LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

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\82X020822W.M
Title : SW846 8260

Signal : TIC: VX027078.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.751	105	109	115	rVB	4608	6458	1.14%	0.126%
2	2.343	199	206	219	rVB2	12181	26403	4.65%	0.513%
3	2.532	231	237	248	rBV	4492	8659	1.53%	0.168%
4	2.788	271	279	282	rBV	12312	25976	4.58%	0.505%
5	2.830	282	286	302	rVB	11512	26135	4.60%	0.508%
6	3.105	323	331	340	rBV	15784	35810	6.31%	0.696%
7	4.220	506	514	524	rVB3	3097	9035	1.59%	0.176%
8	5.141	655	665	678	rBV6	2143	8990	1.58%	0.175%
9	5.391	696	706	720	rBV2	65160	191101	33.67%	3.716%
10	5.556	722	733	757	rVB2	113990	353219	62.22%	6.868%
11	5.842	771	780	789	rVB5	2424	7343	1.29%	0.143%
12	5.964	789	800	809	rVV	72782	191089	33.66%	3.716%
13	6.049	809	814	821	rVV3	4446	11651	2.05%	0.227%
14	6.171	821	834	842	rVV6	8272	33583	5.92%	0.653%
15	6.257	842	848	858	rVV2	9080	27369	4.82%	0.532%
16	6.372	858	867	874	rVB3	3344	8931	1.57%	0.174%
17	6.769	919	932	943	rBV	181758	442529	77.96%	8.605%
18	6.860	943	947	957	rVB2	6676	14758	2.60%	0.287%
19	7.177	991	999	1008	rVB3	3589	7534	1.33%	0.146%
20	7.366	1022	1030	1037	rBV4	3748	8485	1.49%	0.165%
21	7.781	1092	1098	1104	rBV3	3343	6673	1.18%	0.130%
22	7.884	1113	1115	1123	rVV7	2937	6425	1.13%	0.125%
23	7.988	1123	1132	1142	rVB4	5655	15526	2.74%	0.302%
24	8.110	1145	1152	1163	rBV3	11830	27342	4.82%	0.532%
25	8.506	1209	1217	1224	rVB5	5411	12354	2.18%	0.240%
26	8.653	1235	1241	1250	rVB	363828	567652	100.00%	11.038%
27	8.842	1265	1272	1279	rBV4	4667	9329	1.64%	0.181%
28	8.957	1282	1291	1300	rVV	142002	231476	40.78%	4.501%
29	9.049	1300	1306	1312	rVV	165415	250827	44.19%	4.877%
30	9.104	1312	1315	1320	rVV2	5416	9384	1.65%	0.182%
31	9.165	1320	1325	1335	rVB	15238	24454	4.31%	0.476%
32	9.433	1364	1369	1371	rBV	5081	7889	1.39%	0.153%
33	9.457	1371	1373	1377	rVV	7311	10657	1.88%	0.207%
34	9.500	1377	1380	1387	rVV4	3403	7267	1.28%	0.141%
35	9.762	1419	1423	1427	rVB4	3611	6418	1.13%	0.125%
36	10.055	1462	1471	1481	rBV	365151	498335	87.79%	9.690%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-11R

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\82X020822W.M
Title : SW846 8260

37	10.146	1481	1486	1491	rVB4	6105	11105	1.96%	0.216%
38	10.250	1497	1503	1506	rBV4	3672	6072	1.07%	0.118%
39	10.414	1525	1530	1536	rVB5	7033	13061	2.30%	0.254%
40	10.963	1617	1620	1624	rVB	4954	5953	1.05%	0.116%
41	11.085	1634	1640	1645	rBV	281329	353982	62.36%	6.883%
42	11.615	1724	1727	1732	rVB	9052	11136	1.96%	0.217%
43	11.890	1769	1772	1777	rVB	8952	11987	2.11%	0.233%
44	12.024	1789	1794	1800	rBV	359511	430045	75.76%	8.362%
45	12.176	1816	1819	1823	rVB	7148	8711	1.53%	0.169%
46	12.249	1823	1831	1837	rVV2	198835	262551	46.25%	5.105%
47	12.317	1837	1842	1848	rVV	23878	27767	4.89%	0.540%
48	12.408	1852	1857	1864	rVB2	14525	19460	3.43%	0.378%
49	12.530	1873	1877	1882	rVB	8565	10071	1.77%	0.196%
50	12.615	1886	1891	1894	rBV	74969	91146	16.06%	1.772%
51	12.646	1894	1896	1900	rVV	22076	24009	4.23%	0.467%
52	12.701	1900	1905	1912	rVB	207669	246672	43.45%	4.797%
53	12.841	1922	1928	1932	rVB	6963	8378	1.48%	0.163%
54	13.030	1954	1959	1964	rBV2	8970	13379	2.36%	0.260%
55	13.140	1973	1977	1979	rBV	8952	11830	2.08%	0.230%
56	13.170	1979	1982	1992	rVB	76572	96208	16.95%	1.871%
57	13.261	1992	1997	1999	rBV2	4689	7109	1.25%	0.138%
58	13.292	1999	2002	2008	rVV	83660	110150	19.40%	2.142%
59	13.347	2008	2011	2013	rVV3	12912	16469	2.90%	0.320%
60	13.371	2013	2015	2021	rVV	8546	9838	1.73%	0.191%
61	13.426	2021	2024	2029	rVB	5546	6621	1.17%	0.129%
62	13.511	2031	2038	2040	rBV3	7308	10675	1.88%	0.208%
63	13.548	2040	2044	2047	rVV	36809	50746	8.94%	0.987%
64	13.591	2047	2051	2054	rVV4	12973	21774	3.84%	0.423%
65	13.646	2054	2060	2061	rVV2	17733	30592	5.39%	0.595%
66	13.664	2061	2063	2072	rVB2	34197	45156	7.95%	0.878%
67	13.871	2091	2097	2102	rBV2	6825	10590	1.87%	0.206%
68	13.926	2102	2106	2110	rVV	5617	6663	1.17%	0.130%
69	14.078	2126	2131	2136	rBV	7933	9796	1.73%	0.190%
70	14.377	2176	2180	2189	rVB3	4402	5895	1.04%	0.115%

Sum of corrected areas: 5142663

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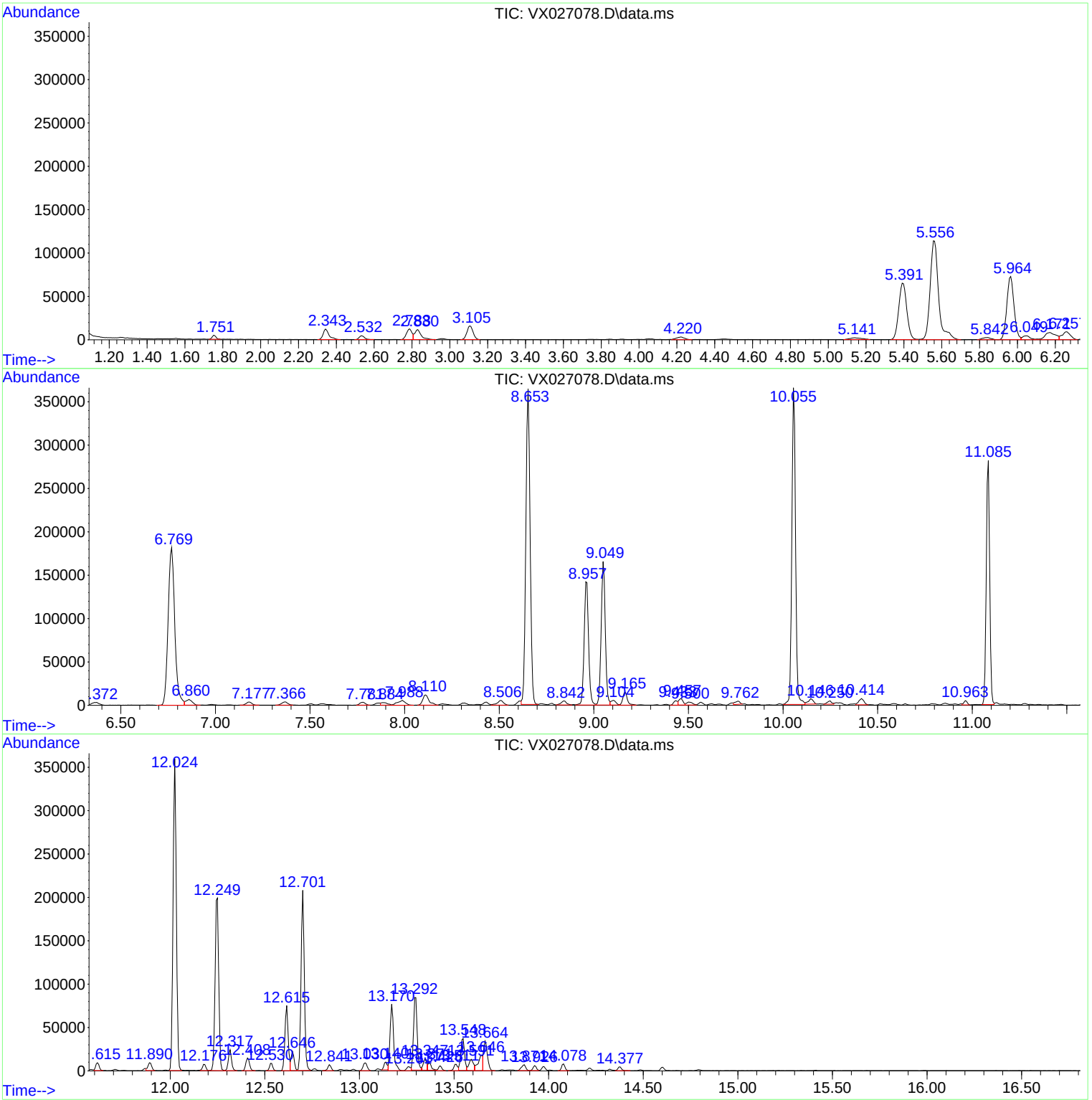
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
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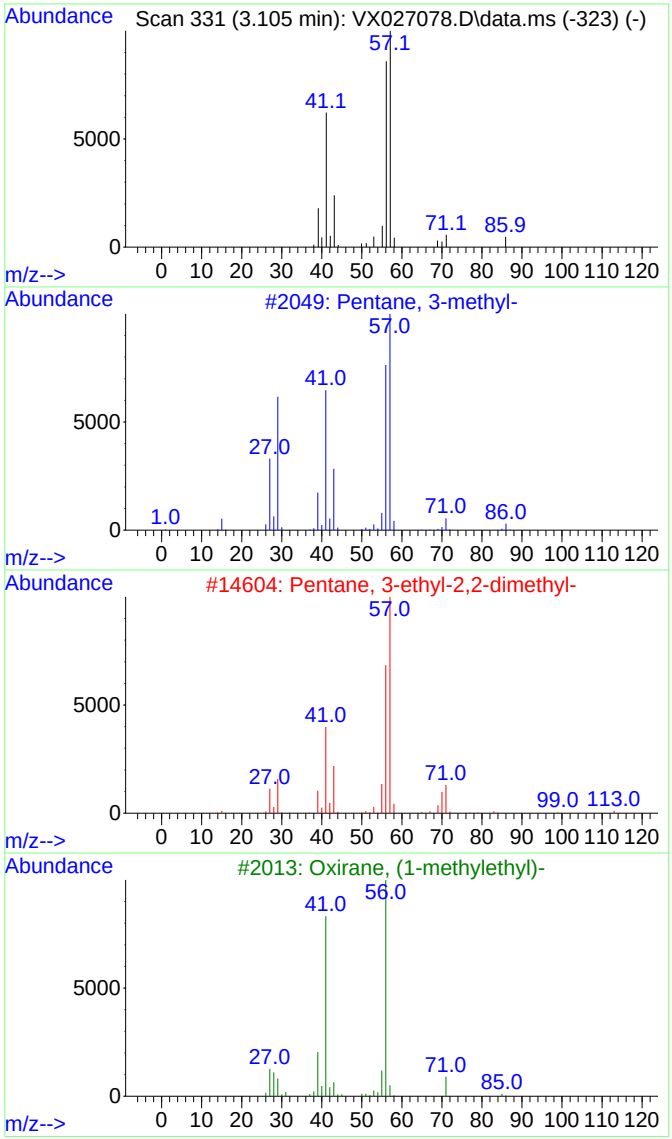
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

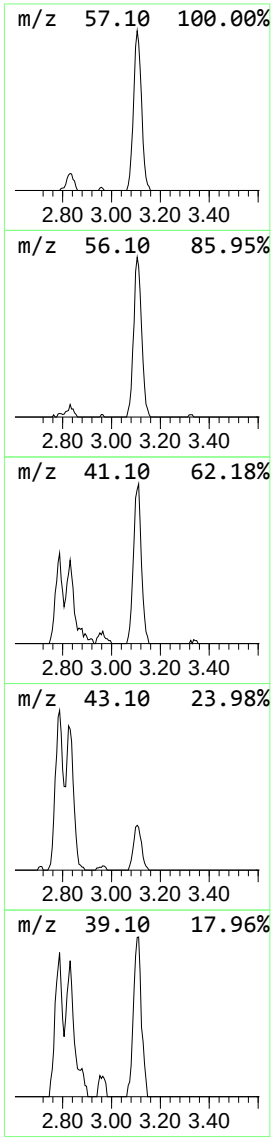
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Pentane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	5.07 ug/l	35810	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	74
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

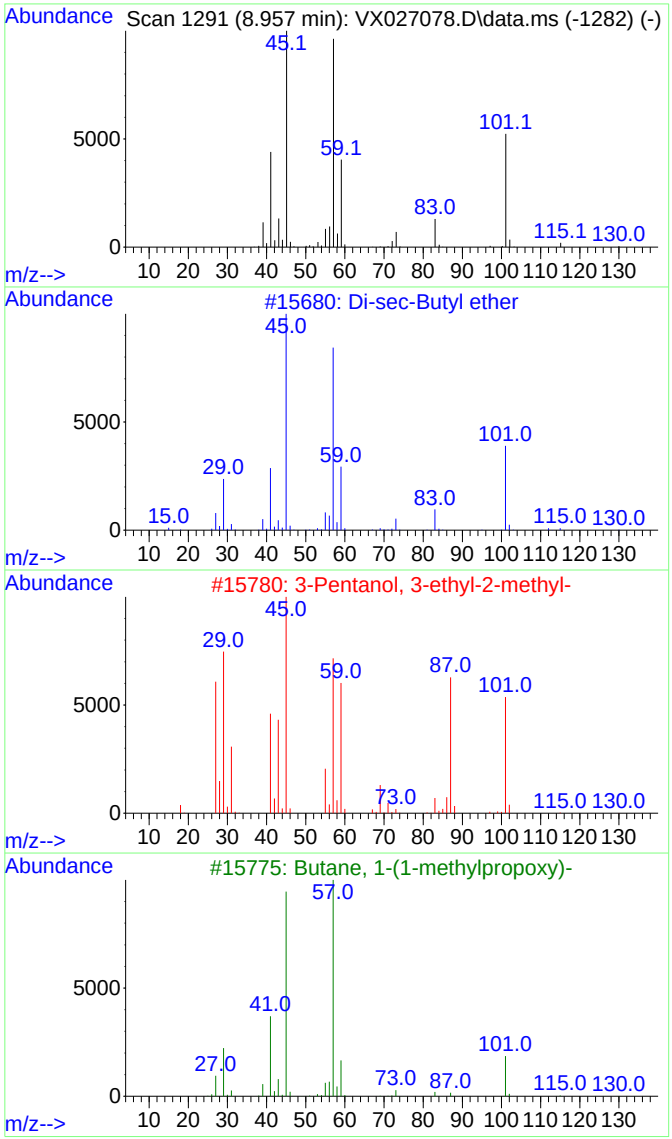
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

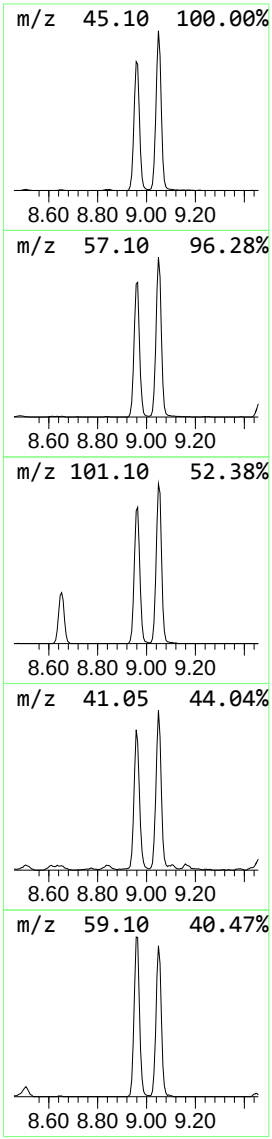
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Di-sec-Butyl ether Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.957	23.22 ug/l	231476	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-Butyl ether	130	C8H18O	006863-58-7	91
2	3-Pentanol, 3-ethyl-2-methyl-	130	C8H18O	000597-05-7	64
3	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50
4	Diethyl carbitol	162	C8H18O3	000112-36-7	47
5	Sulfurous acid, isobutyl 2-methy...	238	C10H22O4S	1000309-20-3	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

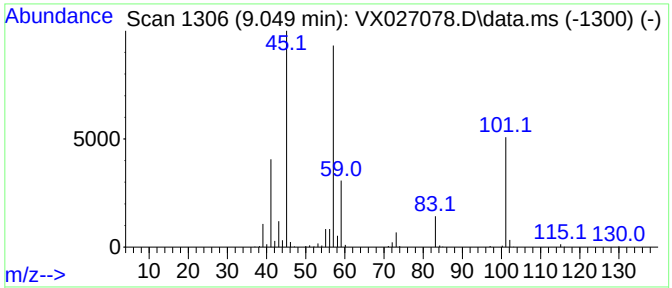
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

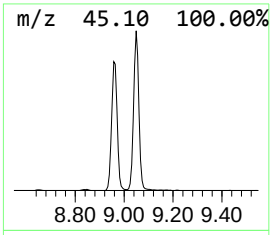
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

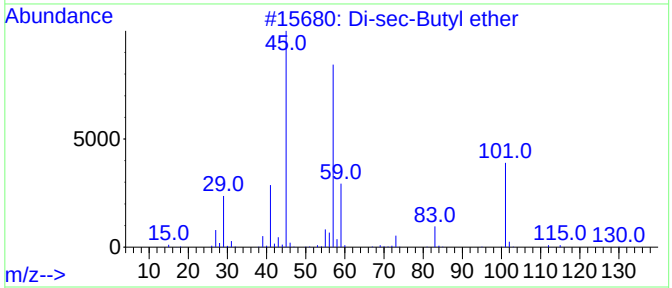
****
Peak Number 3 3-Pentanol, 3-ethyl-2-methyl- Concentration Rank 3

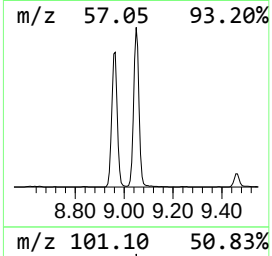
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.049	25.17 ug/l	250827	Chlorobenzene-d5	10.055

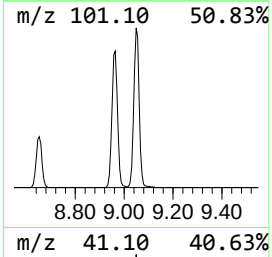
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-Butyl ether	130	C8H18O	006863-58-7	91
2	3-Pentanol, 3-ethyl-2-methyl-	130	C8H18O	000597-05-7	72
3	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	53
4	Diethyl carbitol	162	C8H18O3	000112-36-7	47
5	1-Octene, 3-(methoxymethoxy)-	172	C10H20O2	088738-34-5	40

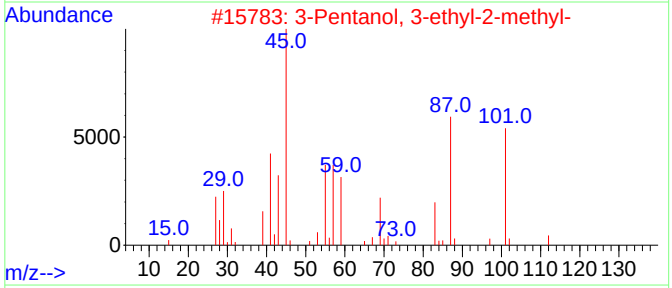


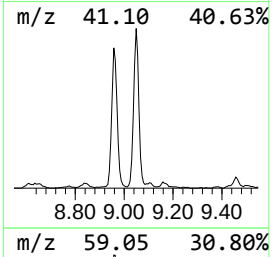


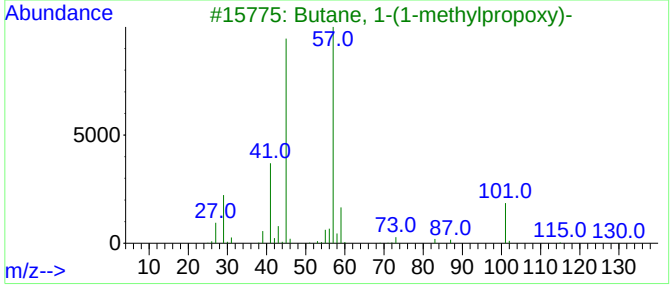


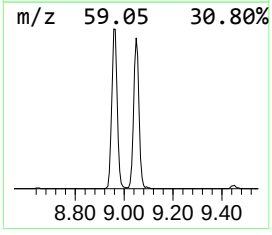












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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

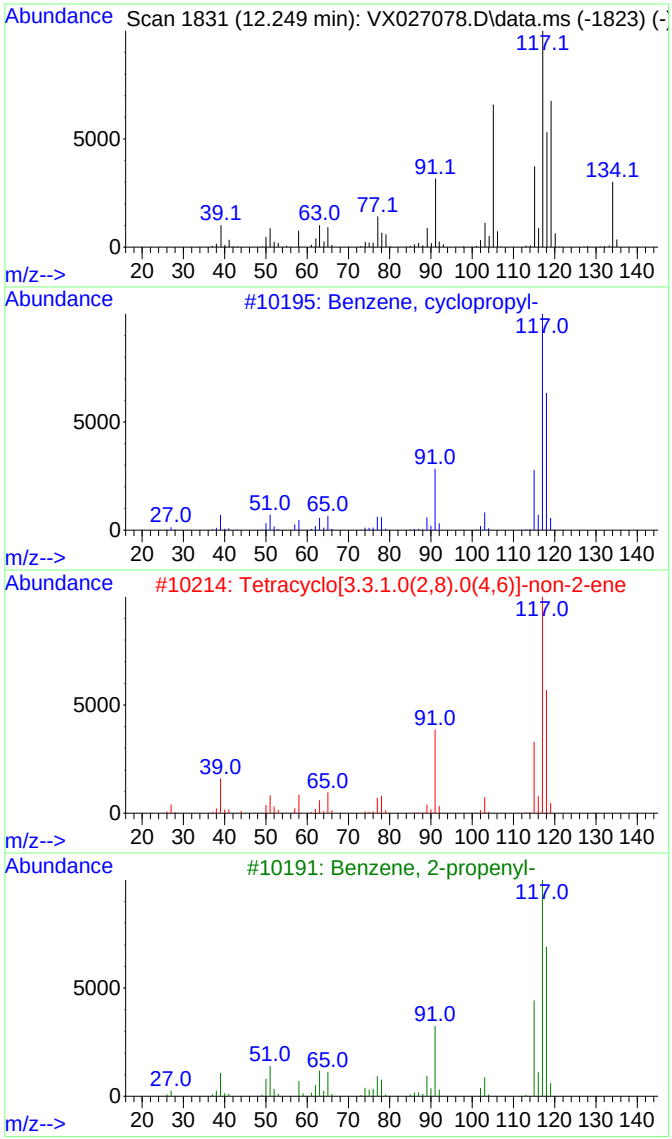
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

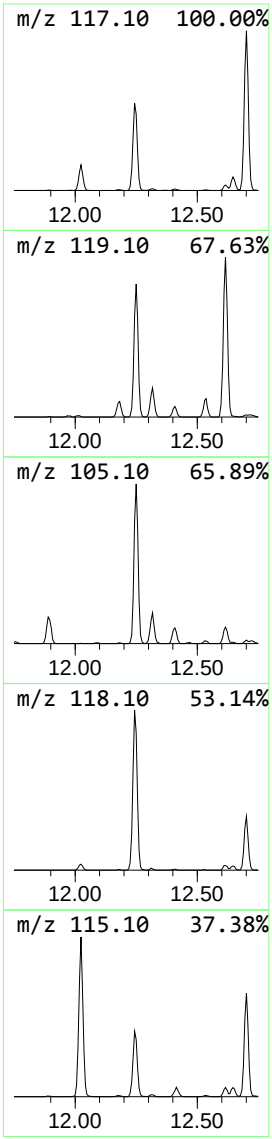
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.249	30.53 ug/l	262551	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	60
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
5	Deltacyclene	118	C9H10	007785-10-6	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

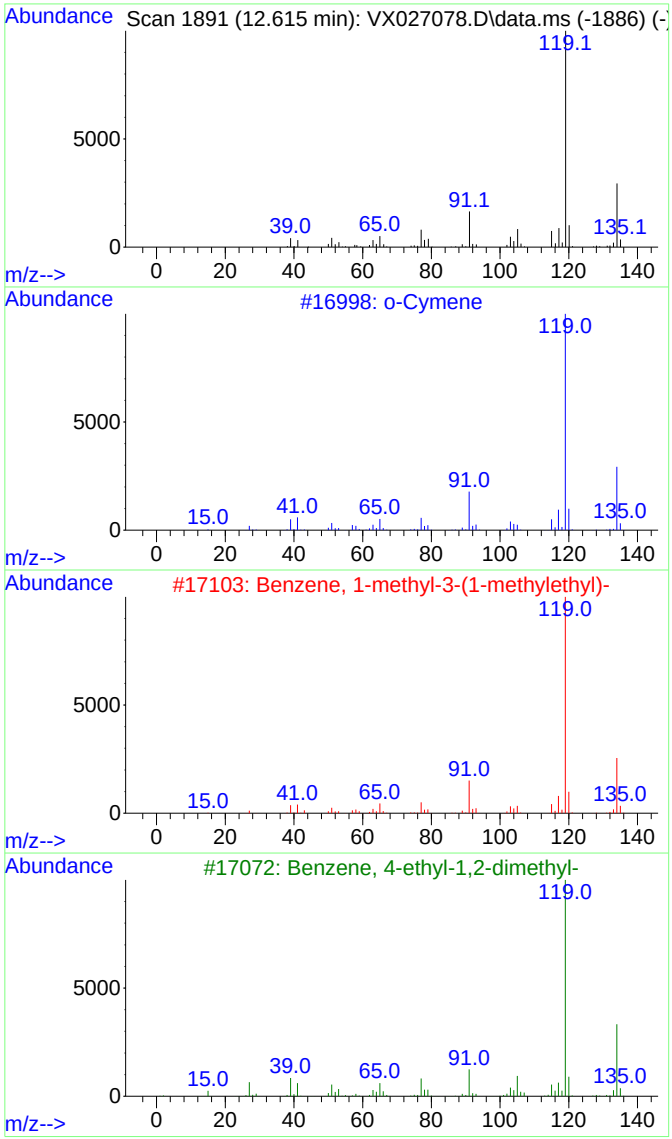
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

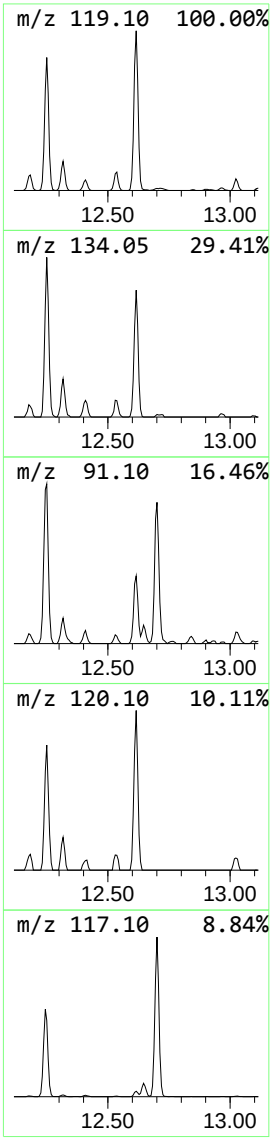
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.615	10.60 ug/l	91146	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	p-Cymene	134	C10H14	000099-87-6	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
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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-11R

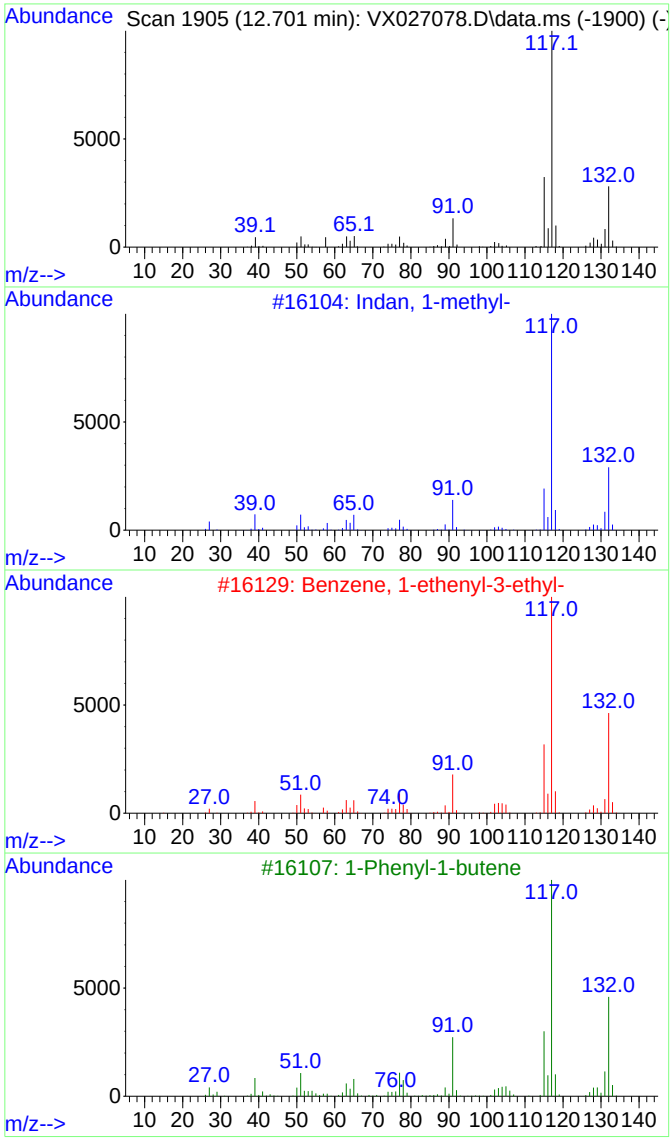
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

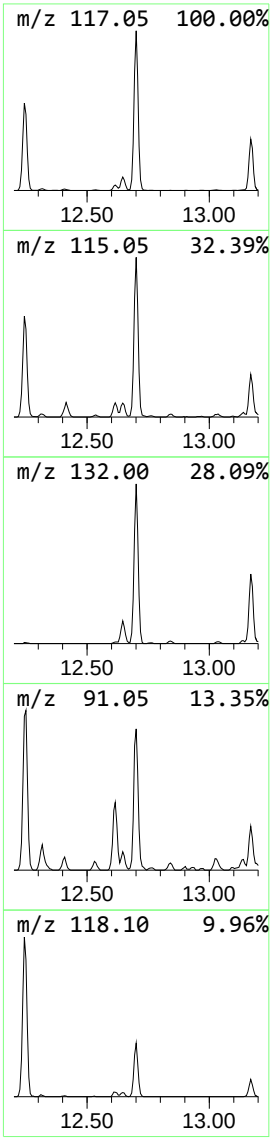
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	28.68 ug/l	246672	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

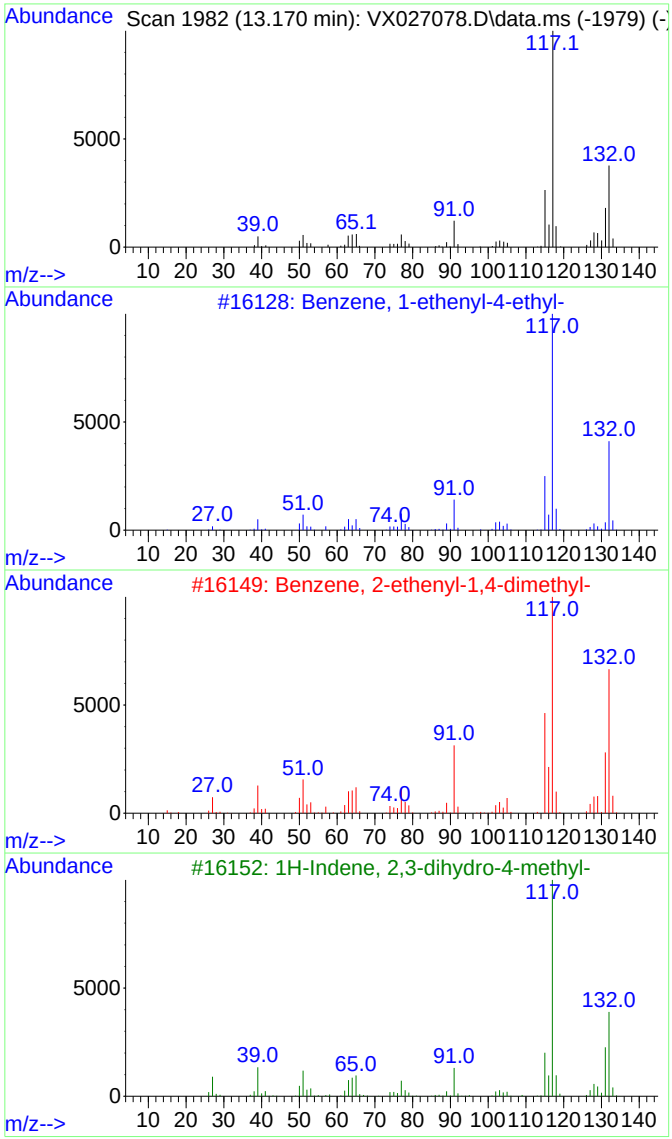
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

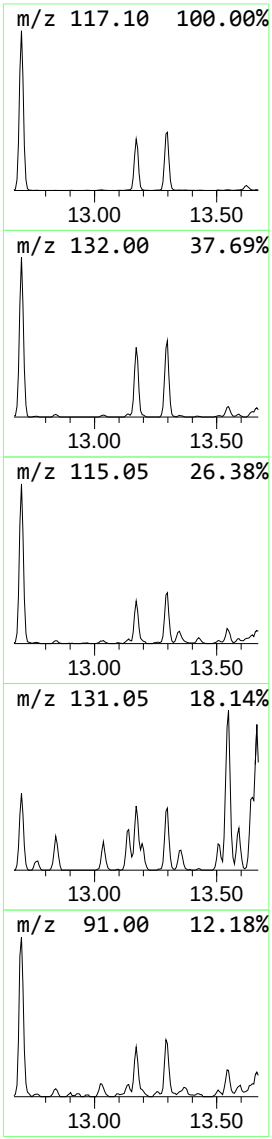
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	11.19 ug/l	96208	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

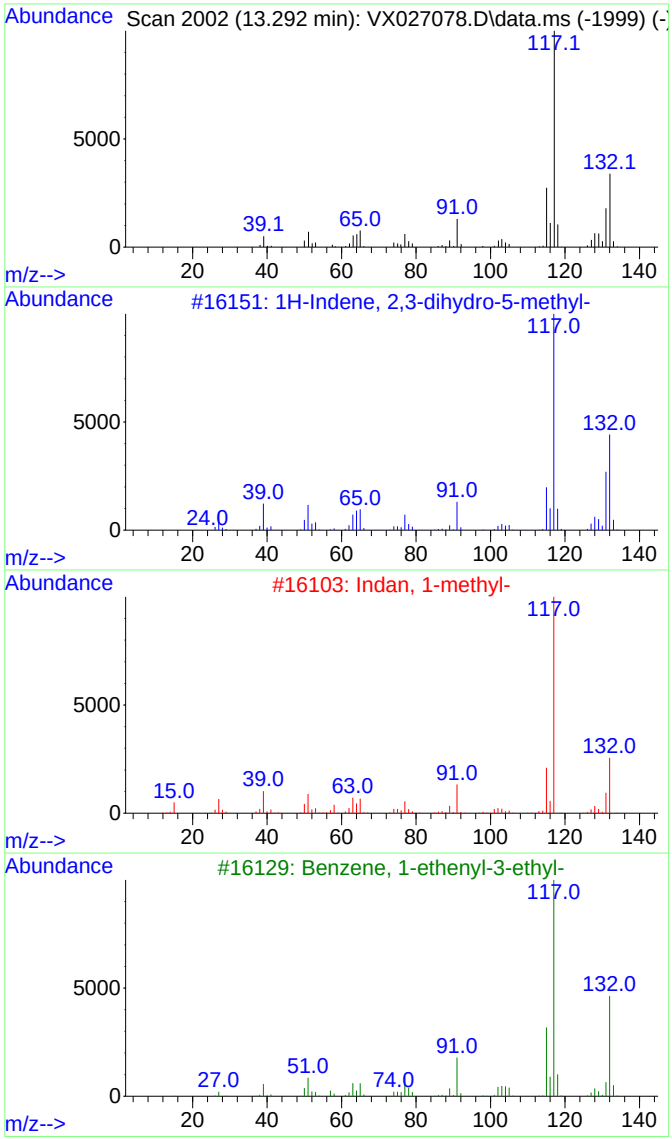
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

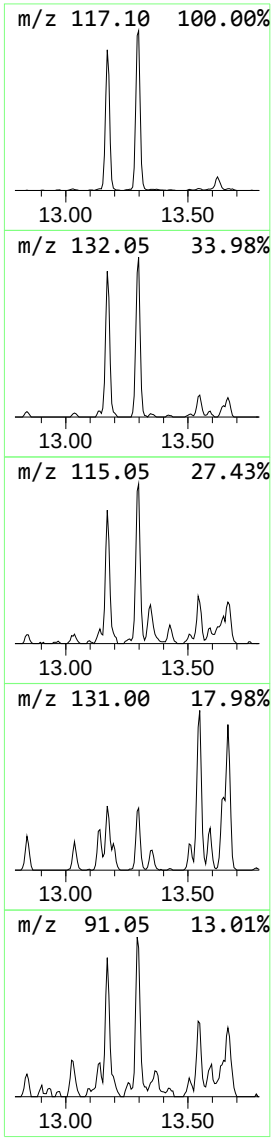
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	12.81 ug/l	110150	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	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

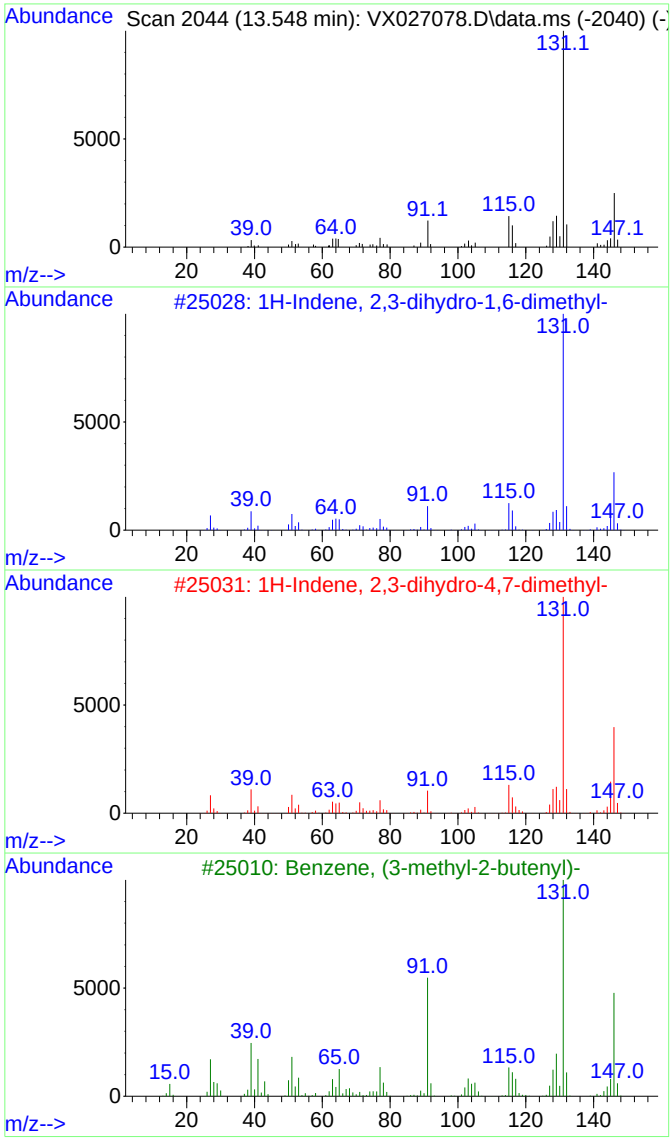
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

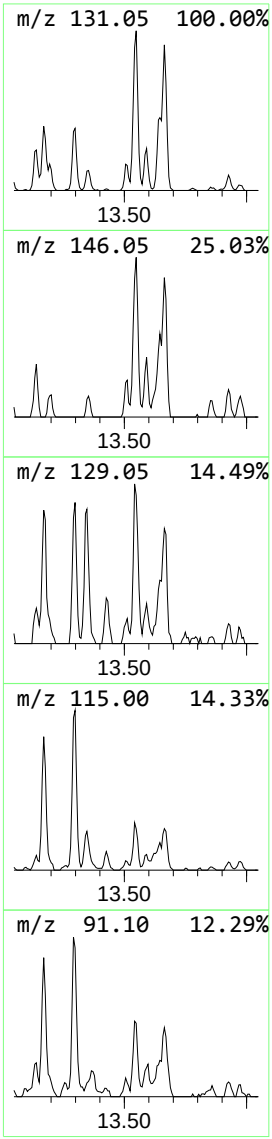
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.548	5.90 ug/l	50746	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	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

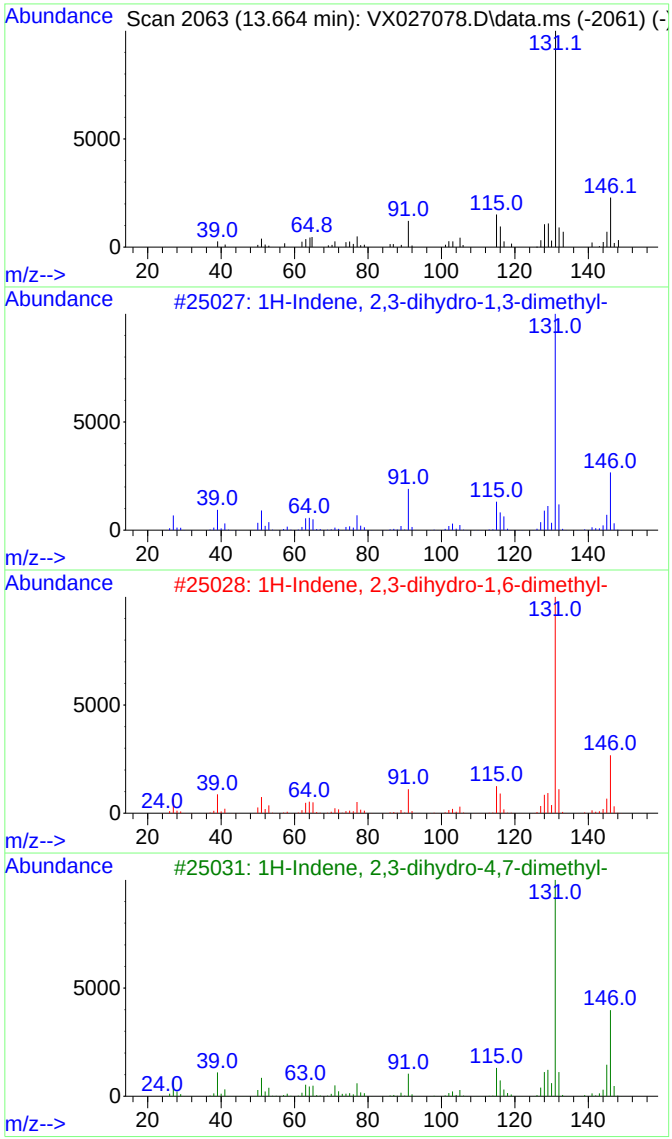
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
Quant Title : SW846 8260

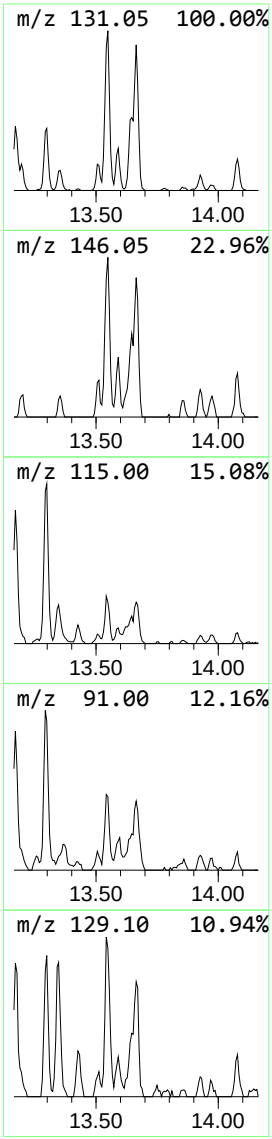
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	5.25 ug/l	45156	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	93
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	92
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021722\
Data File : VX027078.D
Acq On : 17 Feb 2022 15:23
Operator : JC/MD
Sample : N1593-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.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
Pentane, 3-methyl-	3.105	5.1	ug/l	35810	1	5.562	353219	50.0
Di-sec-Butyl ether	8.957	23.2	ug/l	231476	3	10.055	498335	50.0
3-Pentanol, 3-e...	9.049	25.2	ug/l	250827	3	10.055	498335	50.0
Benzene, cyclop...	12.249	30.5	ug/l	262551	4	12.024	430045	50.0
o-Cymene	12.615	10.6	ug/l	91146	4	12.024	430045	50.0
Indan, 1-methyl-	12.701	28.7	ug/l	246672	4	12.024	430045	50.0
Benzene, 1-ethe...	13.170	11.2	ug/l	96208	4	12.024	430045	50.0
1H-Indene, 2,3-...	13.292	12.8	ug/l	110150	4	12.024	430045	50.0
1H-Indene, 2,3-...	13.548	5.9	ug/l	50746	4	12.024	430045	50.0
1H-Indene, 2,3-...	13.664	5.3	ug/l	45156	4	12.024	430045	50.0