LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

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\82N110921W.M
Title : SW846 8260

Signal : TIC: VN069560.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.311	108	120	137	rVV	54025	167867	1.06%	0.064%
2	2.443	160	169	188	rVB	164477	283149	1.78%	0.108%
3	3.140	410	429	461	rBV	1968922	4471970	28.15%	1.713%
4	3.513	550	568	597	rVB	573304	1391965	8.76%	0.533%
5	3.990	725	746	769	rBV3	158404	433199	2.73%	0.166%
6	4.253	817	844	886	rBV4	236157	896832	5.65%	0.343%
7	5.079	1122	1152	1159	rBV3	1108937	3369607	21.21%	1.291%
8	5.618	1319	1353	1400	rBV	2107323	7198666	45.32%	2.757%
9	5.935	1446	1471	1508	rVB4	106297	349497	2.20%	0.134%
10	6.122	1519	1541	1567	rBV5	61223	183458	1.15%	0.070%
11	6.468	1639	1670	1697	rBV3	495017	1506337	9.48%	0.577%
12	6.611	1699	1723	1744	rBV4	333242	1057900	6.66%	0.405%
13	6.903	1808	1832	1864	rBV4	419228	1309027	8.24%	0.501%
14	7.048	1865	1886	1902	rBV2	476123	1393953	8.78%	0.534%
15	7.147	1902	1923	1944	rBV	3384514	9646134	60.72%	3.694%
16	7.327	1977	1990	2021	rVB9	73619	266411	1.68%	0.102%
17	7.670	2095	2118	2139	rVB4	84710	300005	1.89%	0.115%
18	7.783	2139	2160	2169	rBV5	165529	437092	2.75%	0.167%
19	7.847	2171	2184	2205	rVB3	463134	1176475	7.41%	0.451%
20	8.024	2227	2250	2258	rBV	978673	2401122	15.12%	0.920%
21	8.091	2260	2275	2292	rVV4	4008143	10995053	69.22%	4.211%
22	8.177	2293	2307	2332	rVB2	1954871	5137719	32.34%	1.968%
23	8.298	2333	2352	2366	rBV2	1175005	2742665	17.27%	1.050%
24	8.445	2389	2407	2435	rBV3	1381046	4061498	25.57%	1.556%
25	8.574	2436	2455	2463	rBV2	1289854	3105720	19.55%	1.189%
26	8.708	2494	2505	2526	rVB2	1050809	2451466	15.43%	0.939%
27	8.802	2527	2540	2571	rVB4	257002	663722	4.18%	0.254%
28	8.971	2579	2603	2630	rBV3	4185082	10696580	67.34%	4.097%
29	9.126	2650	2661	2675	rVB3	297111	559653	3.52%	0.214%
30	9.201	2675	2689	2696	rVV3	409367	803504	5.06%	0.308%
31	9.247	2698	2706	2729	rVB2	585802	1184402	7.46%	0.454%
32	9.462	2752	2786	2801	rBV2	3192605	9257297	58.28%	3.546%
33	9.596	2824	2836	2841	rBV2	96840	169968	1.07%	0.065%
34	9.641	2841	2853	2868	rVB2	618688	1192626	7.51%	0.457%
35	9.735	2869	2888	2901	rBV5	170967	397875	2.50%	0.152%
36	9.805	2901	2914	2918	rBV3	274055	471242	2.97%	0.180%

LSC Area Percent Report

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Instrument :
MSVOA_N
ClientSampleId :
LB-3R

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\82N110921W.M
Title : SW846 8260

37	9.883	2932	2943	2961	rVB	878416	1801088	11.34%	0.690%
38	9.974	2962	2977	2985	rBV	1719243	3319534	20.90%	1.271%
39	10.017	2987	2993	3009	rVB2	1383143	2517542	15.85%	0.964%
40	10.210	3049	3065	3093	rBV10	249536	766078	4.82%	0.293%
41	10.322	3094	3107	3118	rBV2	761300	1431923	9.01%	0.548%
42	10.443	3135	3152	3167	rBV	5002085	9686476	60.98%	3.710%
43	10.636	3205	3224	3237	rVB6	546114	1279289	8.05%	0.490%
44	10.784	3268	3279	3293	rBV6	247523	508198	3.20%	0.195%
45	10.862	3294	3308	3330	rVB5	1190912	2752151	17.33%	1.054%
46	10.950	3331	3341	3358	rVB5	118971	276453	1.74%	0.106%
47	11.039	3358	3374	3385	rBV5	83348	209871	1.32%	0.080%
48	11.130	3395	3408	3417	rBV3	227373	436011	2.74%	0.167%
49	11.173	3418	3424	3432	rVV2	161190	213626	1.34%	0.082%
50	11.216	3433	3440	3446	rVV3	143550	194441	1.22%	0.074%
51	11.256	3447	3455	3464	rVB4	301982	457107	2.88%	0.175%
52	11.454	3520	3529	3543	rVB6	147163	260862	1.64%	0.100%
53	11.556	3550	3567	3584	rBV7	112245	289079	1.82%	0.111%
54	11.747	3622	3638	3653	rBV	4639914	8378726	52.75%	3.209%
55	11.846	3664	3675	3696	rVB	1180149	2225744	14.01%	0.852%
56	11.956	3702	3716	3741	rVB	614231	1403445	8.83%	0.538%
57	12.280	3826	3837	3857	rVB	250875	453995	2.86%	0.174%
58	12.449	3875	3900	3912	rBV8	90459	255953	1.61%	0.098%
59	12.581	3934	3949	3968	rBV	3322457	5463423	34.39%	2.093%
60	12.731	3990	4005	4026	rBV	3567062	6211777	39.10%	2.379%
61	12.921	4060	4076	4093	rBV	9767367	15885171	100.00%	6.084%
62	13.015	4103	4111	4121	rVB	287128	422681	2.66%	0.162%
63	13.222	4172	4188	4207	rBV	2315277	3838861	24.17%	1.470%
64	13.369	4233	4243	4256	rVB	253330	413564	2.60%	0.158%
65	13.498	4272	4291	4305	rBV3	973069	2380702	14.99%	0.912%
66	13.560	4305	4314	4325	rVB	227415	344298	2.17%	0.132%
67	13.616	4325	4335	4343	rBV	156174	234834	1.48%	0.090%
68	13.675	4343	4357	4381	rVV2	4108418	7717897	48.59%	2.956%
69	13.771	4383	4393	4404	rVB2	153432	243599	1.53%	0.093%
70	13.838	4404	4418	4424	rBV	2739228	4358718	27.44%	1.669%
71	13.879	4424	4433	4443	rVV	6502345	11079693	69.75%	4.244%
72	13.922	4444	4449	4469	rVV3	1333538	3231294	20.34%	1.238%
73	14.002	4469	4479	4492	rVB2	604904	981004	6.18%	0.376%
74	14.069	4492	4504	4516	rBV	324178	513211	3.23%	0.197%
75	14.131	4516	4527	4543	rVB	287755	450685	2.84%	0.173%
76	14.217	4544	4559	4572	rBV	6560251	10489397	66.03%	4.017%

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Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
LB-3R

Integration Parameters: RTEINT.P
Integrator: RTE
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Title : SW846 8260

77	14.281	4572	4583	4590	rVV	1235212	2024893	12.75%	0.776%
78	14.329	4590	4601	4616	rVB2	4895866	8324837	52.41%	3.188%
79	14.396	4616	4626	4639	rVB6	141354	265514	1.67%	0.102%
80	14.466	4641	4652	4661	rVB2	211397	320817	2.02%	0.123%
81	14.538	4662	4679	4701	rBV	4107505	6890115	43.37%	2.639%
82	14.622	4701	4710	4718	rBV	345694	464548	2.92%	0.178%
83	14.764	4744	4763	4769	rBV3	335152	597529	3.76%	0.229%
84	14.809	4769	4780	4792	rVV	3148505	5032394	31.68%	1.927%
85	14.930	4806	4825	4838	rBV3	5877079	12276885	77.29%	4.702%
86	14.984	4839	4845	4863	rVB3	425752	886443	5.58%	0.340%
87	15.086	4871	4883	4898	rBV2	536800	996685	6.27%	0.382%
88	15.196	4900	4924	4933	rBV4	1637935	3976084	25.03%	1.523%
89	15.311	4952	4967	4990	rVB3	1582285	3816123	24.02%	1.462%
90	15.458	5012	5022	5026	rBV2	103247	164909	1.04%	0.063%
91	15.507	5028	5040	5053	rVB	1192972	2123671	13.37%	0.813%
92	15.563	5053	5061	5070	rBV3	175144	257857	1.62%	0.099%
93	15.617	5070	5081	5090	rBV3	384124	754389	4.75%	0.289%
94	15.804	5136	5151	5167	rBV	701451	1372288	8.64%	0.526%
95	15.981	5198	5217	5253	rVB5	462088	1561392	9.83%	0.598%
96	16.110	5254	5265	5277	rBV3	187666	372146	2.34%	0.143%
97	16.191	5282	5295	5327	rVB4	390507	1045822	6.58%	0.401%
98	16.346	5336	5353	5380	rBV4	283362	703044	4.43%	0.269%
99	16.550	5407	5429	5440	rBV5	127240	333082	2.10%	0.128%
100	16.617	5441	5454	5491	rVB3	389574	1023388	6.44%	0.392%

Sum of corrected areas: 261094912

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Instrument :
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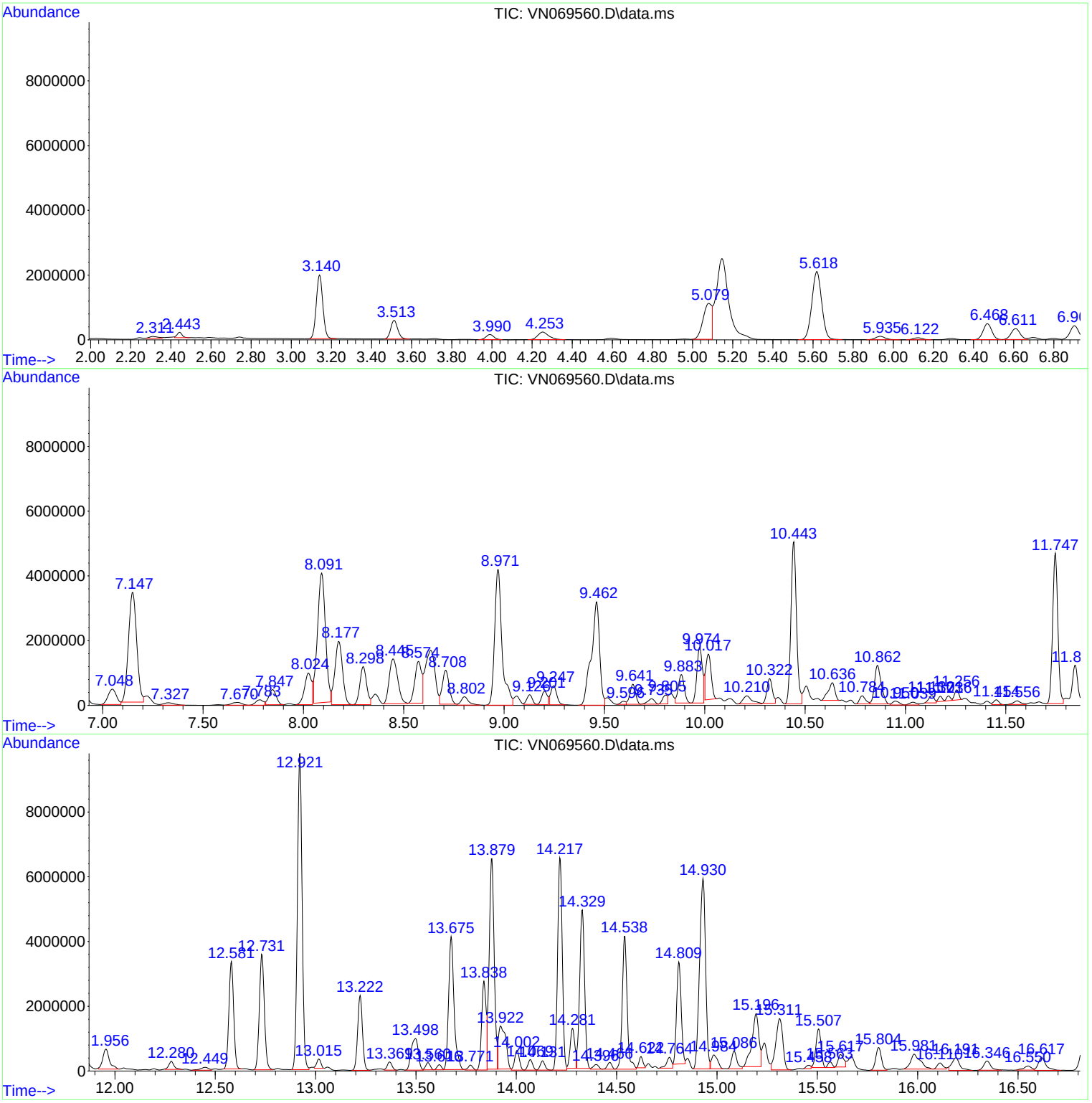
ClientSampleId :
LB-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
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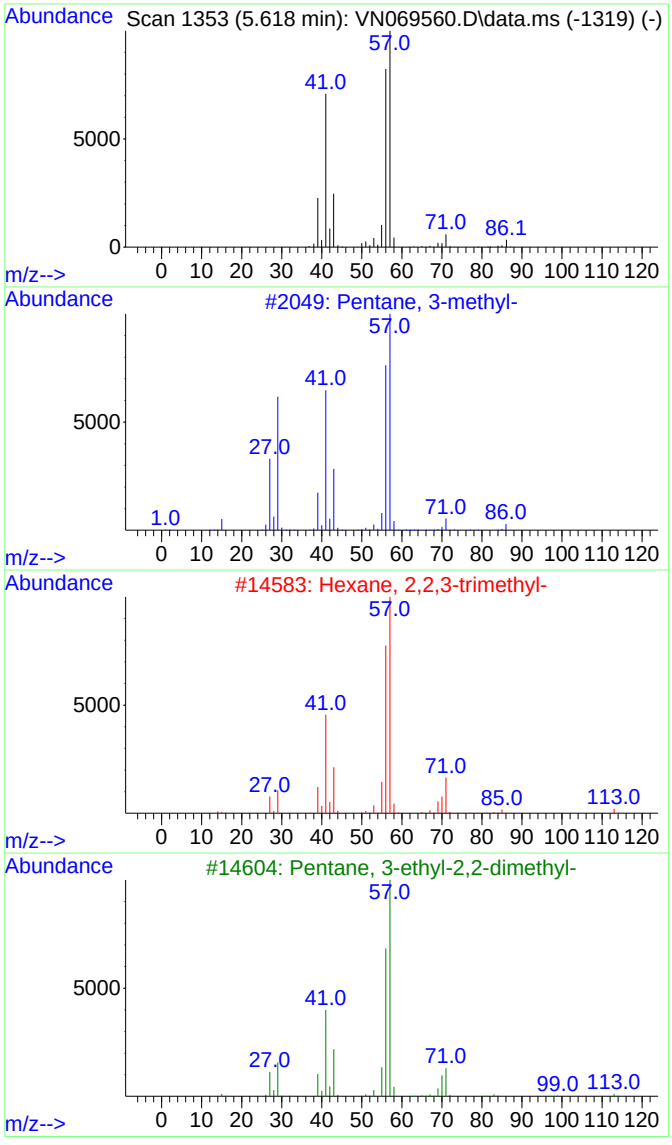
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

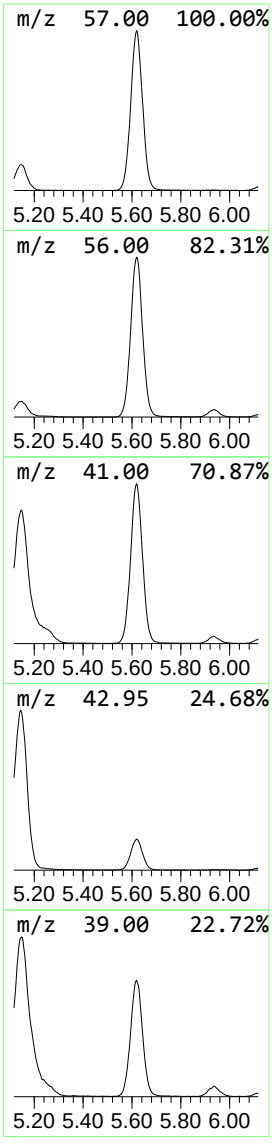
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.618	32.74 ug/l	7198670	Pentafluorobenzene	8.088

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
4	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40





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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

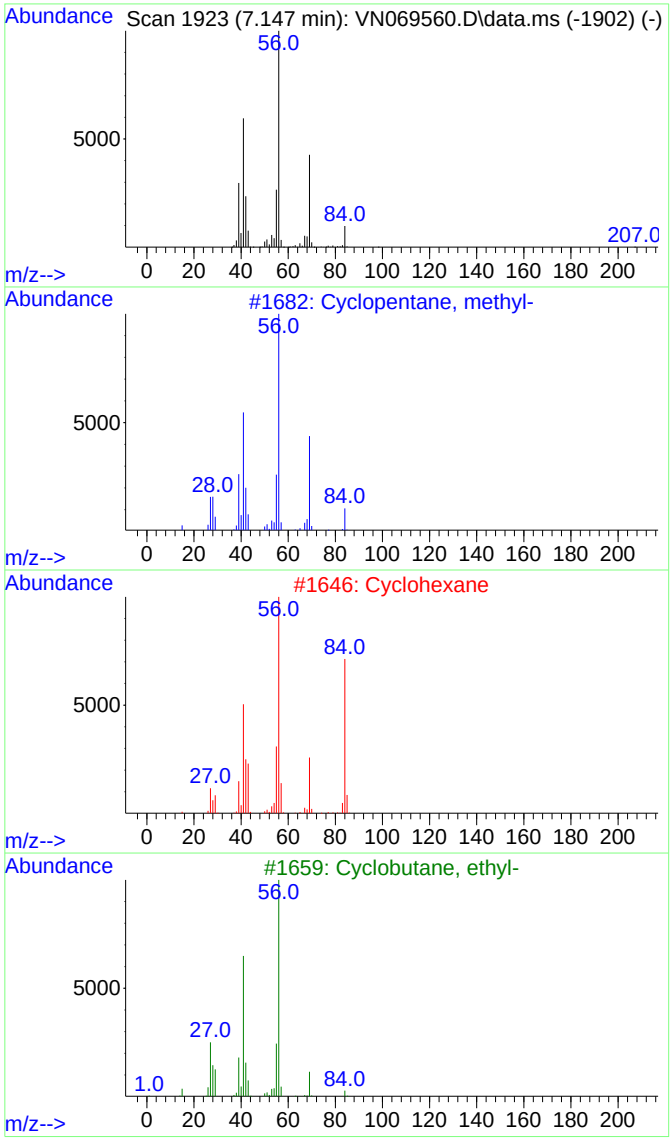
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

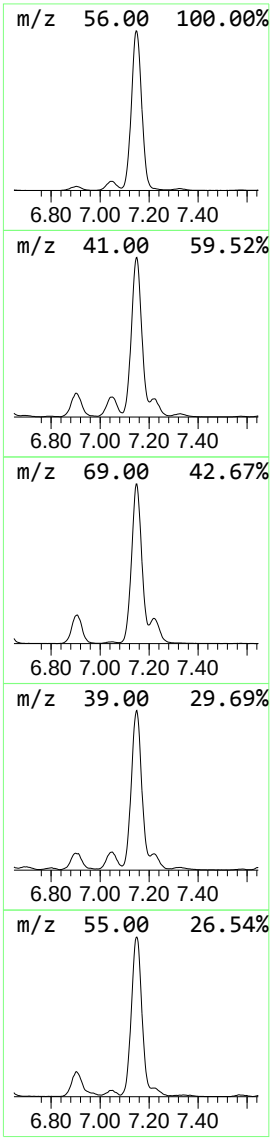
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.147	43.87 ug/l	9646130	Pentafluorobenzene	8.088

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59
5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	59





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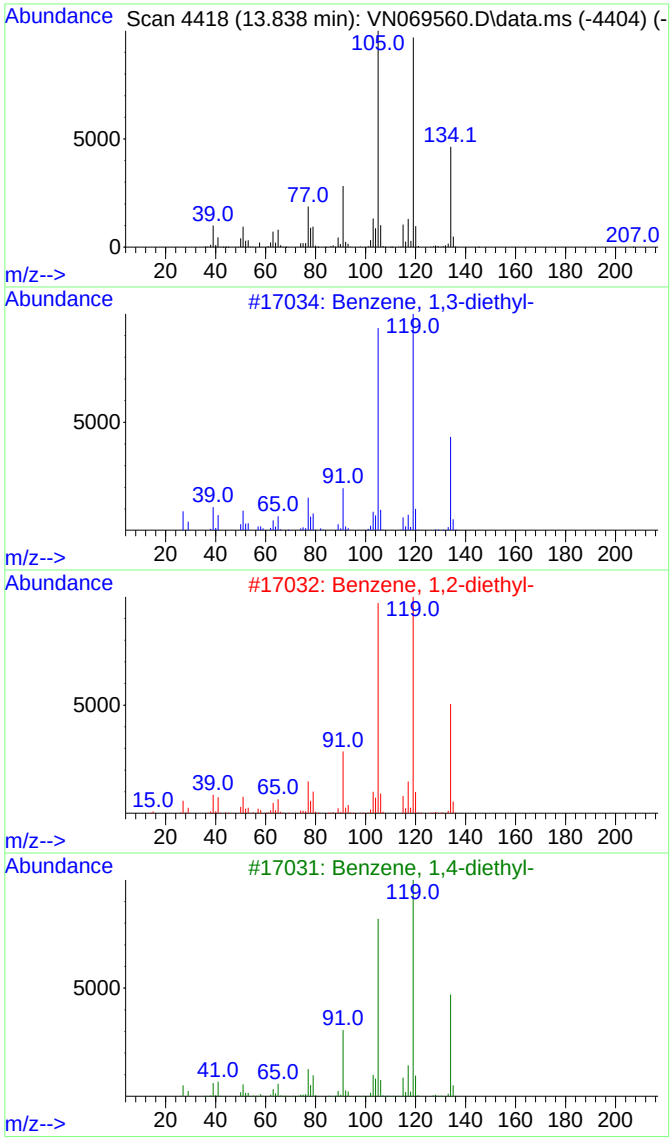
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

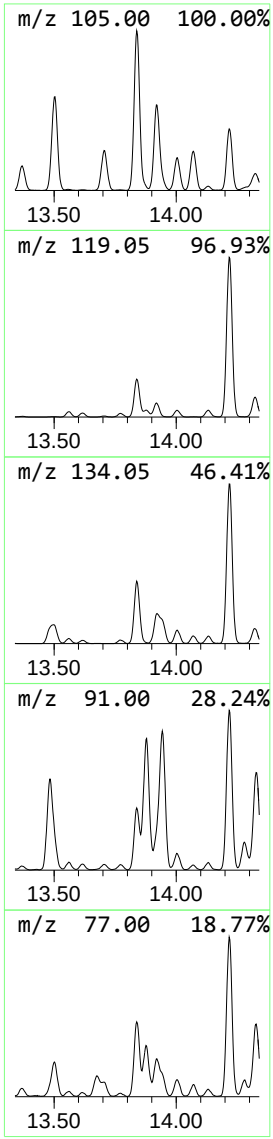
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.839	28.24 ug/l	4358720	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87





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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
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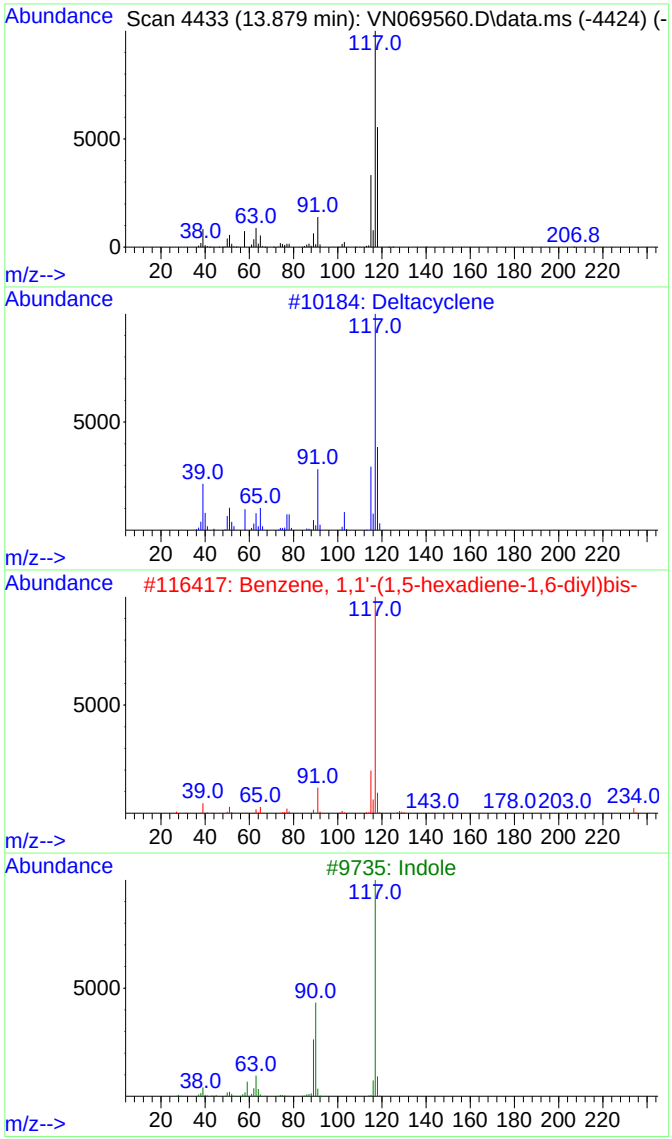
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

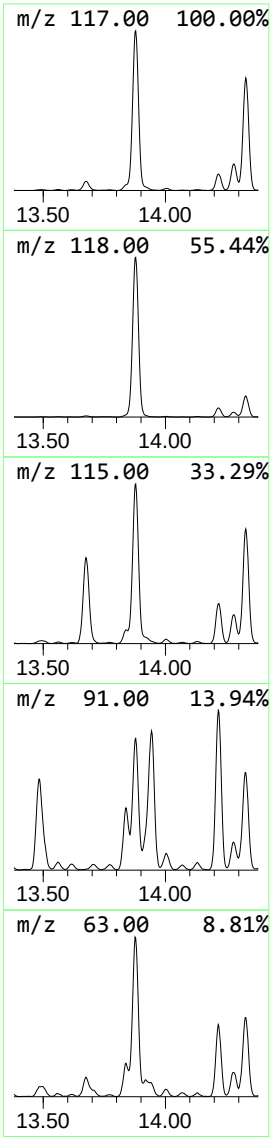
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Deltacyclene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.879	71.78 ug/l	11079700	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Deltacyclene	118	C9H10	007785-10-6	64
2	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
3	Indole	117	C8H7N	000120-72-9	49
4	5H-1-Pyrindine	117	C8H7N	000270-91-7	43
5	4-Ethynylaniline	117	C8H7N	014235-81-5	43





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Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

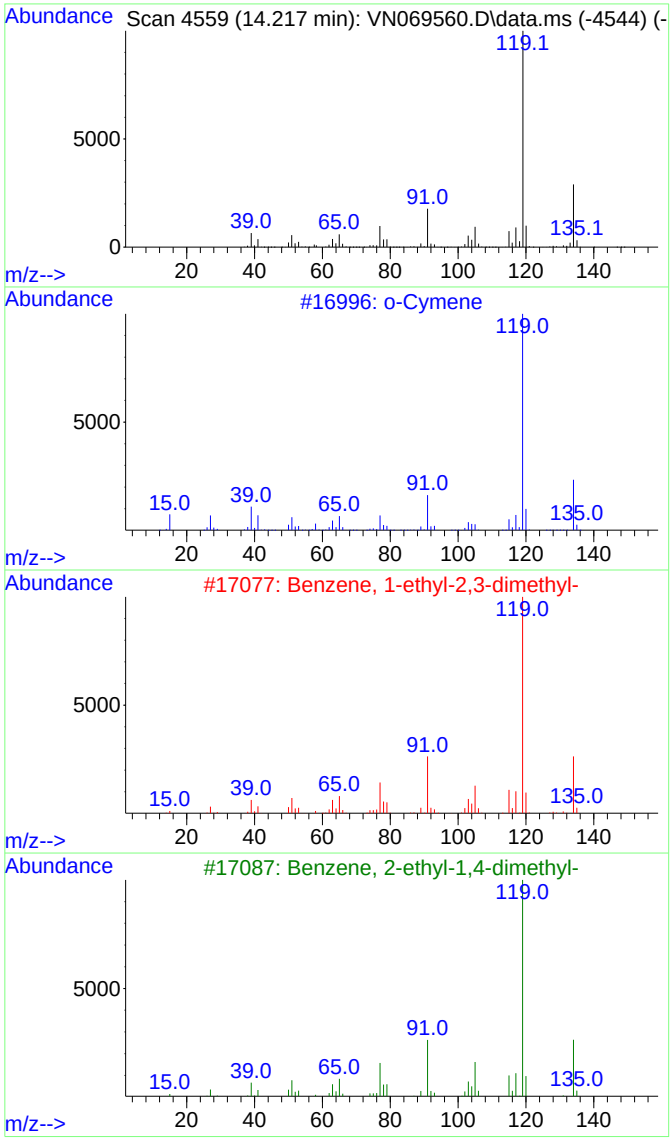
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

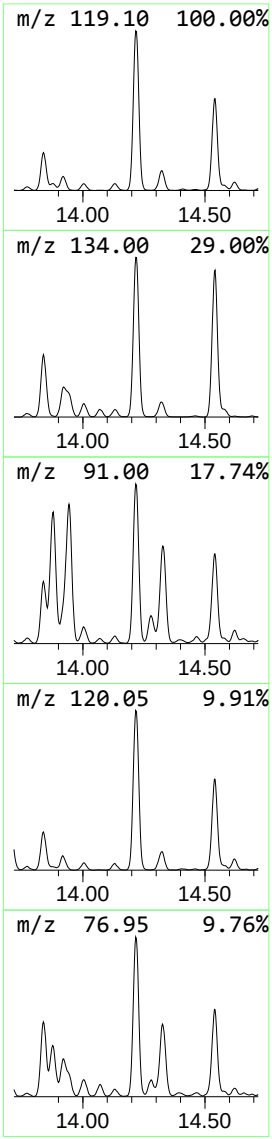
****

Peak Number 5 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.217	67.95 ug/l	10489400	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





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Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

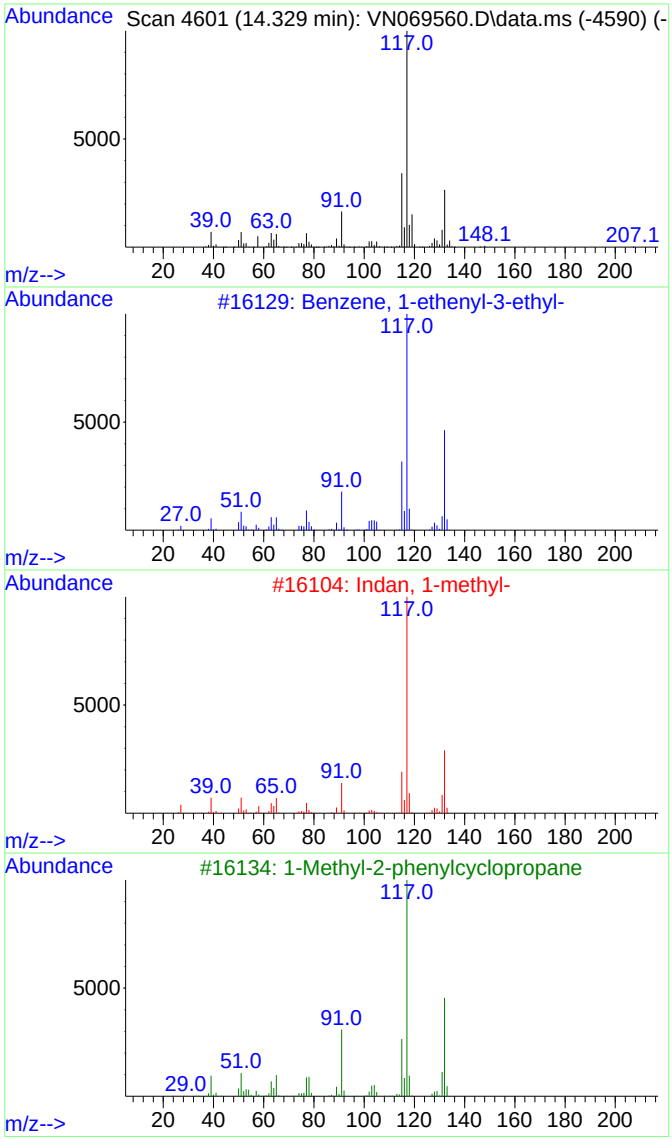
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

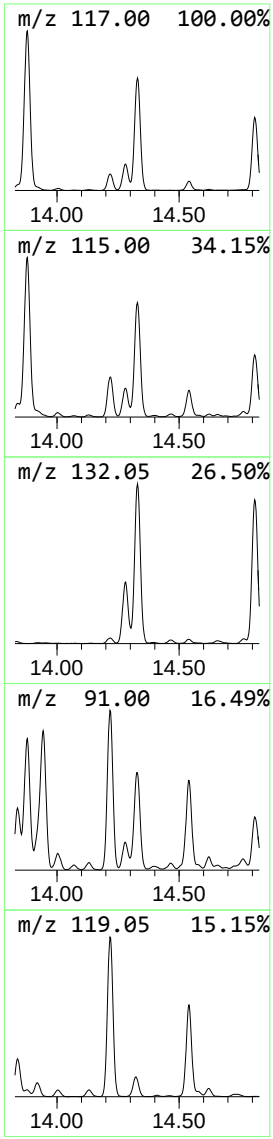
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	53.93 ug/l	8324840	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	72





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Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

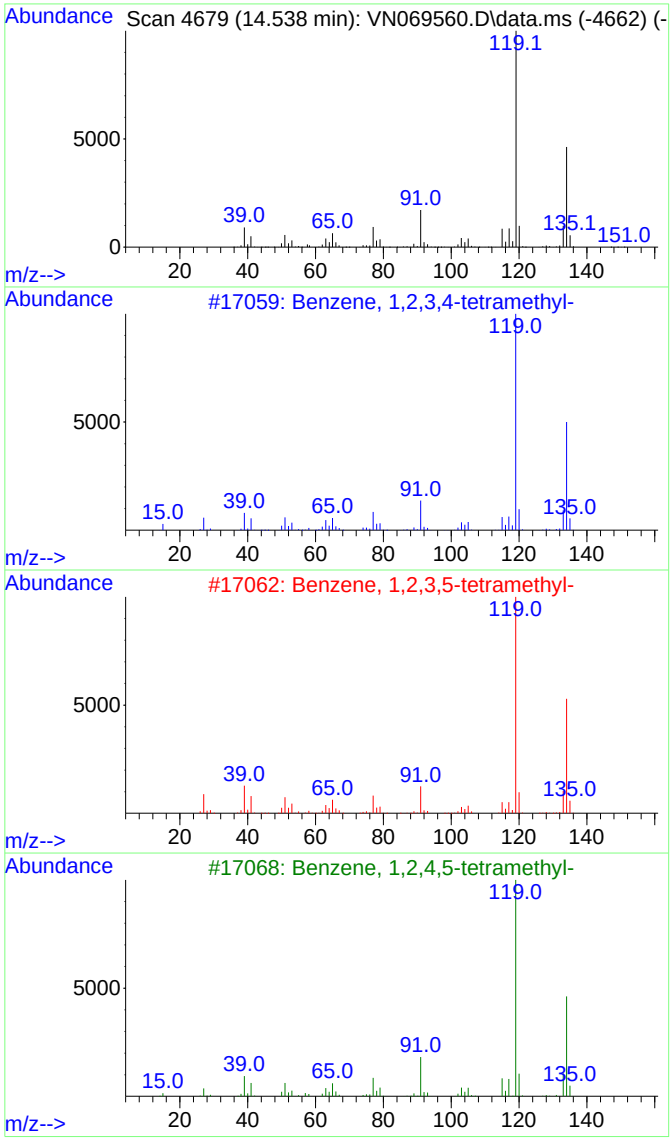
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

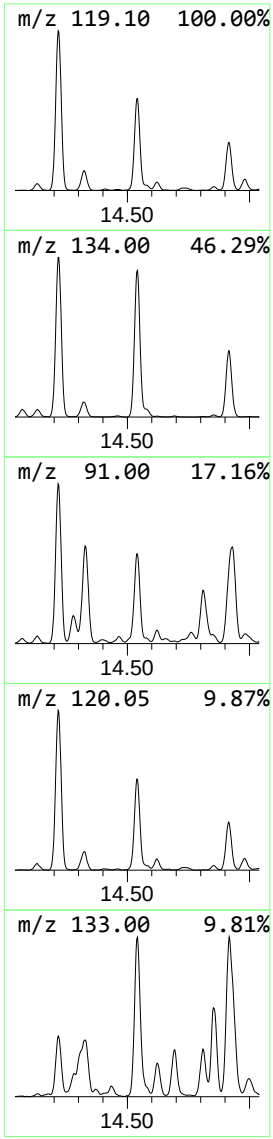
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.539	44.64 ug/l	6890120	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

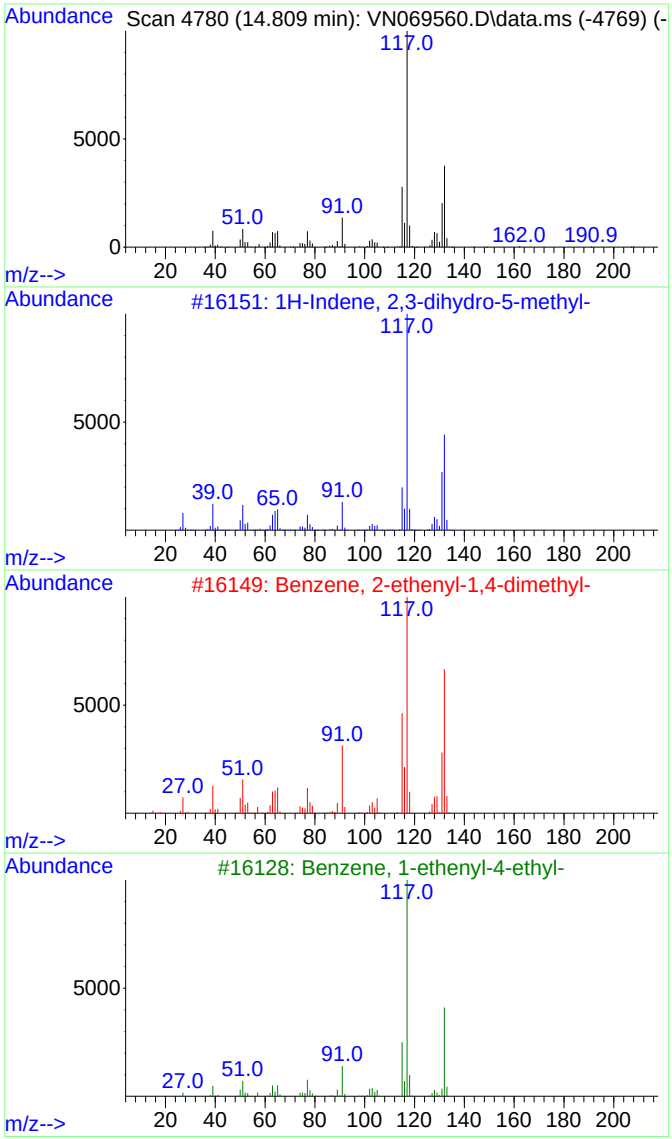
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

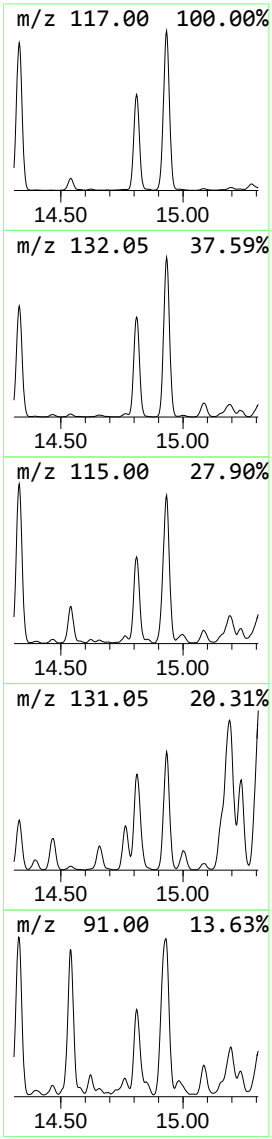
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.809	32.60 ug/l	5032390	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	96
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

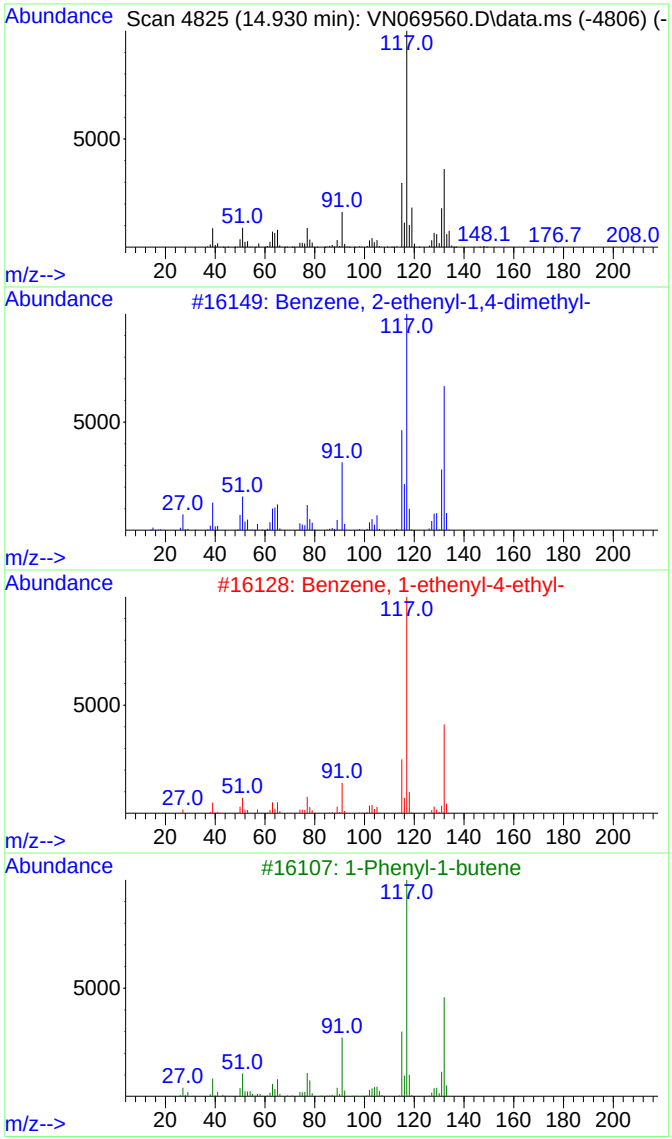
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

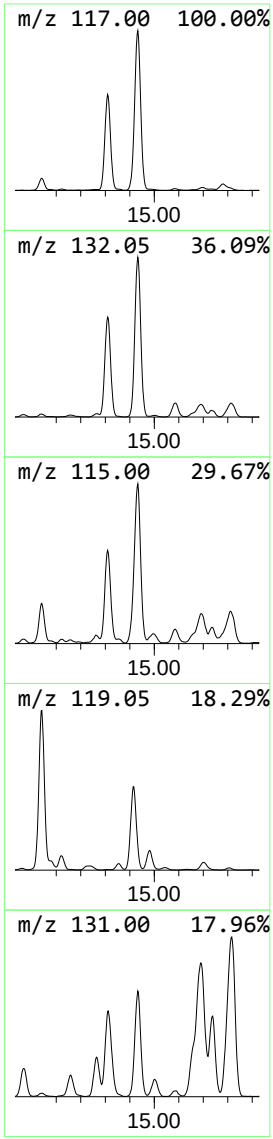
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.930	79.54 ug/l	12276900	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

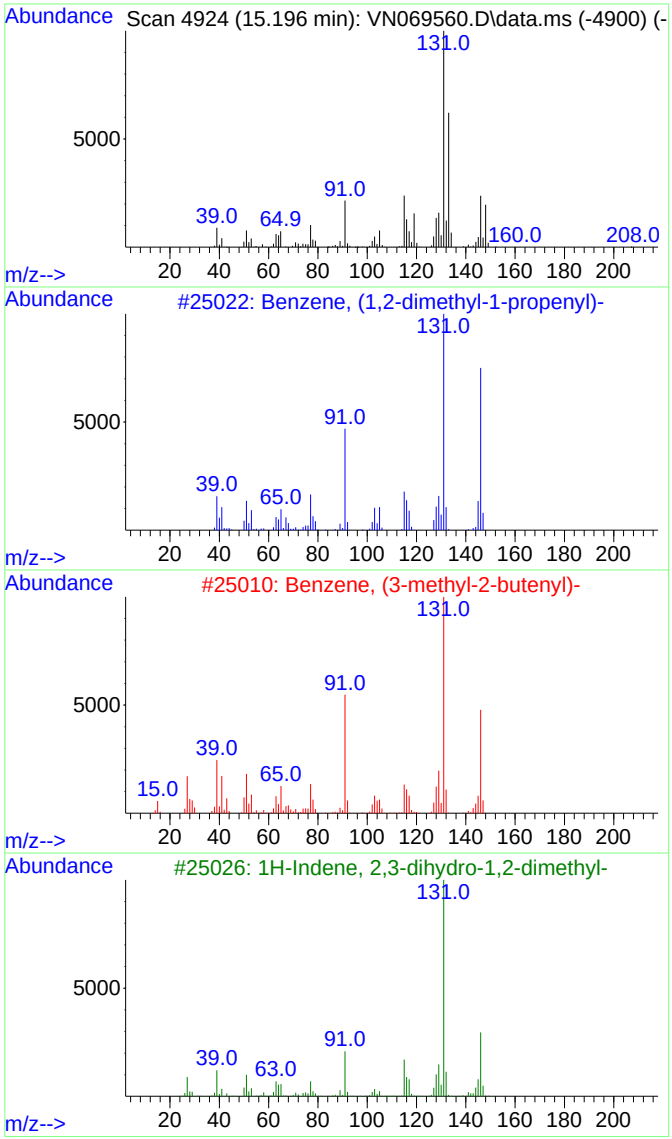
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

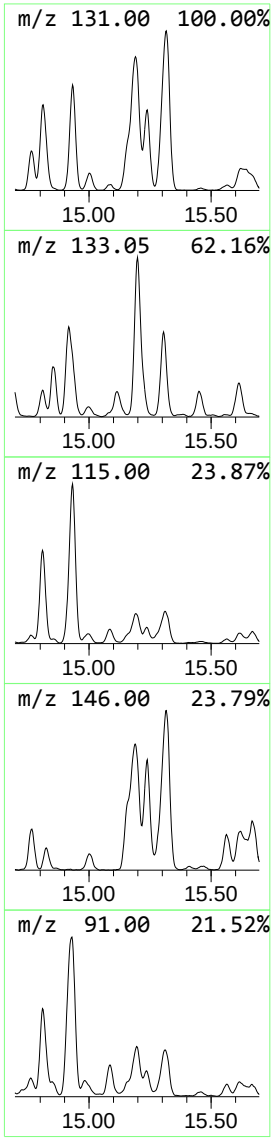
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.196	25.76 ug/l	3976080	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
Data File : VN069560.D
Acq On : 18 Nov 2021 18:17
Operator : JC/MD
Sample : M4739-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
LB-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Pentane, 3-methyl-	5.618	32.7	ug/l	7198670	1	8.088	10995100	50.0
Cyclopentane, m...	7.147	43.9	ug/l	9646130	1	8.088	10995100	50.0
Benzene, 1,3-di...	13.839	28.2	ug/l	4358720	4	13.675	7717900	50.0
Deltacyclene	13.879	71.8	ug/l	11079700	4	13.675	7717900	50.0
o-Cymene	14.217	68.0	ug/l	10489400	4	13.675	7717900	50.0
Benzene, 1-ethe...	14.329	53.9	ug/l	8324840	4	13.675	7717900	50.0
Benzene, 1,2,3,...	14.539	44.6	ug/l	6890120	4	13.675	7717900	50.0
1H-Indene, 2,3-...	14.809	32.6	ug/l	5032390	4	13.675	7717900	50.0
Benzene, 2-ethe...	14.930	79.5	ug/l	12276900	4	13.675	7717900	50.0
Benzene, (1,2-d...	15.196	25.8	ug/l	3976080	4	13.675	7717900	50.0