User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 12:28:24

Sample: AD22919
Analysis: $GCMS GC/MS Scan For Unknowns
Started: 10/5/01 12:27 Ended: 10/5/01 12:27 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D Vial: 9
Acq On : 2 Oct 2001 9:56 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 4 9:54 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	450578	20.00	ug/l	-0.12
19) Naphthalene-d8	15.40	136	1746363	20.00	ug/l	-0.13
31) Acenaphthene-d10	20.69	164	812866	20.00	ug/l	-0.13
49) Phenanthrene-d10	25.11	188	1138444	20.00	ug/l	-0.13
59) Chrysene-d12	33.15	240	817411	20.00	ug/l	-0.13
68) Perylene-d12	37.27	264	613822	20.00	ug/l	-0.15
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	0.00	132	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
7) 1,2-Dichlorobenzene-d4	12.31	152	4531	0.21	ug/l	-0.13
Spiked Amount 50.000			Recovery =	0.42%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.82	172	1833	0.03	ug/l	0.01
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	0.00	79	0	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

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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. 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	21.12	65	197	N.D.	
44)	2,4-Dinitrotoluene	0.00	165	0	N.D.	
45)	Diethylphthalate	22.19	149	316	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	25.10	178	272	N.D.	
56)	Anthracene	25.10	178	272	N.D.	
57)	Di-n-butylphthalate	27.10	149	50302	0.55 ug/l	99
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	33.15	228	1785	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.41	149	5468	N.D.	
67)	Chrysene	33.15	228	1785	N.D.	
69)	Di-n-Octylphthalate	35.21	149	175	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 4 9:54 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	37.28	252	2379	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

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Acq On : 2 Oct 2001 9:56 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
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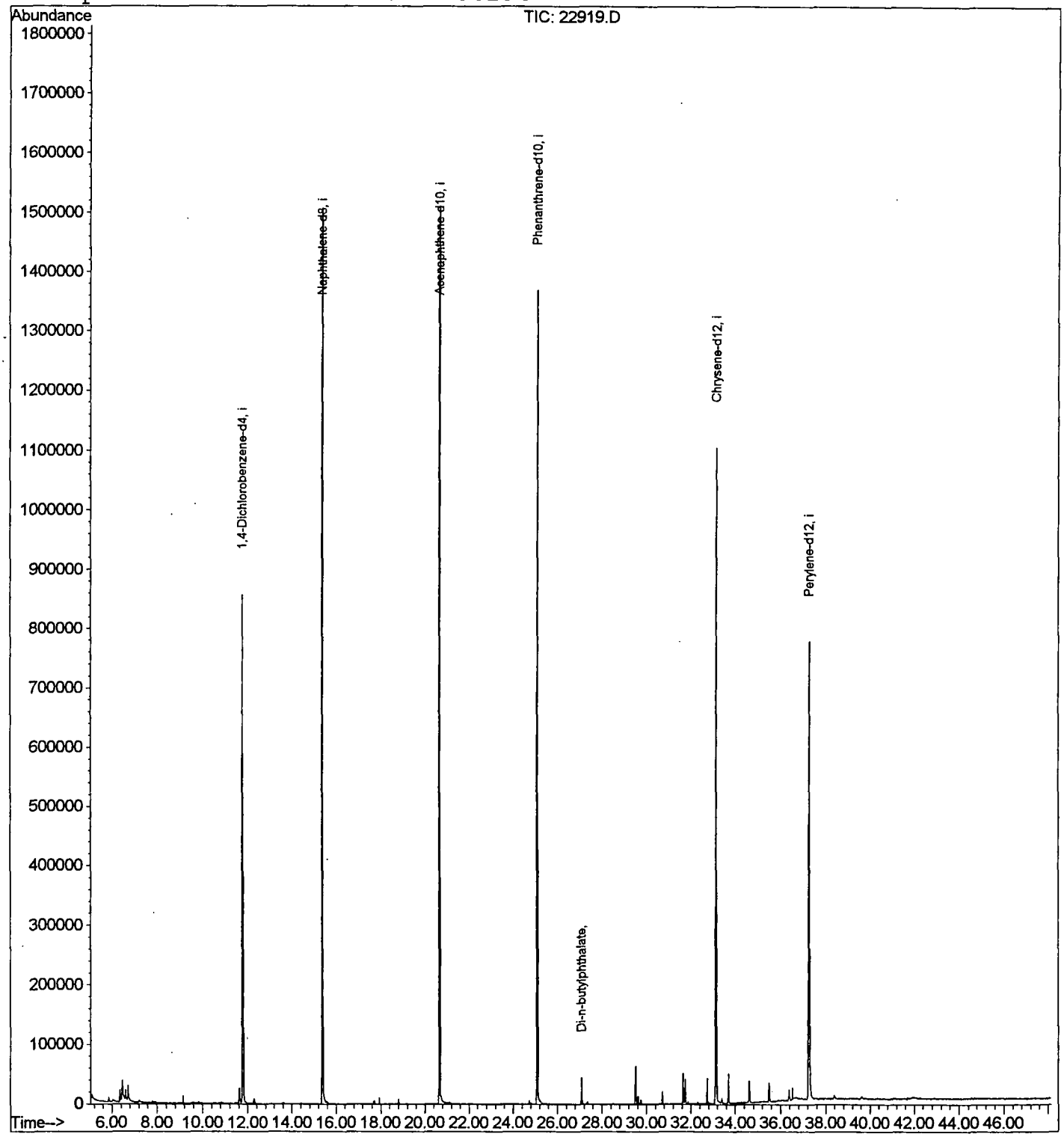
Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

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

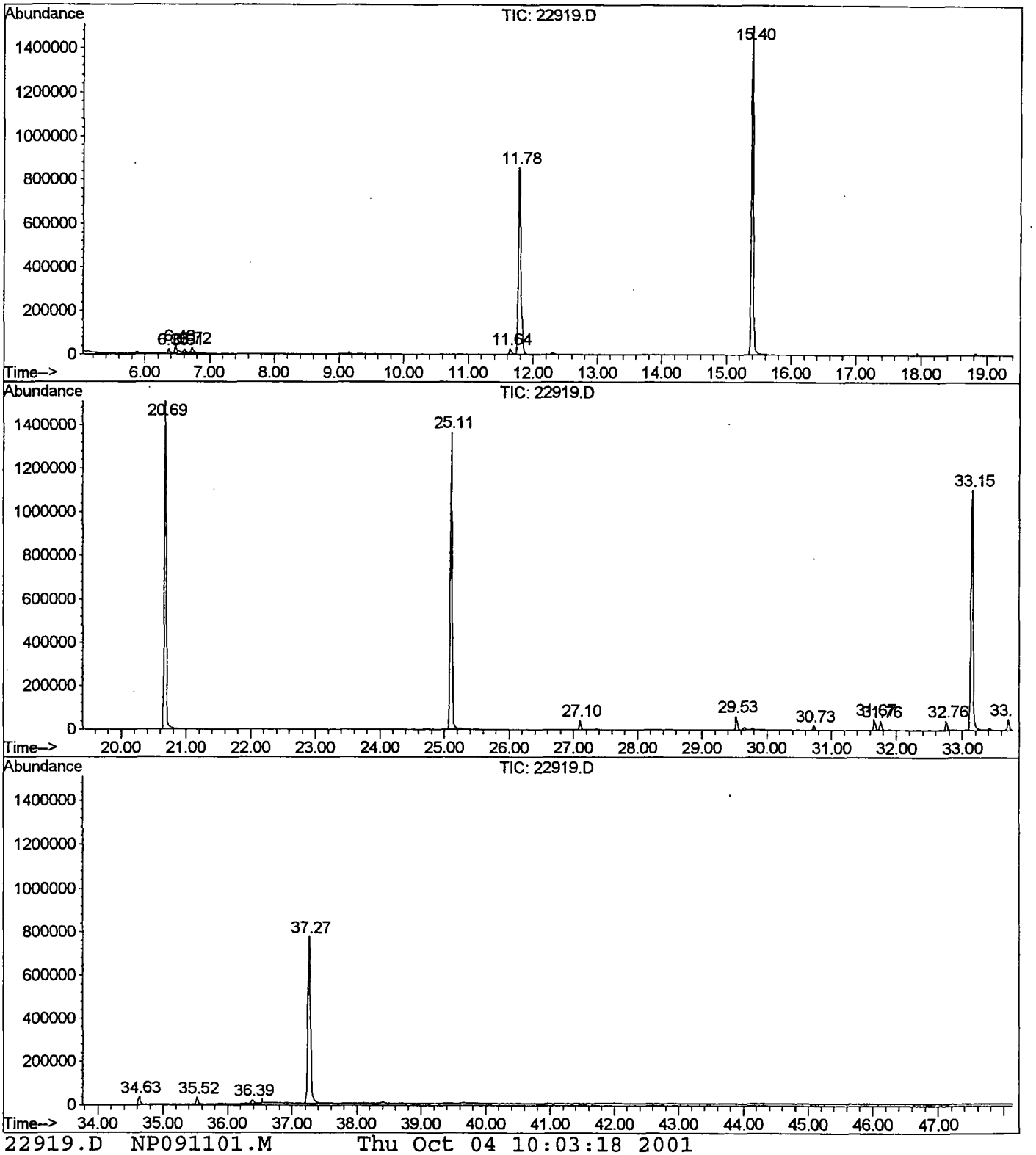
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.375	141	145	150	rBV	20444	45922	1.42%	0.263%
2	6.476	152	156	159	rBV	32200	58819	1.82%	0.337%
3	6.614	167	171	179	rVB	16146	36390	1.13%	0.208%
4	6.724	179	183	191	rVB	23482	48453	1.50%	0.277%
5	11.645	715	719	727	rBV	24714	59990	1.86%	0.343%
6	~~11.783~~	729	734	748	rVV	855123	2328976	72.14%	13.326%
7	~~15.400~~	1122	1128	1144	rBV	1506609	3173456	98.30%	18.158%
8	~~20.687~~	1697	1704	1718	rBV	1510100	3228453	100.00%	18.473%
9	~~25.112~~	2179	2186	2198	rBV	1367909	2897546	89.75%	16.579%
10	27.105	2398	2403	2410	rVB	43557	78734	2.44%	0.451%
11	29.528	2663	2667	2674	rBV	62871	123928	3.84%	0.709%
12	30.731	2793	2798	2804	rBV	20277	38880	1.20%	0.222%
13	31.667	2895	2900	2906	rBV	50300	93285	2.89%	0.534%
14	31.759	2906	2910	2916	rVB	40307	82206	2.55%	0.470%
15	32.760	3014	3019	3025	rBV	42112	81467	2.52%	0.466%
16	~~33.154~~	3055	3062	3082	rBV2	1104125	2538769	78.64%	14.526%
17	33.714	3115	3123	3130	rBV2	49195	105164	3.26%	0.602%
18	34.632	3214	3223	3231	rVB	36120	82102	2.54%	0.470%
19	35.523	3315	3320	3327	rBV	31240	67262	2.08%	0.385%
20	36.386	3409	3414	3421	rBV	18189	49441	1.53%	0.283%
21	~~37.267~~	3501	3510	3525	rBV2	769305	2257665	69.93%	12.918%

Sum of corrected areas: 17476908

22919.D NP091101.M Thu Oct 04 10:03:16 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Operator : 209 GALOS
Acquired : 2 Oct 2001 9:56 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 9
Quant File :NP091101.RES (RTE Integrator)



22919.D NP091101.M Thu Oct 04 10:03:18 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

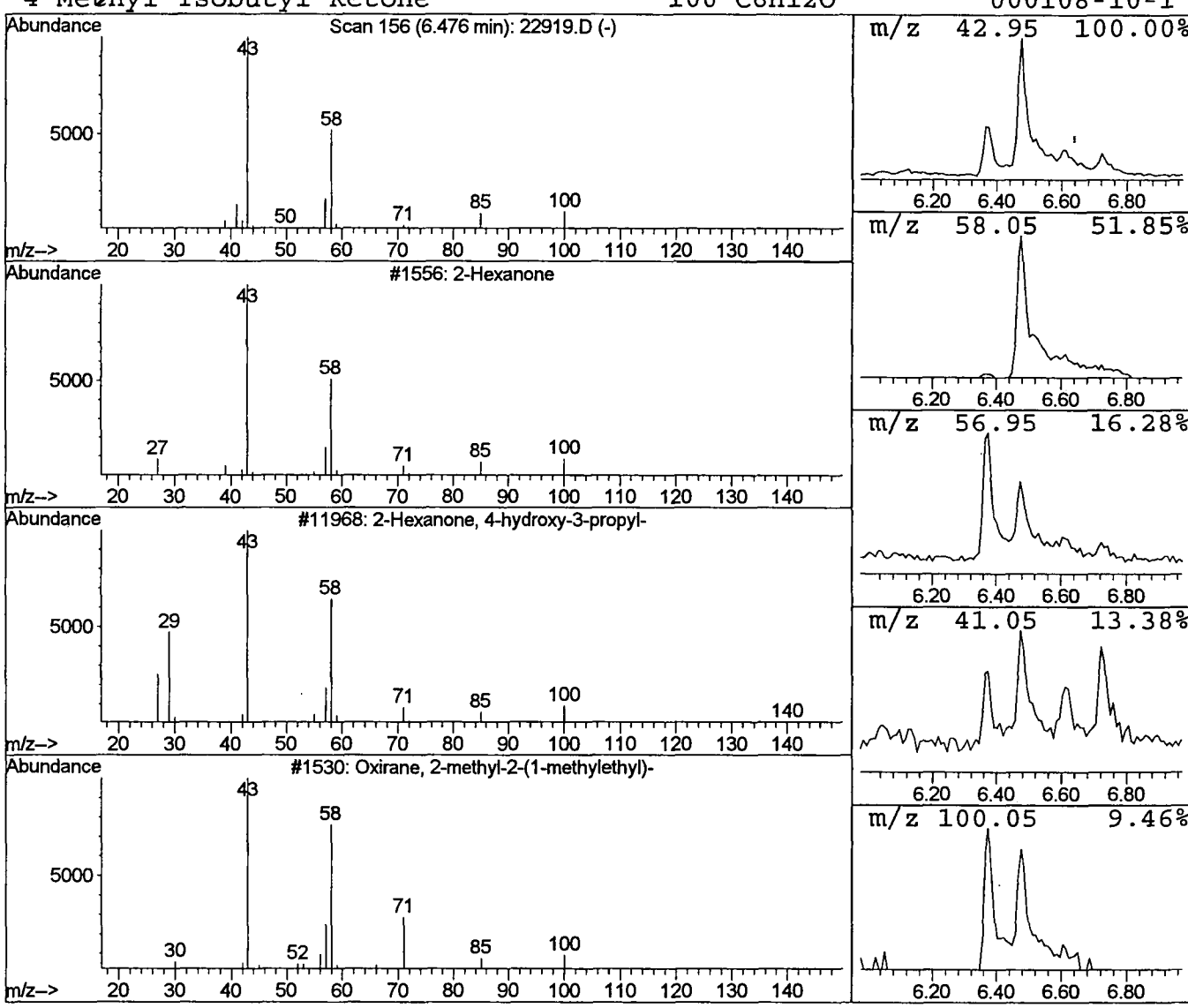
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Hexanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	0.51 ug/l	58819	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	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	74
3	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	64
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	49



Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

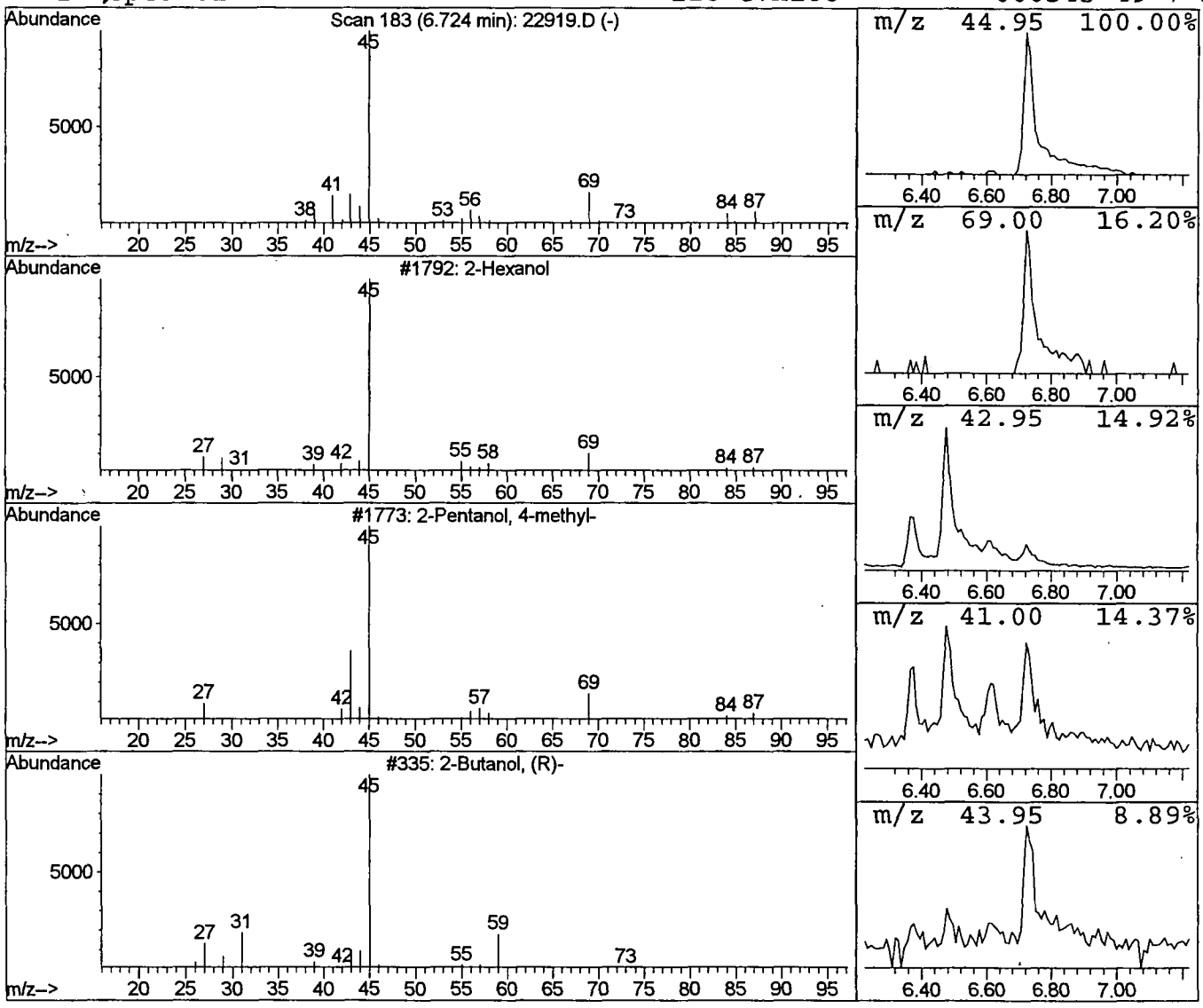
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	0.42 ug/l	48453	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	74
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	64
3	2-Butanol, (R)-	74	C4H10O	014898-79-4	45
4	2-Heptanol	116	C7H16O	000543-49-7	39



Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

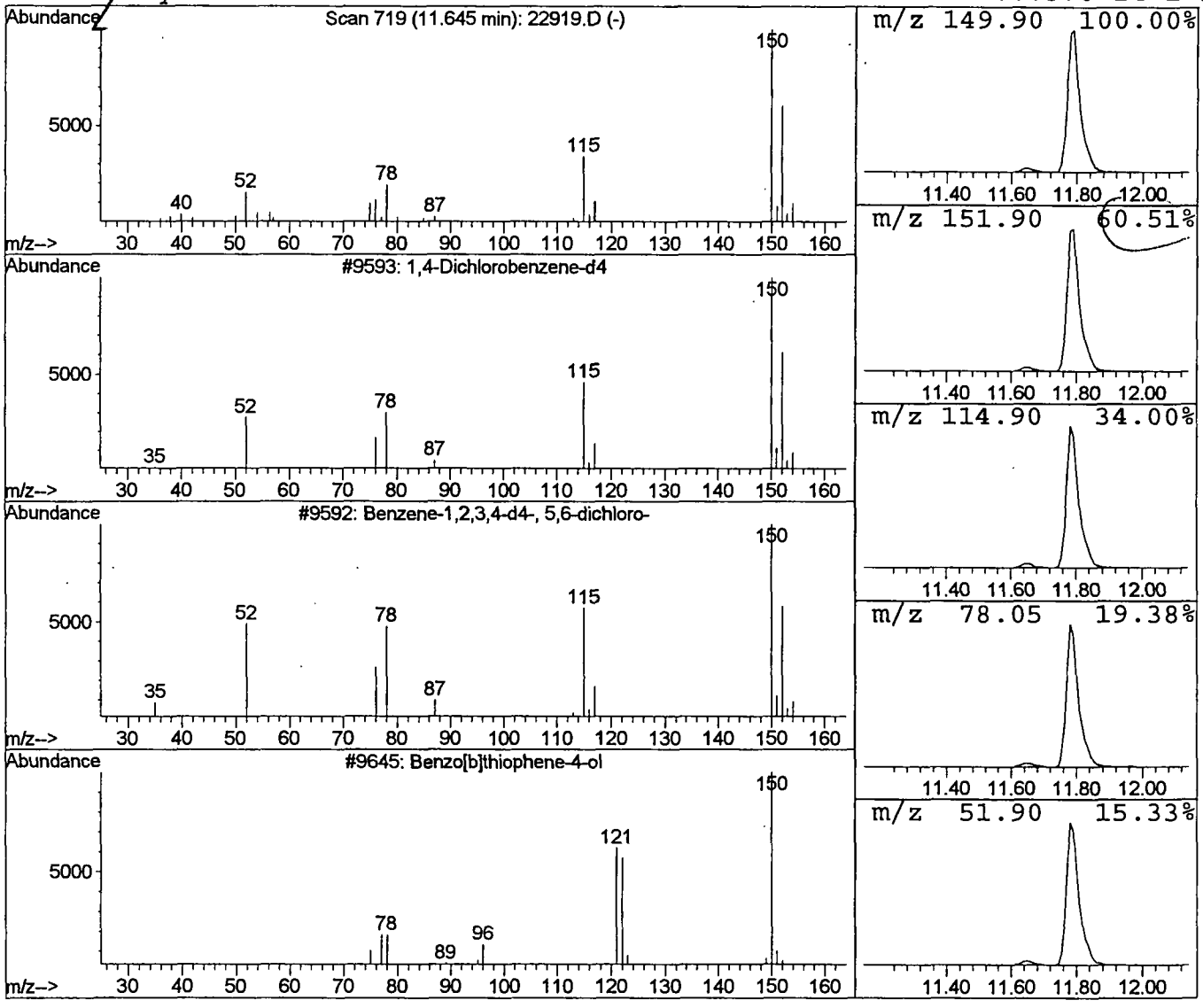
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 1,4-Dichlorobenzene-d4 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	0.52 ug/l	59990	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	33
3	Benzo[b]thiophene-4-ol	150	C8H6OS	003610-02-4	25
4	2-Methyl-6-nitroaniline	152	C7H8N2O2	000570-24-1	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

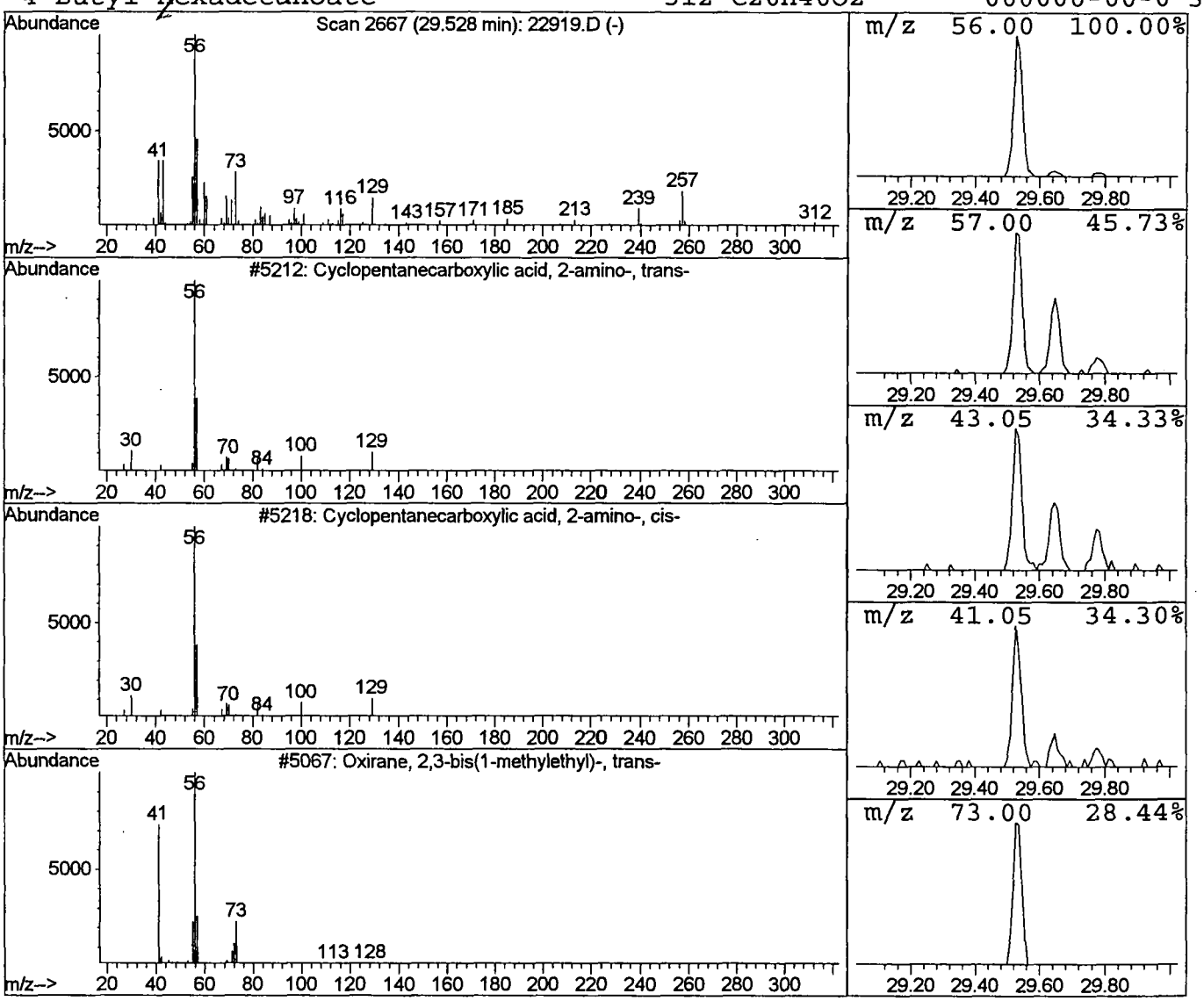
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Cyclopentanecarboxylic acid, 2 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.53	0.98 ug/l	123928	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	38
2	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	037910-65-9	38
3	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	37
4	Butyl hexadecanoate	312	C20H40O2	000000-00-0	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

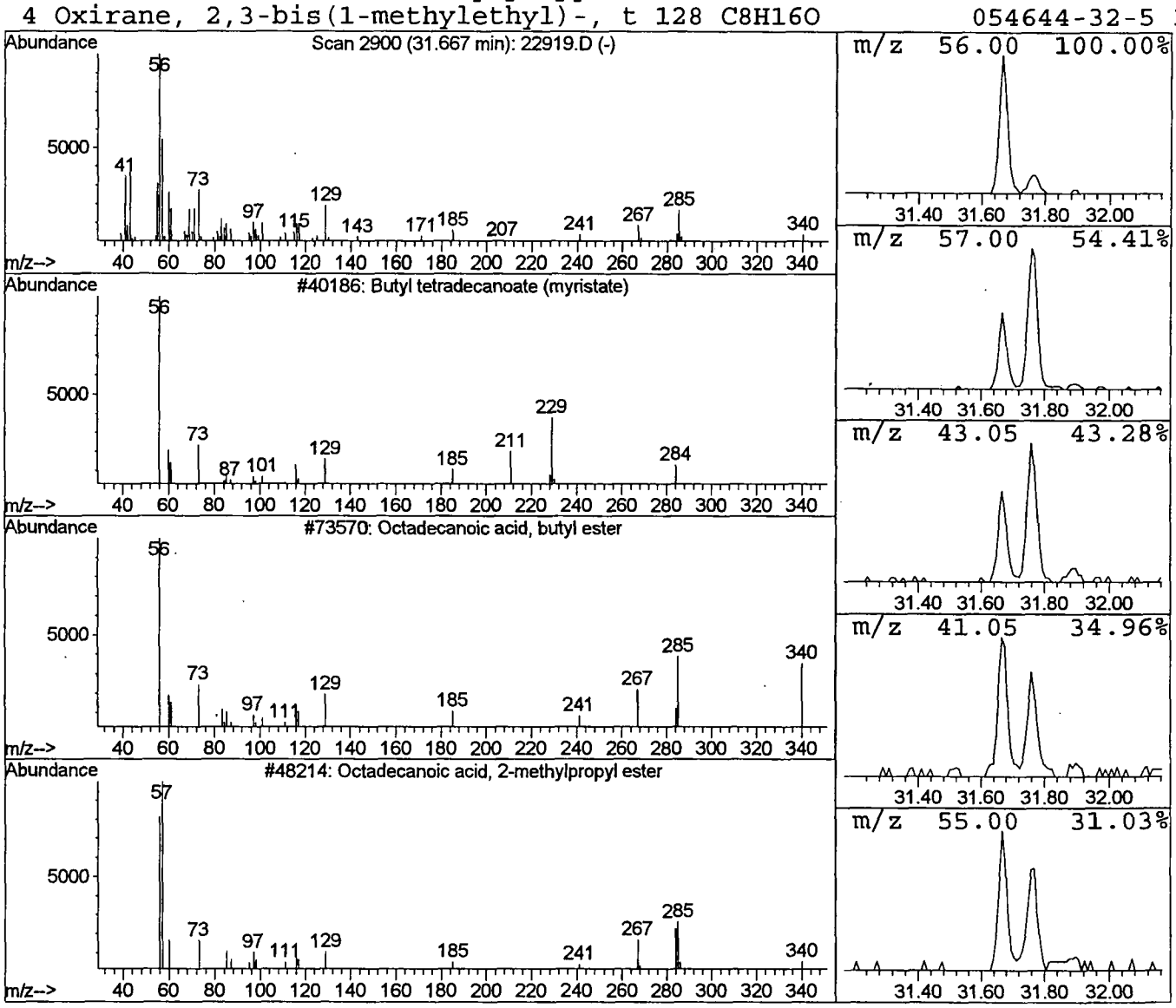
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Butyl tetradecanoate (myristat Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.67	0.73 ug/l	93285	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	80
2	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	64
3	Octadecanoic acid, 2-methylpropyl e	340	C22H44O2	000646-13-9	59
4	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

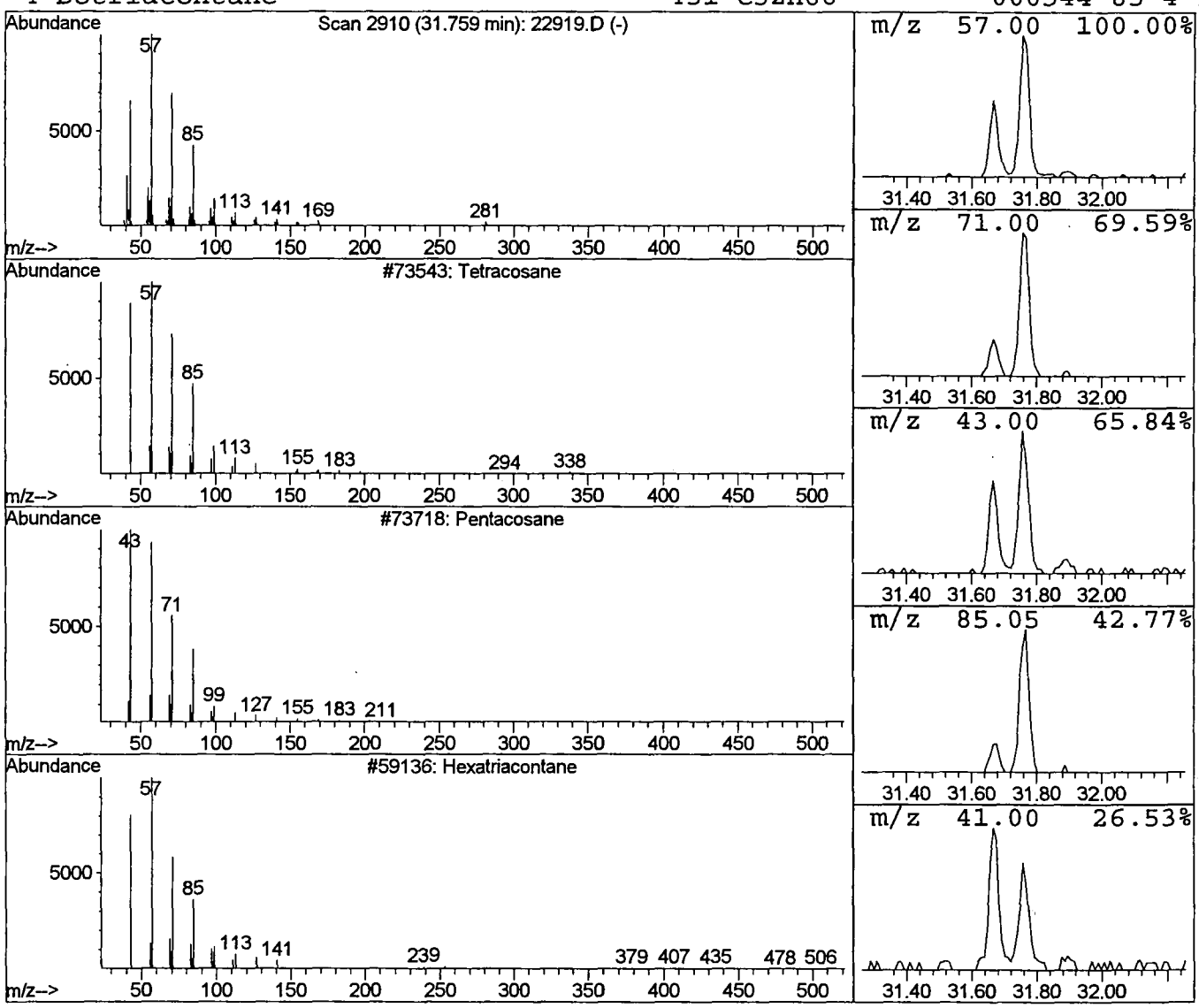
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Tetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.76	0.65 ug/l	82206	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	Hexatriacontane	507	C36H74	000630-06-8	91
4	Dotriacontane	451	C32H66	000544-85-4	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

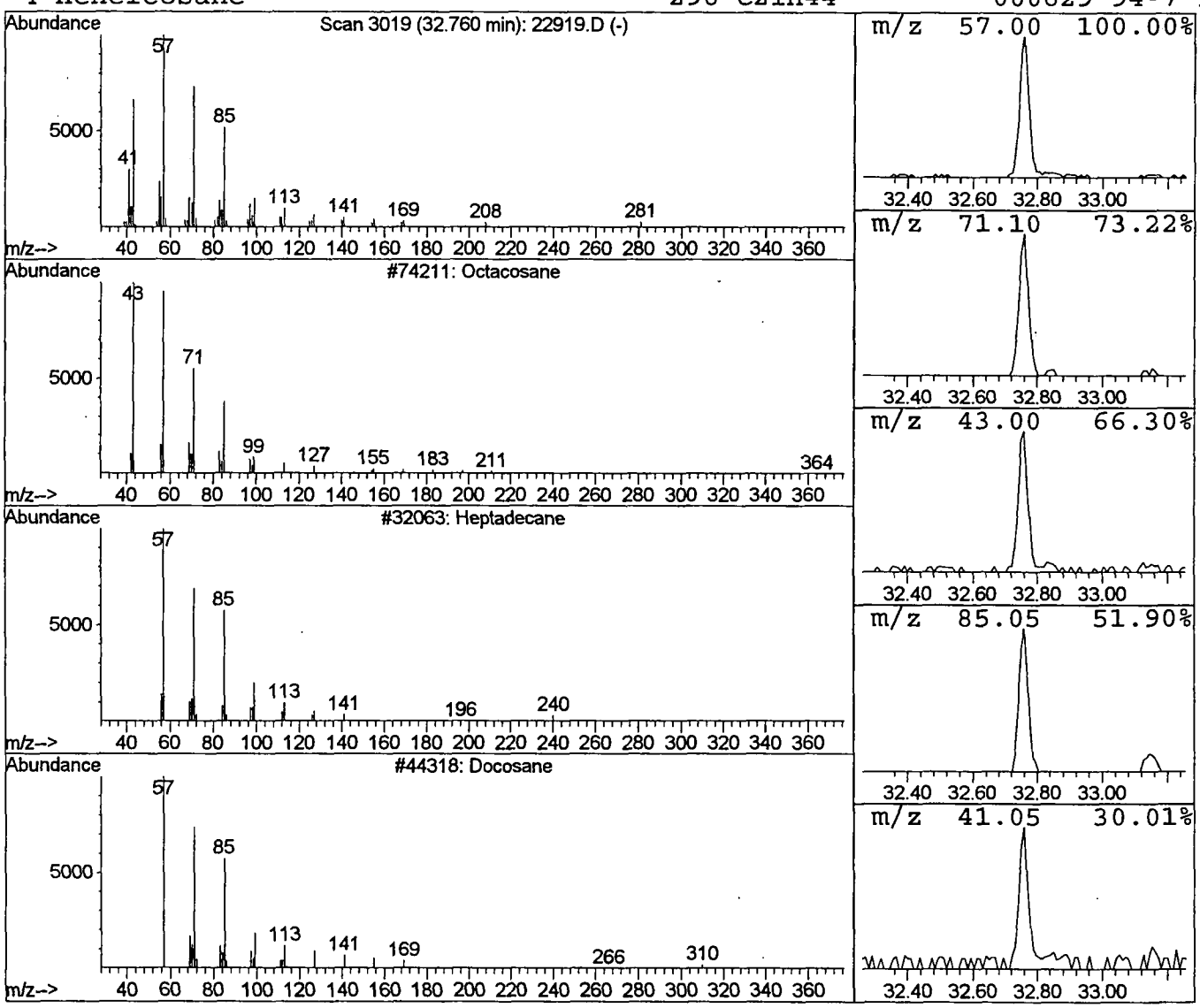
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Octacosane — Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.76	0.64 ug/l	81467	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Docosane	310	C22H46	000629-97-0	91
4	Heneicosane	296	C21H44	000629-94-7	91



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Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

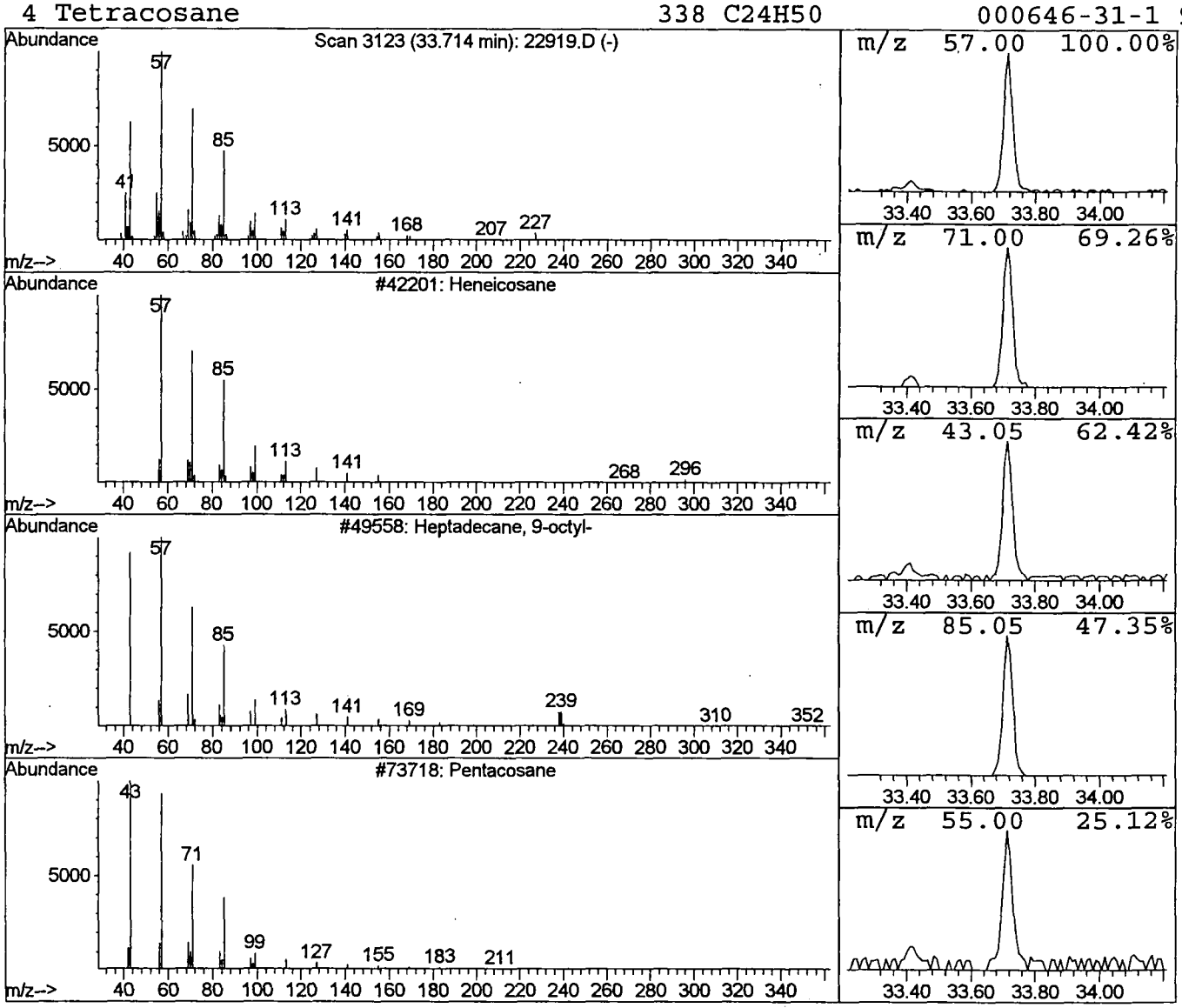
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.71	0.83 ug/l	105164	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Pentacosane	352	C25H52	000629-99-2	91
4	Tetracosane	338	C24H50	000646-31-1	90



Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
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Misc :
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Vial: 9
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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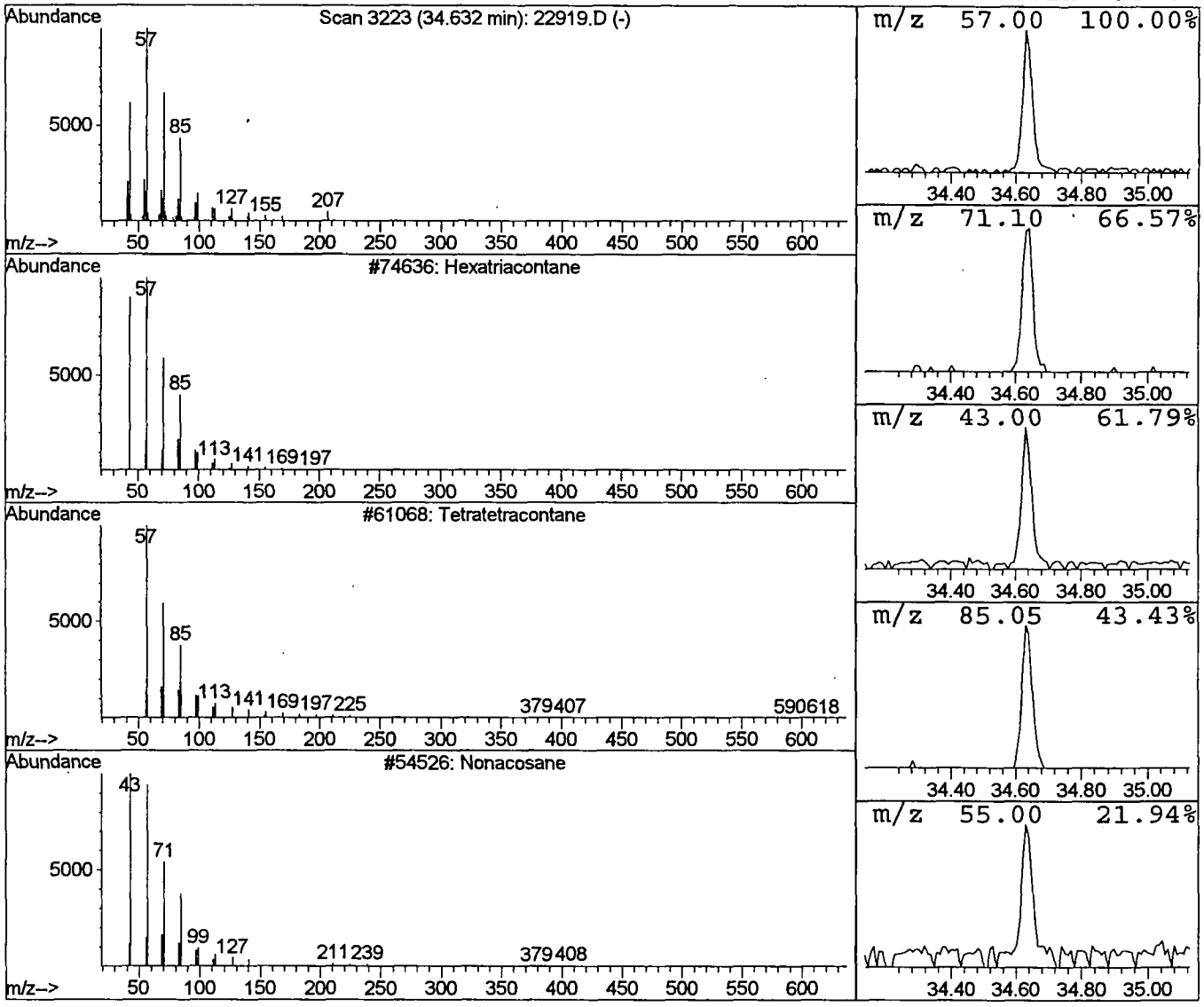
Peak Number 9 Hexatriacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.63	0.65 ug/l	82102	Chrysene-d12	33.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Nonacosane	408	C29H60	000630-03-5	90
4	Eicosane	282	C20H42	000112-95-8	87



Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
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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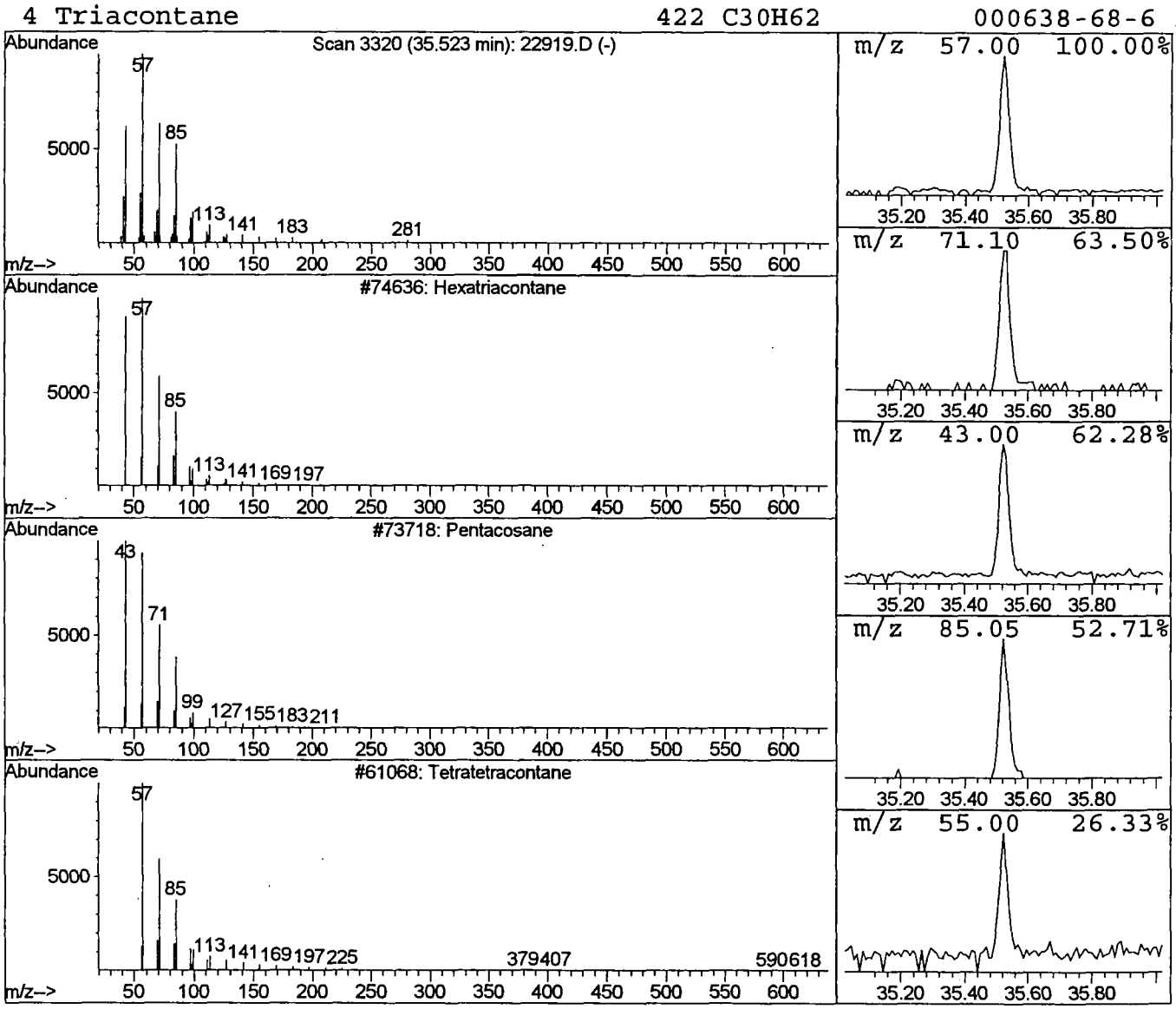
Peak Number 10 Hexatriacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.52	0.60 ug/l	67262	Perylene-d12	37.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	Tetratetracontane	619	C44H90	007098-22-8	90
4	Triacontane	422	C30H62	000638-68-6	90



Data File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Acq On : 2 Oct 2001 9:56 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
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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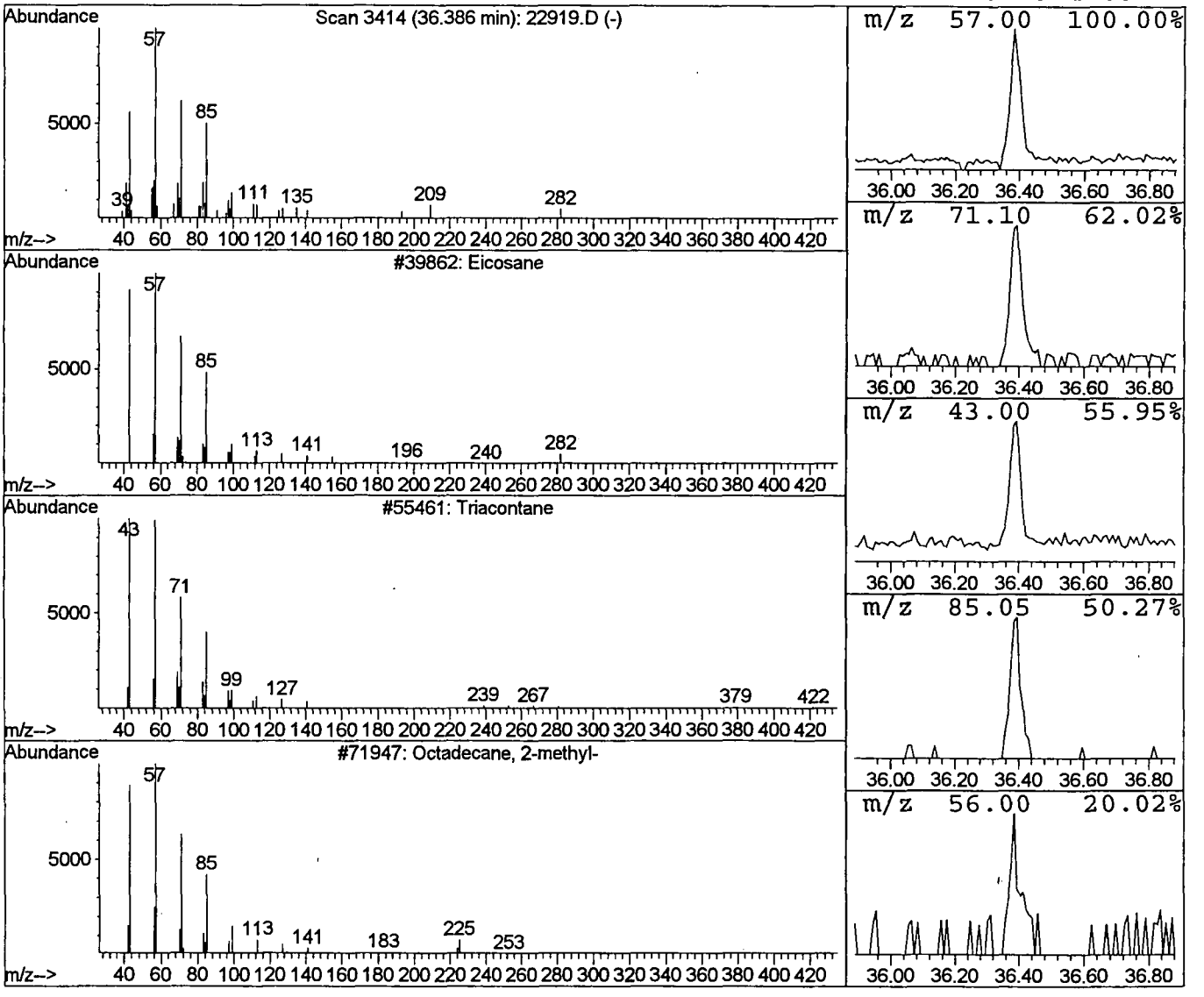
Peak Number 11 Eicosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.39	0.44 ug/l	49441	Perylene-d12	37.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Triacontane	422	C30H62	000638-68-6	86
3	Octadecane, 2-methyl-	268	C19H40	001560-88-9	86
4	Dotriacontane	451	C32H66	000544-85-4	86



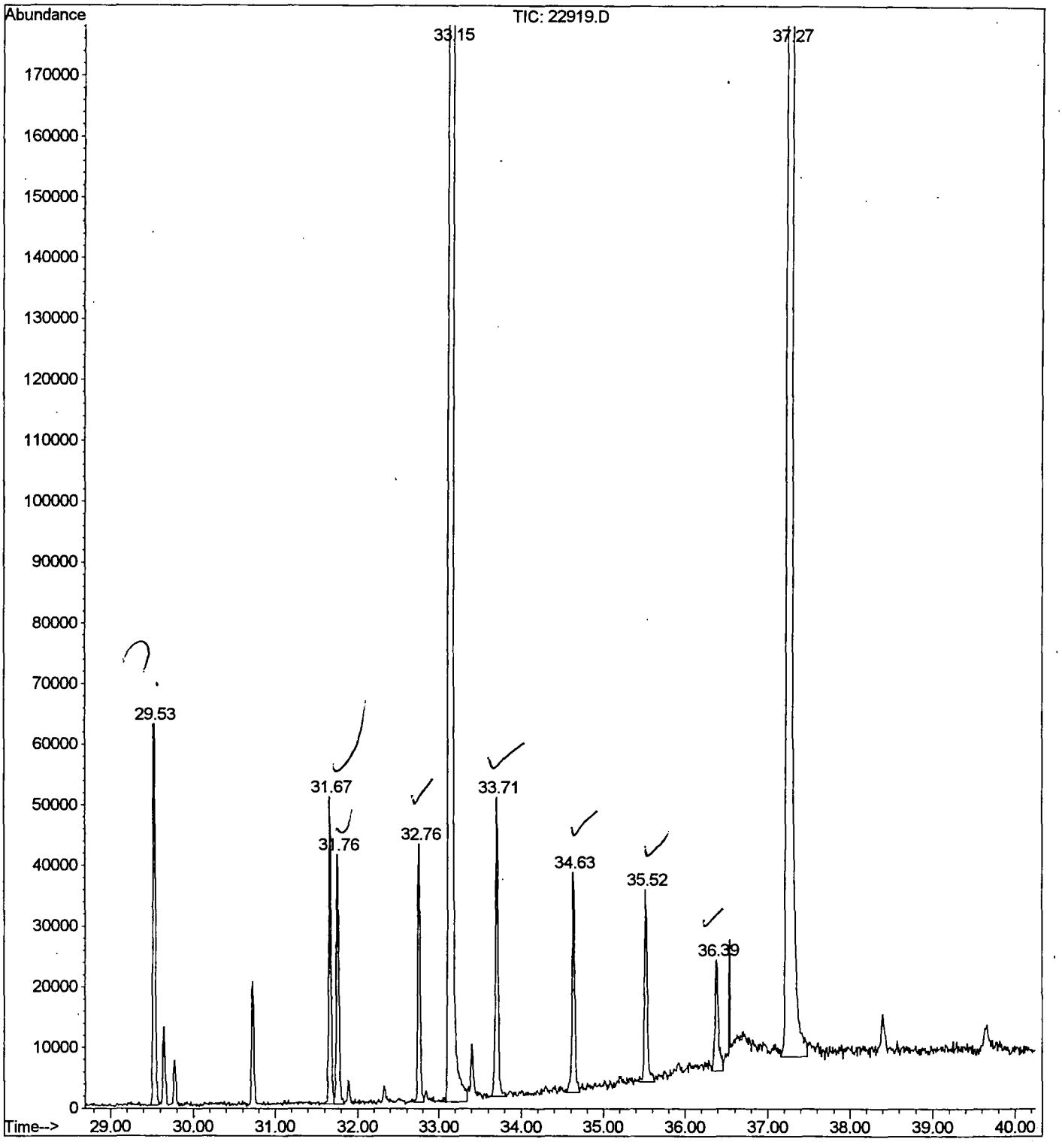
Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 2 Oct 2001 9:56 pm
Data File: C:\HPCHEM\1\DATA\OCT0201\22919.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
2-Hexanone	6.48	0.5	ug/l	58819	ISTD01	11.79	2328980	20.0
2-Hexanol	6.72	0.4	ug/l	48453	ISTD01	11.79	2328980	20.0
1,4-Dichlorobenzene-	11.64	0.5	ug/l	59990	ISTD01	11.79	2328980	20.0
Cyclopentanecarboxyl	29.53	1.0	ug/l	123928	ISTD05	33.15	2538770	20.0
Butyl tetradecanoate	31.67	0.7	ug/l	93285	ISTD05	33.15	2538770	20.0
Tetracosane	31.76	0.6	ug/l	82206	ISTD05	33.15	2538770	20.0
Octacosane	32.76	0.6	ug/l	81467	ISTD05	33.15	2538770	20.0
Heneicosane	33.71	0.8	ug/l	105164	ISTD05	33.15	2538770	20.0
Hexatriacontane	34.63	0.6	ug/l	82102	ISTD05	33.15	2538770	20.0
Hexatriacontane	35.52	0.6	ug/l	67262	ISTD06	37.27	2257670	20.0
Eicosane	36.39	0.4	ug/l	49441	ISTD06	37.27	2257670	20.0

22919.D NP091101.M Thu Oct 04 10:03:45 2001

File : C:\HPCHEM\1\DATA\OCT0201\22919.D
Operator : 209 GALOS
Acquired : 2 Oct 2001 9:56 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 9



Comments for Sample AD22919 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-08-2001 Time: 12:12:40

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 present in the Laboratory Blank.