LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

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\82N060419W.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	2.793	306	317	332	rVB2	49585	99112	3.82%	0.376%
2	3.545	537	551	567	rVB5	14812	40886	1.58%	0.155%
3	3.783	609	625	651	rVB6	8277	32097	1.24%	0.122%
4	4.519	833	854	860	rBV4	35163	101937	3.93%	0.386%
5	5.047	990	1018	1041	rBV3	58111	211328	8.15%	0.801%
6	5.924	1270	1291	1308	rBV3	19433	59849	2.31%	0.227%
7	6.060	1314	1333	1352	rBV4	19444	62258	2.40%	0.236%
8	6.365	1409	1428	1443	rBV5	17460	56338	2.17%	0.214%
9	6.513	1459	1474	1489	rBV4	12201	36415	1.40%	0.138%
10	6.619	1490	1507	1525	rVV2	96507	299069	11.53%	1.134%
11	6.706	1527	1534	1550	rVB7	14408	36893	1.42%	0.140%
12	7.323	1708	1726	1736	rBV2	17747	47613	1.84%	0.180%
13	7.391	1736	1747	1762	rVB4	23414	62077	2.39%	0.235%
14	7.590	1784	1809	1820	rBV2	230078	602950	23.26%	2.285%
15	7.664	1820	1832	1847	rVB	390218	939331	36.23%	3.560%
16	7.870	1882	1896	1908	rBV3	23128	55450	2.14%	0.210%
17	8.027	1928	1945	1970	rBV	236889	567769	21.90%	2.152%
18	8.159	1970	1986	1994	rBV4	64741	163129	6.29%	0.618%
19	8.301	2022	2030	2050	rVB6	22372	54185	2.09%	0.205%
20	8.407	2053	2063	2076	rVV3	11568	26352	1.02%	0.100%
21	8.584	2102	2118	2140	rBV	677461	1540064	59.40%	5.838%
22	8.744	2159	2168	2179	rVB3	18303	33817	1.30%	0.128%
23	8.818	2179	2191	2199	rBV	48218	102314	3.95%	0.388%
24	9.043	2241	2261	2266	rBV3	52167	104901	4.05%	0.398%
25	9.268	2322	2331	2344	rVB3	16410	33870	1.31%	0.128%
26	9.362	2345	2360	2373	rBV6	12812	35847	1.38%	0.136%
27	9.516	2399	2408	2420	rVB4	35344	69050	2.66%	0.262%
28	9.612	2427	2438	2444	rBV3	47373	91251	3.52%	0.346%
29	9.767	2480	2486	2498	rVB7	12965	26703	1.03%	0.101%
30	9.854	2498	2513	2524	rBV7	15305	42705	1.65%	0.162%
31	9.966	2538	2548	2558	rBV2	52342	97819	3.77%	0.371%
32	10.092	2572	2587	2615	rBV	969090	1928652	74.39%	7.310%
33	10.516	2707	2719	2735	rVV3	80335	172886	6.67%	0.655%
34	10.792	2790	2805	2812	rBV7	21269	44312	1.71%	0.168%

LSC Area Percent Report

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Sample : K3345-13
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Title : SW846 8260

35	11.407	2982	2996	3018	rVV	904700	1687836	65.10%	6.398%
36	11.509	3020	3028	3046	rVB2	49275	110682	4.27%	0.420%
37	12.249	3245	3258	3278	rVB	1584003	2592764	100.00%	9.828%
38	12.400	3292	3305	3322	rBV	720703	1255071	48.41%	4.757%
39	12.590	3338	3364	3384	rVB	546191	930376	35.88%	3.527%
40	12.889	3442	3457	3470	rBV2	38855	69631	2.69%	0.264%
41	12.992	3482	3489	3497	rVV	20248	30528	1.18%	0.116%
42	13.040	3497	3504	3516	rVB	22338	36850	1.42%	0.140%
43	13.169	3530	3544	3557	rBV4	139852	280115	10.80%	1.062%
44	13.342	3586	3598	3619	rVB2	762120	1524111	58.78%	5.777%
45	13.442	3620	3629	3638	rVB	42682	66582	2.57%	0.252%
46	13.542	3639	3660	3671	rBV	952207	1827126	70.47%	6.926%
47	13.593	3671	3676	3690	rVV2	61204	144573	5.58%	0.548%
48	13.670	3690	3700	3711	rVB2	84003	142074	5.48%	0.539%
49	13.886	3757	3767	3777	rBV	282516	438523	16.91%	1.662%
50	13.947	3777	3786	3792	rVV	137233	219540	8.47%	0.832%
51	13.995	3792	3801	3813	rVV2	518011	858899	33.13%	3.256%
52	14.066	3816	3823	3831	rVB6	22432	37290	1.44%	0.141%
53	14.130	3832	3843	3851	rBV3	79439	130268	5.02%	0.494%
54	14.210	3851	3868	3885	rVB	245494	450767	17.39%	1.709%
55	14.291	3886	3893	3899	rBV2	27882	42160	1.63%	0.160%
56	14.361	3909	3915	3922	rVB	23356	30451	1.17%	0.115%
57	14.426	3923	3935	3942	rBV3	48830	89491	3.45%	0.339%
58	14.474	3942	3950	3961	rVV2	197963	332009	12.81%	1.258%
59	14.586	3972	3985	3998	rBV3	848438	1771975	68.34%	6.717%
60	14.657	3998	4007	4016	rVB6	44920	75795	2.92%	0.287%
61	14.738	4024	4032	4041	rVB2	97136	143481	5.53%	0.544%
62	14.847	4048	4066	4072	rBV4	111878	253451	9.78%	0.961%
63	14.886	4073	4078	4087	rVV2	95652	163974	6.32%	0.622%
64	14.956	4088	4100	4112	rVB2	167176	387626	14.95%	1.469%
65	15.046	4119	4128	4134	rBV4	23356	40491	1.56%	0.153%
66	15.088	4134	4141	4147	rVV	41389	71224	2.75%	0.270%
67	15.133	4148	4155	4163	rVV	72965	122063	4.71%	0.463%
68	15.188	4163	4172	4180	rVV2	83030	139084	5.36%	0.527%
69	15.259	4180	4194	4229	rVB5	86032	344007	13.27%	1.304%
70	15.413	4230	4242	4254	rBV3	44411	85685	3.30%	0.325%
71	15.484	4254	4264	4275	rBV7	26763	49115	1.89%	0.186%

LSC Area Percent Report

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Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Title : SW846 8260

72	15.574	4277	4292	4296	rBV4	61599	118693	4.58%	0.450%
73	15.612	4298	4304	4320	rVB2	129766	232105	8.95%	0.880%
74	15.699	4320	4331	4341	rBV2	34070	66328	2.56%	0.251%
75	15.767	4342	4352	4362	rBV4	25289	51847	2.00%	0.197%
76	15.911	4383	4397	4410	rBV3	132095	272861	10.52%	1.034%
77	16.101	4440	4456	4464	rBV5	26781	62616	2.42%	0.237%
78	16.159	4464	4474	4506	rVB2	78225	195837	7.55%	0.742%
79	16.352	4517	4534	4547	rBV	233285	526838	20.32%	1.997%

Sum of corrected areas: 26382042

LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
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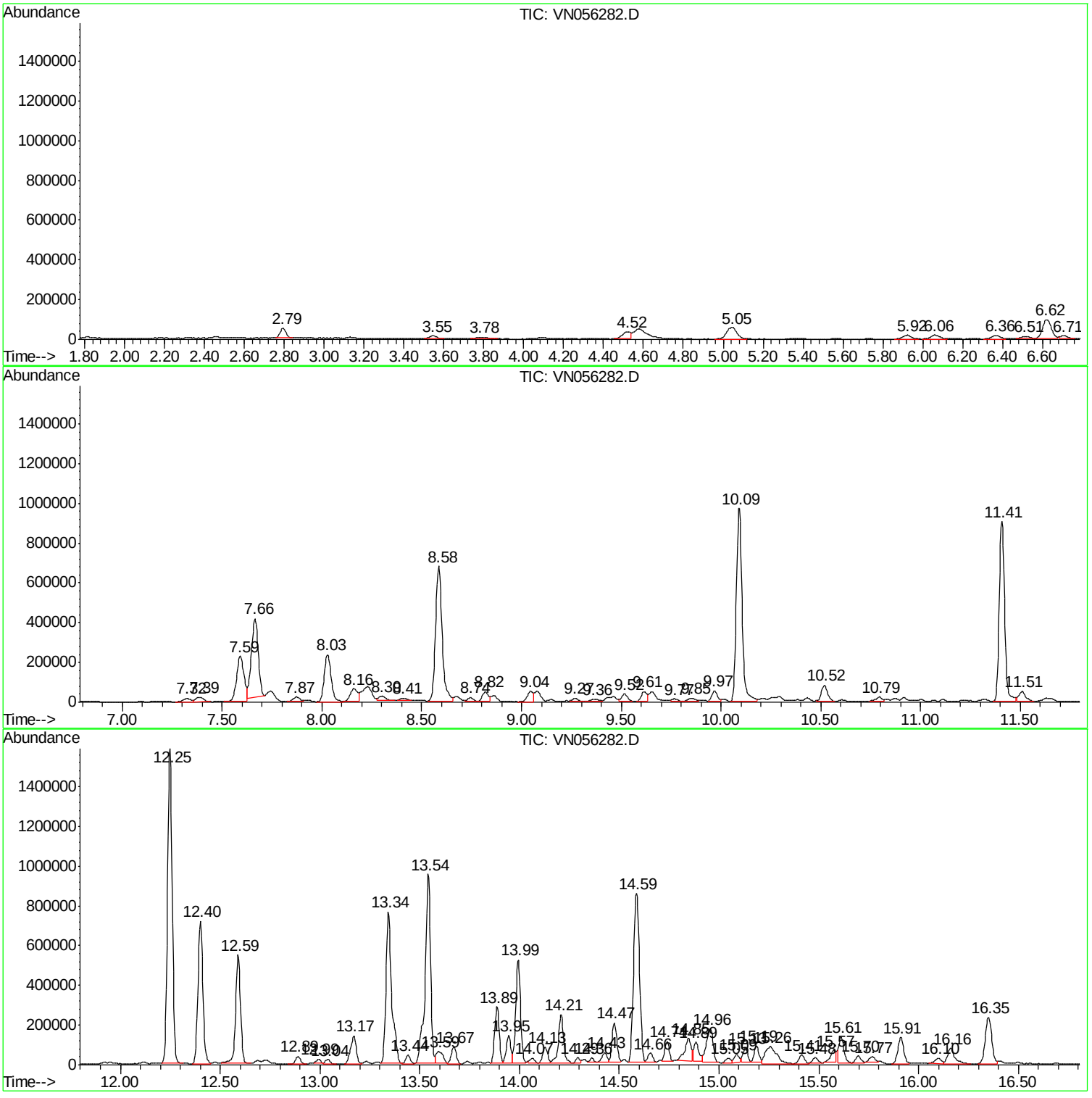
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

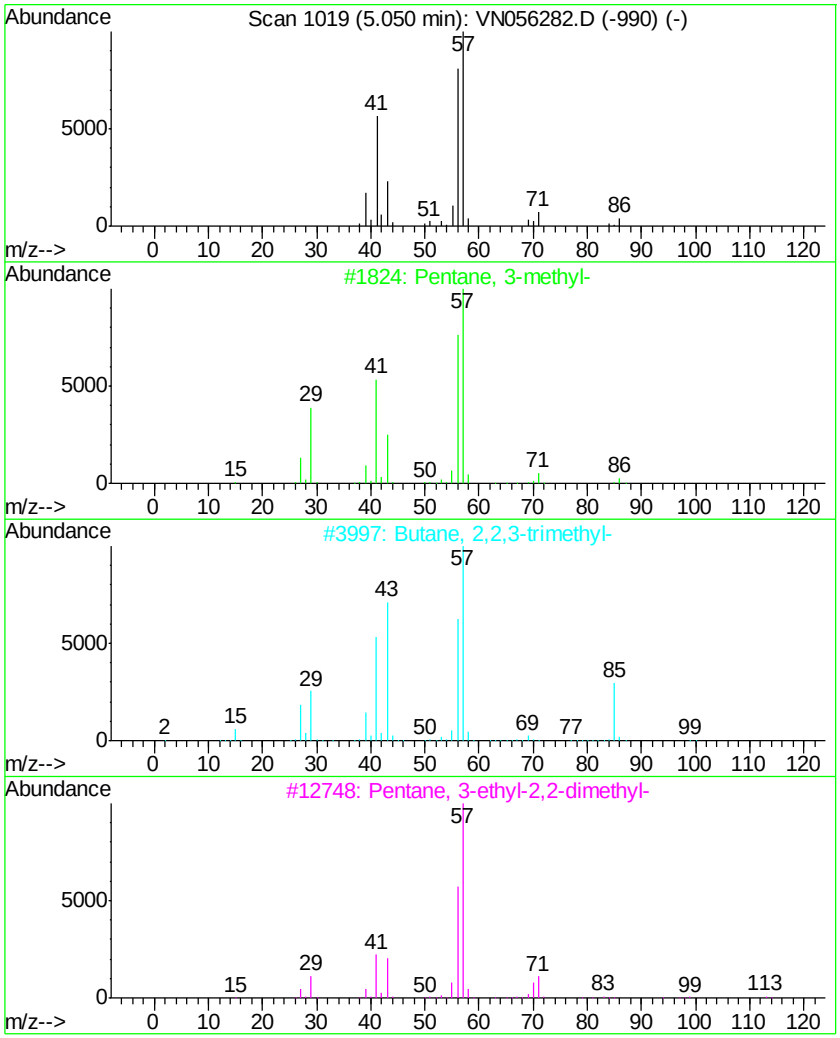
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

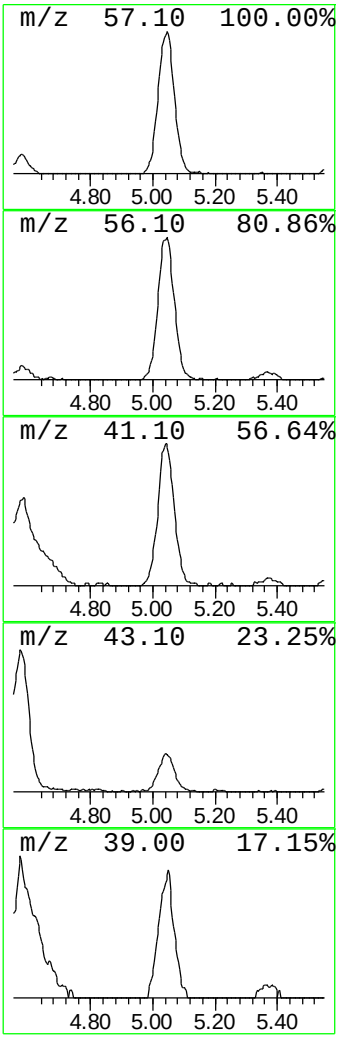
****

Peak Number 1 Pentane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	11.25 ug/l	211328	Pentafluorobenzene	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	78
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
5	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	56





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
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Operator : JC/SP
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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

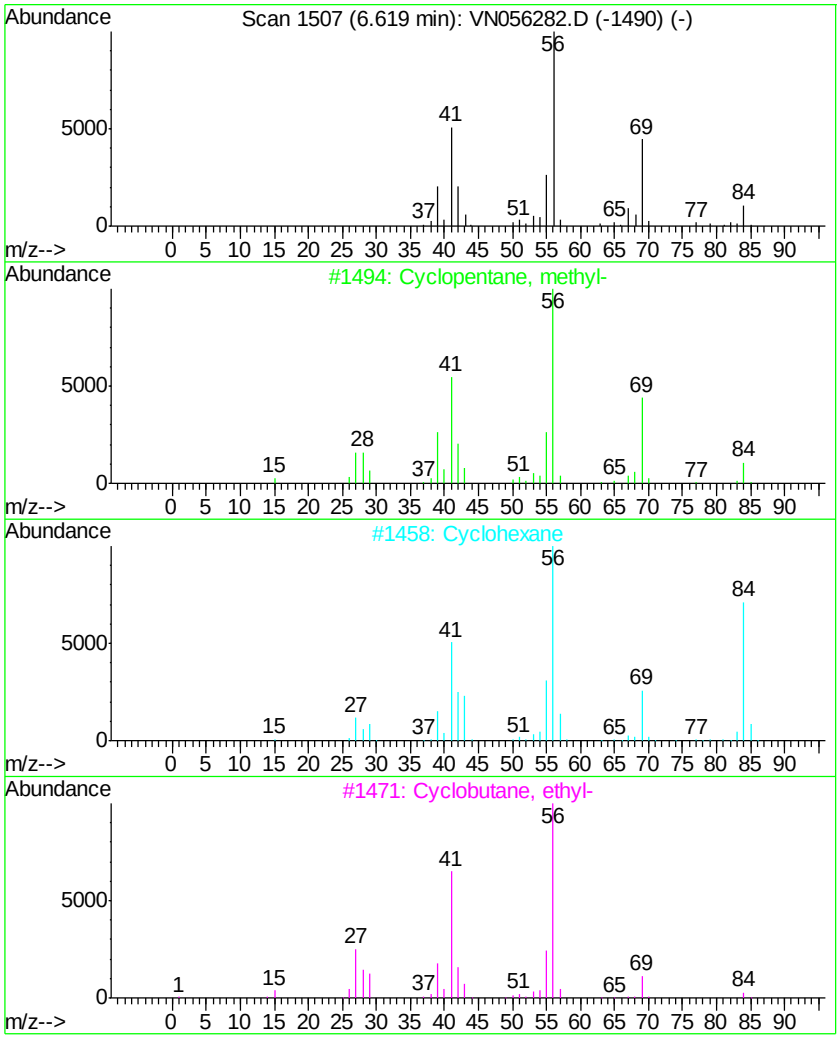
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

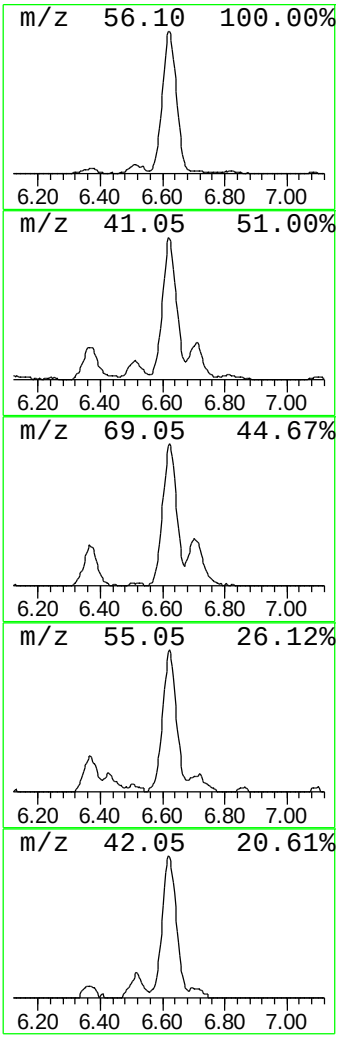
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	15.92 ug/l	299069	Pentafluorobenzene	7.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclohexane	84	C6H12	000110-82-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4	Cyclopropane, propyl-	84	C6H12	002415-72-7	53
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

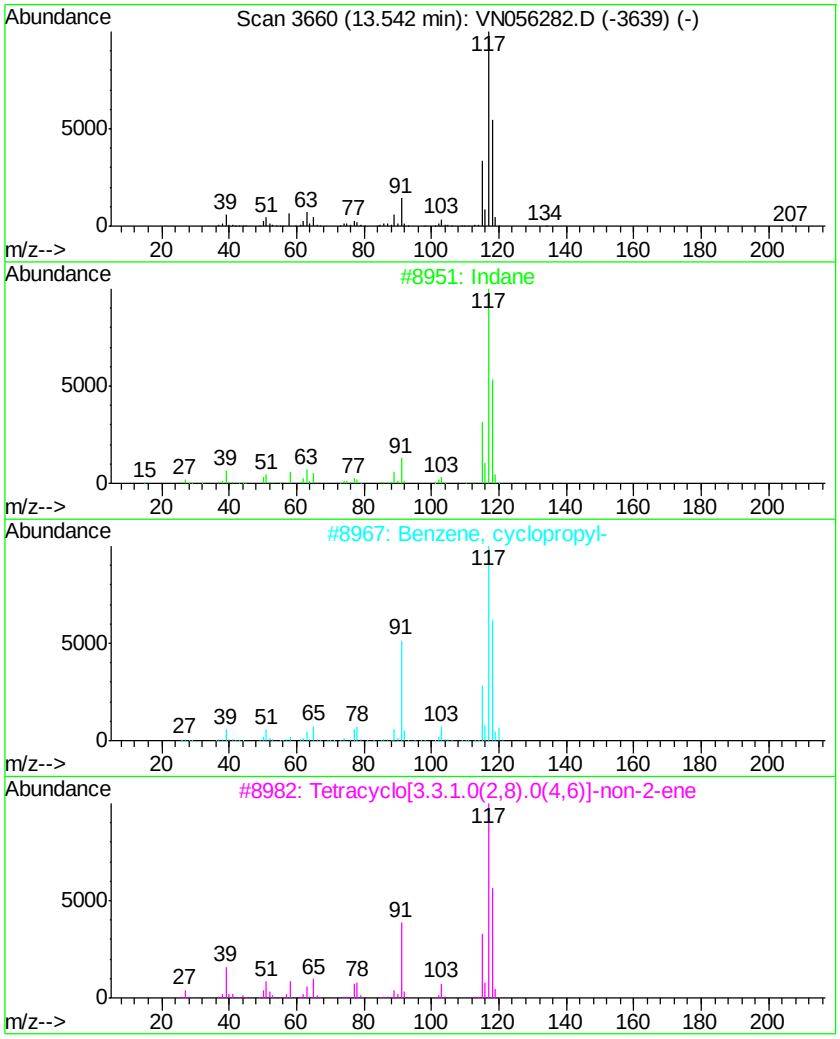
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

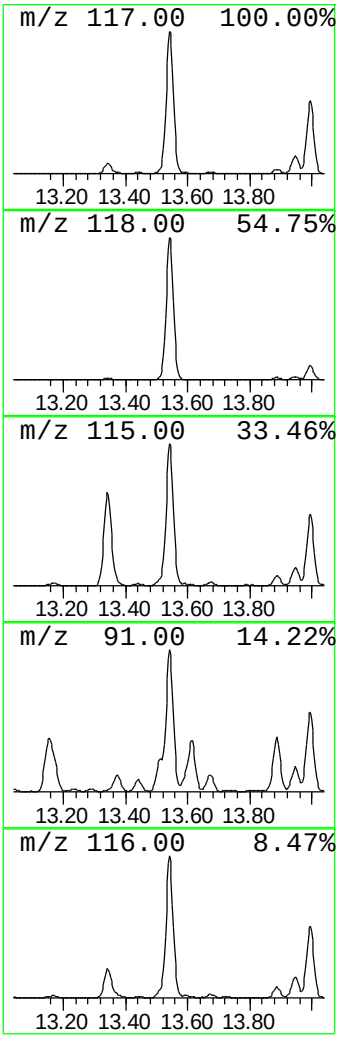
****

Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	59.94 ug/l	1827130	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64





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ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

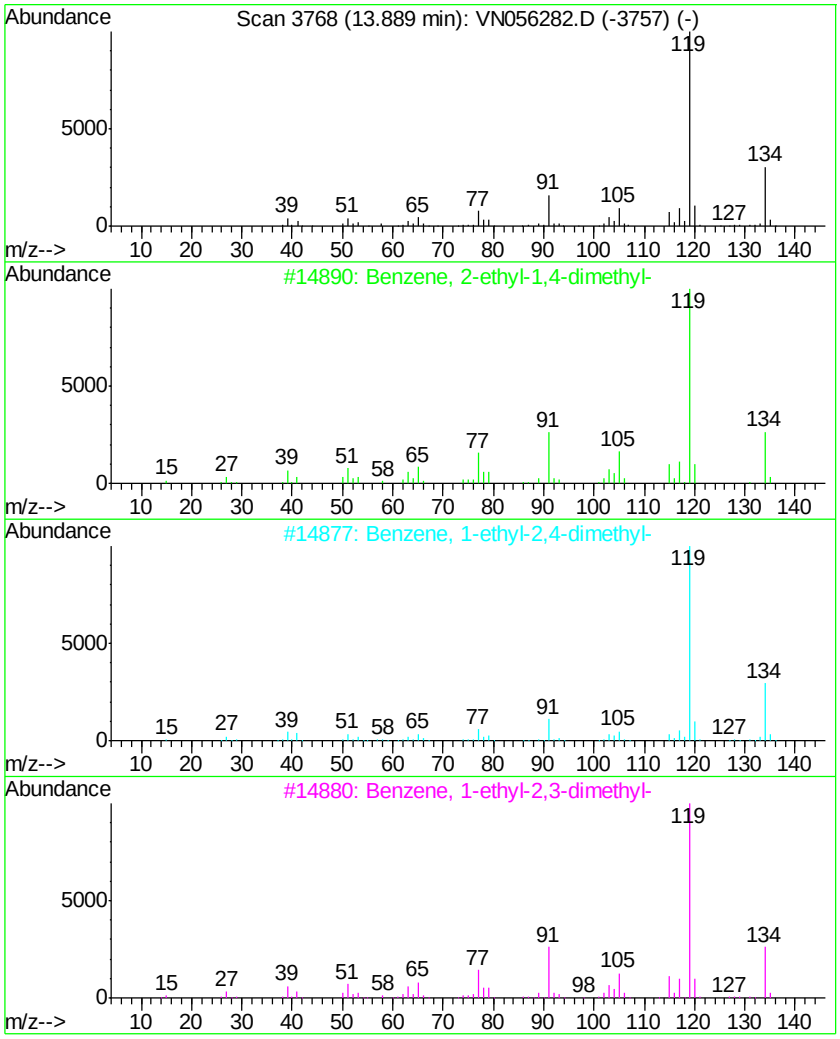
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

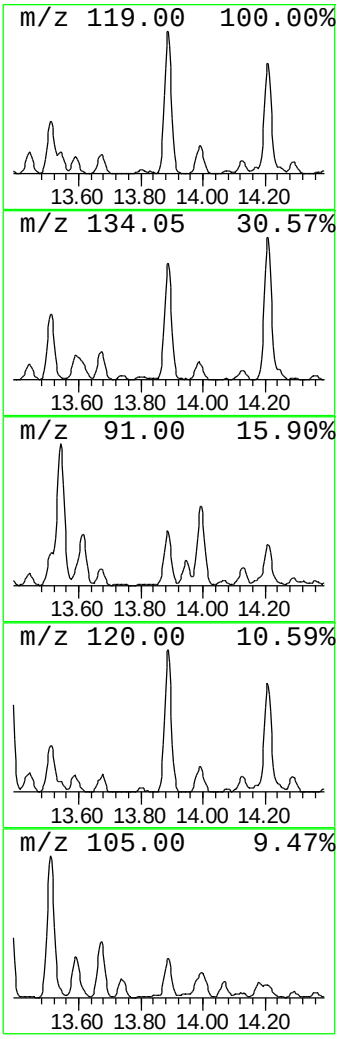
****

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	14.39 ug/l	438523	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	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	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





Library Search Compound Report

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Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

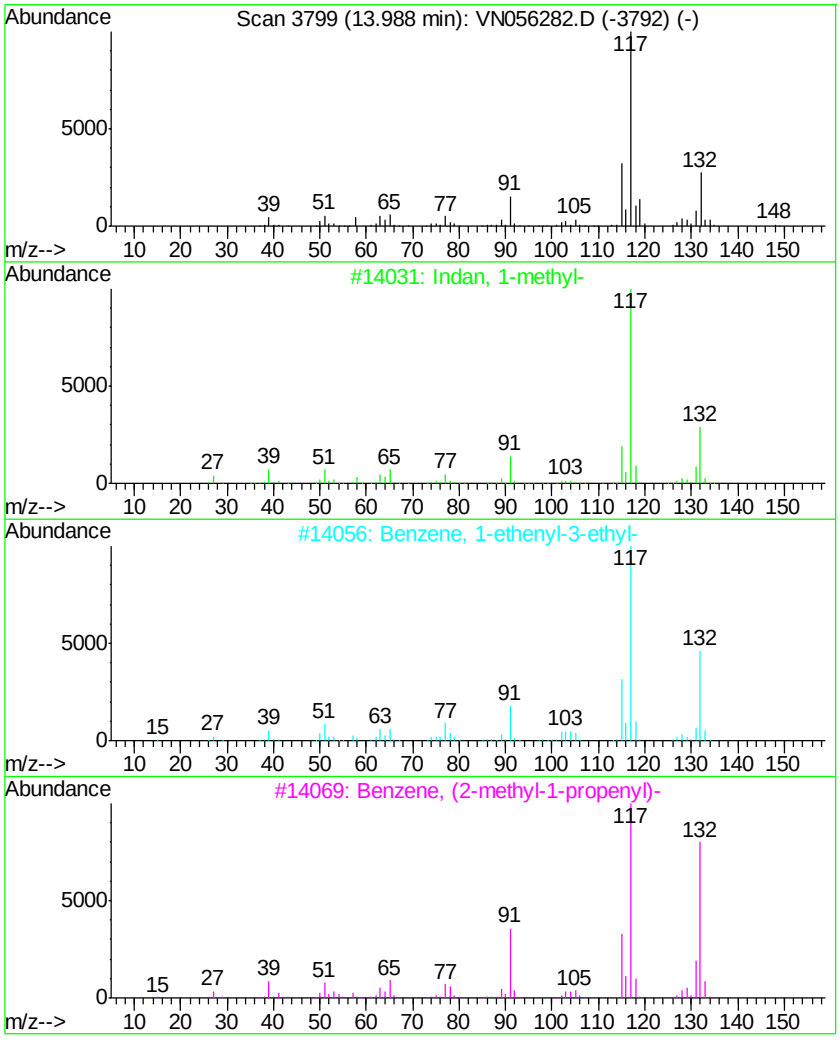
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

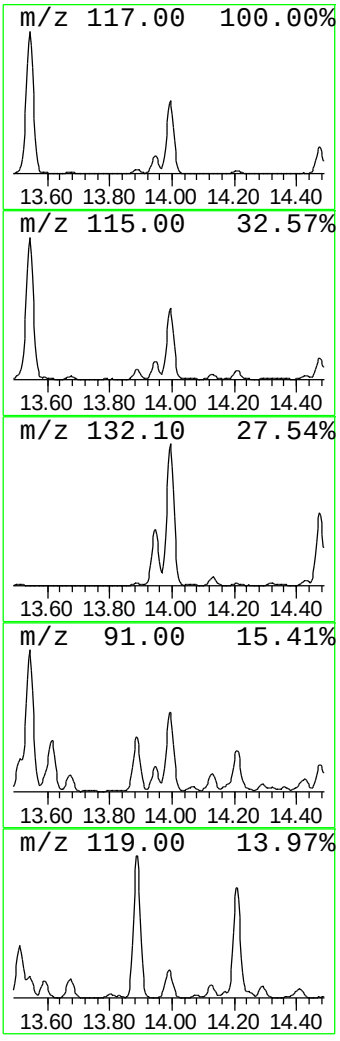
****

Peak Number 5 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	28.18 ug/l	858899	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

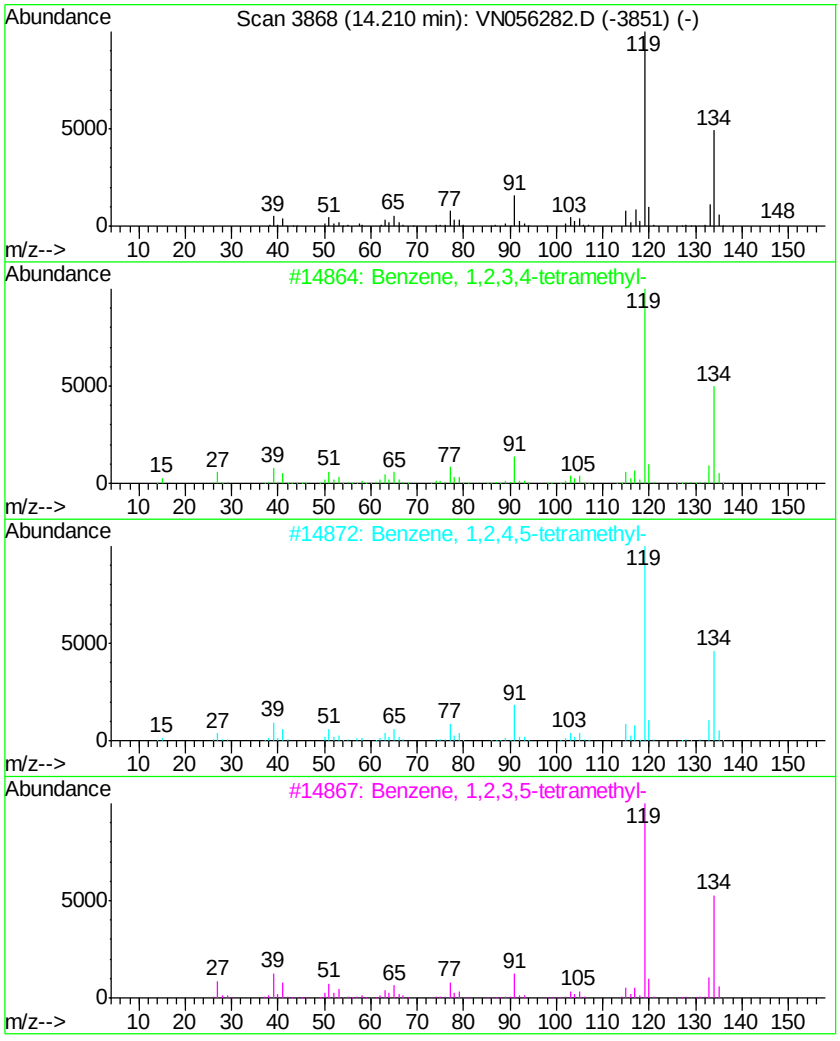
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

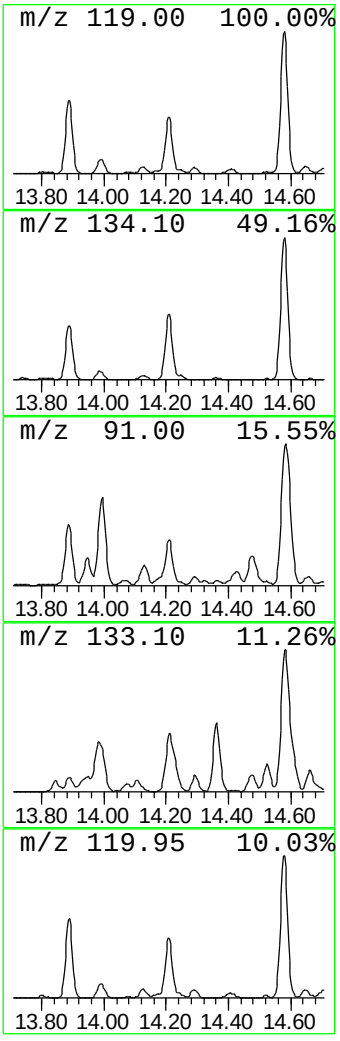
****

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	14.79 ug/l	450767	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





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Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

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

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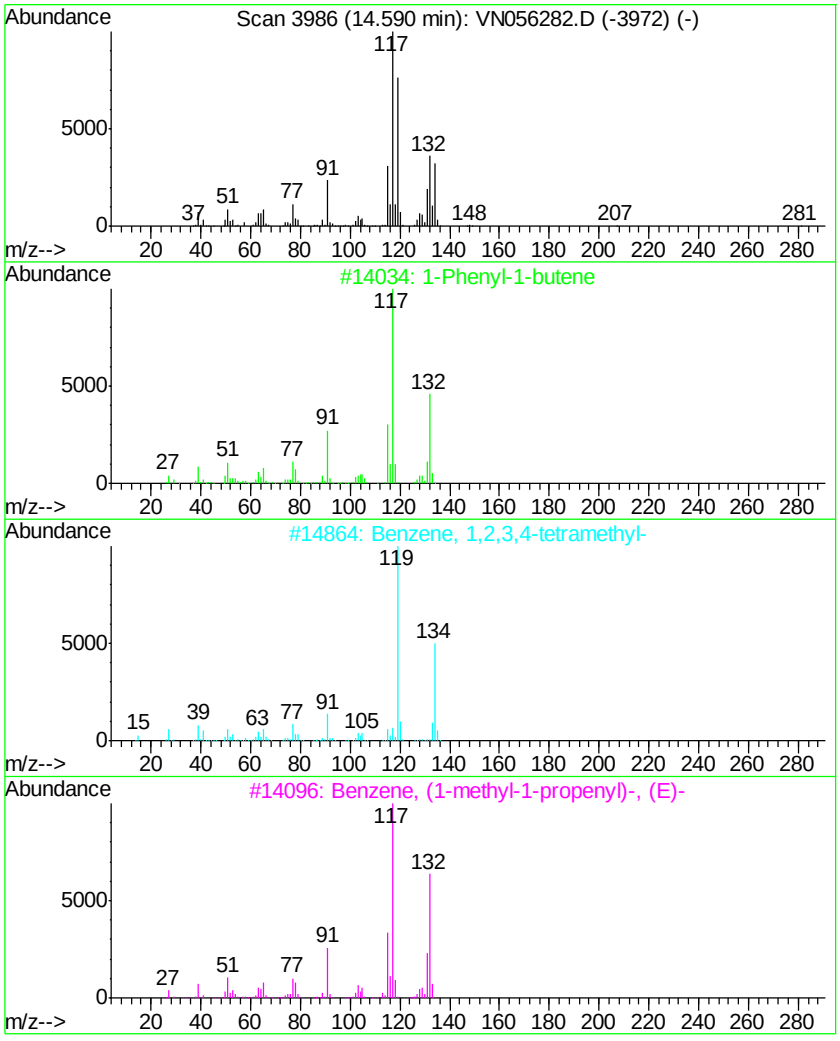
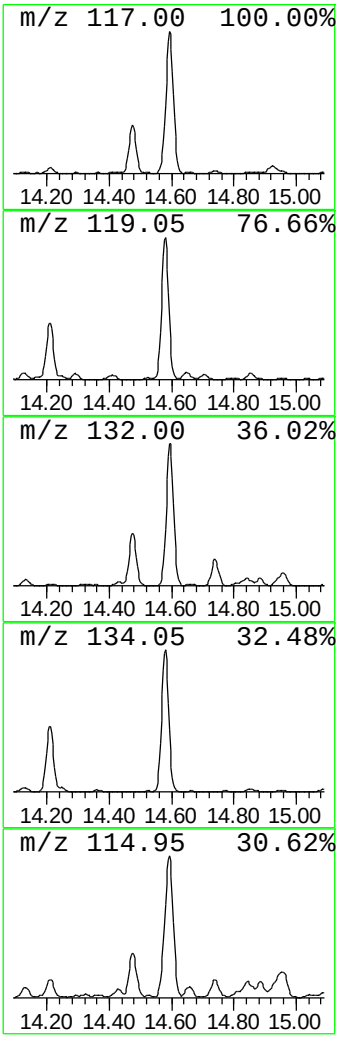
Peak Number 7 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	58.13 ug/l	1771980	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	50
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

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

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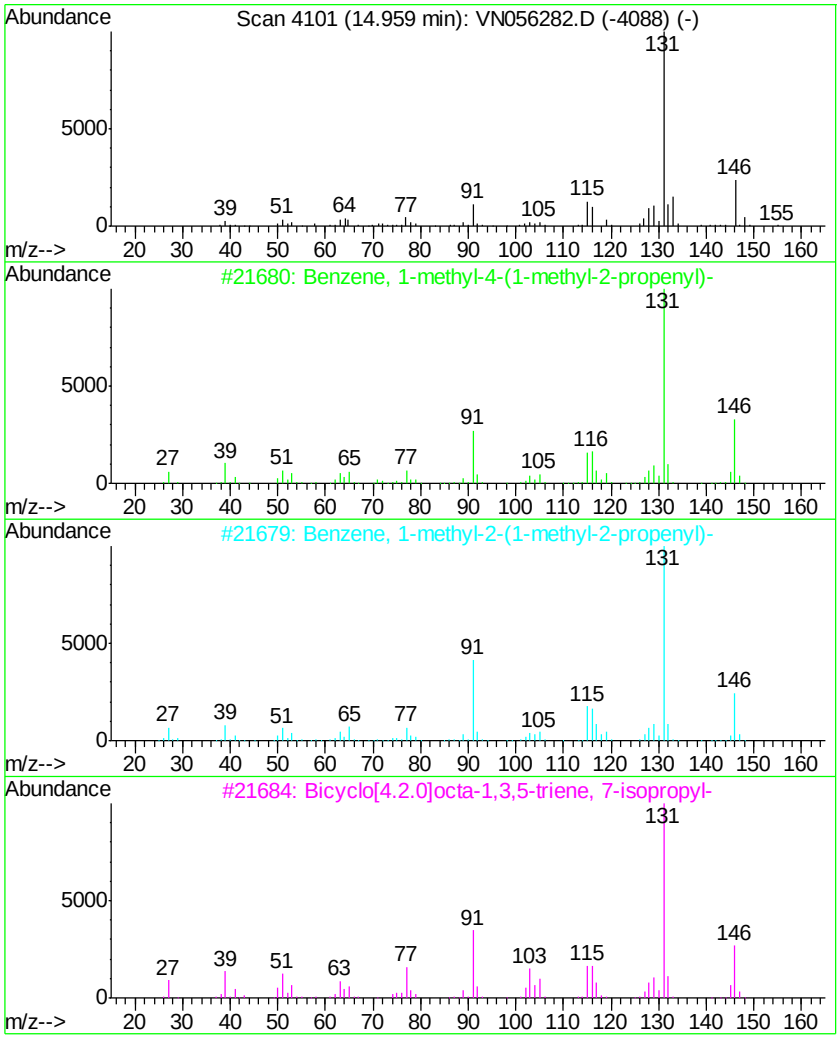
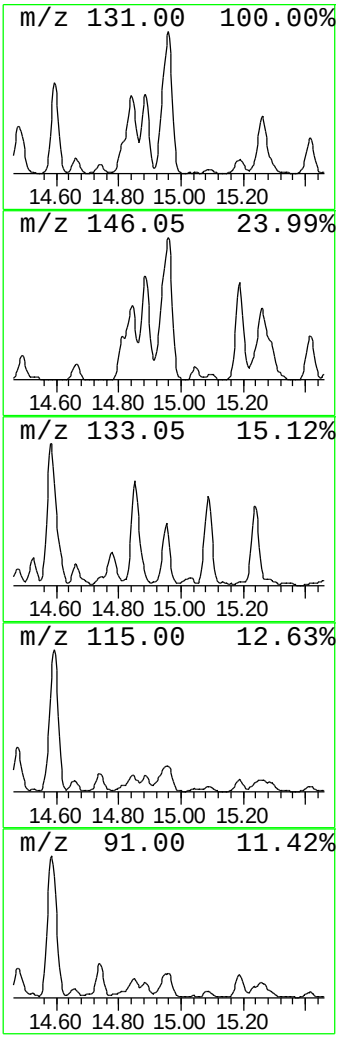
Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	12.72 ug/l	387626	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
2	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	87
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	87
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87





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Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

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

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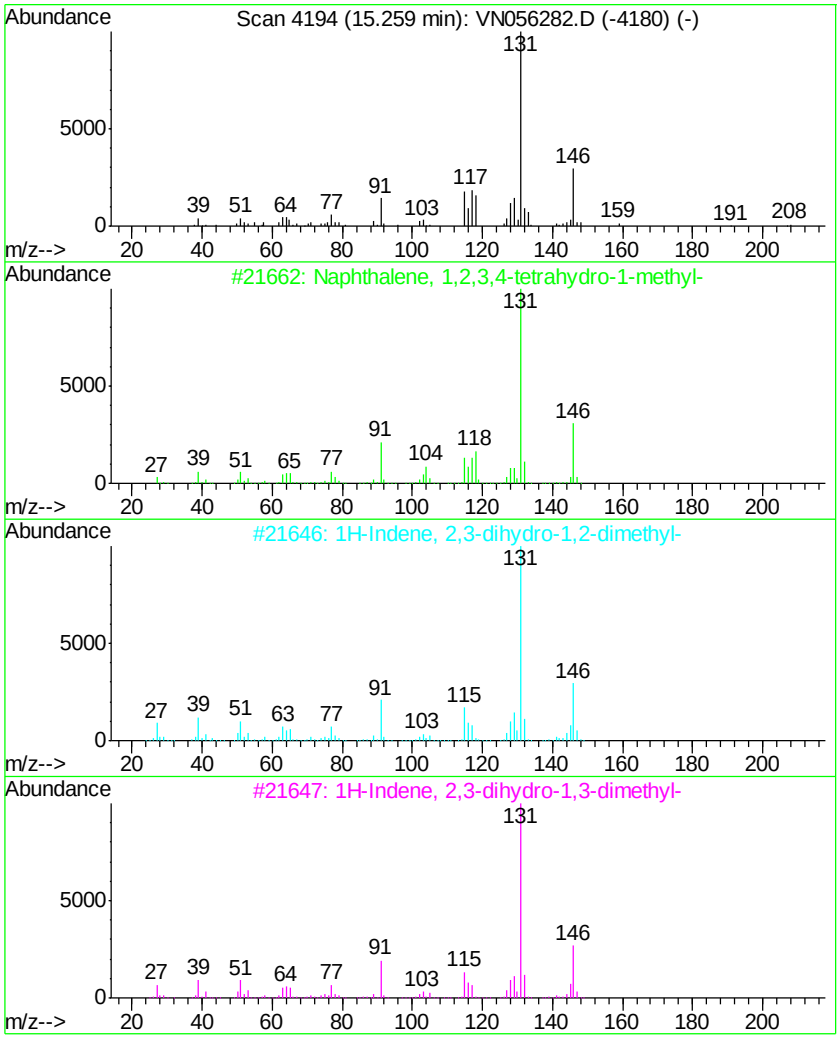
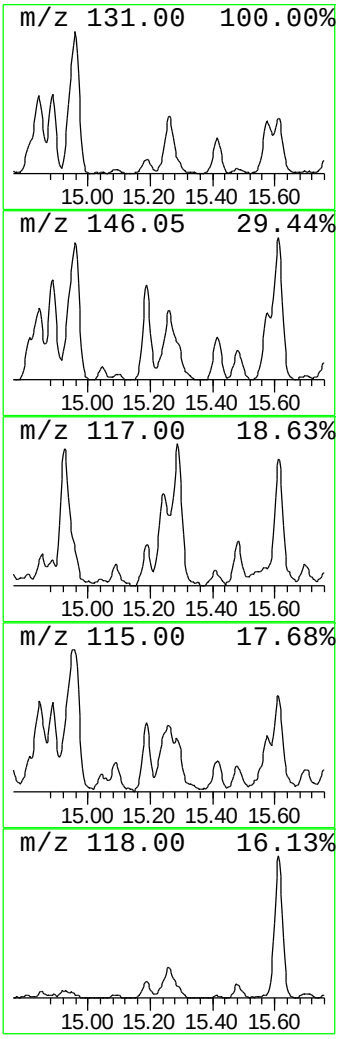
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	11.29 ug/l	344007	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	83
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260

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

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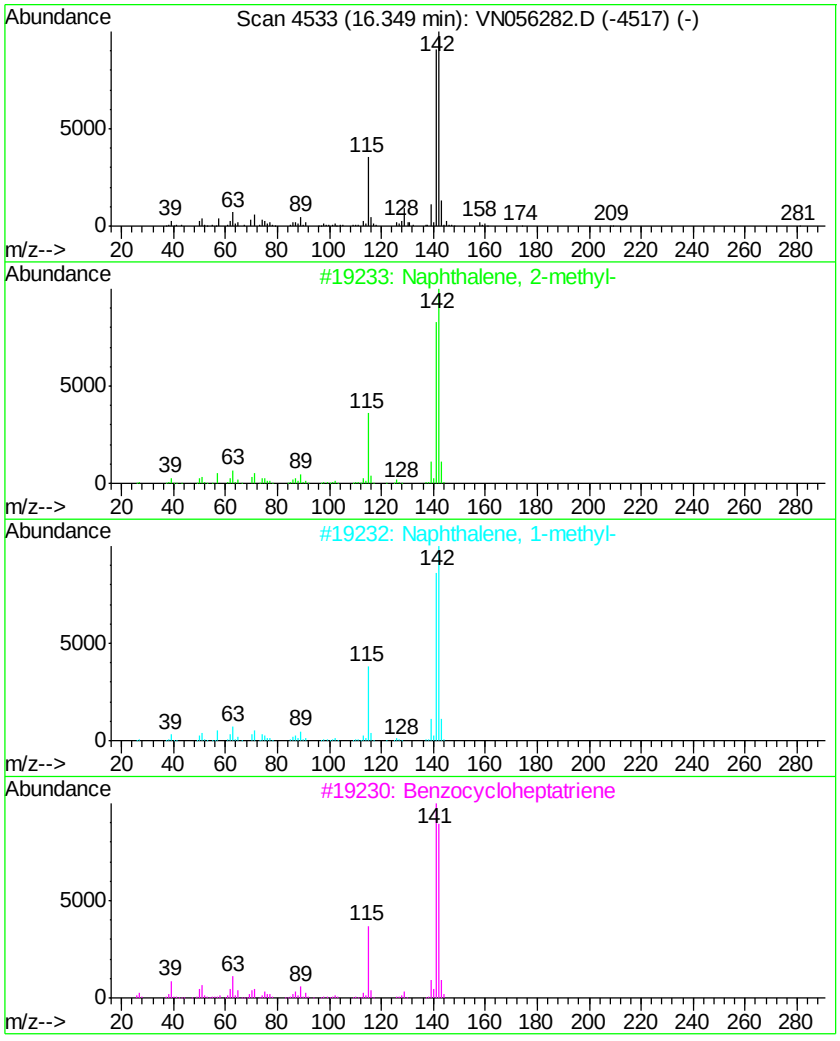
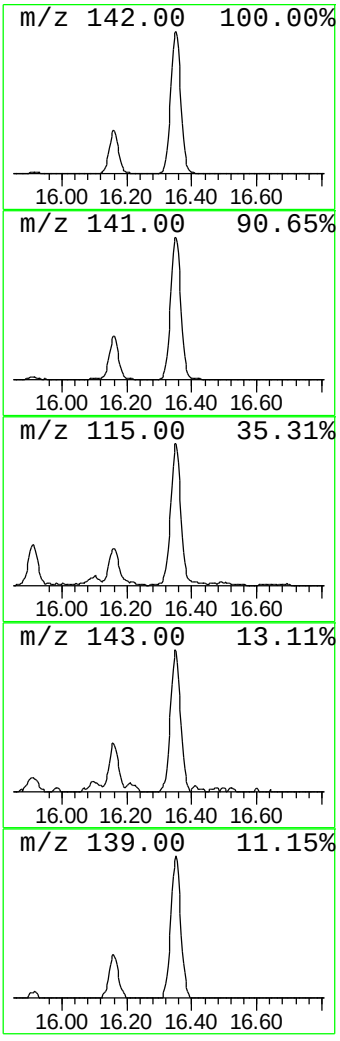
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.35	17.28 ug/l	526838	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN061319\
Data File : VN056282.D
Acq On : 13 Jun 2019 15:31
Operator : JC/SP
Sample : K3345-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : DUP-01_061219

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.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, 3-methyl-	5.05	11.3	ug/l	211328	1	7.66	939331	50.0
Cyclopentane, met...	6.62	15.9	ug/l	299069	1	7.66	939331	50.0
Indane	13.54	59.9	ug/l	1827130	4	13.34	1524110	50.0
Benzene, 2-ethyl-...	13.89	14.4	ug/l	438523	4	13.34	1524110	50.0
Indan, 1-methyl-	13.99	28.2	ug/l	858899	4	13.34	1524110	50.0
Benzene, 1,2,3,4-...	14.21	14.8	ug/l	450767	4	13.34	1524110	50.0
1-Phenyl-1-butene	14.59	58.1	ug/l	1771980	4	13.34	1524110	50.0
Benzene, 1-methyl...	14.96	12.7	ug/l	387626	4	13.34	1524110	50.0
Naphthalene, 1,2,...	15.26	11.3	ug/l	344007	4	13.34	1524110	50.0
Naphthalene, 1-me...	16.35	17.3	ug/l	526838	4	13.34	1524110	50.0