LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

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\82N033022W.M
Title : SW846 8260

Signal : TIC: VN071756.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	5	7	19	rVB3	29499	64027	1.62%	0.136%
2	2.257	38	46	60	rBV3	152740	395309	10.03%	0.837%
3	2.440	72	77	85	rVB	132321	210948	5.35%	0.447%
4	3.134	186	195	218	rVB	1283844	2945161	74.69%	6.237%
5	3.516	251	260	270	rBV2	38662	97460	2.47%	0.206%
6	3.998	329	342	351	rBV3	24374	68163	1.73%	0.144%
7	4.263	371	387	404	rBV2	215644	779818	19.78%	1.651%
8	4.569	429	439	449	rBV3	12861	43396	1.10%	0.092%
9	5.081	510	526	537	rBV2	415612	1675002	42.48%	3.547%
10	5.622	603	618	632	rBV2	167943	579043	14.69%	1.226%
11	6.487	755	765	776	rVB8	12368	46319	1.17%	0.098%
12	6.610	776	786	795	rBV4	18408	56435	1.43%	0.120%
13	6.886	823	833	845	rBV3	37652	132215	3.35%	0.280%
14	7.045	849	860	869	rVV2	143887	453257	11.50%	0.960%
15	7.145	869	877	886	rVV2	216719	633993	16.08%	1.343%
16	7.222	886	890	900	rVV3	43105	111495	2.83%	0.236%
17	7.322	900	907	918	rVB6	19840	64616	1.64%	0.137%
18	7.651	956	963	977	rBV4	17422	58119	1.47%	0.123%
19	7.839	977	995	1005	rBV4	70560	230713	5.85%	0.489%
20	8.022	1015	1026	1031	rBV	442736	1087559	27.58%	2.303%
21	8.086	1031	1037	1045	rVV	887151	2087391	52.94%	4.420%
22	8.175	1045	1052	1064	rVB	452995	1118788	28.37%	2.369%
23	8.357	1076	1083	1088	rBV4	27432	60159	1.53%	0.127%
24	8.451	1088	1099	1109	rVV3	994737	2875980	72.94%	6.090%
25	8.616	1111	1127	1136	rVV	515678	1541300	39.09%	3.264%
26	8.704	1136	1142	1155	rVB	123599	296522	7.52%	0.628%
27	8.963	1177	1186	1198	rBV	1253950	2650388	67.22%	5.612%
28	9.057	1198	1202	1209	rVB3	29941	63859	1.62%	0.135%
29	9.198	1220	1226	1231	rBV2	42335	85238	2.16%	0.180%
30	9.422	1254	1264	1275	rBV3	181890	524915	13.31%	1.112%
31	9.510	1275	1279	1287	rVB6	31734	64042	1.62%	0.136%
32	9.592	1287	1293	1297	rBV	25263	47330	1.20%	0.100%
33	9.727	1306	1316	1323	rBV3	53593	125474	3.18%	0.266%
34	9.880	1336	1342	1351	rVB	226875	449861	11.41%	0.953%
35	9.969	1351	1357	1359	rVV	92211	151564	3.84%	0.321%
36	10.010	1359	1364	1377	rVV	279305	611959	15.52%	1.296%

LSC Area Percent Report

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Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

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\82N033022W.M
Title : SW846 8260

37	10.198	1388	1396	1402	rBV4	72410	152556	3.87%	0.323%
38	10.322	1411	1417	1420	rVV	59523	118196	3.00%	0.250%
39	10.363	1420	1424	1429	rVV4	55962	107928	2.74%	0.229%
40	10.439	1429	1437	1458	rVV	2021028	3943065	100.00%	8.350%
41	10.639	1458	1471	1478	rVV	346387	740161	18.77%	1.567%
42	10.710	1478	1483	1490	rVV	288327	521921	13.24%	1.105%
43	10.774	1490	1494	1500	rVV2	38995	79846	2.02%	0.169%
44	10.857	1500	1508	1517	rVV3	230158	545624	13.84%	1.155%
45	10.939	1517	1522	1532	rVB10	24398	63265	1.60%	0.134%
46	11.127	1539	1554	1558	rBV6	40019	118519	3.01%	0.251%
47	11.251	1570	1575	1581	rVB3	54628	99922	2.53%	0.212%
48	11.451	1604	1609	1615	rVB3	32926	56492	1.43%	0.120%
49	11.551	1619	1626	1634	rBV9	15364	43218	1.10%	0.092%
50	11.663	1636	1645	1650	rBV8	15650	43699	1.11%	0.093%
51	11.739	1650	1658	1669	rBV	1905942	3444201	87.35%	7.293%
52	11.839	1669	1675	1685	rVB	491742	889266	22.55%	1.883%
53	11.957	1689	1695	1698	rBV2	25220	52782	1.34%	0.112%
54	12.574	1793	1800	1810	rBV	507579	840176	21.31%	1.779%
55	12.727	1819	1826	1836	rVB	1433052	2445067	62.01%	5.178%
56	12.916	1851	1858	1868	rVB	510990	833485	21.14%	1.765%
57	13.215	1903	1909	1917	rBV	60886	104144	2.64%	0.221%
58	13.480	1947	1954	1962	rVV2	104079	231156	5.86%	0.489%
59	13.668	1979	1986	1998	rBV	1530019	2767202	70.18%	5.860%
60	13.768	1998	2003	2007	rVB	38580	62529	1.59%	0.132%
61	13.833	2008	2014	2016	rBV	208324	320801	8.14%	0.679%
62	13.874	2016	2021	2027	rVB	1742184	2794170	70.86%	5.917%
63	13.998	2036	2042	2048	rVB	188244	303284	7.69%	0.642%
64	14.068	2048	2054	2061	rVB	145551	241665	6.13%	0.512%
65	14.145	2061	2067	2073	rBV	137641	235908	5.98%	0.500%
66	14.210	2073	2078	2083	rVB	54666	84962	2.15%	0.180%
67	14.274	2083	2089	2092	rBV	83953	140549	3.56%	0.298%
68	14.327	2092	2098	2104	rVV	515201	876415	22.23%	1.856%
69	14.404	2104	2111	2116	rVV4	64831	141613	3.59%	0.300%
70	14.462	2116	2121	2125	rVV	65486	96820	2.46%	0.205%
71	14.533	2125	2133	2139	rVV	129192	234544	5.95%	0.497%
72	14.651	2149	2153	2158	rVB	44082	64580	1.64%	0.137%
73	14.757	2163	2171	2176	rBV2	69983	110778	2.81%	0.235%
74	14.815	2176	2181	2191	rVB4	69262	154702	3.92%	0.328%
75	14.927	2191	2200	2207	rBV6	37407	93713	2.38%	0.198%
76	14.998	2207	2212	2217	rBV3	41069	70509	1.79%	0.149%

LSC Area Percent Report

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Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

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\82N033022W.M
Title : SW846 8260

77	15.192	2239	2245	2250	rBV2	44866	86941	2.20%	0.184%
78	15.298	2254	2263	2274	rVB5	70522	213355	5.41%	0.452%
79	15.562	2301	2308	2313	rBV4	21844	44521	1.13%	0.094%
80	15.633	2313	2320	2330	rVB5	25309	67094	1.70%	0.142%
81	16.109	2394	2401	2407	rBV2	25806	55193	1.40%	0.117%

Sum of corrected areas: 47224355

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
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Operator : JC\MD
Sample : N2090-05
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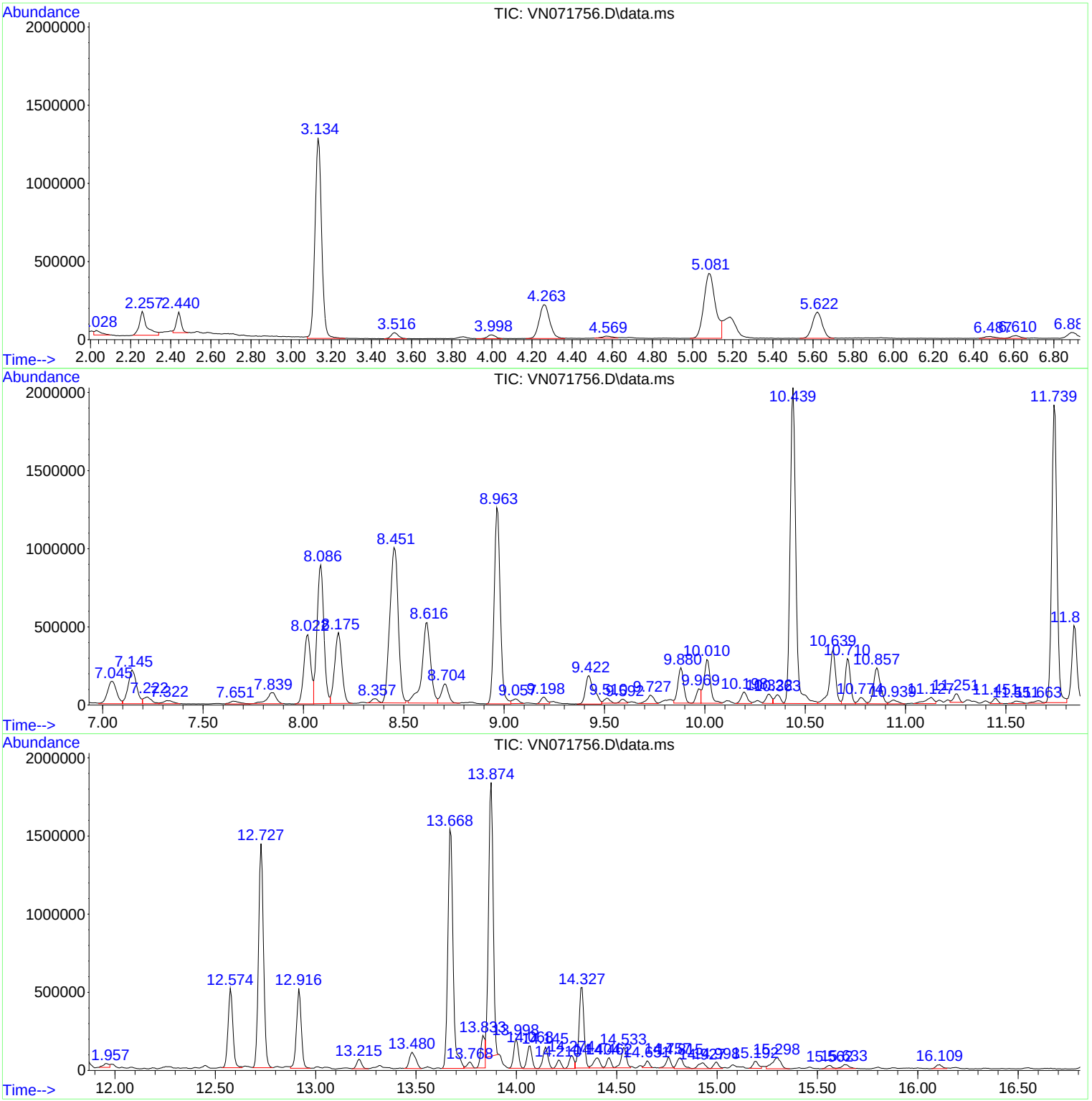
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

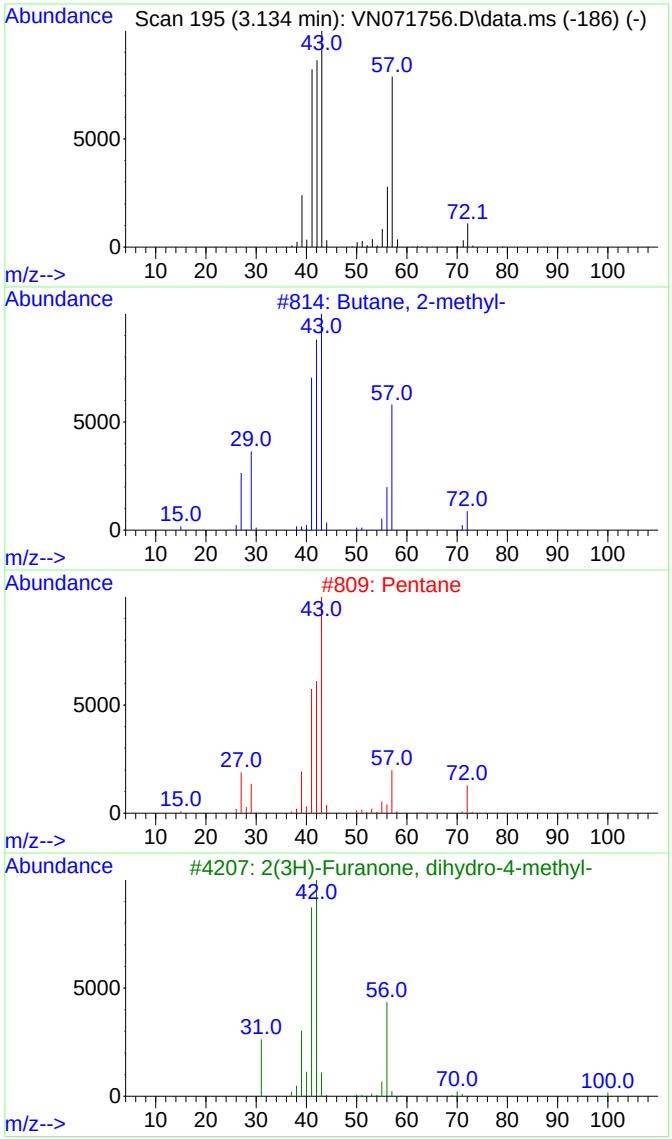
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

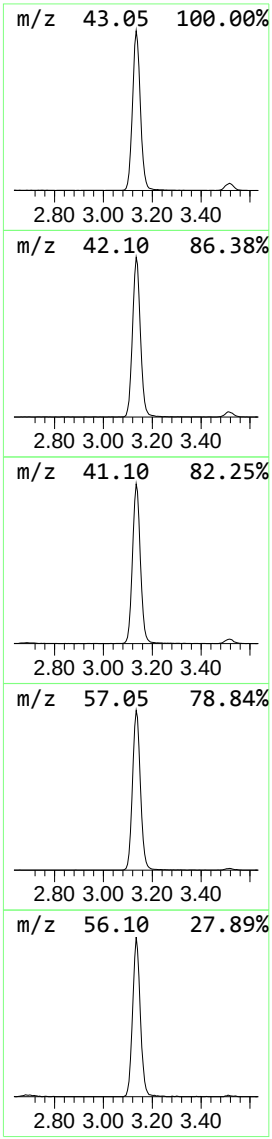
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.134	70.55 ug/l	2945160	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	86
2	Pentane	72	C5H12	000109-66-0	46
3	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	36
4	3-Buten-1-ol	72	C4H8O	000627-27-0	25
5	1-Butanesulfonyl chloride	156	C4H9ClO2S	002386-60-9	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
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Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

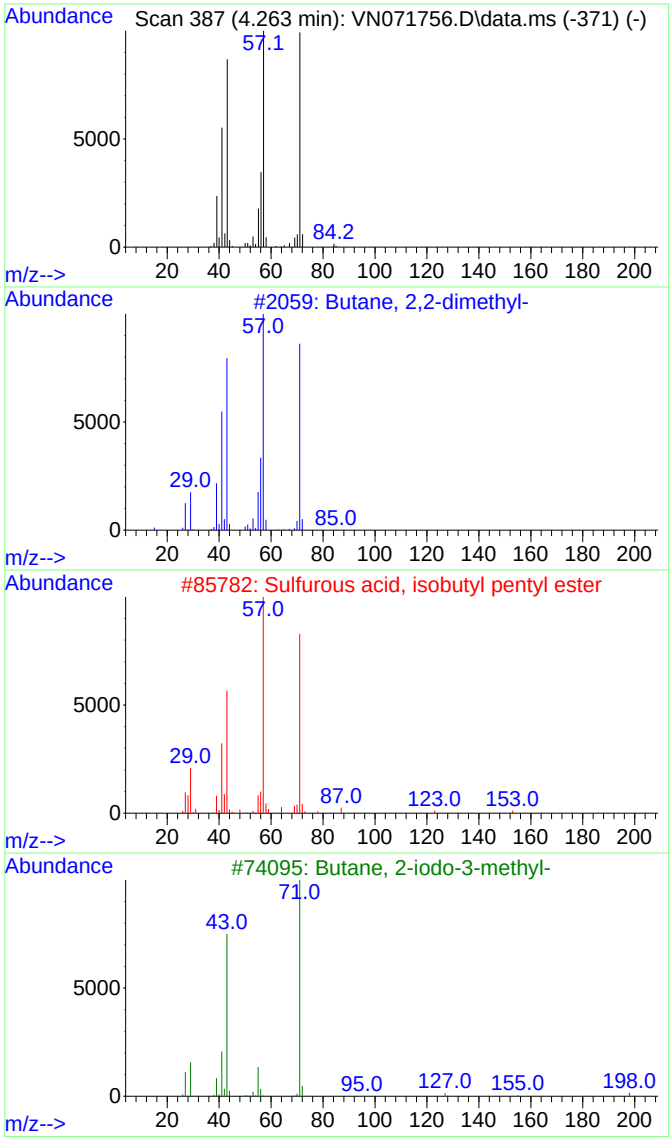
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

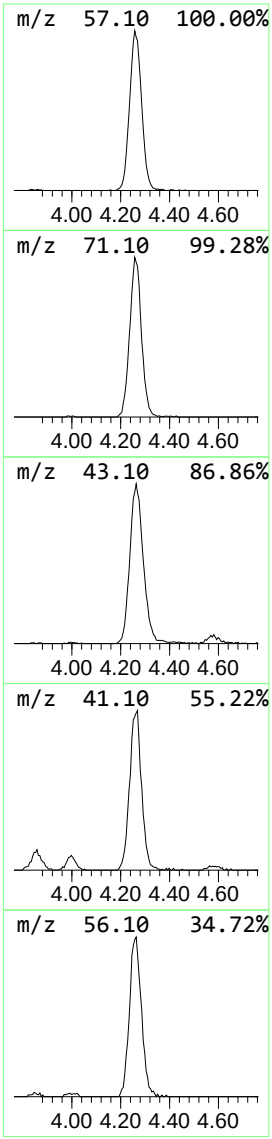
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.263	18.68 ug/l	779818	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	90
2	Sulfurous acid, isobutyl pentyl ...	208	C9H20O3S	1000309-13-8	64
3	Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	42
4	Pentane, 3-bromo-	150	C5H11Br	001809-10-5	40
5	Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
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Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

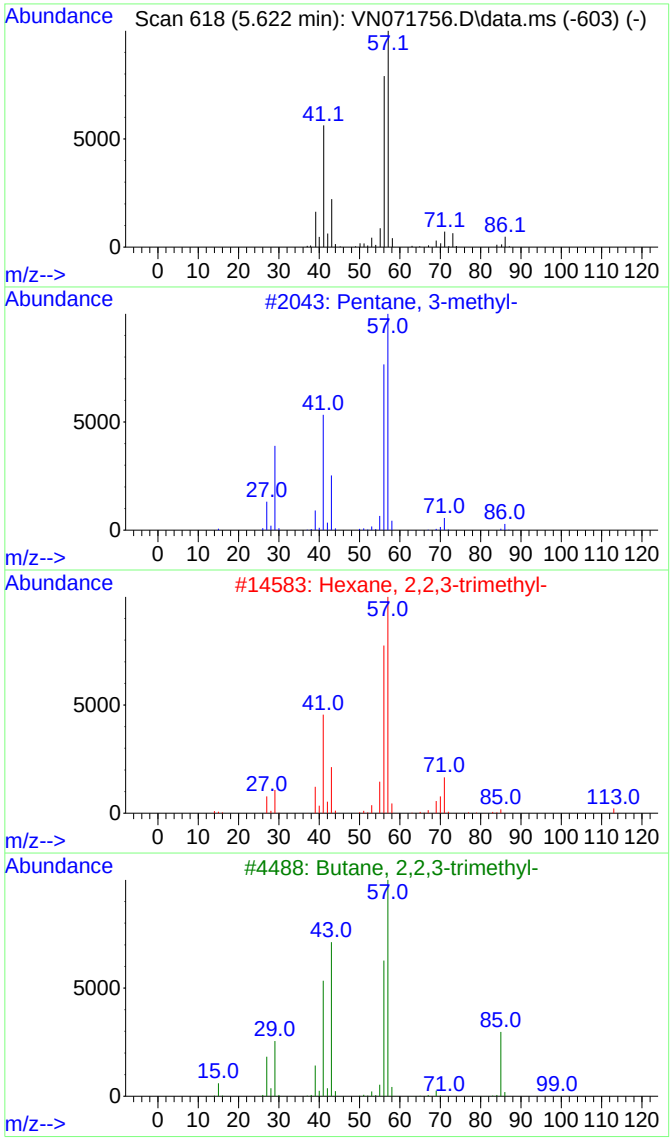
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

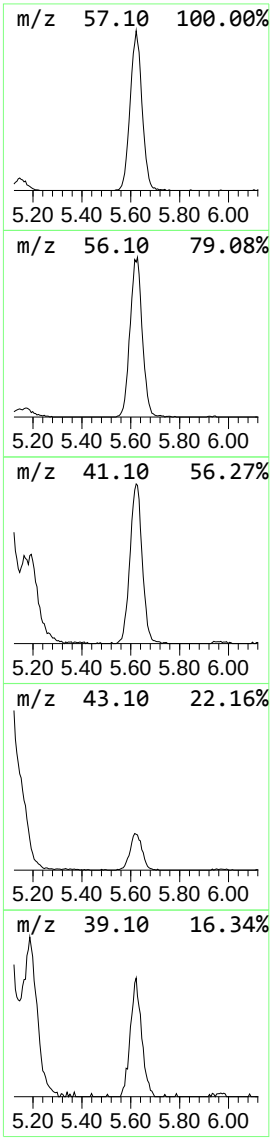
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.622	13.87 ug/l	579043	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	72
3	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	72
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	42
5	Butane, 1-chloro-	92	C4H9Cl	000109-69-3	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
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Instrument :
MSVOA_N
ClientSampleId :
MW-50

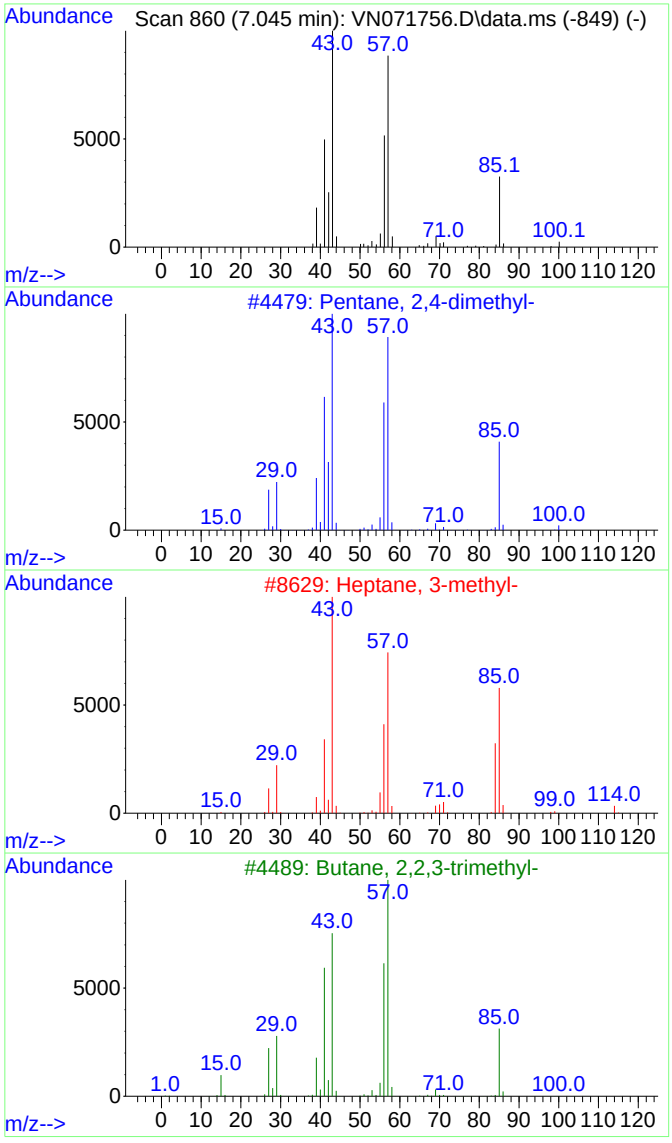
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

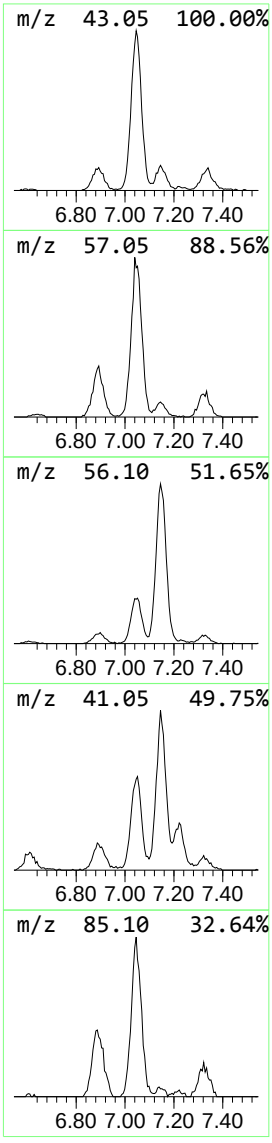
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.045	10.86 ug/l	453257	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Heptane, 3-methyl-	114	C8H18	000589-81-1	72
3	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59
4	Hexane, 2-methyl-	100	C7H16	000591-76-4	50
5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	43





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Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

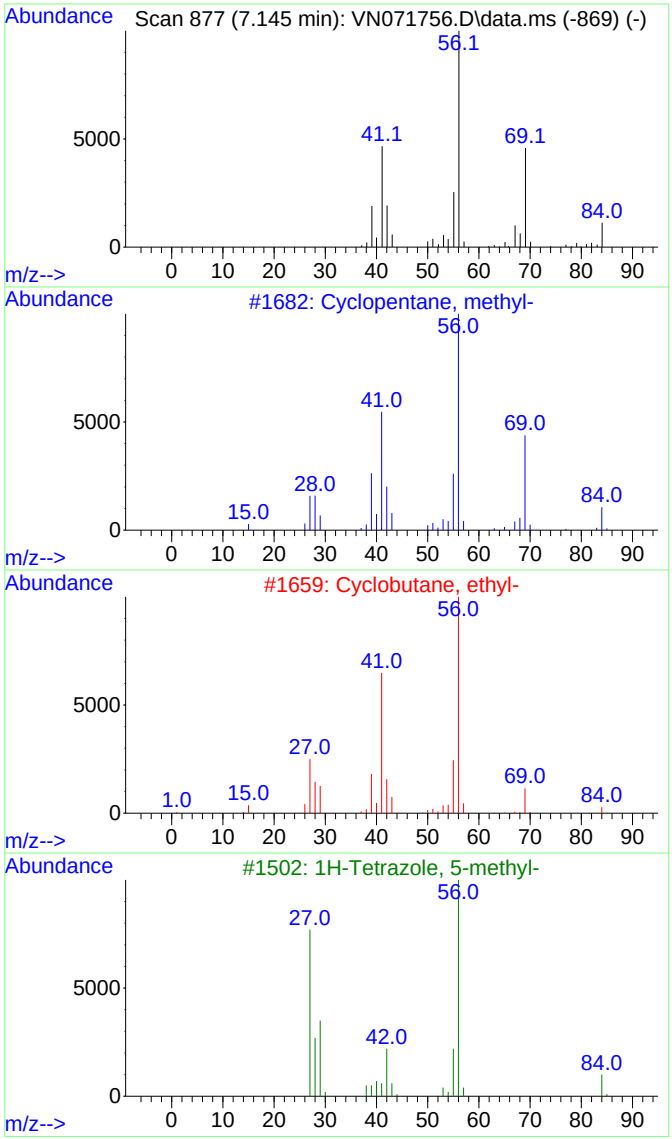
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

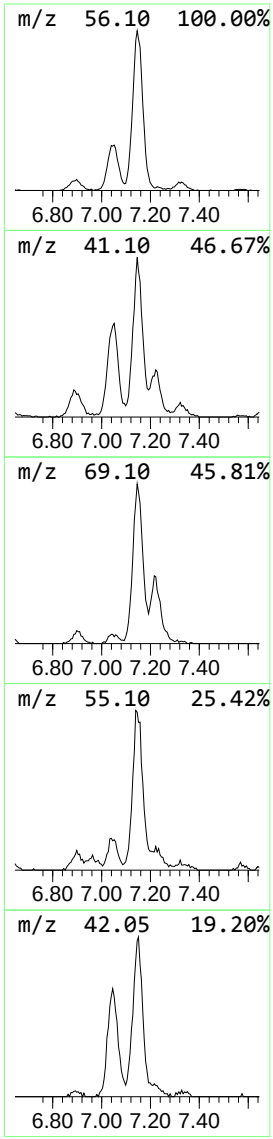
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.145	15.19 ug/l	633993	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43
5	1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

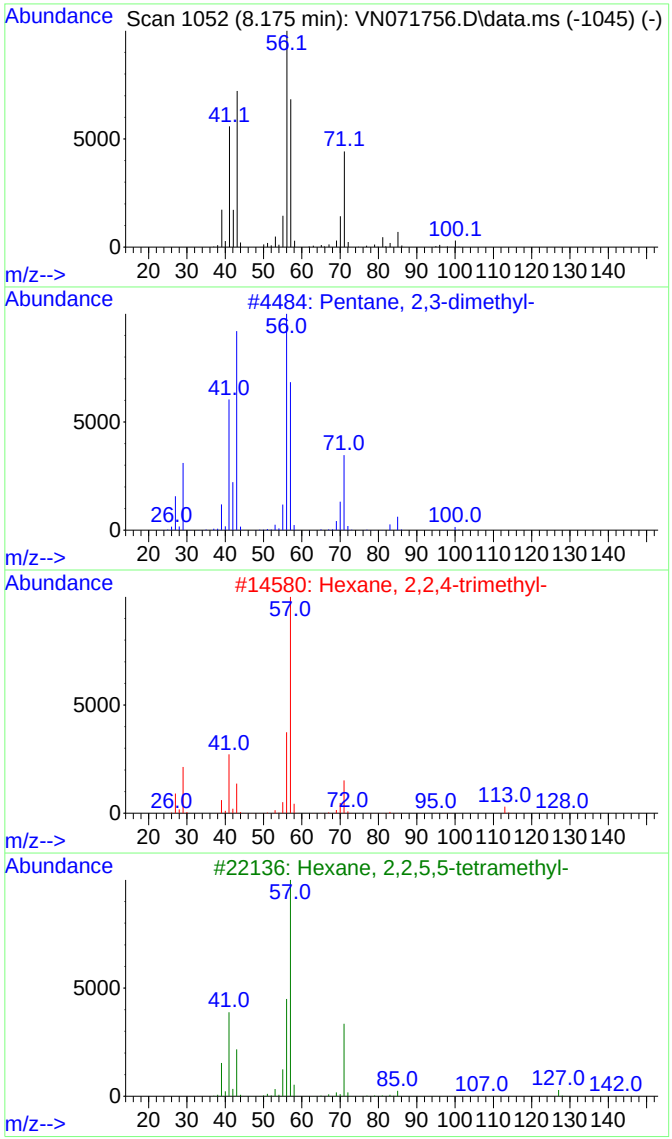
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

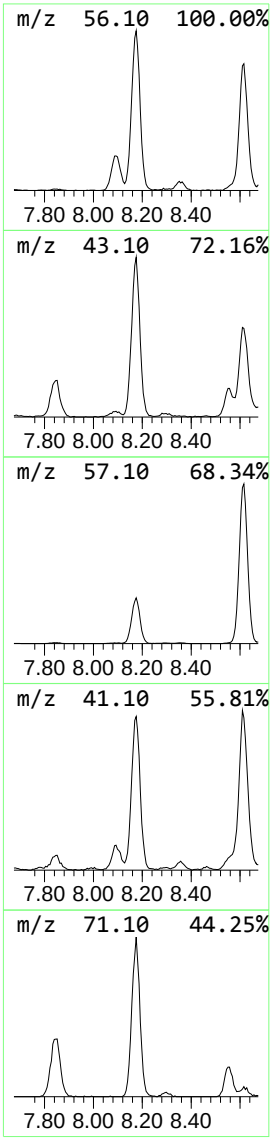
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.175	26.80 ug/l	1118790	Pentafluorobenzene	8.086

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	50
3	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	50
4	Heptane	100	C7H16	000142-82-5	43
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	40





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Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

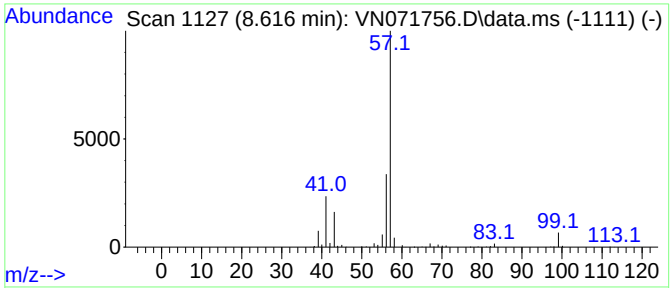
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

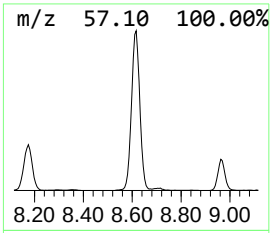
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

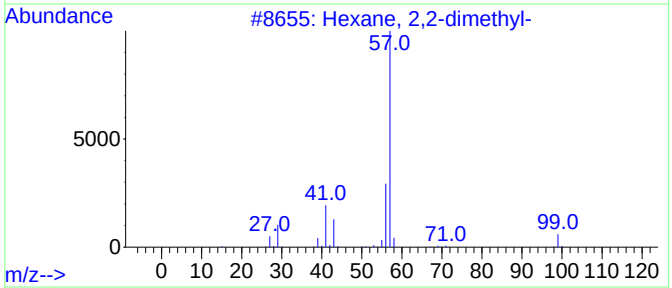
****
Peak Number 7 Hexane, 2,2-dimethyl- Concentration Rank 3

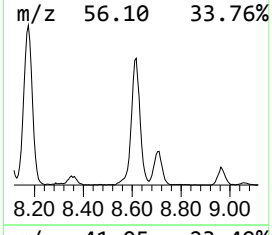
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.616	29.08 ug/l	1541300	1,4-Difluorobenzene	8.963

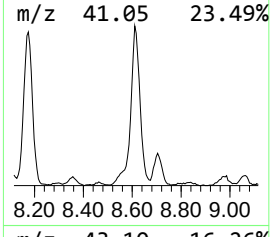
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59

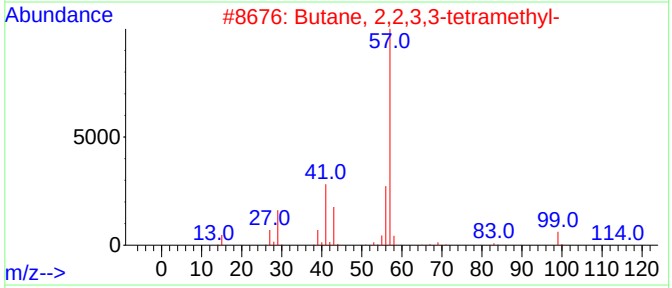


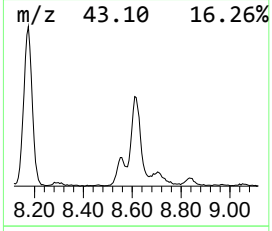


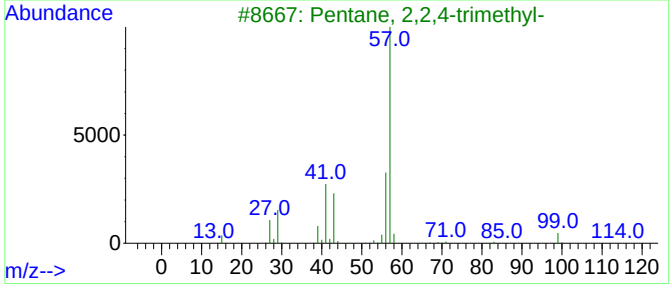


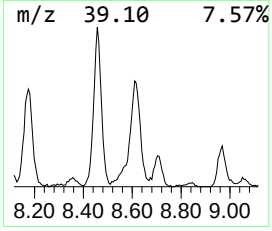












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Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

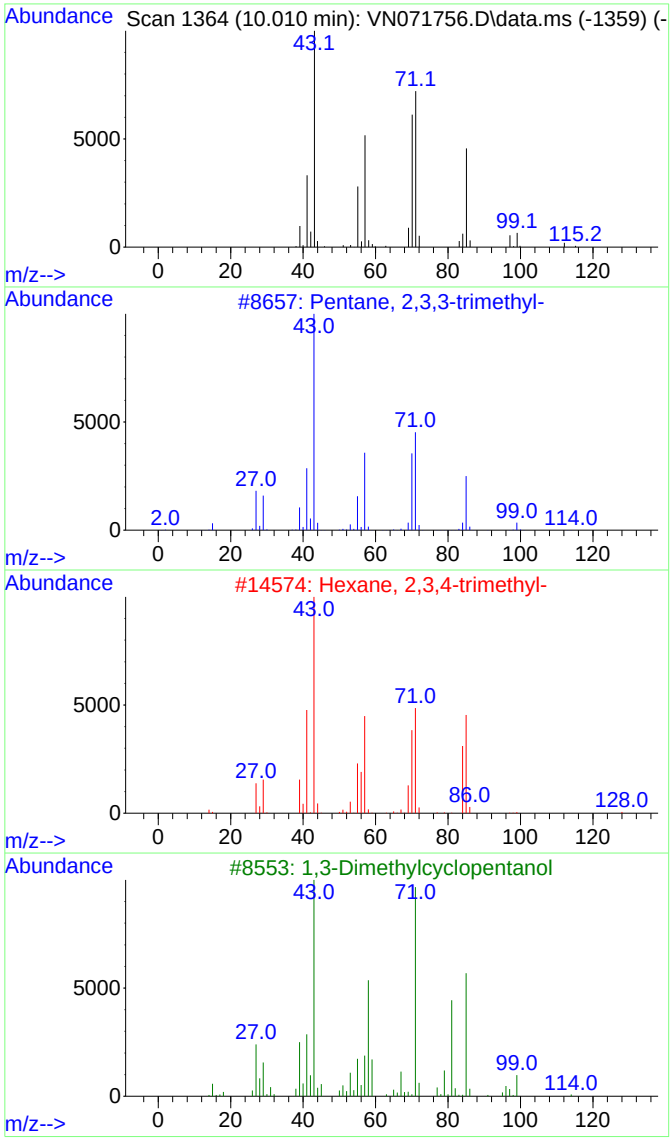
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

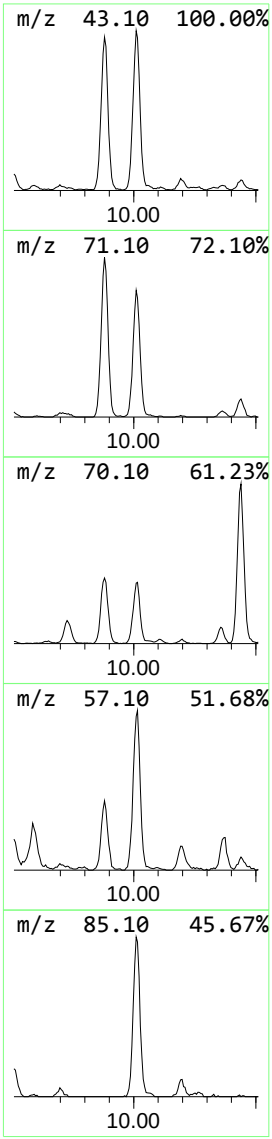
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Pentane, 2,3,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.010	11.54 ug/l	611959	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	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3	1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	53
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5	Propanoic acid, 2-methyl-, 2-eth...	200	C12H24O2	035061-61-1	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

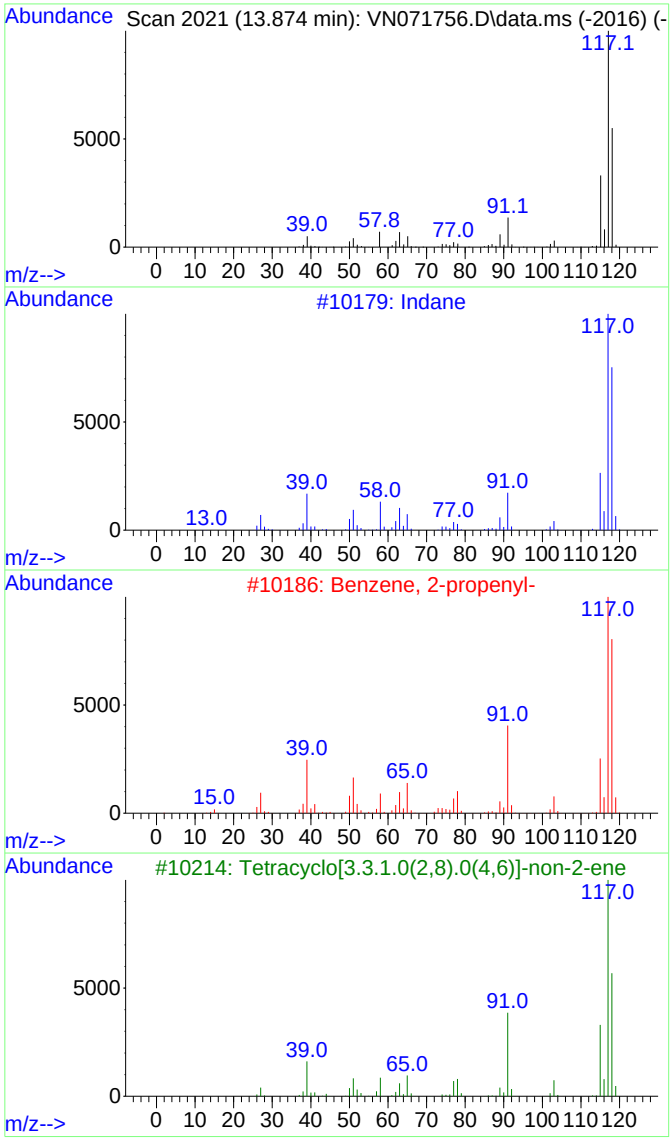
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

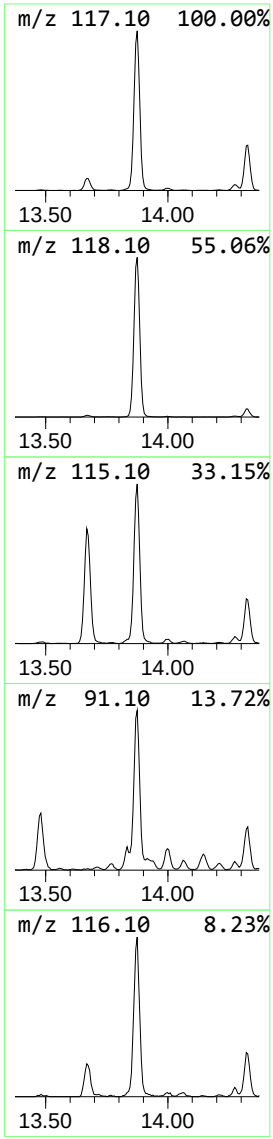
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	50.49 ug/l	2794170	1,4-Dichlorobenzene-d4	13.668

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

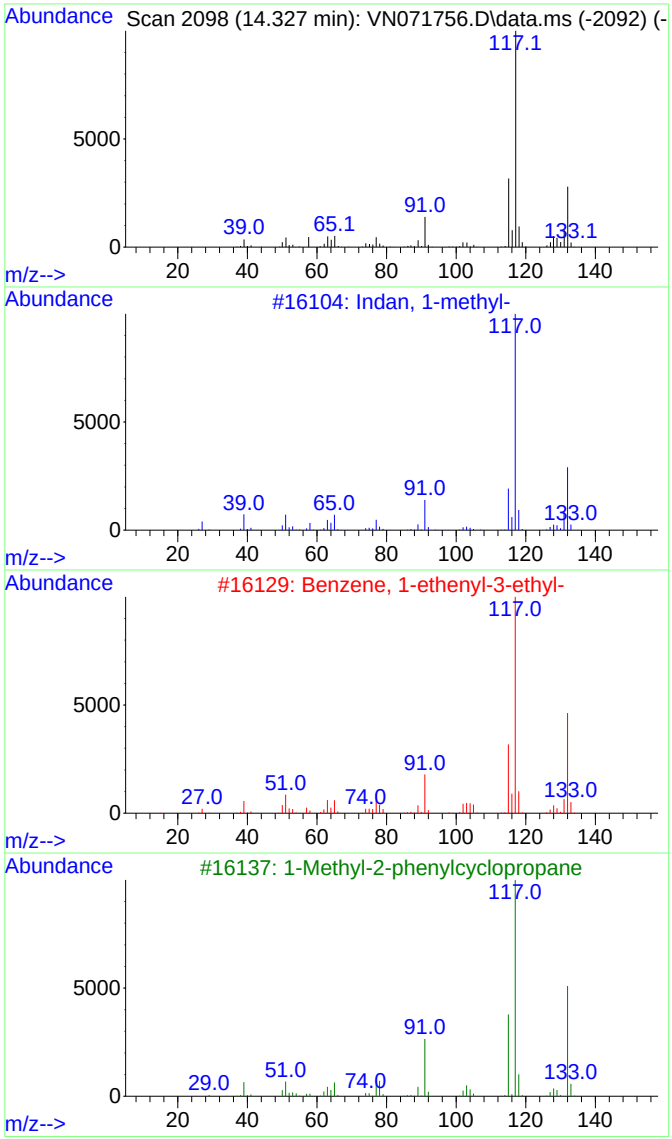
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

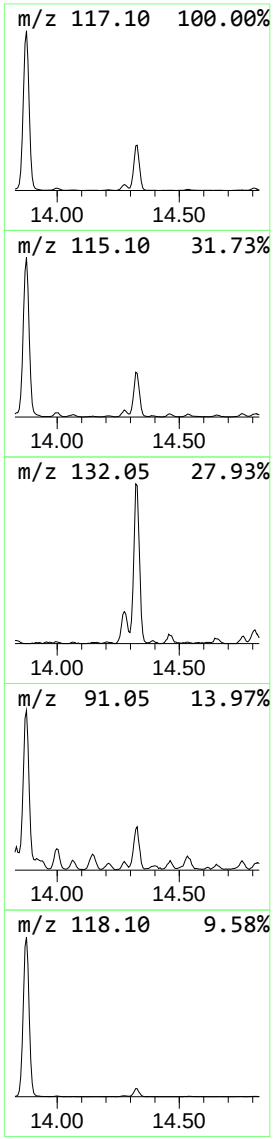
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.327	15.84 ug/l	876415	1,4-Dichlorobenzene-d4	13.668

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	86
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	80
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071756.D
Acq On : 01 Apr 2022 16:31
Operator : JC\MD
Sample : N2090-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-50

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.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-	3.134	70.5	ug/l	2945160	1	8.086	2087390	50.0
Butane, 2,2-dim...	4.263	18.7	ug/l	779818	1	8.086	2087390	50.0
Pentane, 3-methyl-	5.622	13.9	ug/l	579043	1	8.086	2087390	50.0
Pentane, 2,4-di...	7.045	10.9	ug/l	453257	1	8.086	2087390	50.0
Cyclopentane, m...	7.145	15.2	ug/l	633993	1	8.086	2087390	50.0
Pentane, 2,3-di...	8.175	26.8	ug/l	1118790	1	8.086	2087390	50.0
Hexane, 2,2-dim...	8.616	29.1	ug/l	1541300	2	8.963	2650390	50.0
Pentane, 2,3,3-...	10.010	11.5	ug/l	611959	2	8.963	2650390	50.0
Indane	13.874	50.5	ug/l	2794170	4	13.668	2767200	50.0
Indan, 1-methyl-	14.327	15.8	ug/l	876415	4	13.668	2767200	50.0