LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

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\82X112222W.M
Title : SW846 8260

Signal : TIC: VX033055.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.374	40	47	53	rVV	297664	346924	3.16%	0.350%
2	1.752	103	109	122	rVB	468102	614378	5.59%	0.620%
3	1.965	138	144	156	rVB2	220533	458762	4.18%	0.463%
4	2.069	156	161	169	rVV	226635	327059	2.98%	0.330%
5	2.166	169	177	181	rVV	141573	212625	1.94%	0.214%
6	2.215	181	185	196	rVB	346659	533994	4.86%	0.539%
7	2.752	259	273	281	rBV2	173349	442170	4.03%	0.446%
8	2.867	287	292	303	rVB2	221922	520981	4.74%	0.525%
9	2.989	303	312	324	rVV	100041	378362	3.45%	0.382%
10	3.111	324	332	350	rVB2	156006	382545	3.48%	0.386%
11	3.733	428	434	444	rVB3	117422	298679	2.72%	0.301%
12	3.837	444	451	457	rBV	67964	165893	1.51%	0.167%
13	3.904	457	462	472	rBV2	42704	110570	1.01%	0.112%
14	4.105	486	495	506	rBV2	78664	204707	1.86%	0.206%
15	4.306	517	528	539	rVV	284127	806831	7.35%	0.814%
16	5.196	663	674	687	rVB	80025	249344	2.27%	0.251%
17	5.391	692	706	712	rVV	71464	222889	2.03%	0.225%
18	5.477	712	720	726	rVV	127452	409006	3.72%	0.413%
19	5.550	726	732	763	rVB	145636	559925	5.10%	0.565%
20	5.958	790	799	803	rBV	70354	178960	1.63%	0.181%
21	6.044	803	813	826	rVB	4069130	10619689	96.69%	10.711%
22	6.202	826	839	844	rBV3	86432	278192	2.53%	0.281%
23	6.367	857	866	876	rVB3	35253	113411	1.03%	0.114%
24	6.763	923	931	938	rBV	225654	560216	5.10%	0.565%
25	6.842	941	944	953	rVB4	66579	170205	1.55%	0.172%
26	7.171	990	998	1008	rVB2	55324	128870	1.17%	0.130%
27	7.379	1018	1032	1045	rBV5	153196	409150	3.73%	0.413%
28	8.141	1153	1157	1163	rVB	151829	252586	2.30%	0.255%
29	8.507	1210	1217	1227	rVB	119055	212213	1.93%	0.214%
30	8.647	1234	1240	1246	rVB	344397	534909	4.87%	0.540%
31	8.720	1246	1252	1259	rVB	649941	981241	8.93%	0.990%
32	9.458	1364	1373	1377	rBV3	60488	116429	1.06%	0.117%
33	10.055	1461	1471	1476	rBV	484665	657882	5.99%	0.664%
34	10.195	1488	1494	1500	rBV	8040563	10423466	94.91%	10.513%
35	10.305	1506	1512	1518	rVB	8328564	10982910	100.00%	11.078%
36	10.646	1562	1568	1578	rVB	5126831	6787765	61.80%	6.846%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

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\82X112222W.M
Title : SW846 8260

37	10.963	1614	1620	1626	rBV	660172	810846	7.38%	0.818%
38	11.079	1634	1639	1645	rBV	327374	415188	3.78%	0.419%
39	11.305	1668	1676	1683	rBV	1689210	1975845	17.99%	1.993%
40	11.378	1683	1688	1690	rVV	1452712	2006755	18.27%	2.024%
41	11.402	1690	1692	1696	rVV	1862746	1884119	17.16%	1.900%
42	11.451	1696	1700	1706	rVB	1556094	1826793	16.63%	1.843%
43	11.616	1721	1727	1736	rBV	2038654	2574999	23.45%	2.597%
44	11.756	1744	1750	1755	rVB	7578162	9159834	83.40%	9.239%
45	11.969	1779	1785	1789	rBV	90911	117629	1.07%	0.119%
46	12.024	1789	1794	1798	rVV2	570432	758913	6.91%	0.765%
47	12.085	1798	1804	1809	rVB	2483588	3125571	28.46%	3.152%
48	12.244	1824	1830	1838	rBV	3516536	4556529	41.49%	4.596%
49	12.317	1838	1842	1850	rVB3	599067	883572	8.04%	0.891%
50	12.408	1852	1857	1862	rBV2	260178	361854	3.29%	0.365%
51	12.463	1862	1866	1872	rVB	239627	291795	2.66%	0.294%
52	12.530	1872	1877	1879	rBV	454005	543086	4.94%	0.548%
53	12.555	1879	1881	1886	rVB	343446	362122	3.30%	0.365%
54	12.616	1886	1891	1894	rBV	892666	1123399	10.23%	1.133%
55	12.646	1894	1896	1900	rVB	217067	218537	1.99%	0.220%
56	12.701	1900	1905	1913	rBV2	671172	955395	8.70%	0.964%
57	12.847	1924	1929	1934	rVB2	224099	286399	2.61%	0.289%
58	12.920	1934	1941	1945	rBV	572912	699930	6.37%	0.706%
59	12.963	1945	1948	1954	rVB	660262	749928	6.83%	0.756%
60	13.170	1978	1982	1987	rVB	625152	721499	6.57%	0.728%
61	13.292	1997	2002	2007	rVB2	1429834	1834420	16.70%	1.850%
62	13.420	2018	2023	2032	rVB	409744	521735	4.75%	0.526%
63	13.542	2040	2043	2046	rBV	126057	146895	1.34%	0.148%
64	13.585	2046	2050	2055	rVB3	133028	203800	1.86%	0.206%
65	13.664	2060	2063	2069	rVB2	151952	236346	2.15%	0.238%
66	13.774	2073	2081	2090	rBV	2893233	3676138	33.47%	3.708%
67	13.865	2090	2096	2100	rBV2	86959	145461	1.32%	0.147%
68	13.926	2100	2106	2110	rBV2	97124	131114	1.19%	0.132%
69	14.073	2126	2130	2134	rBV	110817	144807	1.32%	0.146%
70	14.225	2149	2155	2160	rBV2	145789	296494	2.70%	0.299%
71	14.481	2192	2197	2202	rBV2	102429	145204	1.32%	0.146%
72	14.634	2210	2222	2233	rBV	1366356	1851103	16.85%	1.867%
73	14.780	2241	2246	2255	rBV	1057243	1329127	12.10%	1.341%
74	15.207	2306	2316	2320	rBV2	59909	121168	1.10%	0.122%
75	15.371	2335	2343	2347	rBV	69891	114173	1.04%	0.115%
76	15.475	2353	2360	2365	rBV	195788	304133	2.77%	0.307%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

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\82X112222W.M
Title : SW846 8260

77	15.615	2373	2383	2386	rBV	273109	438991	4.00%	0.443%
78	15.652	2386	2389	2395	rVB	167458	234427	2.13%	0.236%
79	15.835	2409	2419	2437	rVB	76428	206260	1.88%	0.208%

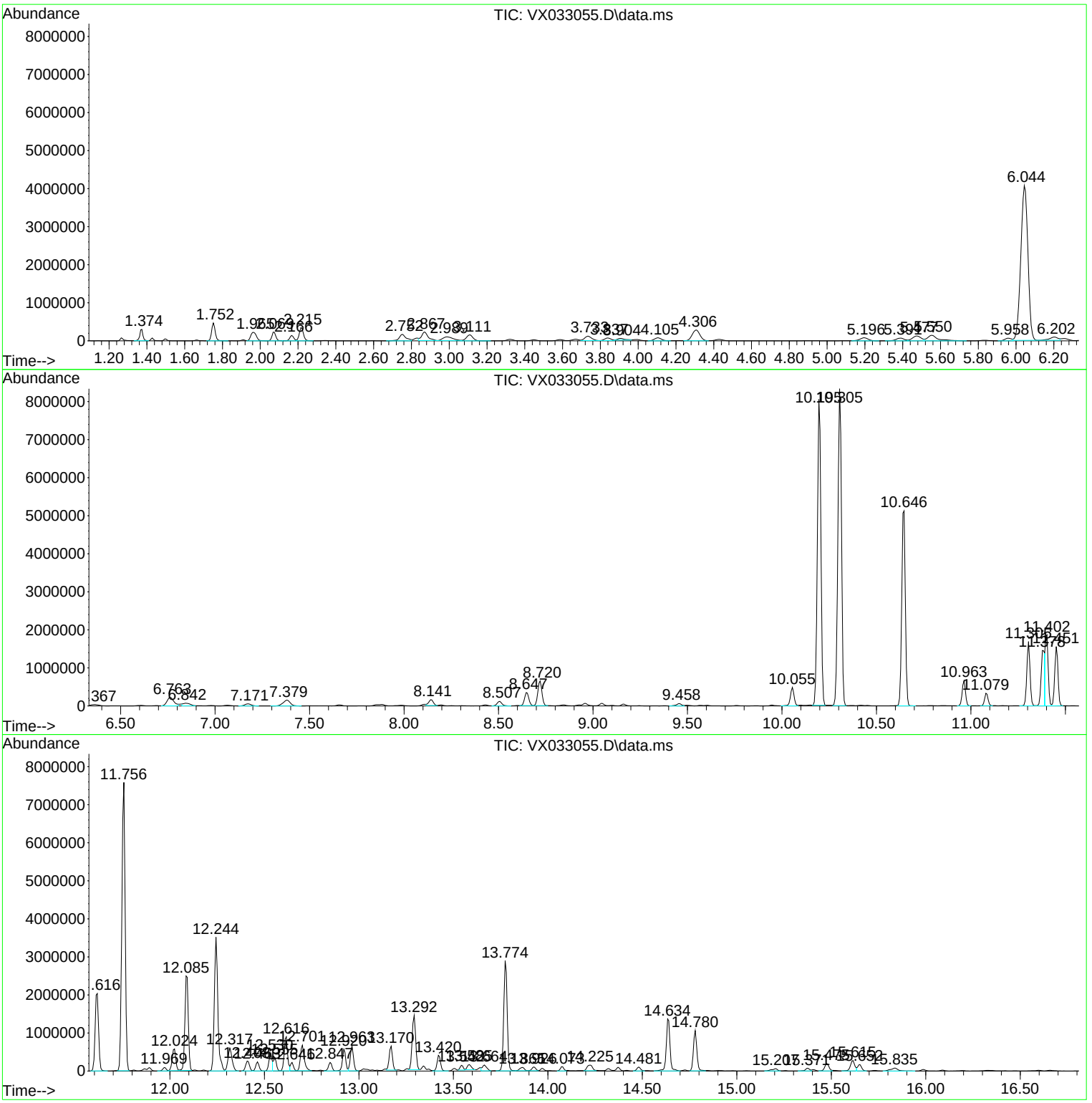
Sum of corrected areas: 99146071

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

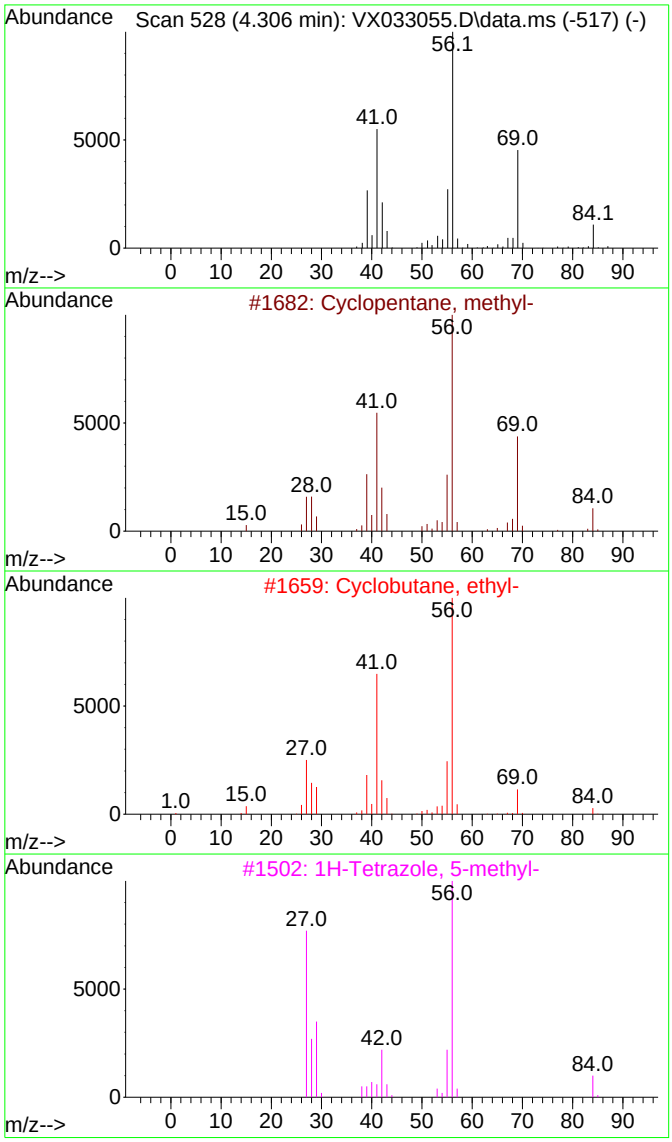
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

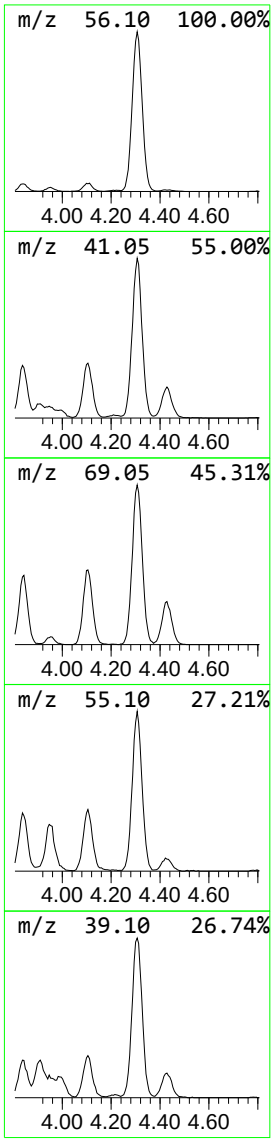
****

Peak Number 1 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	72.05 ug/l	806831	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	56
5	2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

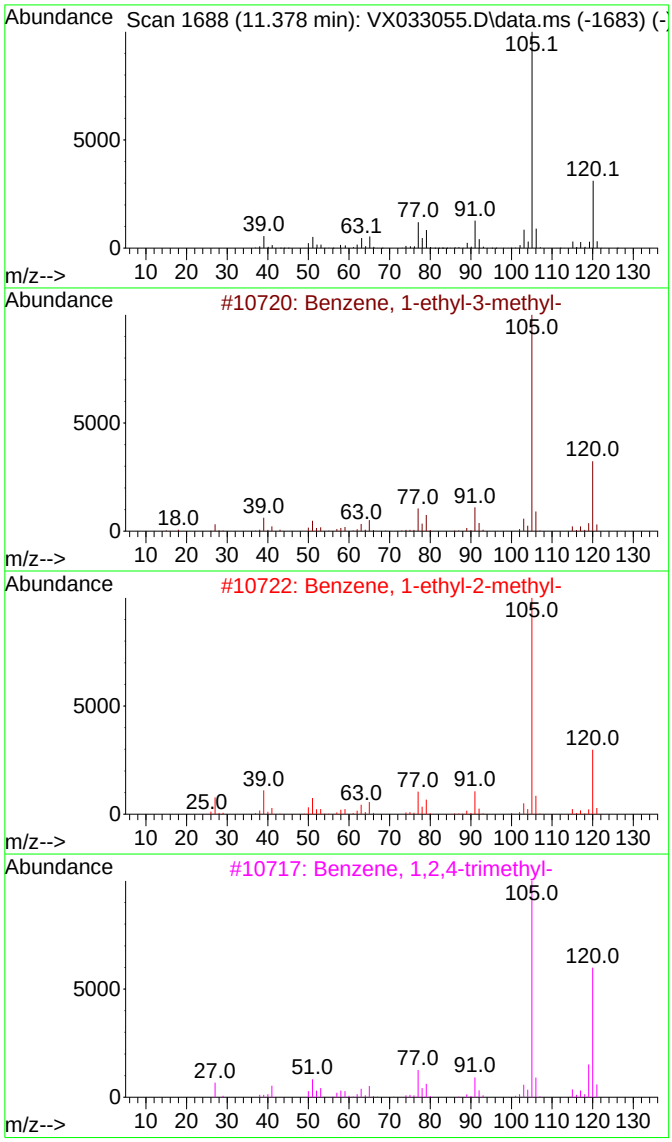
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

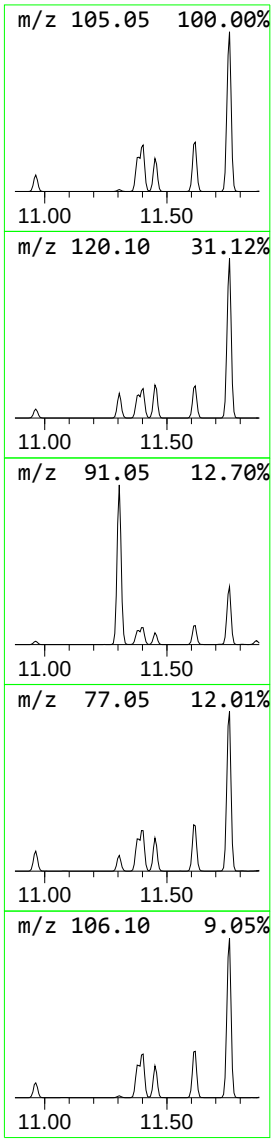
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	132.21 ug/l	2006760	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

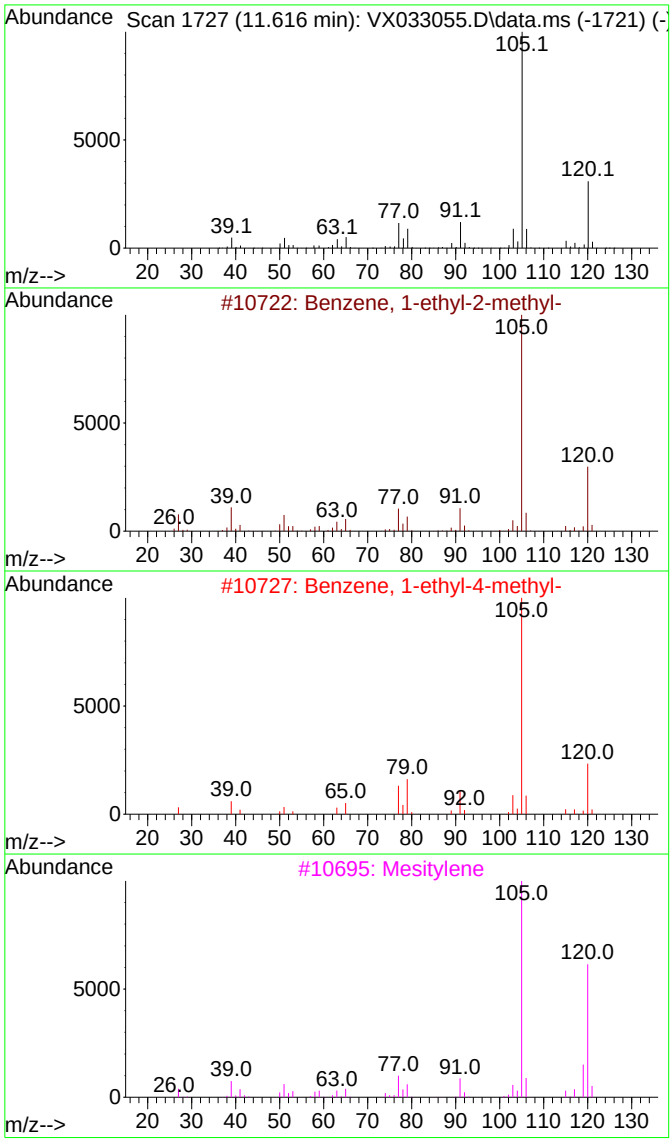
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

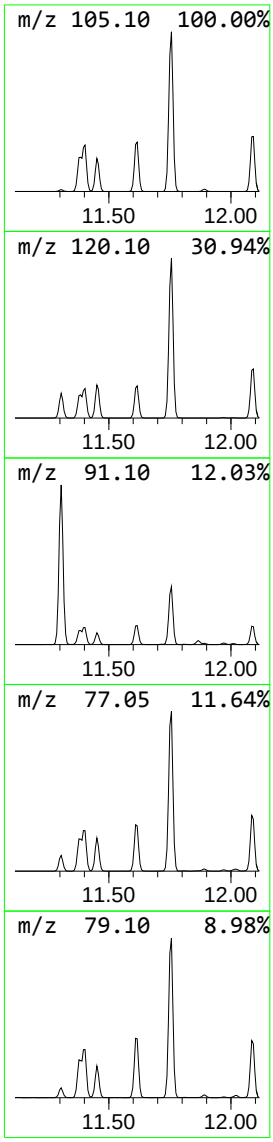
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	169.65 ug/l	2575000	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

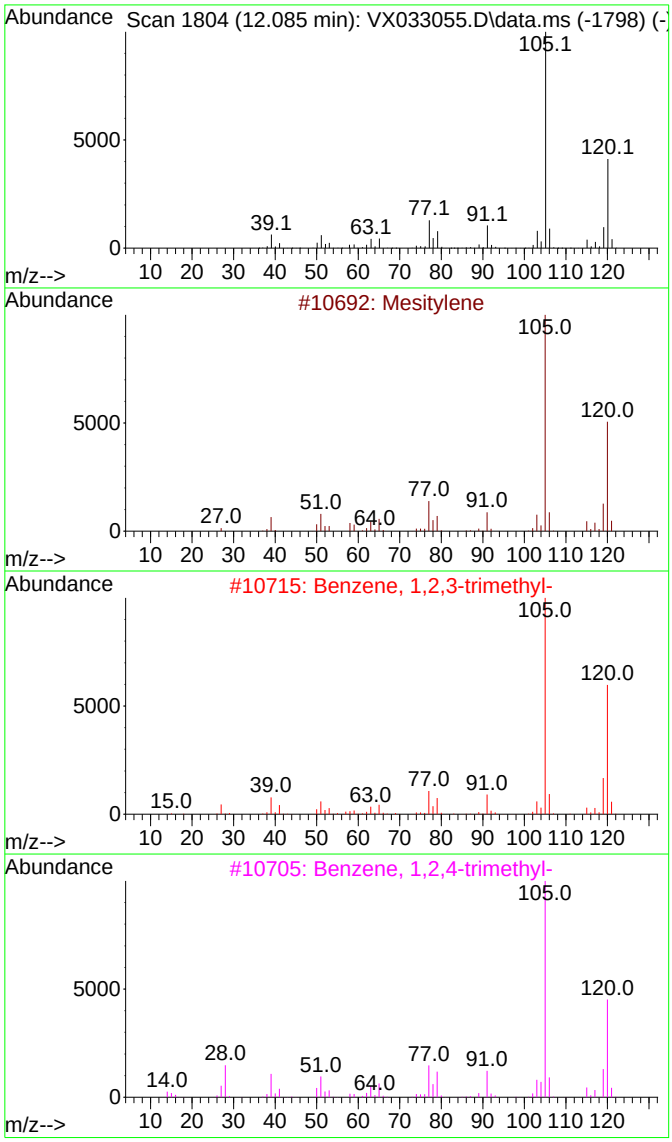
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

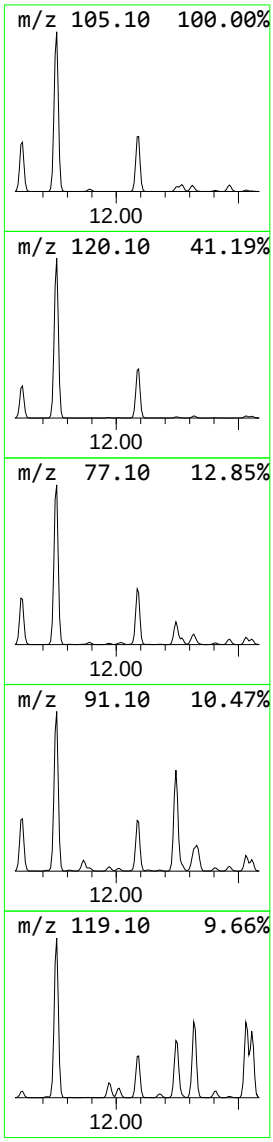
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	205.92 ug/l	3125570	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

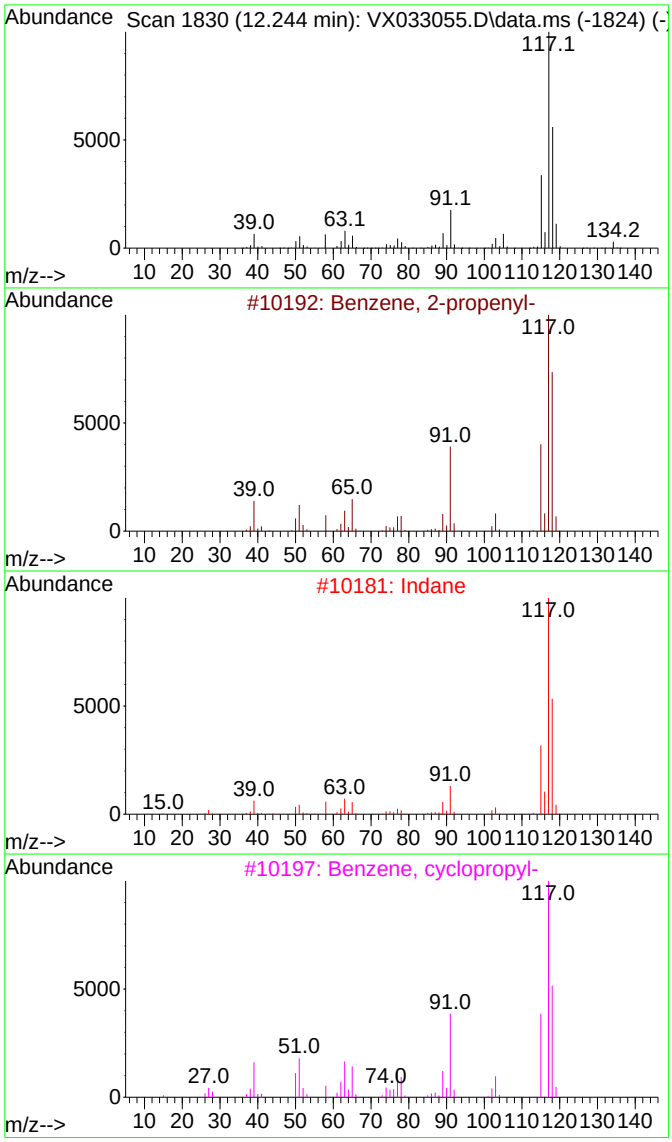
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

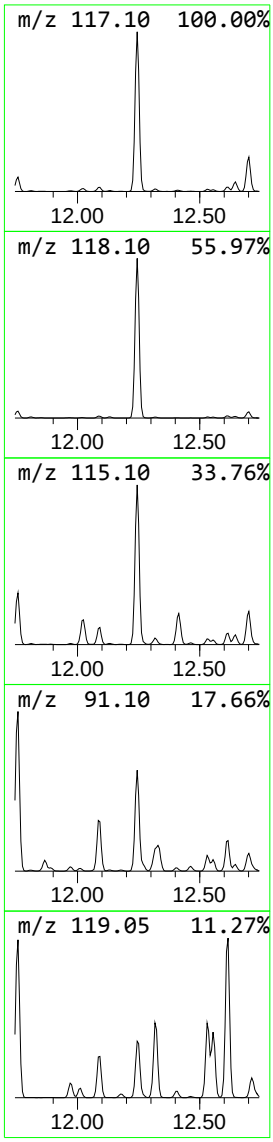
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 2-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	300.20 ug/l	4556530	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
2	Indane	118	C9H10	000496-11-7	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Deltacyclene	118	C9H10	007785-10-6	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

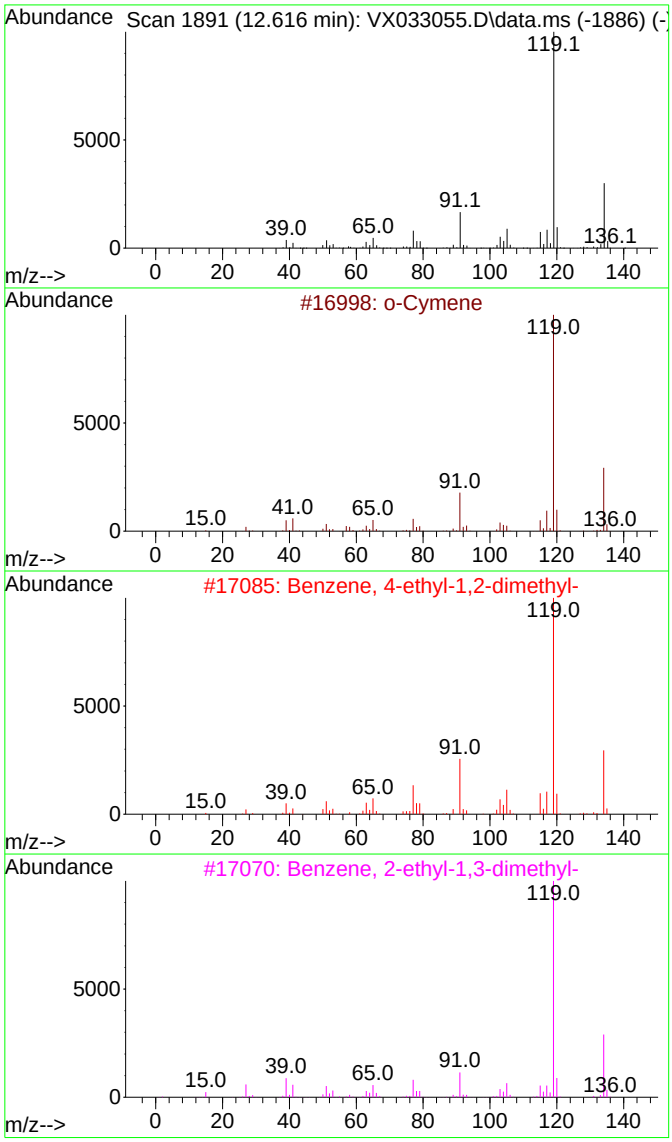
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

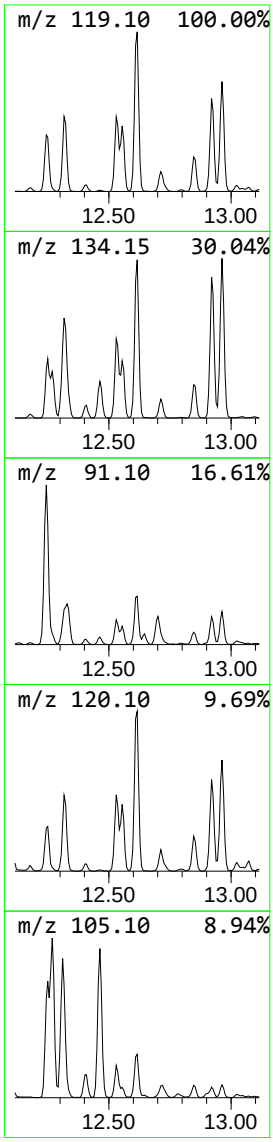
****

Peak Number 6 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	74.01 ug/l	1123400	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, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

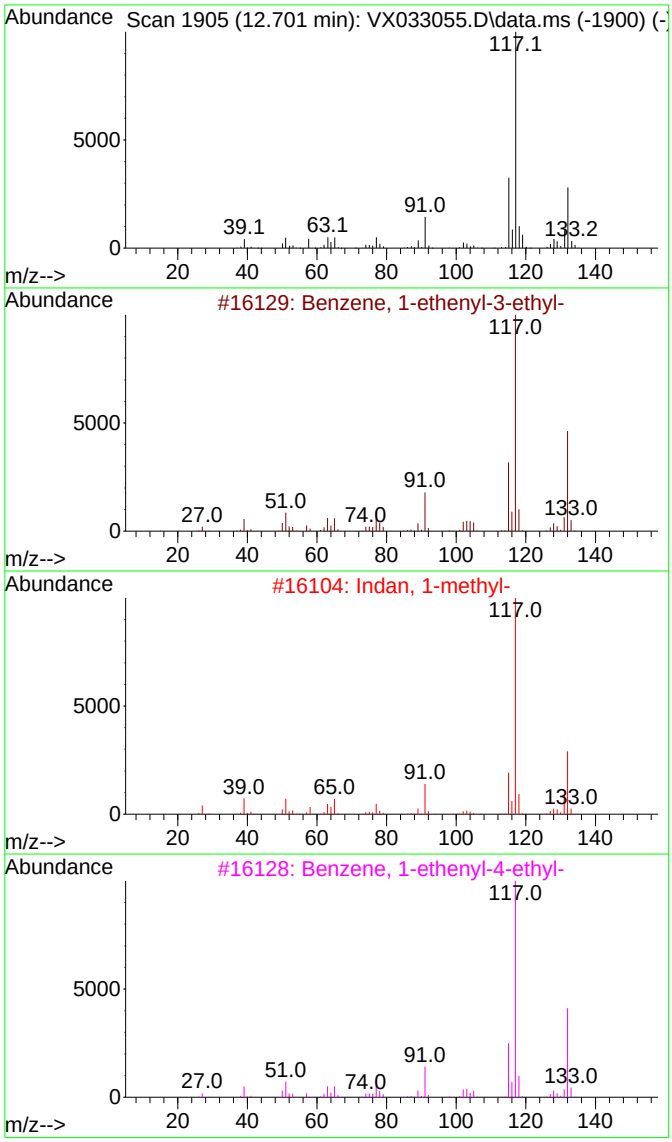
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

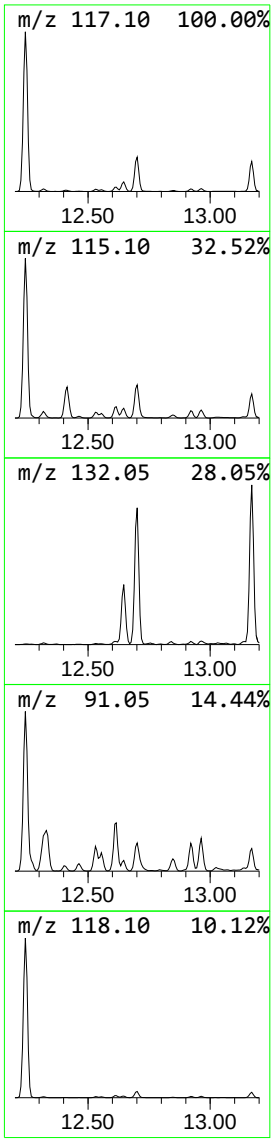
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	62.95 ug/l	955395	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	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	83
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

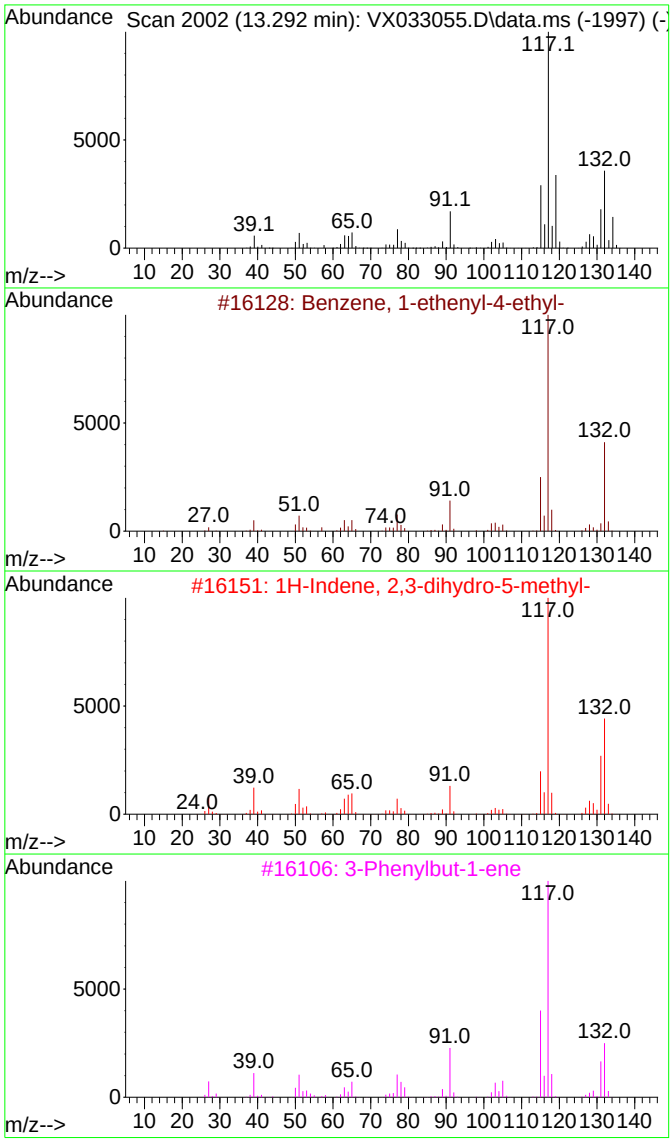
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

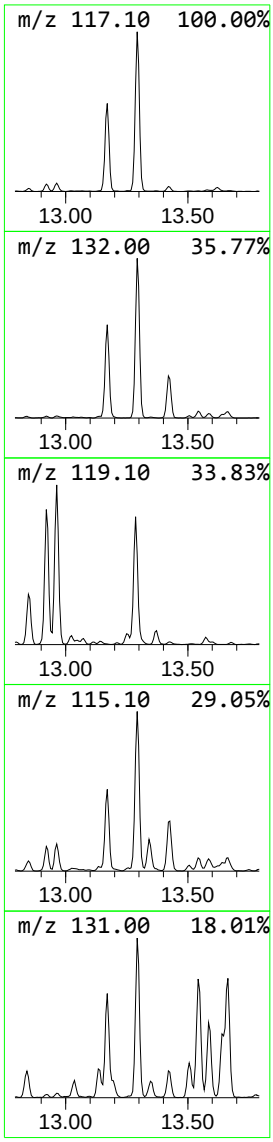
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	120.86 ug/l	1834420	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	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

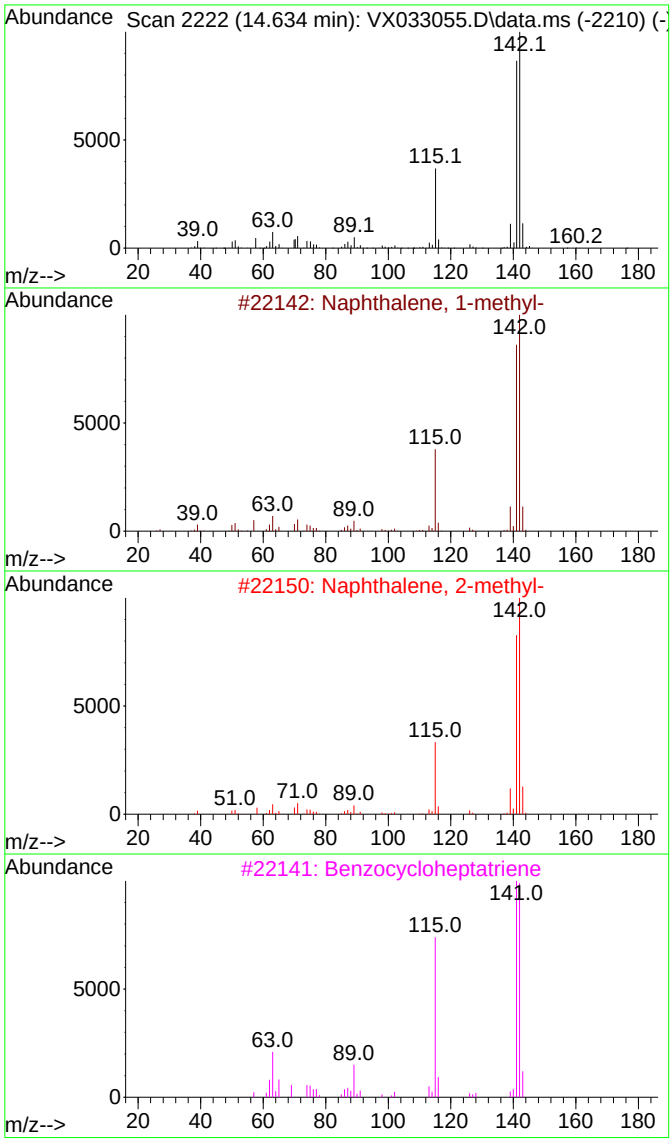
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

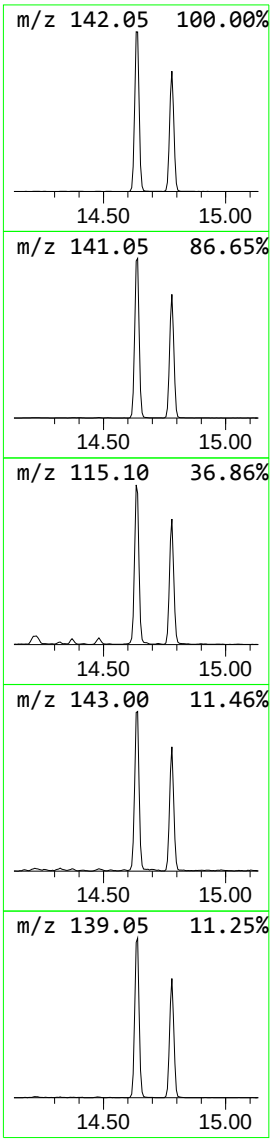
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzocycloheptatriene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	121.96 ug/l	1851100	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

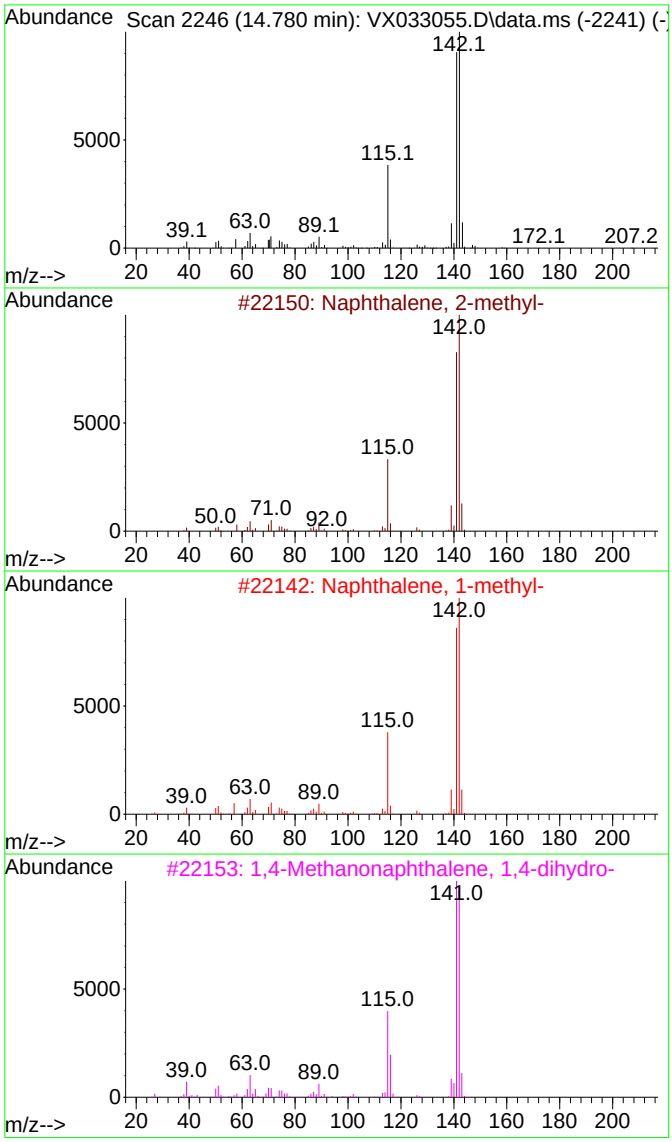
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

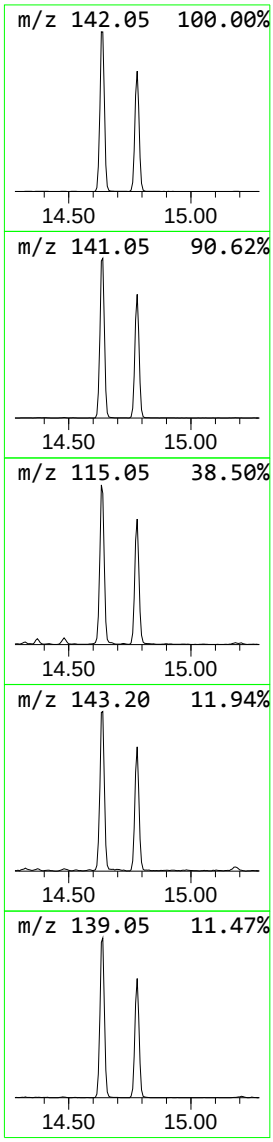
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	87.57 ug/l	1329130	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033055.D
Acq On : 24 Nov 2022 11:25
Operator : JC/MD
Sample : N5747-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.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
Cyclopentane, m...	4.306	72.0	ug/l	806831	1	5.556	559925	50.0
Benzene, 1-ethy...	11.378	132.2	ug/l	2006760	4	12.024	758913	50.0
Benzene, 1-ethy...	11.616	169.7	ug/l	2575000	4	12.024	758913	50.0
Benzene, 1,2,3-...	12.085	205.9	ug/l	3125570	4	12.024	758913	50.0
Benzene, 2-prop...	12.244	300.2	ug/l	4556530	4	12.024	758913	50.0
o-Cymene	12.616	74.0	ug/l	1123400	4	12.024	758913	50.0
Benzene, 1-ethe...	12.701	63.0	ug/l	955395	4	12.024	758913	50.0
Benzene, 1-ethe...	13.292	120.9	ug/l	1834420	4	12.024	758913	50.0
Benzocyclohepta...	14.634	122.0	ug/l	1851100	4	12.024	758913	50.0
1,4-Methanonaph...	14.780	87.6	ug/l	1329130	4	12.024	758913	50.0