LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 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\82N072822W.M
Title : SW846 8260

Signal : TIC: VN073736.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.269	48	65	67	rBV2	157047	497411	6.34%	0.797%
2	2.399	82	87	102	rVB	262010	471284	6.01%	0.755%
3	3.087	195	204	222	rVB	880779	2027627	25.84%	3.249%
4	3.451	256	266	284	rVB4	46193	127620	1.63%	0.204%
5	3.928	336	347	362	rBV2	84411	230849	2.94%	0.370%
6	4.187	378	391	408	rBV5	30422	122250	1.56%	0.196%
7	4.998	517	529	536	rVV3	119302	465331	5.93%	0.746%
8	5.104	537	547	567	rVV3	169855	700462	8.93%	1.122%
9	5.534	605	620	634	rBV4	74211	257847	3.29%	0.413%
10	6.410	752	769	780	rBV5	40685	153369	1.95%	0.246%
11	7.063	870	880	890	rVV2	269504	811043	10.34%	1.300%
12	7.139	890	893	905	rVB2	34683	80996	1.03%	0.130%
13	7.775	994	1001	1012	rVB3	32709	82440	1.05%	0.132%
14	7.957	1019	1032	1036	rBV	261106	633002	8.07%	1.014%
15	8.022	1036	1043	1051	rVV2	662303	1595401	20.33%	2.556%
16	8.104	1051	1057	1071	rVB3	89949	236852	3.02%	0.380%
17	8.392	1095	1106	1118	rBV	3278600	7847401	100.00%	12.574%
18	8.639	1143	1148	1157	rVB3	45806	106993	1.36%	0.171%
19	8.904	1183	1193	1205	rBV	661141	1416796	18.05%	2.270%
20	9.363	1261	1271	1273	rBV2	71459	144586	1.84%	0.232%
21	9.392	1273	1276	1283	rVB2	80084	155363	1.98%	0.249%
22	9.822	1344	1349	1357	rVB2	44247	88142	1.12%	0.141%
23	9.910	1357	1364	1368	rBV2	103443	208512	2.66%	0.334%
24	9.951	1368	1371	1381	rVB3	87271	171994	2.19%	0.276%
25	10.139	1395	1403	1411	rBV3	36039	90686	1.16%	0.145%
26	10.263	1418	1424	1429	rBV	62901	120700	1.54%	0.193%
27	10.380	1437	1444	1451	rBV	976395	1754678	22.36%	2.812%
28	10.445	1451	1455	1461	rVB	52524	94894	1.21%	0.152%
29	10.580	1468	1478	1485	rVB3	105163	219778	2.80%	0.352%
30	10.657	1485	1491	1497	rBV2	92643	153536	1.96%	0.246%
31	10.798	1508	1515	1524	rVB3	78662	177965	2.27%	0.285%
32	11.686	1657	1666	1677	rBV	960728	1707902	21.76%	2.737%
33	11.786	1677	1683	1693	rVB2	86238	187931	2.39%	0.301%
34	12.521	1802	1808	1816	rBV	754858	1205439	15.36%	1.932%
35	12.674	1826	1834	1844	rBV	824671	1381686	17.61%	2.214%
36	12.863	1858	1866	1875	rBV	1925622	2976254	37.93%	4.769%

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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-1

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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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\82N072822W.M
Title : SW846 8260

37	13.163	1911	1917	1924	rBV	105148	177069	2.26%	0.284%
38	13.310	1936	1942	1949	rVB	94945	158967	2.03%	0.255%
39	13.427	1954	1962	1970	rBV3	263991	641333	8.17%	1.028%
40	13.616	1987	1994	2005	rBV	1025959	1737851	22.15%	2.785%
41	13.710	2005	2010	2015	rVB	70374	106328	1.35%	0.170%
42	13.780	2015	2022	2024	rBV	1105418	1715807	21.86%	2.749%
43	13.815	2024	2028	2033	rVV	2539854	4157377	52.98%	6.662%
44	13.857	2033	2035	2045	rVB2	495730	1086387	13.84%	1.741%
45	13.945	2045	2050	2056	rVB	339132	522495	6.66%	0.837%
46	14.068	2066	2071	2079	rBV	83788	151654	1.93%	0.243%
47	14.157	2079	2086	2092	rBV	1981753	2980468	37.98%	4.776%
48	14.215	2092	2096	2099	rBV	225804	351820	4.48%	0.564%
49	14.268	2099	2105	2112	rVB2	1903809	3122281	39.79%	5.003%
50	14.333	2112	2116	2122	rVB	74610	121301	1.55%	0.194%
51	14.404	2122	2128	2132	rBV	96385	144160	1.84%	0.231%
52	14.480	2132	2141	2146	rBV	1963818	3151758	40.16%	5.050%
53	14.557	2151	2154	2158	rVV	196514	274731	3.50%	0.440%
54	14.598	2158	2161	2164	rVB	77900	87760	1.12%	0.141%
55	14.662	2169	2172	2174	rVV	59224	85173	1.09%	0.136%
56	14.704	2174	2179	2182	rVV	175369	325718	4.15%	0.522%
57	14.751	2182	2187	2192	rVV2	1136071	1963706	25.02%	3.147%
58	14.792	2192	2194	2199	rVB2	180811	234012	2.98%	0.375%
59	14.868	2199	2207	2212	rBV3	1103218	2294059	29.23%	3.676%
60	14.921	2212	2216	2223	rVB3	254494	498184	6.35%	0.798%
61	15.021	2228	2233	2236	rVV	82752	146537	1.87%	0.235%
62	15.127	2237	2251	2256	rVV3	843568	2021869	25.76%	3.240%
63	15.168	2256	2258	2263	rVV2	236922	413760	5.27%	0.663%
64	15.245	2263	2271	2282	rVB3	660391	1586930	20.22%	2.543%
65	15.392	2291	2296	2299	rBV3	58175	103486	1.32%	0.166%
66	15.433	2299	2303	2308	rVB	91187	160990	2.05%	0.258%
67	15.492	2308	2313	2317	rBV2	55641	88780	1.13%	0.142%
68	15.545	2317	2322	2327	rVV2	160607	332197	4.23%	0.532%
69	15.598	2328	2331	2346	rVB3	105390	232045	2.96%	0.372%
70	15.733	2346	2354	2361	rBV	196928	405691	5.17%	0.650%
71	15.904	2374	2383	2393	rBV3	92028	320354	4.08%	0.513%
72	16.033	2399	2405	2410	rBV	153341	315166	4.02%	0.505%
73	16.104	2411	2417	2424	rVB6	94681	262680	3.35%	0.421%
74	16.256	2434	2443	2448	rBV4	37488	87333	1.11%	0.140%
75	16.462	2467	2478	2485	rBV3	57194	143002	1.82%	0.229%
76	16.739	2515	2525	2533	rBV	421329	992458	12.65%	1.590%

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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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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\82N072822W.M
Title : SW846 8260

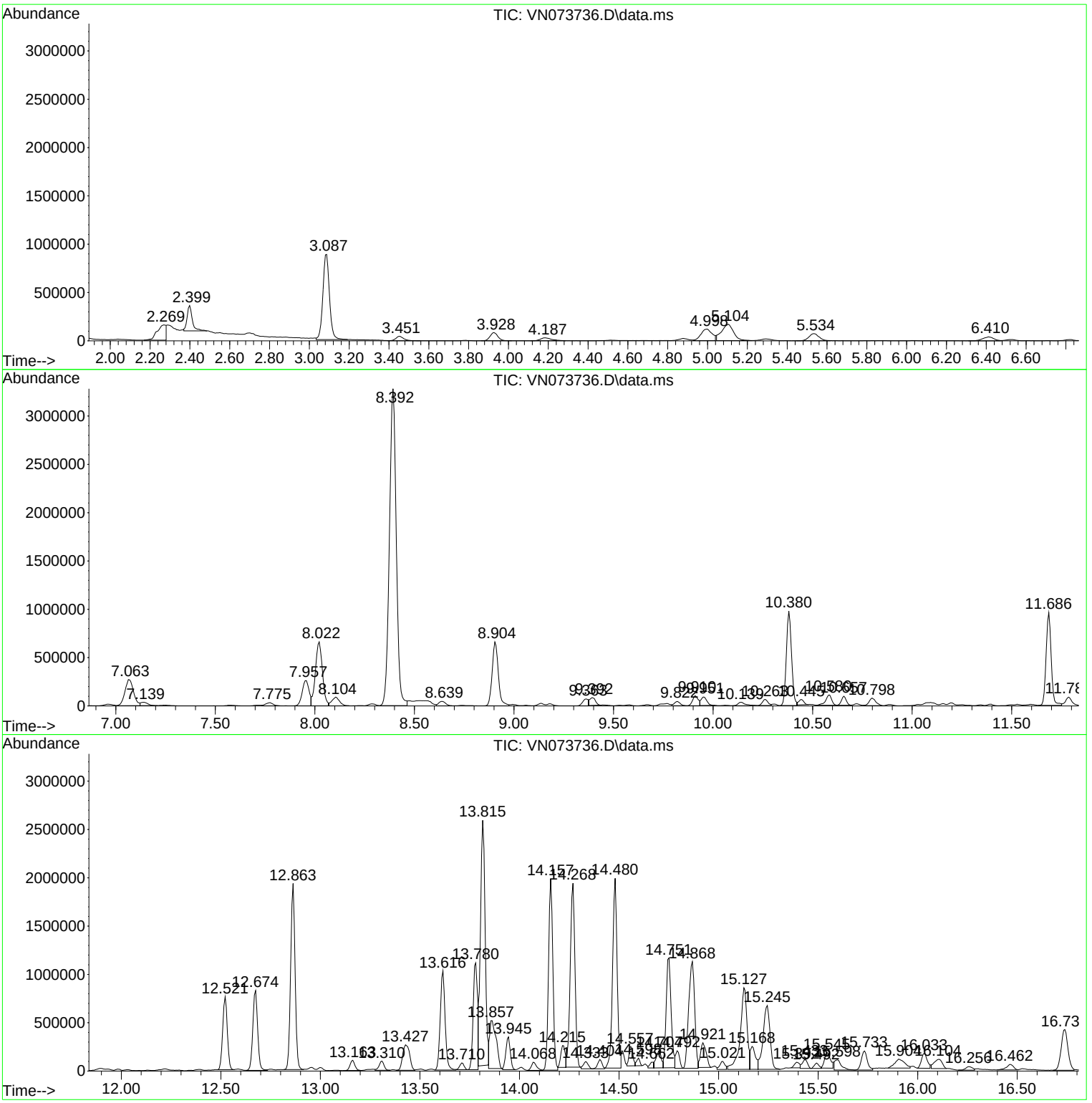
Sum of corrected areas: 62408697

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
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MSVOA_N
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

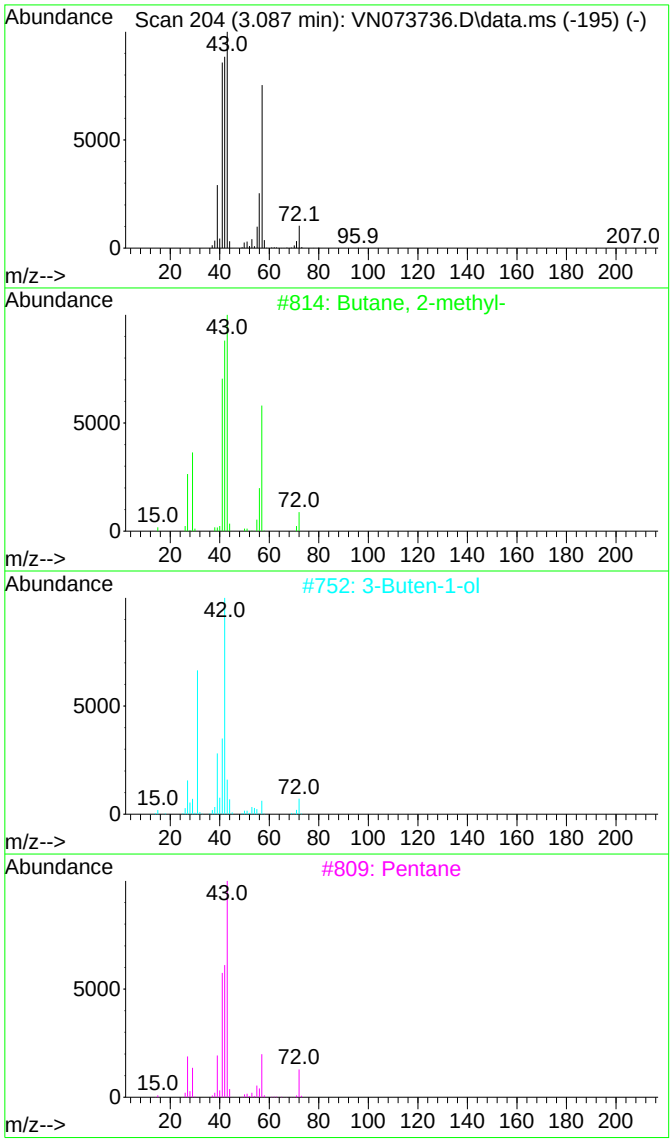
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

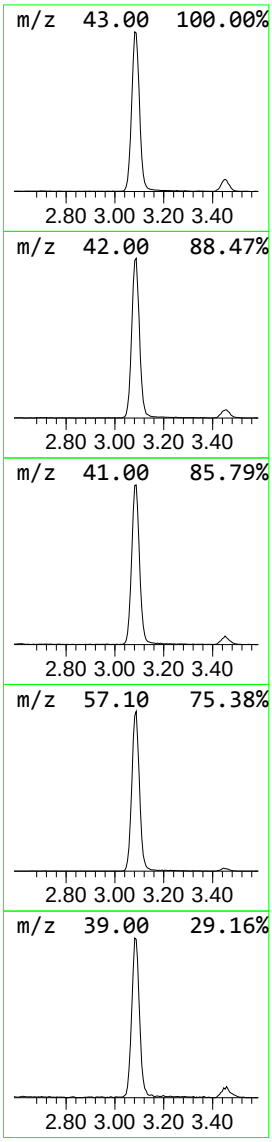
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.087	63.55 ug/l	2027630	Pentafluorobenzene	8.022

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	87
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	28
4	1-Butanesulfonyl chloride	156	C4H9ClO2S	002386-60-9	12
5	1-Butene	56	C4H8	000106-98-9	10





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Data File : VN073736.D
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Instrument :
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ClientSampleId :
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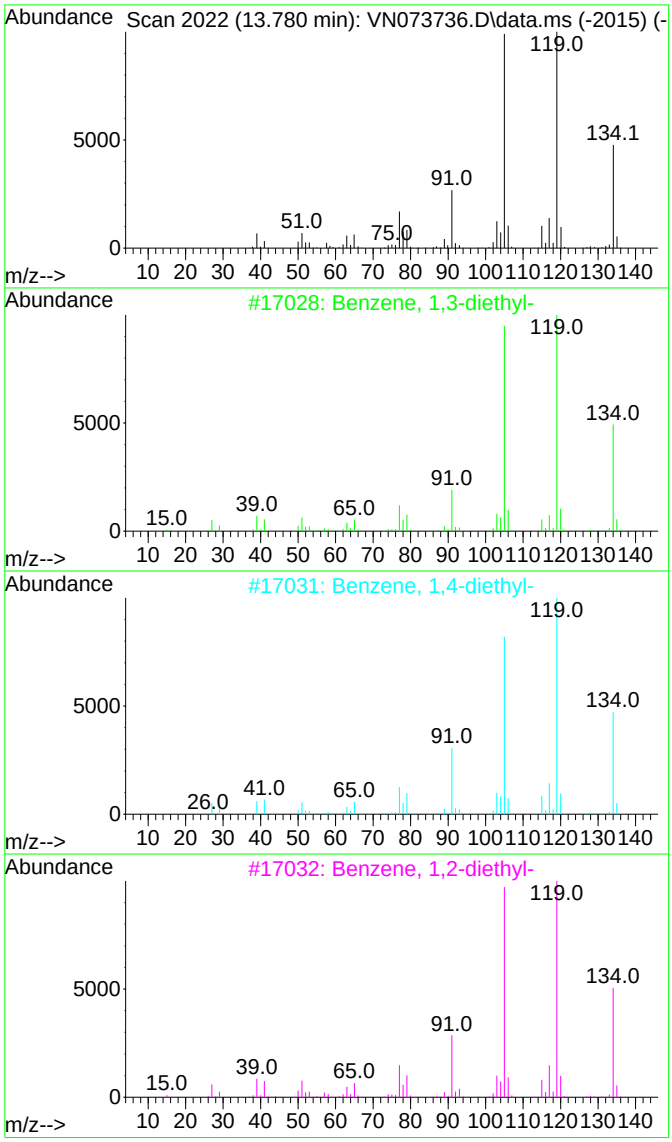
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

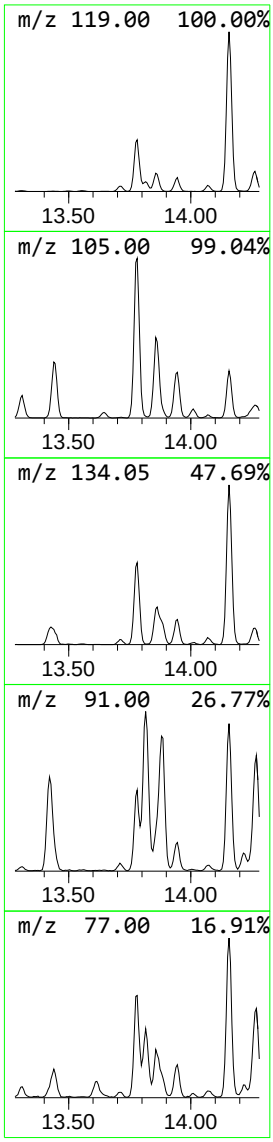
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,3-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	49.37 ug/l	1715810	1,4-Dichlorobenzene-d4	13.616

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	97
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-1,3-dimethyl-	134	C10H14	002870-04-4	83





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Data File : VN073736.D
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Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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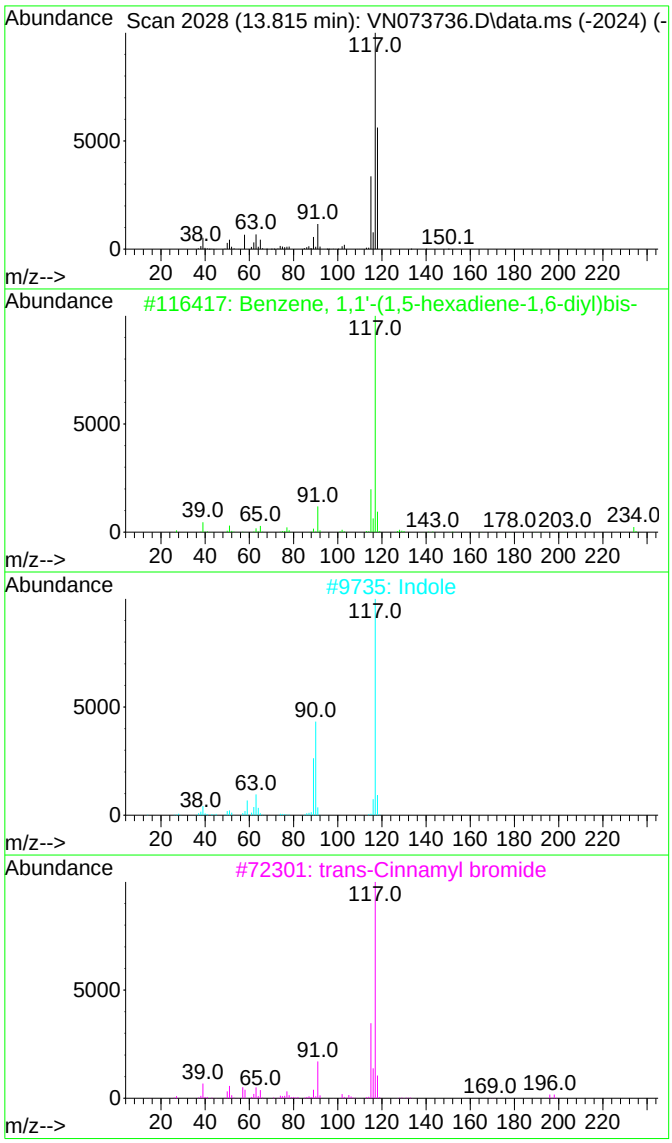
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

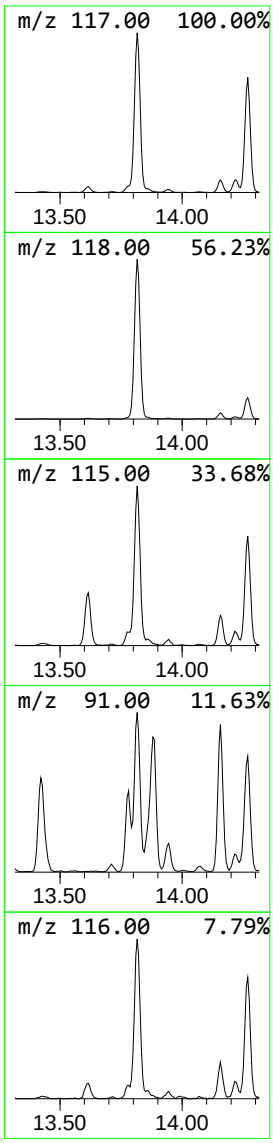
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	119.61 ug/l	4157380	1,4-Dichlorobenzene-d4	13.616

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	Indole	117	C8H7N	000120-72-9	46
3	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42
4	Indolizine	117	C8H7N	000274-40-8	37
5	Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	36





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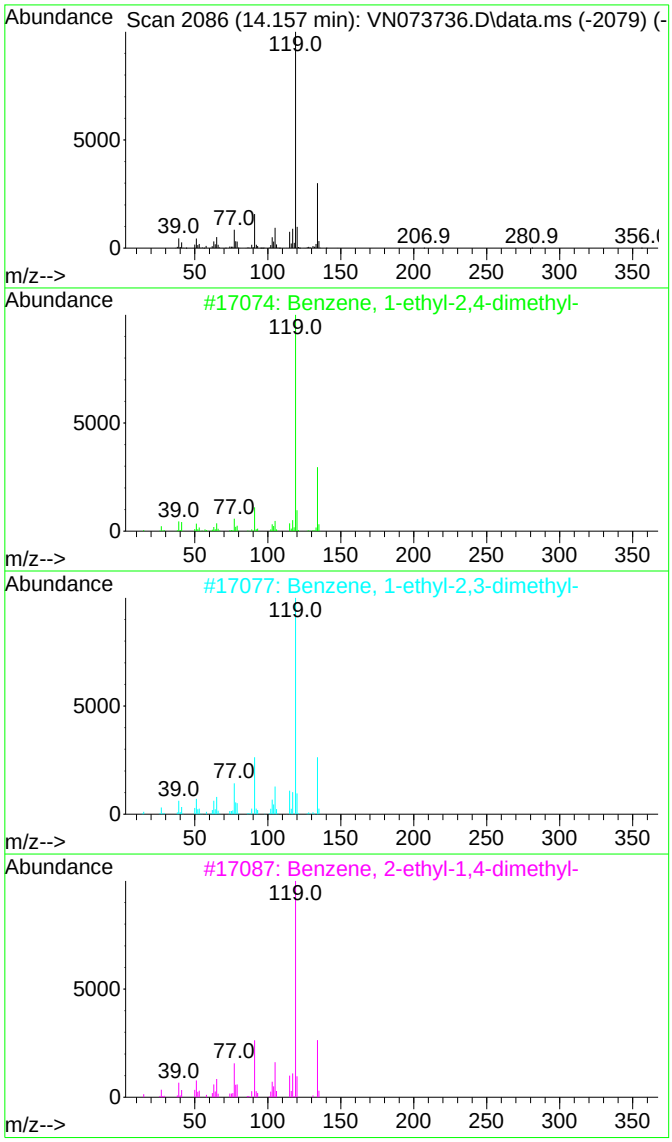
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

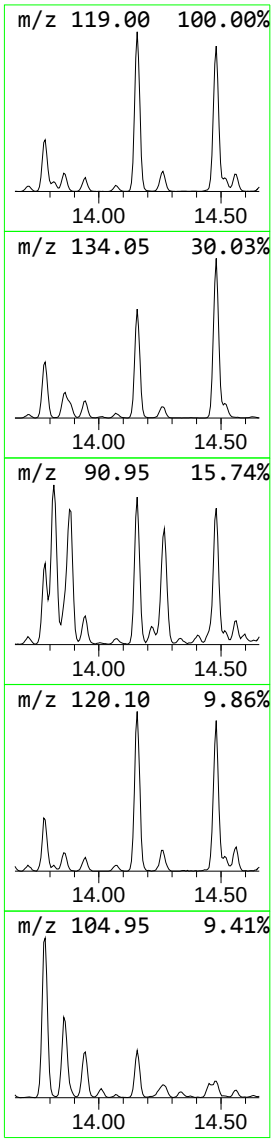
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.157	85.75 ug/l	2980470	1,4-Dichlorobenzene-d4	13.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
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





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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

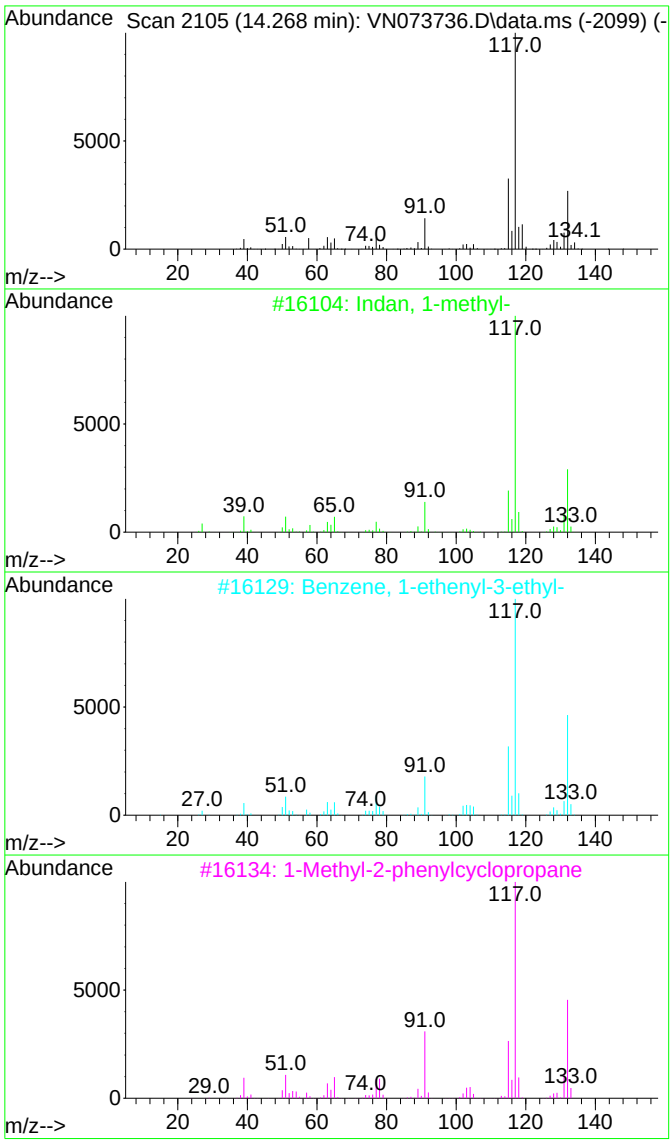
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

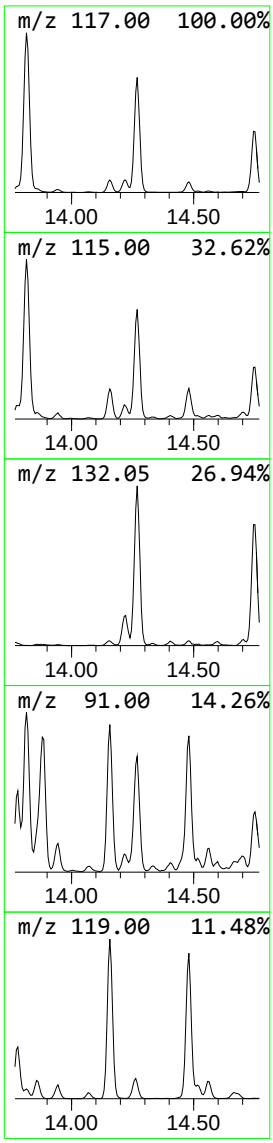
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.268	89.83 ug/l	3122280	1,4-Dichlorobenzene-d4	13.616

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-methyl-1-propenyl)-	132	C10H12	000768-49-0	72
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	72





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Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

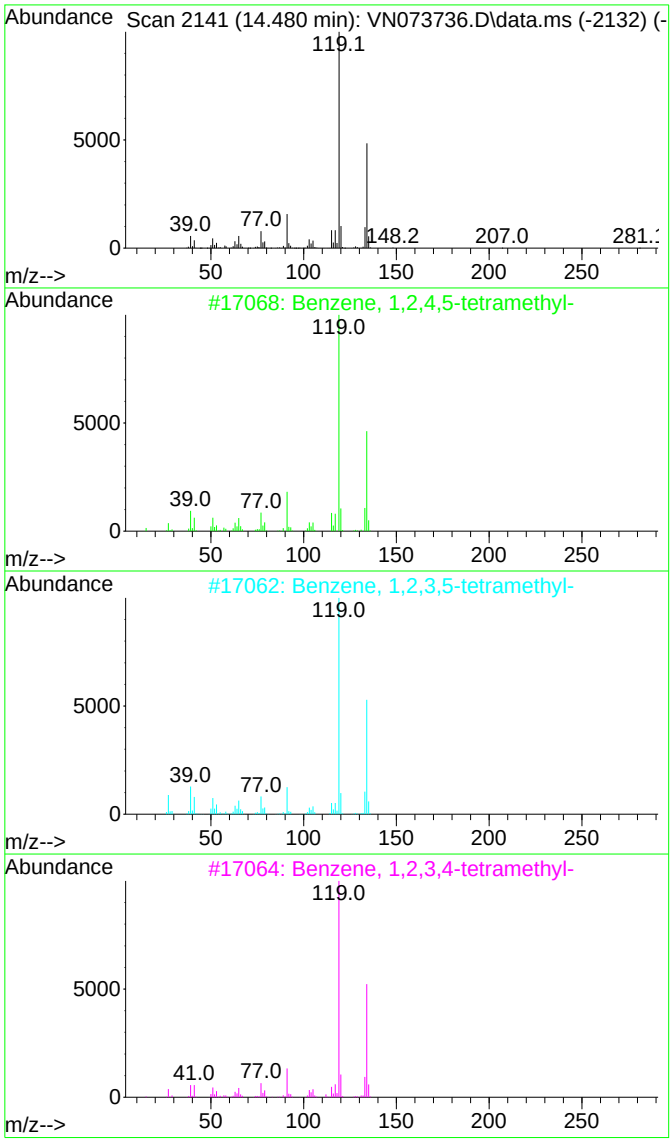
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

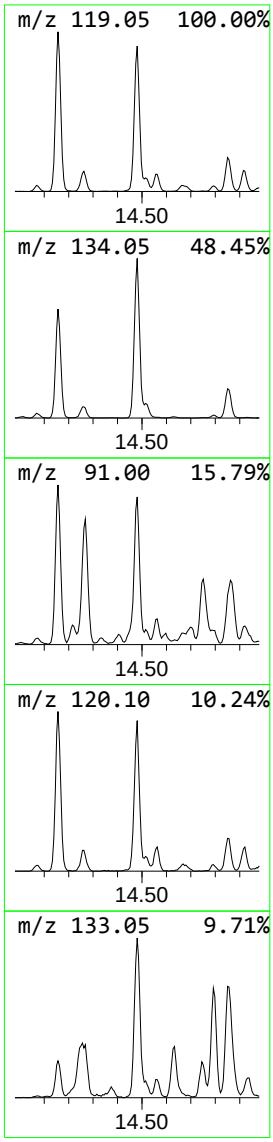
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	90.68 ug/l	3151760	1,4-Dichlorobenzene-d4	13.616

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	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





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Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

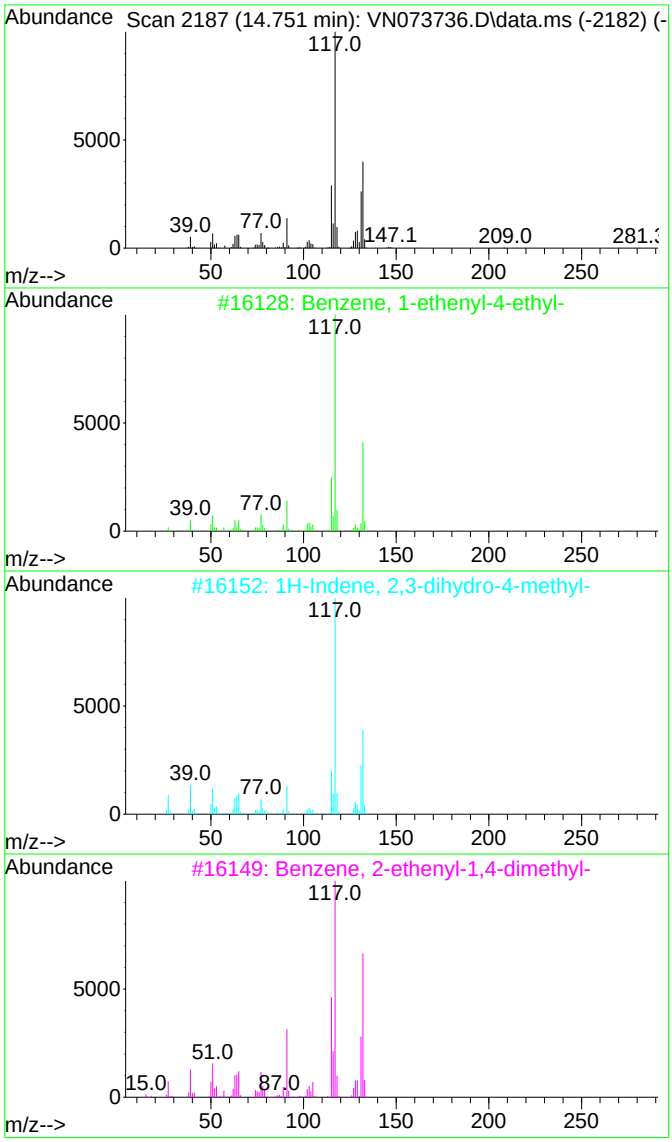
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

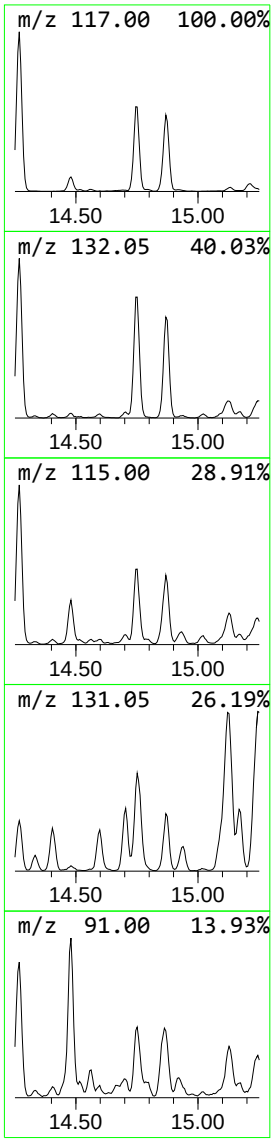
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.751	56.50 ug/l	1963710	1,4-Dichlorobenzene-d4	13.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

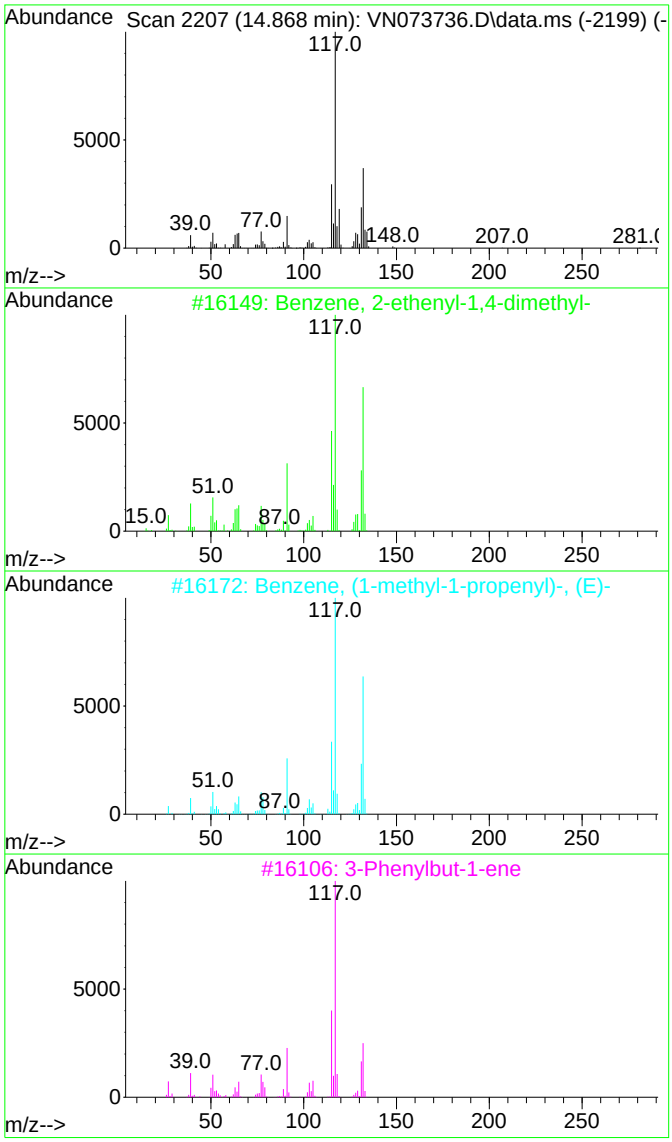
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

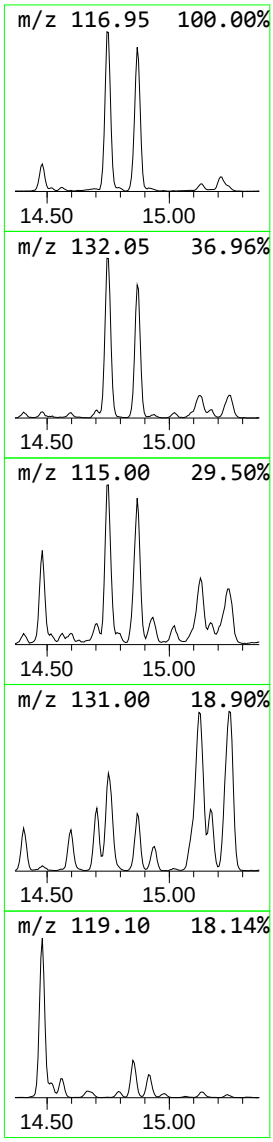
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	66.00 ug/l	2294060	1,4-Dichlorobenzene-d4	13.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

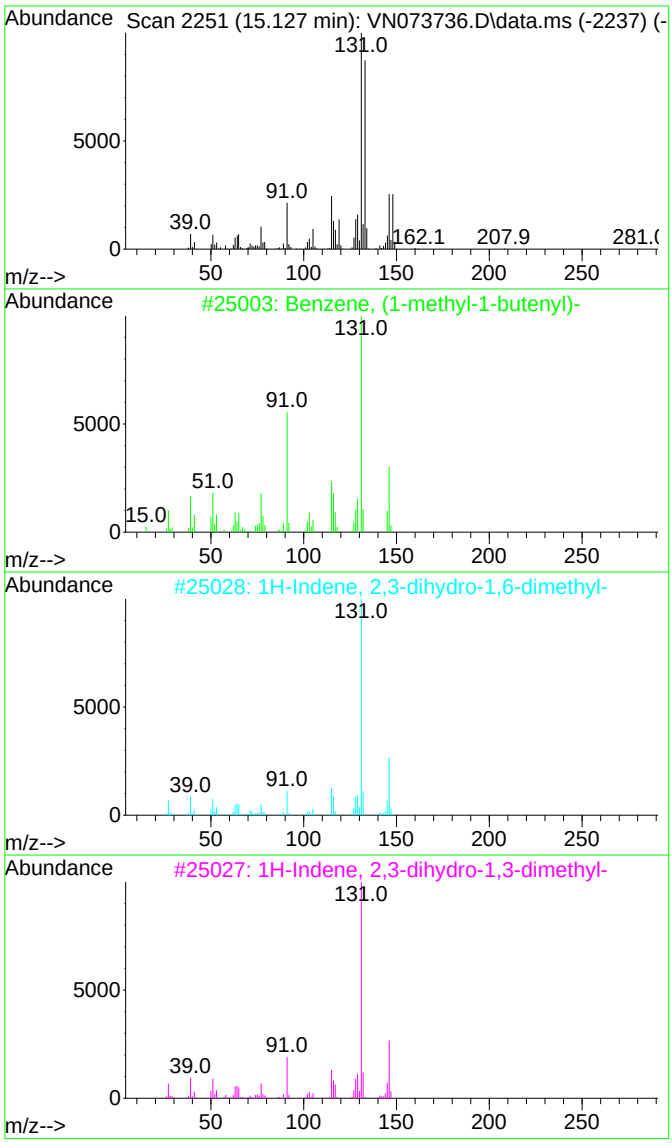
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

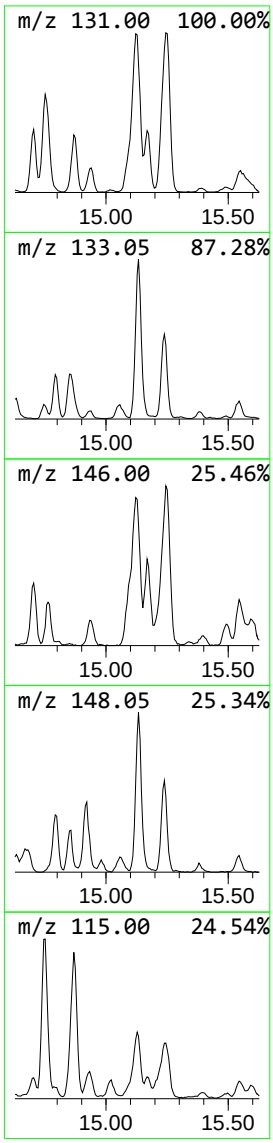
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, (1-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.127	58.17 ug/l	2021870	1,4-Dichlorobenzene-d4	13.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	86
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	86
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64





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Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

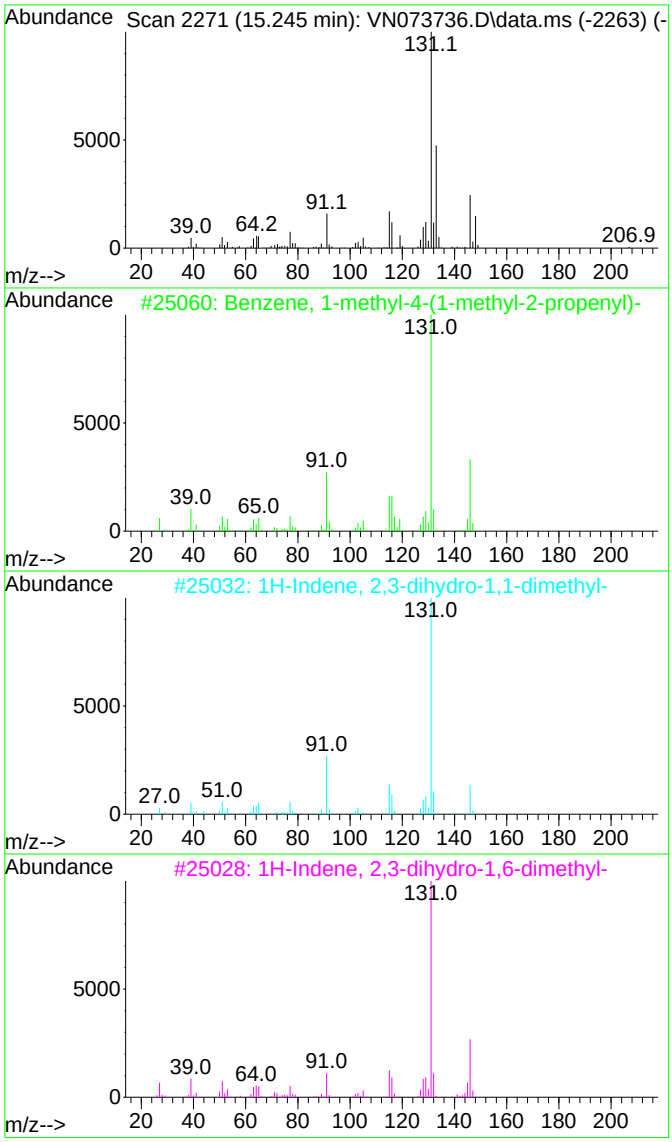
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260

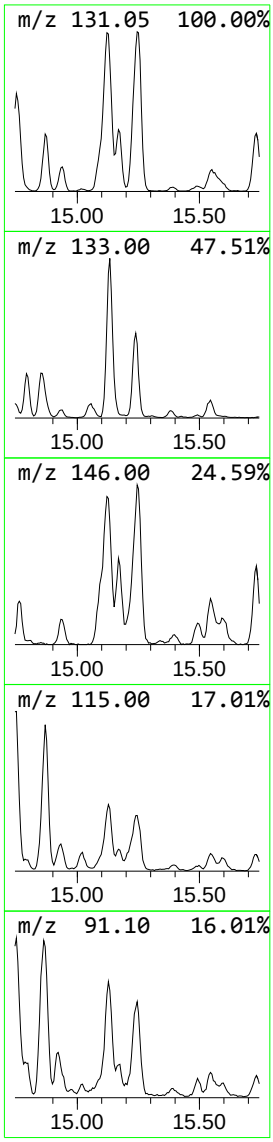
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-methyl-4-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.245	45.66 ug/l	1586930	1,4-Dichlorobenzene-d4	13.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	86
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	70
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	70
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
Data File : VN073736.D
Acq On : 04 Aug 2022 15:05
Operator : JC\MD
Sample : N3887-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.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-	3.087	63.5	ug/l	2027630	1	8.022	1595400	50.0
Benzene, 1,3-di...	13.780	49.4	ug/l	1715810	4	13.616	1737850	50.0
Benzene, 1,1'-(...	13.816	119.6	ug/l	4157380	4	13.616	1737850	50.0
Benzene, 1-ethy...	14.157	85.8	ug/l	2980470	4	13.616	1737850	50.0
Indan, 1-methyl-	14.268	89.8	ug/l	3122280	4	13.616	1737850	50.0
Benzene, 1,2,4,...	14.480	90.7	ug/l	3151760	4	13.616	1737850	50.0
Benzene, 1-ethe...	14.751	56.5	ug/l	1963710	4	13.616	1737850	50.0
Benzene, 2-ethe...	14.868	66.0	ug/l	2294060	4	13.616	1737850	50.0
Benzene, (1-met...	15.127	58.2	ug/l	2021870	4	13.616	1737850	50.0
Benzene, 1-meth...	15.245	45.7	ug/l	1586930	4	13.616	1737850	50.0