LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

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\82X041222W.M
Title : SW846 8260

Signal : TIC: VX028108.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	428183	590856	1.95%	0.289%
2	2.782	270	278	282	rBV	340885	761059	2.52%	0.372%
3	2.825	282	285	289	rVB	198520	321943	1.06%	0.158%
4	3.111	321	332	346	rVB3	555268	1398146	4.62%	0.684%
5	3.733	425	434	443	rBV	159101	380051	1.26%	0.186%
6	4.105	481	495	504	rBV	114630	315371	1.04%	0.154%
7	4.306	519	528	539	rVB	579629	1628183	5.38%	0.797%
8	5.190	649	673	690	rVB2	152170	610985	2.02%	0.299%
9	5.391	691	706	712	rBV	187109	555050	1.84%	0.272%
10	5.477	712	720	727	rVV3	217064	741386	2.45%	0.363%
11	5.556	727	733	739	rVV2	257595	788702	2.61%	0.386%
12	5.623	739	744	765	rVB3	202972	690459	2.28%	0.338%
13	5.958	790	799	806	rBV	179635	465117	1.54%	0.228%
14	6.251	824	847	857	rVV6	308750	1450189	4.79%	0.709%
15	6.367	857	866	877	rVB2	117487	337707	1.12%	0.165%
16	6.763	921	931	938	rBV	427824	1125146	3.72%	0.550%
17	6.848	941	945	957	rVB5	197640	543441	1.80%	0.266%
18	7.178	990	999	1008	rBV	168250	380771	1.26%	0.186%
19	7.379	1017	1032	1044	rBV4	316502	914887	3.03%	0.448%
20	7.988	1123	1132	1143	rVB	181133	418859	1.38%	0.205%
21	8.110	1143	1152	1155	rBV	334026	700110	2.31%	0.343%
22	8.147	1155	1158	1163	rVB	369700	560653	1.85%	0.274%
23	8.507	1210	1217	1227	rVB	439112	759678	2.51%	0.372%
24	8.653	1235	1241	1247	rVB	907094	1426234	4.72%	0.698%
25	8.720	1247	1252	1258	rVB	281070	431131	1.43%	0.211%
26	9.458	1364	1373	1377	rBV4	143003	328062	1.08%	0.160%
27	10.055	1467	1471	1476	rVB	954814	1242293	4.11%	0.608%
28	10.201	1489	1495	1500	rBV	23112613	30243962	100.00%	14.796%
29	10.305	1506	1512	1518	rVB	20707062	28246104	93.39%	13.819%
30	10.646	1562	1568	1578	rVB	5642222	6712989	22.20%	3.284%
31	10.963	1613	1620	1625	rBV	1822436	2221340	7.34%	1.087%
32	11.085	1635	1640	1644	rBV	964558	1204557	3.98%	0.589%
33	11.305	1672	1676	1683	rVB	3882929	4701682	15.55%	2.300%
34	11.402	1684	1692	1696	rVV	4312733	5445859	18.01%	2.664%
35	11.457	1696	1701	1707	rVB	2459514	3159784	10.45%	1.546%
36	11.616	1721	1727	1735	rBV	5599481	6649382	21.99%	3.253%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

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\82X041222W.M
Title : SW846 8260

37	11.756	1745	1750	1755	rVB	10669509	12459739	41.20%	6.096%
38	11.866	1763	1768	1770	rBV	226128	310604	1.03%	0.152%
39	11.890	1770	1772	1779	rVB	352060	430021	1.42%	0.210%
40	12.024	1789	1794	1799	rBV	1592862	1966456	6.50%	0.962%
41	12.091	1799	1805	1809	rBV	5562081	6755043	22.34%	3.305%
42	12.250	1825	1831	1838	rBV	21385951	27282900	90.21%	13.348%
43	12.317	1838	1842	1850	rVB2	1447974	2100778	6.95%	1.028%
44	12.414	1852	1858	1862	rBV2	1658584	2196635	7.26%	1.075%
45	12.469	1862	1867	1872	rVB2	562479	800767	2.65%	0.392%
46	12.530	1872	1877	1880	rBV	703505	945331	3.13%	0.462%
47	12.616	1886	1891	1894	rBV	4466639	5258777	17.39%	2.573%
48	12.646	1894	1896	1900	rVV	734963	818261	2.71%	0.400%
49	12.701	1900	1905	1912	rVB	3008032	3940628	13.03%	1.928%
50	12.847	1924	1929	1934	rVB2	358821	488845	1.62%	0.239%
51	12.920	1934	1941	1945	rBV	1569347	2028725	6.71%	0.993%
52	12.963	1945	1948	1954	rVV	1023837	1228369	4.06%	0.601%
53	13.024	1954	1958	1967	rVB2	221836	356698	1.18%	0.175%
54	13.170	1978	1982	1992	rVB	2080487	2583654	8.54%	1.264%
55	13.292	1997	2002	2007	rVV2	3260346	4324909	14.30%	2.116%
56	13.341	2007	2010	2019	rVV3	622341	954217	3.16%	0.467%
57	13.426	2019	2024	2030	rVB	1205851	1481734	4.90%	0.725%
58	13.542	2040	2043	2047	rBV	284451	376384	1.24%	0.184%
59	13.585	2047	2050	2054	rVV3	251302	368281	1.22%	0.180%
60	13.640	2054	2059	2061	rVV2	231084	335635	1.11%	0.164%
61	13.664	2061	2063	2070	rVB2	324424	457225	1.51%	0.224%
62	13.780	2073	2082	2091	rBV	9709922	11974743	39.59%	5.858%
63	13.871	2091	2097	2101	rBV	479688	596341	1.97%	0.292%
64	14.640	2219	2223	2233	rVB	1425449	1791122	5.92%	0.876%
65	14.780	2241	2246	2261	rVB2	984154	1335351	4.42%	0.653%

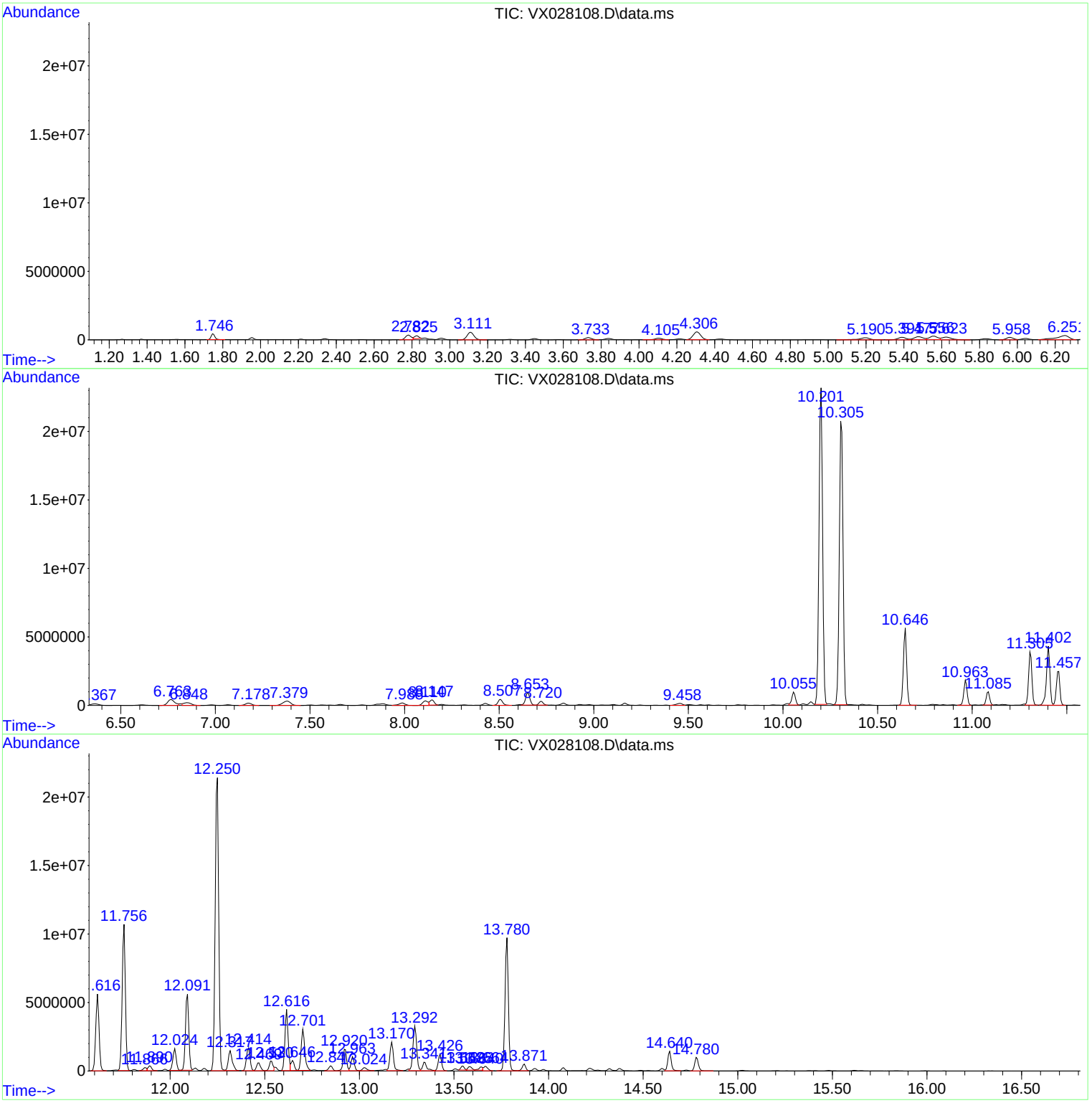
Sum of corrected areas: 204400301

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

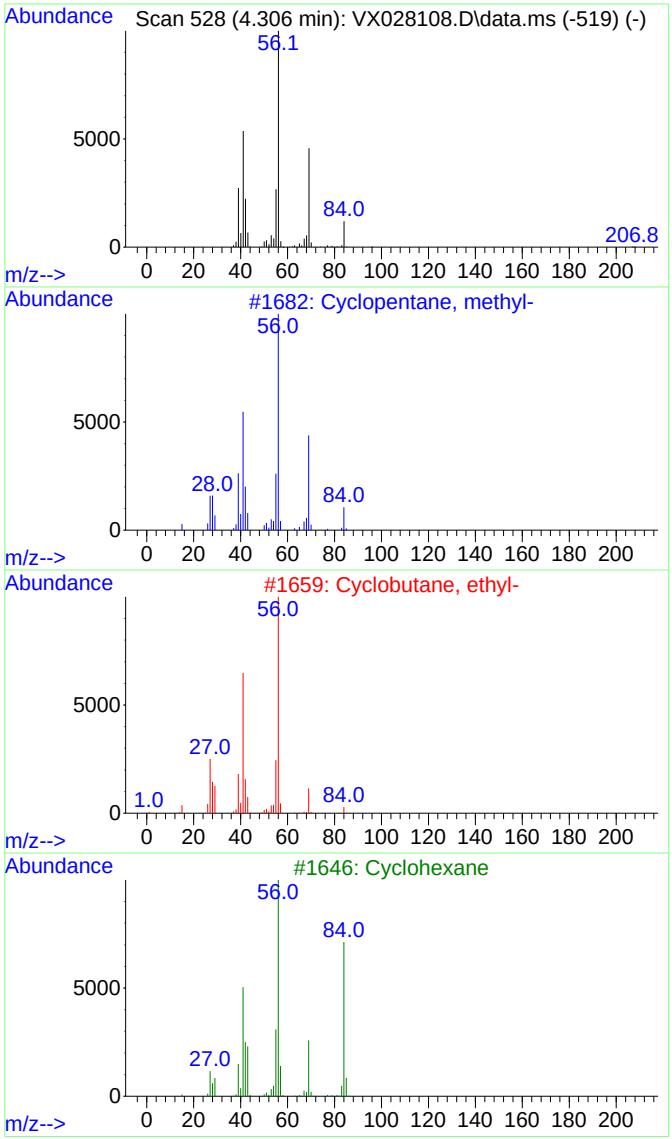
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

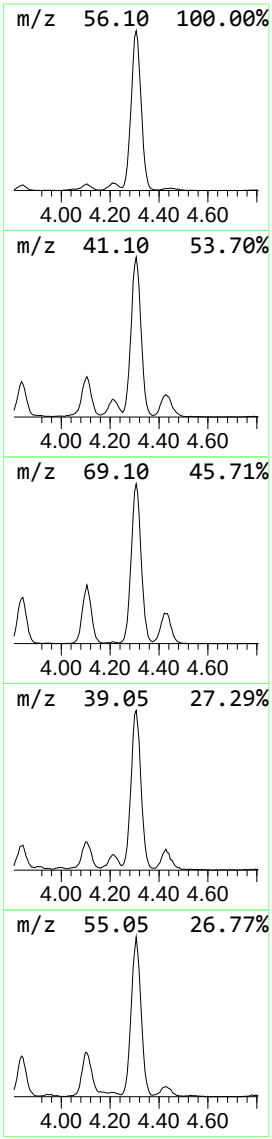
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	103.22 ug/l	1628180	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3	Cyclohexane	84	C6H12	000110-82-7	64
4	1-Hexene	84	C6H12	000592-41-6	50
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

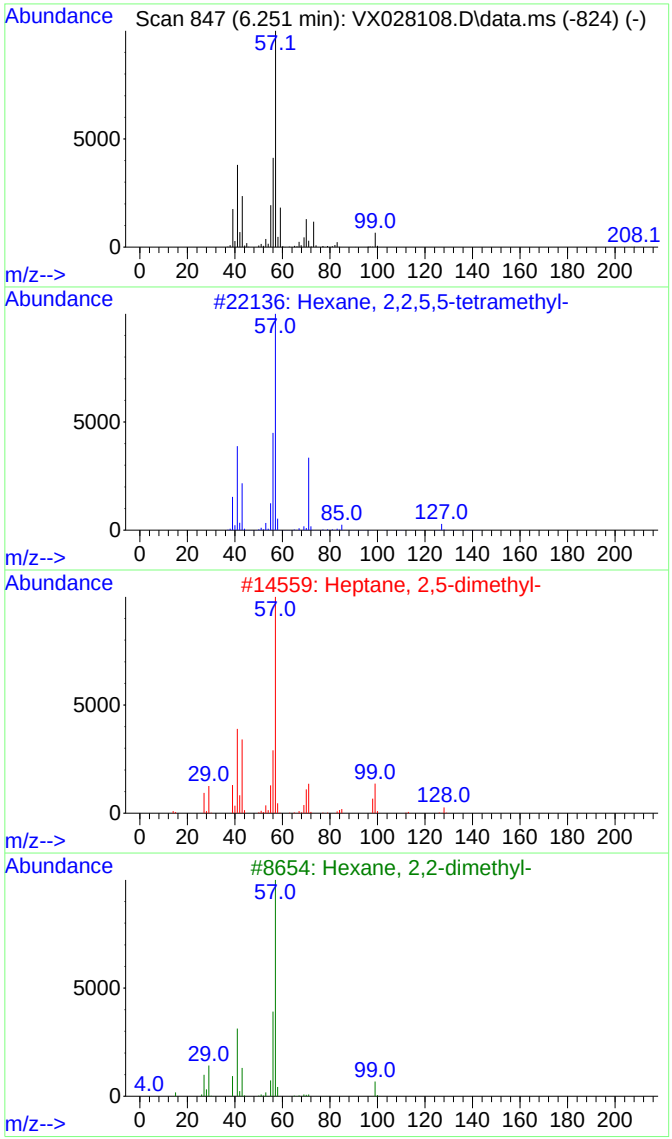
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

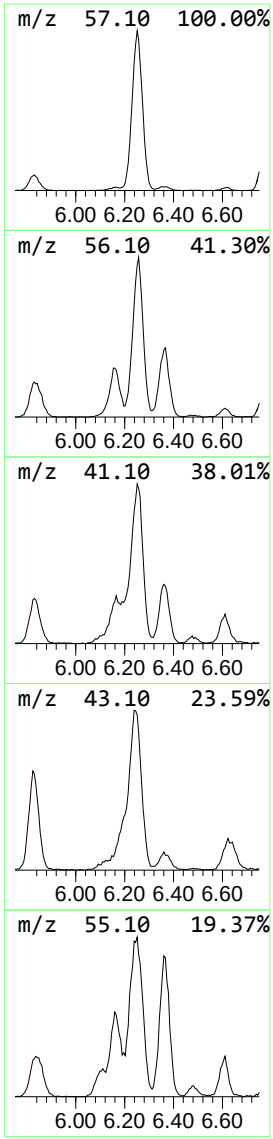
****

Peak Number 2 unknown6.251 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	64.44 ug/l	1450190	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	47
2	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	47
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	42
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	40
5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

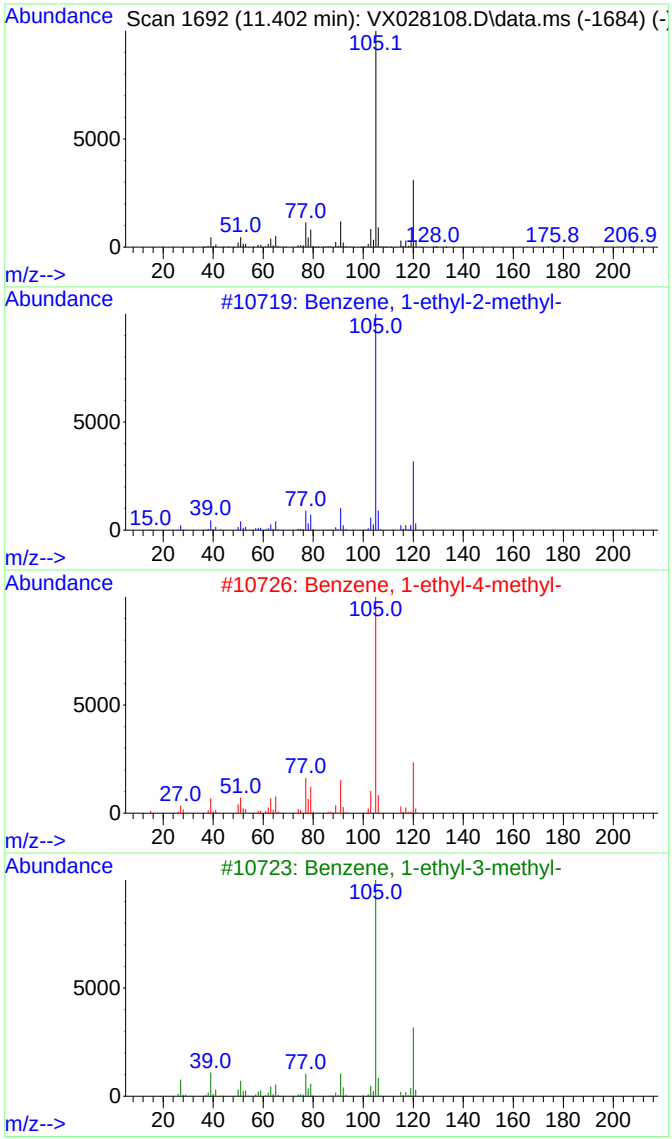
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

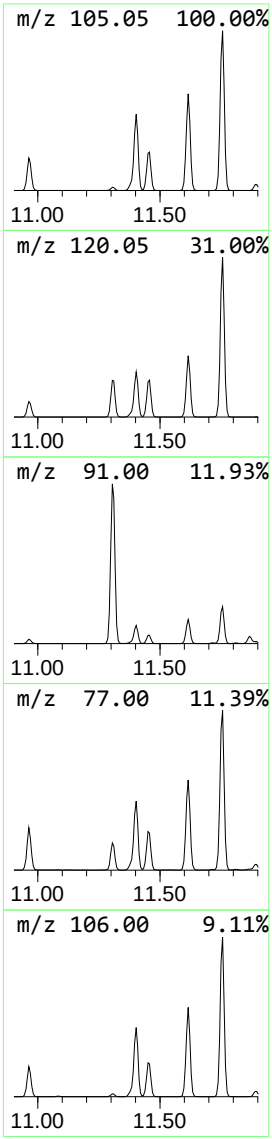
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.402	138.47 ug/l	5445860	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	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

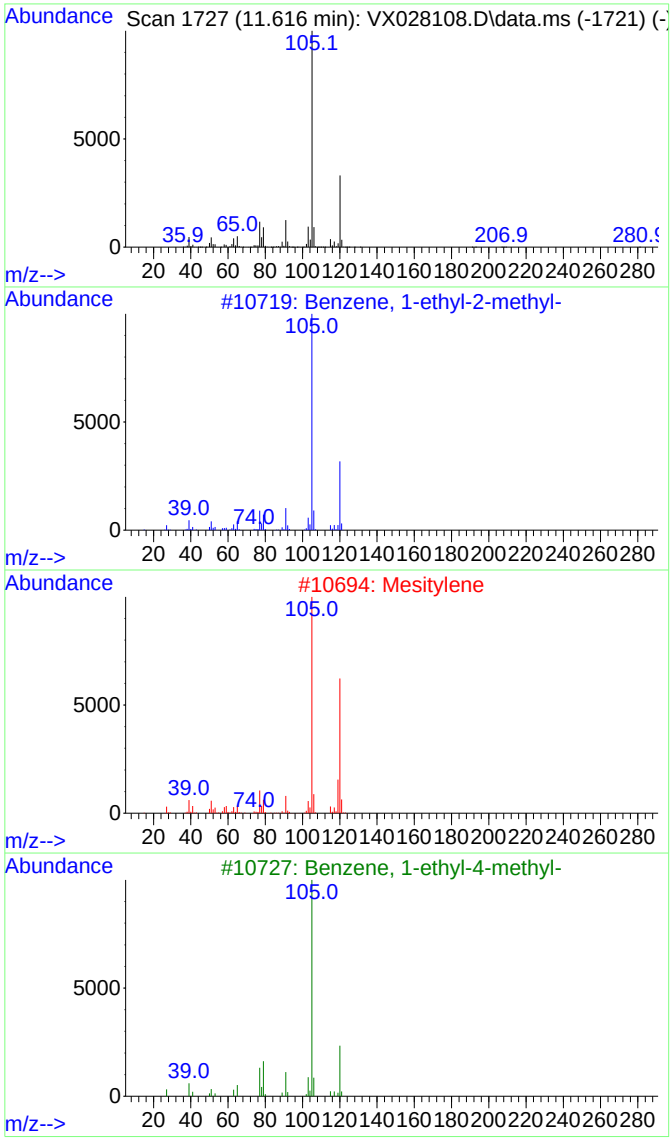
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

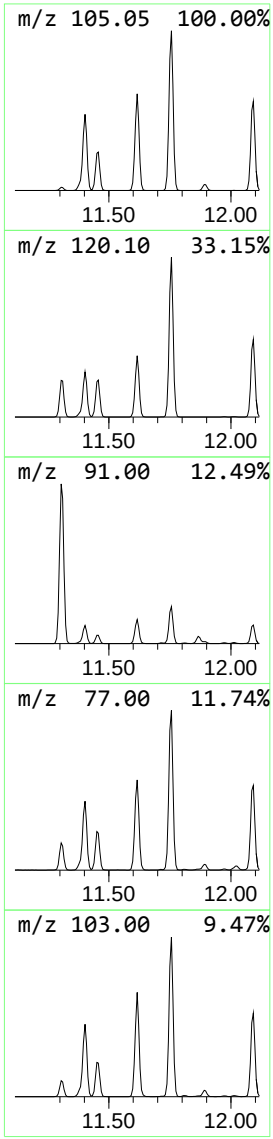
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	169.07 ug/l	6649380	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	Mesitylene	120	C9H12	000108-67-8	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	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

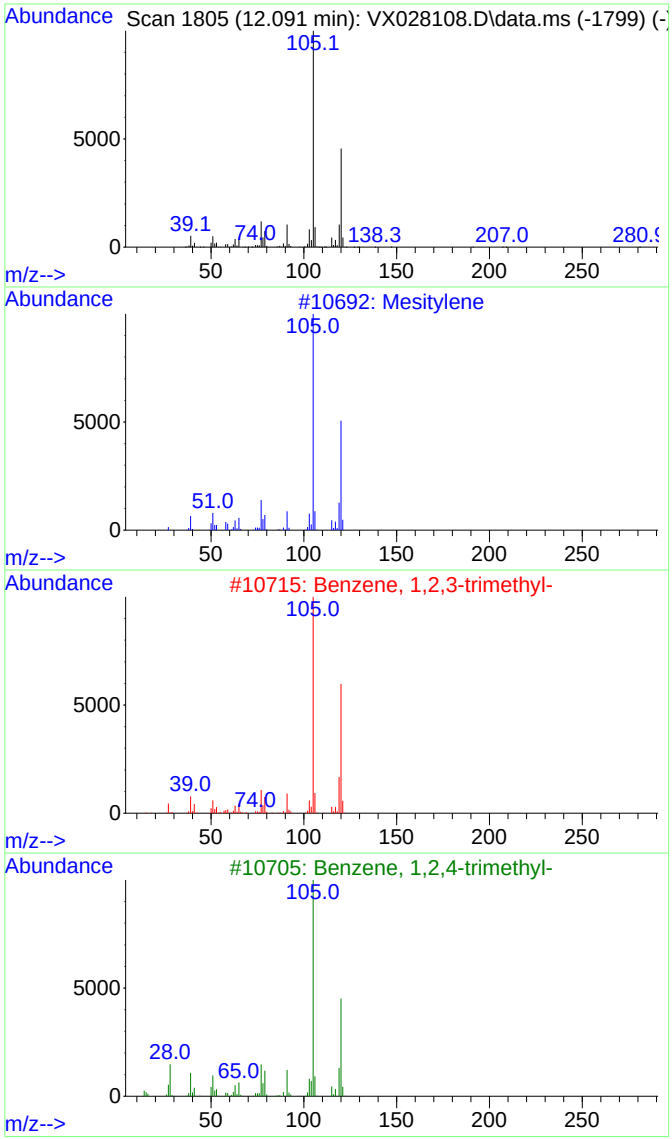
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

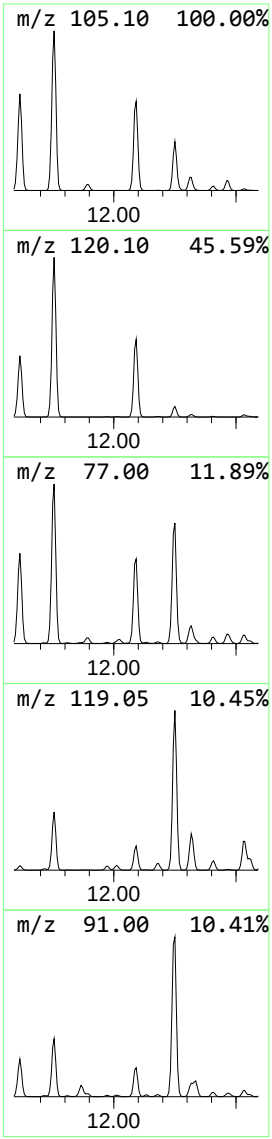
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	171.76 ug/l	6755040	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\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

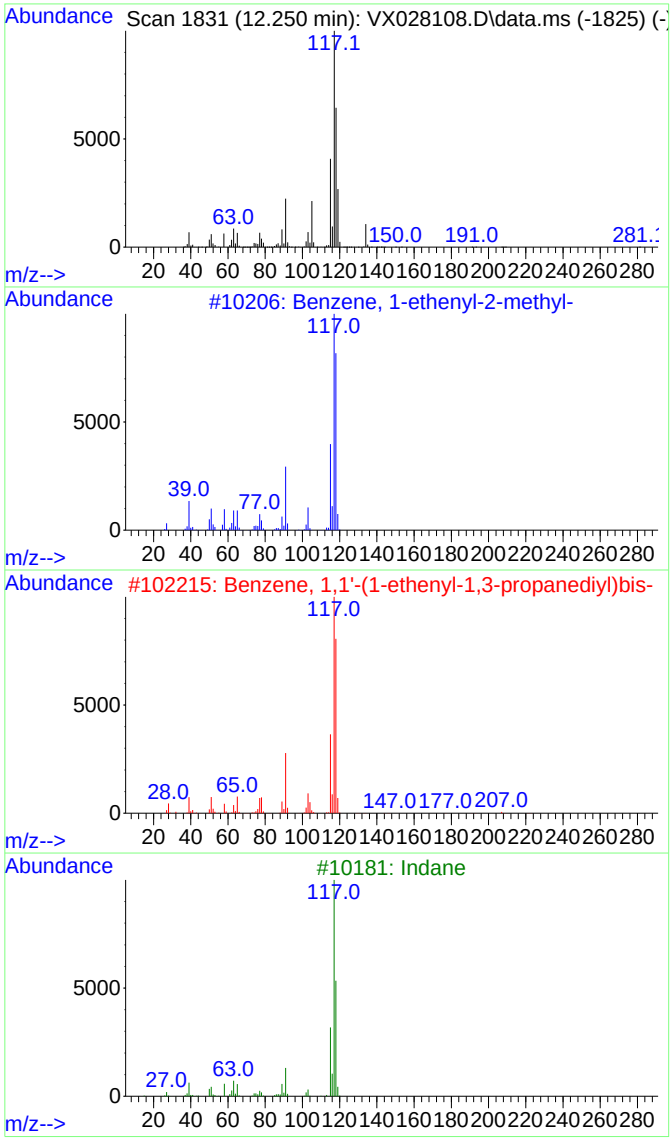
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

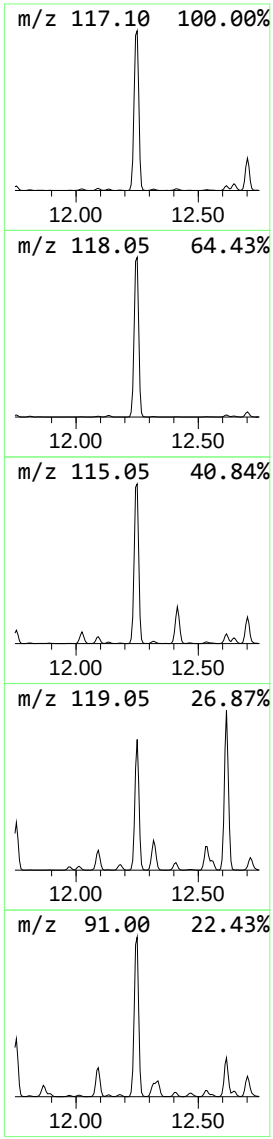
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	693.71 ug/l	27282900	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83
2	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
3	Indane	118	C9H10	000496-11-7	60
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

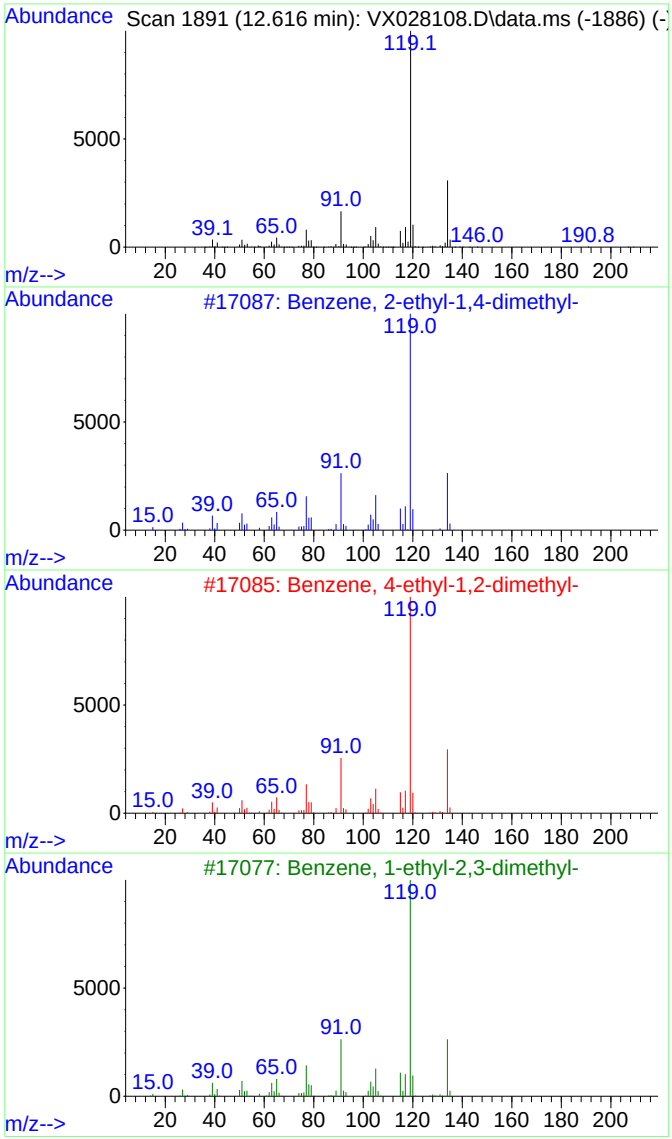
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

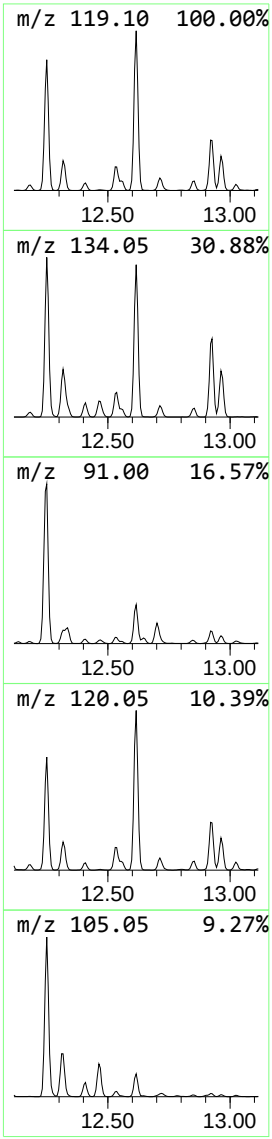
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	133.71 ug/l	5258780	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

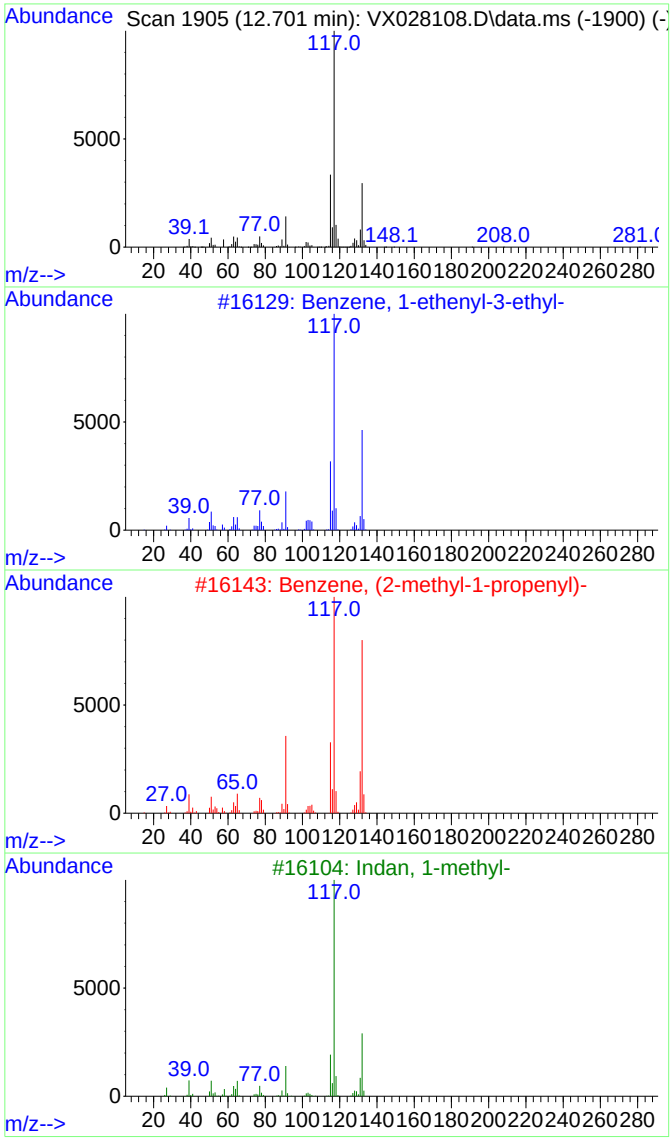
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

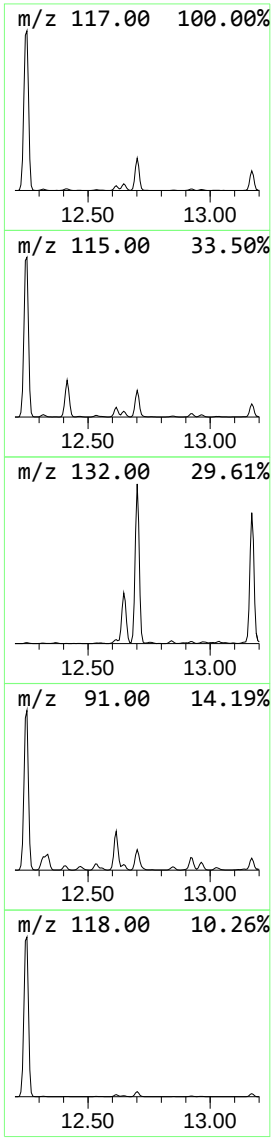
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	100.20 ug/l	3940630	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	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

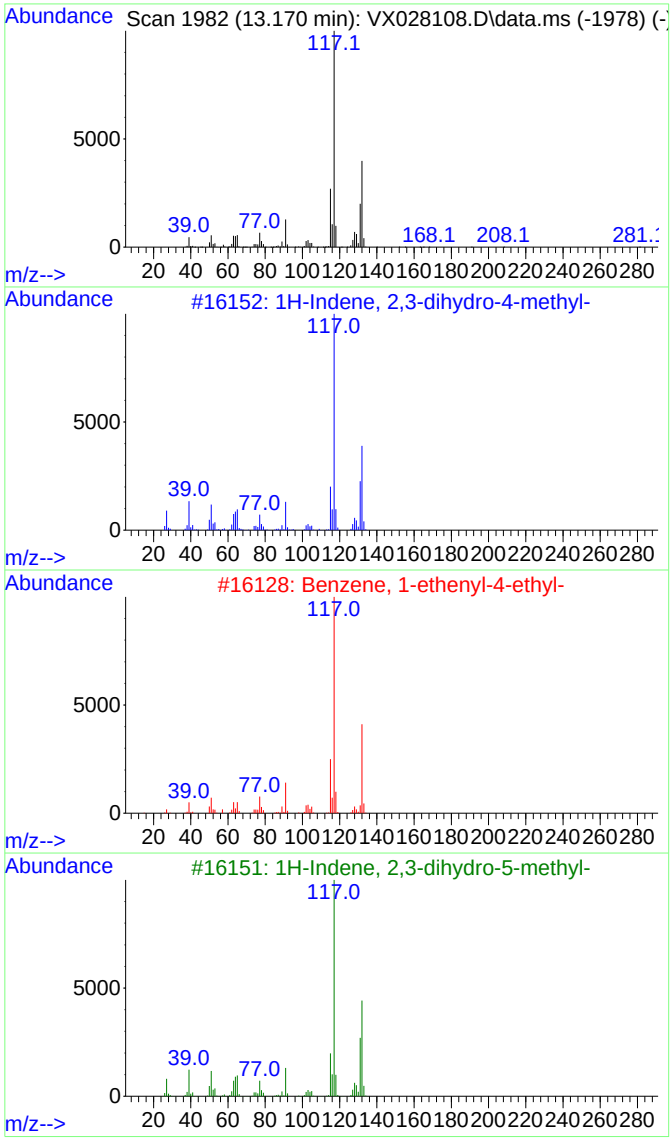
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

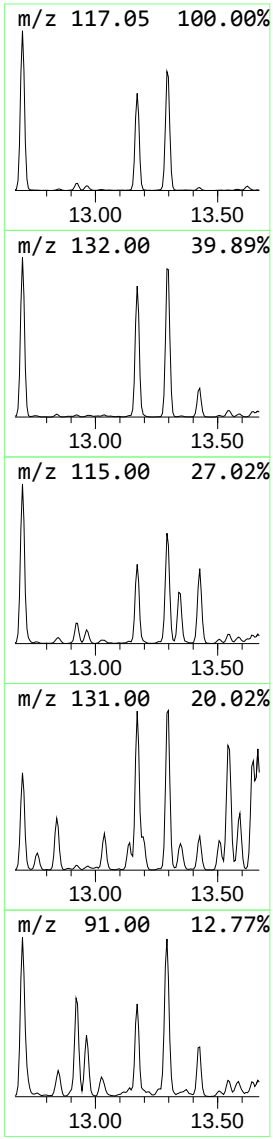
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	65.69 ug/l	2583650	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	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

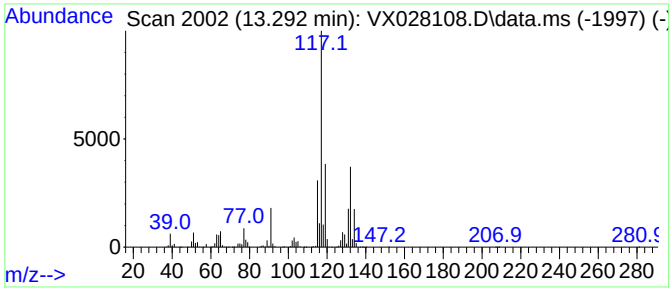
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

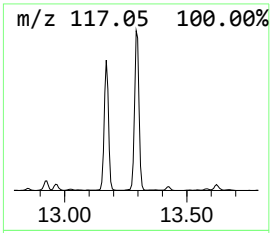
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

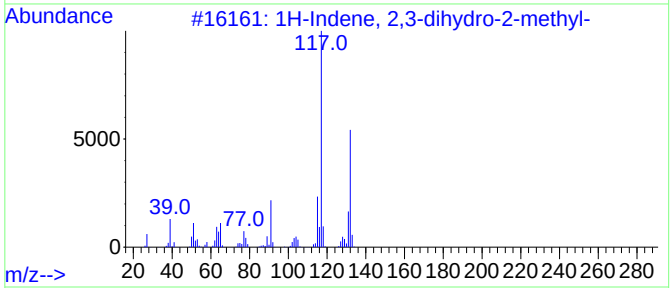
****
Peak Number 10 1H-Indene, 2,3-dihydro-2-me... Concentration Rank 6

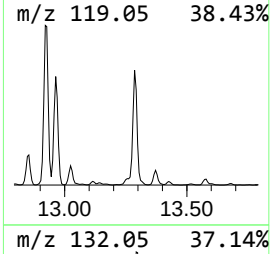
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	109.97 ug/l	4324910	1,4-Dichlorobenzene-d4	12.024

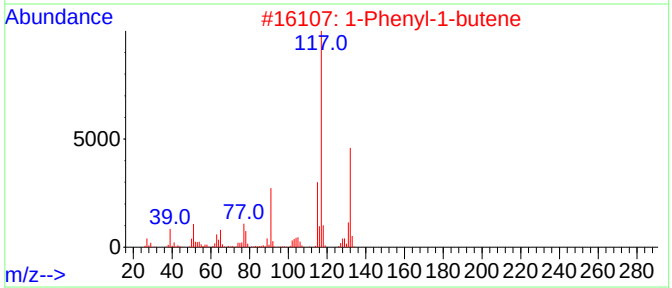
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	76
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	76
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	76
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76

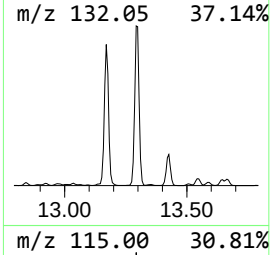


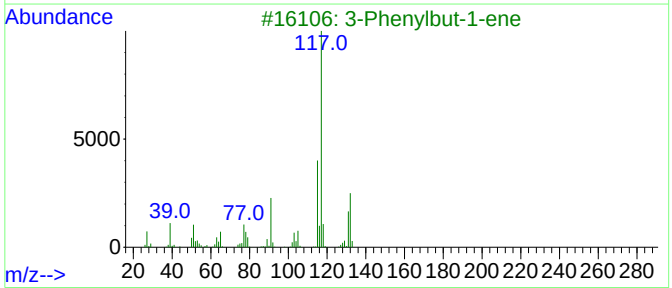


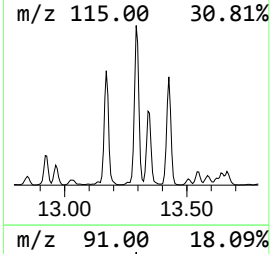


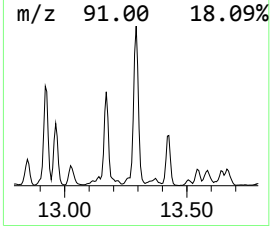












Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028108.D
Acq On : 14 Apr 2022 15:38
Operator : JC/MD
Sample : N2403-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-47

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.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	103.2	ug/l	1628180	1	5.556	788702	50.0
unknown6.251	6.251	64.4	ug/l	1450190	2	6.763	1125150	50.0
Benzene, 1-ethy...	11.402	138.5	ug/l	5445860	4	12.024	1966460	50.0
Benzene, 1-ethy...	11.616	169.1	ug/l	6649380	4	12.024	1966460	50.0
Benzene, 1,2,3-...	12.091	171.8	ug/l	6755040	4	12.024	1966460	50.0
Benzene, 1-ethe...	12.250	693.7	ug/l	27282900	4	12.024	1966460	50.0
Benzene, 2-ethy...	12.616	133.7	ug/l	5258780	4	12.024	1966460	50.0
Benzene, 1-ethe...	12.701	100.2	ug/l	3940630	4	12.024	1966460	50.0
1H-Indene, 2,3-...	13.170	65.7	ug/l	2583650	4	12.024	1966460	50.0
1H-Indene, 2,3-...	13.292	110.0	ug/l	4324910	4	12.024	1966460	50.0