LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

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\82N060822W.M
Title : SW846 8260

Signal : TIC: VN072764.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.028	3	7	14	rVB2	21422	31418	1.45%	0.189%
2	2.581	95	101	109	rBV2	29572	53502	2.47%	0.322%
3	2.740	124	128	136	rVB5	9783	23428	1.08%	0.141%
4	4.281	375	390	401	rBV2	22096	78594	3.63%	0.474%
5	4.528	423	432	437	rBV	9810	28843	1.33%	0.174%
6	5.081	509	526	547	rBV3	29999	130893	6.04%	0.789%
7	5.622	604	618	627	rBV5	7421	25100	1.16%	0.151%
8	6.892	823	834	846	rVB6	6982	25020	1.15%	0.151%
9	7.039	846	859	874	rBV3	108229	348559	16.08%	2.101%
10	7.328	896	908	920	rBV4	17990	60906	2.81%	0.367%
11	7.833	983	994	1005	rVB6	20189	63708	2.94%	0.384%
12	8.016	1016	1025	1031	rBV	258659	637223	29.40%	3.840%
13	8.080	1031	1036	1044	rVV	268203	601354	27.75%	3.624%
14	8.169	1044	1051	1065	rVV2	178595	439792	20.29%	2.651%
15	8.433	1087	1096	1109	rBV	316980	708555	32.69%	4.270%
16	8.610	1109	1126	1137	rBV	892600	2167249	100.00%	13.061%
17	8.963	1178	1186	1195	rVV	403735	837231	38.63%	5.046%
18	9.051	1195	1201	1209	rVB2	38740	83825	3.87%	0.505%
19	9.427	1256	1265	1266	rBV3	28069	55938	2.58%	0.337%
20	9.451	1266	1269	1274	rVV3	43580	83605	3.86%	0.504%
21	9.504	1274	1278	1286	rVV3	58183	126028	5.82%	0.760%
22	9.592	1286	1293	1305	rVB	46961	104418	4.82%	0.629%
23	9.733	1305	1317	1322	rBV4	29153	73036	3.37%	0.440%
24	9.880	1332	1342	1353	rBV	595545	1160714	53.56%	6.995%
25	10.010	1355	1364	1377	rBV	574960	1175189	54.22%	7.083%
26	10.110	1377	1381	1387	rVB5	14284	26715	1.23%	0.161%
27	10.198	1387	1396	1403	rBV2	30989	72353	3.34%	0.436%
28	10.369	1412	1425	1430	rBV2	81876	169366	7.81%	1.021%
29	10.439	1430	1437	1451	rVB	899272	1655495	76.39%	9.977%
30	10.610	1459	1466	1468	rBV4	17715	32266	1.49%	0.194%
31	10.651	1469	1473	1478	rVB	28020	45960	2.12%	0.277%
32	10.880	1504	1512	1518	rBV6	22125	58216	2.69%	0.351%
33	11.398	1595	1600	1605	rBV6	12326	21908	1.01%	0.132%
34	11.463	1607	1611	1618	rVB5	13146	25246	1.16%	0.152%
35	11.545	1620	1625	1635	rVB7	11893	32603	1.50%	0.196%
36	11.739	1647	1658	1668	rBV	550137	995405	45.93%	5.999%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
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Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

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\82N060822W.M
Title : SW846 8260

37	11.874	1672	1681	1688	rVB3	46883	93196	4.30%	0.562%
38	11.992	1694	1701	1711	rVB8	21055	58435	2.70%	0.352%
39	12.257	1739	1746	1753	rBV8	17948	44909	2.07%	0.271%
40	12.451	1774	1779	1784	rVV5	20945	41207	1.90%	0.248%
41	12.510	1785	1789	1795	rVB6	16790	34131	1.57%	0.206%
42	12.727	1818	1826	1835	rVB	638047	1092895	50.43%	6.587%
43	12.892	1849	1854	1862	rVB7	43409	91590	4.23%	0.552%
44	12.986	1862	1870	1878	rBV7	21670	63383	2.92%	0.382%
45	13.063	1878	1883	1891	rVV10	15298	42305	1.95%	0.255%
46	13.286	1913	1921	1922	rVV6	14997	28525	1.32%	0.172%
47	13.304	1922	1924	1933	rVB8	16085	28200	1.30%	0.170%
48	13.480	1946	1954	1955	rBV	42724	61435	2.83%	0.370%
49	13.668	1979	1986	1993	rBV	455440	778896	35.94%	4.694%
50	13.815	2008	2011	2021	rVB9	19654	45981	2.12%	0.277%
51	13.915	2024	2028	2037	rVV6	23199	59223	2.73%	0.357%
52	13.998	2037	2042	2048	rVB4	50008	84573	3.90%	0.510%
53	14.145	2061	2067	2073	rBV4	35254	65136	3.01%	0.393%
54	14.209	2073	2078	2086	rVB8	17982	36323	1.68%	0.219%
55	14.327	2091	2098	2103	rBV6	25838	47856	2.21%	0.288%
56	14.398	2103	2110	2114	rBV7	28616	76724	3.54%	0.462%
57	14.509	2124	2129	2131	rBV4	37740	64619	2.98%	0.389%
58	14.533	2131	2133	2140	rVB3	56204	87238	4.03%	0.526%
59	14.615	2143	2147	2150	rBV4	16165	25476	1.18%	0.154%
60	14.662	2151	2155	2162	rVV6	15523	33223	1.53%	0.200%
61	14.739	2162	2168	2176	rVB5	37177	101792	4.70%	0.613%
62	14.815	2176	2181	2186	rBV3	23171	39424	1.82%	0.238%
63	14.909	2193	2197	2205	rBV6	23526	55159	2.55%	0.332%
64	14.974	2205	2208	2215	rVV5	22119	48115	2.22%	0.290%
65	15.133	2230	2235	2240	rBV7	20061	49134	2.27%	0.296%
66	15.192	2240	2245	2249	rVV2	40413	60340	2.78%	0.364%
67	15.239	2249	2253	2257	rVV3	20952	48272	2.23%	0.291%
68	15.298	2259	2263	2270	rVB3	42730	86001	3.97%	0.518%
69	15.404	2275	2281	2286	rVV3	21223	45614	2.10%	0.275%
70	15.456	2286	2290	2296	rVB5	25936	46953	2.17%	0.283%
71	15.556	2301	2307	2312	rBV4	42765	99747	4.60%	0.601%
72	15.798	2340	2348	2354	rBV8	24039	73388	3.39%	0.442%
73	15.886	2358	2363	2367	rBV8	12983	25497	1.18%	0.154%
74	15.956	2367	2375	2376	rBV4	17662	39274	1.81%	0.237%
75	16.103	2395	2400	2406	rBV4	41514	79508	3.67%	0.479%
76	16.186	2408	2414	2420	rBV5	11190	33366	1.54%	0.201%

LSC Area Percent Report

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Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

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\82N060822W.M
Title : SW846 8260

77	16.339	2431	2440	2447	rBV6	44243	113722	5.25%	0.685%
78	16.545	2465	2475	2480	rBV6	22276	67149	3.10%	0.405%
79	16.615	2481	2487	2503	rVB6	34195	135183	6.24%	0.815%

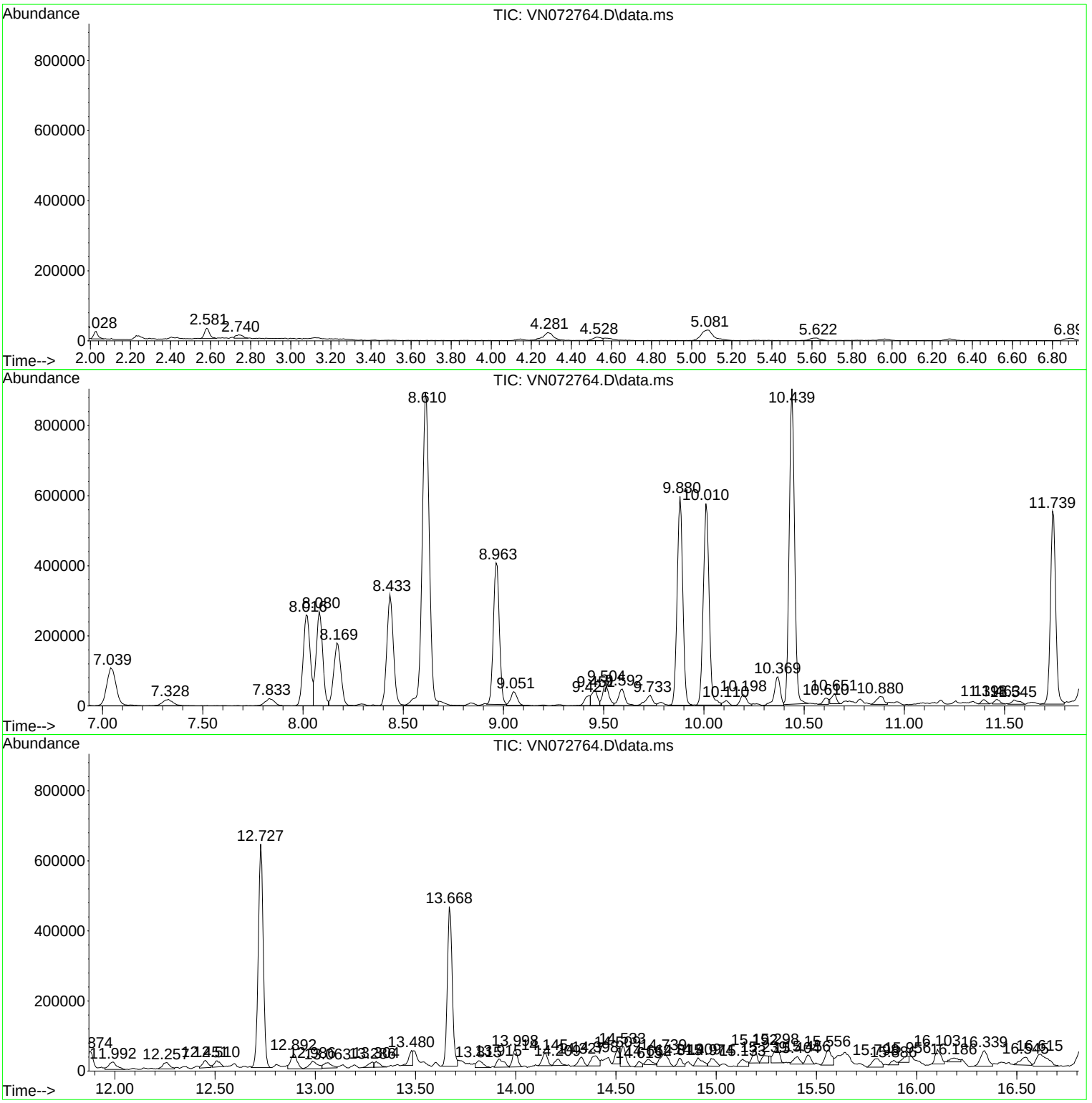
Sum of corrected areas: 16592778

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

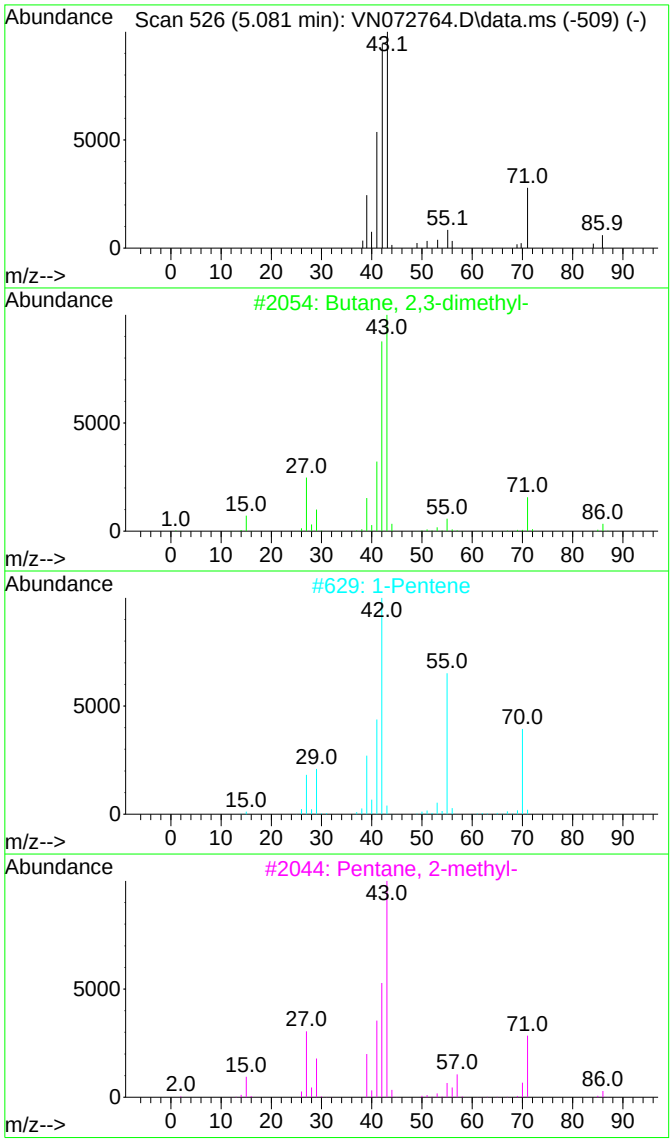
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

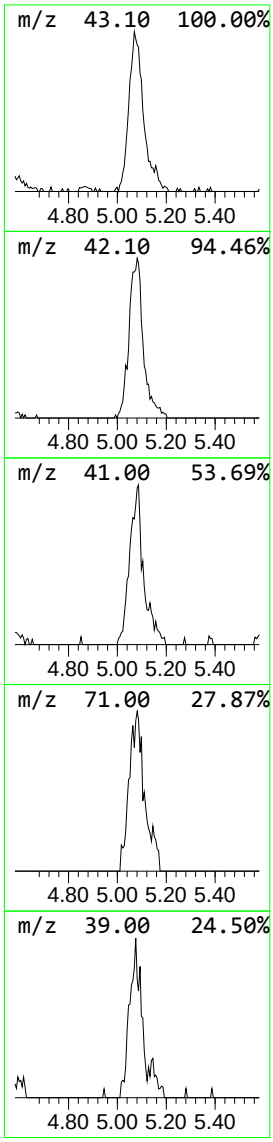
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.081	10.88 ug/l	130893	Pentafluorobenzene	8.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	1-Pentene	70	C5H10	000109-67-1	33
3	Pentane, 2-methyl-	86	C6H14	000107-83-5	33
4	Cyclobutane, methyl-	70	C5H10	000598-61-8	9
5	Pentane, 2-bromo-	150	C5H11Br	000107-81-3	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

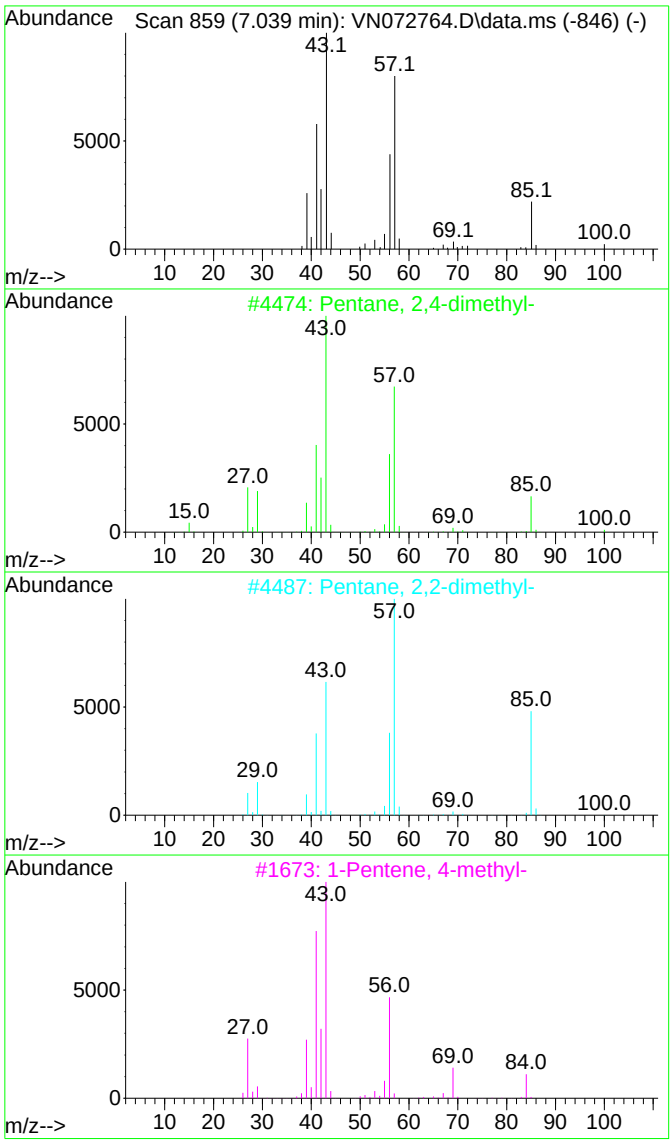
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

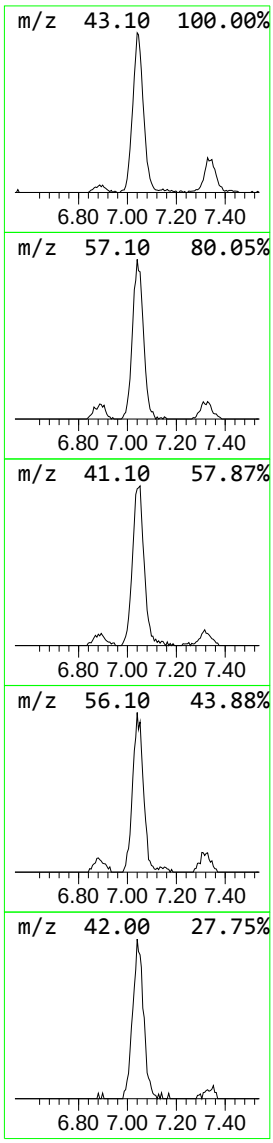
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.039	28.98 ug/l	348559	Pentafluorobenzene	8.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64
3	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	47
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	45
5	Hexane, 2-methyl-	100	C7H16	000591-76-4	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

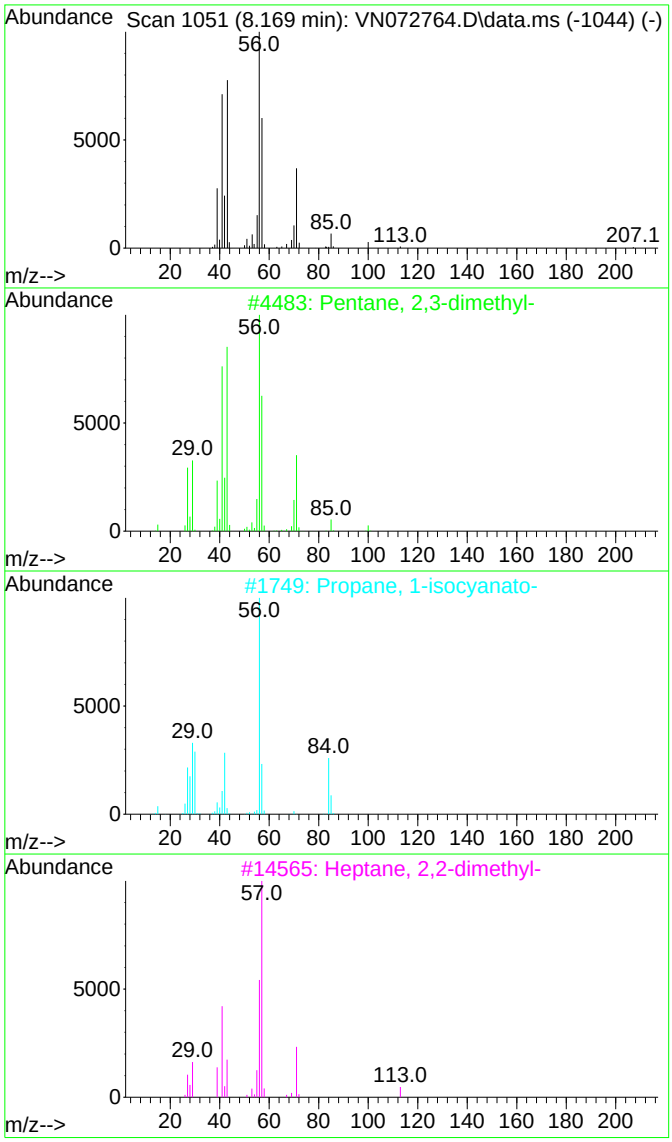
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

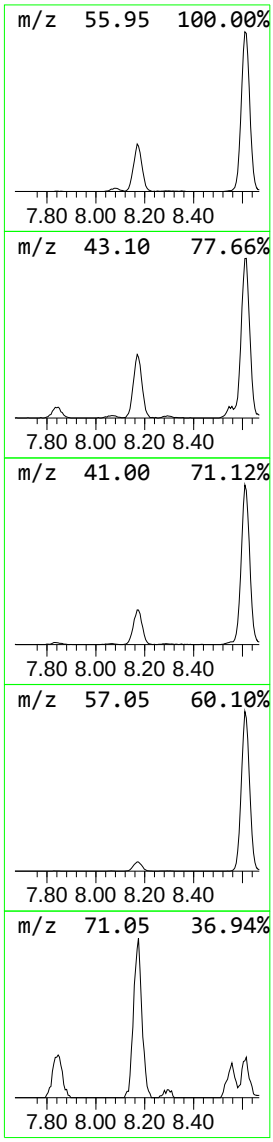
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.169	36.57 ug/l	439792	Pentafluorobenzene	8.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	50
3	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	40
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	40
5	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

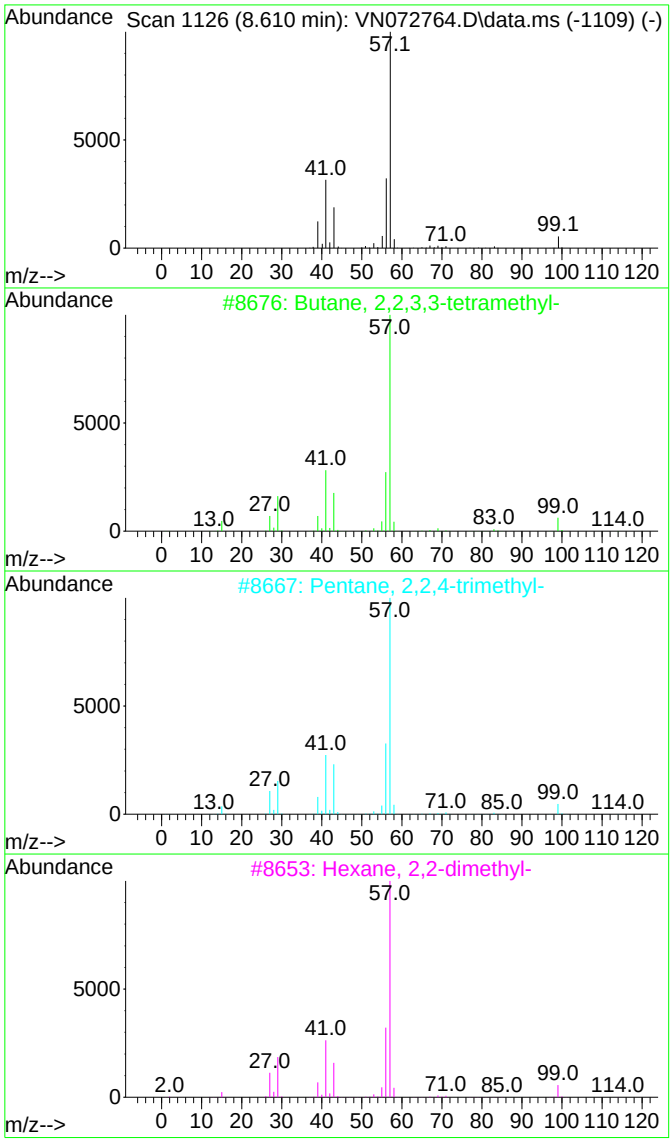
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

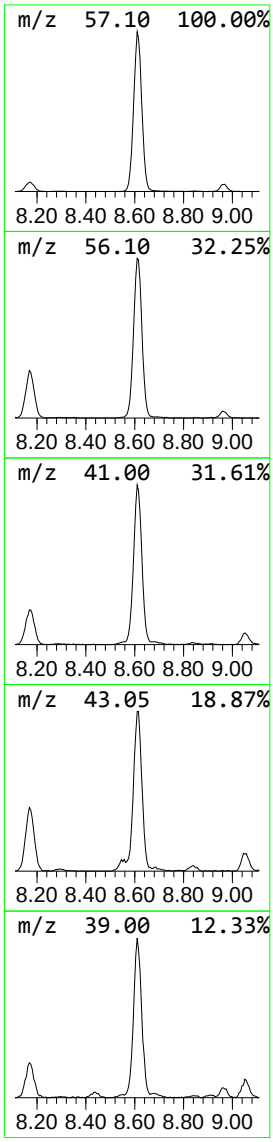
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.610	129.43 ug/l	2167250	1,4-Difluorobenzene	8.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	53
5	Decane, 2,2,4-trimethyl-	184	C13H28	062237-98-3	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
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Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

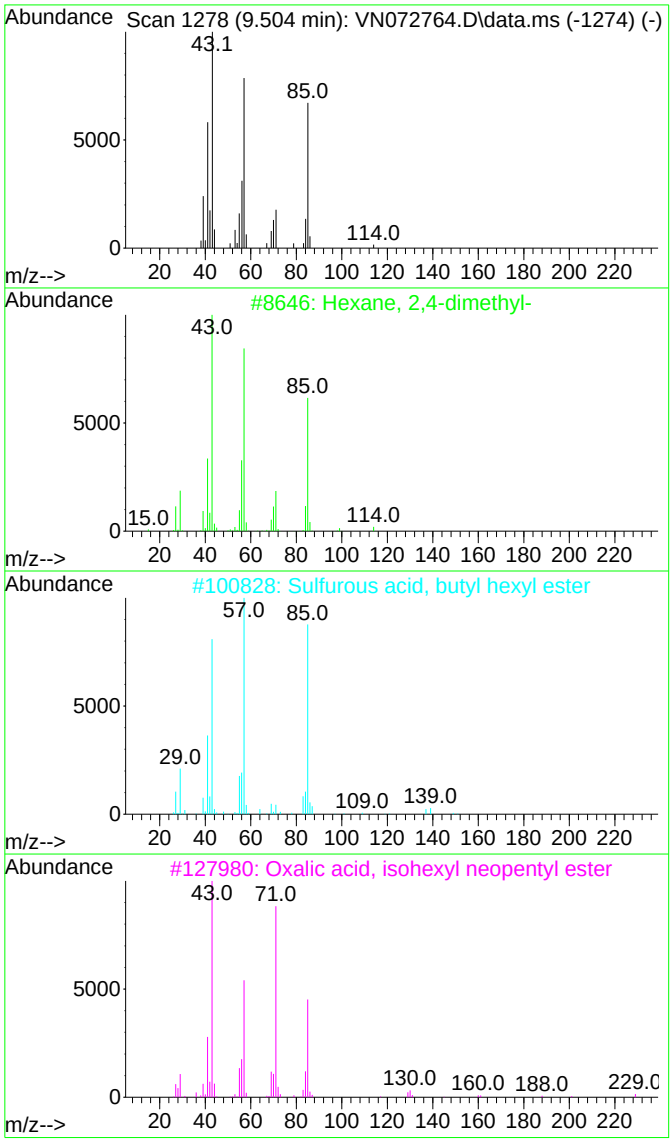
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

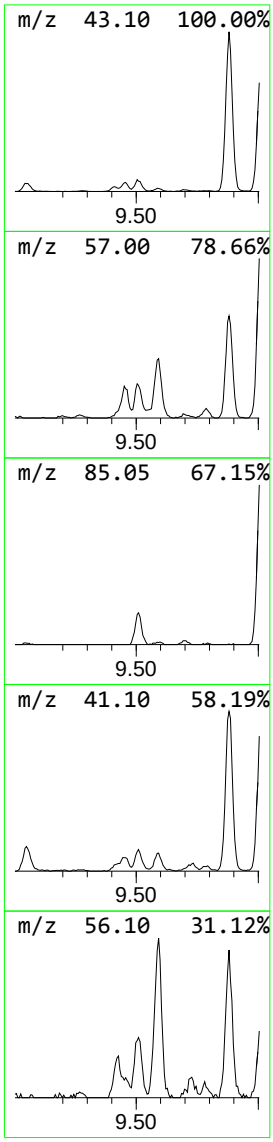
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexane, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.504	7.53 ug/l	126028	1,4-Difluorobenzene	8.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
2	Sulfurous acid, butyl hexyl ester	222	C10H22O3S	1000309-17-3	64
3	Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	59
4	Pentanoic acid, 2,2-dimethyl-, e...	156	C9H16O2	044970-05-0	53
5	Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

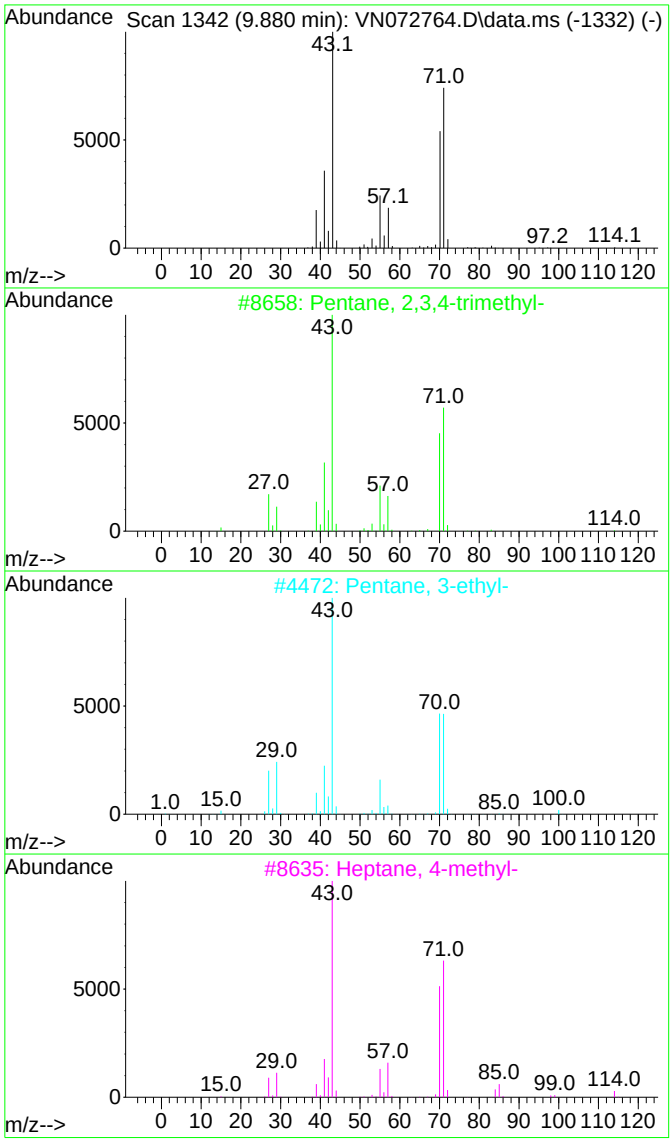
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

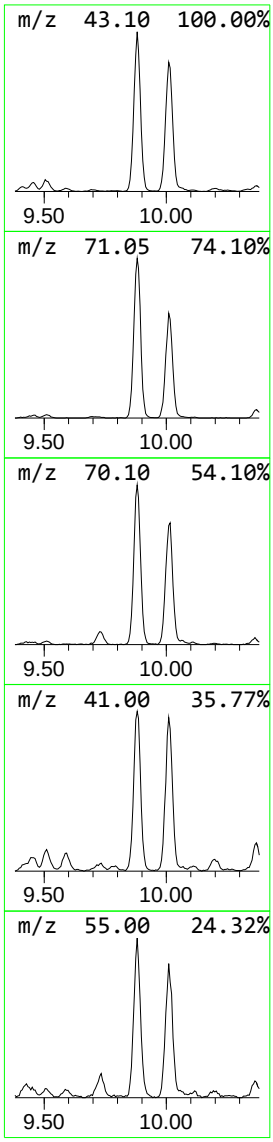
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.880	69.32 ug/l	1160710	1,4-Difluorobenzene	8.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	74
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64
5	Ethanedioic acid, bis(3-methylbu...	230	C12H22O4	002051-00-5	59





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Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

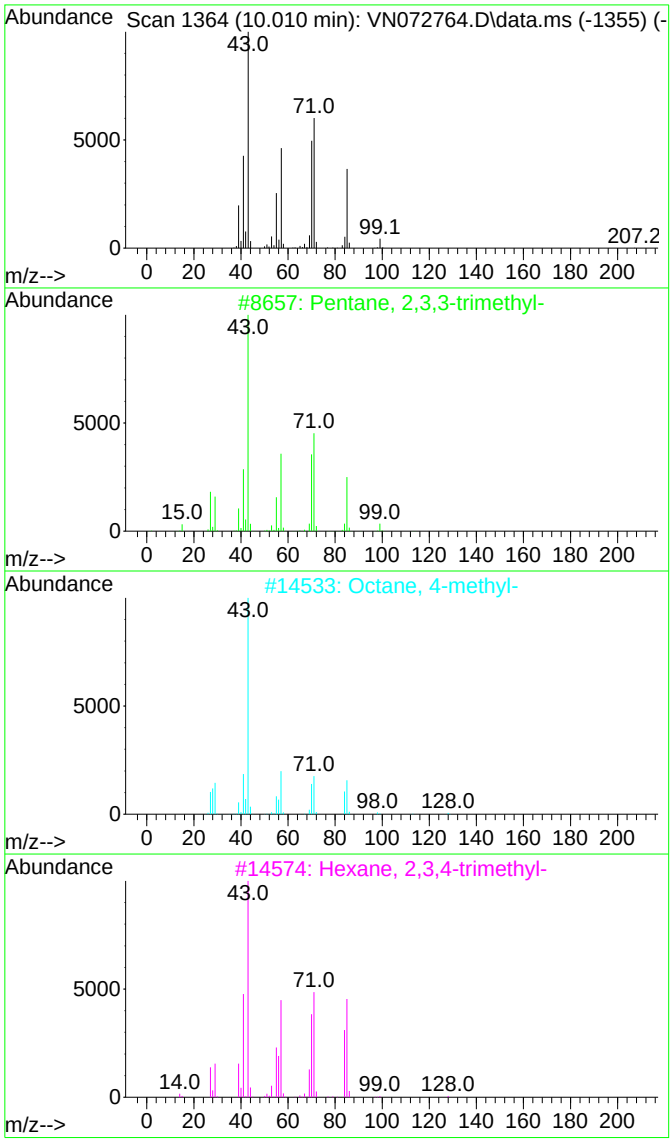
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

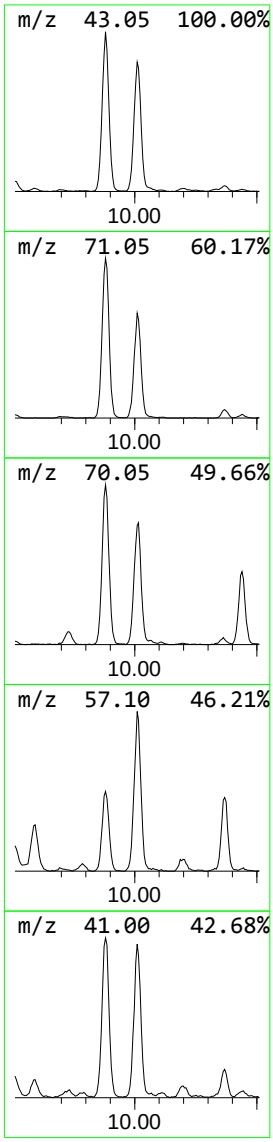
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.010	70.18 ug/l	1175190	1,4-Difluorobenzene	8.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2	Octane, 4-methyl-	128	C9H20	002216-34-4	78
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59





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Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

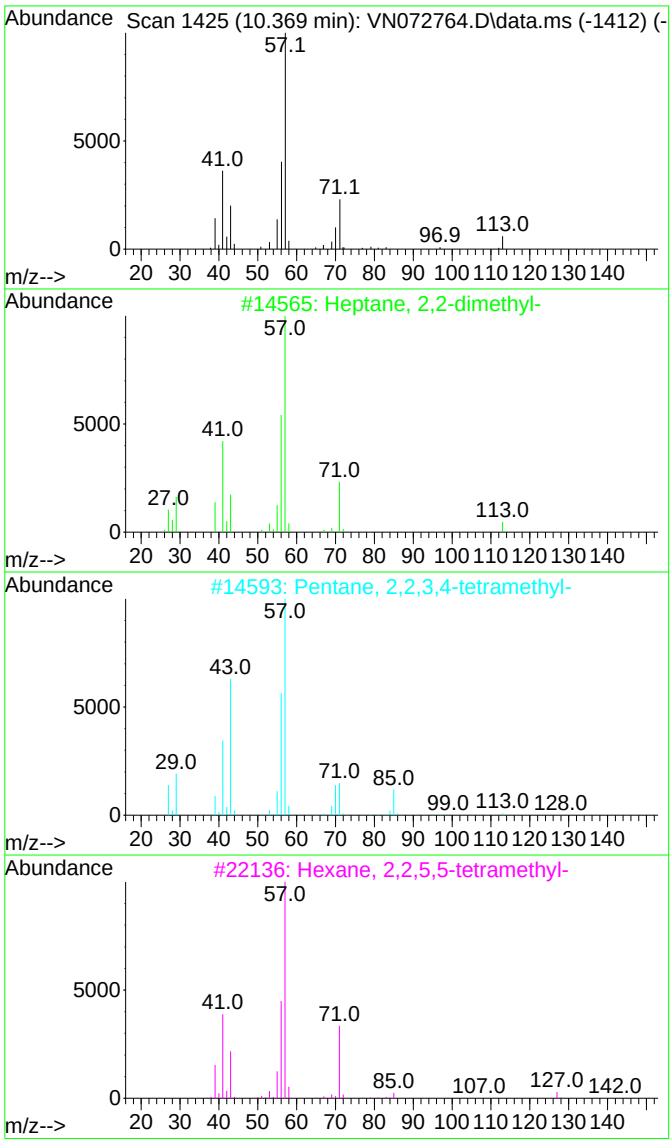
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

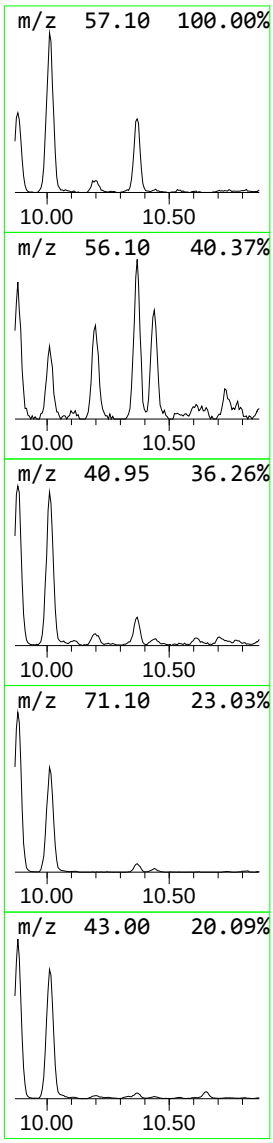
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Heptane, 2,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.369	8.51 ug/l	169366	Chlorobenzene-d5	11.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	72
2	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	64
3	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	64
4	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	56
5	Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	50





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Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

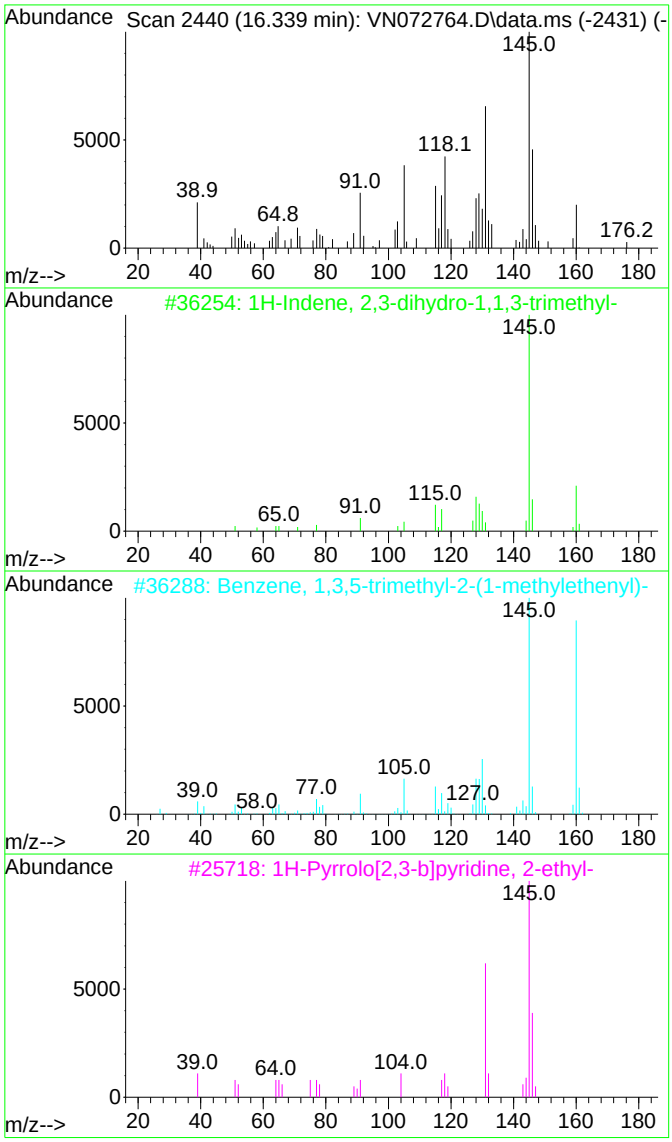
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

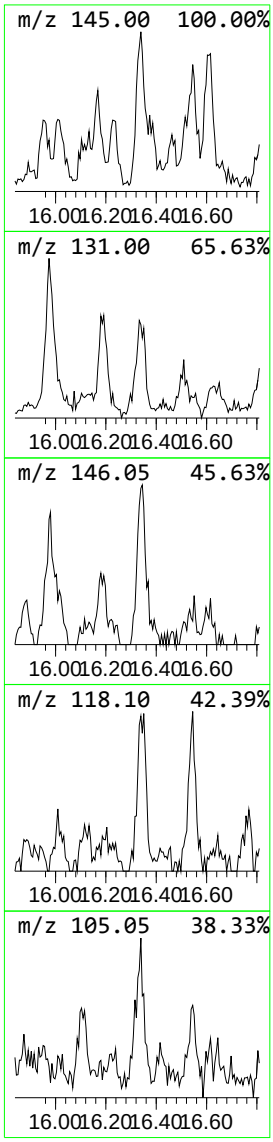
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.339	7.30 ug/l	113722	1,4-Dichlorobenzene-d4	13.668

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	55
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	55
3	1H-Pyrrolo[2,3-b]pyridine, 2-ethyl-	146	C9H10N2	023612-49-9	53
4	1H-Indene, 2,3-dihydro-1,4,7-tri...	160	C12H16	054340-87-3	50
5	1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	49





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Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

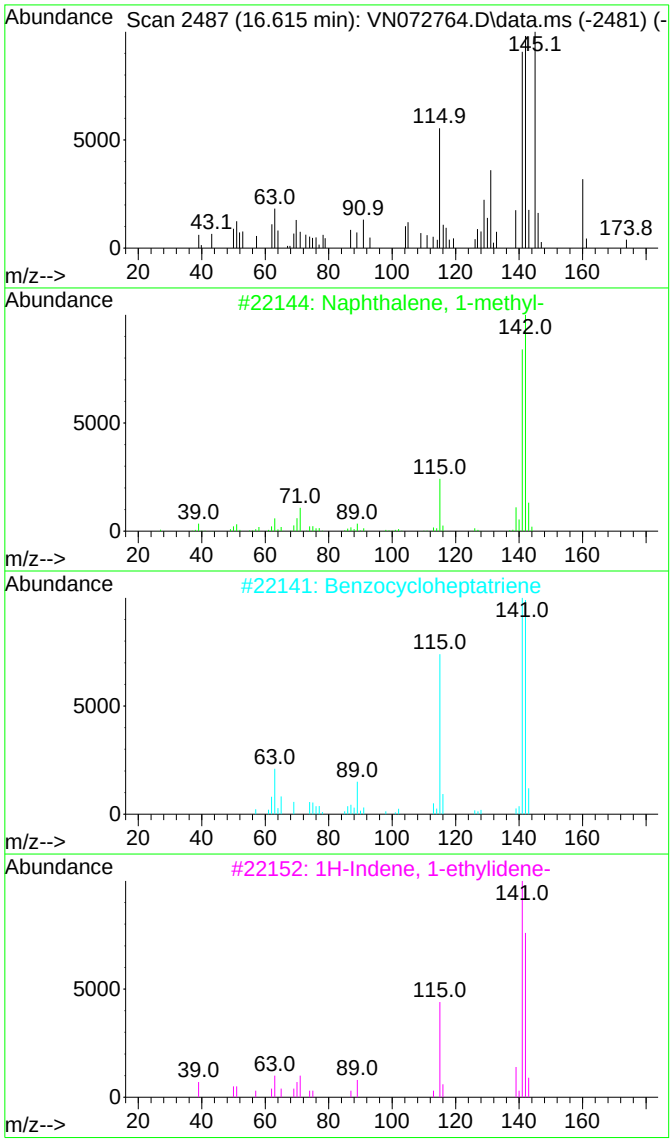
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

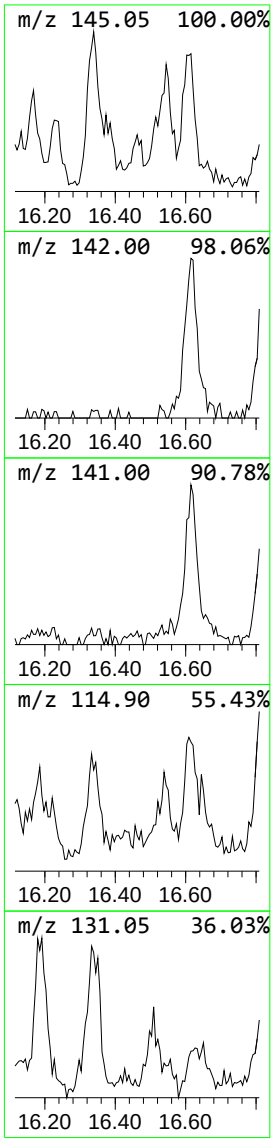
****

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.615	8.68 ug/l	135183	1,4-Dichlorobenzene-d4	13.668

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	86
2	Benzocycloheptatriene	142	C11H10	000264-09-5	60
3	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	53
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	50
5	Ethanone, 1-(2,3-dihydro-1H-inde...	160	C11H12O	004228-10-8	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072764.D
Acq On : 10 Jun 2022 14:57
Operator : JC\MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.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, 2,3-dim...	5.081	10.9	ug/l	130893	1	8.080	601354	50.0
Pentane, 2,4-di...	7.039	29.0	ug/l	348559	1	8.080	601354	50.0
Pentane, 2,3-di...	8.169	36.6	ug/l	439792	1	8.080	601354	50.0
Butane, 2,2,3,3...	8.610	129.4	ug/l	2167250	2	8.963	837231	50.0
Hexane, 2,4-dim...	9.504	7.5	ug/l	126028	2	8.963	837231	50.0
Pentane, 2,3,4-...	9.880	69.3	ug/l	1160710	2	8.963	837231	50.0
Pentane, 2,3,3-...	10.010	70.2	ug/l	1175190	2	8.963	837231	50.0
Heptane, 2,2-di...	10.369	8.5	ug/l	169366	3	11.739	995405	50.0
1H-Indene, 2,3-...	16.339	7.3	ug/l	113722	4	13.668	778896	50.0
Naphthalene, 1-...	16.615	8.7	ug/l	135183	4	13.668	778896	50.0