LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-24

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: VX029771.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.264	23	29	40	rBV2	18263	28032	2.23%	0.226%
2	1.746	104	108	117	rVB	49614	68640	5.45%	0.553%
3	2.337	196	205	219	rVB	30169	64769	5.14%	0.522%
4	2.782	270	278	283	rBV	26797	58203	4.62%	0.469%
5	2.965	301	308	318	rVB	10170	23146	1.84%	0.186%
6	3.105	321	331	342	rBV3	33253	79596	6.32%	0.641%
7	3.757	429	438	451	rBV2	15403	42547	3.38%	0.343%
8	4.208	500	512	522	rBV3	6029	18588	1.48%	0.150%
9	5.391	695	706	719	rBV2	136719	401934	31.91%	3.237%
10	5.556	722	733	742	rBV	264394	753725	59.84%	6.071%
11	5.964	789	800	807	rVV	143498	383897	30.48%	3.092%
12	6.044	807	813	823	rVV	39645	108065	8.58%	0.870%
13	6.159	824	832	839	rVV4	17650	52957	4.20%	0.427%
14	6.251	839	847	858	rVV6	39368	125874	9.99%	1.014%
15	6.361	858	865	873	rVB8	8650	24845	1.97%	0.200%
16	6.763	921	931	945	rBV	397216	940843	74.70%	7.578%
17	7.366	1021	1030	1038	rBV2	12924	31341	2.49%	0.252%
18	7.775	1091	1097	1105	rVB2	11574	22260	1.77%	0.179%
19	7.988	1122	1132	1141	rVB	37297	86010	6.83%	0.693%
20	8.110	1143	1152	1160	rBV	81381	167379	13.29%	1.348%
21	8.311	1179	1185	1193	rBV6	8036	17457	1.39%	0.141%
22	8.519	1212	1219	1225	rVB2	11393	26712	2.12%	0.215%
23	8.604	1225	1233	1235	rBV2	23711	41035	3.26%	0.331%
24	8.653	1235	1241	1254	rVB	791725	1259480	100.00%	10.144%
25	8.775	1254	1261	1265	rBV3	11813	19494	1.55%	0.157%
26	8.927	1282	1286	1289	rVV2	9131	12986	1.03%	0.105%
27	8.976	1289	1294	1299	rVB2	15846	28216	2.24%	0.227%
28	9.104	1307	1315	1320	rVV3	35755	64522	5.12%	0.520%
29	9.165	1320	1325	1336	rVB	70912	119665	9.50%	0.964%
30	9.433	1364	1369	1372	rBV2	11099	18882	1.50%	0.152%
31	9.506	1377	1381	1387	rVB4	23322	45063	3.58%	0.363%
32	9.659	1403	1406	1416	rVB2	13910	24190	1.92%	0.195%
33	9.762	1416	1423	1428	rBV2	24997	41923	3.33%	0.338%
34	10.055	1466	1471	1477	rBV	771941	1031737	81.92%	8.310%
35	10.128	1477	1483	1487	rVB2	13259	20928	1.66%	0.169%
36	10.274	1503	1507	1509	rBV2	14447	18944	1.50%	0.153%

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

Instrument :
MSVOA_X
ClientSampleId :
MW-24

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	10.305	1509	1512	1518	rVB4	16780	26846	2.13%	0.216%
38	10.378	1518	1524	1528	rBV8	5906	14374	1.14%	0.116%
39	10.585	1550	1558	1565	rBV4	17452	34251	2.72%	0.276%
40	10.781	1582	1590	1593	rBV3	11703	22461	1.78%	0.181%
41	10.860	1598	1603	1611	rVB8	6712	16733	1.33%	0.135%
42	10.969	1616	1621	1626	rBV4	10722	22699	1.80%	0.183%
43	11.085	1635	1640	1645	rBV	605015	778182	61.79%	6.268%
44	11.128	1645	1647	1655	rVB4	15743	20117	1.60%	0.162%
45	11.311	1674	1677	1685	rVB5	8936	16213	1.29%	0.131%
46	11.616	1720	1727	1736	rBV	60698	82221	6.53%	0.662%
47	11.713	1736	1743	1747	rBV4	13377	23105	1.83%	0.186%
48	11.890	1768	1772	1778	rVB3	14324	20393	1.62%	0.164%
49	12.024	1789	1794	1799	rBV	815559	998083	79.25%	8.039%
50	12.177	1815	1819	1825	rVB	31889	42393	3.37%	0.341%
51	12.250	1825	1831	1837	rBV	602328	735098	58.37%	5.921%
52	12.317	1837	1842	1848	rVB	135323	161264	12.80%	1.299%
53	12.408	1852	1857	1862	rBV2	64160	83654	6.64%	0.674%
54	12.475	1862	1868	1872	rVB2	35589	49574	3.94%	0.399%
55	12.646	1891	1896	1900	rBV	193118	224574	17.83%	1.809%
56	12.701	1900	1905	1912	rVV	945026	1167898	92.73%	9.407%
57	12.762	1912	1915	1921	rVB2	17348	21309	1.69%	0.172%
58	12.841	1922	1928	1933	rVB	44058	57354	4.55%	0.462%
59	12.902	1933	1938	1941	rBV3	13388	17750	1.41%	0.143%
60	13.024	1953	1958	1966	rVB2	76656	122365	9.72%	0.986%
61	13.140	1968	1977	1979	rBV2	63467	121137	9.62%	0.976%
62	13.170	1979	1982	1985	rBV	87278	91952	7.30%	0.741%
63	13.237	1990	1993	1997	rVB2	20509	24455	1.94%	0.197%
64	13.298	1997	2003	2007	rBV2	75510	103372	8.21%	0.833%
65	13.347	2007	2011	2015	rBV3	59995	81114	6.44%	0.653%
66	13.390	2015	2018	2020	rBV2	12169	13354	1.06%	0.108%
67	13.426	2020	2024	2029	rVB	59687	80409	6.38%	0.648%
68	13.506	2032	2037	2039	rBV2	15046	24641	1.96%	0.198%
69	13.542	2039	2043	2047	rBV	64134	82455	6.55%	0.664%
70	13.640	2052	2059	2068	rVB3	54663	129054	10.25%	1.039%
71	13.804	2080	2086	2091	rVV4	9278	27177	2.16%	0.219%
72	13.853	2091	2094	2099	rVB3	21641	29196	2.32%	0.235%
73	13.926	2099	2106	2110	rBV	61350	88010	6.99%	0.709%
74	13.969	2110	2113	2118	rVB2	13017	17819	1.41%	0.144%
75	14.079	2126	2131	2135	rBV	62742	84269	6.69%	0.679%
76	14.219	2148	2154	2158	rBV3	22002	41013	3.26%	0.330%

LSC Area Percent Report

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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

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
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Title : SW846 8260

77	14.323	2166	2171	2176	rBV	48058	71824	5.70%	0.578%
78	14.371	2176	2179	2185	rVB	52743	68874	5.47%	0.555%
79	14.487	2193	2198	2202	rBV2	11597	20778	1.65%	0.167%
80	14.603	2210	2217	2222	rBV4	28507	63741	5.06%	0.513%
81	14.676	2227	2229	2233	rVB2	13972	15120	1.20%	0.122%
82	14.792	2243	2248	2252	rBV3	15825	22521	1.79%	0.181%
83	14.908	2261	2267	2273	rBV10	7003	15221	1.21%	0.123%
84	15.182	2306	2312	2320	rBV	15036	30098	2.39%	0.242%
85	15.408	2346	2349	2354	rVB	13189	18077	1.44%	0.146%
86	15.987	2439	2444	2454	rVB4	9022	17302	1.37%	0.139%

Sum of corrected areas: 12415779

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MSVOA_X
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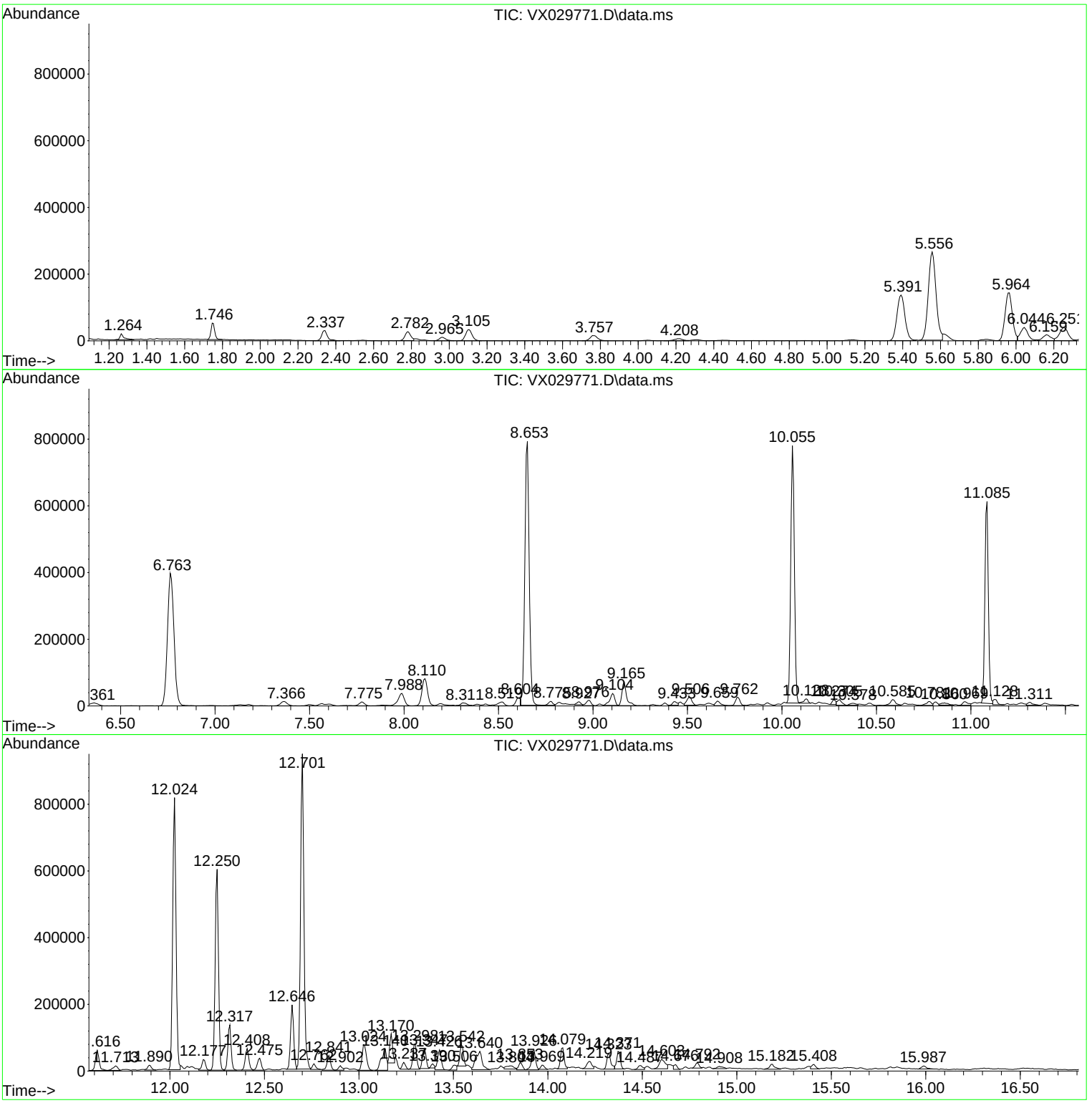
MW-24

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



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Instrument :
MSVOA_X
ClientSampleId :
MW-24

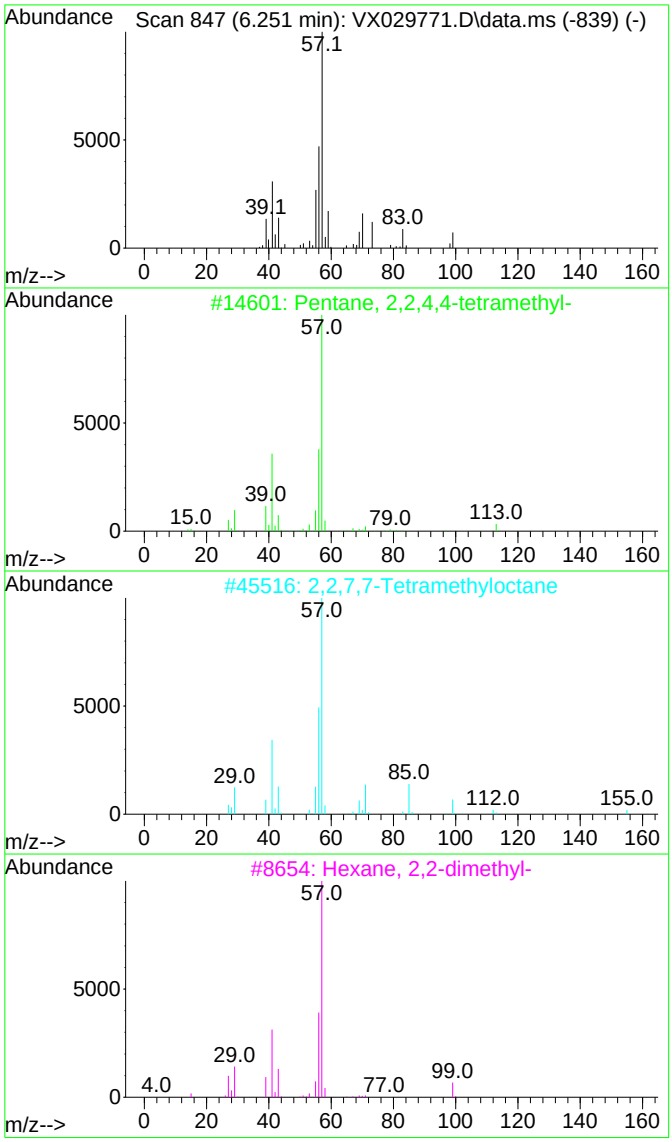
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

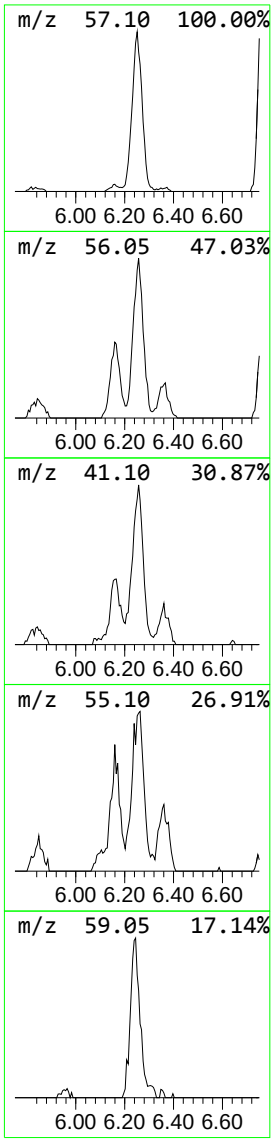
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	6.69 ug/l	125874	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	43
2	2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	40
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	38
4	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	36
5	Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

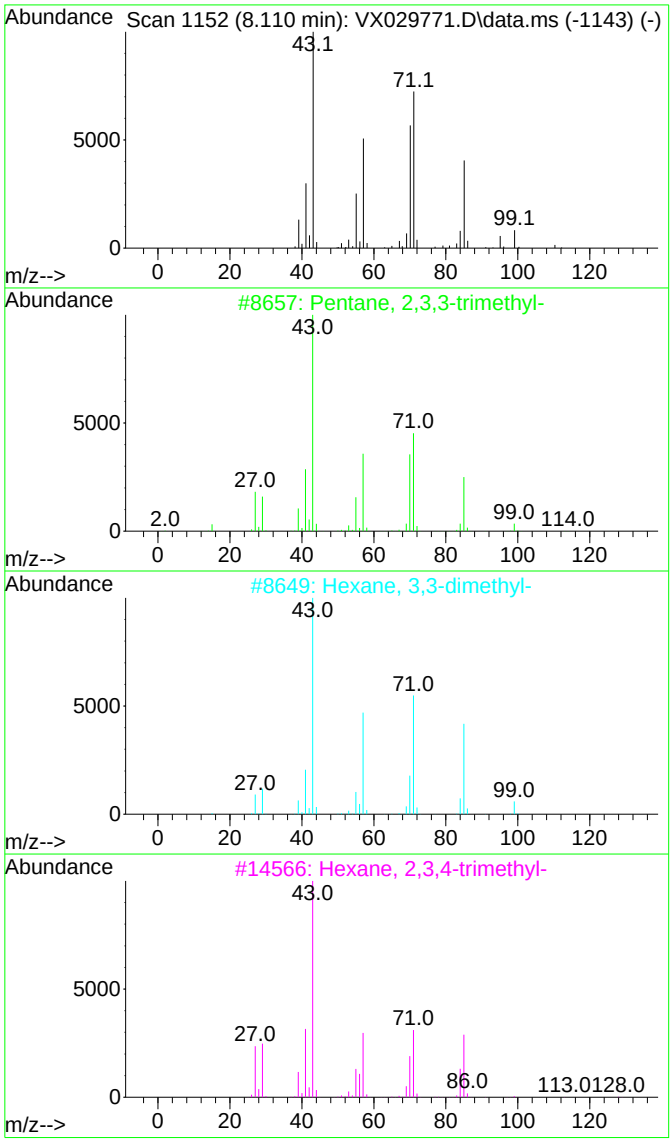
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

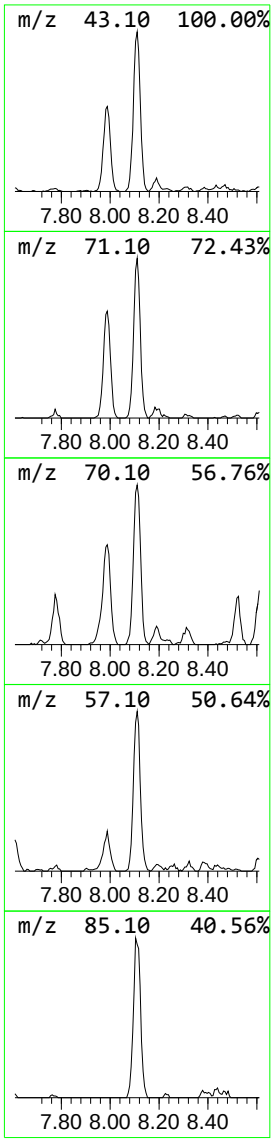
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	8.90 ug/l	167379	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, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	50





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ALS Vial : 13 Sample Multiplier: 1

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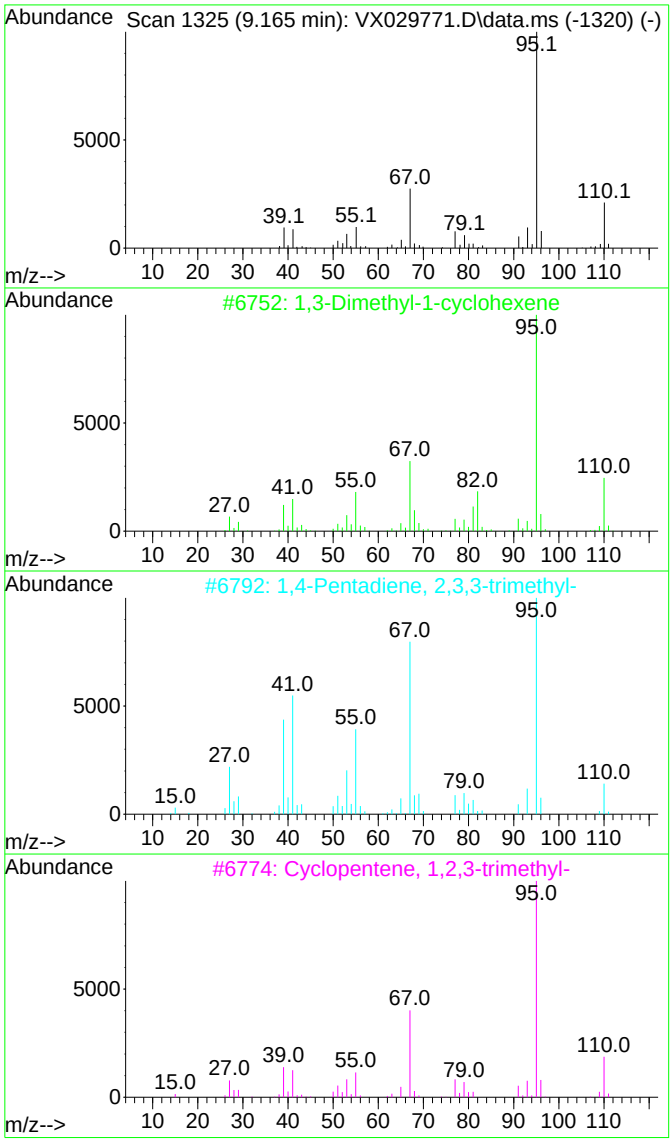
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

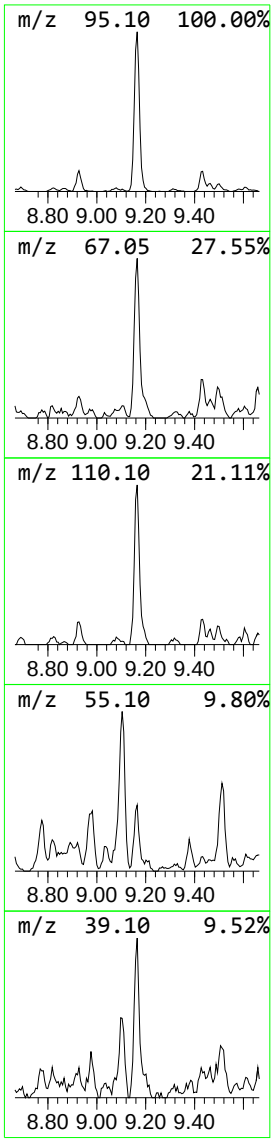
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,3-Dimethyl-1-cyclohexene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.165	5.80 ug/l	119665	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	91
2	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	91
3	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	91
4	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1010142-17-5	80
5	5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	78





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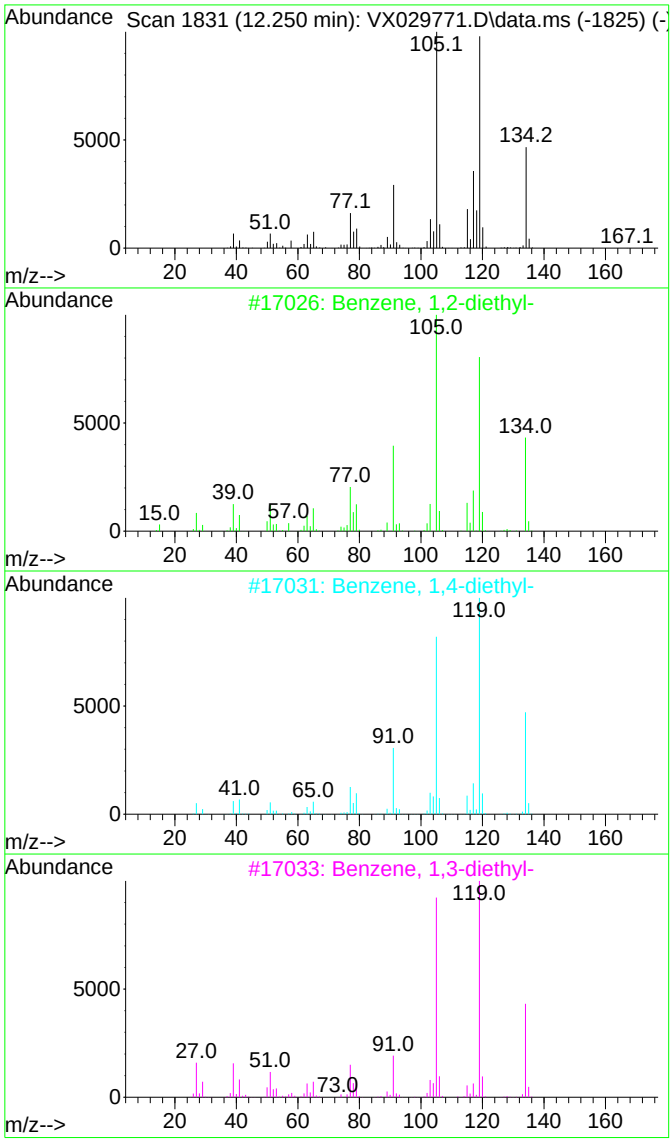
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

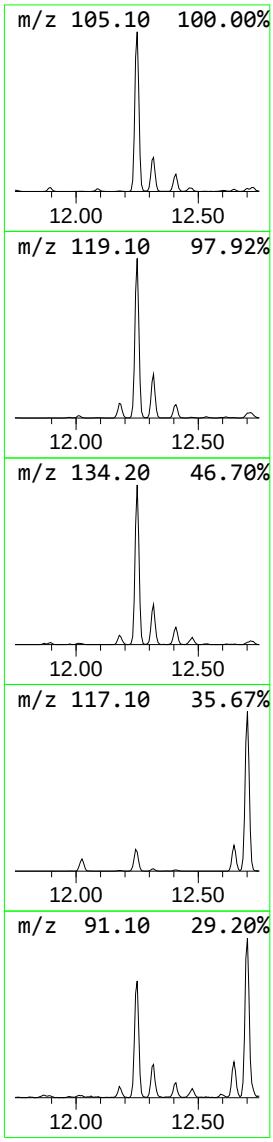
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	36.83 ug/l	735098	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	92
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
4	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	86
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83





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

Instrument :
MSVOA_X
ClientSampleId :
MW-24

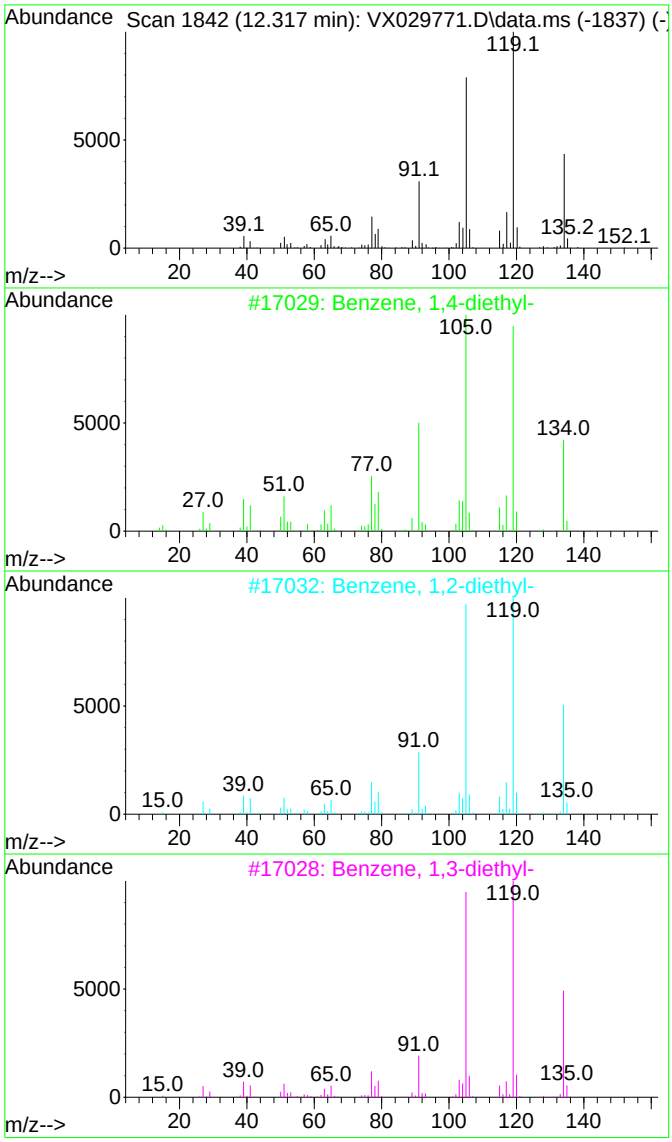
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

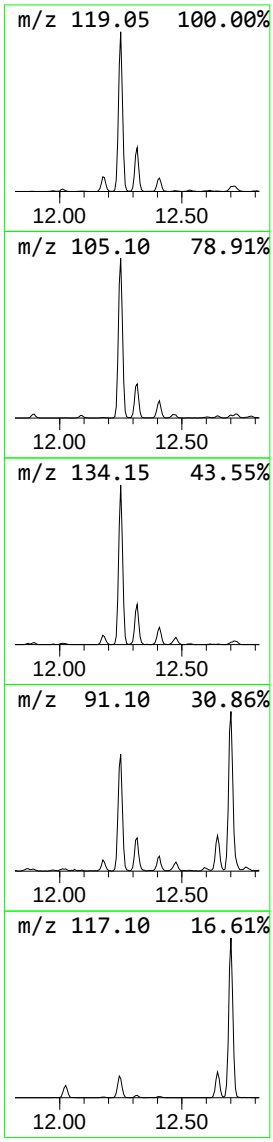
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,4-diethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	8.08 ug/l	161264	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5	2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	87





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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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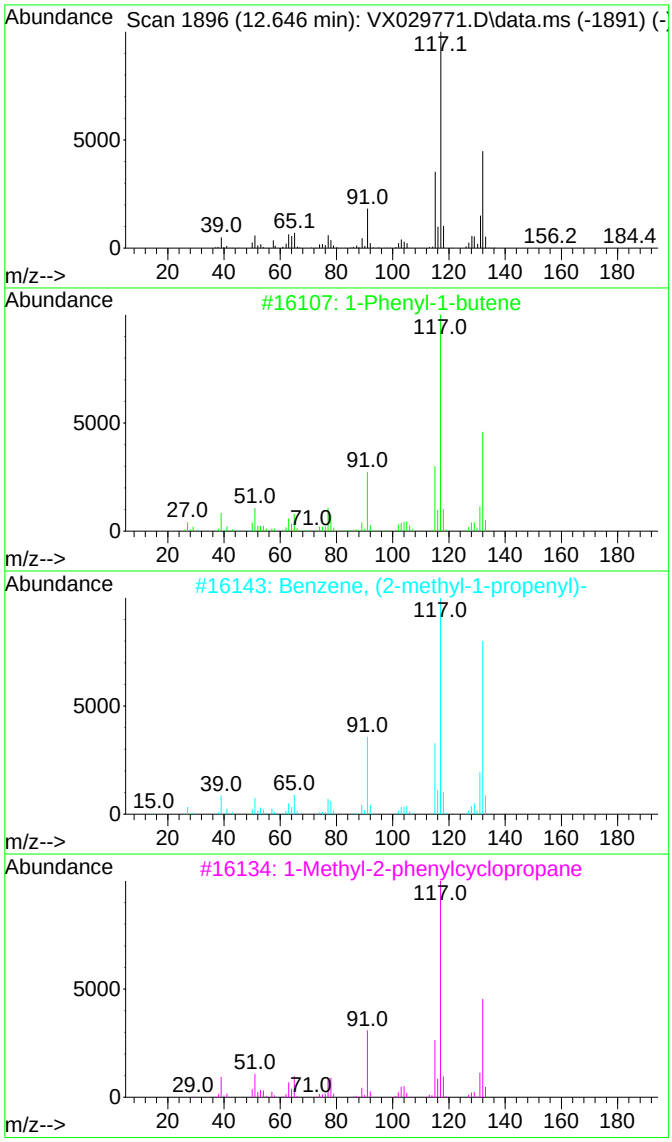
Quant Title : SW846 8260

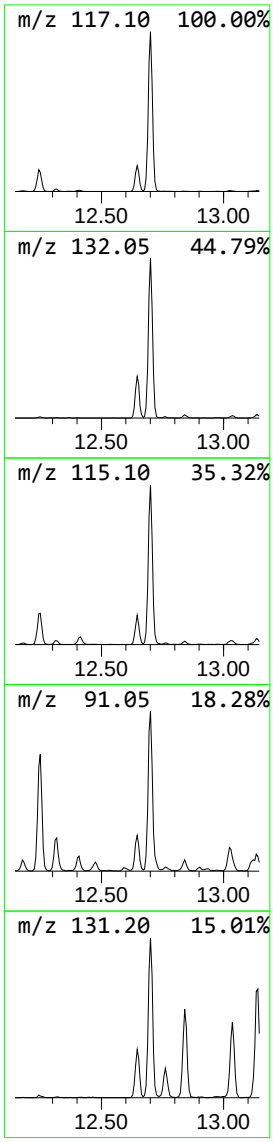
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	11.25 ug/l	224574	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	94
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	91
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

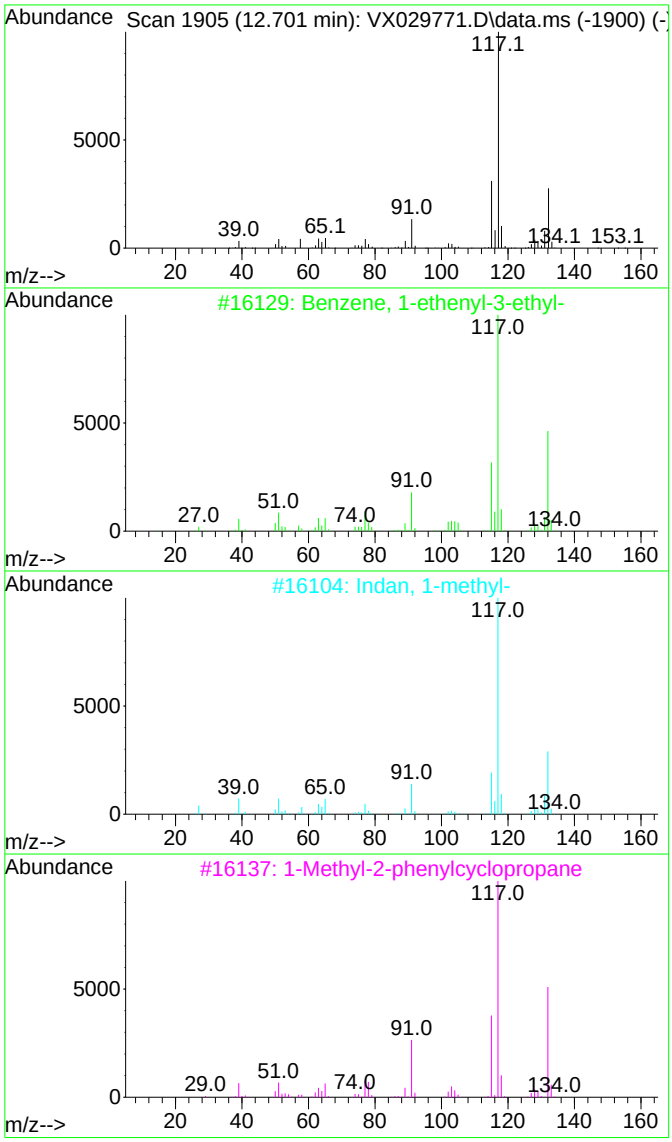
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

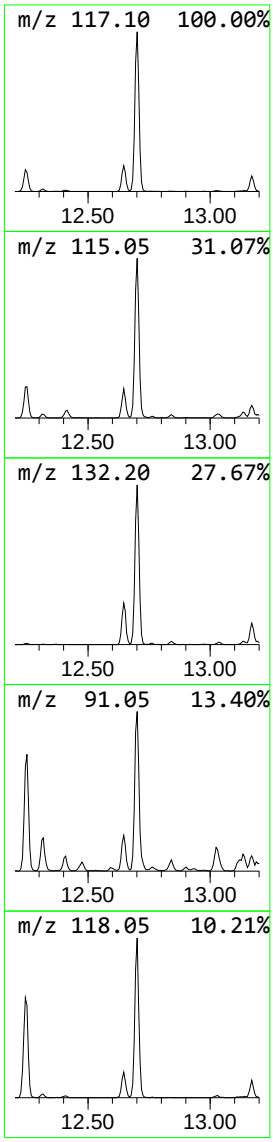
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	58.51 ug/l	1167900	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	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

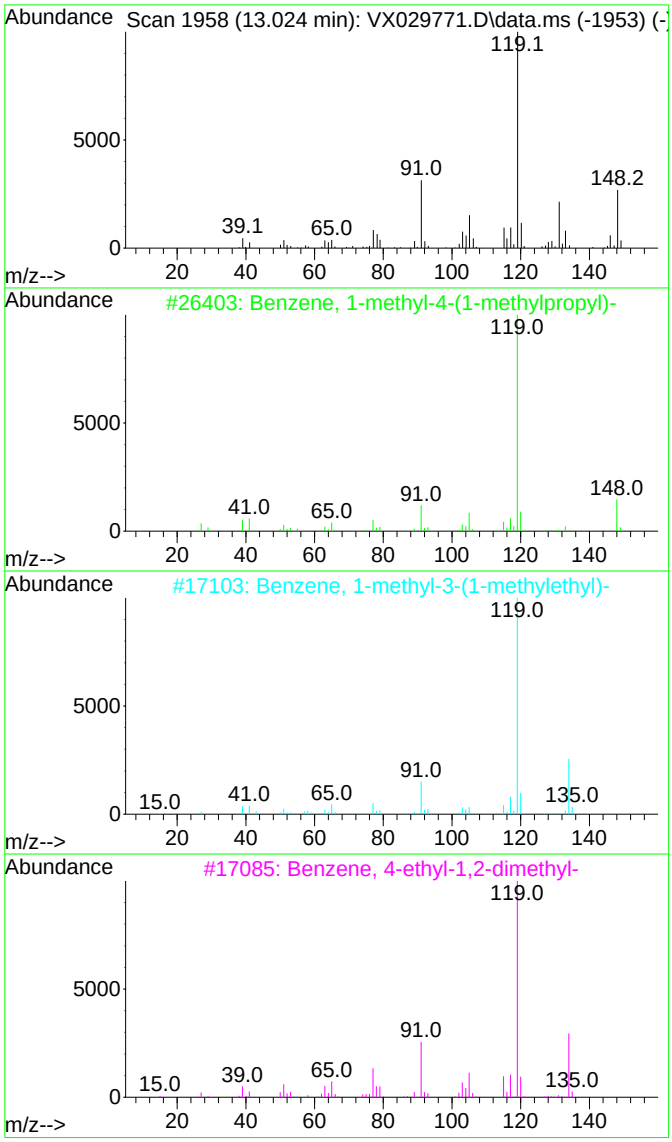
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

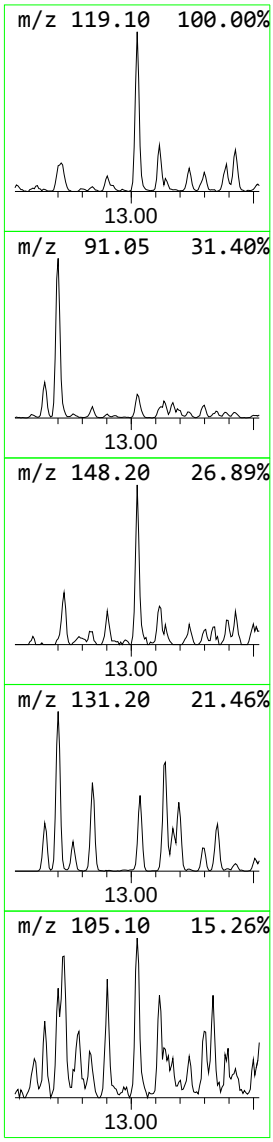
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	6.13 ug/l	122365	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	81
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	64
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	62





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

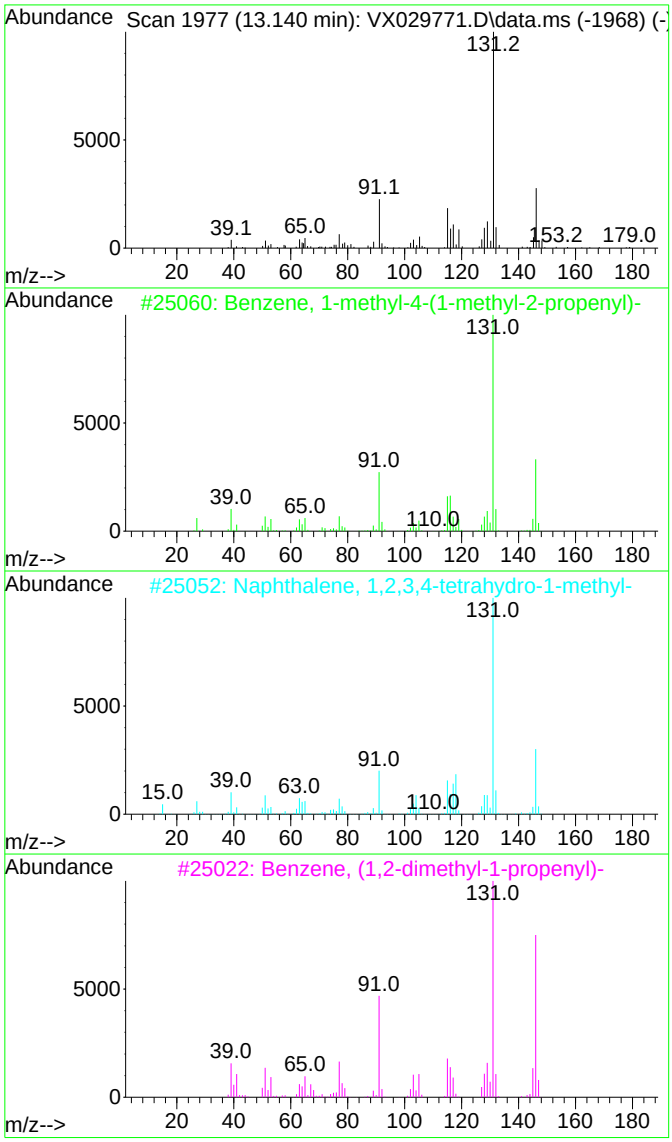
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

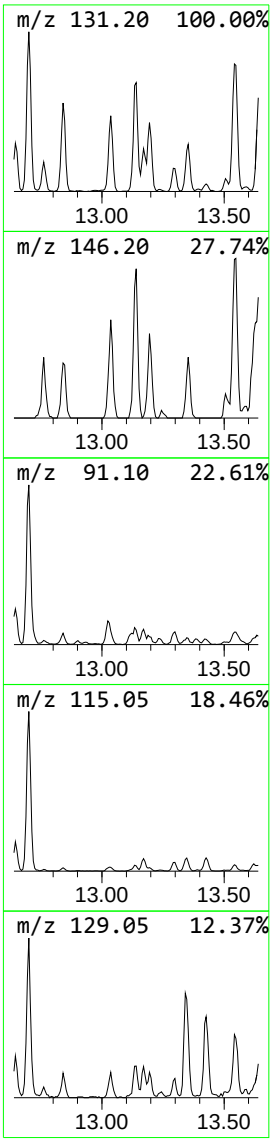
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-methyl-4-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.140	6.07 ug/l	121137	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

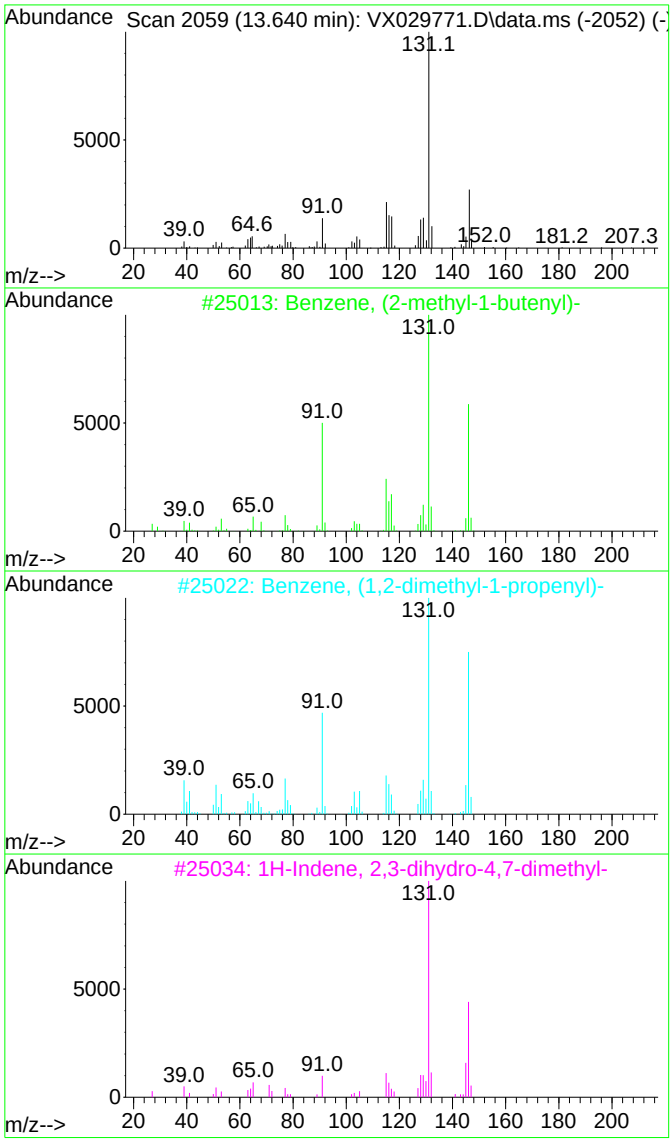
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

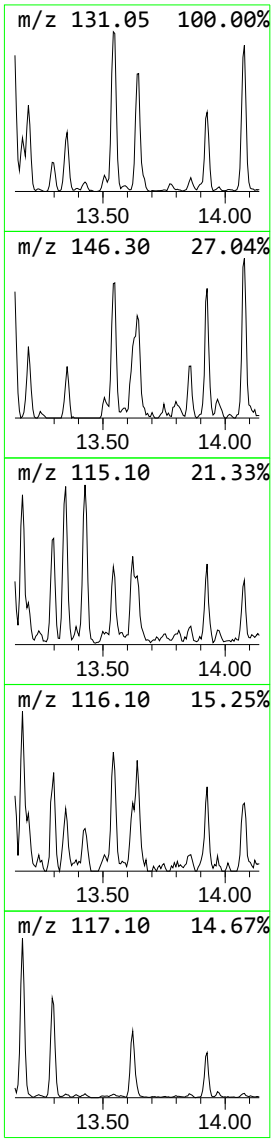
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	6.47 ug/l	129054	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	96
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
Data File : VX029771.D
Acq On : 24 Jun 2022 13:36
Operator : JC/MD
Sample : N3366-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

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	6.7	ug/l	125874	2	6.763	940843	50.0
Pentane, 2,3,3-...	8.110	8.9	ug/l	167379	2	6.763	940843	50.0
1,3-Dimethyl-1-...	9.165	5.8	ug/l	119665	3	10.055	1031740	50.0
Benzene, 1,2-di...	12.250	36.8	ug/l	735098	4	12.024	998083	50.0
Benzene, 1,4-di...	12.317	8.1	ug/l	161264	4	12.024	998083	50.0
1-Phenyl-1-butene	12.646	11.3	ug/l	224574	4	12.024	998083	50.0
Benzene, 1-ethe...	12.701	58.5	ug/l	1167900	4	12.024	998083	50.0
Benzene, 1-meth...	13.024	6.1	ug/l	122365	4	12.024	998083	50.0
Benzene, 1-meth...	13.140	6.1	ug/l	121137	4	12.024	998083	50.0
Benzene, (2-met...	13.640	6.5	ug/l	129054	4	12.024	998083	50.0