compound Report

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Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

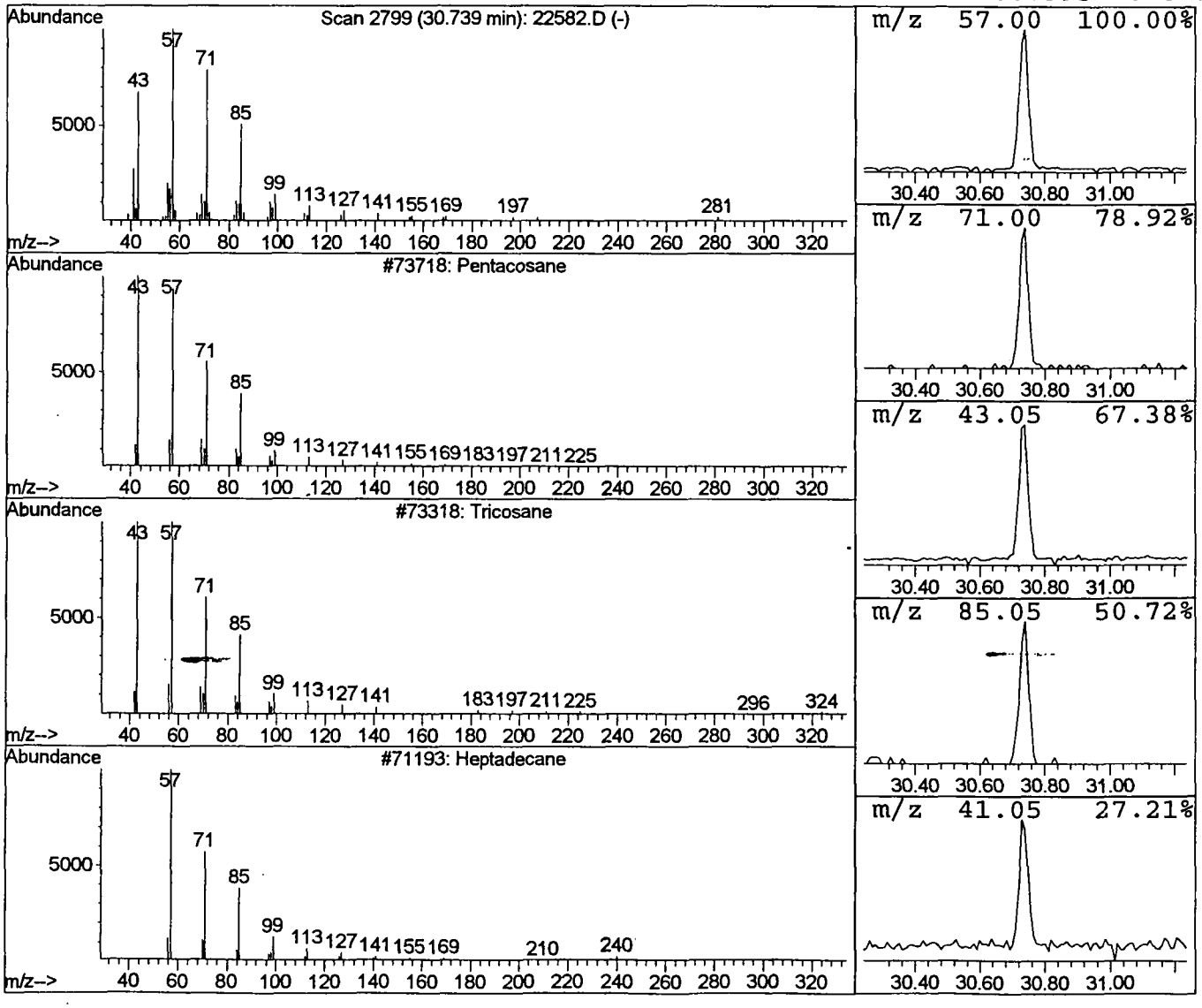
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Pentacosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.74	0.60 ug/l	82051	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Tricosane	324	C23H48	000638-67-5	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Octadecane	254	C18H38	000593-45-3	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

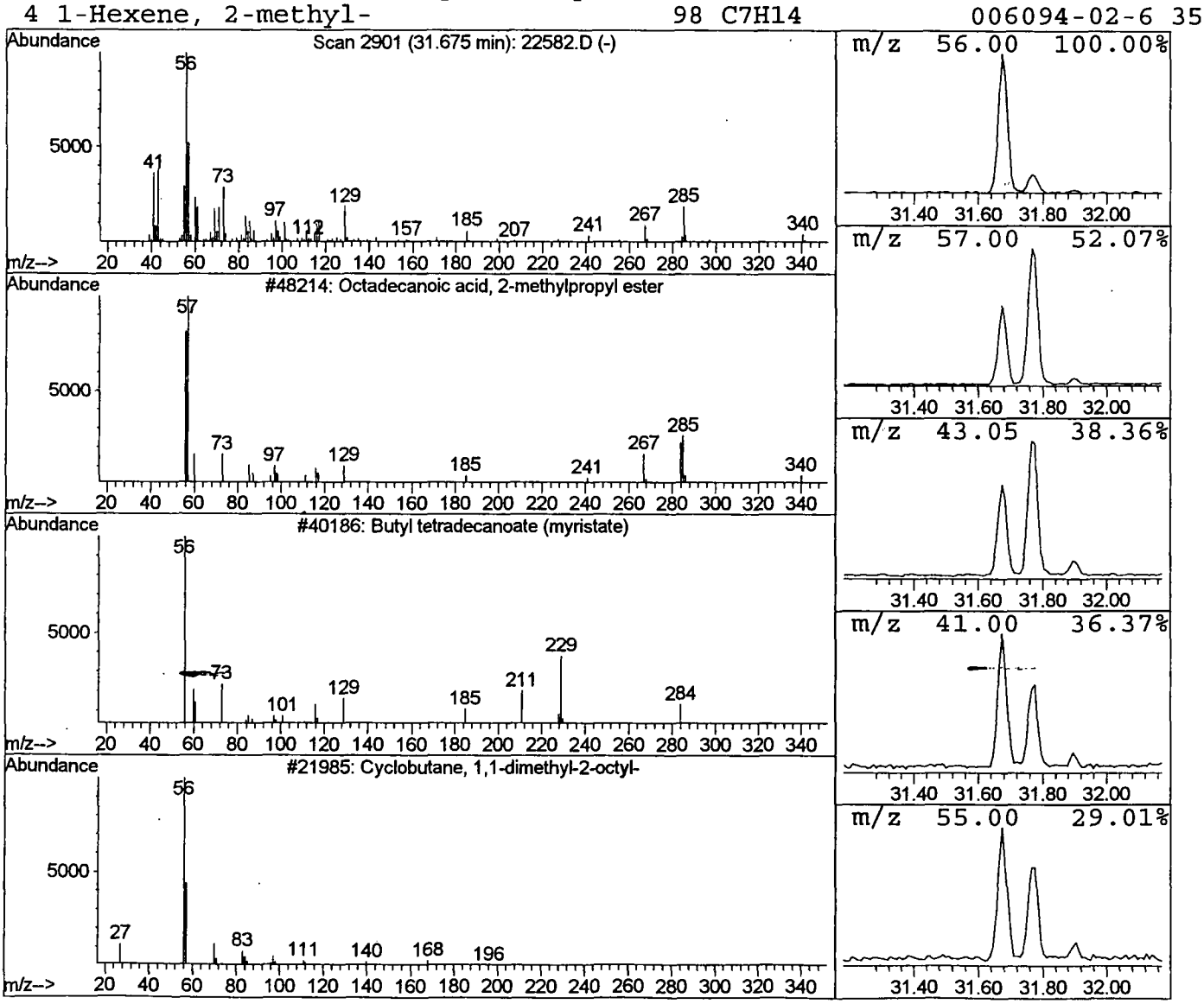
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Octadecanoic acid, 2-methylpro Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.68	1.52 ug/l	208581	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, 2-methylpropyl e	340	C22H44O2	000646-13-9	84
2	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	72
3	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	35



Library Search Compound Report

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Acq On : 28 Sep 2001 9:15 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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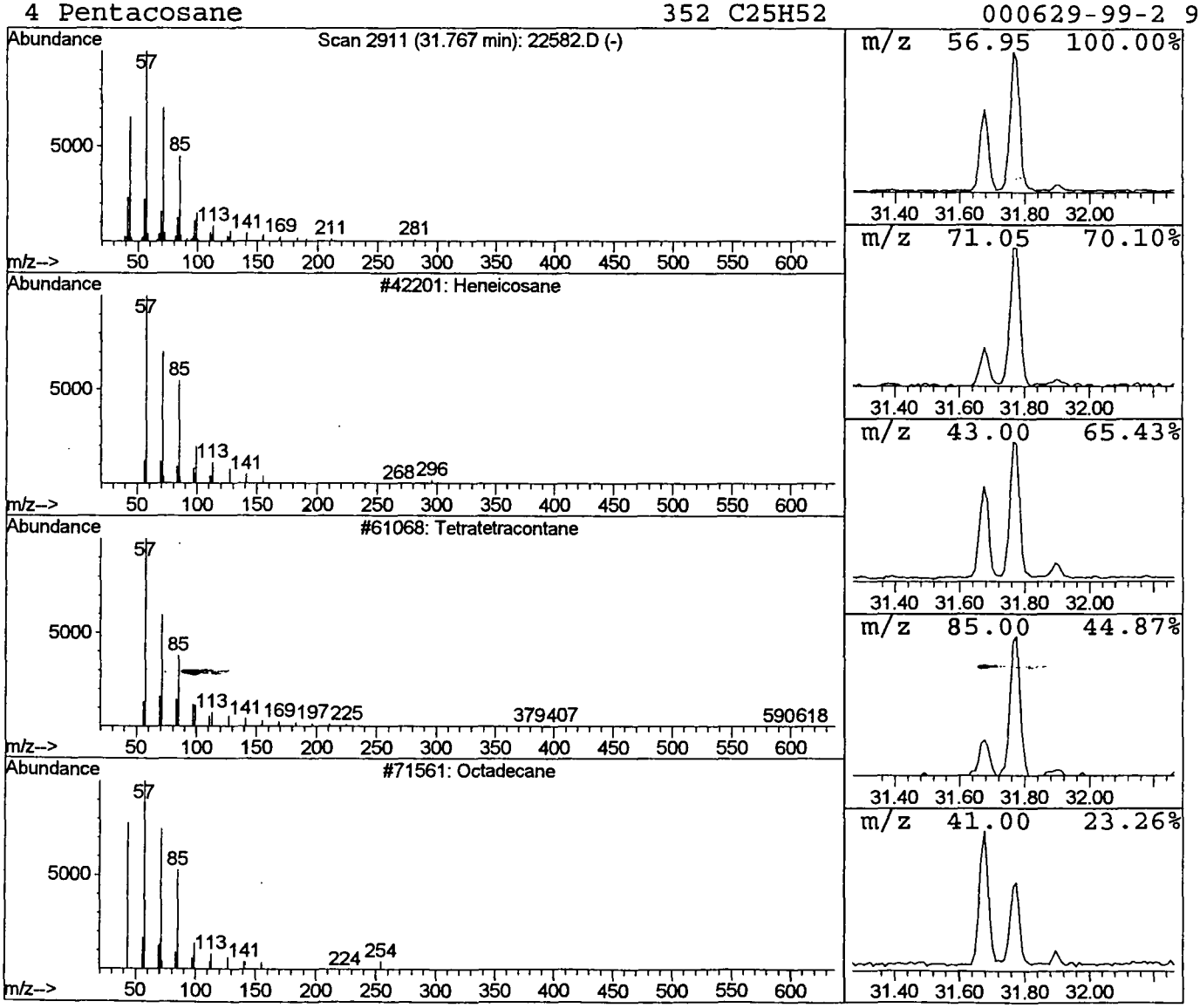
Peak Number 8 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.77	1.18 ug/l	162025	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Octadecane	254	C18H38	000593-45-3	91
4	Pentacosane	352	C25H52	000629-99-2	91



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Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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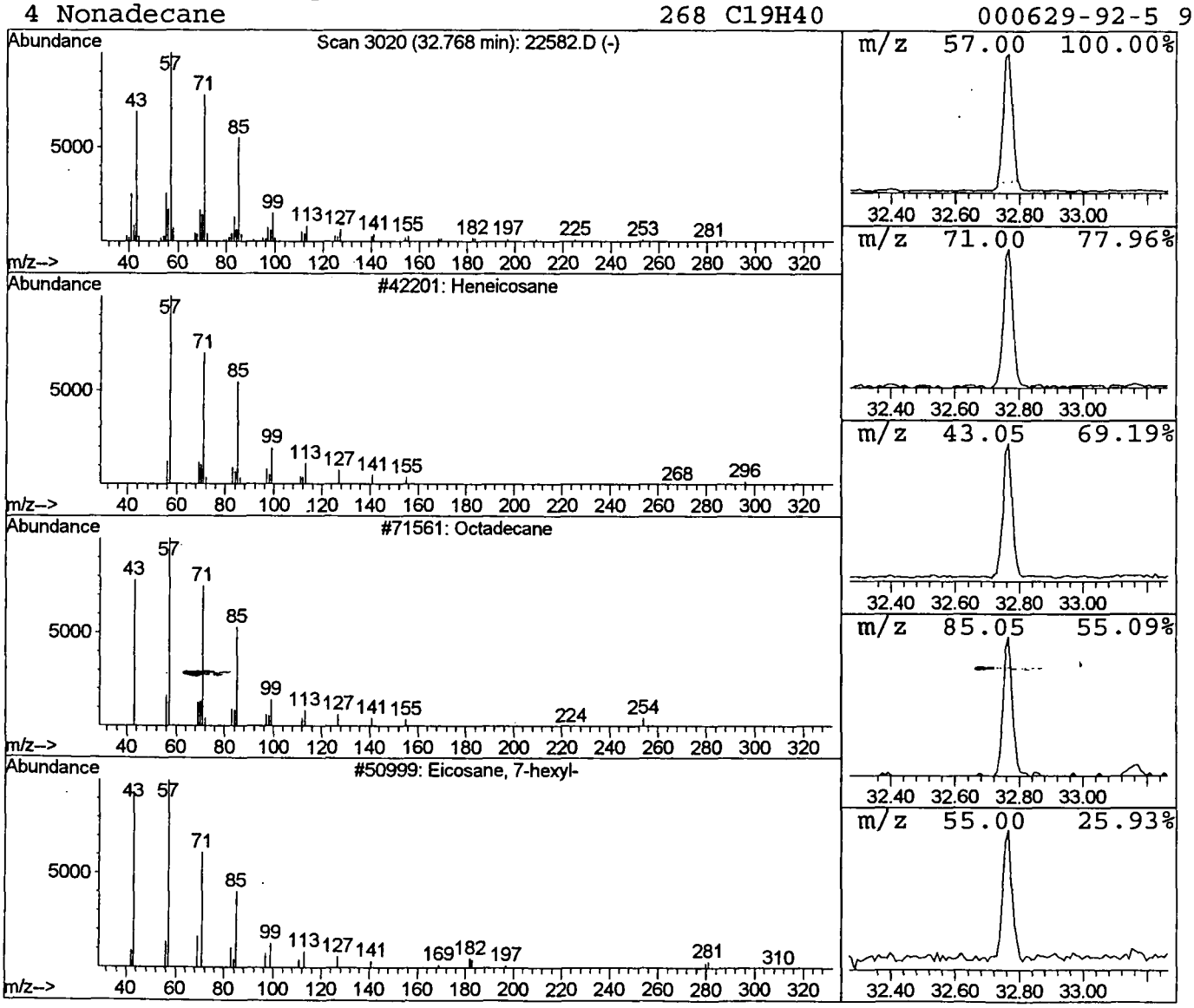
Peak Number 9 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.77	1.27 ug/l	174249	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Octadecane	254	C18H38	000593-45-3	91
3	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4	Nonadecane	268	C19H40	000629-92-5	91



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Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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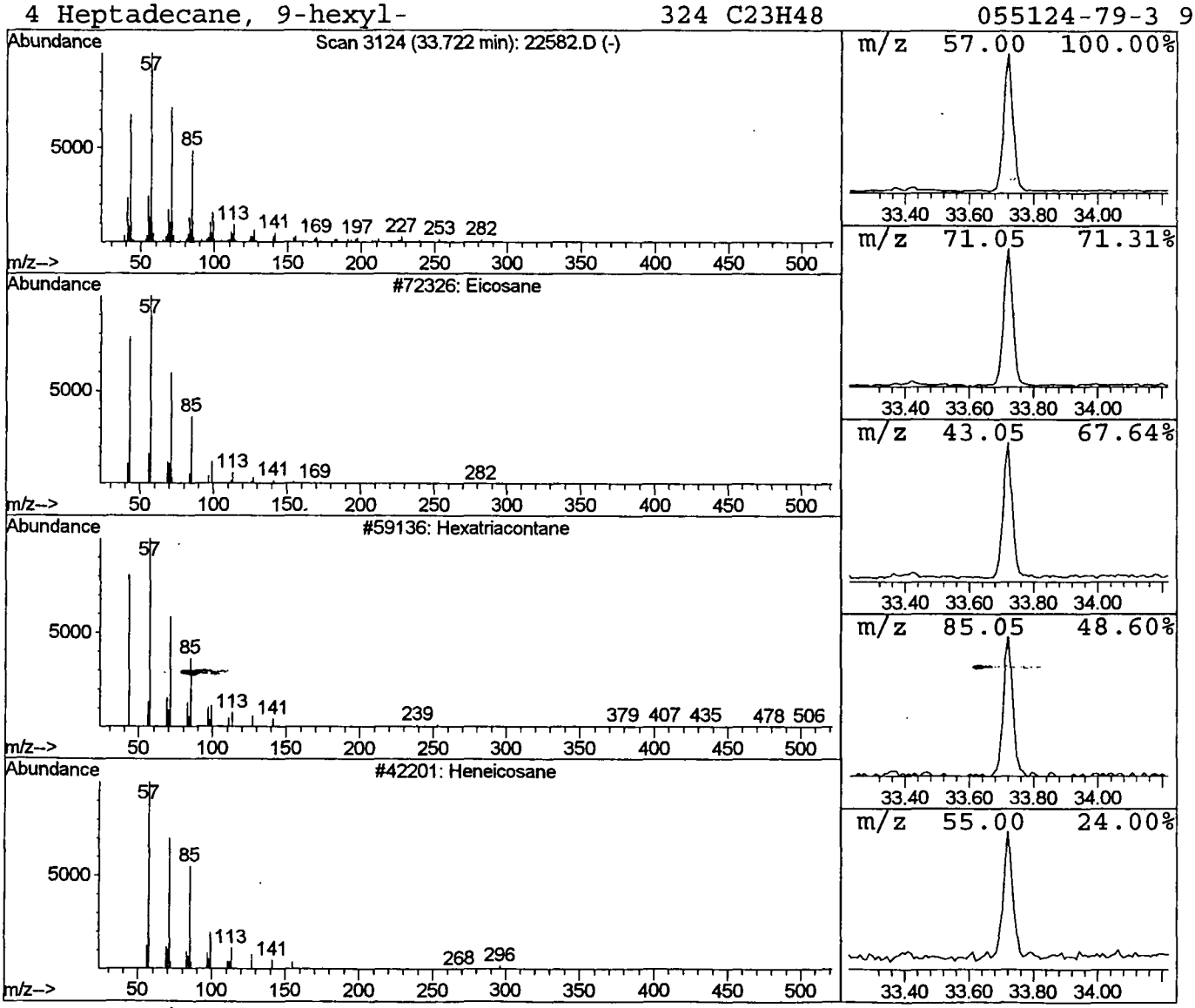
Peak Number 10 Eicosane — Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.72	1.57 ug/l	215720	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91



Library Search Compound Report

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Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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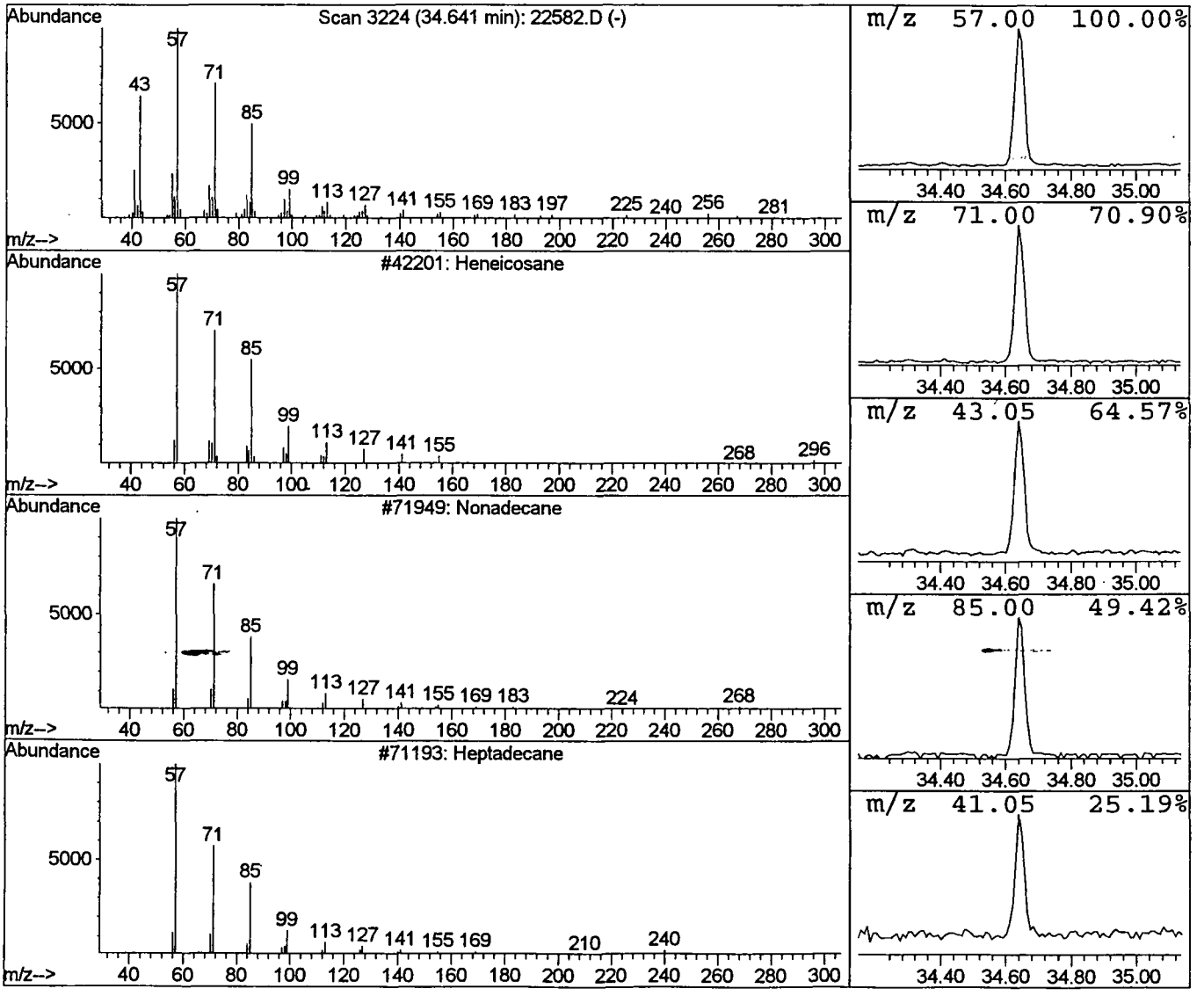
Peak Number 11 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.64	1.07 ug/l	147673	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Nonadecane	268	C19H40	000629-92-5	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Pentacosane	352	C25H52	000629-99-2	91



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Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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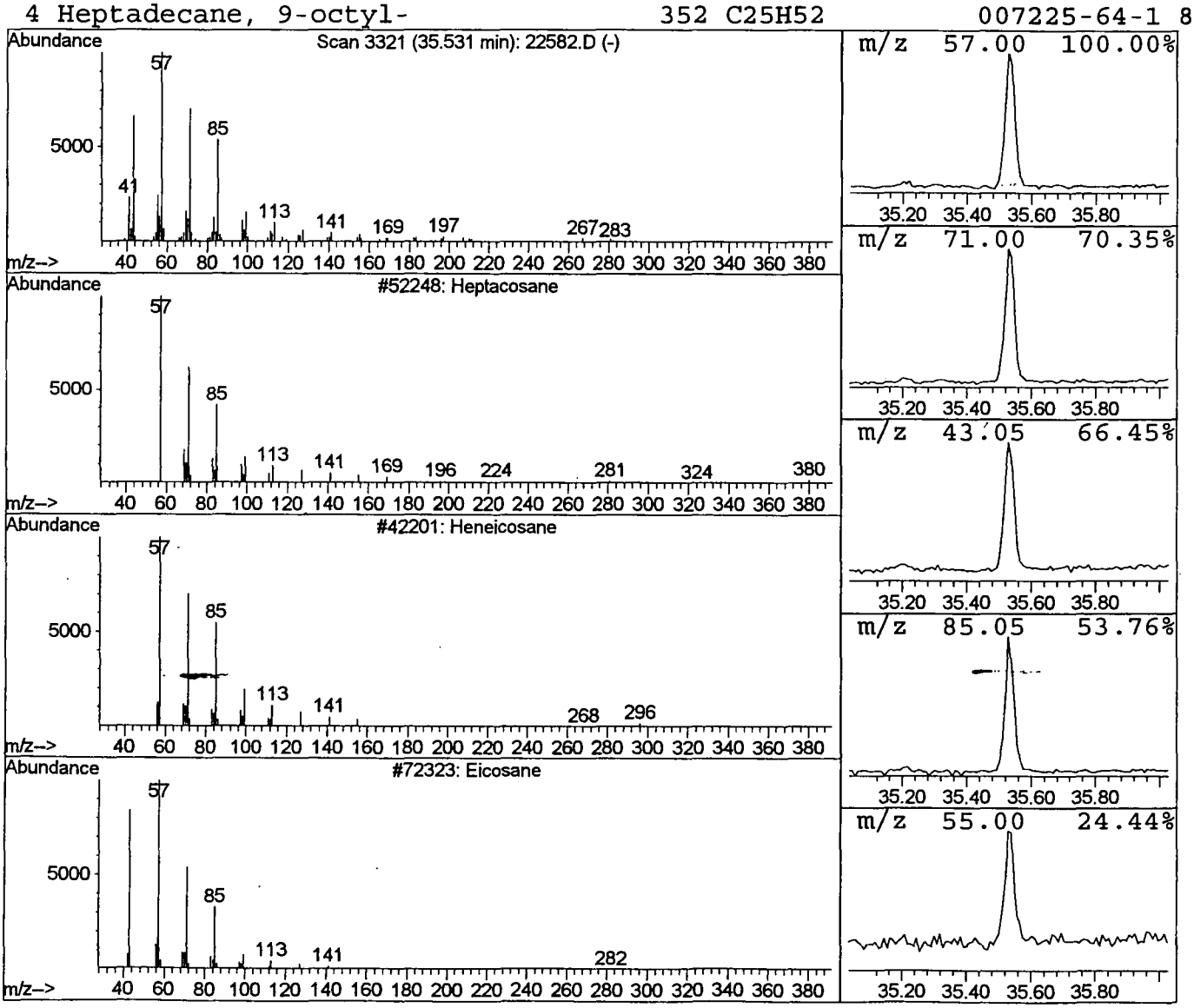
Peak Number 12 Heptacosane — Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.53	1.14 ug/l	138906	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Eicosane	282	C20H42	000112-95-8	89
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



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Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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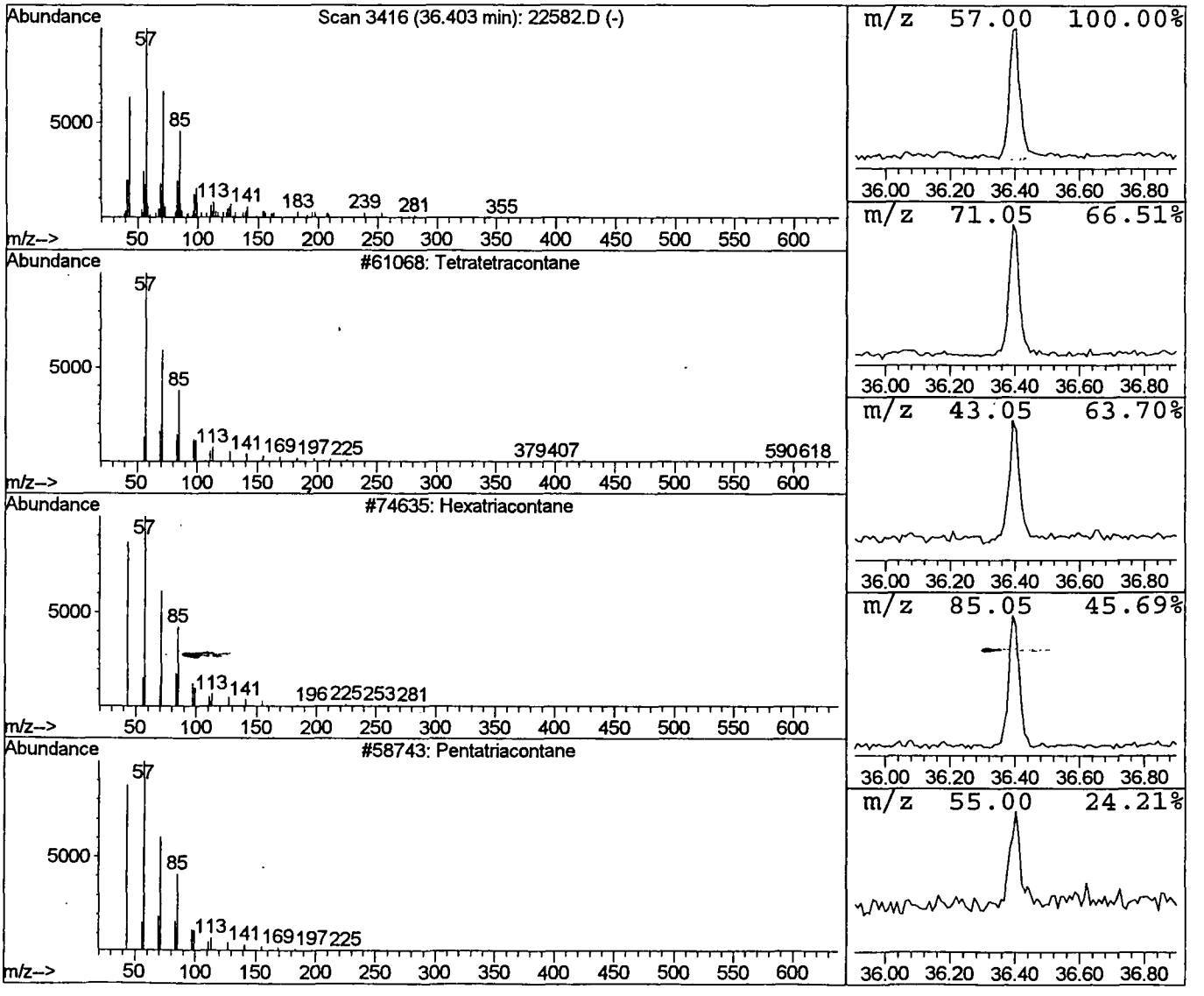
Peak Number 13 Tetratetracontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.40	0.74 ug/l	89450	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Pentatriacontane	493	C35H72	000630-07-9	91
4	Dotriacontane	451	C32H66	000544-85-4	91



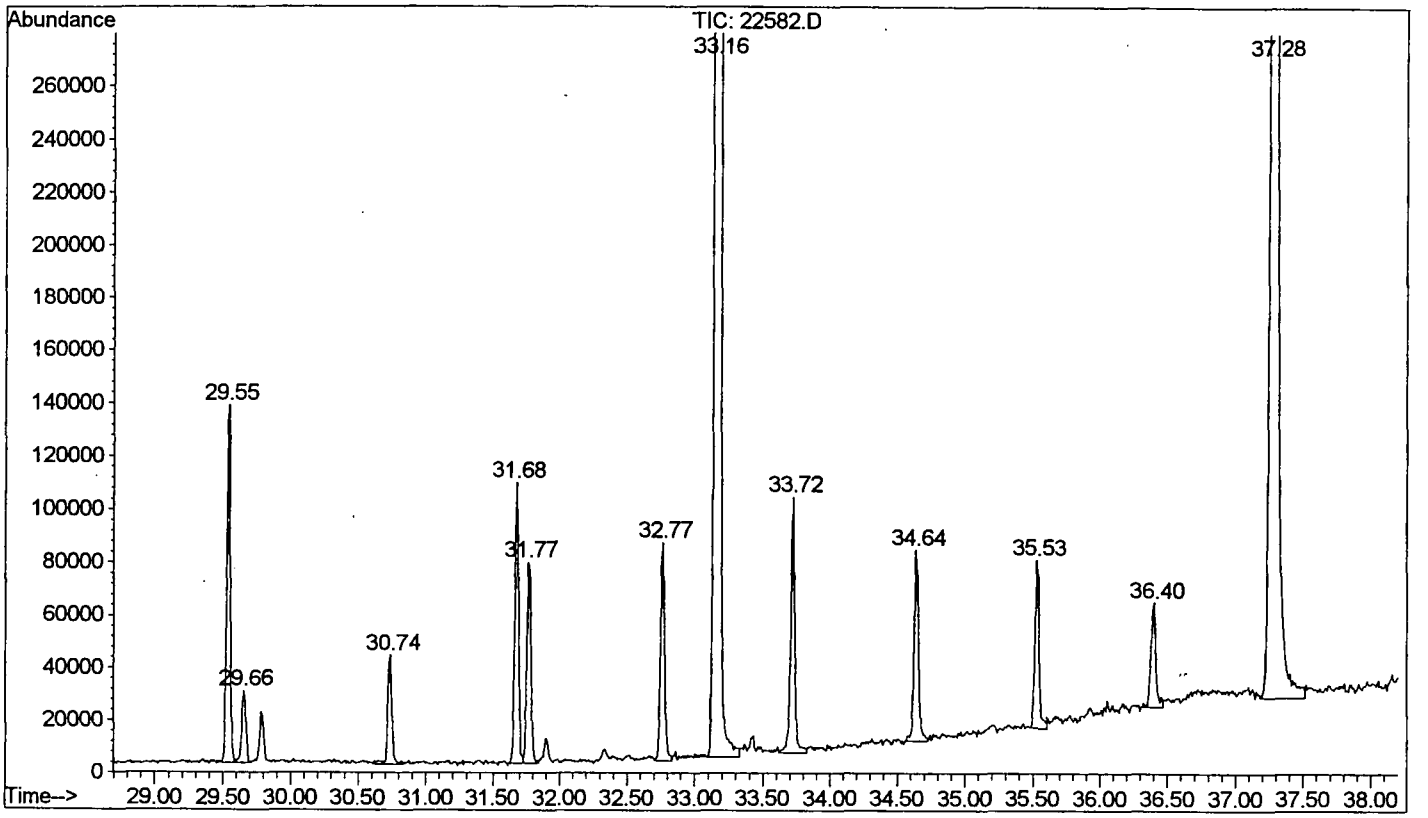
Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 28 Sep 2001 9:15 pm
Data File: C:\HPCHEM\1\DATA\SEP2801\22582.D
Name: gc/ms unknown
Misc:
Method: C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title: 209 625/TCLP calibration table 7/16/01
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
3-Hexanone	6.38	0.4	ug/l	53564	ISTD01	11.79	2529920	20.0
2-Hexanone	6.49	0.6	ug/l	82023	ISTD01	11.79	2529920	20.0
2-Hexanol	6.74	0.4	ug/l	55422	ISTD01	11.79	2529920	20.0
1,4-Dichlorobenzene-	11.64	0.6	ug/l	70536	ISTD01	11.79	2529920	20.0
Cyclobutane, 1,1-dim	29.55	2.0	ug/l	268733	ISTD05	33.16	2752480	20.0
Pentacosane	30.74	0.6	ug/l	82051	ISTD05	33.16	2752480	20.0
Octadecanoic acid, 2	31.68	1.5	ug/l	208581	ISTD05	33.16	2752480	20.0
Heneicosane	31.77	1.2	ug/l	162025	ISTD05	33.16	2752480	20.0
Heneicosane	32.77	1.3	ug/l	174249	ISTD05	33.16	2752480	20.0
Eicosane	33.72	1.6	ug/l	215720	ISTD05	33.16	2752480	20.0
Heneicosane	34.64	1.1	ug/l	147673	ISTD05	33.16	2752480	20.0
Heptacosane	35.53	1.1	ug/l	138906	ISTD06	37.28	2426590	20.0
Tetratetracontane	36.40	0.7	ug/l	89450	ISTD06	37.28	2426590	20.0

22582.D NP091101.M Wed Oct 03 12:09:25 2001

File : C:\HPCHEM\1\DATA\SEP2801\22582.D
Operator : 209 GALOS
Acquired : 28 Sep 2001 9:15 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 10



Comments for Sample AD22582 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-12-2001 Time: 13:47:26

A scan of the extract prepared for the 508 analysis, from 35 - 550 amu using a mass spectrometer, does not reveal the presence of any non-target compounds generally different than those in the Laboratory Blank, except that the levels are higher in the sample than in the Laboratory Blank.