LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

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_X\Method\82X061822W.M
Title : SW846 8260

Signal : TIC: VX029756.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.746	103	108	120	rVB	211301	299566	3.10%	0.464%
2	2.337	195	205	210	rBV	60395	120601	1.25%	0.187%
3	2.782	267	278	281	rBV	134680	297942	3.09%	0.462%
4	2.825	281	285	300	rVV	235485	563591	5.84%	0.874%
5	2.959	300	307	320	rVV	125544	278859	2.89%	0.432%
6	3.105	322	331	345	rVB2	339590	802069	8.31%	1.244%
7	3.763	425	439	446	rBV3	64401	227160	2.35%	0.352%
8	4.215	502	513	520	rBV2	67754	198623	2.06%	0.308%
9	4.306	520	528	540	rVB	191061	547699	5.68%	0.849%
10	5.391	692	706	713	rBV	149864	446068	4.62%	0.692%
11	5.556	713	733	739	rVV3	279256	1270870	13.17%	1.970%
12	5.623	739	744	765	rVB2	170613	549969	5.70%	0.853%
13	5.830	765	778	790	rBV2	159073	483647	5.01%	0.750%
14	5.958	790	799	805	rVV	152227	397826	4.12%	0.617%
15	6.044	805	813	823	rVV	720191	1937111	20.08%	3.003%
16	6.159	824	832	838	rVV2	141205	454218	4.71%	0.704%
17	6.251	838	847	857	rVV5	375138	1324807	13.73%	2.054%
18	6.361	857	865	878	rVB2	110051	326221	3.38%	0.506%
19	6.763	922	931	939	rBV	410386	1003340	10.40%	1.556%
20	6.848	939	945	957	rVB2	80700	239940	2.49%	0.372%
21	7.178	991	999	1006	rBV	87625	191776	1.99%	0.297%
22	7.366	1020	1030	1042	rBV4	211967	624779	6.48%	0.969%
23	7.659	1073	1078	1090	rVB2	41165	97853	1.01%	0.152%
24	7.775	1090	1097	1104	rVB2	52681	105853	1.10%	0.164%
25	7.885	1112	1115	1123	rVB5	66584	141880	1.47%	0.220%
26	7.988	1123	1132	1143	rVB	211781	473889	4.91%	0.735%
27	8.110	1143	1152	1156	rBV	381929	814379	8.44%	1.263%
28	8.190	1162	1165	1175	rVB4	56390	98923	1.03%	0.153%
29	8.311	1178	1185	1191	rBV4	61235	124888	1.29%	0.194%
30	8.507	1211	1217	1226	rVB3	207041	397082	4.12%	0.616%
31	8.604	1226	1233	1236	rBV3	130682	260699	2.70%	0.404%
32	8.653	1236	1241	1249	rVB	843469	1356367	14.06%	2.103%
33	8.824	1265	1269	1278	rVV3	66700	165464	1.72%	0.257%
34	8.927	1282	1286	1290	rVV	82085	135932	1.41%	0.211%
35	9.104	1307	1315	1320	rBV	136201	249309	2.58%	0.387%
36	9.165	1320	1325	1330	rBV	294384	450883	4.67%	0.699%

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MSVOA_X
ClientSampleId :
MW-14

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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Title : SW846 8260

37	9.458	1370	1373	1377	rVV2	107536	174883	1.81%	0.271%
38	9.506	1377	1381	1387	rVB3	113621	223123	2.31%	0.346%
39	9.762	1416	1423	1430	rBV3	89048	155970	1.62%	0.242%
40	10.055	1467	1471	1476	rVB	810401	1040797	10.79%	1.614%
41	10.134	1481	1484	1490	rVV3	60643	145629	1.51%	0.226%
42	10.195	1490	1494	1499	rVB	1569525	1947852	20.19%	3.020%
43	10.305	1509	1512	1518	rVB	188318	250796	2.60%	0.389%
44	10.585	1550	1558	1562	rBV2	60605	103053	1.07%	0.160%
45	10.781	1583	1590	1593	rBV6	49988	104535	1.08%	0.162%
46	10.963	1615	1620	1626	rBV	879165	1114517	11.55%	1.728%
47	11.085	1635	1640	1644	rBV	626031	792472	8.21%	1.229%
48	11.305	1672	1676	1685	rVB	876892	1109906	11.50%	1.721%
49	11.402	1686	1692	1696	rBV	243386	309498	3.21%	0.480%
50	11.616	1721	1727	1735	rBV	523296	659397	6.83%	1.022%
51	11.866	1763	1768	1770	rBV	136336	181652	1.88%	0.282%
52	11.890	1770	1772	1778	rVB	194678	225491	2.34%	0.350%
53	12.024	1789	1794	1798	rBV	924914	1122116	11.63%	1.740%
54	12.091	1802	1805	1816	rVB3	62887	123180	1.28%	0.191%
55	12.183	1816	1820	1825	rVB	85178	105809	1.10%	0.164%
56	12.244	1825	1830	1837	rBV	7664574	9647902	100.00%	14.958%
57	12.317	1837	1842	1852	rVB2	818526	1229371	12.74%	1.906%
58	12.408	1852	1857	1862	rBV2	358205	458623	4.75%	0.711%
59	12.463	1862	1866	1872	rVB2	152515	218377	2.26%	0.339%
60	12.530	1872	1877	1884	rBV	154715	200517	2.08%	0.311%
61	12.616	1886	1891	1893	rBV	954720	1079333	11.19%	1.673%
62	12.646	1893	1896	1900	rVB	775368	876932	9.09%	1.360%
63	12.701	1900	1905	1913	rBV	2863596	3745355	38.82%	5.807%
64	12.841	1924	1928	1933	rVB	118034	158079	1.64%	0.245%
65	12.920	1933	1941	1946	rBV	1299400	1681546	17.43%	2.607%
66	13.024	1954	1958	1967	rVB2	273801	436523	4.52%	0.677%
67	13.140	1967	1977	1978	rBV3	213290	391293	4.06%	0.607%
68	13.170	1978	1982	1991	rVB	1562863	1951448	20.23%	3.025%
69	13.292	1997	2002	2007	rVB2	3761058	4826496	50.03%	7.483%
70	13.347	2007	2011	2014	rBV3	137368	162101	1.68%	0.251%
71	13.426	2020	2024	2031	rVB	420206	549590	5.70%	0.852%
72	13.506	2031	2037	2040	rBV2	215284	319784	3.31%	0.496%
73	13.542	2040	2043	2047	rVV	700535	1021167	10.58%	1.583%
74	13.585	2047	2050	2054	rVV3	565729	833319	8.64%	1.292%
75	13.640	2054	2059	2060	rVV2	342645	481868	4.99%	0.747%
76	13.664	2060	2063	2069	rVB2	623570	954181	9.89%	1.479%

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Integration Parameters: RTEINT.P
Integrator: RTE
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Title : SW846 8260

77	13.774	2074	2081	2091	rVV	1126522	1575766	16.33%	2.443%
78	13.853	2091	2094	2099	rVB	89959	125292	1.30%	0.194%
79	13.926	2099	2106	2110	rBV2	334408	474627	4.92%	0.736%
80	13.969	2110	2113	2117	rVB	261897	310107	3.21%	0.481%
81	14.079	2126	2131	2137	rBV	524256	676028	7.01%	1.048%
82	14.219	2149	2154	2158	rBV	284247	450939	4.67%	0.699%
83	14.323	2167	2171	2176	rBV	247605	330090	3.42%	0.512%
84	14.371	2176	2179	2184	rVB	346778	427513	4.43%	0.663%
85	14.481	2192	2197	2201	rBV2	120859	178039	1.85%	0.276%
86	14.609	2210	2218	2220	rBV3	82490	169923	1.76%	0.263%
87	14.640	2220	2223	2227	rVB	215572	263740	2.73%	0.409%
88	14.780	2241	2246	2251	rBV	611007	795683	8.25%	1.234%
89	15.615	2374	2383	2387	rBV	90485	148418	1.54%	0.230%

Sum of corrected areas: 64500729

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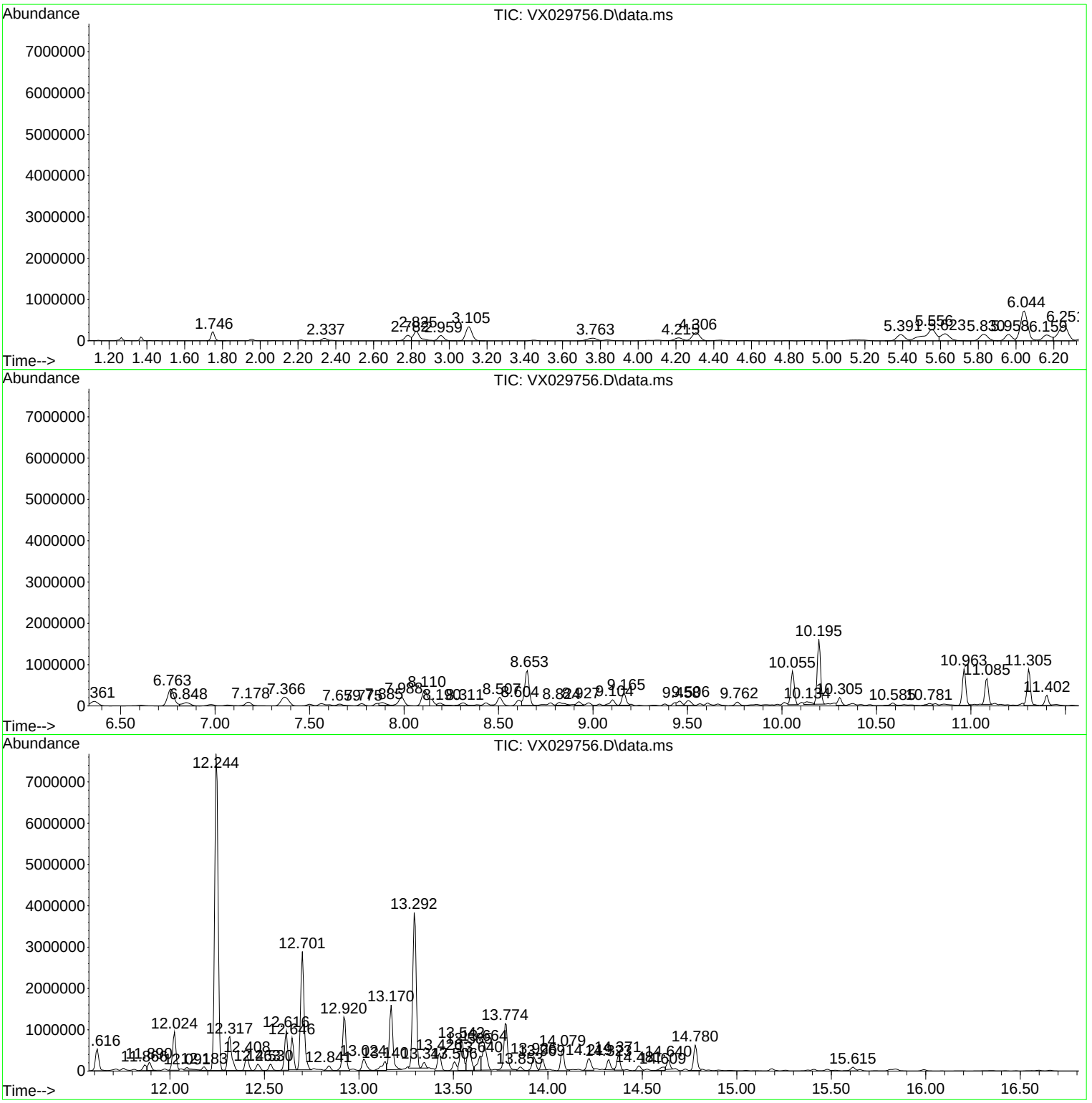
MSVOA_X
ClientSampleId :
MW-14

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
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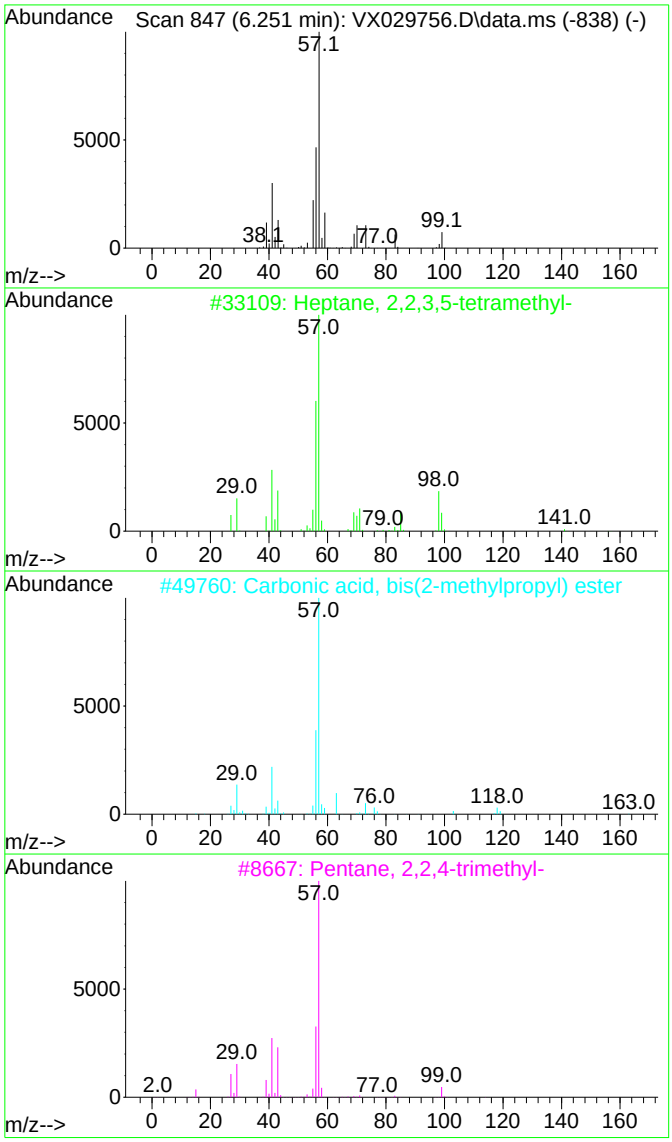
Quant Title : SW846 8260

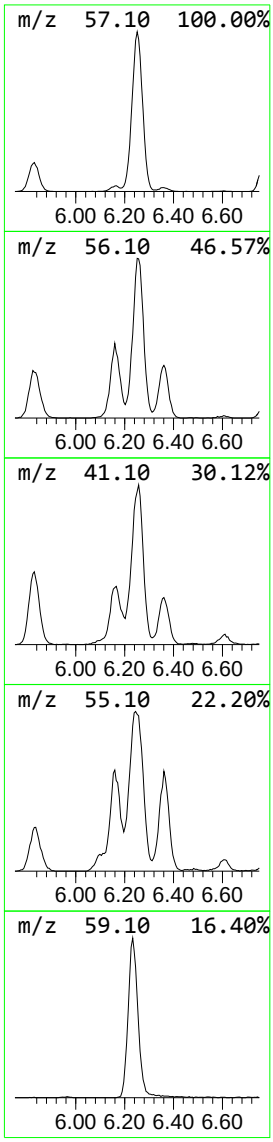
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 unknown6.251 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	66.02 ug/l	1324810	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	45
2	Carbonic acid, bis(2-methylpropy...	174	C9H18O3	000539-92-4	40
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	38
4	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	38
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	37





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Data File : VX029756.D
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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

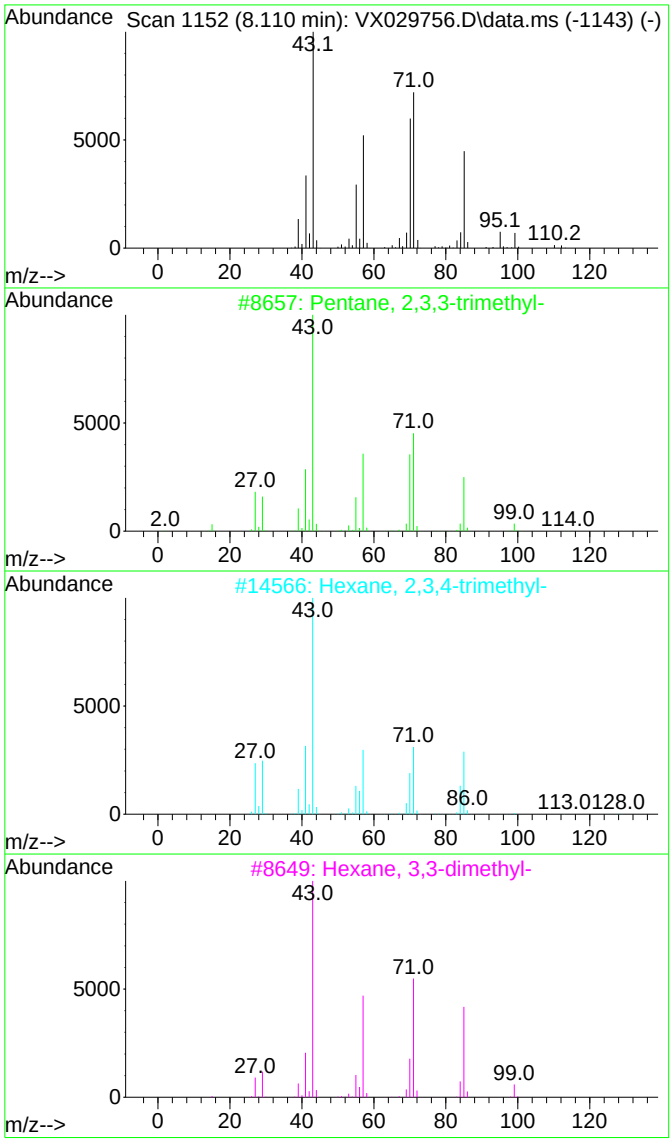
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

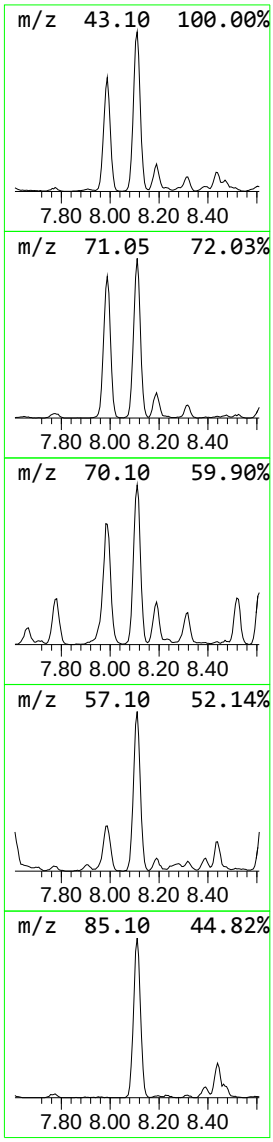
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2,3,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	40.58 ug/l	814379	1,4-Difluorobenzene	6.763

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	59
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	50





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Instrument :
MSVOA_X
ClientSampleId :
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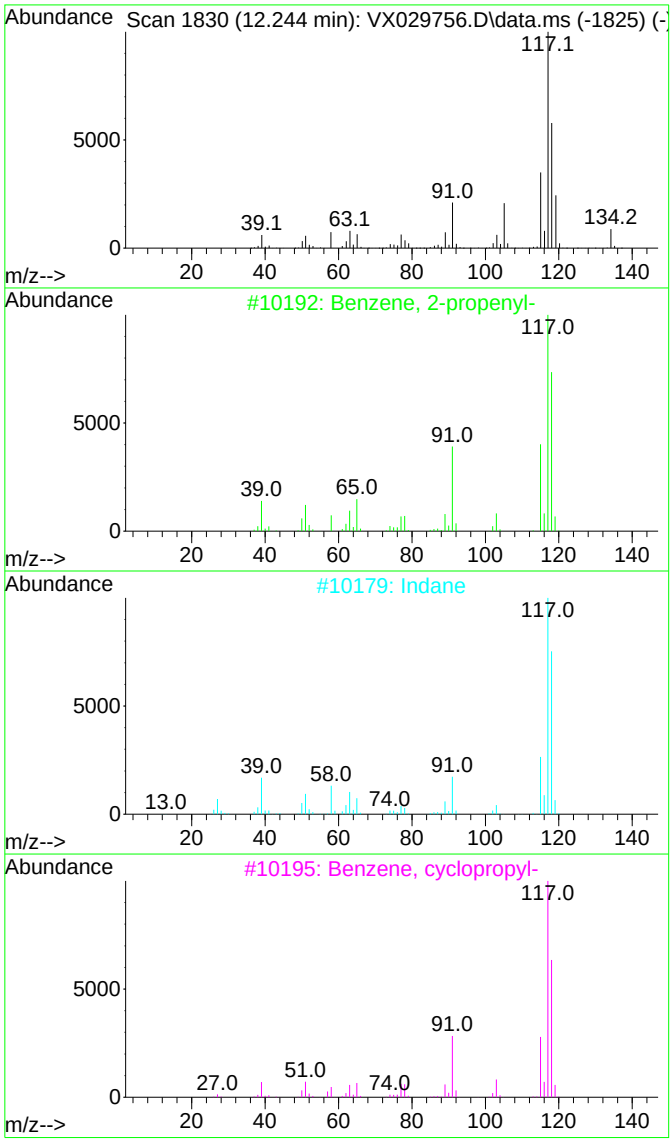
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

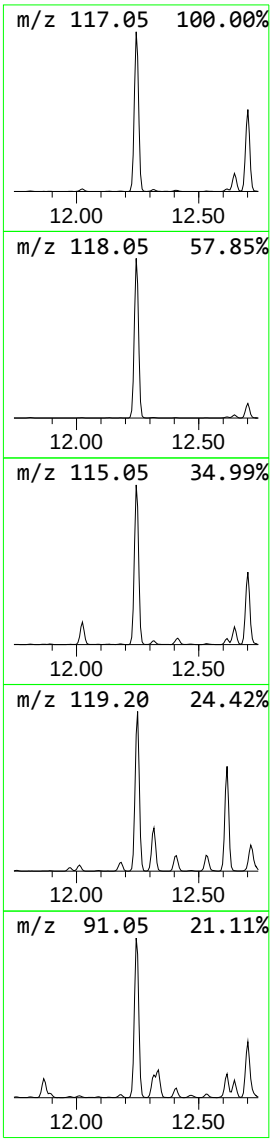
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 2-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	429.90 ug/l	9647900	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
2	Indane	118	C9H10	000496-11-7	70
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4	Deltacyclene	118	C9H10	007785-10-6	52
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	46





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ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
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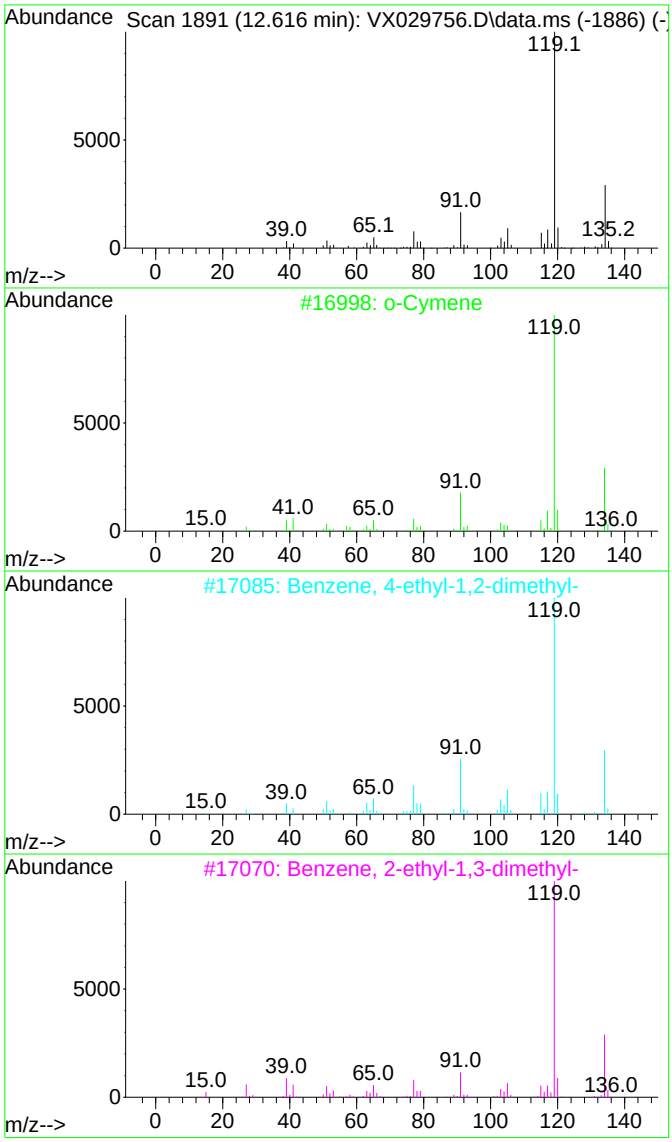
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

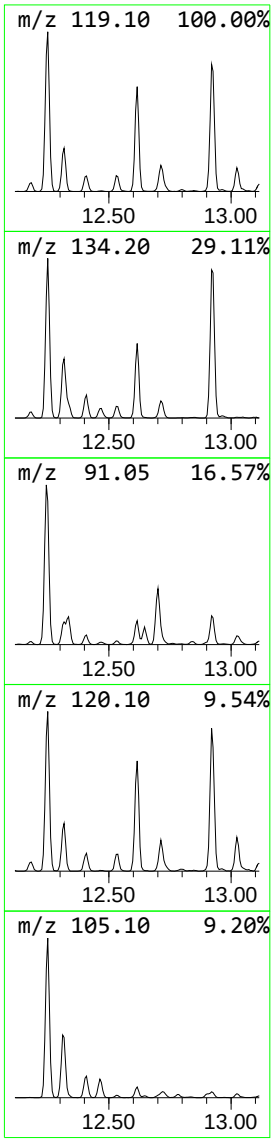
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	48.09 ug/l	1079330	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

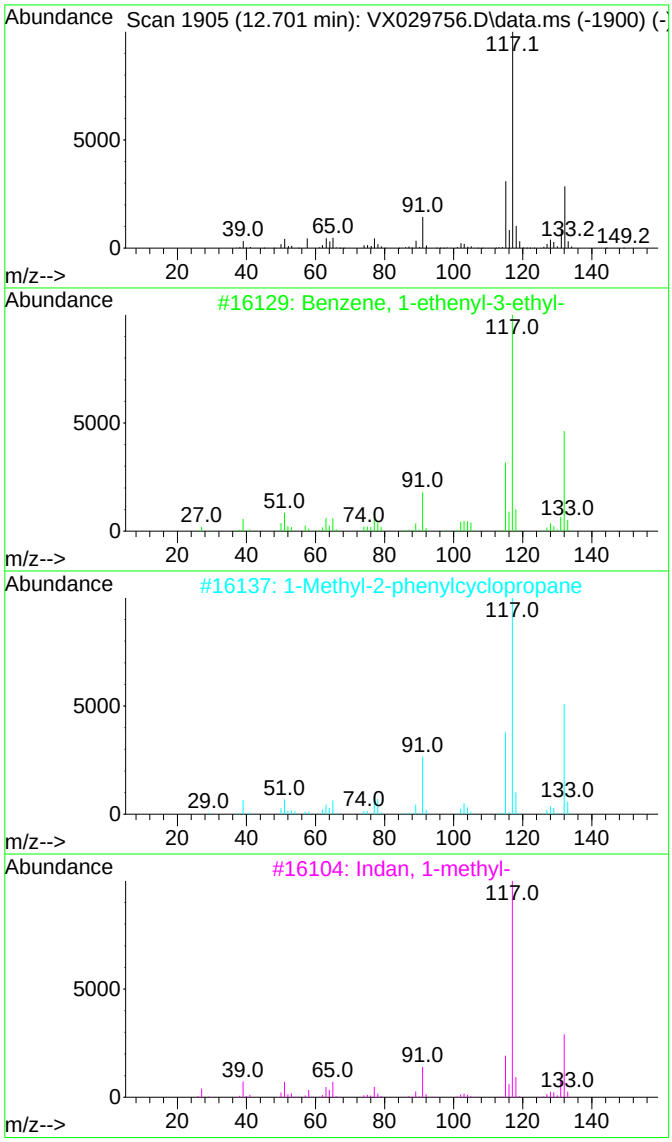
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

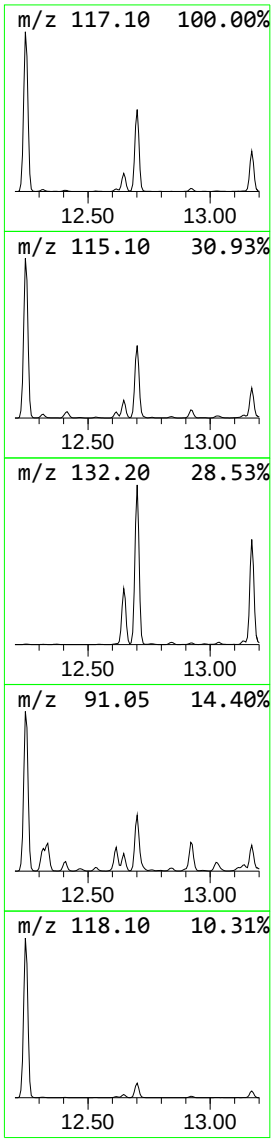
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	166.89 ug/l	3745360	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	90
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
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ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

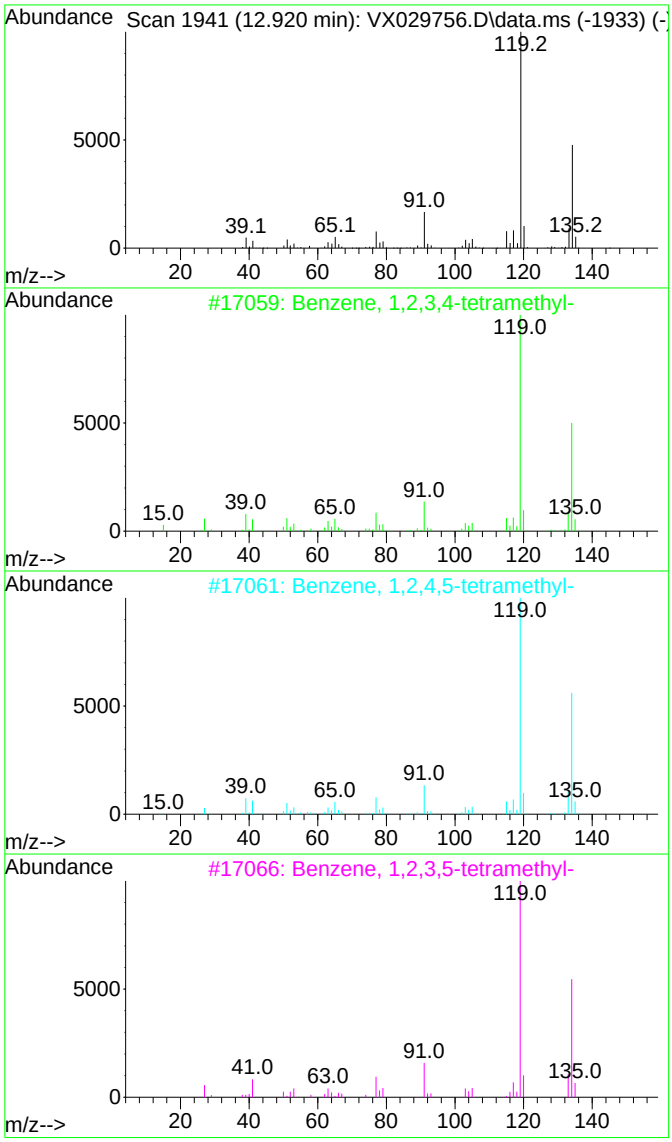
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

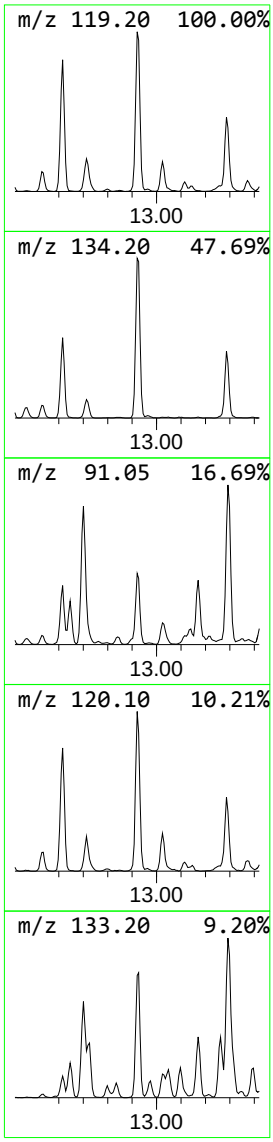
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	74.93 ug/l	1681550	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

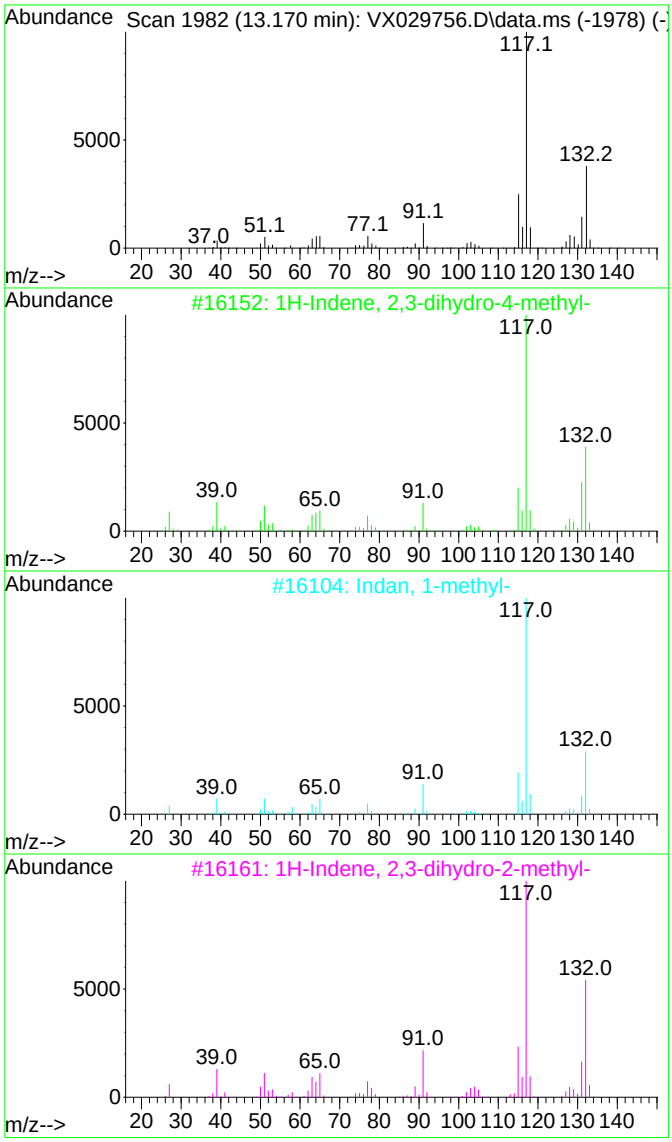
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

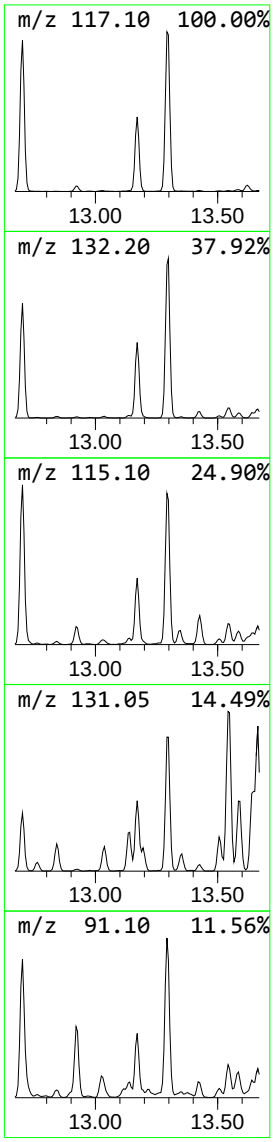
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	86.95 ug/l	1951450	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

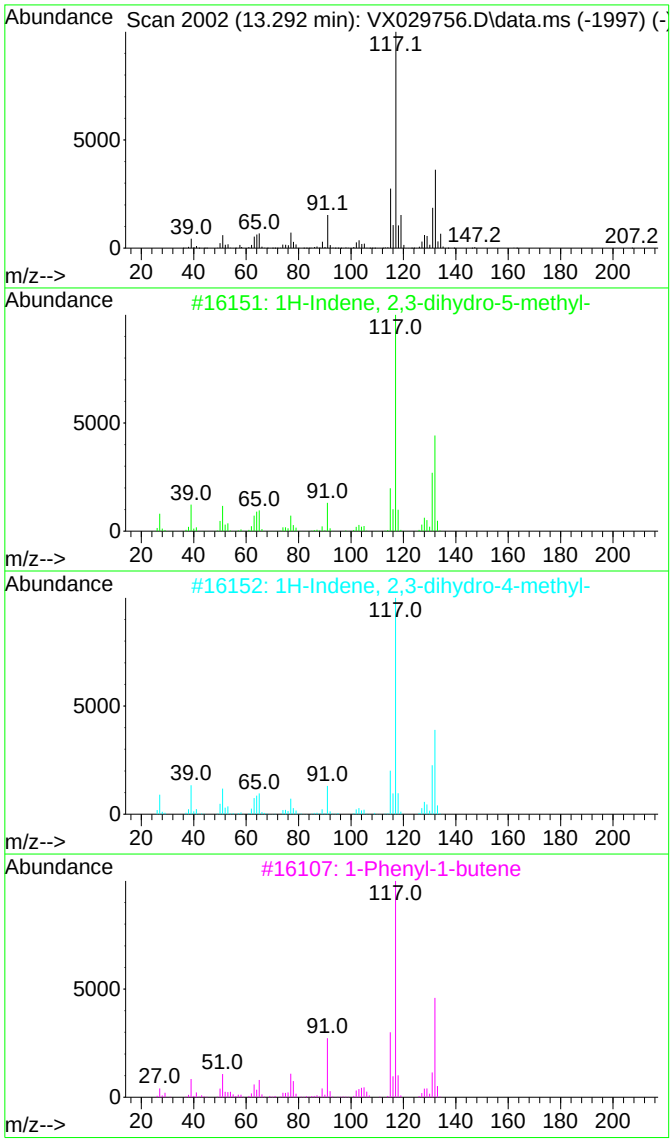
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

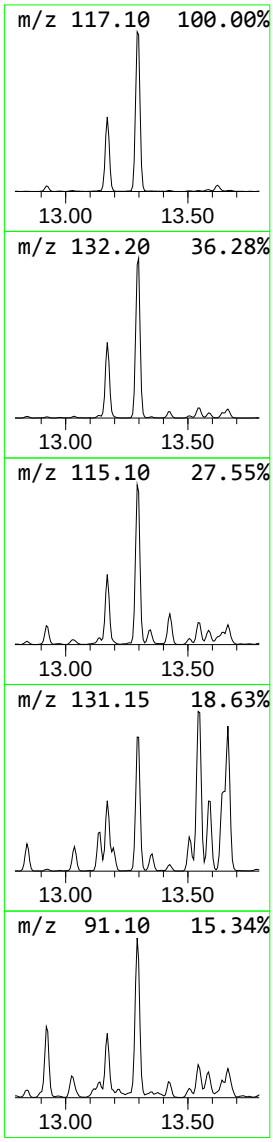
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	215.06 ug/l	4826500	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

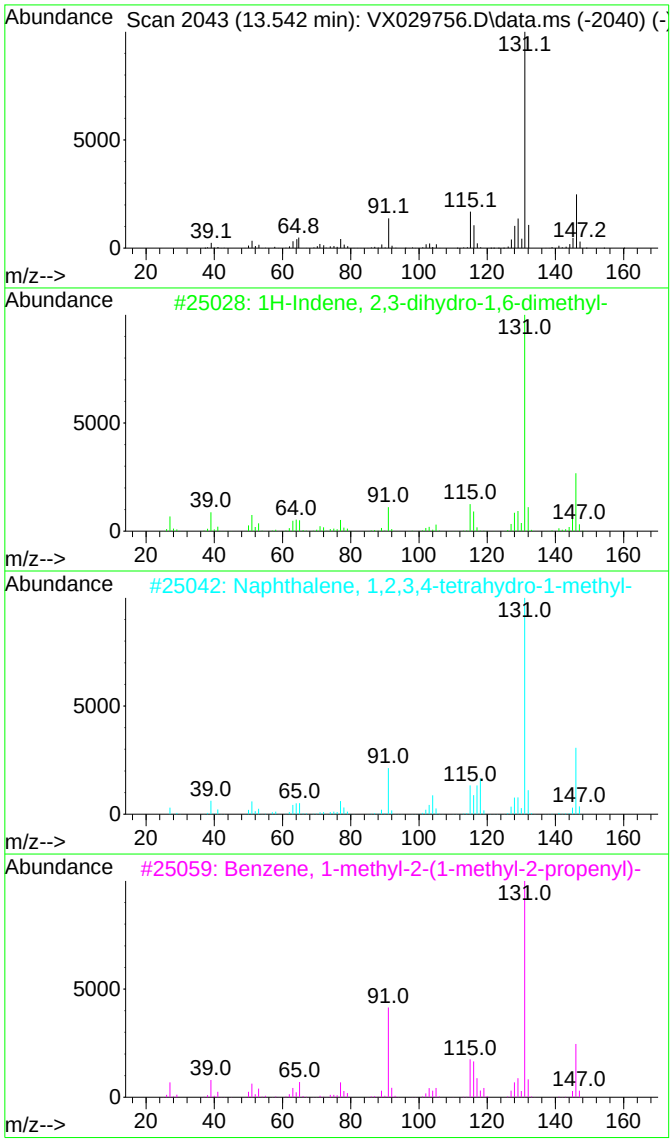
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

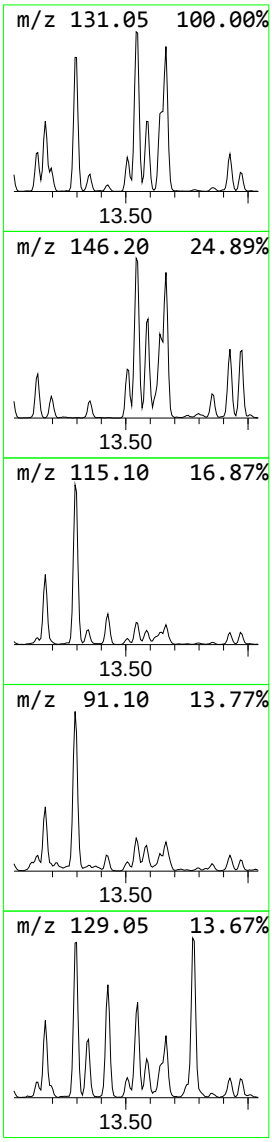
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	45.50 ug/l	1021170	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
3	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91





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Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

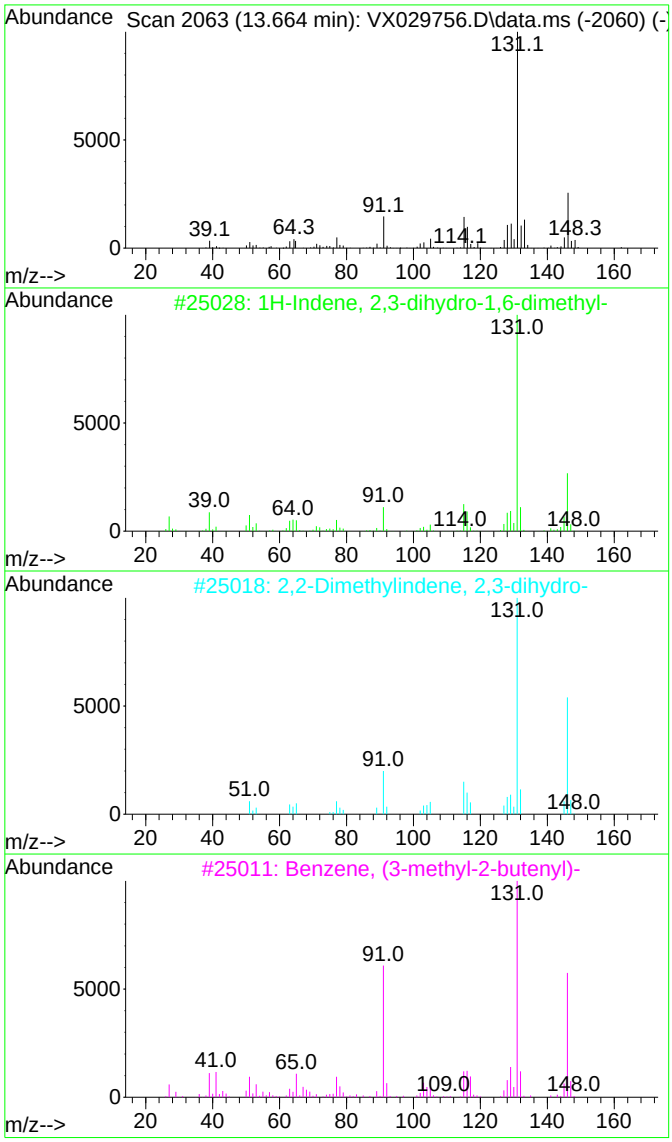
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

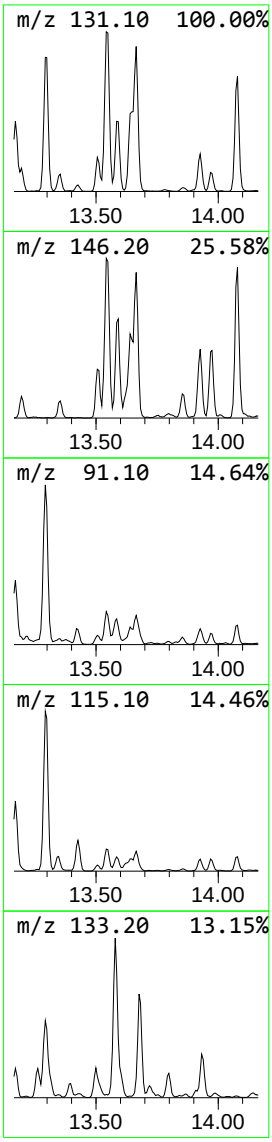
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2,2-Dimethylindene, 2,3-dih... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	42.52 ug/l	954181	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062322\
Data File : VX029756.D
Acq On : 23 Jun 2022 19:57
Operator : JC/MD
Sample : N3366-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.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
unknown6.251	6.251	66.0	ug/l	1324810	2	6.763	1003340	50.0
Pentane, 2,3,3-...	8.110	40.6	ug/l	814379	2	6.763	1003340	50.0
Benzene, 2-prop...	12.244	429.9	ug/l	9647900	4	12.024	1122120	50.0
o-Cymene	12.616	48.1	ug/l	1079330	4	12.024	1122120	50.0
Benzene, 1-ethe...	12.701	166.9	ug/l	3745360	4	12.024	1122120	50.0
Benzene, 1,2,3,...	12.920	74.9	ug/l	1681550	4	12.024	1122120	50.0
1H-Indene, 2,3-...	13.170	87.0	ug/l	1951450	4	12.024	1122120	50.0
1H-Indene, 2,3-...	13.292	215.1	ug/l	4826500	4	12.024	1122120	50.0
1H-Indene, 2,3-...	13.542	45.5	ug/l	1021170	4	12.024	1122120	50.0
2,2-Dimethylind...	13.664	42.5	ug/l	954181	4	12.024	1122120	50.0