LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

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\82N061722W.M
Title : SW846 8260

Signal : TIC: VN073124.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	192	202	214	rBV	63697	150813	7.85%	0.736%
2	3.451	257	266	279	rVB2	30241	78895	4.11%	0.385%
3	3.651	294	300	308	rBV4	10321	22876	1.19%	0.112%
4	3.922	338	346	355	rVB2	12905	34082	1.77%	0.166%
5	5.004	517	530	531	rBV4	18245	53250	2.77%	0.260%
6	5.057	532	539	552	rVB3	52688	187129	9.74%	0.914%
7	5.534	605	620	636	rBV3	37380	139547	7.27%	0.681%
8	5.851	663	674	687	rVB7	6875	28668	1.49%	0.140%
9	6.040	694	706	718	rVB3	23016	72675	3.78%	0.355%
10	6.316	744	753	758	rBV4	7597	21386	1.11%	0.104%
11	6.816	829	838	850	rVB5	11860	34152	1.78%	0.167%
12	6.963	853	863	871	rBV3	29844	86619	4.51%	0.423%
13	7.063	871	880	890	rVV2	31421	87563	4.56%	0.427%
14	7.239	900	910	924	rBV	40729	117412	6.11%	0.573%
15	7.775	993	1001	1007	rVB4	10739	25454	1.33%	0.124%
16	7.951	1020	1031	1036	rBV2	236714	595249	31.00%	2.906%
17	8.016	1036	1042	1051	rVV	496787	1183057	61.60%	5.776%
18	8.104	1051	1057	1069	rVB	84510	205631	10.71%	1.004%
19	8.228	1069	1078	1087	rBV3	59338	140059	7.29%	0.684%
20	8.369	1093	1102	1116	rBV	244147	570851	29.73%	2.787%
21	8.545	1116	1132	1144	rVV	276704	749599	39.03%	3.659%
22	8.645	1144	1149	1158	rVB5	12927	31502	1.64%	0.154%
23	8.757	1158	1168	1179	rVB4	27427	71572	3.73%	0.349%
24	8.904	1183	1193	1208	rBV	631568	1383310	72.03%	6.753%
25	9.169	1226	1238	1249	rBV10	6771	28774	1.50%	0.140%
26	9.392	1262	1276	1281	rBV3	68467	176840	9.21%	0.863%
27	9.445	1281	1285	1293	rVV2	44658	85639	4.46%	0.418%
28	9.575	1302	1307	1312	rVB4	14409	27466	1.43%	0.134%
29	9.663	1314	1322	1330	rVB6	9745	23960	1.25%	0.117%
30	9.816	1335	1348	1359	rVV2	122555	252797	13.16%	1.234%
31	9.951	1359	1371	1377	rVV	129163	283632	14.77%	1.385%
32	10.016	1377	1382	1396	rVB3	39750	101620	5.29%	0.496%
33	10.151	1396	1405	1415	rBV4	42000	109023	5.68%	0.532%
34	10.304	1425	1431	1436	rBV3	38406	67234	3.50%	0.328%
35	10.375	1436	1443	1451	rBV	858123	1602035	83.42%	7.821%
36	10.439	1451	1454	1460	rVB	40410	71283	3.71%	0.348%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
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Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

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\82N061722W.M
Title : SW846 8260

37	10.563	1468	1475	1480	rVB7	34177	70833	3.69%	0.346%
38	10.816	1509	1518	1526	rVV10	16343	47943	2.50%	0.234%
39	11.092	1560	1565	1574	rVB3	10767	23070	1.20%	0.113%
40	11.463	1620	1628	1635	rVV3	28357	64523	3.36%	0.315%
41	11.569	1639	1646	1654	rVB2	23102	43454	2.26%	0.212%
42	11.680	1654	1665	1673	rBV	917517	1620093	84.36%	7.909%
43	11.780	1675	1682	1691	rVB	170876	296634	15.45%	1.448%
44	11.886	1691	1700	1714	rBV	599159	1171382	61.00%	5.719%
45	12.216	1746	1756	1764	rBV	243410	424821	22.12%	2.074%
46	12.516	1802	1807	1812	rBV2	18734	33898	1.77%	0.165%
47	12.669	1823	1833	1843	rVV	783395	1350819	70.34%	6.595%
48	12.857	1860	1865	1870	rVB	70275	106312	5.54%	0.519%
49	12.921	1870	1876	1879	rVV	284771	462194	24.07%	2.256%
50	12.951	1879	1881	1885	rVV	120450	185601	9.66%	0.906%
51	12.998	1885	1889	1896	rVB	205420	327396	17.05%	1.598%
52	13.157	1910	1916	1924	rVB	105309	184048	9.58%	0.898%
53	13.304	1935	1941	1948	rBV	560767	865344	45.06%	4.224%
54	13.433	1956	1963	1967	rBV4	11577	31274	1.63%	0.153%
55	13.498	1971	1974	1978	rVV	18198	27963	1.46%	0.137%
56	13.610	1985	1993	2013	rVB2	1018318	1920416	100.00%	9.375%
57	13.774	2016	2021	2022	rBV2	37241	45282	2.36%	0.221%
58	13.810	2022	2027	2030	rVV2	126356	231052	12.03%	1.128%
59	13.845	2030	2033	2044	rVB2	127593	240242	12.51%	1.173%
60	13.933	2044	2048	2053	rVB6	15347	21147	1.10%	0.103%
61	14.010	2053	2061	2064	rBV2	27369	44572	2.32%	0.218%
62	14.068	2066	2071	2073	rBV	60590	97811	5.09%	0.478%
63	14.092	2073	2075	2081	rVV	60557	93843	4.89%	0.458%
64	14.151	2081	2085	2093	rVB	104337	158754	8.27%	0.775%
65	14.263	2099	2104	2109	rVB	44271	70196	3.66%	0.343%
66	14.392	2121	2126	2132	rBV	23969	43284	2.25%	0.211%
67	14.474	2132	2140	2143	rVV	65096	123194	6.41%	0.601%
68	14.515	2143	2147	2151	rVV	91011	144725	7.54%	0.707%
69	14.557	2151	2154	2158	rVV3	38275	58411	3.04%	0.285%
70	14.598	2158	2161	2164	rVV	21003	30604	1.59%	0.149%
71	14.627	2164	2166	2172	rVV5	12757	24971	1.30%	0.122%
72	14.698	2174	2178	2182	rVV2	18251	28606	1.49%	0.140%
73	14.745	2182	2186	2190	rVV	51749	92408	4.81%	0.451%
74	14.786	2190	2193	2198	rVV3	30939	48613	2.53%	0.237%
75	14.863	2198	2206	2212	rVV2	74030	195957	10.20%	0.957%
76	14.921	2212	2216	2222	rVB2	22671	39505	2.06%	0.193%

LSC Area Percent Report

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Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

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\82N061722W.M
Title : SW846 8260

77	15.127	2247	2251	2261	rVB3	29322	65637	3.42%	0.320%
78	15.233	2263	2269	2270	rBV5	21224	29683	1.55%	0.145%
79	15.433	2291	2303	2311	rVB2	60840	131521	6.85%	0.642%
80	15.633	2333	2337	2345	rVB2	15618	34927	1.82%	0.171%
81	15.727	2345	2353	2359	rBV4	17424	47168	2.46%	0.230%
82	16.109	2411	2418	2424	rVB9	10811	27372	1.43%	0.134%
83	16.527	2482	2489	2495	rBV	43329	97785	5.09%	0.477%
84	16.733	2517	2524	2533	rVB3	23595	56997	2.97%	0.278%

Sum of corrected areas: 20483939

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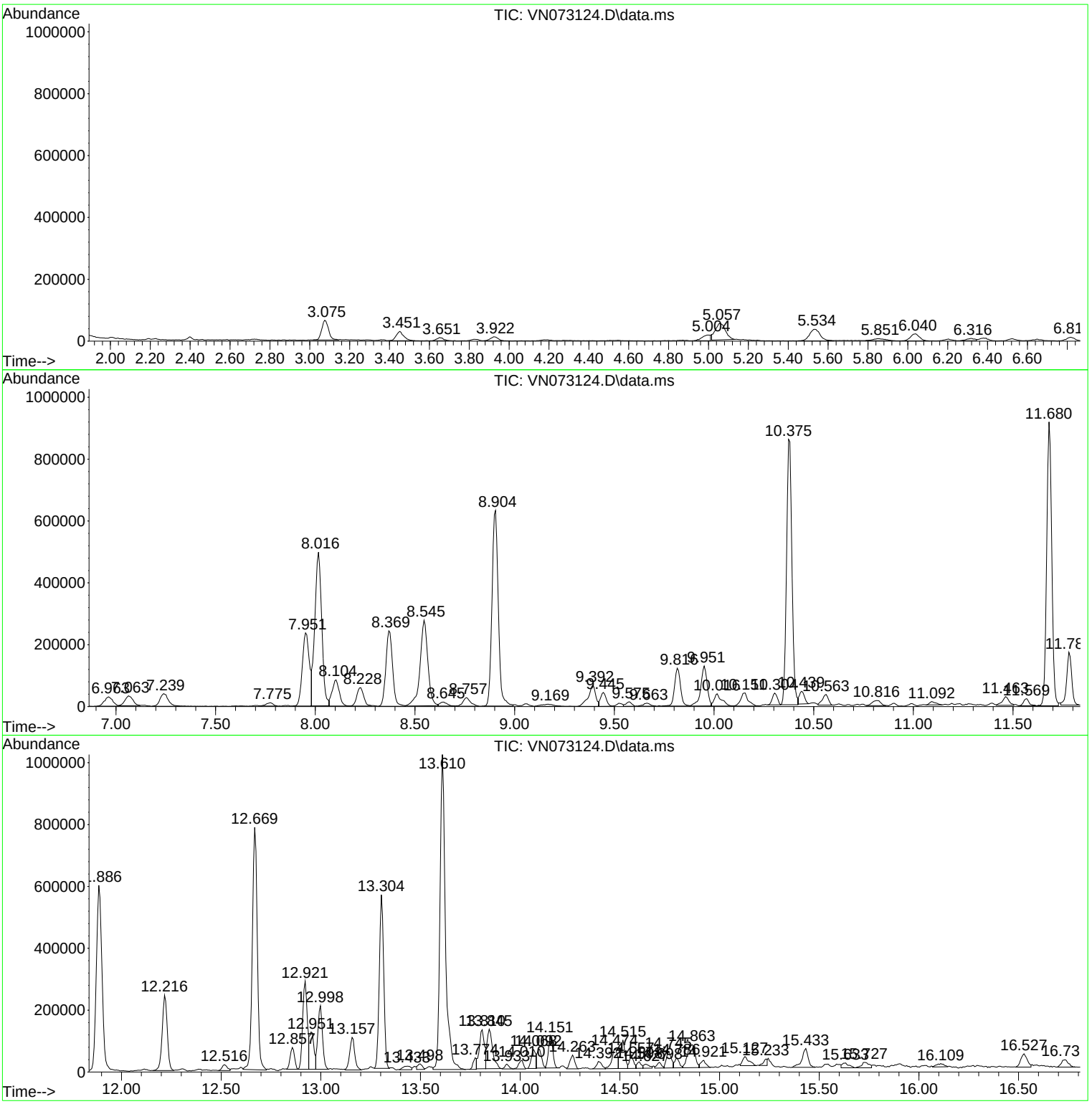
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
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Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

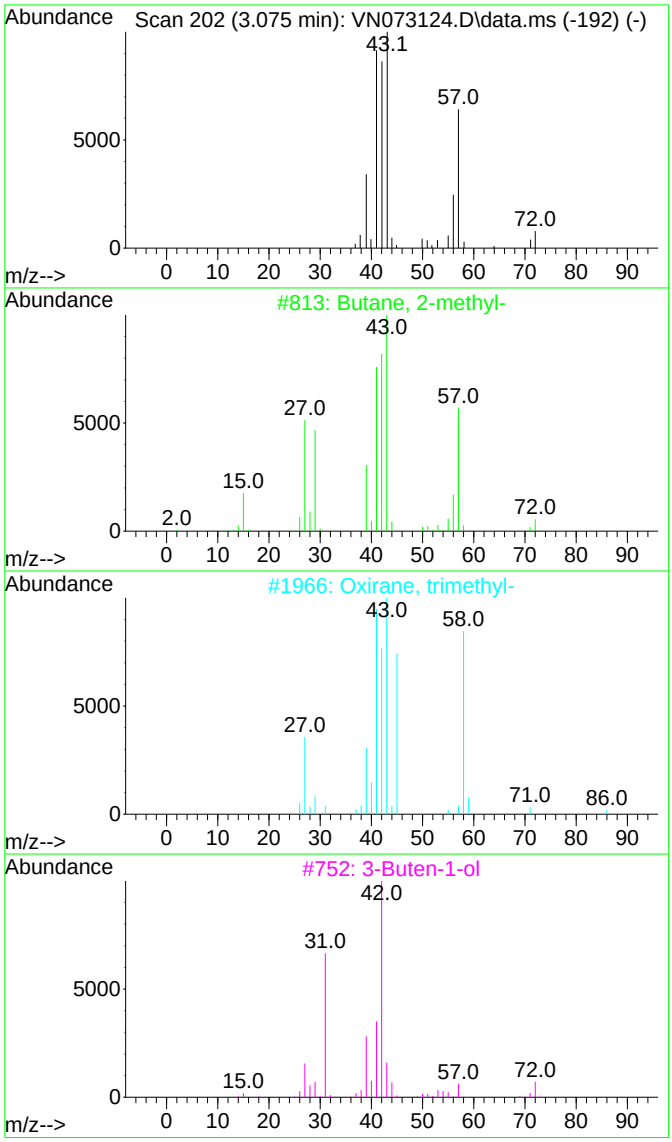
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

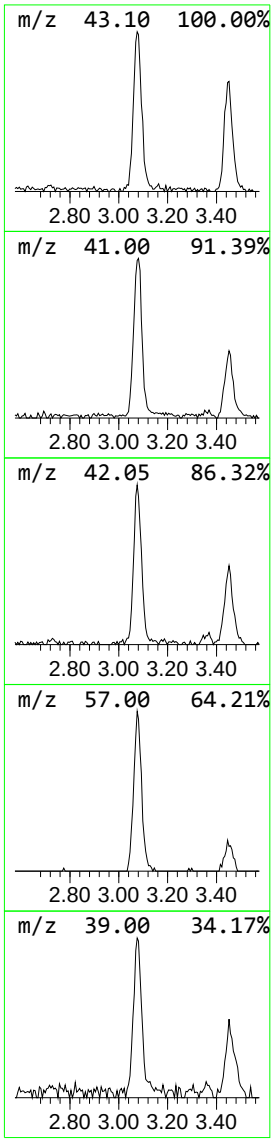
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	6.37 ug/l	150813	Pentafluorobenzene	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Oxirane, trimethyl-	86	C5H10O	005076-19-7	42
3	3-Buten-1-ol	72	C4H8O	000627-27-0	38
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	33
5	Pentane, 2-bromo-	150	C5H11Br	000107-81-3	17





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

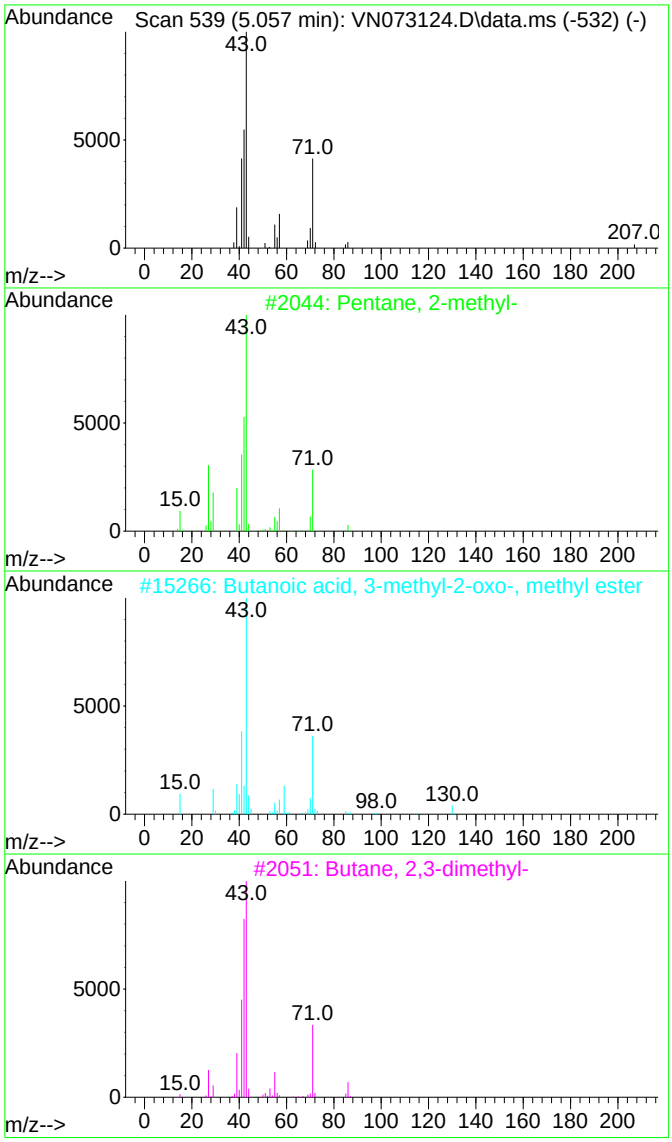
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

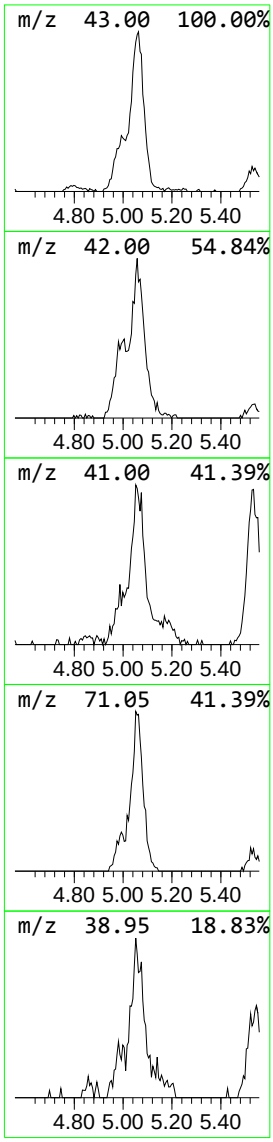
****

Peak Number 2 Pentane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.057	7.91 ug/l	187129	Pentafluorobenzene	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	87
2	Butanoic acid, 3-methyl-2-oxo-, ...	130	C6H10O3	003952-67-8	53
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38
5	Pentane	72	C5H12	000109-66-0	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

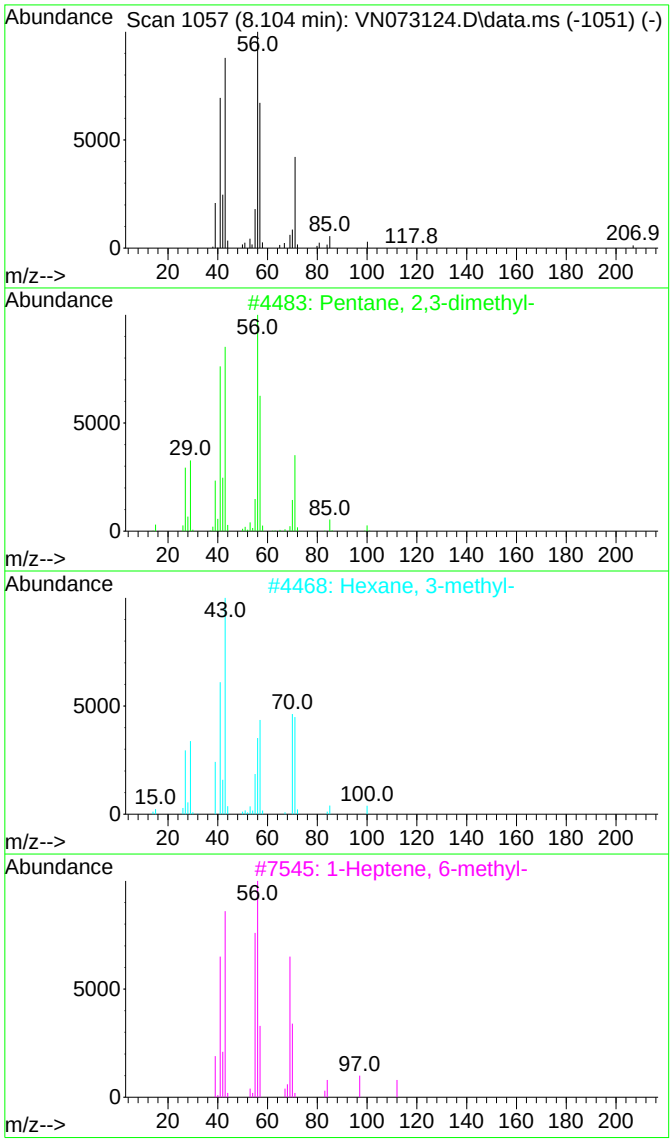
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

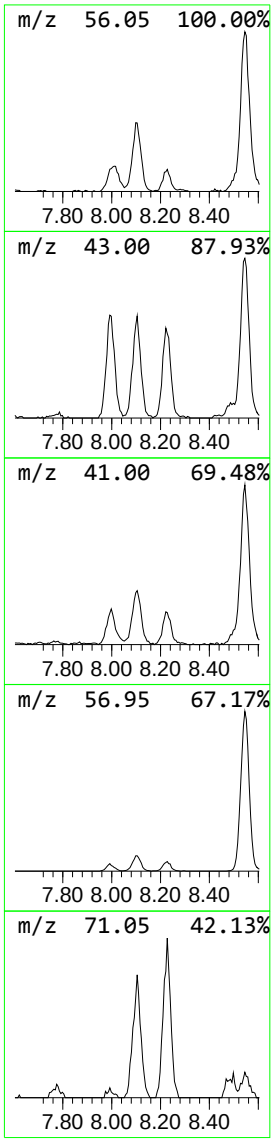
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	8.69 ug/l	205631	Pentafluorobenzene	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Hexane, 3-methyl-	100	C7H16	000589-34-4	46
3	1-Heptene, 6-methyl-	112	C8H16	005026-76-6	40
4	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	38
5	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	38





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Operator : JC\MD
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Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

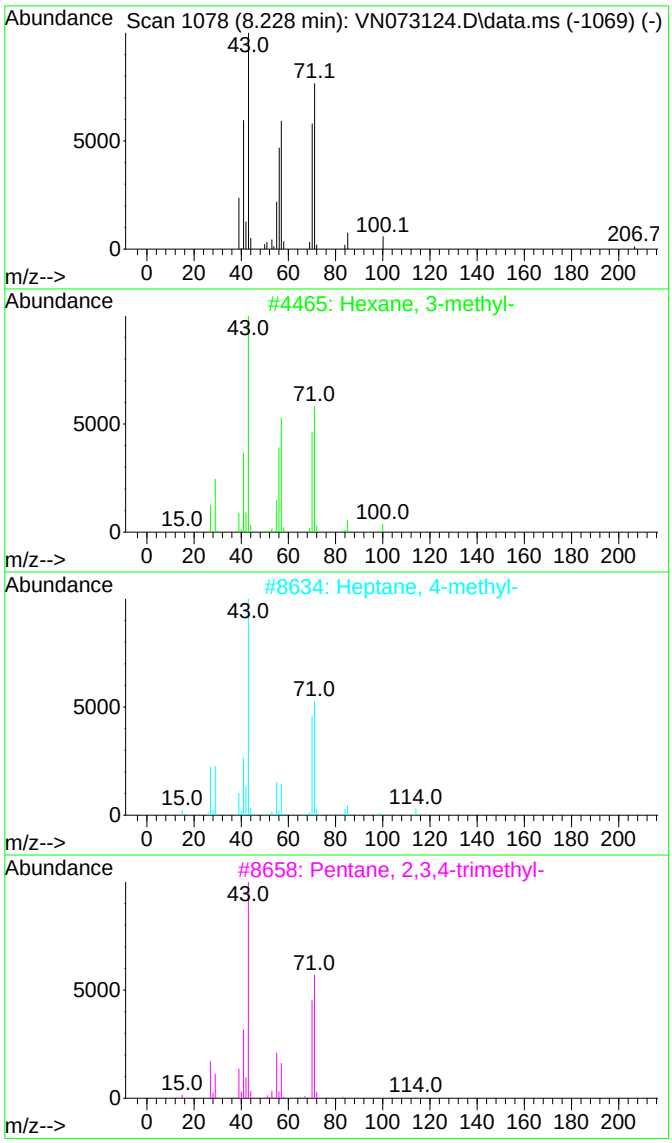
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

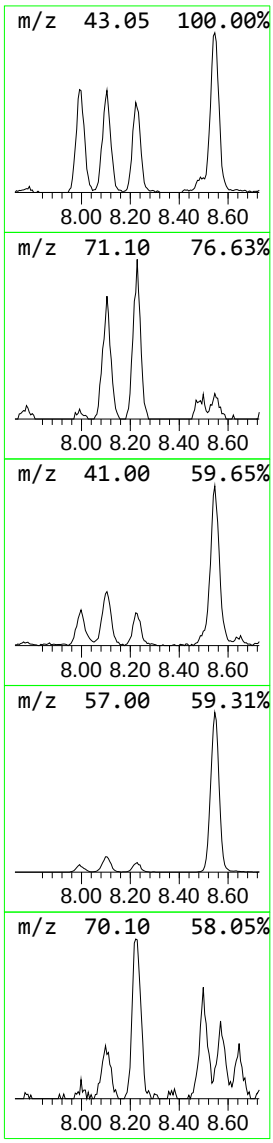
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Hexane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.228	5.92 ug/l	140059	Pentafluorobenzene	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
4	Pentane, 3-ethyl-	100	C7H16	000617-78-7	50
5	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	50





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Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

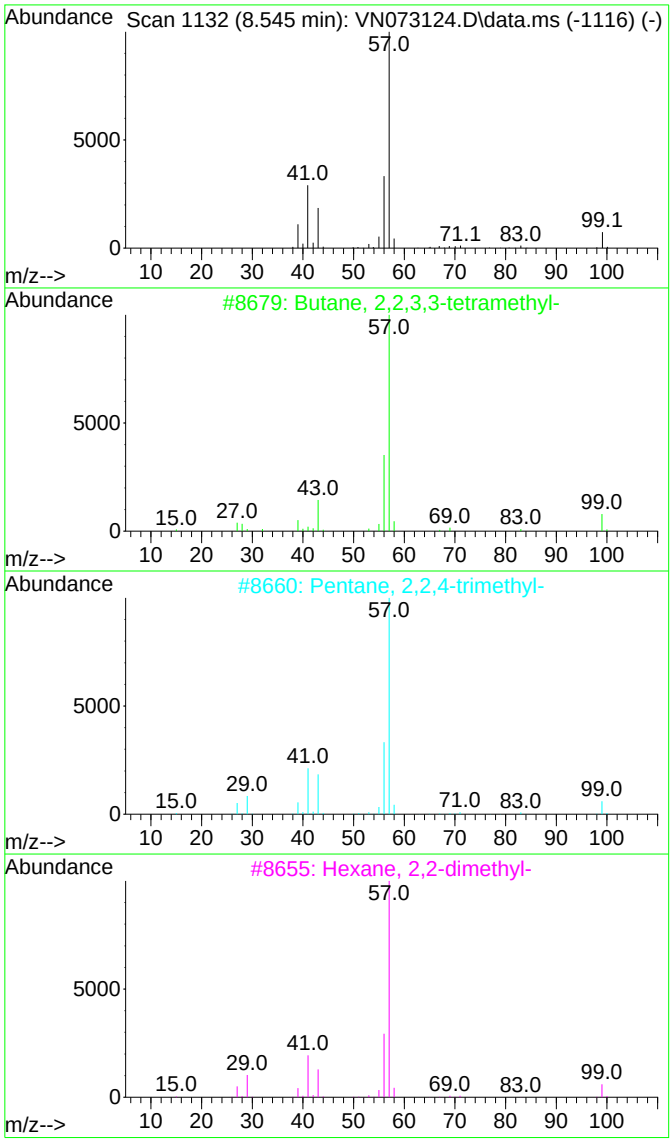
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

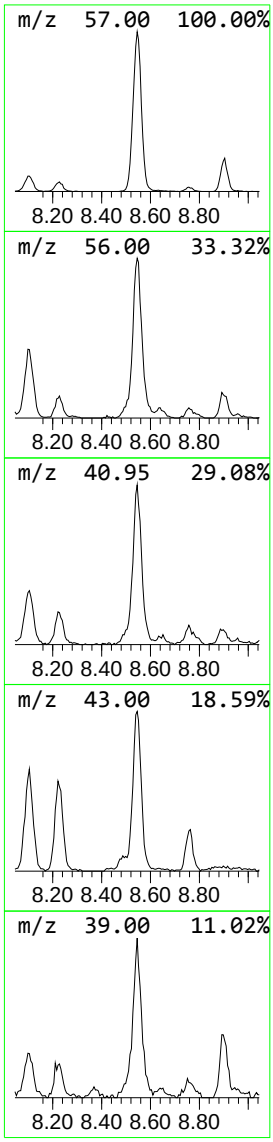
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.545	27.09 ug/l	749599	1,4-Difluorobenzene	8.904

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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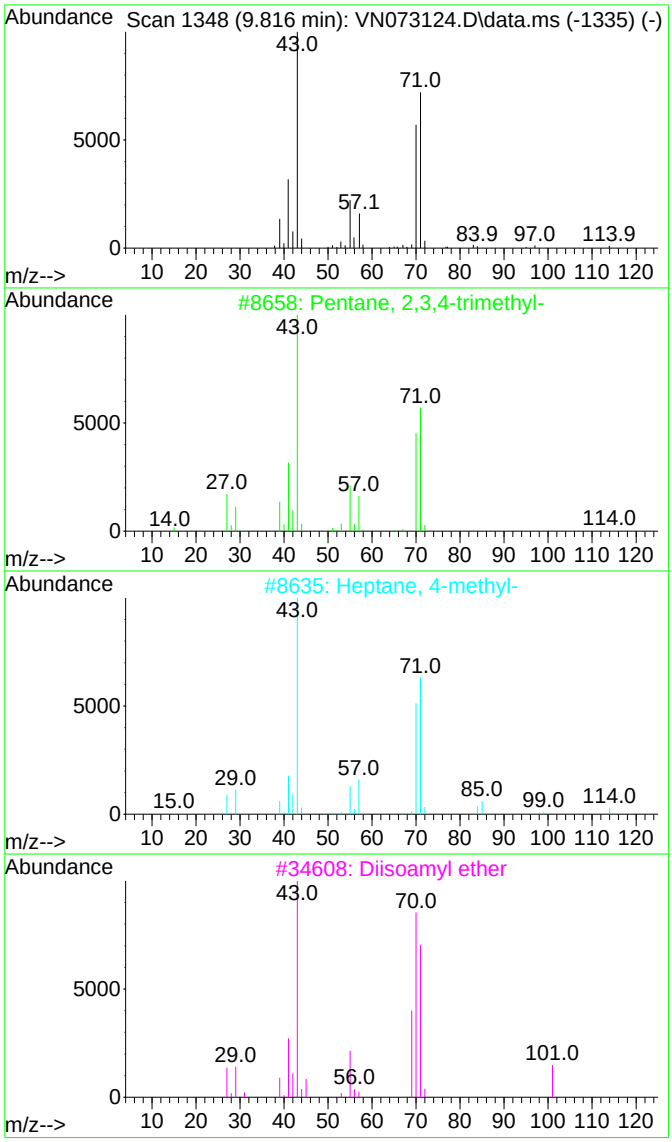
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

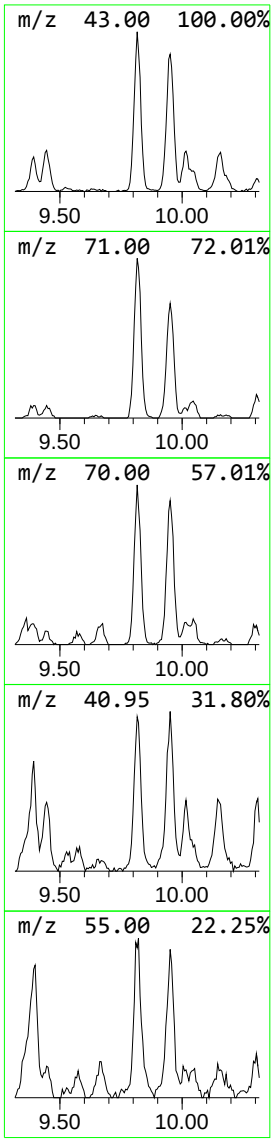
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 2,3,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.816	9.14 ug/l	252797	1,4-Difluorobenzene	8.904

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	90
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3	Diisoamyl ether	158	C10H22O	000544-01-4	78
4	Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

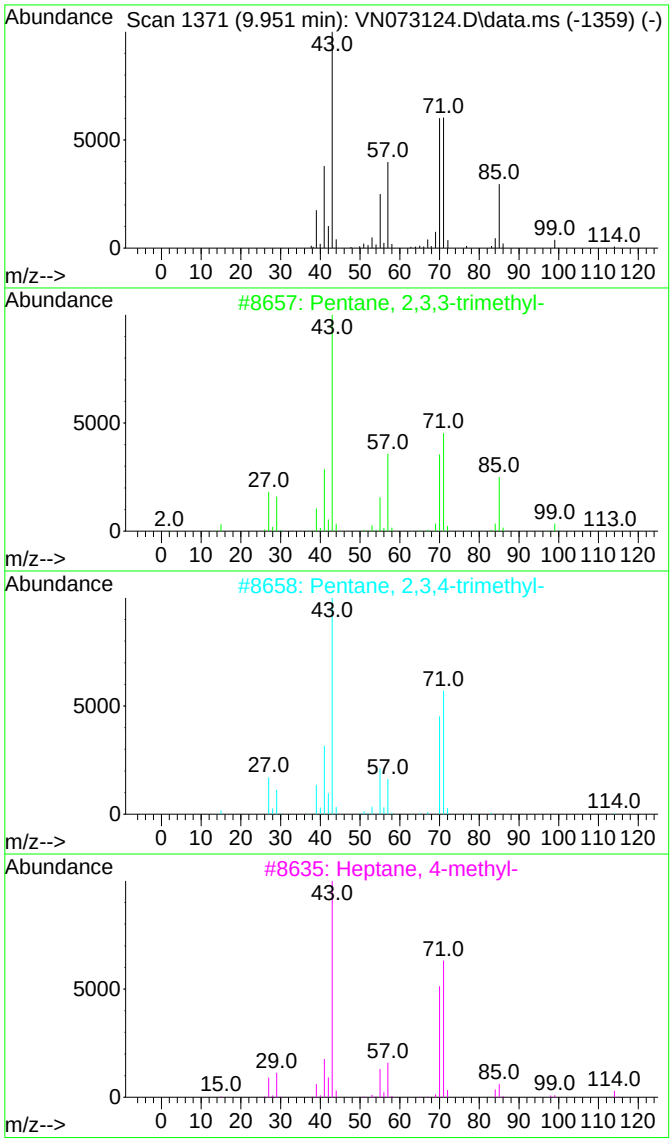
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

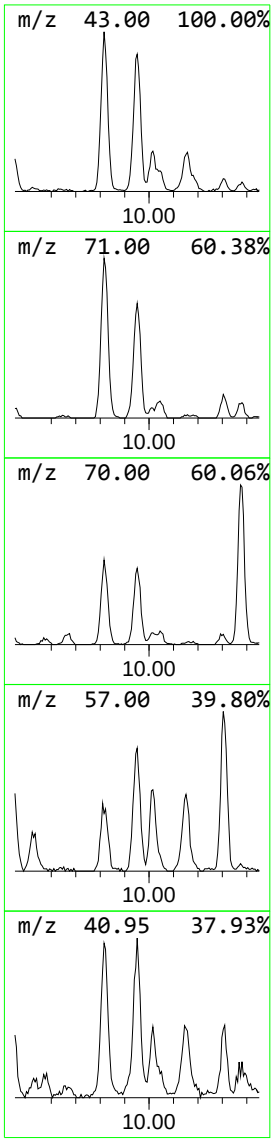
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentane, 2,3,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.951	10.25 ug/l	283632	1,4-Difluorobenzene	8.904

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	81
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	72
4	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
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ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

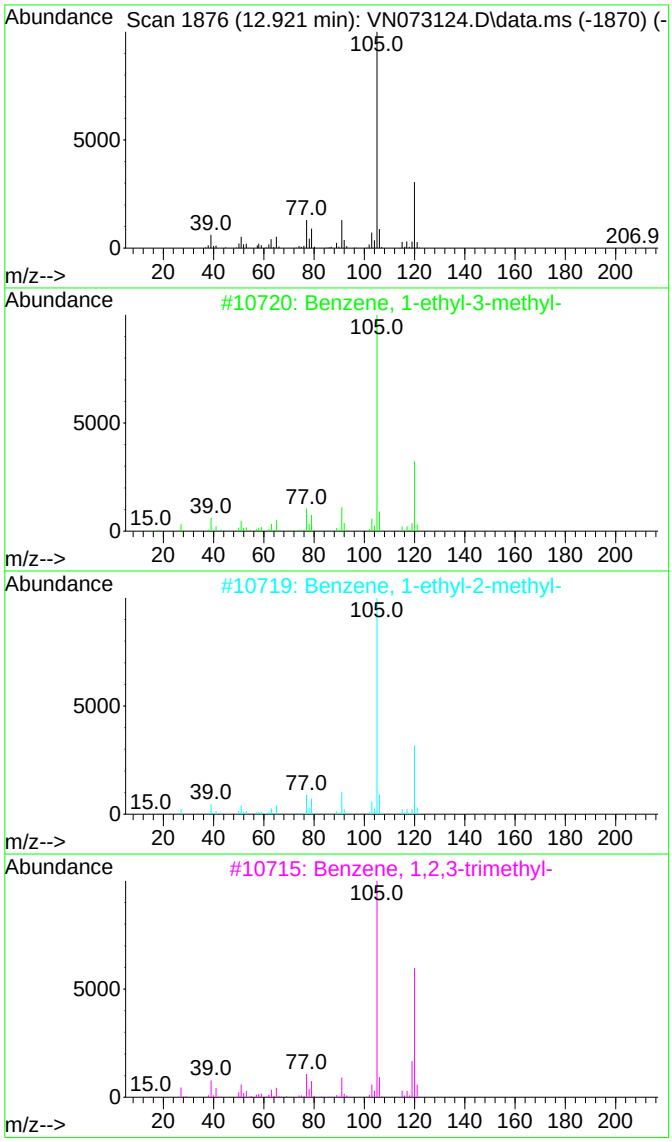
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

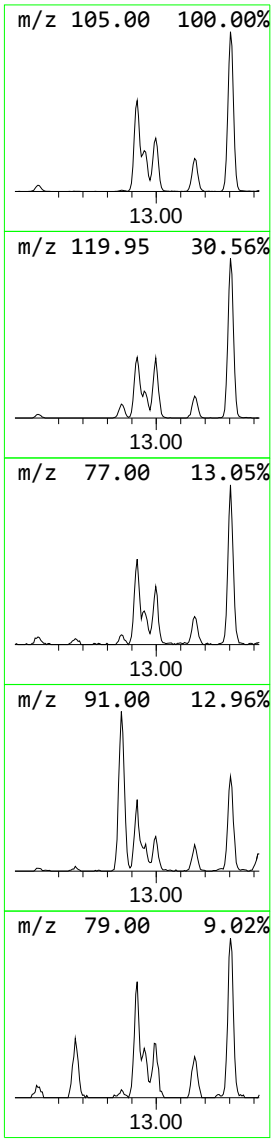
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.922	12.03 ug/l	462194	1,4-Dichlorobenzene-d4	13.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

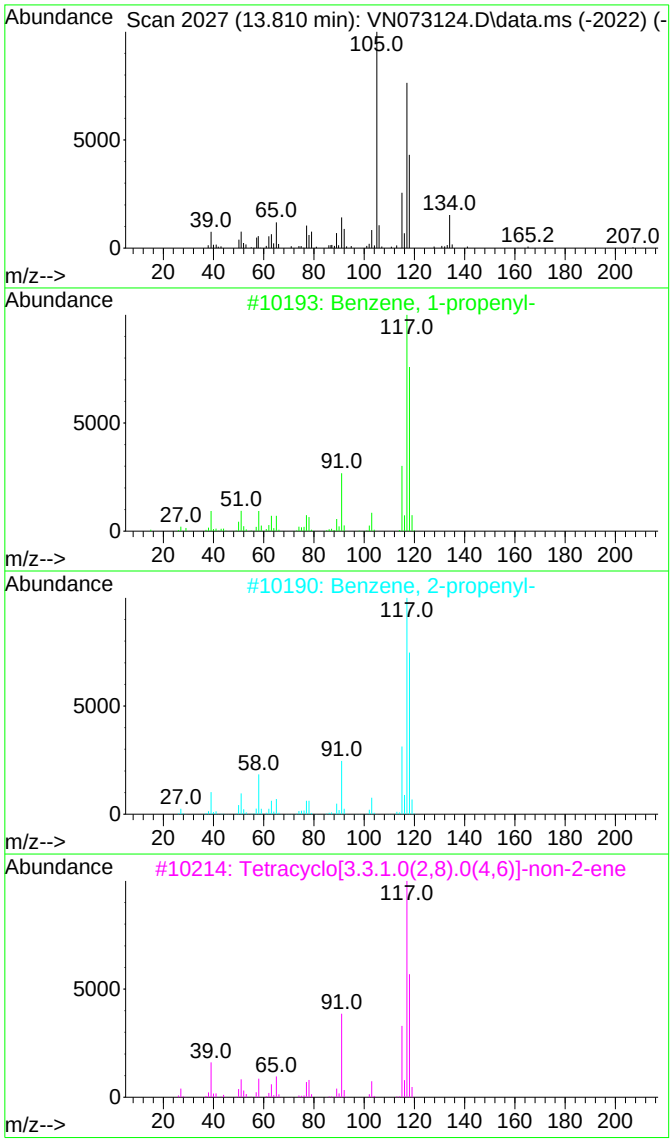
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

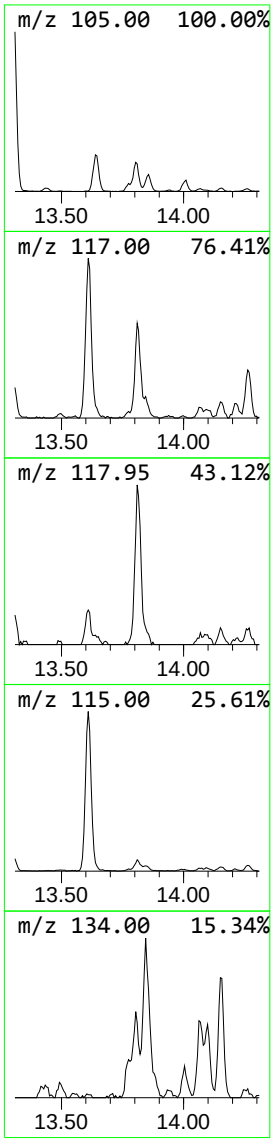
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	6.02 ug/l	231052	1,4-Dichlorobenzene-d4	13.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	70
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4	Indane	118	C9H10	000496-11-7	60
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	58





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Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

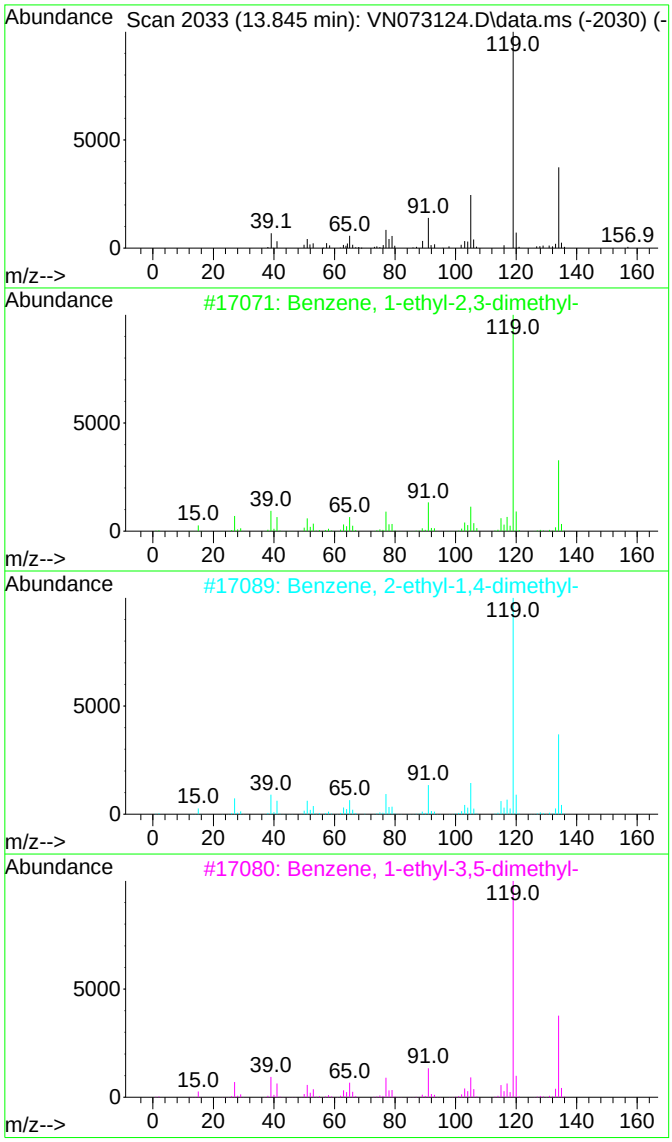
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260

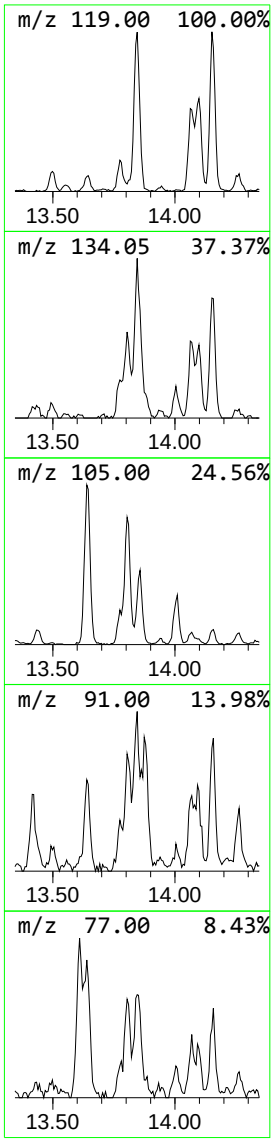
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.845	6.25 ug/l	240242	1,4-Dichlorobenzene-d4	13.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	83
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	80
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	72
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	68
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
Data File : VN073124.D
Acq On : 25 Jun 2022 20:48
Operator : JC\MD
Sample : N3434-07 40X
Misc : 6.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B2206-29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.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
Butane, 2-methyl-	3.075	6.4	ug/l	150813	1	8.016	1183060	50.0
Pentane, 2-methyl-	5.057	7.9	ug/l	187129	1	8.016	1183060	50.0
Pentane, 2,3-di...	8.104	8.7	ug/l	205631	1	8.016	1183060	50.0
Hexane, 3-methyl-	8.228	5.9	ug/l	140059	1	8.016	1183060	50.0
Butane, 2,2,3,3...	8.545	27.1	ug/l	749599	2	8.904	1383310	50.0
Pentane, 2,3,4-...	9.816	9.1	ug/l	252797	2	8.904	1383310	50.0
Pentane, 2,3,3-...	9.951	10.3	ug/l	283632	2	8.904	1383310	50.0
Benzene, 1-ethy...	12.922	12.0	ug/l	462194	4	13.610	1920420	50.0
Benzene, 1-prop...	13.810	6.0	ug/l	231052	4	13.610	1920420	50.0
Benzene, 1-ethy...	13.845	6.3	ug/l	240242	4	13.610	1920420	50.0