LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN020320\
Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.564	827	868	892	rBV2	41369	172869	13.58%	1.061%
2	5.030	989	1013	1039	rVB3	37688	135614	10.66%	0.833%
3	5.545	1154	1173	1195	rBB2	26811	86048	6.76%	0.528%
4	6.329	1399	1417	1434	rBV6	4896	17745	1.39%	0.109%
5	6.500	1450	1470	1486	rBV2	23099	70673	5.55%	0.434%
6	6.603	1487	1502	1522	rVB3	16092	48237	3.79%	0.296%
7	7.304	1704	1720	1731	rBV3	8728	23311	1.83%	0.143%
8	7.368	1731	1740	1756	rVB5	7058	18700	1.47%	0.115%
9	7.567	1779	1802	1810	rBV	125209	333228	26.19%	2.046%
10	7.725	1840	1851	1871	rVB2	77983	207357	16.29%	1.273%
11	7.857	1875	1892	1923	rBV	147512	379371	29.81%	2.329%
12	8.008	1923	1939	1964	rVV	130940	319639	25.12%	1.963%
13	8.136	1964	1979	1990	rVV5	61409	163114	12.82%	1.002%
14	8.214	1992	2003	2016	rVV5	61975	185451	14.57%	1.139%
15	8.284	2016	2025	2047	rVV5	40142	102004	8.02%	0.626%
16	8.419	2047	2067	2084	rVB2	90071	197990	15.56%	1.216%
17	8.567	2096	2113	2135	rBV	382685	847313	66.58%	5.203%
18	8.661	2136	2142	2153	rVB5	7359	13505	1.06%	0.083%
19	8.731	2153	2164	2175	rVB4	14019	26839	2.11%	0.165%
20	8.805	2175	2187	2196	rBV4	24180	50720	3.99%	0.311%
21	9.066	2244	2268	2280	rBV3	190038	506762	39.82%	3.112%
22	9.130	2280	2288	2303	rVB2	53581	105219	8.27%	0.646%
23	9.255	2316	2327	2339	rVV2	36587	72335	5.68%	0.444%
24	9.352	2339	2357	2373	rVV5	40701	109447	8.60%	0.672%
25	9.503	2391	2404	2414	rBV4	60196	124939	9.82%	0.767%
26	9.603	2426	2435	2440	rBV2	28244	48415	3.80%	0.297%
27	9.651	2441	2450	2459	rVB6	52956	98978	7.78%	0.608%
28	9.718	2460	2471	2477	rBV	129631	243322	19.12%	1.494%
29	9.857	2494	2514	2536	rBV	163546	427779	33.61%	2.627%
30	9.956	2536	2545	2549	rBV7	10073	17098	1.34%	0.105%
31	10.005	2550	2560	2570	rVV4	35124	76570	6.02%	0.470%
32	10.079	2570	2583	2605	rVB	638245	1272584	100.00%	7.814%
33	10.204	2606	2622	2628	rBV5	40219	90844	7.14%	0.558%
34	10.275	2628	2644	2656	rVV5	137936	335187	26.34%	2.058%

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ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Title : SW846 8260

35	10.336	2656	2663	2670	rVV4	15226	25782	2.03%	0.158%
36	10.378	2670	2676	2682	rVV2	17210	24131	1.90%	0.148%
37	10.423	2682	2690	2699	rVB3	42668	69680	5.48%	0.428%
38	10.513	2707	2718	2740	rVB6	104605	241949	19.01%	1.486%
39	10.615	2742	2750	2761	rVB5	38861	72833	5.72%	0.447%
40	10.712	2761	2780	2789	rBV2	70924	137978	10.84%	0.847%
41	10.815	2790	2812	2823	rBV2	69644	188015	14.77%	1.154%
42	10.873	2824	2830	2835	rVV7	24243	38367	3.01%	0.236%
43	10.908	2836	2841	2846	rVV3	42643	66721	5.24%	0.410%
44	10.943	2847	2852	2860	rVV	56925	95066	7.47%	0.584%
45	10.998	2861	2869	2879	rVV3	61565	119054	9.36%	0.731%
46	11.056	2881	2887	2896	rVV9	30535	56415	4.43%	0.346%
47	11.117	2899	2906	2913	rVV3	43493	68860	5.41%	0.423%
48	11.191	2914	2929	2950	rVB5	176157	453579	35.64%	2.785%
49	11.294	2950	2961	2981	rBV	151357	294217	23.12%	1.807%
50	11.400	2983	2994	3009	rBV	511987	897828	70.55%	5.513%
51	11.586	3044	3052	3060	rBV6	26651	48954	3.85%	0.301%
52	11.651	3061	3072	3081	rBV5	48618	101724	7.99%	0.625%
53	11.744	3094	3101	3110	rBV8	14572	25699	2.02%	0.158%
54	11.808	3113	3121	3126	rBV5	12757	21036	1.65%	0.129%
55	11.853	3128	3135	3143	rVV7	25280	42098	3.31%	0.258%
56	11.921	3145	3156	3168	rVV8	46798	106604	8.38%	0.655%
57	12.043	3187	3194	3202	rVB	48990	69170	5.44%	0.425%
58	12.114	3209	3216	3223	rBV7	33018	49831	3.92%	0.306%
59	12.156	3224	3229	3238	rVB5	34929	54160	4.26%	0.333%
60	12.397	3296	3304	3314	rVV	388912	623502	48.99%	3.828%
61	12.442	3314	3318	3326	rVB	62025	76091	5.98%	0.467%
62	12.586	3356	3363	3373	rVB	138661	213970	16.81%	1.314%
63	12.873	3445	3452	3461	rBV5	17548	35602	2.80%	0.219%
64	12.924	3464	3468	3472	rVV4	24840	30527	2.40%	0.187%
65	12.959	3473	3479	3492	rVB5	46856	77982	6.13%	0.479%
66	13.088	3512	3519	3529	rVB2	33562	53573	4.21%	0.329%
67	13.149	3530	3538	3549	rBV7	38244	96466	7.58%	0.592%
68	13.284	3573	3580	3585	rBV3	43050	58189	4.57%	0.357%
69	13.342	3586	3598	3616	rVV2	522290	1018814	80.06%	6.256%
70	13.509	3642	3650	3658	rBV	113660	181036	14.23%	1.112%
71	13.593	3668	3676	3690	rVB3	77547	181305	14.25%	1.113%

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Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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72	13.670	3691	3700	3707	rBV6	43231	73373	5.77%	0.451%
73	13.738	3714	3721	3732	rVB	77152	117127	9.20%	0.719%
74	13.885	3757	3767	3777	rBV	357373	555070	43.62%	3.408%
75	13.940	3778	3784	3791	rVV3	59010	93458	7.34%	0.574%
76	13.992	3792	3800	3811	rVB	180746	288674	22.68%	1.773%
77	14.062	3817	3822	3831	rVB5	29580	43273	3.40%	0.266%
78	14.127	3834	3842	3847	rBV2	57631	97346	7.65%	0.598%
79	14.207	3859	3867	3886	rVB	197503	419618	32.97%	2.577%
80	14.294	3886	3894	3901	rBV3	97289	151270	11.89%	0.929%
81	14.474	3942	3950	3959	rBV2	134746	234001	18.39%	1.437%
82	14.590	3975	3986	3996	rBV5	239038	504108	39.61%	3.095%
83	14.644	3997	4003	4017	rVB3	101702	210164	16.51%	1.290%
84	14.850	4060	4067	4073	rVB4	102831	139869	10.99%	0.859%
85	14.956	4092	4100	4112	rVB3	142470	286163	22.49%	1.757%
86	15.416	4236	4243	4255	rVB2	71691	126533	9.94%	0.777%

Sum of corrected areas: 16286062

LSC Report - Integrated Chromatogram

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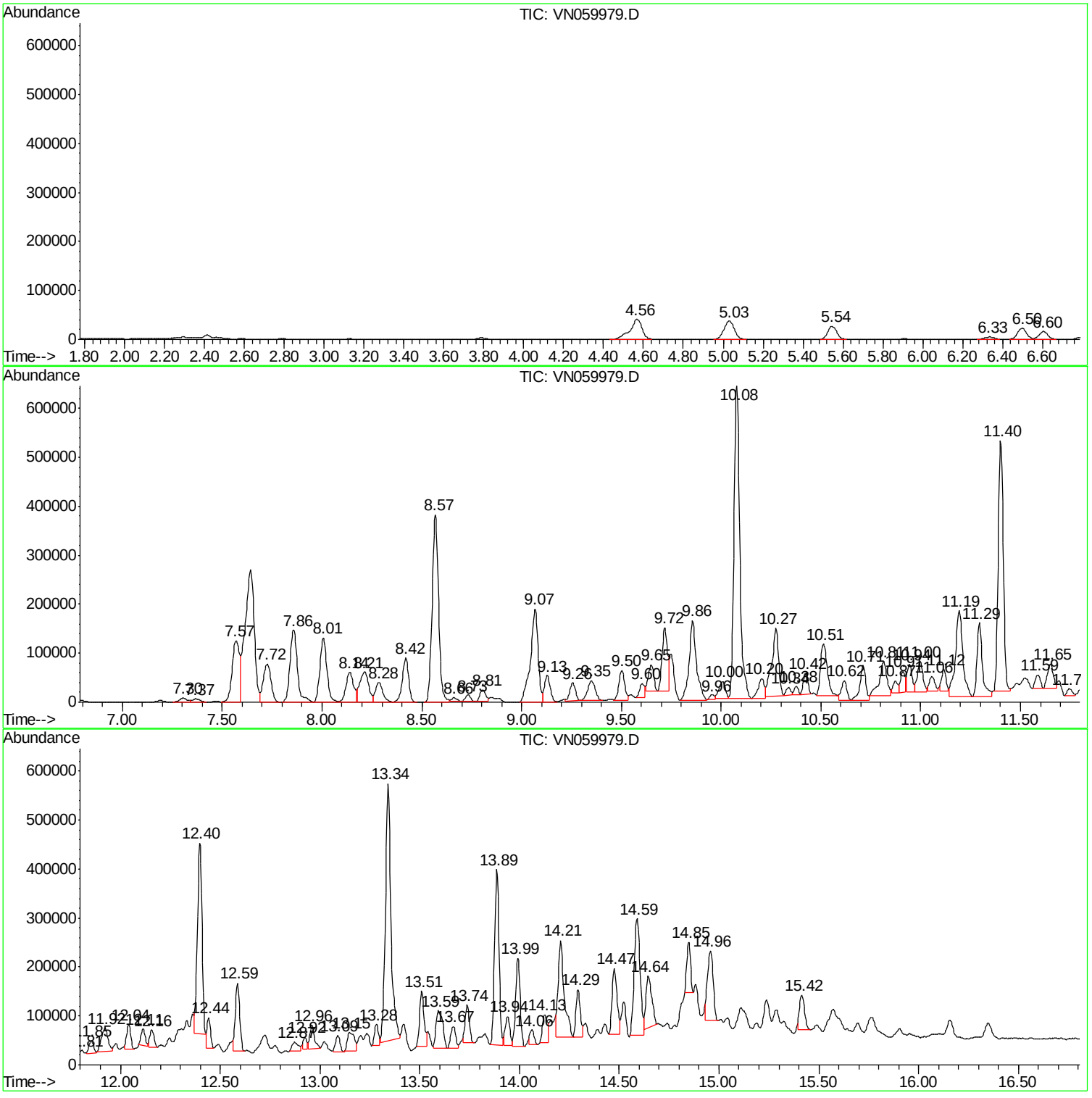
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN020320\
Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
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Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

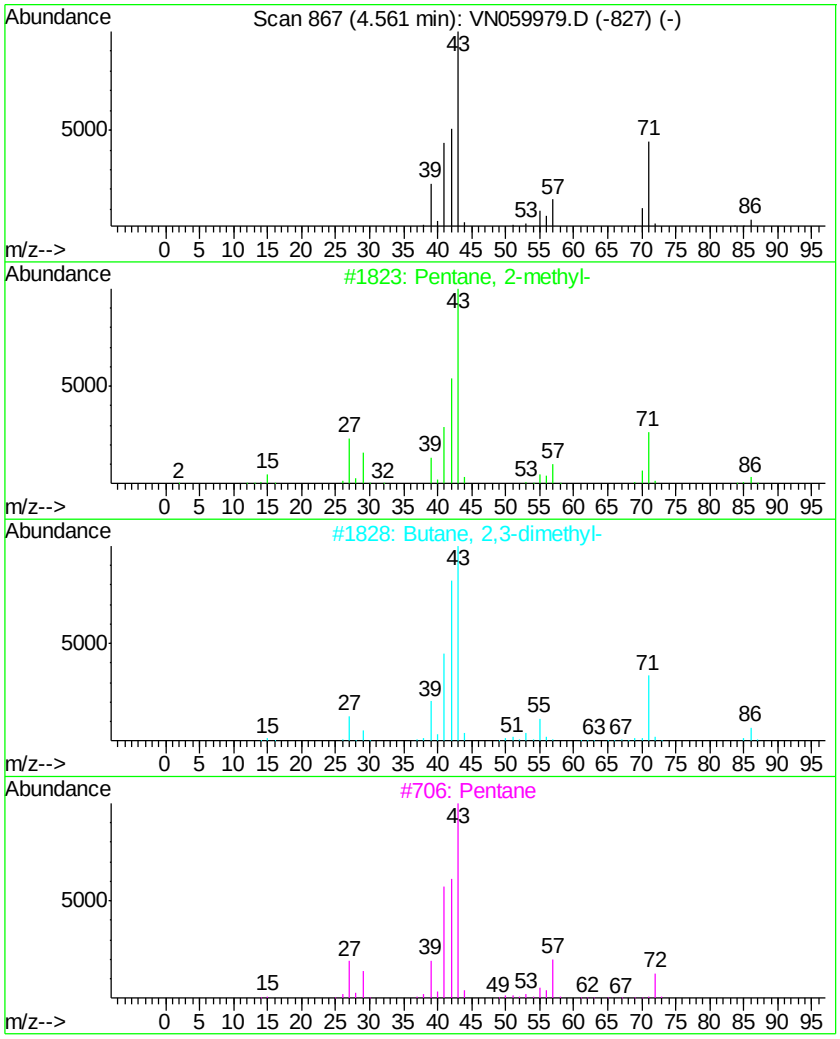
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

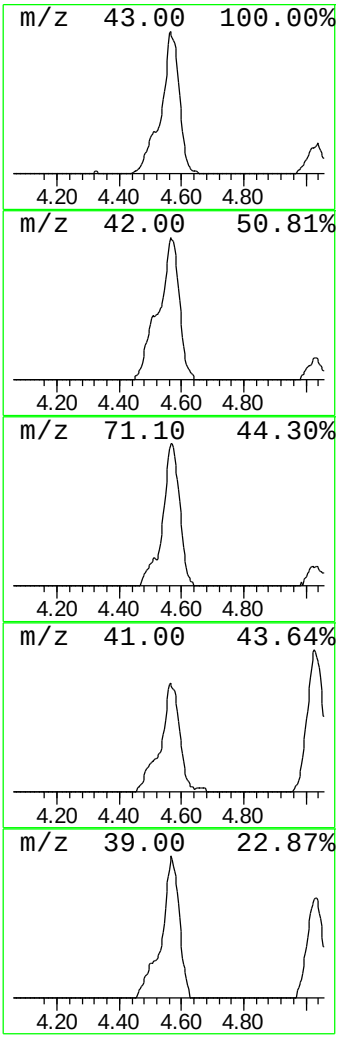
****

Peak Number 1 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	25.94 ug/l	172869	Pentafluorobenzene	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	49
3	Pentane	72	C5H12	000109-66-0	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	42





Library Search Compound Report

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Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

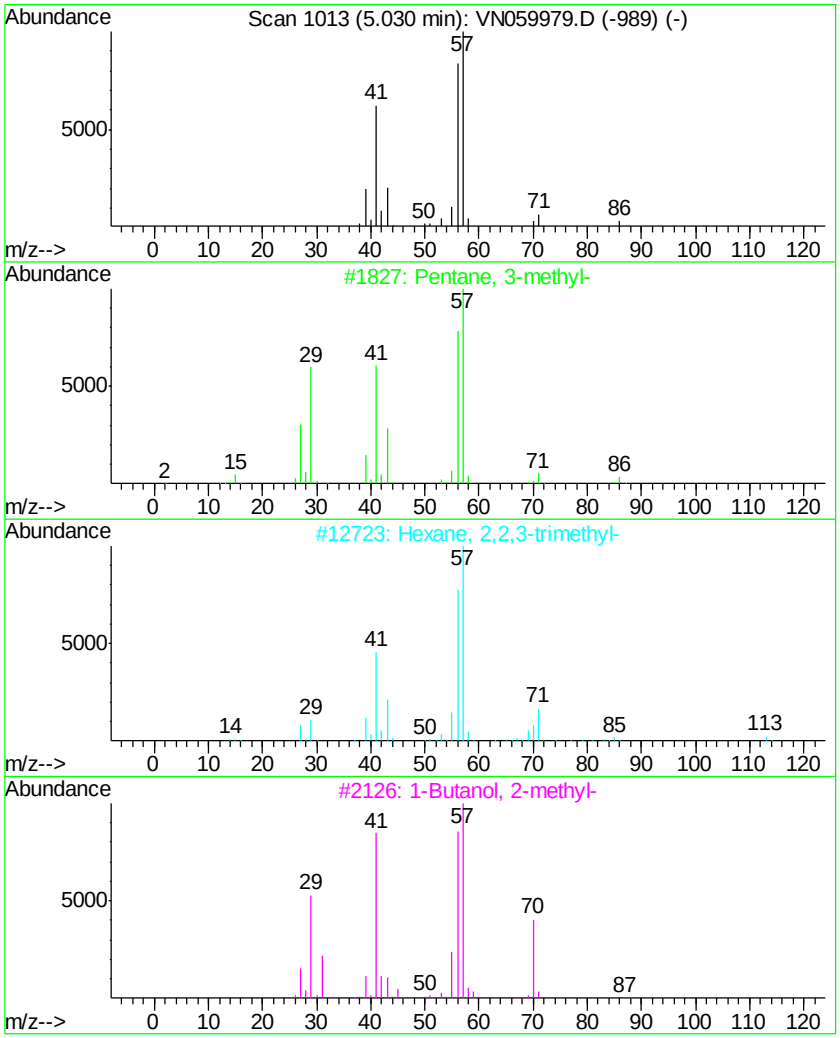
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

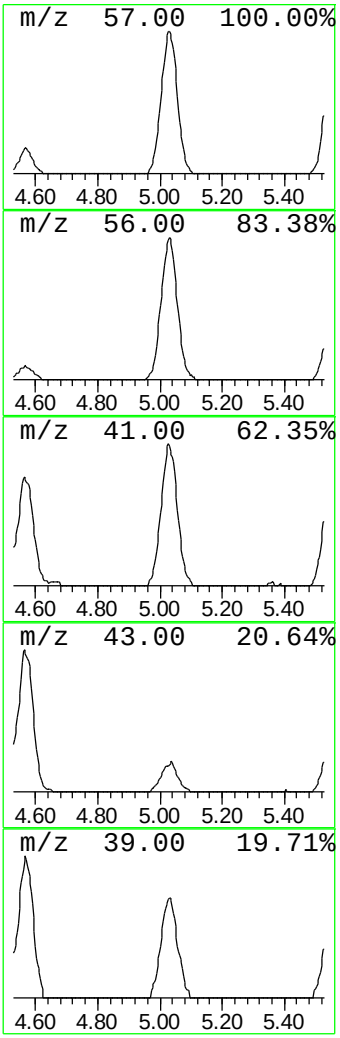
****

Peak Number 2 Pentane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	20.35 ug/l	135614	Pentafluorobenzene	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	83
3	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40
4	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40





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Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

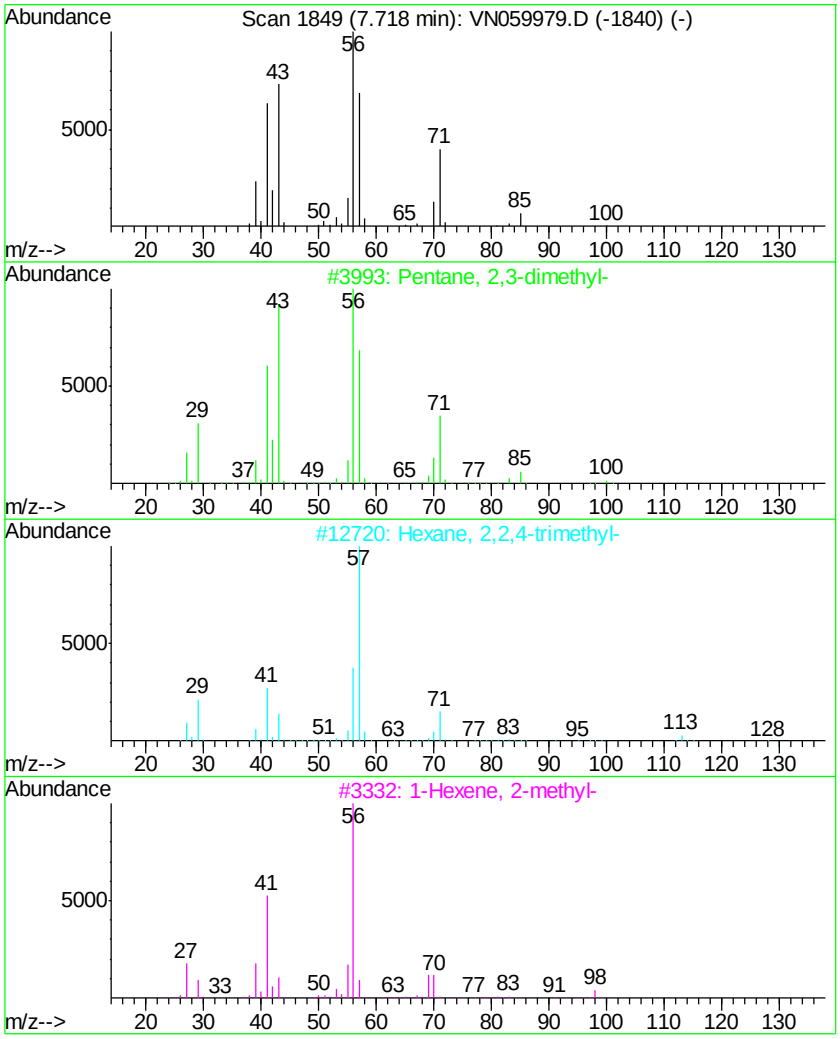
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

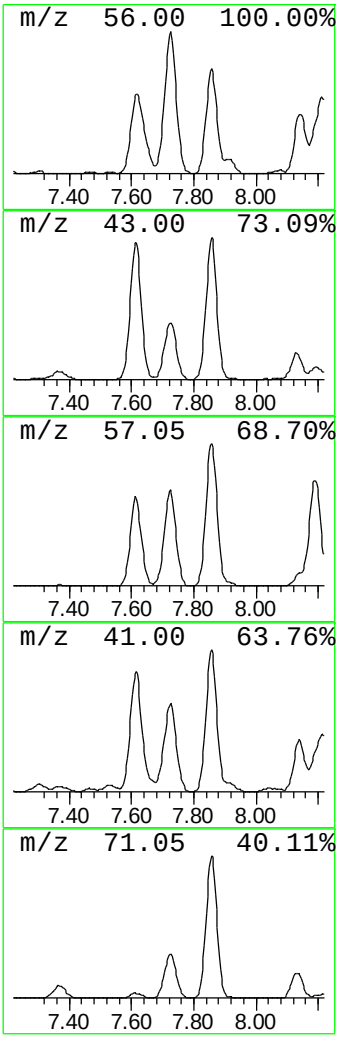
****

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	31.11 ug/l	207357	Pentafluorobenzene	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	56
3	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	47
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	37
5	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	37





Library Search Compound Report

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ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

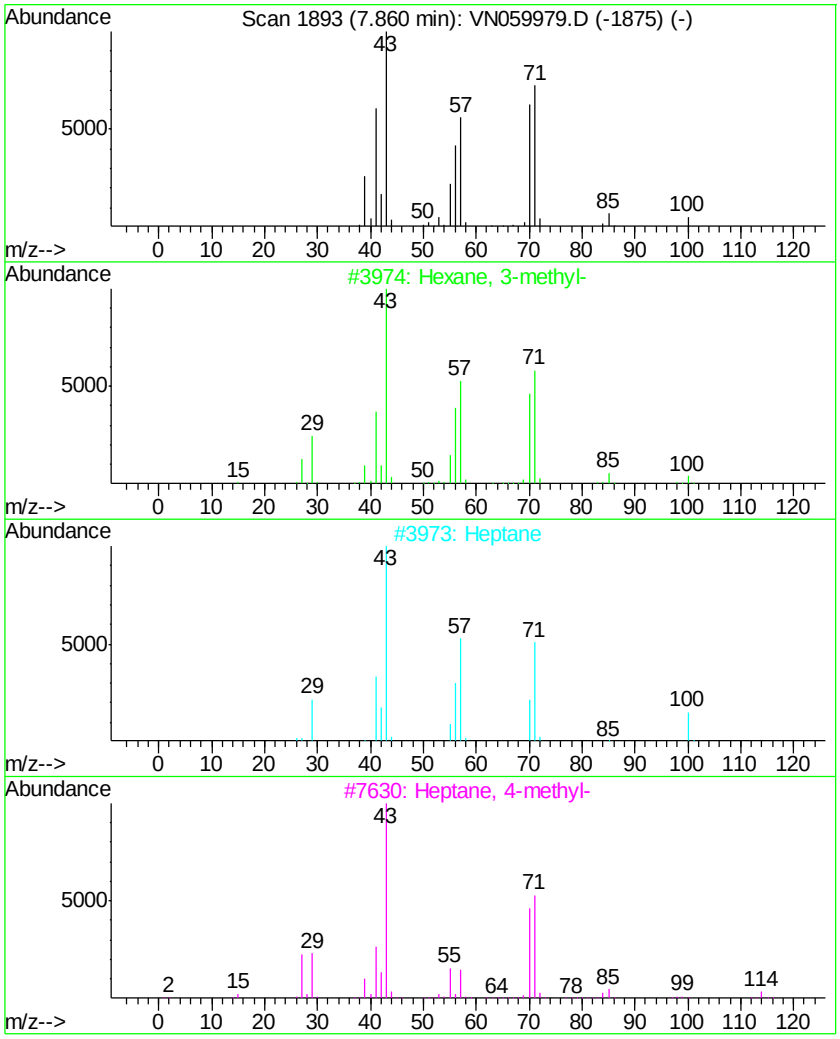
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

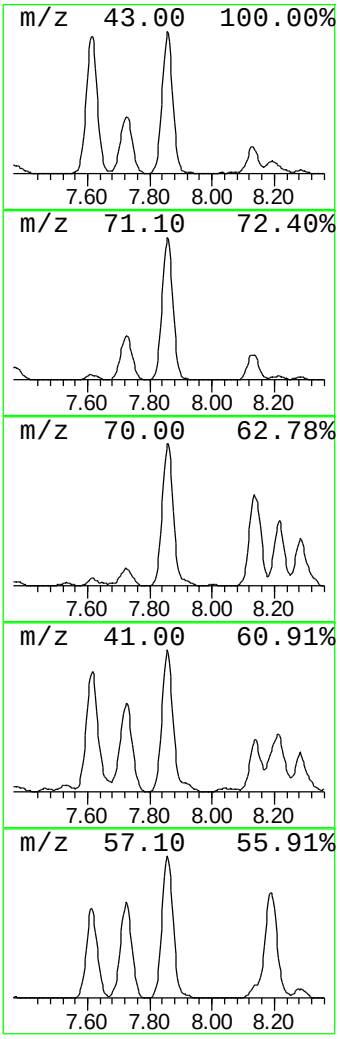
****

Peak Number 4 Hexane, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	56.92 ug/l	379371	Pentafluorobenzene	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2	Heptane	100	C7H16	000142-82-5	80
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	53





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Operator : JC/MD
Sample : L1346-04 40X
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ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

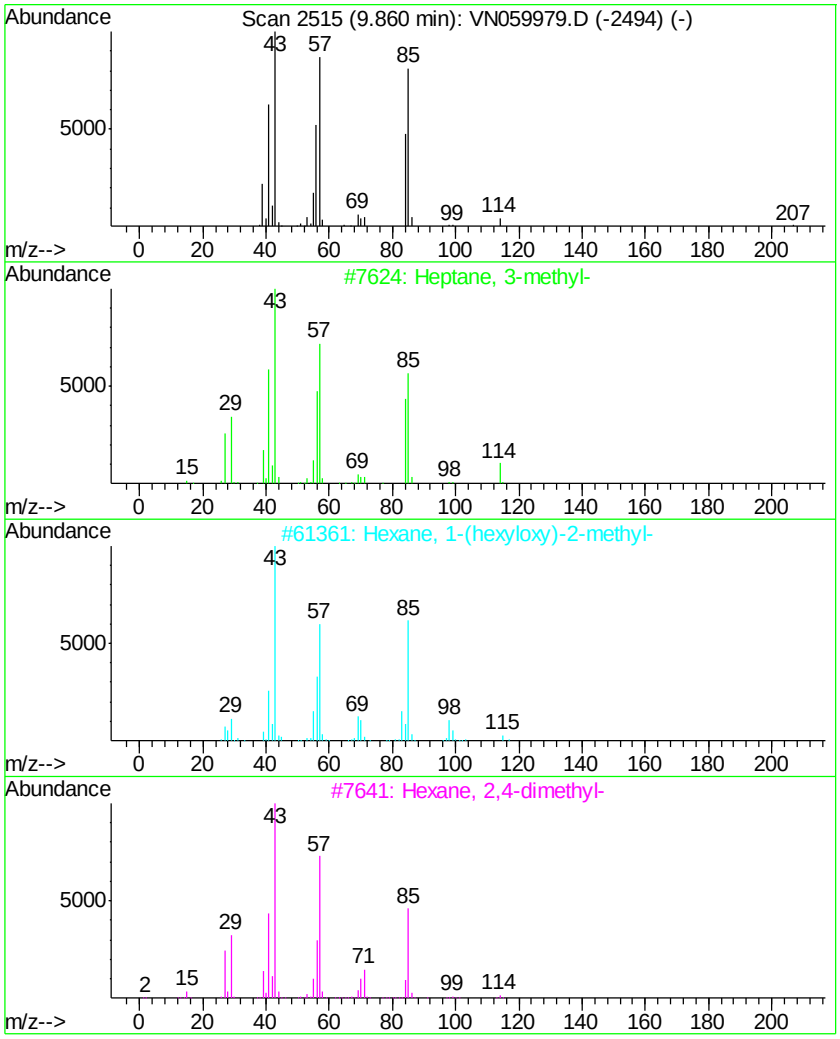
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

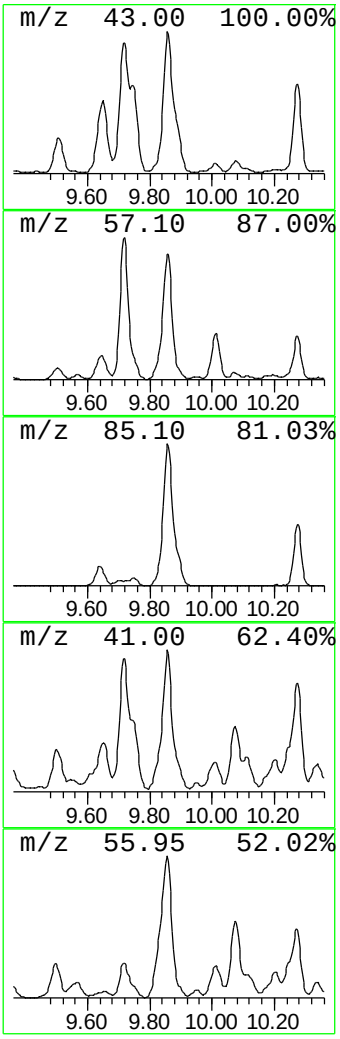
****

Peak Number 5 Heptane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	25.24 ug/l	427779	1,4-Difluorobenzene	8.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2	Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	64
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
4	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59
5	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59





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ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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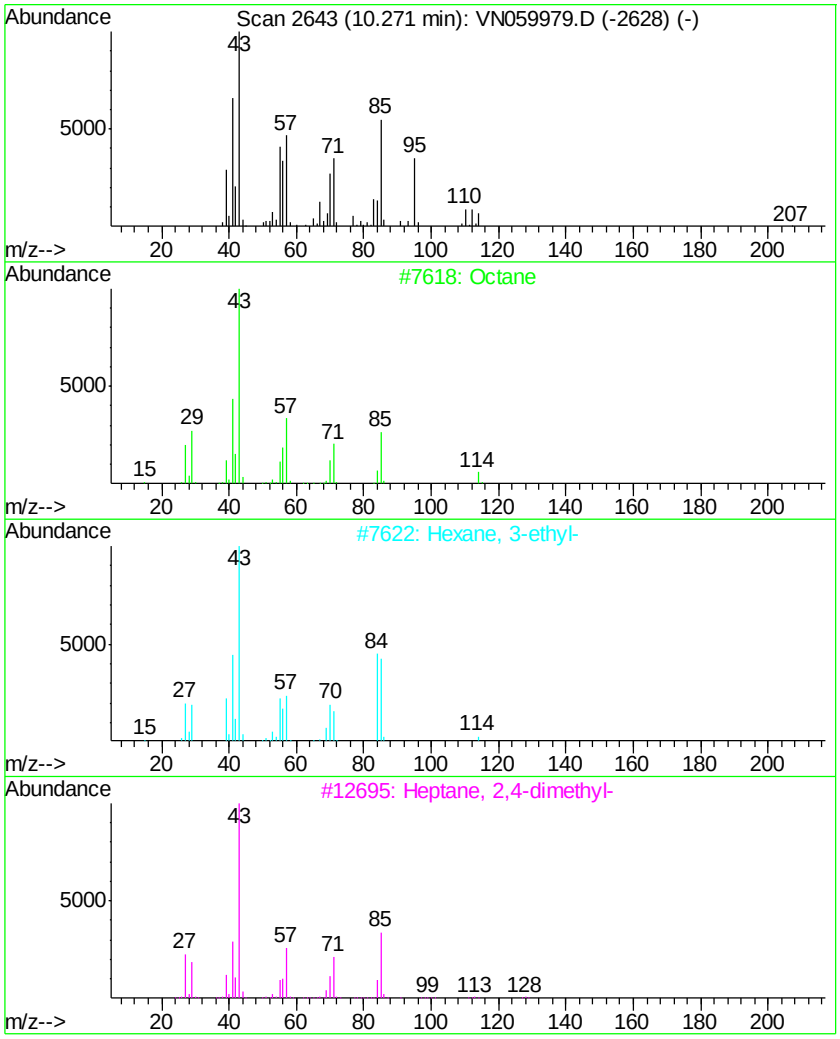
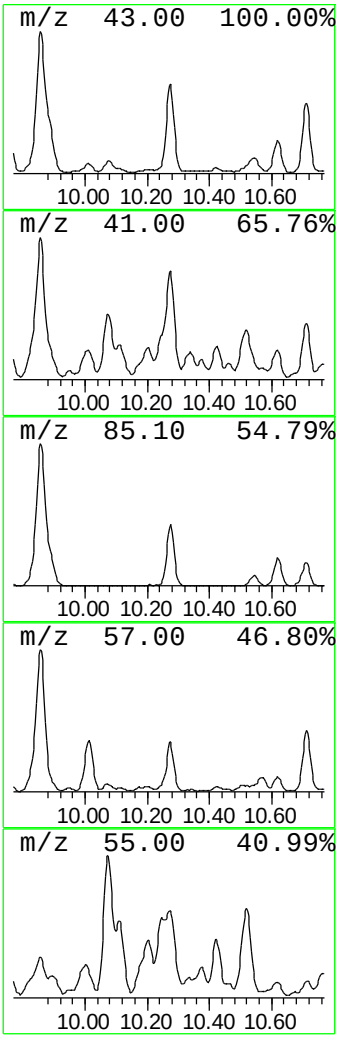
Peak Number 6 Octane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	18.67 ug/l	335187	Chlorobenzene-d5	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	64
2	Hexane, 3-ethyl-	114	C8H18	000619-99-8	49
3	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	43
4	Oxalic acid, diisohexyl ester	258	C14H26O4	1000309-32-9	35
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	35





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Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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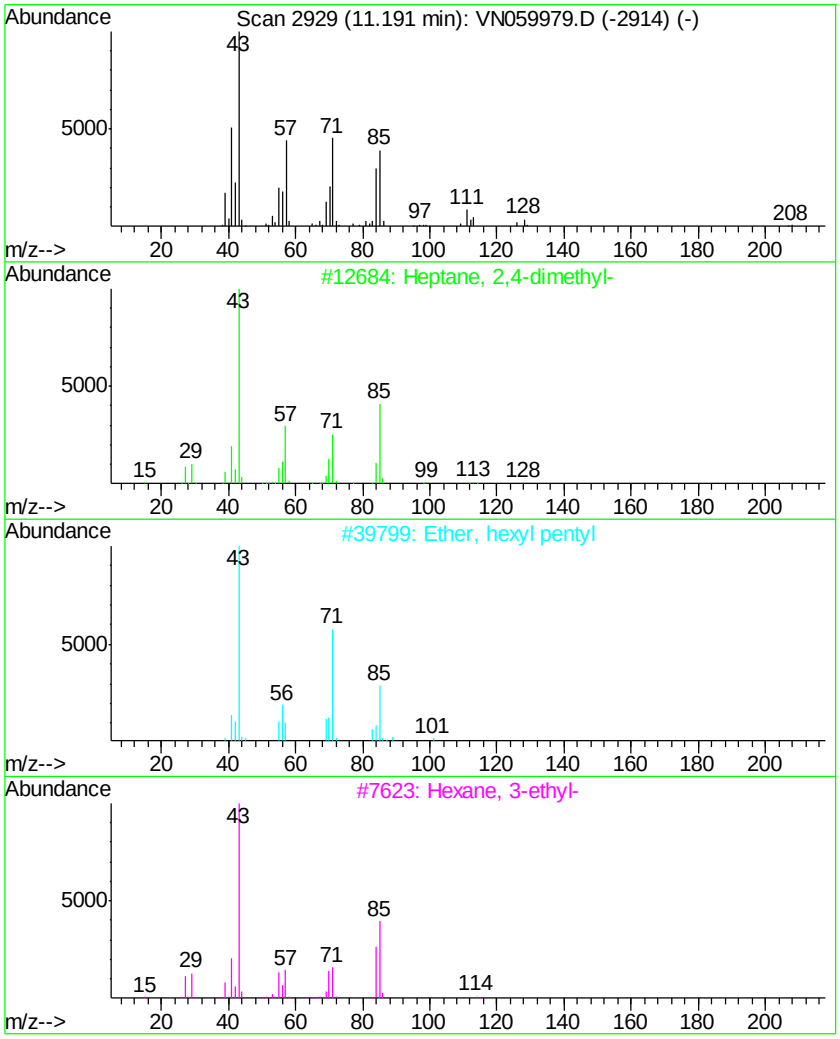
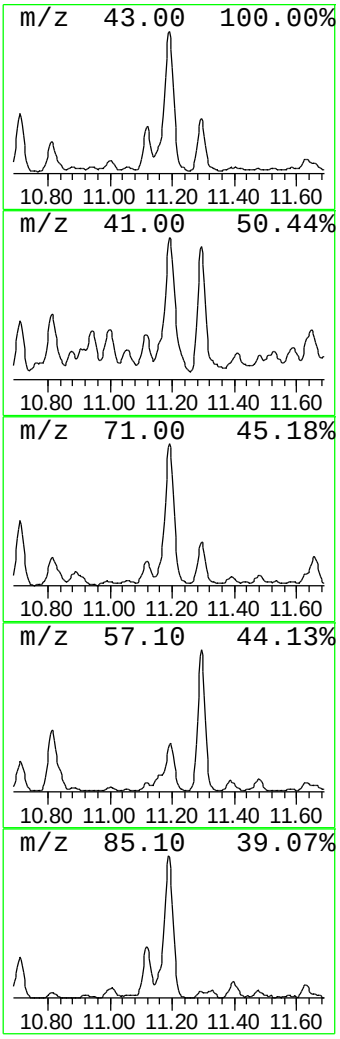
Peak Number 7 Heptane, 2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	25.26 ug/l	453579	Chlorobenzene-d5	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	72
2	Ether, hexyl pentyl	172	C11H24O	032357-83-8	53
3	Hexane, 3-ethyl-	114	C8H18	000619-99-8	53
4	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	53
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN020320\
Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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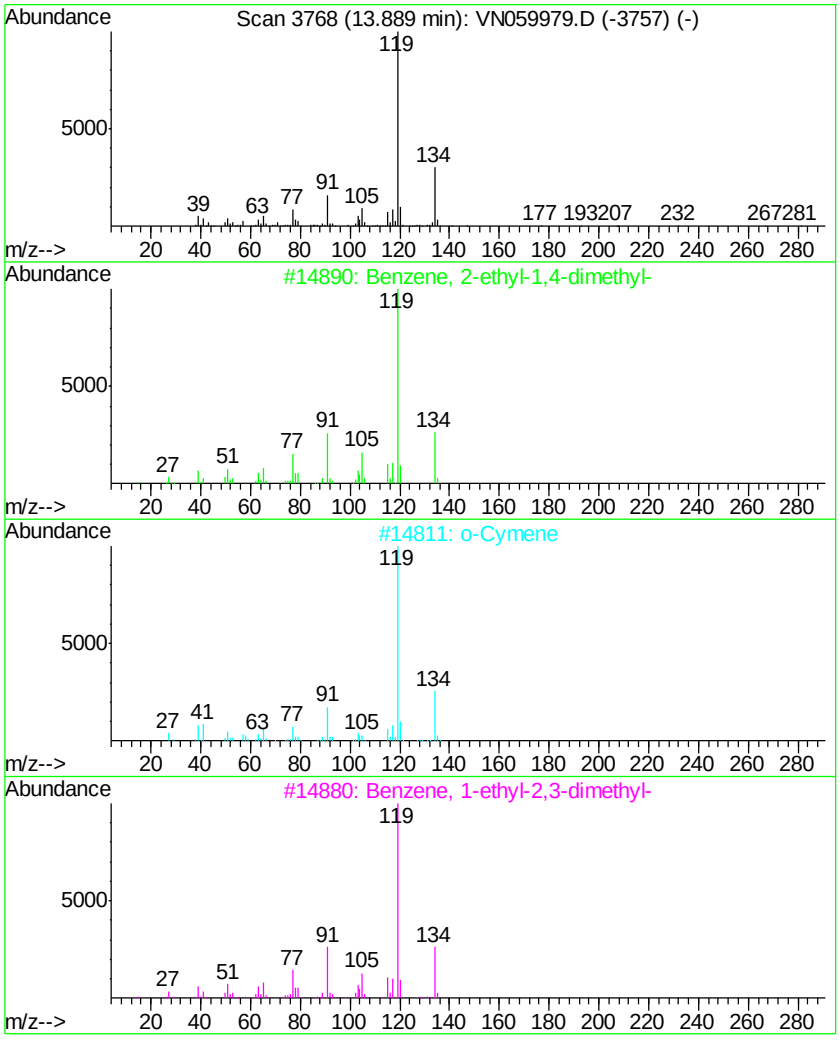
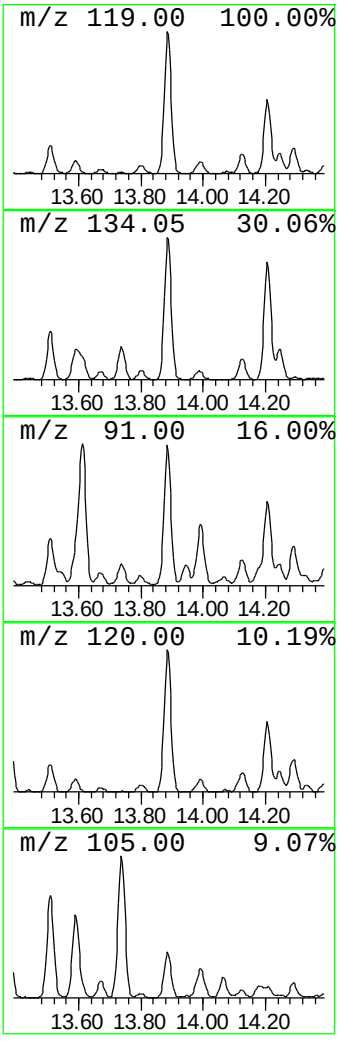
Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	27.24 ug/l	555070	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN020320\
Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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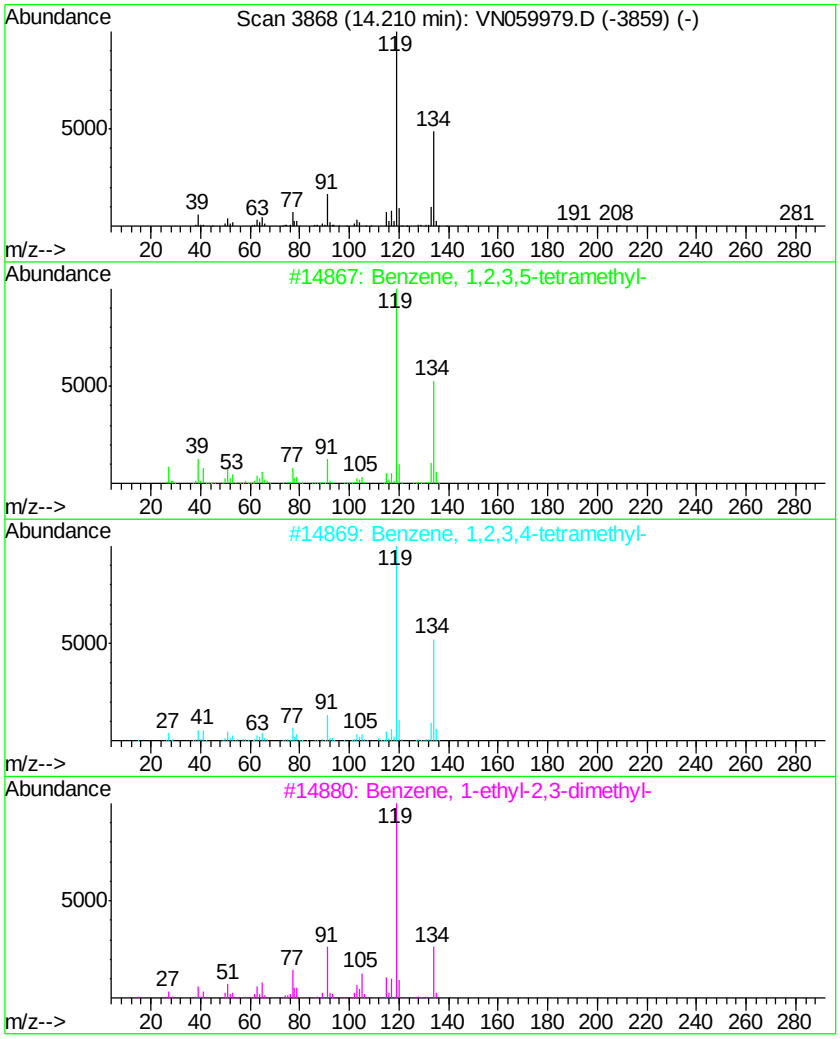
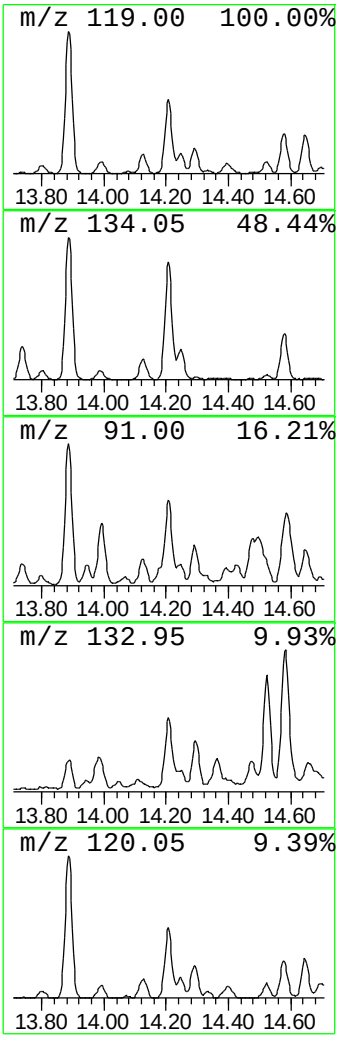
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	20.59 ug/l	419618	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





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Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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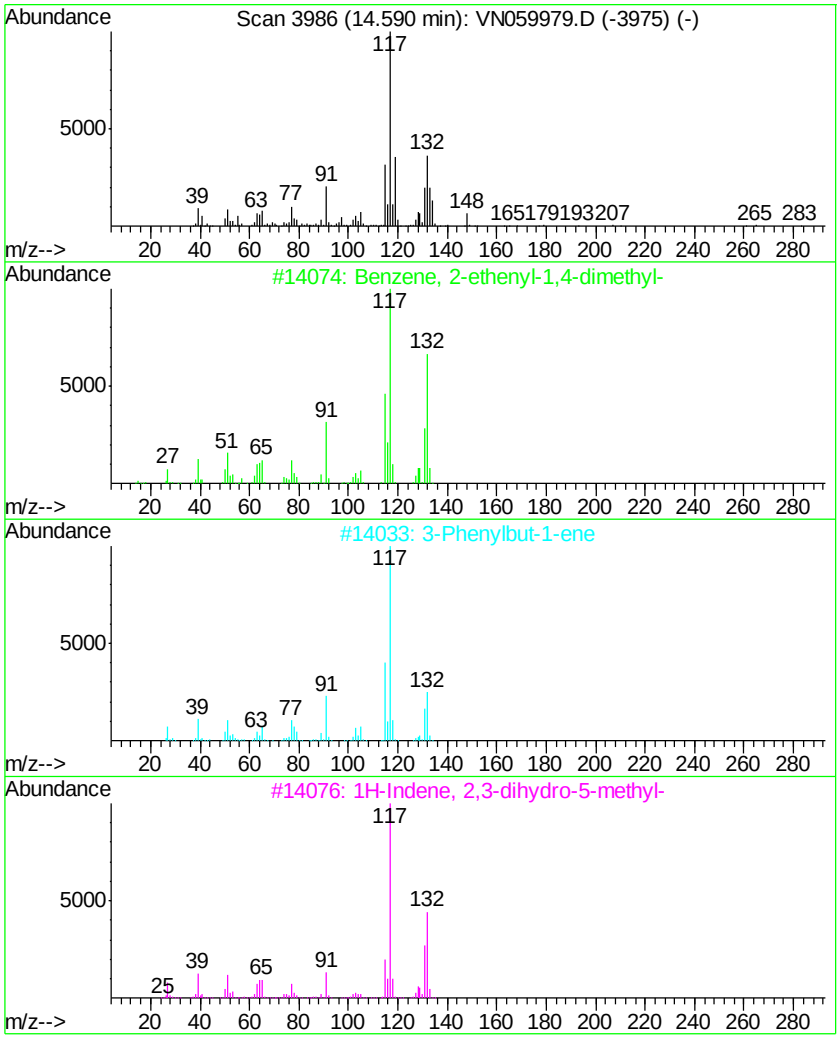
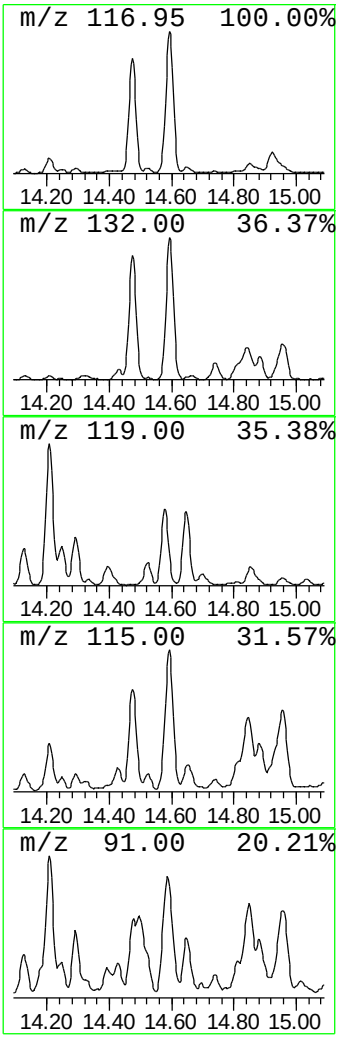
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	24.74 ug/l	504108	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN020320\
Data File : VN059979.D
Acq On : 3 Feb 2020 16:56
Operator : JC/MD
Sample : L1346-04 40X
Misc : 7.24g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : PD-3-(30-32)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Pentane, 2-methyl-	4.56	25.9	ug/l	172869	1	7.64	333228	50.0
Pentane, 3-methyl-	5.03	20.4	ug/l	135614	1	7.64	333228	50.0
Pentane, 2,3-dime...	7.72	31.1	ug/l	207357	1	7.64	333228	50.0
Hexane, 3-methyl-	7.86	56.9	ug/l	379371	1	7.64	333228	50.0
Heptane, 3-methyl-	9.86	25.2	ug/l	427779	2	8.57	847313	50.0
Octane	10.27	18.7	ug/l	335187	3	11.40	897828	50.0
Heptane, 2,4-dime...	11.19	25.3	ug/l	453579	3	11.40	897828	50.0
Benzene, 2-ethyl-...	13.89	27.2	ug/l	555070	4	13.34	1018810	50.0
Benzene, 1,2,3,5-...	14.21	20.6	ug/l	419618	4	13.34	1018810	50.0
Benzene, 2-etheny...	14.59	24.7	ug/l	504108	4	13.34	1018810	50.0