LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

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\82X091922W.M
Title : SW846 8260

Signal : TIC: VX031529.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	119	rVB	313717	406552	2.57%	0.381%
2	1.953	137	142	152	rVB2	173121	259040	1.64%	0.243%
3	2.069	155	161	167	rBV	118691	166170	1.05%	0.156%
4	2.209	180	184	194	rVB	125504	188879	1.19%	0.177%
5	2.776	258	277	280	rBV2	122191	382609	2.42%	0.359%
6	2.825	280	285	289	rVV	268369	575290	3.64%	0.540%
7	2.867	289	292	306	rVB2	171389	409111	2.59%	0.384%
8	3.105	320	331	352	rVB2	269442	666649	4.21%	0.625%
9	3.447	378	387	400	rVB	87834	195279	1.23%	0.183%
10	3.733	426	434	444	rVB	93334	226631	1.43%	0.213%
11	4.306	518	528	540	rVB	615259	1722282	10.88%	1.616%
12	5.385	691	705	712	rBV	94576	305610	1.93%	0.287%
13	5.477	712	720	729	rBV4	180346	623863	3.94%	0.585%
14	5.617	739	743	764	rVB4	109556	391851	2.48%	0.368%
15	5.830	766	778	791	rBV2	111565	334415	2.11%	0.314%
16	5.964	791	800	803	rBV	88273	216480	1.37%	0.203%
17	6.050	803	814	825	rVB	6275386	15825682	100.00%	14.846%
18	6.257	844	848	857	rVB3	176049	448767	2.84%	0.421%
19	6.361	857	865	877	rVB3	84024	253756	1.60%	0.238%
20	6.763	922	931	938	rBV	138264	359553	2.27%	0.337%
21	6.848	938	945	956	rVB4	84196	266605	1.68%	0.250%
22	7.379	1019	1032	1043	rBV4	384900	1055750	6.67%	0.990%
23	7.659	1072	1078	1091	rVB	93029	184762	1.17%	0.173%
24	7.988	1122	1132	1144	rVB	78578	179569	1.13%	0.168%
25	8.110	1144	1152	1154	rBV	139875	265285	1.68%	0.249%
26	8.147	1154	1158	1163	rVB	176785	290000	1.83%	0.272%
27	8.507	1210	1217	1226	rVB2	218650	383246	2.42%	0.360%
28	8.653	1235	1241	1247	rVB	515195	822807	5.20%	0.772%
29	8.720	1247	1252	1258	rBV	263966	392673	2.48%	0.368%
30	8.842	1264	1272	1281	rVB2	100944	193658	1.22%	0.182%
31	8.958	1281	1291	1300	rBV2	312628	546563	3.45%	0.513%
32	9.049	1300	1306	1311	rBV	338604	502929	3.18%	0.472%
33	10.055	1468	1471	1475	rVB	288732	360216	2.28%	0.338%
34	10.201	1489	1495	1500	rBV	10637198	13431683	84.87%	12.600%
35	10.305	1506	1512	1517	rVB	1575111	2020663	12.77%	1.896%
36	10.640	1562	1567	1577	rVB	321452	436505	2.76%	0.409%

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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

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\82X091922W.M
Title : SW846 8260

37	10.964	1613	1620	1626	rVB	1667806	2035990	12.87%	1.910%
38	11.085	1635	1640	1644	rVB	483457	608205	3.84%	0.571%
39	11.305	1666	1676	1682	rBV	4899591	5705978	36.06%	5.353%
40	11.402	1682	1692	1696	rVV	472656	906579	5.73%	0.850%
41	11.451	1696	1700	1707	rVB	355925	455540	2.88%	0.427%
42	11.616	1721	1727	1735	rBV	1326147	1588017	10.03%	1.490%
43	11.756	1744	1750	1755	rVB	1333651	1567970	9.91%	1.471%
44	11.866	1763	1768	1770	rBV	143831	190691	1.20%	0.179%
45	11.890	1770	1772	1778	rVB	214268	240131	1.52%	0.225%
46	12.024	1789	1794	1799	rVB	410263	518649	3.28%	0.487%
47	12.091	1799	1805	1809	rBV	1278432	1548785	9.79%	1.453%
48	12.244	1824	1830	1837	rBV	8900763	10815669	68.34%	10.146%
49	12.317	1837	1842	1852	rVB2	611675	1086285	6.86%	1.019%
50	12.408	1852	1857	1862	rBV2	306529	401414	2.54%	0.377%
51	12.463	1862	1866	1873	rVB	348349	405233	2.56%	0.380%
52	12.530	1873	1877	1880	rBV	242403	313995	1.98%	0.295%
53	12.616	1885	1891	1894	rBV	3289766	3739417	23.63%	3.508%
54	12.646	1894	1896	1900	rVB	539769	553447	3.50%	0.519%
55	12.701	1900	1905	1913	rBV2	1780011	2508878	15.85%	2.354%
56	12.847	1924	1929	1934	rVB3	155129	218661	1.38%	0.205%
57	12.921	1934	1941	1945	rBV	1926216	2335946	14.76%	2.191%
58	12.963	1945	1948	1953	rVB	360855	420972	2.66%	0.395%
59	13.024	1953	1958	1967	rBV2	120287	198376	1.25%	0.186%
60	13.170	1978	1982	1992	rVB	1663864	1982075	12.52%	1.859%
61	13.292	1997	2002	2007	rVV2	3875662	5168825	32.66%	4.849%
62	13.341	2007	2010	2013	rVV4	272860	361280	2.28%	0.339%
63	13.427	2019	2024	2030	rVB	505111	651981	4.12%	0.612%
64	13.506	2030	2037	2040	rBV2	136763	194568	1.23%	0.183%
65	13.542	2040	2043	2046	rVV	260620	325675	2.06%	0.306%
66	13.585	2046	2050	2055	rVB4	295276	439994	2.78%	0.413%
67	13.664	2060	2063	2069	rVB2	309342	497795	3.15%	0.467%
68	13.780	2074	2082	2091	rBV	6345867	8291328	52.39%	7.778%
69	13.926	2100	2106	2110	rBV2	164223	224532	1.42%	0.211%
70	14.079	2126	2131	2138	rBV	212048	280342	1.77%	0.263%
71	14.213	2148	2153	2163	rBV2	198551	384249	2.43%	0.360%
72	14.371	2175	2179	2184	rVB	136917	175332	1.11%	0.164%
73	14.640	2218	2223	2233	rVB	1354985	1716532	10.85%	1.610%
74	14.780	2241	2246	2255	rBV	1426428	1808554	11.43%	1.697%
75	15.475	2354	2360	2367	rBV	111746	185429	1.17%	0.174%
76	15.615	2374	2383	2386	rBV	162920	253281	1.60%	0.238%

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ALS Vial : 21 Sample Multiplier: 1

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ClientSampleId :
MW-6R

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_X\Method\82X091922W.M
Title : SW846 8260

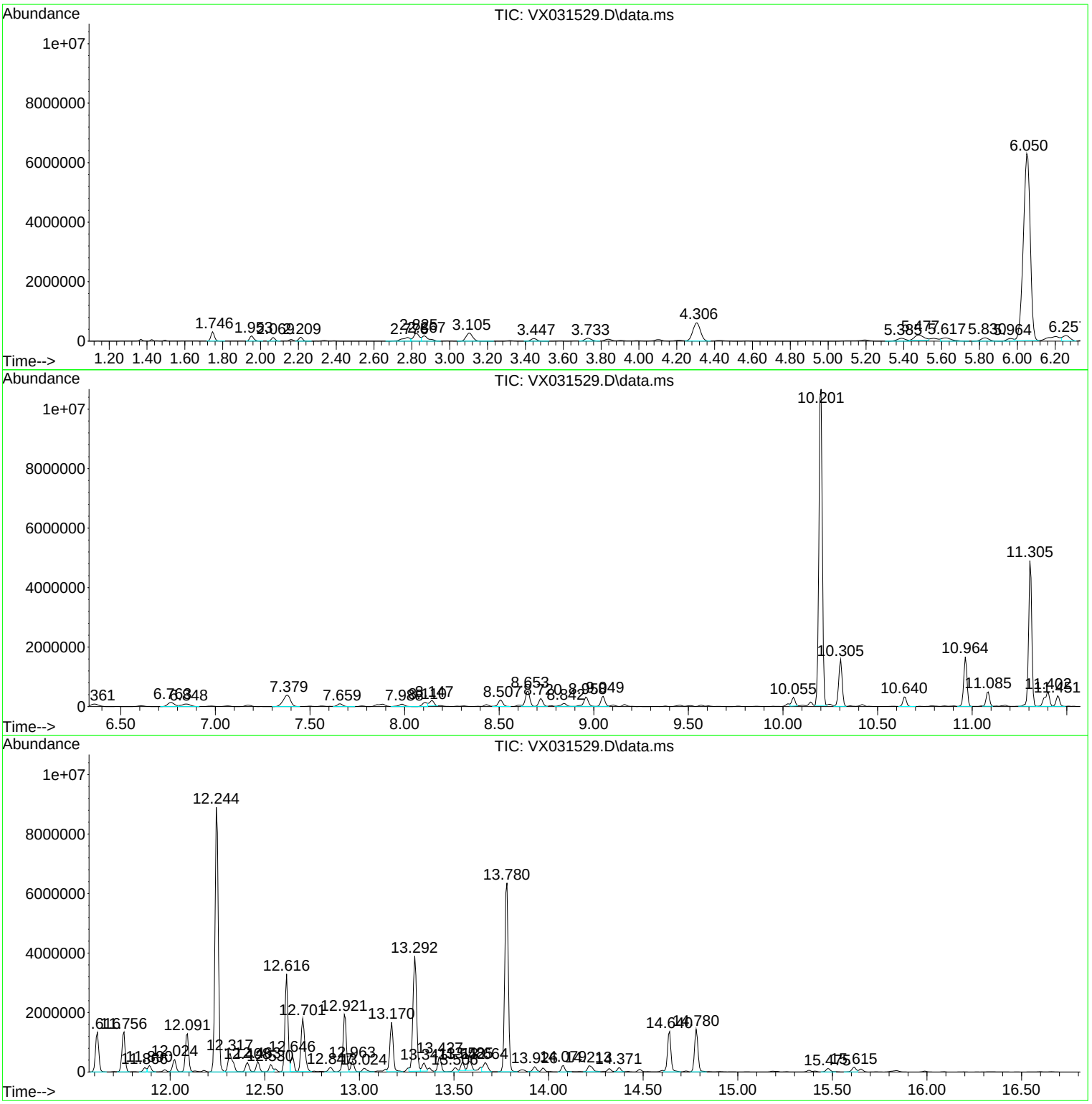
Sum of corrected areas: 106597983

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
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Instrument :
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ClientSampleId :
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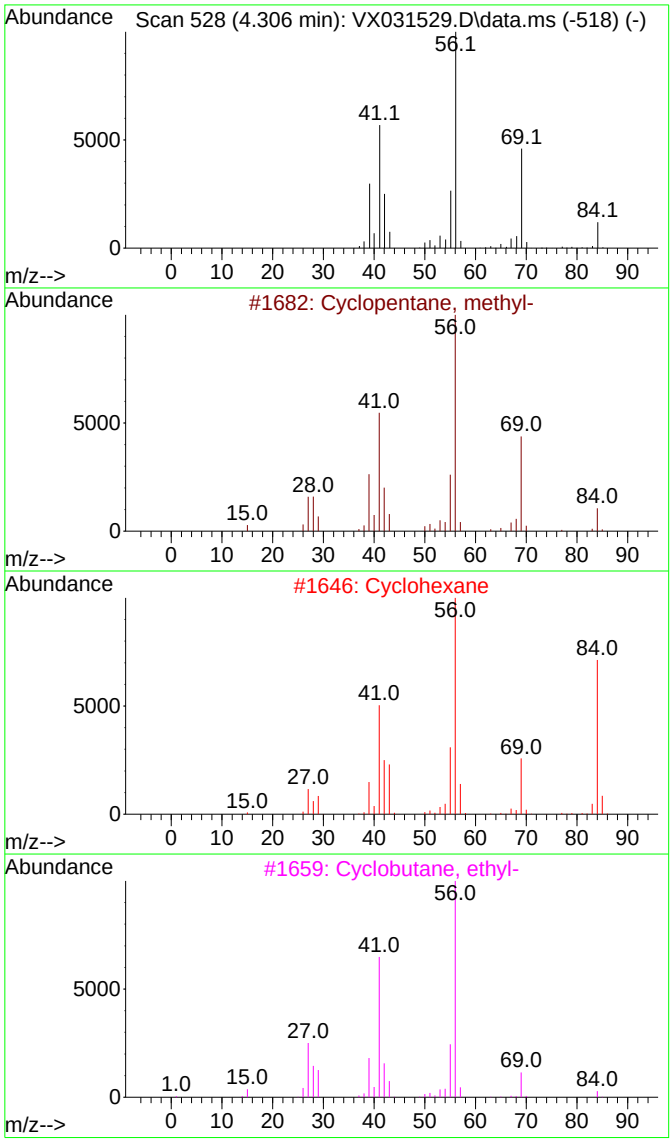
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

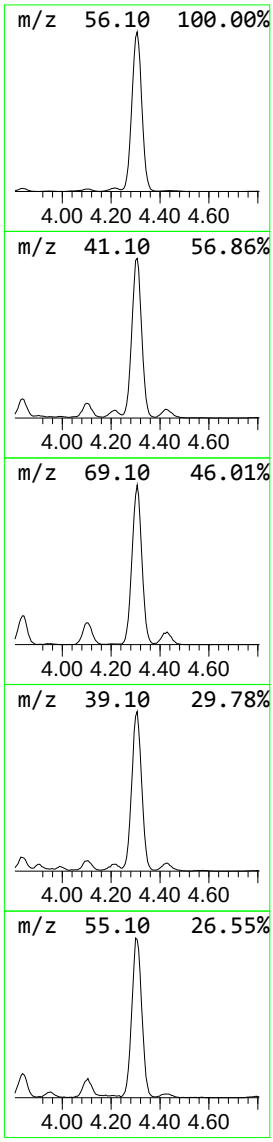
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	219.76 ug/l	1722280	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclohexane	84	C6H12	000110-82-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59
5	Cyclobutane	56	C4H8	000287-23-0	50





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Operator : JC/MD
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Instrument :
MSVOA_X
ClientSampleId :
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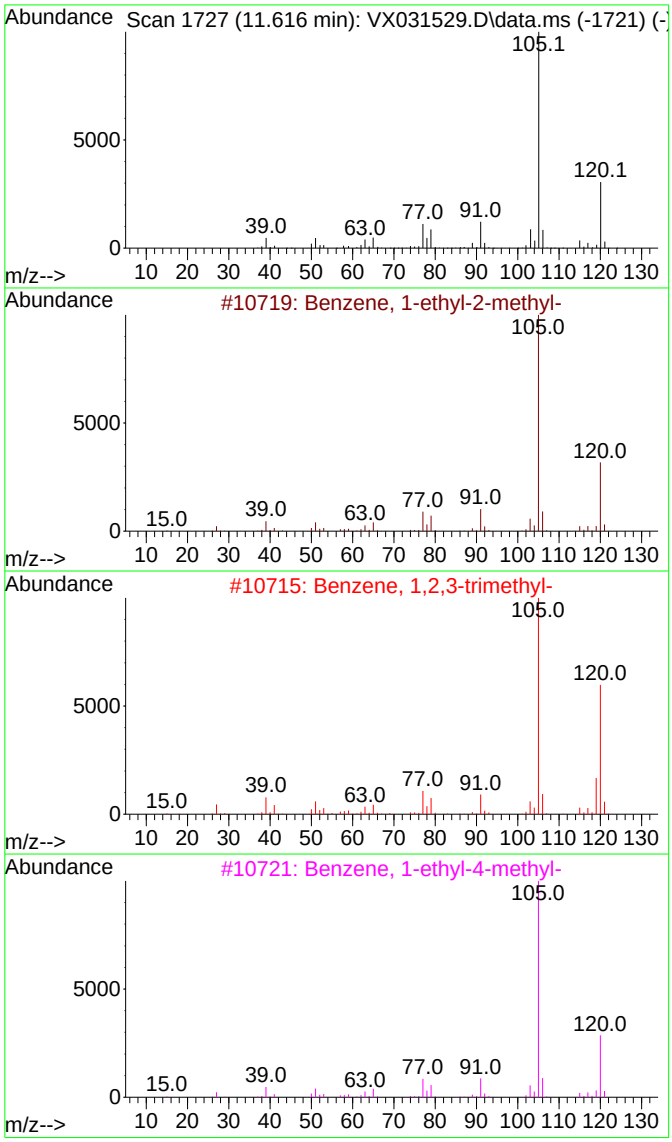
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

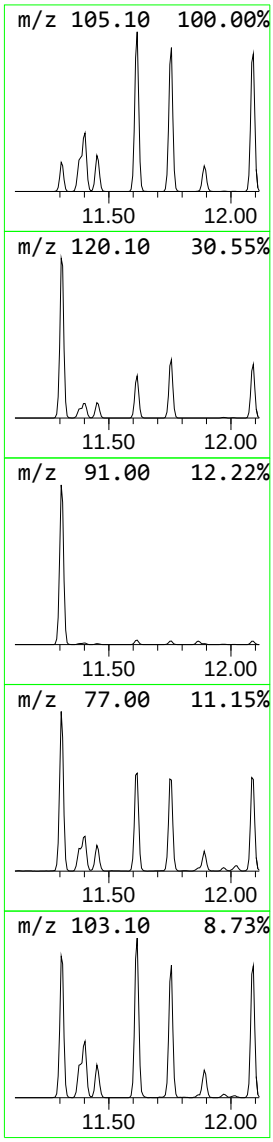
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	153.09 ug/l	1588020	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

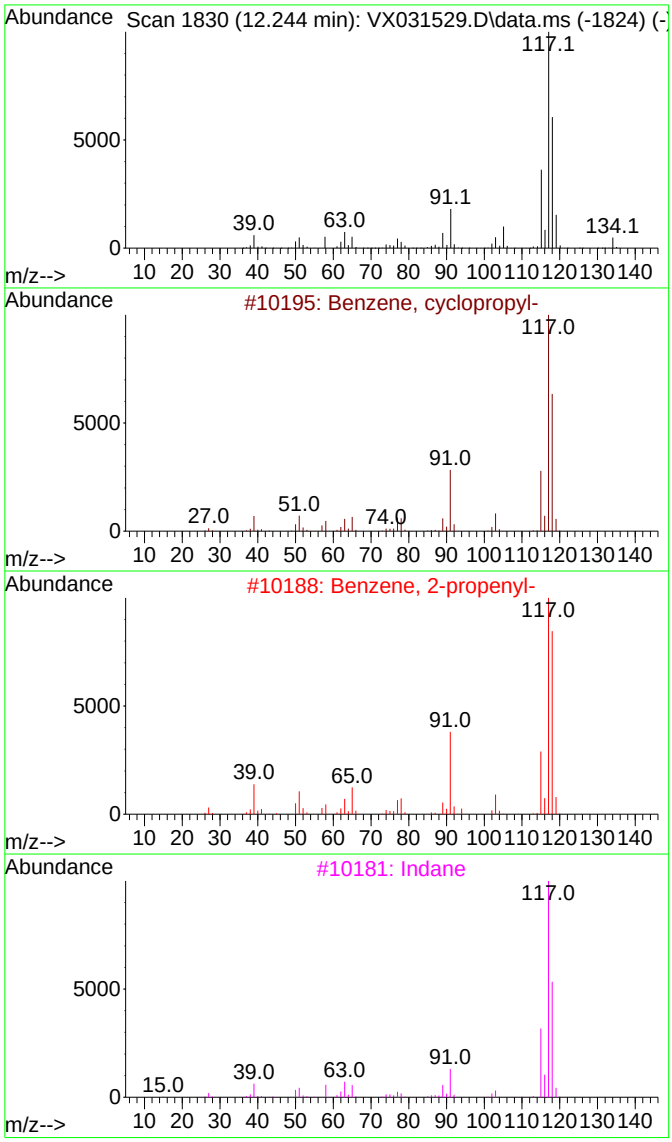
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

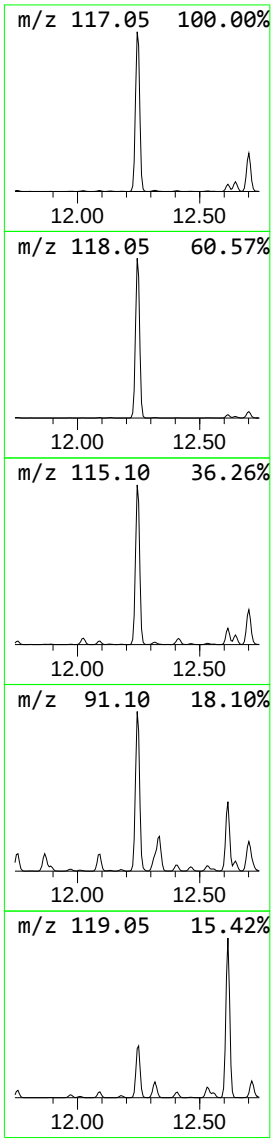
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	1042.68 ug/l	10815700	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
3	Indane	118	C9H10	000496-11-7	76
4	1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	72
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	68





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Instrument :
MSVOA_X
ClientSampleId :
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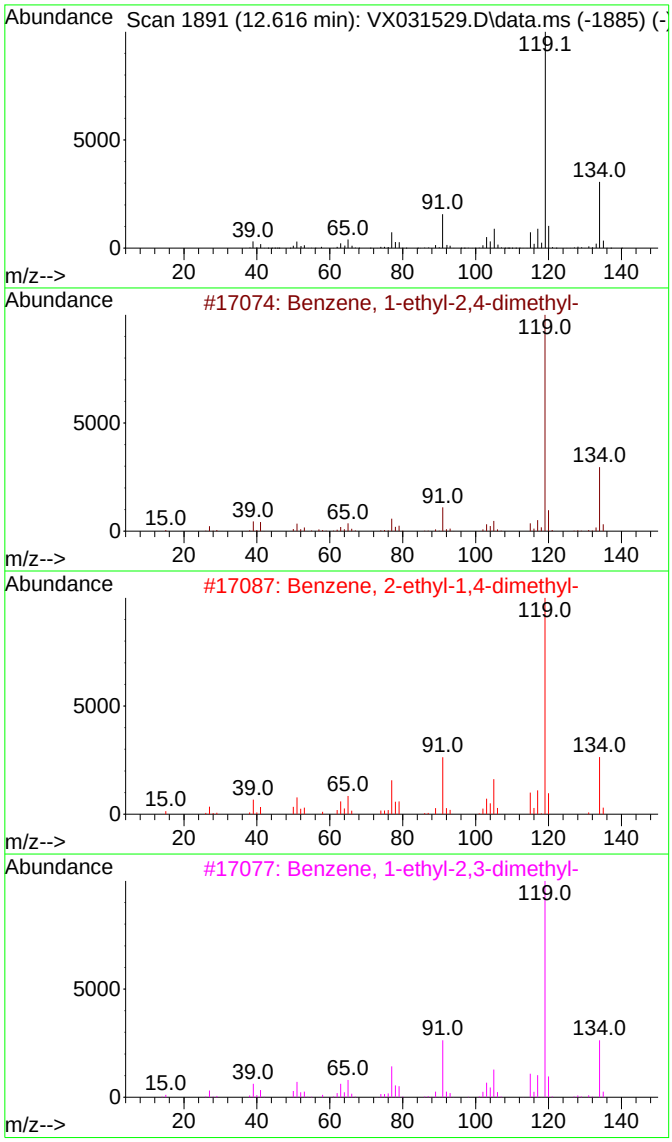
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

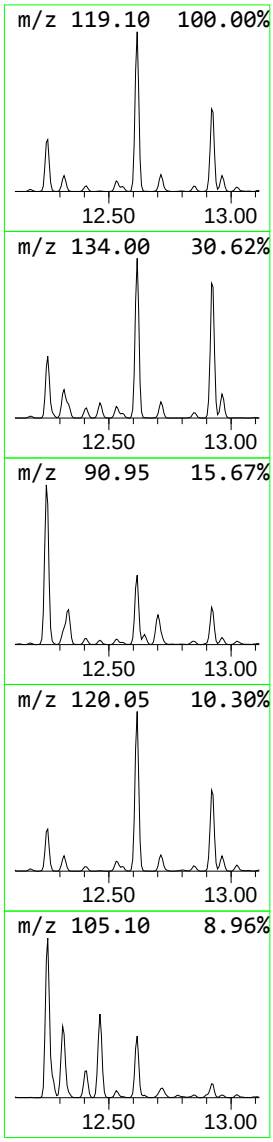
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	360.50 ug/l	3739420	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





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Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

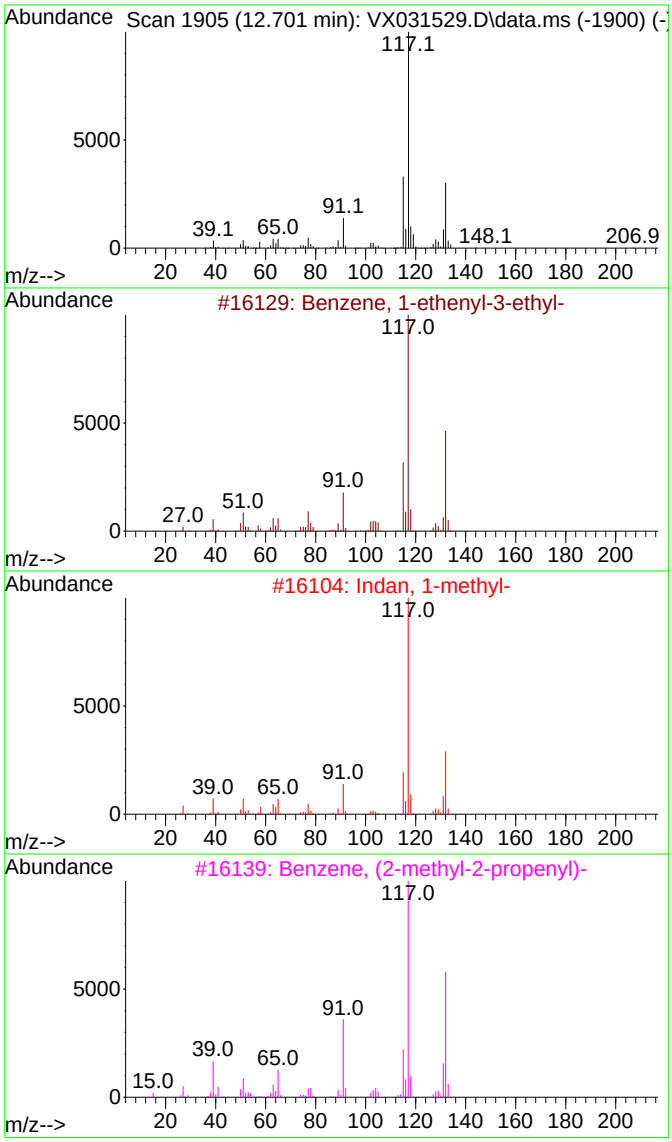
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

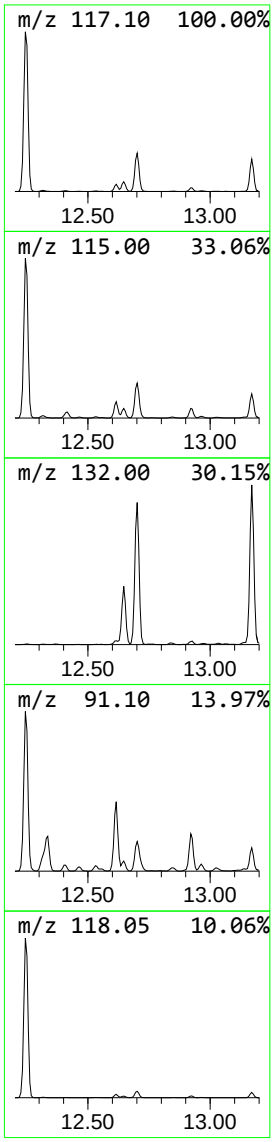
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	241.87 ug/l	2508880	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	90
3	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86





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Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

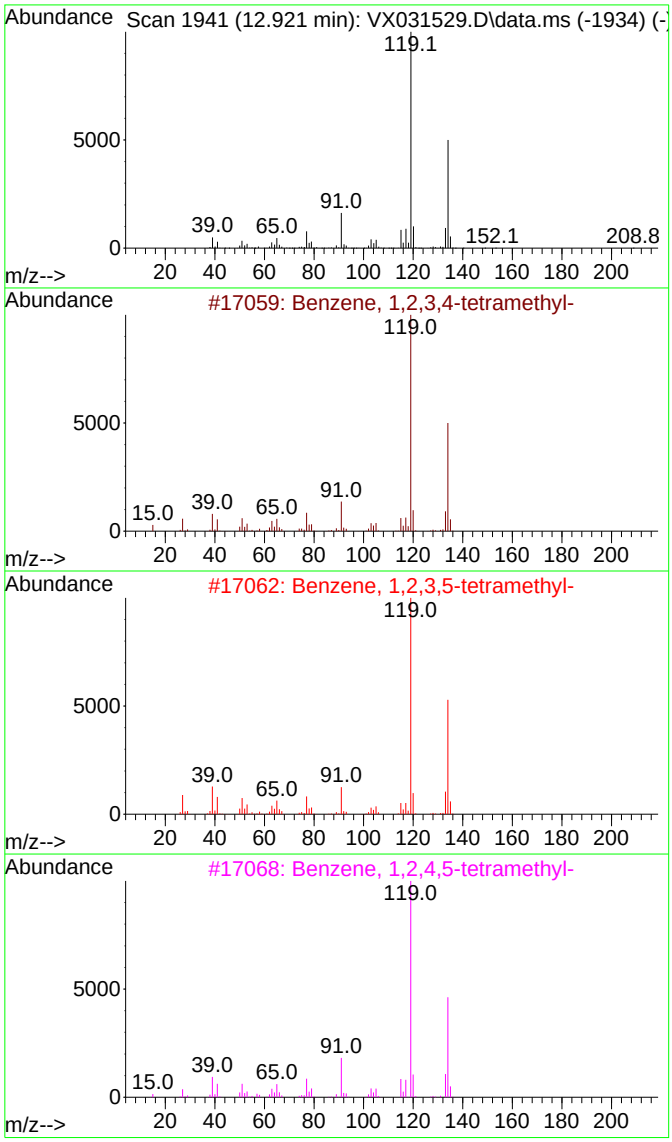
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

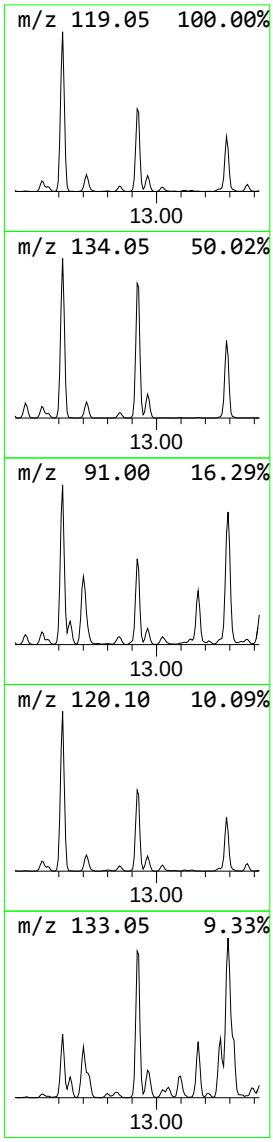
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.921	225.19 ug/l	2335950	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	o-Cymene	134	C10H14	000527-84-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

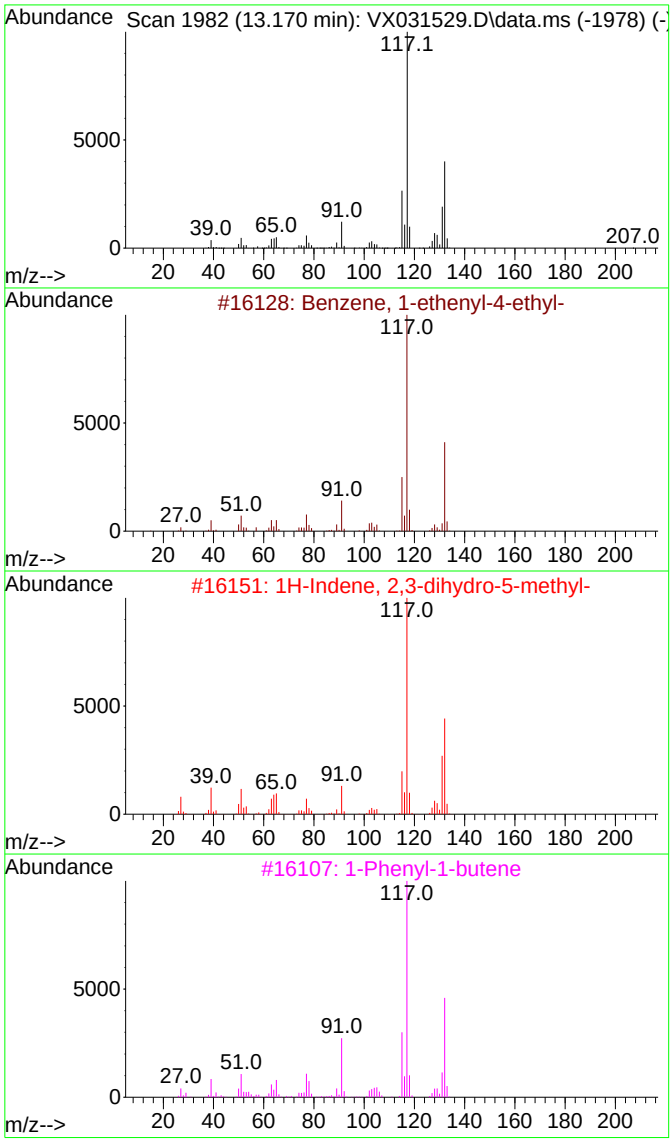
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

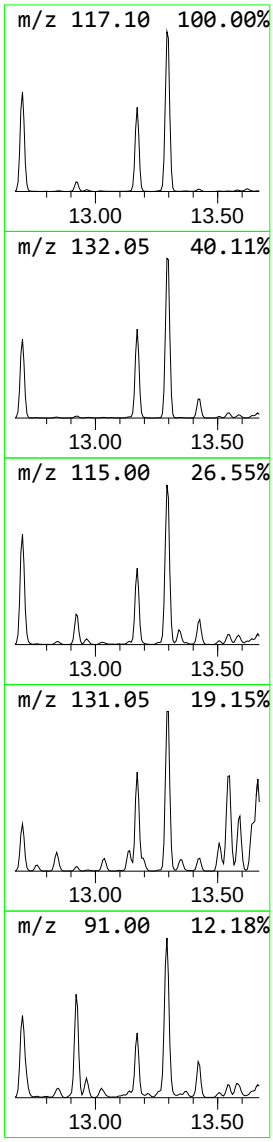
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.171	191.08 ug/l	1982080	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	93
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

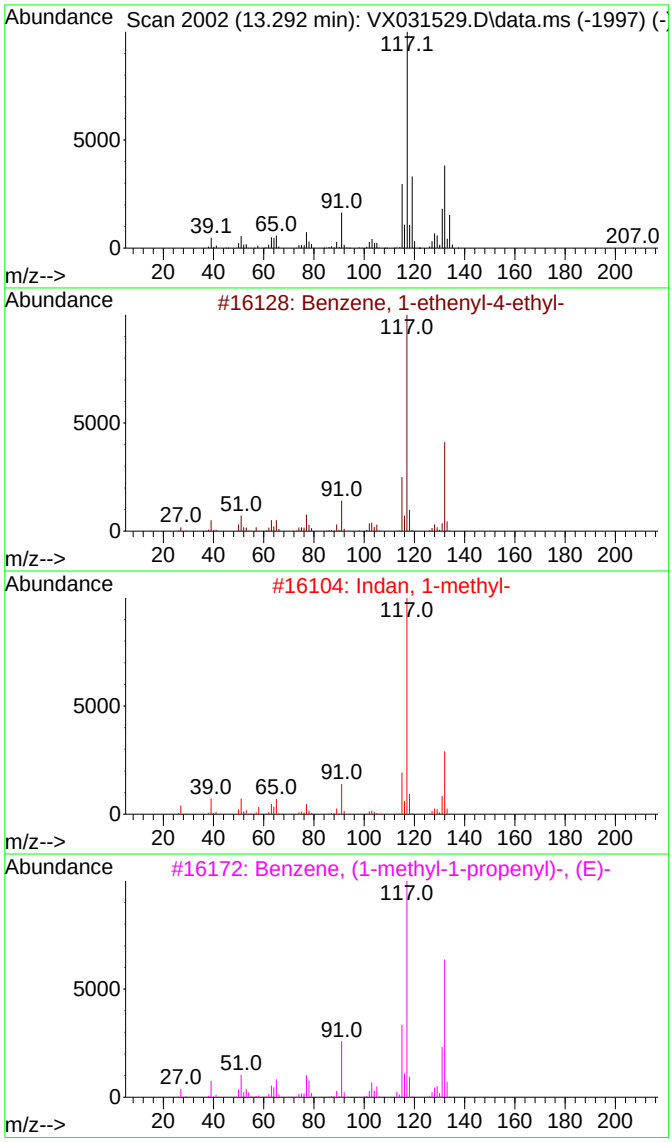
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

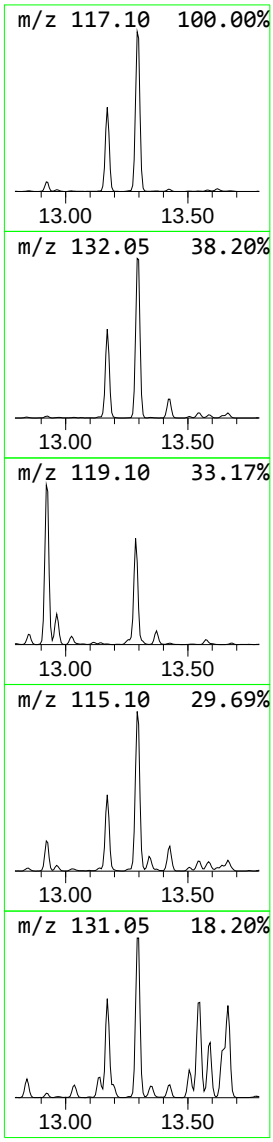
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	498.30 ug/l	5168830	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	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

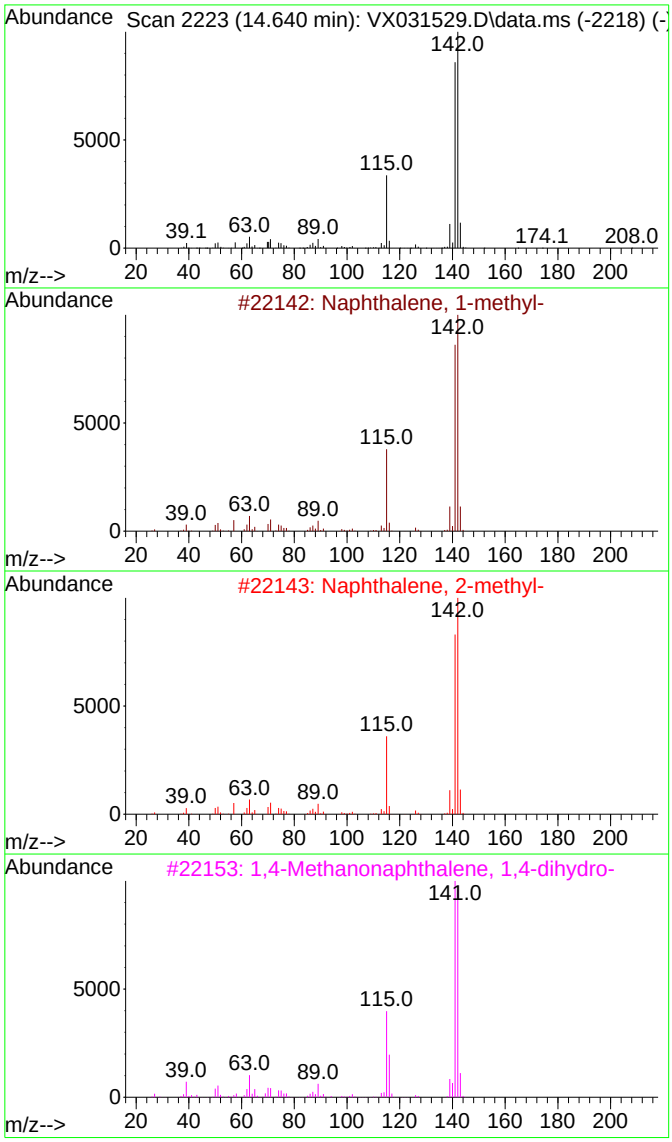
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

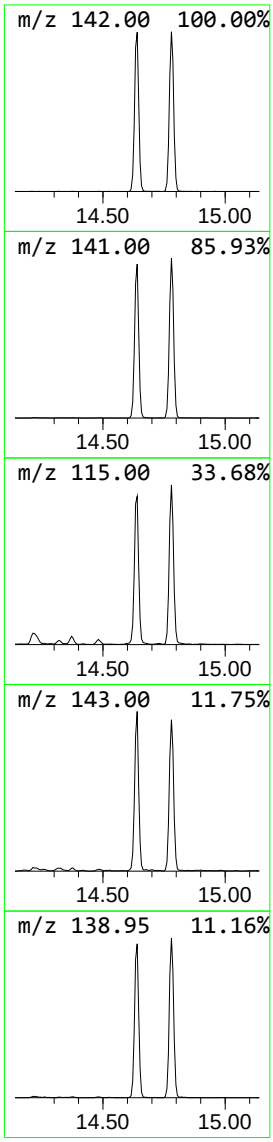
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	165.48 ug/l	1716530	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	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

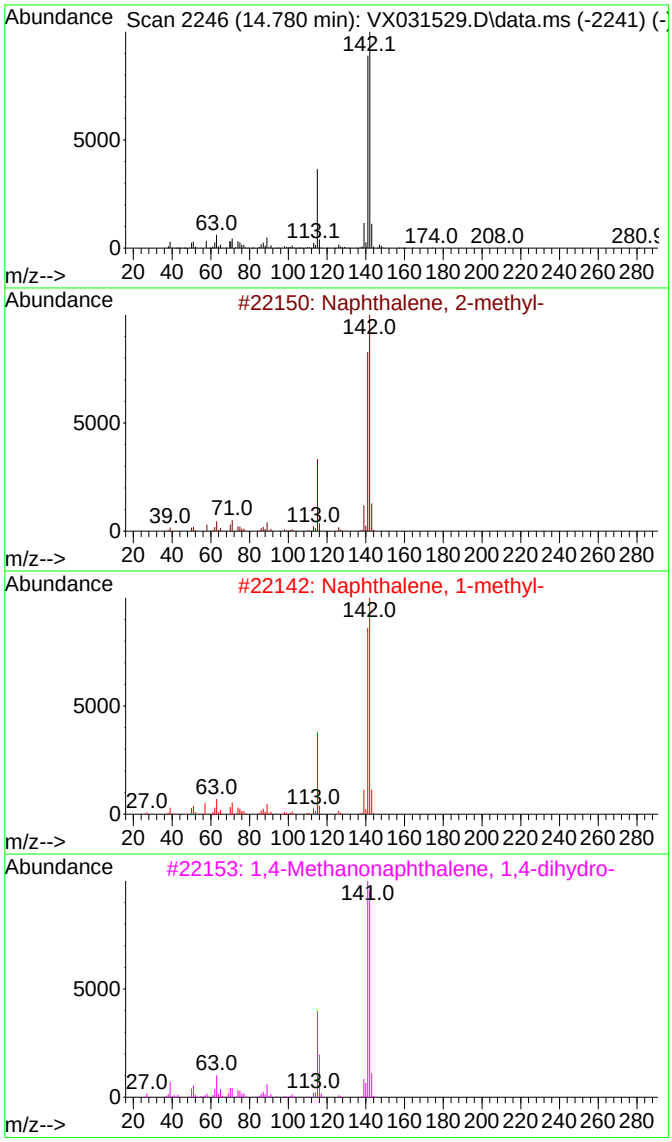
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

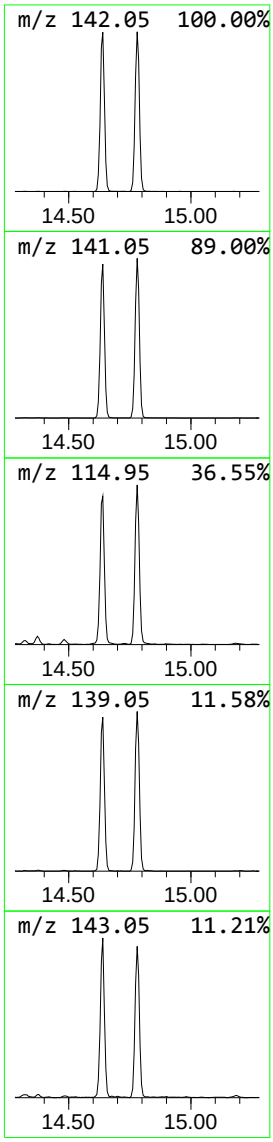
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	174.35 ug/l	1808550	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	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031529.D
Acq On : 21 Sep 2022 17:03
Operator : JC/MD
Sample : N4614-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.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	219.8	ug/l	1722280	1	5.556	391851	50.0
Benzene, 1-ethy...	11.616	153.1	ug/l	1588020	4	12.024	518649	50.0
Benzene, cyclop...	12.244	1042.7	ug/l	10815700	4	12.024	518649	50.0
Benzene, 1-ethy...	12.616	360.5	ug/l	3739420	4	12.024	518649	50.0
Benzene, 1-ethe...	12.701	241.9	ug/l	2508880	4	12.024	518649	50.0
Benzene, 1,2,3,...	12.921	225.2	ug/l	2335950	4	12.024	518649	50.0
Benzene, 1-ethe...	13.171	191.1	ug/l	1982080	4	12.024	518649	50.0
Indan, 1-methyl-	13.292	498.3	ug/l	5168830	4	12.024	518649	50.0
Naphthalene, 1-...	14.640	165.5	ug/l	1716530	4	12.024	518649	50.0
Naphthalene, 2-...	14.780	174.3	ug/l	1808550	4	12.024	518649	50.0