LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

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\82N042423W.M
Title : SW846 8260

Signal : TIC: VN077479.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.519	92	99	110	rBV	1024139	1708836	3.99%	0.458%
2	3.254	213	224	241	rBV	19288999	42802658	100.00%	11.477%
3	3.654	280	292	310	rBV	2854243	7129622	16.66%	1.912%
4	4.154	367	377	388	rVB3	172834	487006	1.14%	0.131%
5	4.437	409	425	442	rBV	1862324	6585627	15.39%	1.766%
6	5.295	555	571	572	rVV2	3082096	8594405	20.08%	2.304%
7	5.354	574	581	617	rVB	7730918	29410659	68.71%	7.886%
8	5.825	644	661	685	rBV	6118152	21218123	49.57%	5.689%
9	6.137	701	714	732	rVB3	126993	444849	1.04%	0.119%
10	6.325	732	746	761	rBV2	2394006	7342793	17.15%	1.969%
11	6.666	783	804	816	rBV2	421977	1332206	3.11%	0.357%
12	6.801	816	827	838	rBV2	297526	932407	2.18%	0.250%
13	7.089	865	876	889	rBV3	581105	1847828	4.32%	0.495%
14	7.231	890	900	909	rVV	1746315	5166372	12.07%	1.385%
15	7.337	909	918	939	rVV	4132181	12358223	28.87%	3.314%
16	7.501	939	946	960	rVB3	149883	474654	1.11%	0.127%
17	7.936	1012	1020	1026	rBV2	228151	611289	1.43%	0.164%
18	8.013	1026	1033	1041	rVB	692555	1747603	4.08%	0.469%
19	8.172	1051	1060	1063	rBV3	566291	1308911	3.06%	0.351%
20	8.231	1063	1070	1081	rVV3	3663286	12249156	28.62%	3.284%
21	8.336	1081	1088	1099	rVB	4960746	12499767	29.20%	3.352%
22	8.454	1099	1108	1115	rBV	3651897	8407129	19.64%	2.254%
23	8.519	1115	1119	1125	rVB	440382	858825	2.01%	0.230%
24	8.583	1125	1130	1140	rVB	406372	859454	2.01%	0.230%
25	8.725	1144	1154	1158	rBV2	2038730	5467071	12.77%	1.466%
26	8.783	1158	1164	1172	rVV4	2136015	6796171	15.88%	1.822%
27	8.860	1172	1177	1186	rVB	1535974	3390701	7.92%	0.909%
28	8.972	1186	1196	1206	rVB2	1285245	2997912	7.00%	0.804%
29	9.107	1209	1219	1226	rBV3	2666985	6702040	15.66%	1.797%
30	9.266	1240	1246	1252	rVB	376990	722060	1.69%	0.194%
31	9.336	1252	1258	1263	rBV	376903	798048	1.86%	0.214%
32	9.389	1263	1267	1287	rVB2	473110	1066968	2.49%	0.286%
33	9.601	1287	1303	1319	rBV3	4703459	14502487	33.88%	3.889%
34	9.778	1328	1333	1340	rVB	992941	1883066	4.40%	0.505%
35	9.872	1340	1349	1355	rVB3	473714	1099795	2.57%	0.295%
36	10.013	1367	1373	1382	rVB2	1618881	3479868	8.13%	0.933%

LSC Area Percent Report

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Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

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\82N042423W.M
Title : SW846 8260

37	10.107	1382	1389	1392	rBV2	1081841	2148418	5.02%	0.576%
38	10.148	1392	1396	1402	rVV	2122900	3866214	9.03%	1.037%
39	10.342	1420	1429	1441	rBV3	929526	2301110	5.38%	0.617%
40	10.454	1441	1448	1451	rBV	452077	851786	1.99%	0.228%
41	10.495	1451	1455	1461	rVB2	580241	1060682	2.48%	0.284%
42	10.572	1461	1468	1474	rBV	2927340	5839378	13.64%	1.566%
43	10.760	1491	1500	1505	rVB5	860885	2066925	4.83%	0.554%
44	10.919	1520	1527	1532	rBV3	509991	1055594	2.47%	0.283%
45	10.983	1532	1538	1549	rVB3	1661508	4235296	9.89%	1.136%
46	11.083	1549	1555	1561	rVB4	176411	432184	1.01%	0.116%
47	11.166	1561	1569	1577	rBV6	139698	435827	1.02%	0.117%
48	11.260	1577	1585	1588	rBV4	370038	828750	1.94%	0.222%
49	11.377	1601	1605	1609	rVV3	403748	759196	1.77%	0.204%
50	11.424	1609	1613	1618	rVB3	335650	564860	1.32%	0.151%
51	11.866	1681	1688	1695	rBV	2032957	3653534	8.54%	0.980%
52	11.966	1700	1705	1715	rVB	1308524	2341573	5.47%	0.628%
53	12.077	1715	1724	1727	rBV2	253720	639876	1.49%	0.172%
54	12.695	1820	1829	1836	rBV	2153973	3734191	8.72%	1.001%
55	12.848	1847	1855	1865	rVB	1449617	2749838	6.42%	0.737%
56	13.036	1877	1887	1894	rVV	6951614	11336363	26.49%	3.040%
57	13.130	1894	1903	1908	rVV	664813	1284881	3.00%	0.345%
58	13.336	1931	1938	1946	rVB	750416	1280196	2.99%	0.343%
59	13.607	1976	1984	1991	rBV3	1264260	3083951	7.21%	0.827%
60	13.795	2009	2016	2027	rVB2	1754130	3264392	7.63%	0.875%
61	13.954	2037	2043	2047	rBV	3277323	5190015	12.13%	1.392%
62	13.995	2047	2050	2054	rBV	1145809	1467761	3.43%	0.394%
63	14.036	2054	2057	2059	rBV	919863	1326685	3.10%	0.356%
64	14.118	2066	2071	2077	rVB	667988	1076848	2.52%	0.289%
65	14.189	2077	2083	2088	rBV	861779	1400915	3.27%	0.376%
66	14.248	2088	2093	2100	rVB	274605	474139	1.11%	0.127%
67	14.330	2100	2107	2114	rBV	7208215	11557521	27.00%	3.099%
68	14.401	2114	2119	2122	rBV	943476	1638075	3.83%	0.439%
69	14.448	2122	2127	2134	rVB2	3920102	6766057	15.81%	1.814%
70	14.654	2154	2162	2172	rBV	4443380	7797655	18.22%	2.091%
71	14.736	2172	2176	2180	rBV	606374	826238	1.93%	0.222%
72	14.883	2196	2201	2204	rBV	376515	495718	1.16%	0.133%
73	14.930	2204	2209	2214	rBV	2824542	4753729	11.11%	1.275%
74	15.048	2221	2229	2234	rBV3	4008757	9106640	21.28%	2.442%
75	15.095	2234	2237	2250	rVV2	662482	1422188	3.32%	0.381%
76	15.201	2250	2255	2261	rVV2	421994	775707	1.81%	0.208%

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Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

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\82N042423W.M
Title : SW846 8260

77	15.301	2261	2272	2276	rVV2	2291501	5650441	13.20%	1.515%
78	15.342	2276	2279	2283	rVV	985353	1600508	3.74%	0.429%
79	15.412	2283	2291	2301	rVB3	1924145	4618537	10.79%	1.238%
80	15.583	2315	2320	2326	rVB	998522	1702849	3.98%	0.457%
81	15.677	2332	2336	2340	rBV2	496835	787940	1.84%	0.211%
82	15.724	2340	2344	2357	rVB	533460	1115588	2.61%	0.299%
83	15.842	2357	2364	2370	rBV	896722	1527704	3.57%	0.410%
84	15.983	2380	2388	2401	rVV4	388841	1296439	3.03%	0.348%
85	16.154	2410	2417	2430	rVB3	326289	862712	2.02%	0.231%
86	16.495	2469	2475	2486	rVB	836897	1636113	3.82%	0.439%
87	16.659	2494	2503	2513	rVB	480126	972375	2.27%	0.261%

Sum of corrected areas: 372946701

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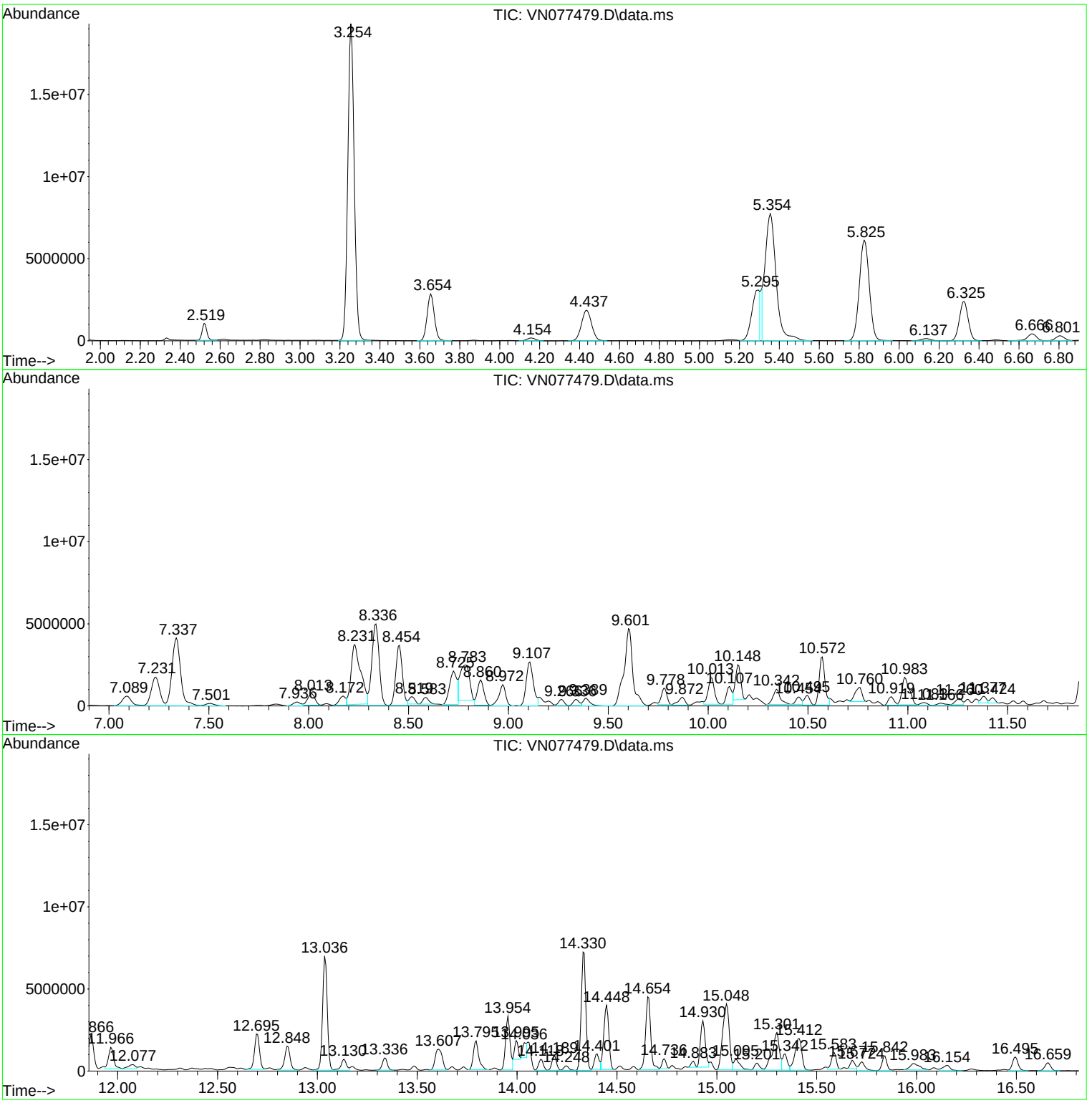
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
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MSVOA_N
ClientSampleId :
MW-1

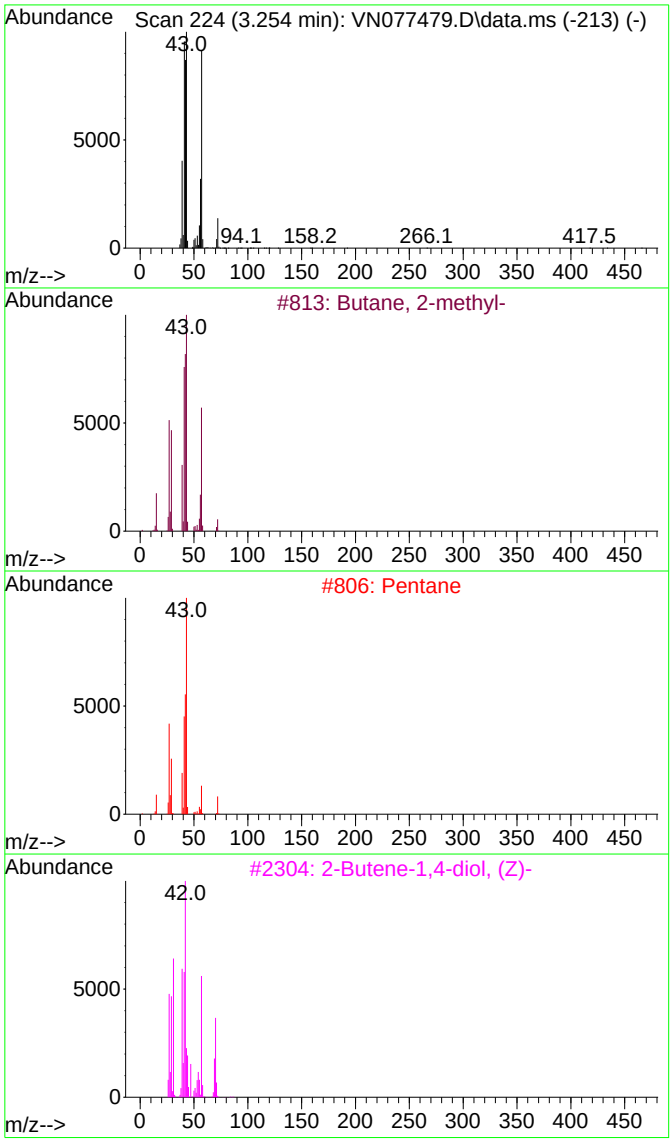
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

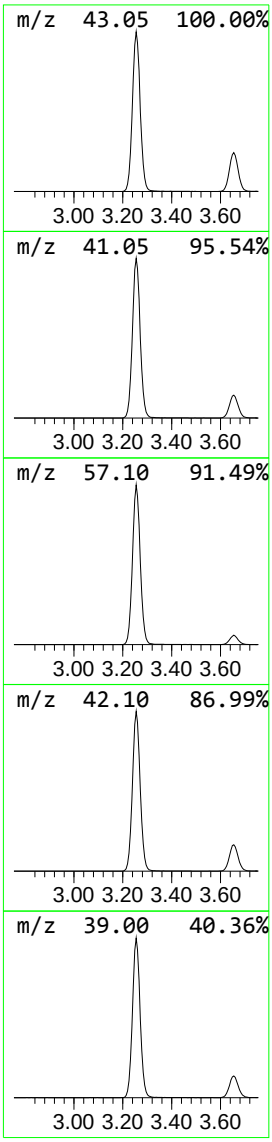
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.254	174.72 ug/l	42802700	Pentafluorobenzene	8.231

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	64
2	Pentane	72	C5H12	000109-66-0	36
3	2-Butene-1,4-diol, (Z)-	88	C4H8O2	006117-80-2	28
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	27
5	Oxirane, trimethyl-	86	C5H10O	005076-19-7	25





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Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

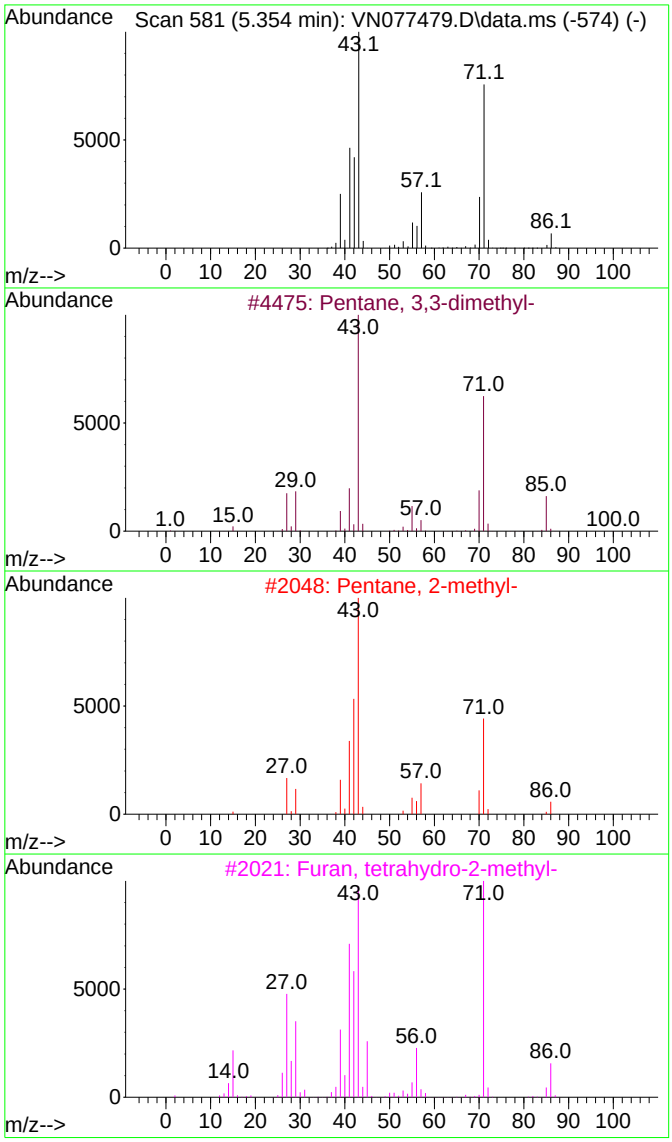
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

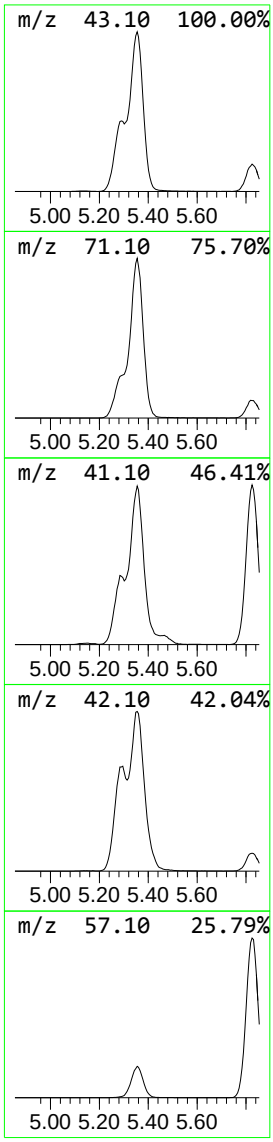
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.354	120.05 ug/l	29410700	Pentafluorobenzene	8.231

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	53
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	49
3	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	45
4	3-Penten-2-ol	86	C5H10O	001569-50-2	43
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	28





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Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

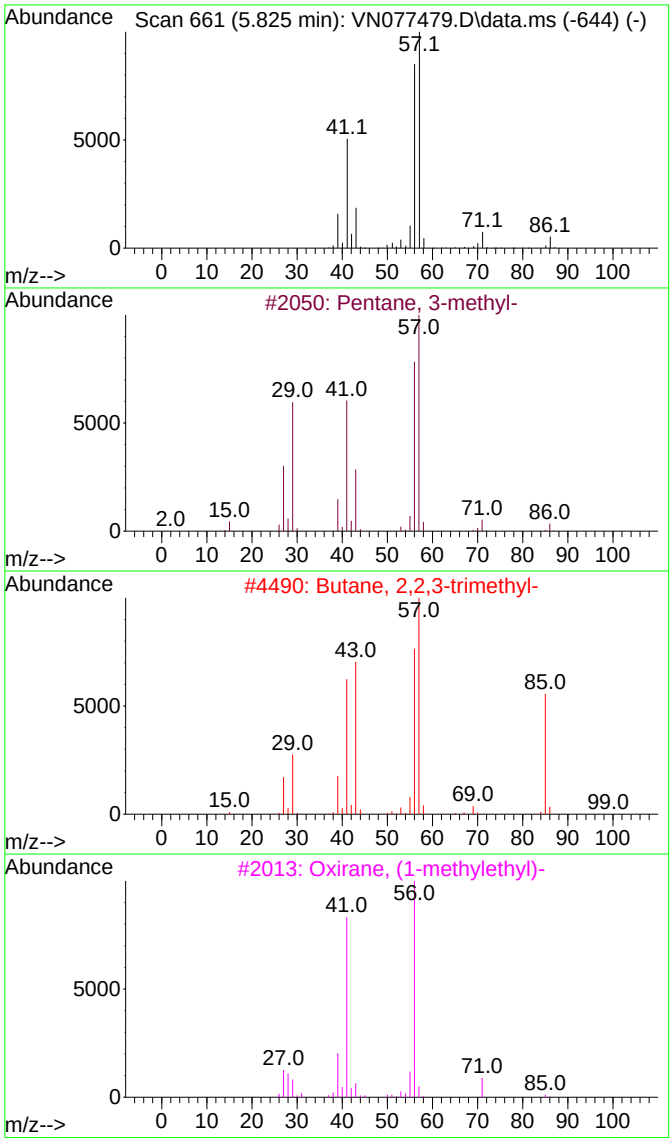
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

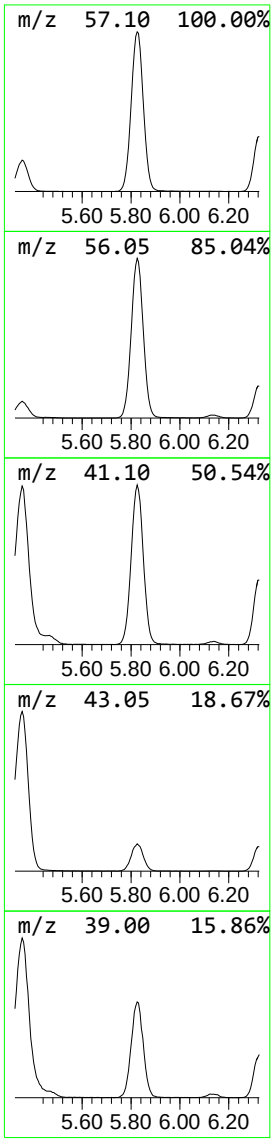
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.825	86.61 ug/l	21218100	Pentafluorobenzene	8.231

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	72
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
4	Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	43
5	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

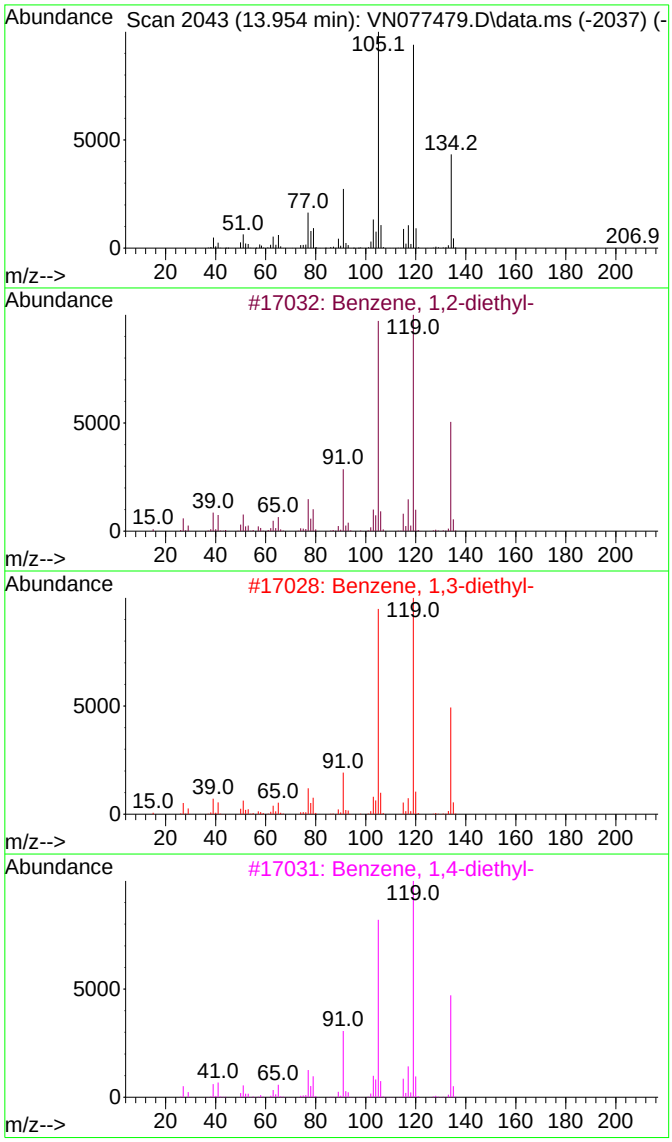
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

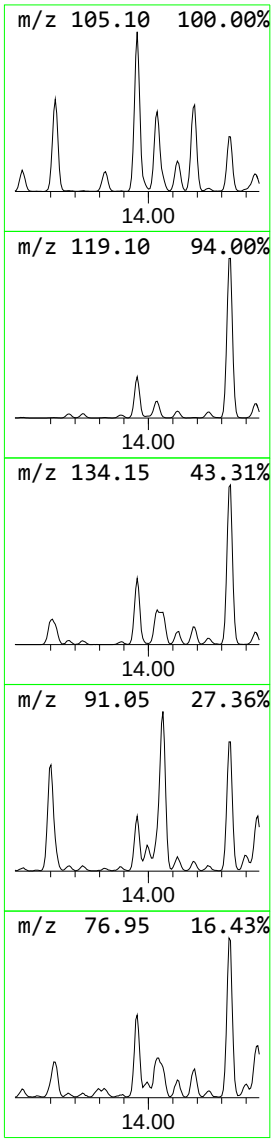
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.954	79.49 ug/l	5190020	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5	o-Cymene	134	C10H14	000527-84-4	81





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Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

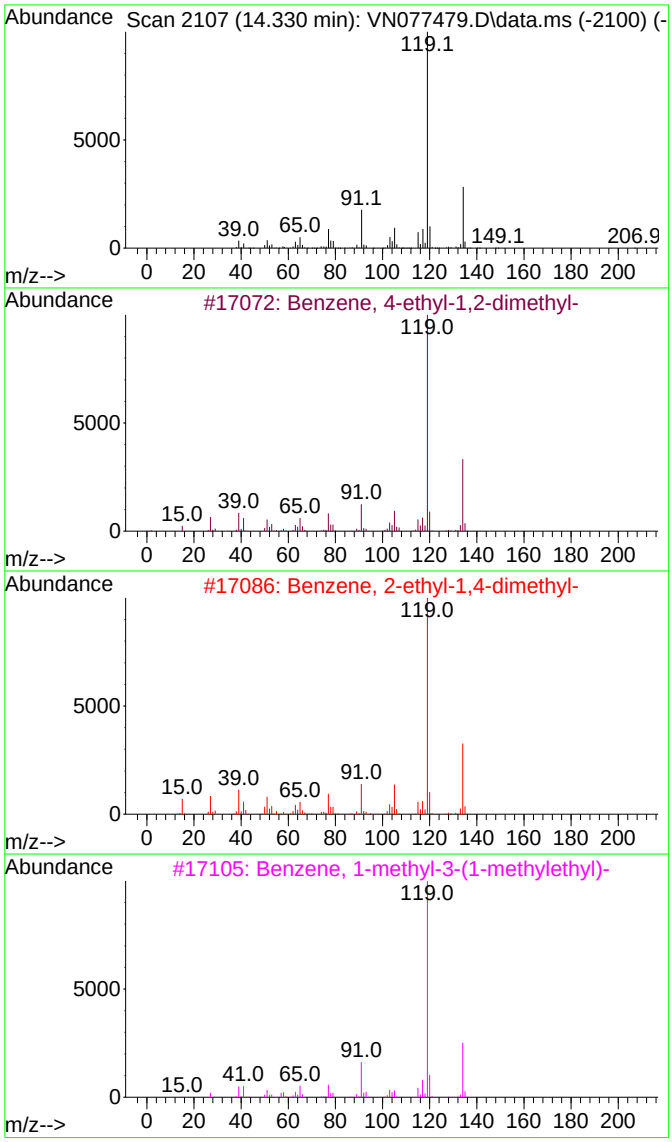
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

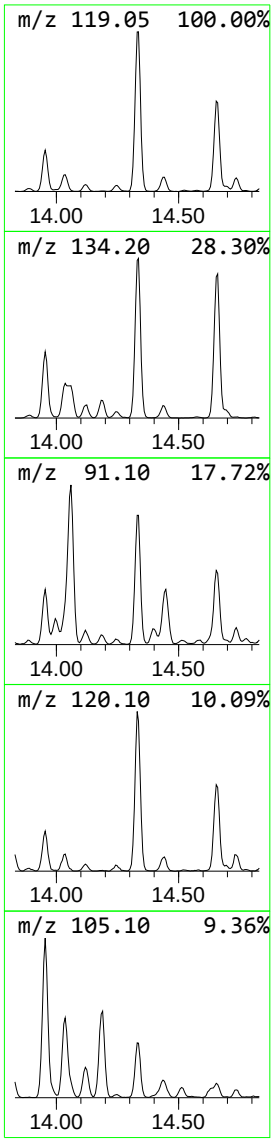
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.330	177.02 ug/l	11557500	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

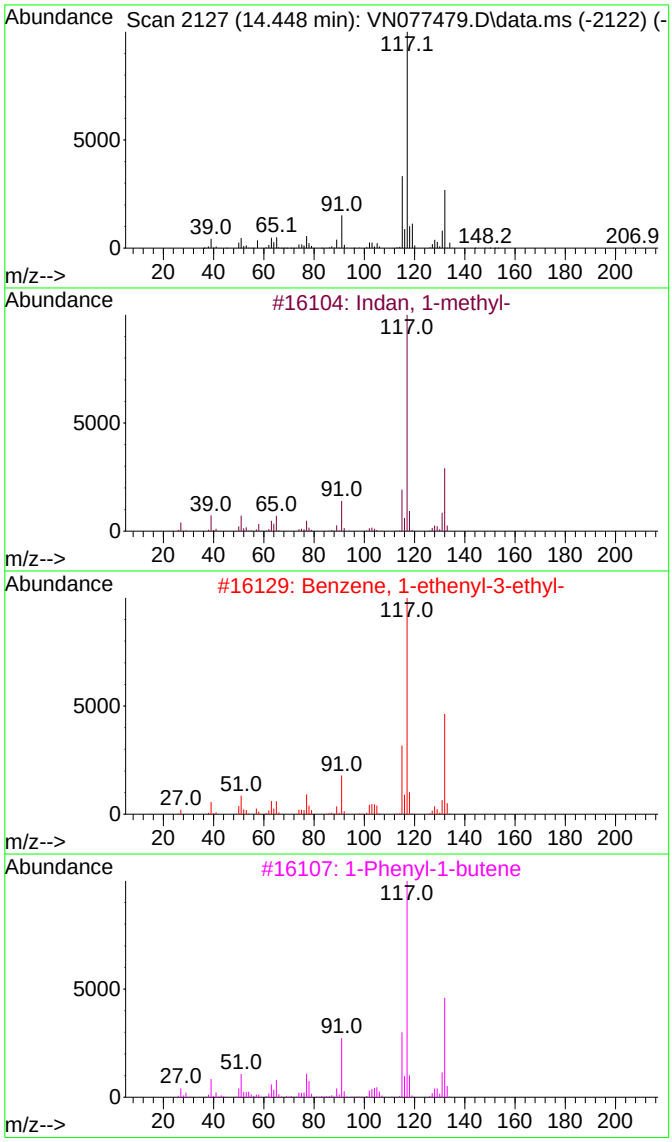
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

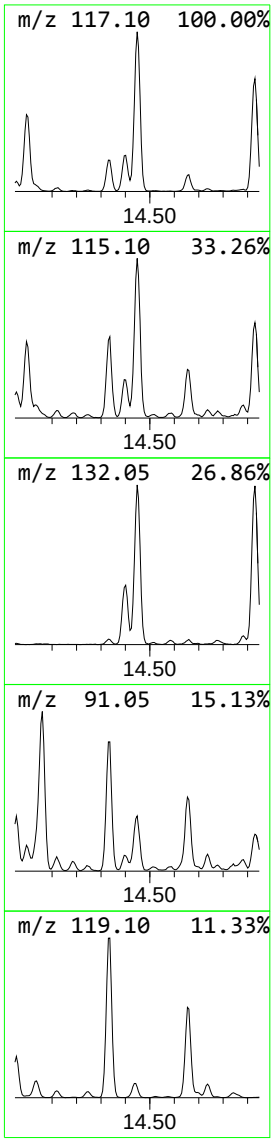
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.448	103.63 ug/l	6766060	1,4-Dichlorobenzene-d4	13.795

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	87
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	83
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

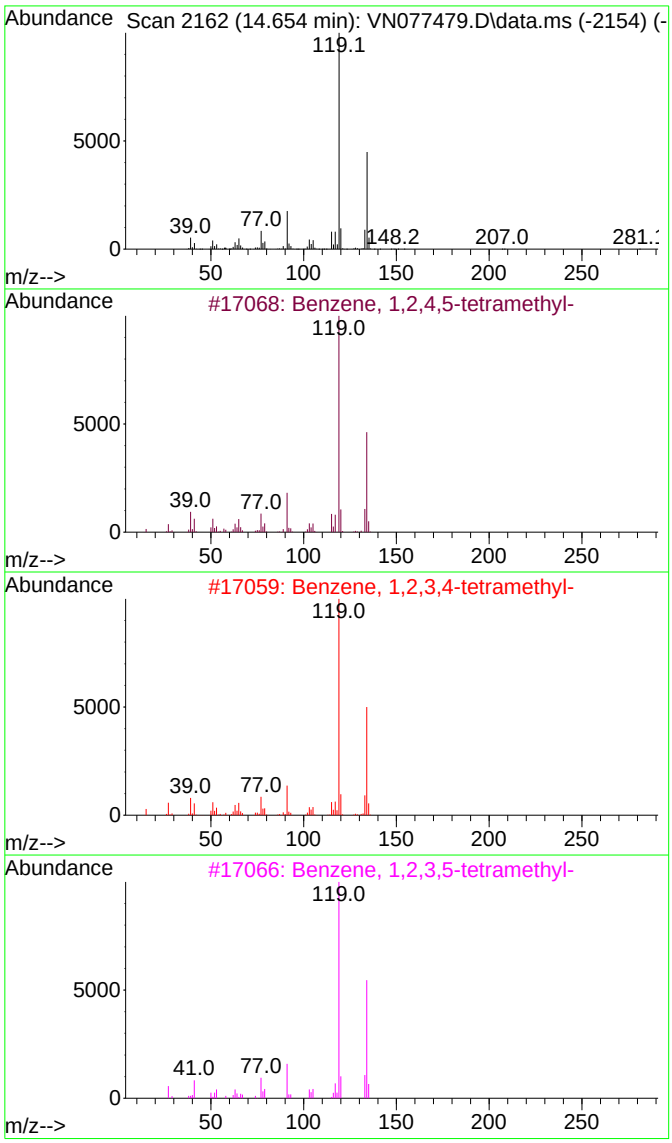
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

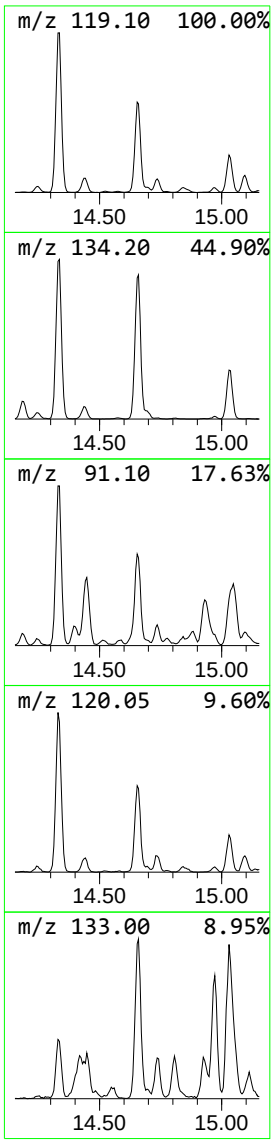
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.654	119.44 ug/l	7797660	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

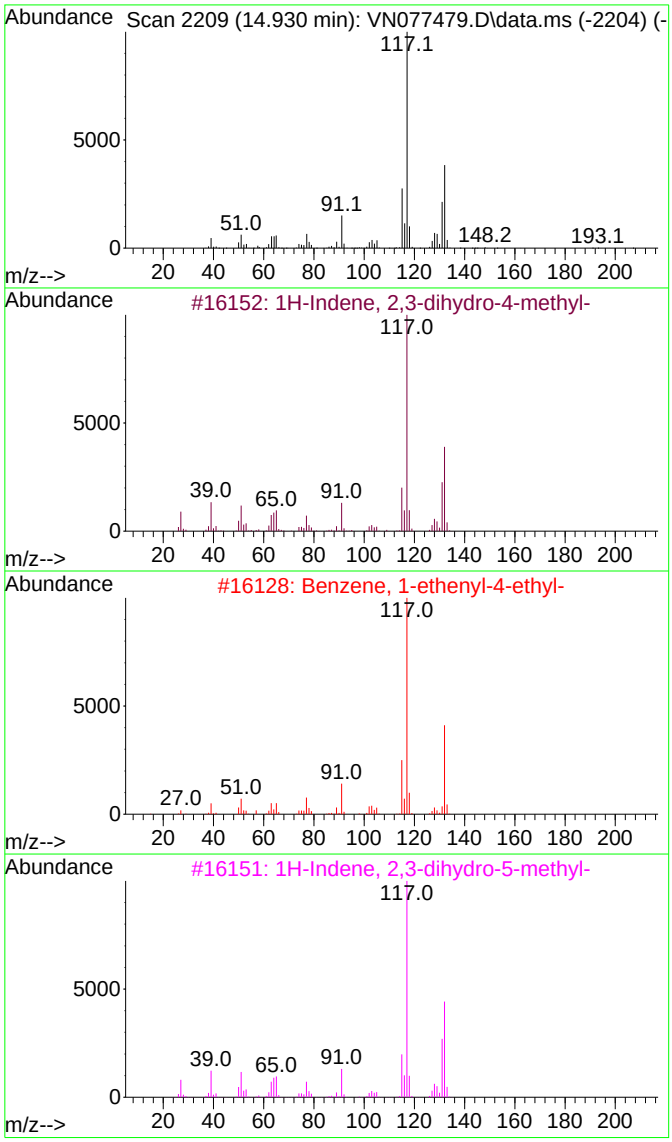
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

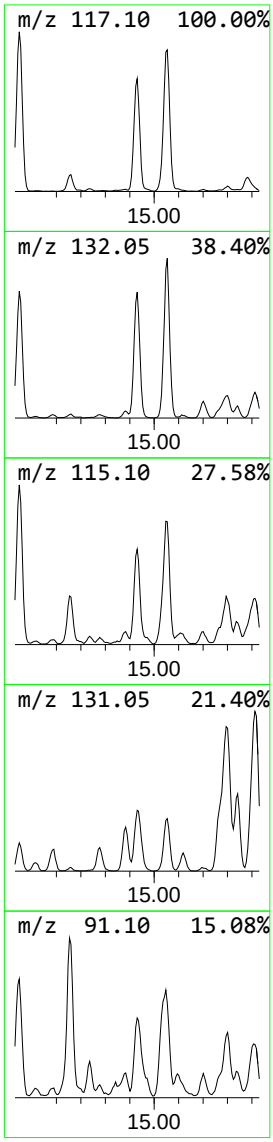
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.930	72.81 ug/l	4753730	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

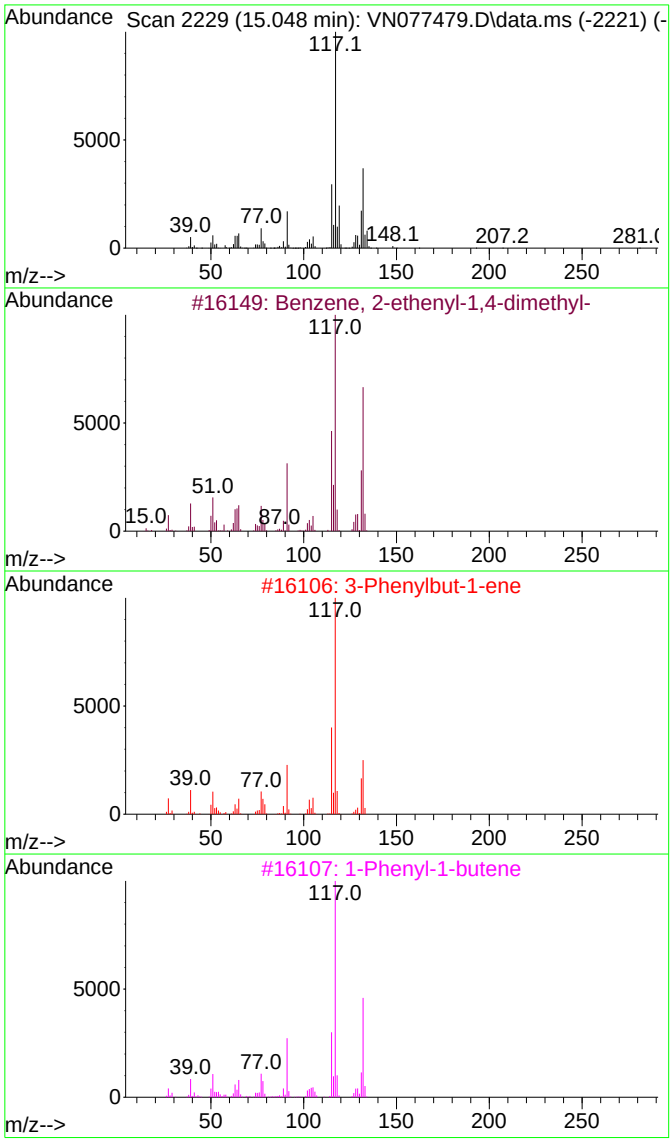
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

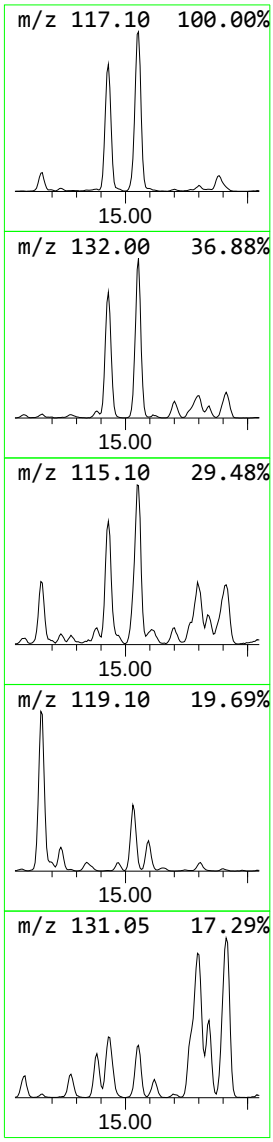
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.048	139.48 ug/l	9106640	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

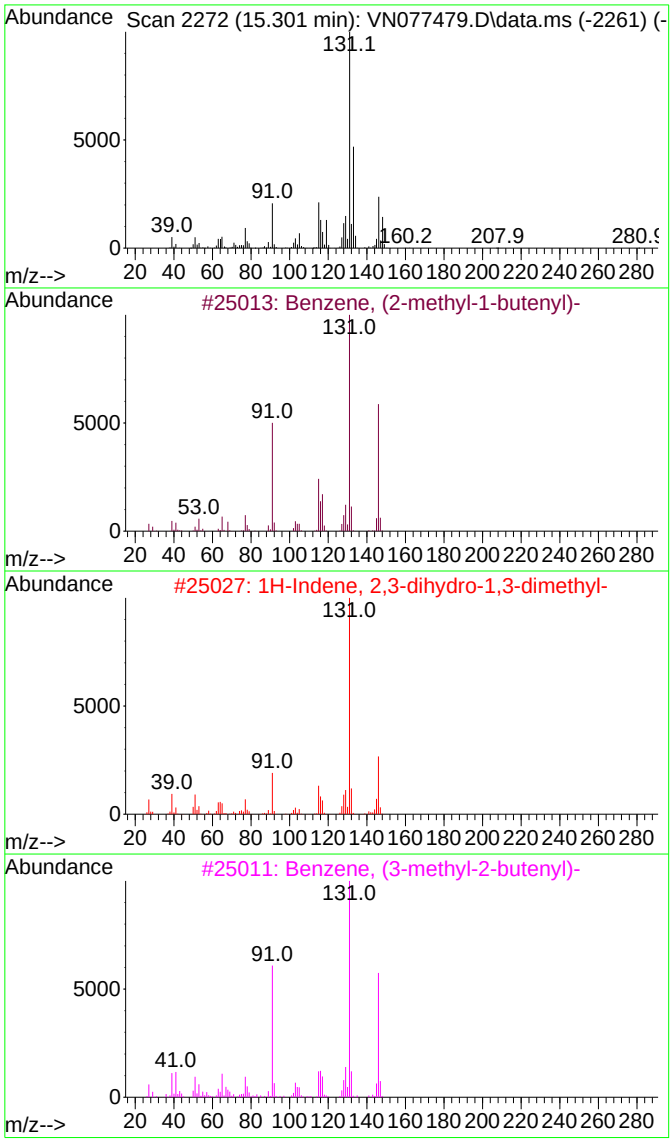
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.M
Quant Title : SW846 8260

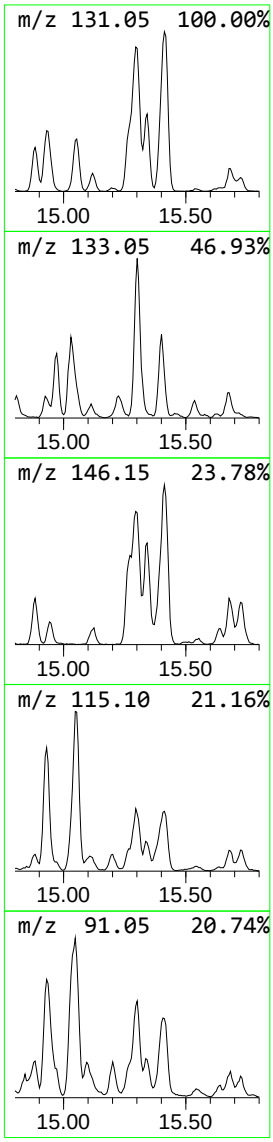
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, (2-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.301	86.55 ug/l	5650440	1,4-Dichlorobenzene-d4	13.795

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	83
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042523\
Data File : VN077479.D
Acq On : 25 Apr 2023 19:09
Operator : JC\MD
Sample : 02460-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042423W.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-	3.254	174.7	ug/l	42802700	1	8.231	12249200	50.0
Pentane, 3,3-di...	5.354	120.1	ug/l	29410700	1	8.231	12249200	50.0
Pentane, 3-methyl-	5.825	86.6	ug/l	21218100	1	8.231	12249200	50.0
Benzene, 1,2-di...	13.954	79.5	ug/l	5190020	4	13.795	3264390	50.0
Benzene, 4-ethy...	14.330	177.0	ug/l	11557500	4	13.795	3264390	50.0
Indan, 1-methyl-	14.448	103.6	ug/l	6766060	4	13.795	3264390	50.0
Benzene, 1,2,4,...	14.654	119.4	ug/l	7797660	4	13.795	3264390	50.0
1H-Indene, 2,3-...	14.930	72.8	ug/l	4753730	4	13.795	3264390	50.0
Benzene, 2-ethe...	15.048	139.5	ug/l	9106640	4	13.795	3264390	50.0
Benzene, (2-met...	15.301	86.5	ug/l	5650440	4	13.795	3264390	50.0