LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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: VN071814.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.234	35	42	62	rBV	425089	1669509	21.16%	1.878%
2	2.440	73	77	89	rVB2	100186	184067	2.33%	0.207%
3	3.134	186	195	208	rVB	814277	1858908	23.56%	2.091%
4	3.516	249	260	272	rBV3	190169	502086	6.36%	0.565%
5	3.999	331	342	356	rVB2	107464	298804	3.79%	0.336%
6	4.263	373	387	407	rBV2	108965	409656	5.19%	0.461%
7	5.081	513	526	530	rBV	145438	465600	5.90%	0.524%
8	5.151	531	538	564	rVV5	299999	1406049	17.82%	1.582%
9	5.363	564	574	590	rVB	89869	322572	4.09%	0.363%
10	5.622	604	618	634	rBV	335695	1167515	14.80%	1.314%
11	6.122	689	703	714	rBV2	53164	163168	2.07%	0.184%
12	6.469	747	762	775	rBV2	170162	521319	6.61%	0.587%
13	6.604	775	785	798	rVV2	87622	273255	3.46%	0.307%
14	6.898	823	835	851	rBV3	94346	301312	3.82%	0.339%
15	7.051	851	861	867	rBV3	39802	114912	1.46%	0.129%
16	7.145	867	877	887	rBV	562932	1655364	20.98%	1.862%
17	7.781	975	985	989	rBV	41977	105324	1.33%	0.118%
18	7.845	989	996	1005	rVB	141712	335699	4.25%	0.378%
19	8.022	1015	1026	1030	rBV	425841	1034481	13.11%	1.164%
20	8.087	1030	1037	1046	rVV2	1078875	2723514	34.51%	3.064%
21	8.175	1046	1052	1061	rVB3	141065	361338	4.58%	0.407%
22	8.298	1063	1073	1080	rBV2	130322	303642	3.85%	0.342%
23	8.445	1087	1098	1109	rBV3	679818	2104538	26.67%	2.368%
24	8.569	1109	1119	1122	rVV5	156510	443812	5.62%	0.499%
25	8.622	1122	1128	1137	rVV2	261965	913352	11.57%	1.028%
26	8.704	1137	1142	1152	rVB	86378	197968	2.51%	0.223%
27	8.804	1152	1159	1174	rVB5	40699	134371	1.70%	0.151%
28	8.963	1174	1186	1199	rBV	1365006	3101152	39.30%	3.489%
29	9.057	1199	1202	1209	rVV2	49656	94153	1.19%	0.106%
30	9.198	1219	1226	1230	rVV2	150537	316314	4.01%	0.356%
31	9.245	1230	1234	1246	rVB	114930	228206	2.89%	0.257%
32	9.457	1255	1270	1277	rBV4	295258	896690	11.36%	1.009%
33	9.639	1296	1301	1308	rVB	83917	157795	2.00%	0.178%
34	9.798	1322	1328	1336	rBV5	48051	139406	1.77%	0.157%
35	9.880	1336	1342	1350	rVB	178220	358452	4.54%	0.403%
36	9.969	1350	1357	1360	rBV	266606	483024	6.12%	0.543%

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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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

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Title : SW846 8260

37	10.016	1360	1365	1372	rVB3	377614	759595	9.63%	0.855%
38	10.204	1390	1397	1402	rBV6	59040	173209	2.20%	0.195%
39	10.322	1411	1417	1421	rBV	131254	224544	2.85%	0.253%
40	10.363	1421	1424	1429	rVB3	61464	102733	1.30%	0.116%
41	10.439	1429	1437	1443	rBV	1951636	3605038	45.69%	4.056%
42	10.504	1443	1448	1455	rVB	437246	782900	9.92%	0.881%
43	10.604	1460	1465	1467	rBV2	47196	83611	1.06%	0.094%
44	10.857	1501	1508	1517	rVB3	104826	242942	3.08%	0.273%
45	10.939	1517	1522	1532	rVB3	45497	98113	1.24%	0.110%
46	11.057	1532	1542	1547	rBV	57829	125209	1.59%	0.141%
47	11.127	1547	1554	1564	rVB	301494	557006	7.06%	0.627%
48	11.739	1650	1658	1668	rBV	1816025	3291377	41.71%	3.703%
49	11.839	1668	1675	1685	rVV	2901980	4916095	62.30%	5.531%
50	11.951	1686	1694	1706	rVB	2243041	4286195	54.32%	4.822%
51	12.274	1738	1749	1758	rVB2	102247	222389	2.82%	0.250%
52	12.574	1792	1800	1815	rVV	4803215	7890821	100.00%	8.878%
53	12.727	1818	1826	1837	rVB	1438533	2387208	30.25%	2.686%
54	12.916	1850	1858	1866	rBV	2440075	3943985	49.98%	4.437%
55	13.010	1866	1874	1880	rVV	342528	556078	7.05%	0.626%
56	13.216	1902	1909	1920	rBV	434490	712907	9.03%	0.802%
57	13.363	1927	1934	1947	rVB	1031394	1679997	21.29%	1.890%
58	13.492	1947	1956	1963	rBV3	195441	461901	5.85%	0.520%
59	13.668	1980	1986	1998	rBV2	1542786	3170250	40.18%	3.567%
60	13.833	2008	2014	2016	rBV	808572	1211249	15.35%	1.363%
61	13.874	2016	2021	2026	rVV	4607978	7730094	97.96%	8.697%
62	13.916	2026	2028	2037	rVV2	344324	626289	7.94%	0.705%
63	13.998	2037	2042	2048	rVV	154028	235293	2.98%	0.265%
64	14.063	2048	2053	2059	rVB3	80068	138169	1.75%	0.155%
65	14.127	2059	2064	2072	rBV2	95612	185010	2.34%	0.208%
66	14.216	2072	2079	2084	rBV	1308263	2058784	26.09%	2.316%
67	14.274	2084	2089	2092	rVV	199283	320351	4.06%	0.360%
68	14.321	2092	2097	2104	rVB2	788757	1324165	16.78%	1.490%
69	14.533	2124	2133	2138	rBV	1039106	1708838	21.66%	1.923%
70	14.574	2138	2140	2144	rVV	112035	141786	1.80%	0.160%
71	14.804	2173	2179	2185	rBV	623919	981980	12.44%	1.105%
72	14.927	2191	2200	2206	rBV2	1392223	2794721	35.42%	3.144%
73	14.986	2206	2210	2218	rVB3	72029	148518	1.88%	0.167%
74	15.080	2220	2226	2233	rBV2	137958	236594	3.00%	0.266%
75	15.192	2235	2245	2249	rVV3	136803	310321	3.93%	0.349%
76	15.233	2249	2252	2256	rVV	50971	78974	1.00%	0.089%

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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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

77	15.304	2256	2264	2277	rVB3	120941	295175	3.74%	0.332%
78	15.504	2290	2298	2310	rVB	1175503	2248051	28.49%	2.529%
79	15.798	2340	2348	2357	rBV	43776	89572	1.14%	0.101%

Sum of corrected areas: 88885364

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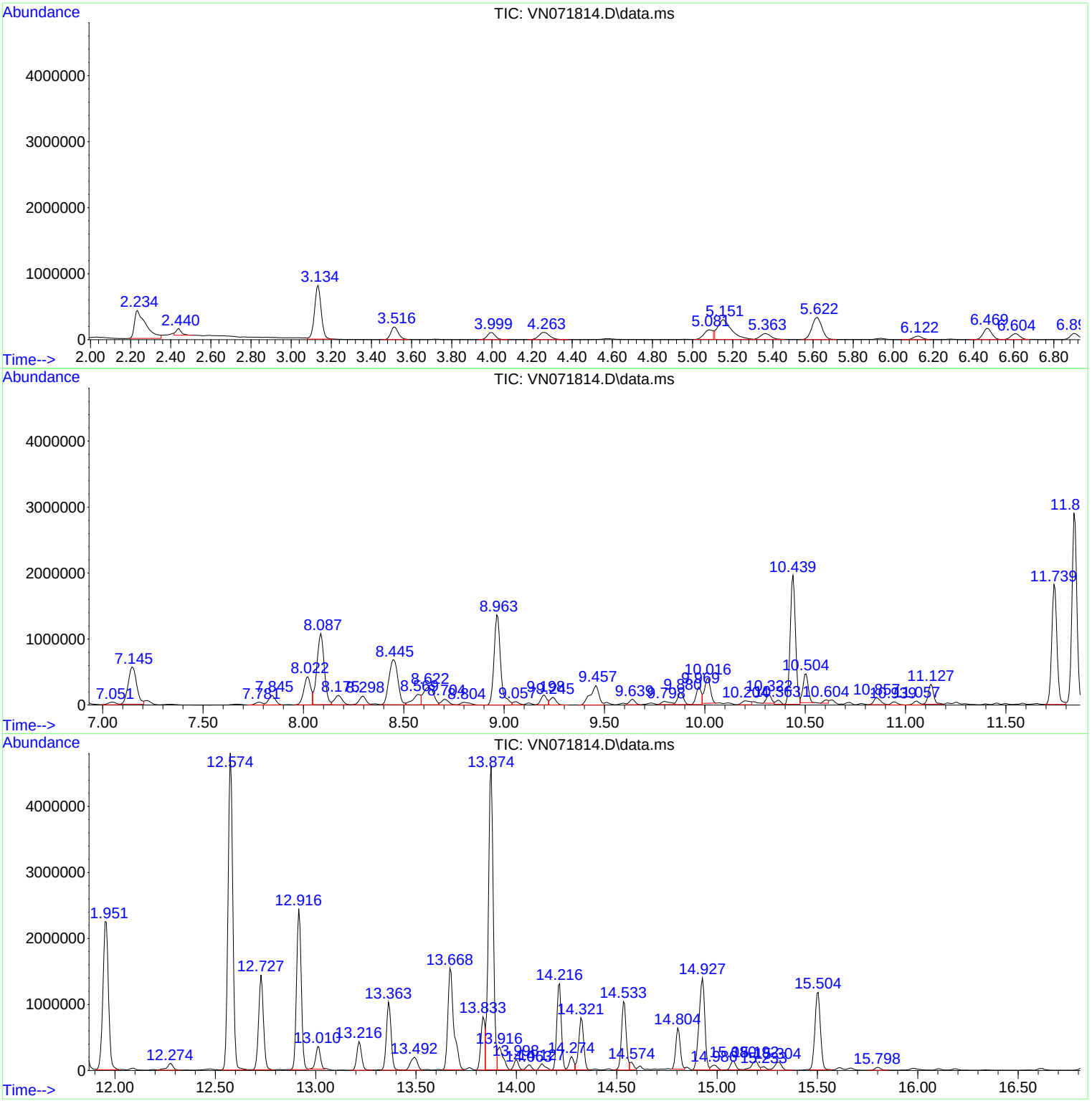
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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\VN040222\
Data File : VN071814.D
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Operator : JC\MD
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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

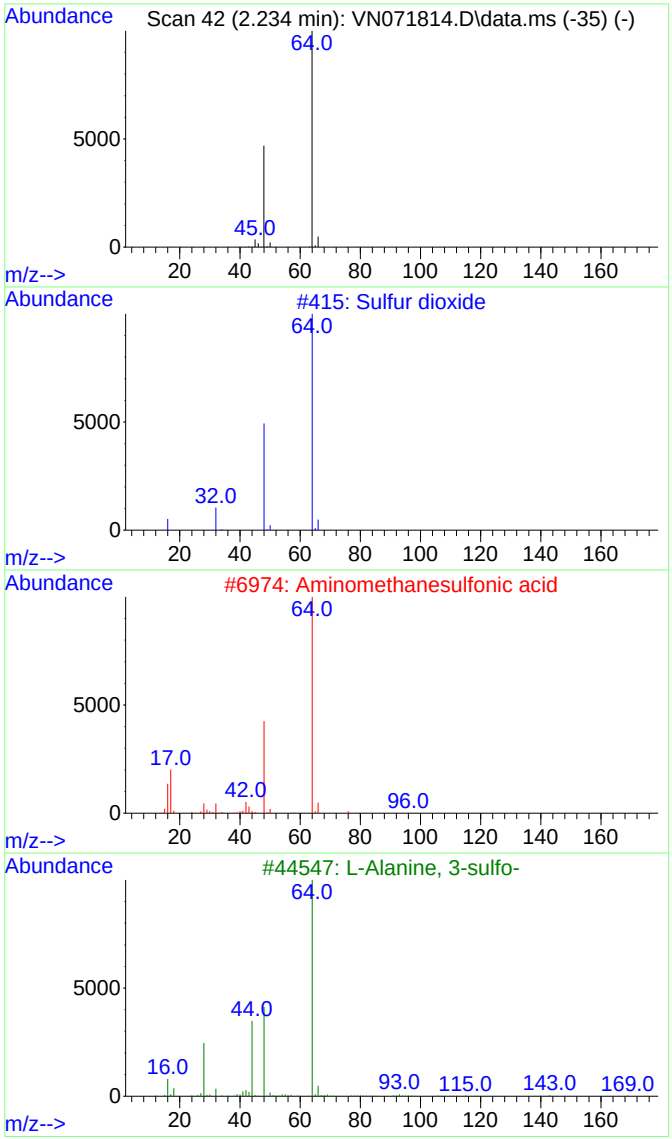
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

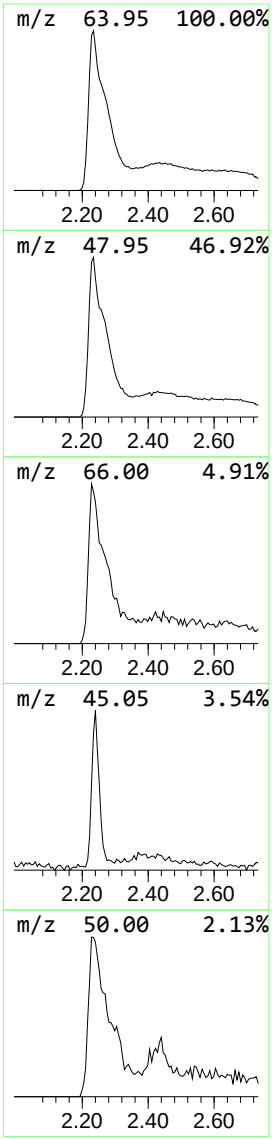
****

Peak Number 1 Sulfur dioxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.234	30.65 ug/l	1669510	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	7
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	3





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Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

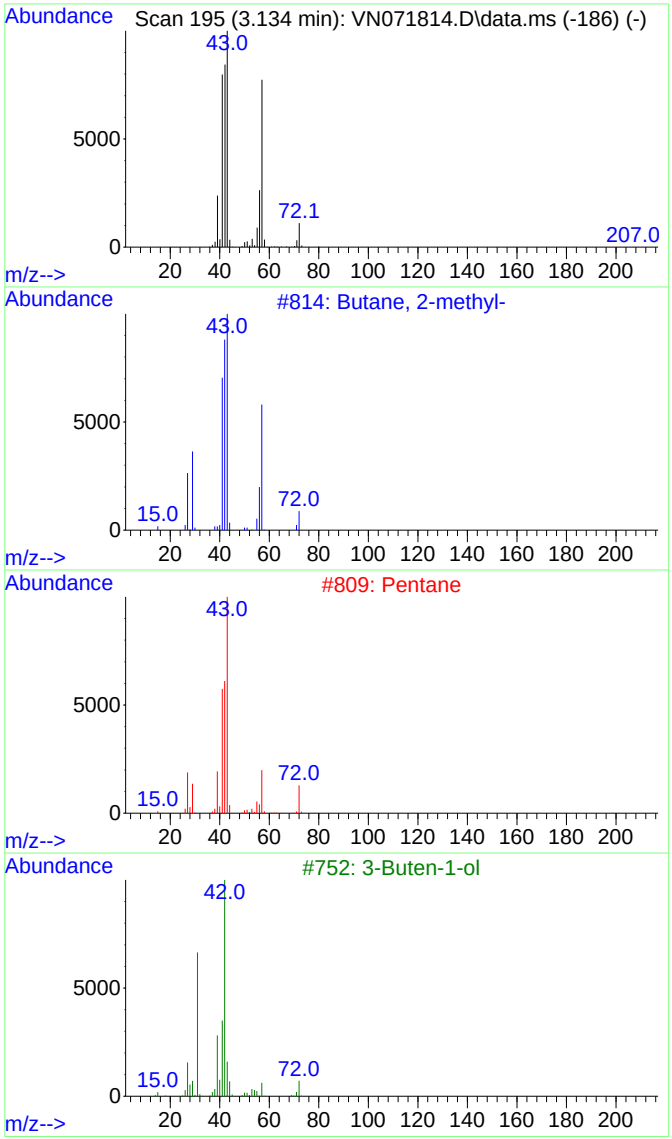
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

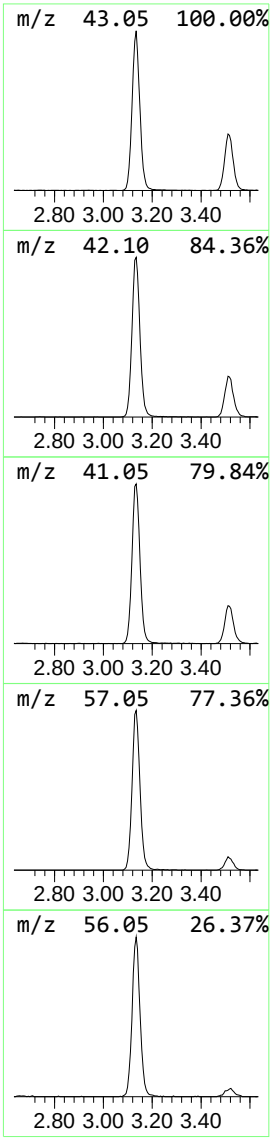
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.134	34.13 ug/l	1858910	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	87
2	Pentane	72	C5H12	000109-66-0	46
3	3-Buten-1-ol	72	C4H8O	000627-27-0	43
4	1-Butanesulfonyl chloride	156	C4H9ClO2S	002386-60-9	12
5	1-Butene	56	C4H8	000106-98-9	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
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Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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

TIC Library : C:\Database\NIST20.L
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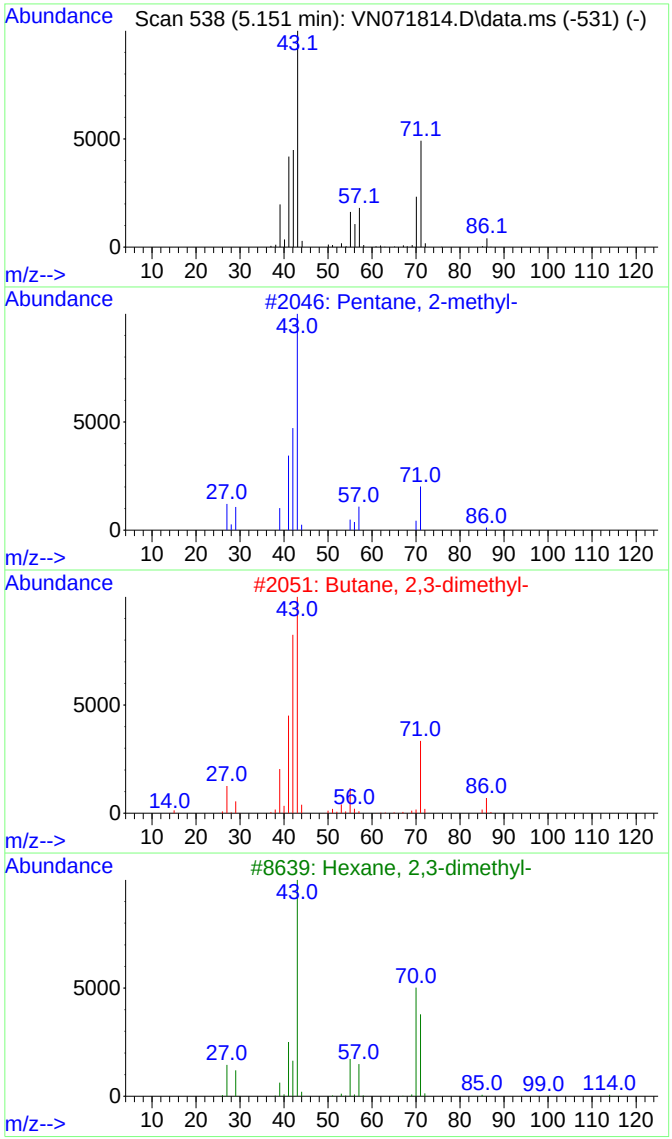
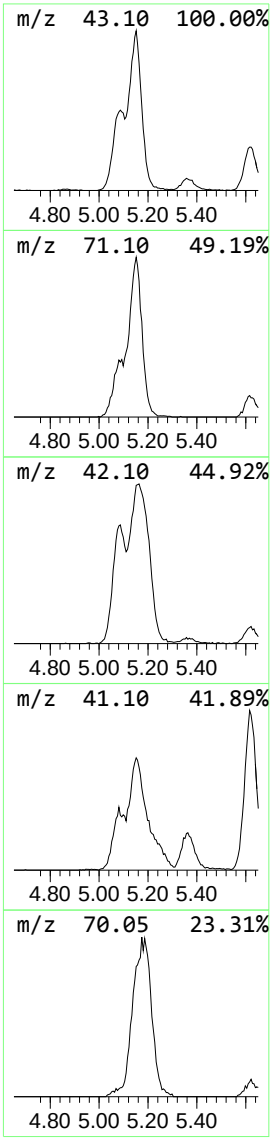
****
Peak Number 3 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.151	25.81 ug/l	1406050	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	78
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	37
3	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	25
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	25
5	Propanoic acid, 2-methyl-, 2-pro...	128	C7H12O2	015727-77-2	23





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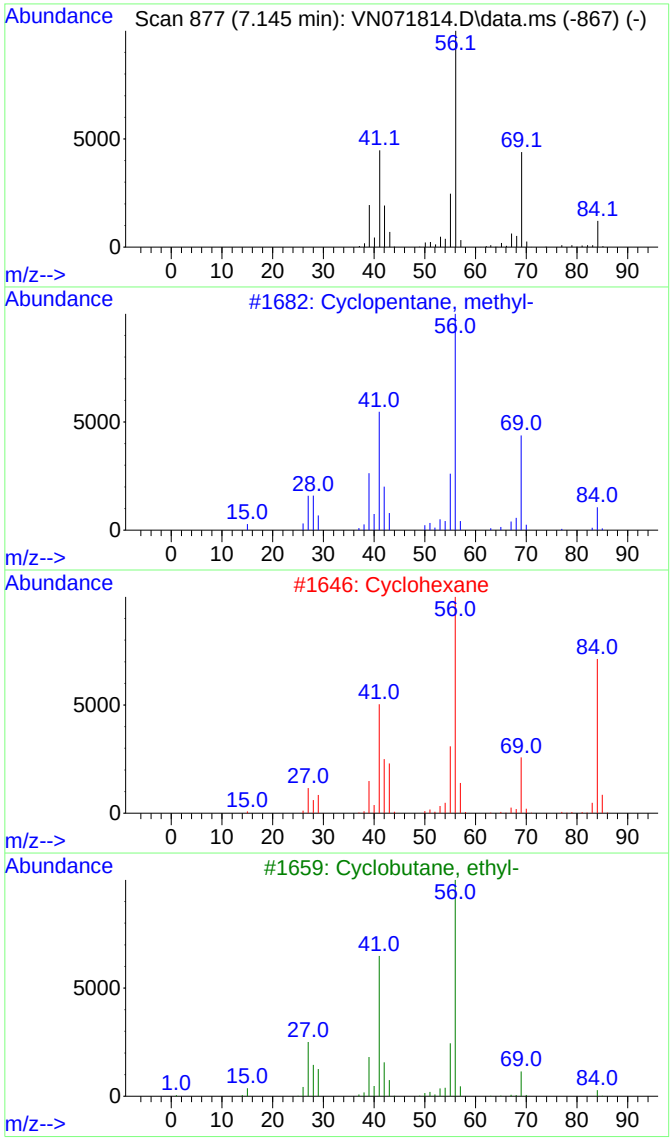
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

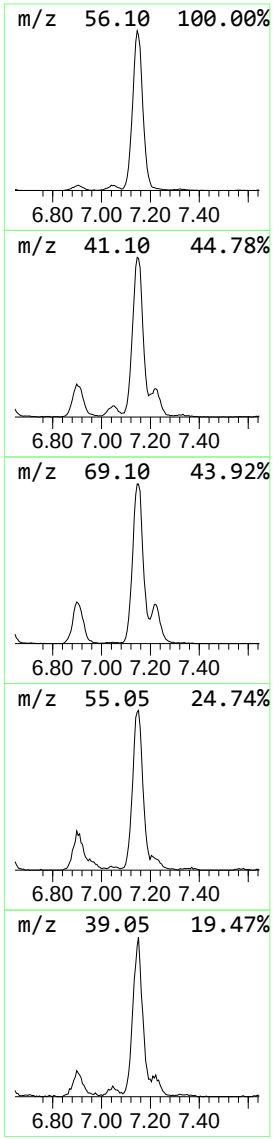
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.145	30.39 ug/l	1655360	Pentafluorobenzene	8.086

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	83
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	47





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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

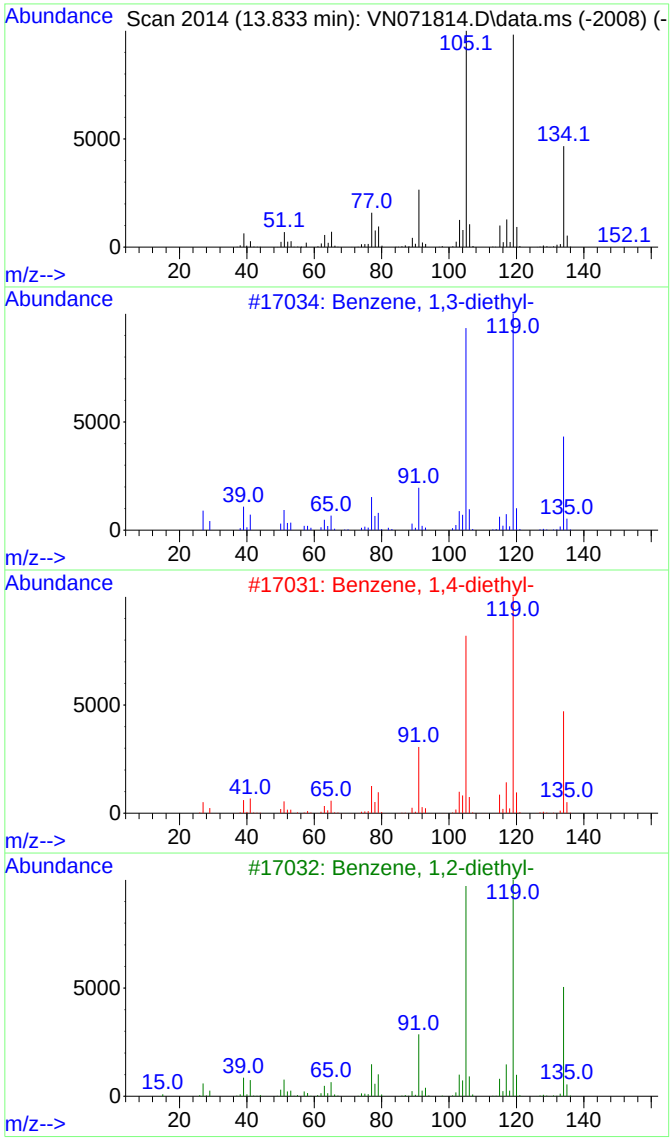
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

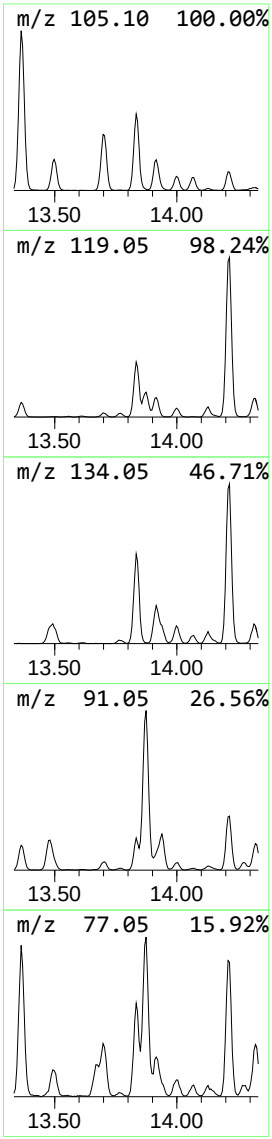
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,3-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.833	19.10 ug/l	1211250	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

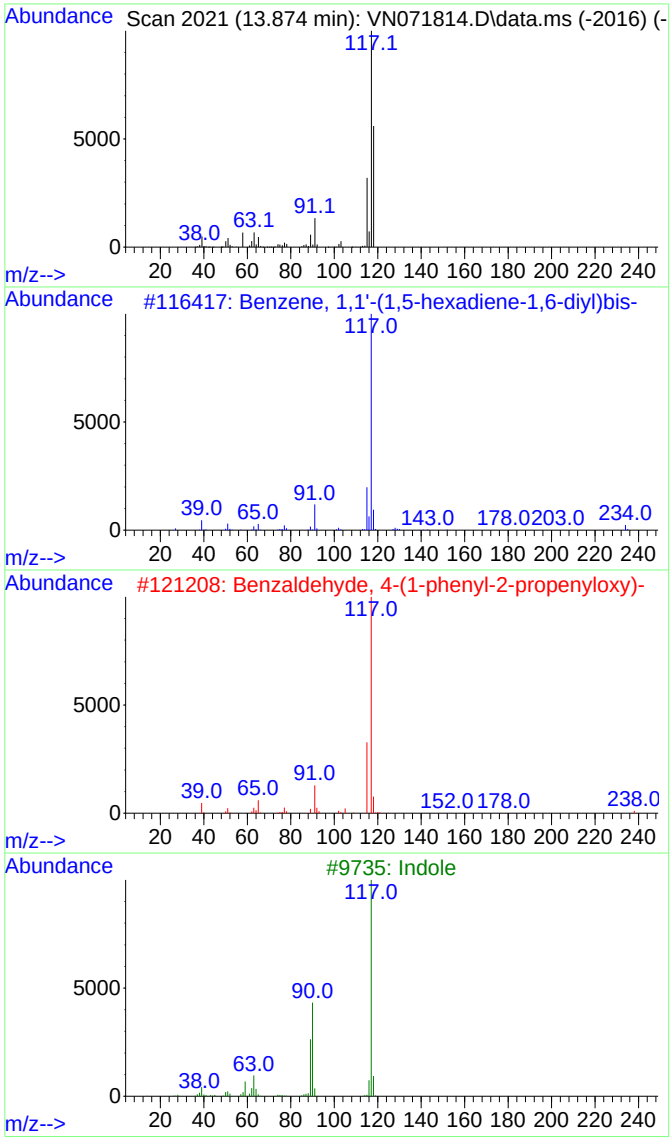
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

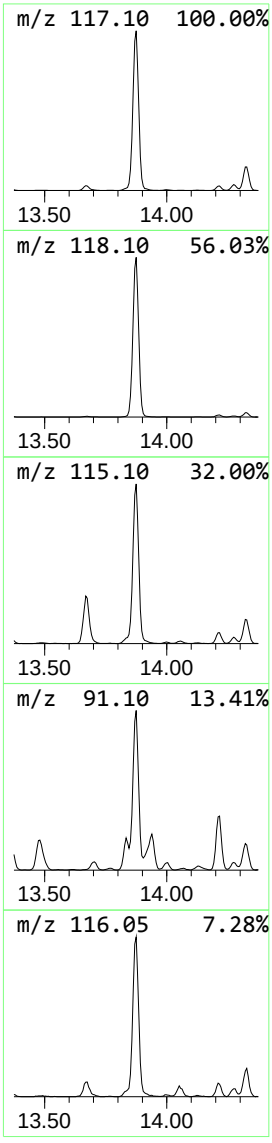
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	121.92 ug/l	7730090	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
2	Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	53
3	Indole	117	C8H7N	000120-72-9	46
4	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

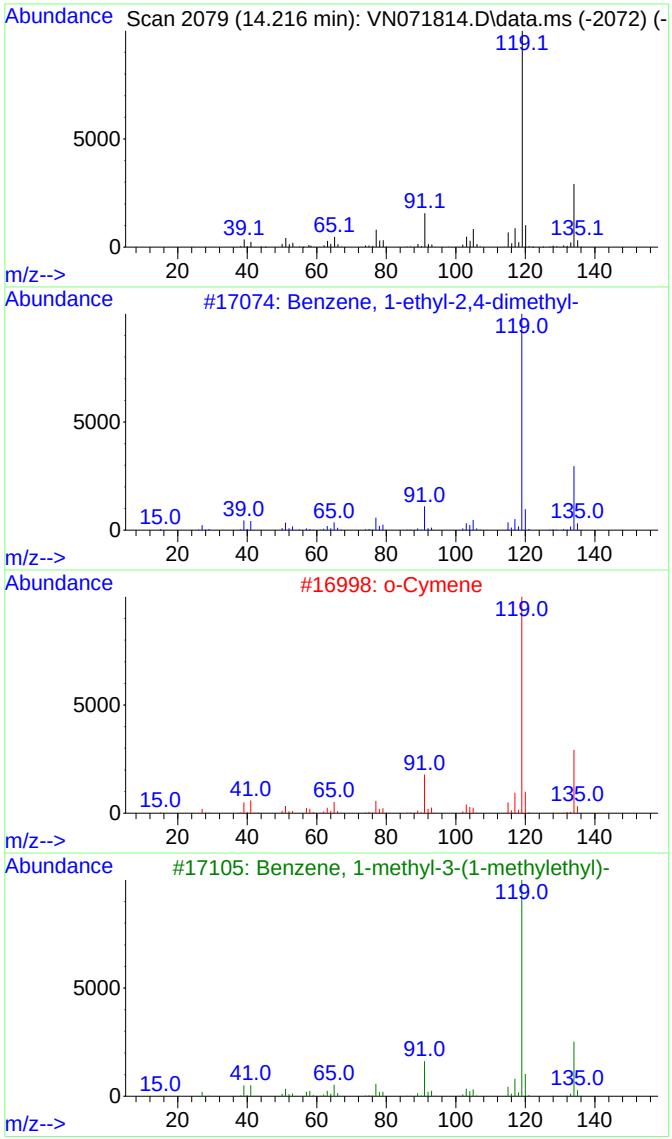
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

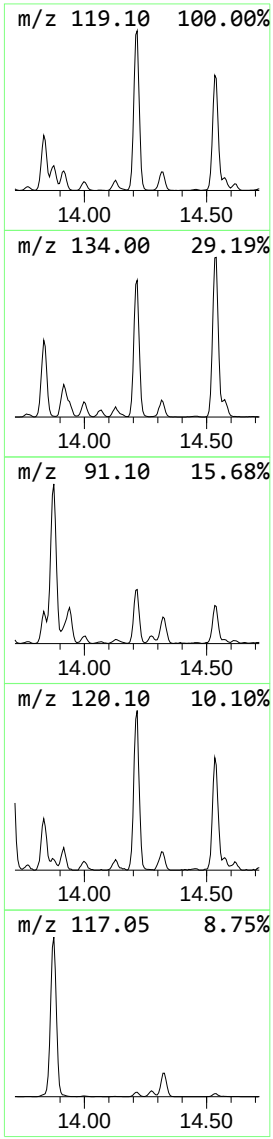
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.216	32.47 ug/l	2058780	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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

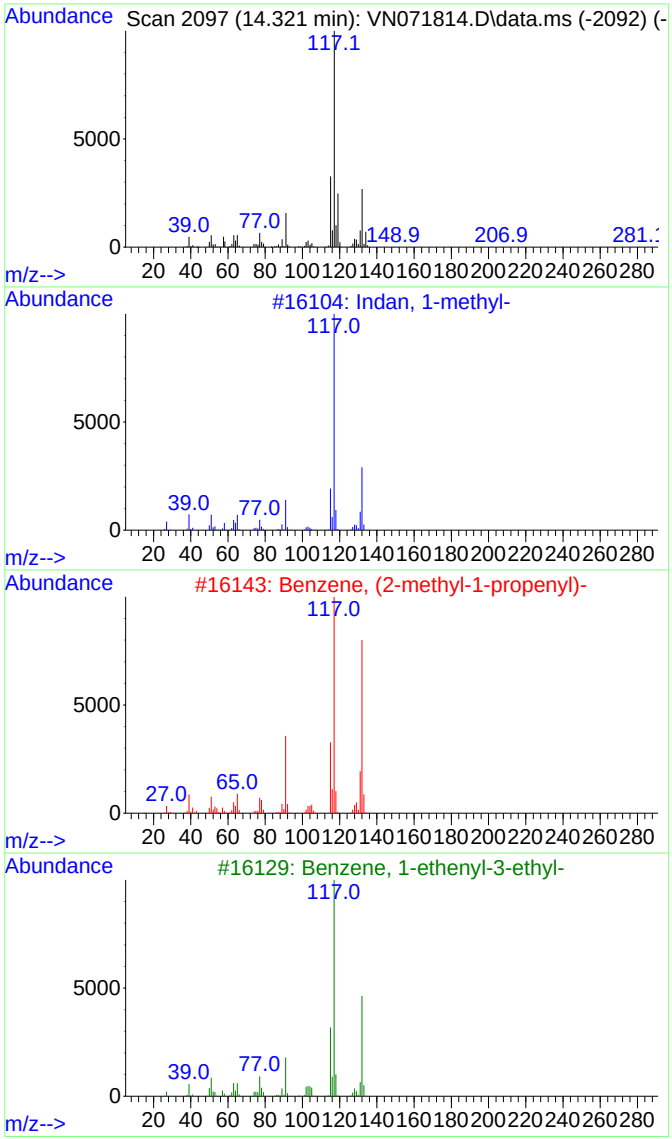
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

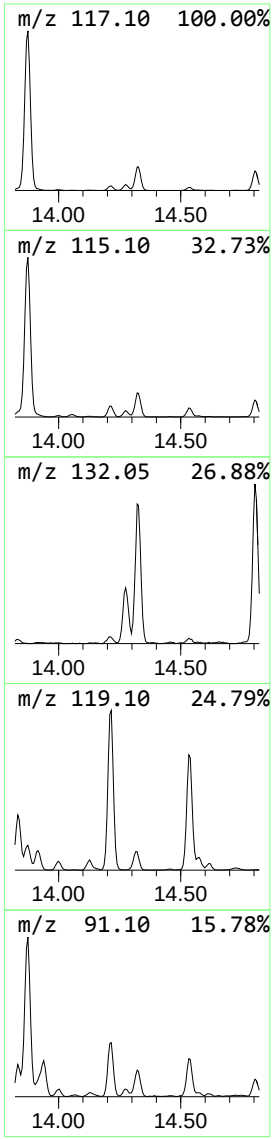
****

Peak Number 8 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.321	20.88 ug/l	1324170	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	76
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	72
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

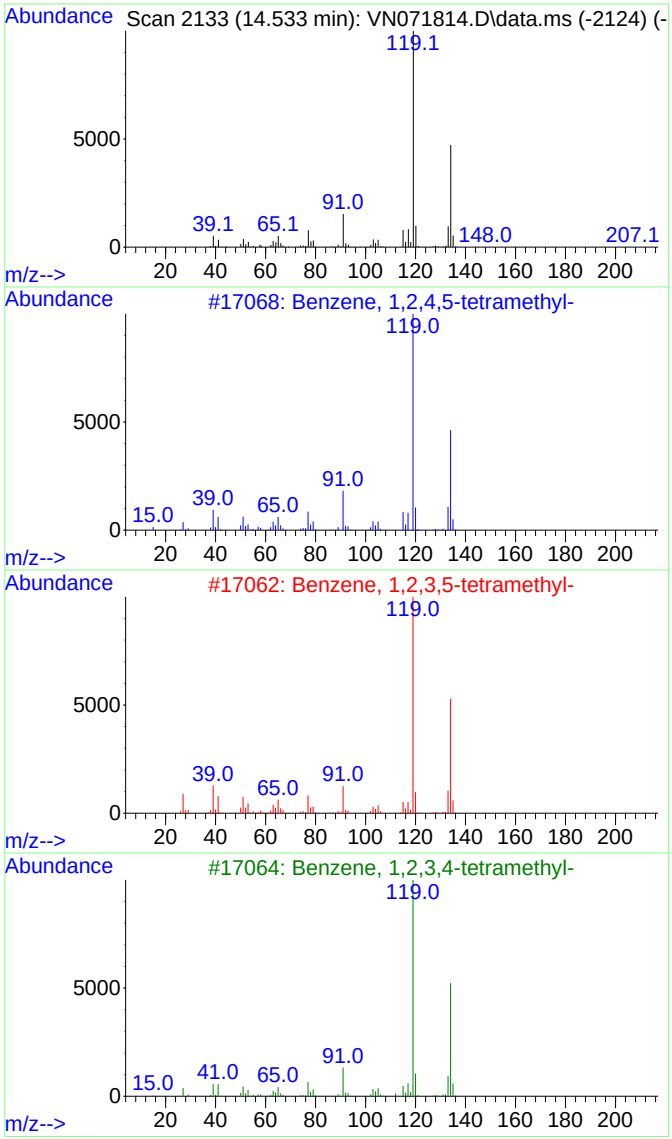
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

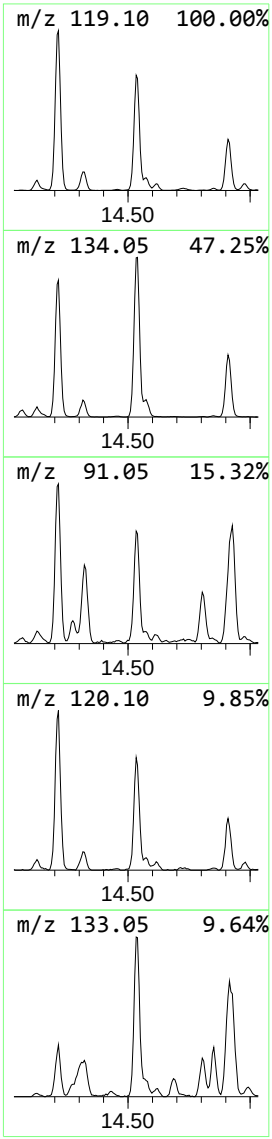
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.533	26.95 ug/l	1708840	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
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\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

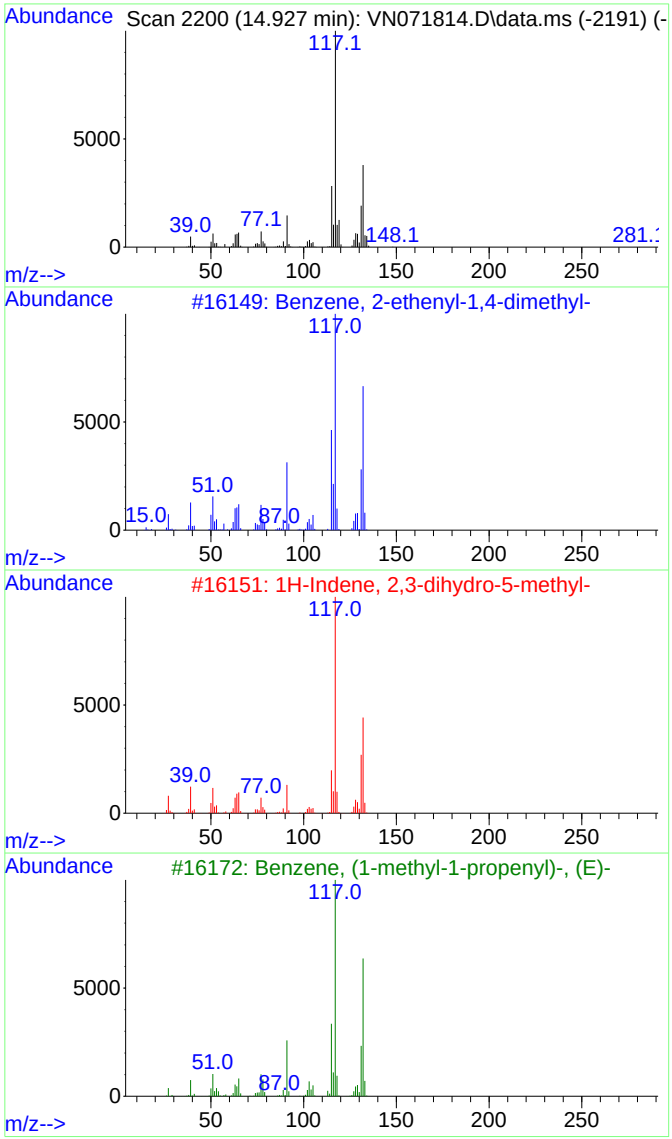
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

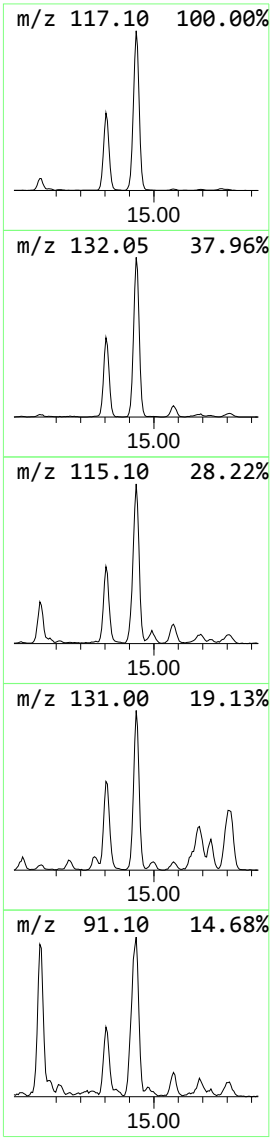
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	44.08 ug/l	2794720	1,4-Dichlorobenzene-d4	13.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040222\
Data File : VN071814.D
Acq On : 02 Apr 2022 21:47
Operator : JC\MD
Sample : N2244-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3-BD

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
Sulfur dioxide	2.234	30.6	ug/l	1669510	1	8.086	2723510	50.0
Butane, 2-methyl-	3.134	34.1	ug/l	1858910	1	8.086	2723510	50.0
Pentane, 2-methyl-	5.151	25.8	ug/l	1406050	1	8.086	2723510	50.0
Cyclopentane, m...	7.145	30.4	ug/l	1655360	1	8.086	2723510	50.0
Benzene, 1,3-di...	13.833	19.1	ug/l	1211250	4	13.669	3170250	50.0
Benzene, 1,1'-(...	13.874	121.9	ug/l	7730090	4	13.669	3170250	50.0
Benzene, 1-ethy...	14.216	32.5	ug/l	2058780	4	13.669	3170250	50.0
Indan, 1-methyl-	14.321	20.9	ug/l	1324170	4	13.669	3170250	50.0
Benzene, 1,2,4,...	14.533	26.9	ug/l	1708840	4	13.669	3170250	50.0
Benzene, 2-ethe...	14.927	44.1	ug/l	2794720	4	13.669	3170250	50.0