LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

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\82X072123W.M
Title : SW846 8260

Signal : TIC: VX036632.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.605	76	85	94	rBV	5871	18651	2.01%	0.135%
2	2.386	207	213	223	rVB	27101	46902	5.04%	0.339%
3	2.550	236	240	250	rVB	6626	16415	1.77%	0.119%
4	4.574	564	572	581	rBV2	3930	11599	1.25%	0.084%
5	5.391	692	706	721	rBV	104210	308647	33.19%	2.229%
6	5.556	721	733	746	rVB	178408	504769	54.28%	3.645%
7	5.958	789	799	817	rBV	115373	306111	32.92%	2.210%
8	6.763	921	931	944	rBV	290590	684909	73.65%	4.945%
9	8.574	1223	1228	1233	rBV	9038	14593	1.57%	0.105%
10	8.647	1234	1240	1248	rVV	586859	929923	100.00%	6.714%
11	8.720	1248	1252	1259	rVB2	6540	11457	1.23%	0.083%
12	9.519	1378	1383	1388	rBV	15608	23300	2.51%	0.168%
13	10.055	1465	1471	1488	rVB	604685	812666	87.39%	5.868%
14	10.195	1488	1494	1502	rBV	18231	26212	2.82%	0.189%
15	10.299	1505	1511	1519	rBV	95098	130895	14.08%	0.945%
16	10.640	1562	1567	1583	rVB	109807	146485	15.75%	1.058%
17	10.854	1598	1602	1612	rVV2	15984	22223	2.39%	0.160%
18	10.963	1615	1620	1625	rBV	17587	22745	2.45%	0.164%
19	11.079	1634	1639	1652	rBV	474168	612207	65.83%	4.420%
20	11.305	1670	1676	1682	rBV	34521	42894	4.61%	0.310%
21	11.378	1682	1688	1696	rVV	189728	331656	35.66%	2.395%
22	11.451	1696	1700	1710	rVB	80034	99745	10.73%	0.720%
23	11.610	1722	1726	1733	rVB	161465	203192	21.85%	1.467%
24	11.750	1744	1749	1758	rVB	318406	388144	41.74%	2.803%
25	11.835	1758	1763	1766	rBV3	16263	27253	2.93%	0.197%
26	11.890	1766	1772	1776	rVB2	25979	38827	4.18%	0.280%
27	11.969	1781	1785	1788	rBV	44432	50808	5.46%	0.367%
28	12.018	1788	1793	1799	rVV	585945	745665	80.19%	5.384%
29	12.085	1799	1804	1812	rVB	272952	340979	36.67%	2.462%
30	12.244	1822	1830	1833	rBV2	274856	416321	44.77%	3.006%
31	12.317	1838	1842	1851	rVB4	121633	189133	20.34%	1.366%
32	12.402	1851	1856	1861	rBV	26376	35893	3.86%	0.259%
33	12.463	1861	1866	1871	rVB	79663	96731	10.40%	0.698%
34	12.530	1871	1877	1879	rBV	85450	109893	11.82%	0.793%
35	12.555	1879	1881	1885	rVB	72251	74712	8.03%	0.539%
36	12.609	1885	1890	1893	rBV	131306	155759	16.75%	1.125%

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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

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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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Title : SW846 8260

37	12.646	1893	1896	1900	rVB	51695	59770	6.43%	0.432%
38	12.695	1900	1904	1913	rVB2	150039	244358	26.28%	1.764%
39	12.798	1917	1921	1925	rVB	26658	33069	3.56%	0.239%
40	12.847	1925	1929	1934	rVB2	74373	97846	10.52%	0.706%
41	12.920	1934	1941	1944	rBV	100296	119906	12.89%	0.866%
42	12.963	1944	1948	1954	rVB	97339	121270	13.04%	0.876%
43	13.024	1954	1958	1960	rBV2	24344	33601	3.61%	0.243%
44	13.134	1973	1976	1978	rVV2	24044	30430	3.27%	0.220%
45	13.164	1978	1981	1986	rVV	133760	173318	18.64%	1.251%
46	13.213	1986	1989	1992	rVB	22972	26385	2.84%	0.191%
47	13.256	1992	1996	1997	rBV2	22818	29517	3.17%	0.213%
48	13.292	1997	2002	2007	rVV2	491830	657838	70.74%	4.750%
49	13.329	2007	2008	2012	rVV3	13823	17847	1.92%	0.129%
50	13.365	2012	2014	2017	rVB	13616	13901	1.49%	0.100%
51	13.420	2017	2023	2033	rVB	581645	696338	74.88%	5.028%
52	13.506	2033	2037	2040	rBV2	27392	38686	4.16%	0.279%
53	13.542	2040	2043	2046	rVV	44790	58268	6.27%	0.421%
54	13.585	2046	2050	2054	rVV3	84378	138688	14.91%	1.001%
55	13.658	2056	2062	2068	rVV2	87835	191742	20.62%	1.384%
56	13.713	2068	2071	2074	rVV2	20076	25917	2.79%	0.187%
57	13.774	2074	2081	2086	rVV	240685	324048	34.85%	2.340%
58	13.823	2086	2089	2090	rVV2	15372	18634	2.00%	0.135%
59	13.853	2090	2094	2098	rVB	145624	186605	20.07%	1.347%
60	13.926	2098	2106	2110	rBV	163598	224509	24.14%	1.621%
61	13.969	2110	2113	2117	rVV2	51198	65606	7.05%	0.474%
62	14.073	2126	2130	2133	rBV	44988	54215	5.83%	0.391%
63	14.109	2133	2136	2141	rVB5	9604	14495	1.56%	0.105%
64	14.225	2148	2155	2161	rBV	215881	354881	38.16%	2.562%
65	14.316	2167	2170	2175	rVB2	30527	35137	3.78%	0.254%
66	14.371	2175	2179	2184	rBV	70334	88160	9.48%	0.637%
67	14.420	2184	2187	2191	rVB2	12250	15785	1.70%	0.114%
68	14.475	2191	2196	2202	rBV2	236346	311677	33.52%	2.250%
69	14.633	2214	2222	2226	rBV2	239734	365799	39.34%	2.641%
70	14.670	2226	2228	2234	rVB3	55699	72046	7.75%	0.520%
71	14.725	2234	2237	2241	rBV2	13326	17872	1.92%	0.129%
72	14.780	2241	2246	2249	rVV	154923	201873	21.71%	1.458%
73	14.816	2249	2252	2255	rVV2	34929	51798	5.57%	0.374%
74	14.847	2255	2257	2260	rVB2	24004	24073	2.59%	0.174%
75	14.902	2260	2266	2272	rVB3	31465	61869	6.65%	0.447%
76	15.048	2283	2290	2297	rBV	16609	28677	3.08%	0.207%

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Integrator: RTE
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Title : SW846 8260

77	15.182	2304	2312	2313	rBV2	36896	52440	5.64%	0.379%
78	15.207	2313	2316	2320	rVV	45272	62039	6.67%	0.448%
79	15.249	2320	2323	2330	rVB3	16422	29993	3.23%	0.217%
80	15.322	2330	2335	2338	rBV6	5925	9340	1.00%	0.067%
81	15.371	2338	2343	2347	rBV	22812	34479	3.71%	0.249%
82	15.475	2353	2360	2365	rBV2	46450	76768	8.26%	0.554%
83	15.524	2365	2368	2373	rVV	12268	21392	2.30%	0.154%
84	15.609	2376	2382	2386	rVV	57801	102237	10.99%	0.738%
85	15.646	2386	2388	2396	rVB	34973	50124	5.39%	0.362%
86	15.841	2411	2420	2426	rVB	15736	36213	3.89%	0.261%
87	15.987	2434	2444	2448	rBV3	7438	15784	1.70%	0.114%
88	16.084	2455	2460	2469	rVV2	10192	18279	1.97%	0.132%
89	16.475	2519	2524	2530	rBV	8899	15768	1.70%	0.114%

Sum of corrected areas: 13849624

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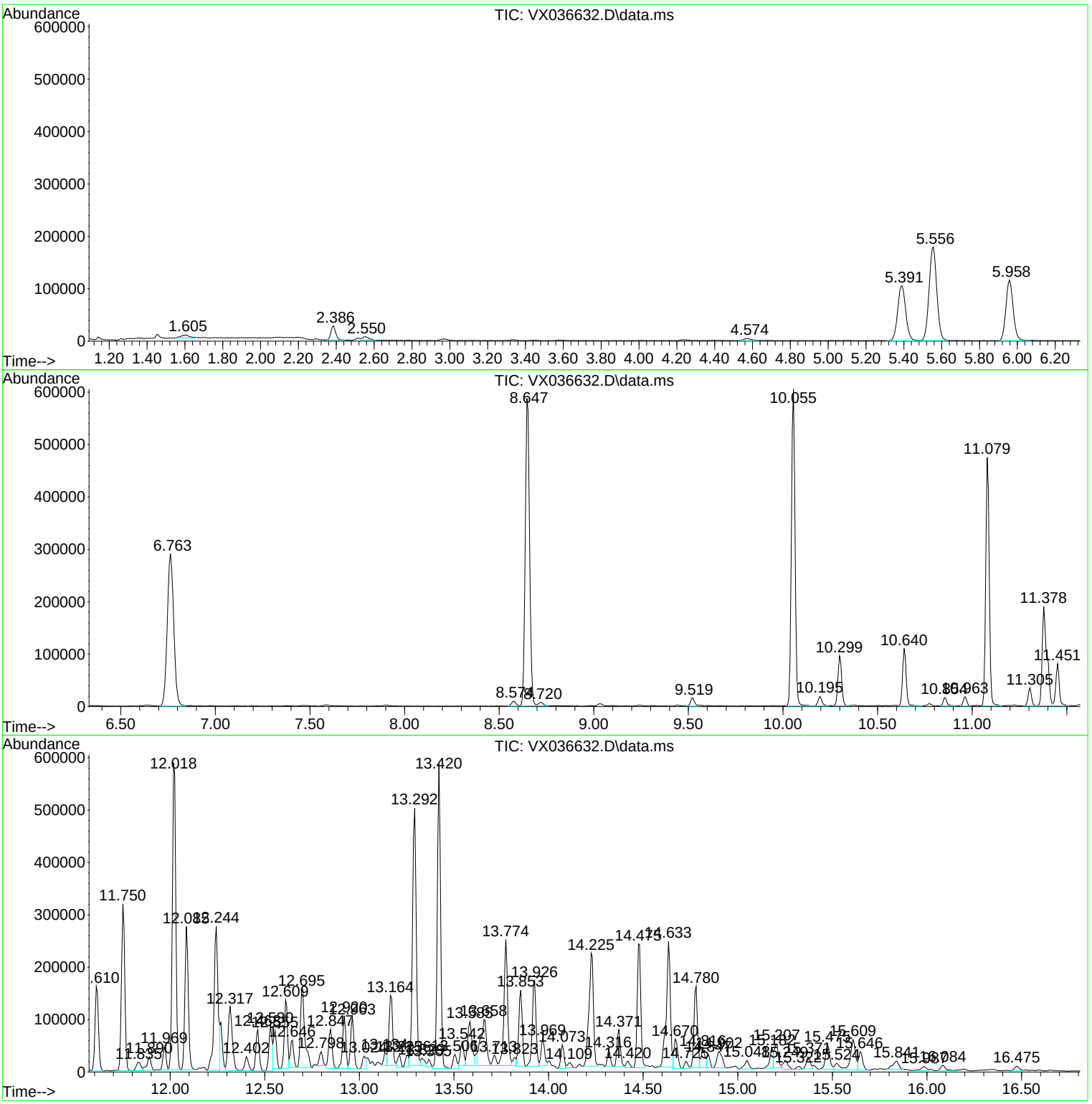
ClientSampleId :
32399

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
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ClientSampleId :
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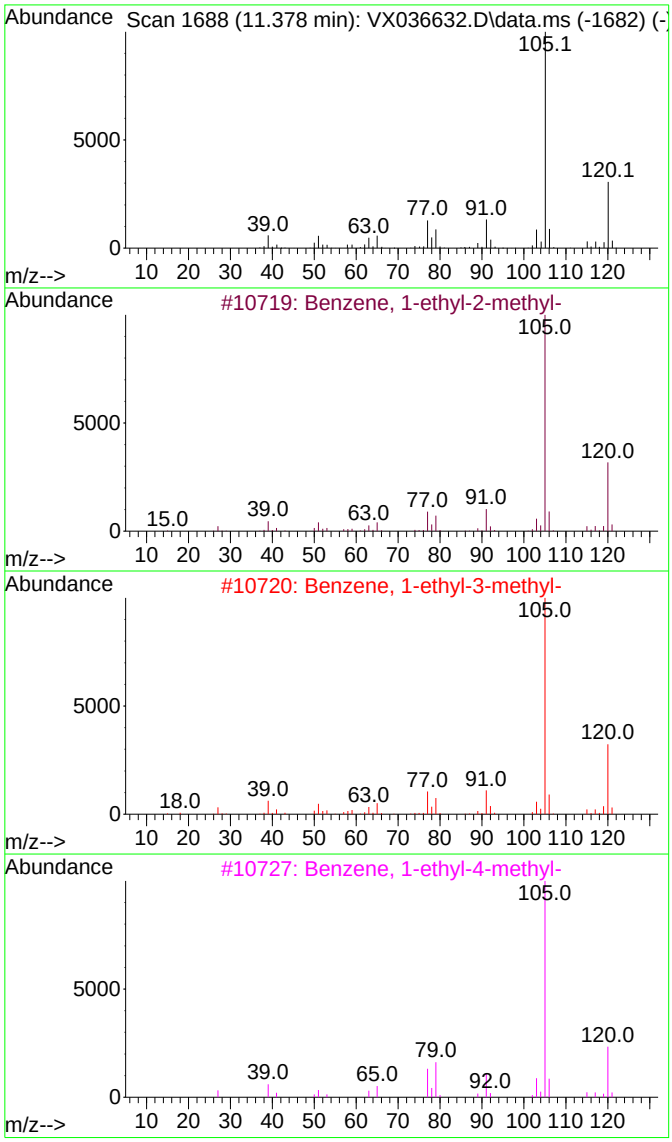
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

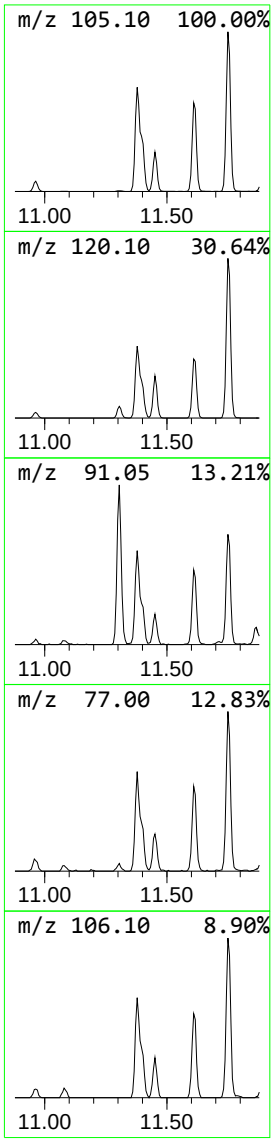
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	22.24 ug/l	331656	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-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

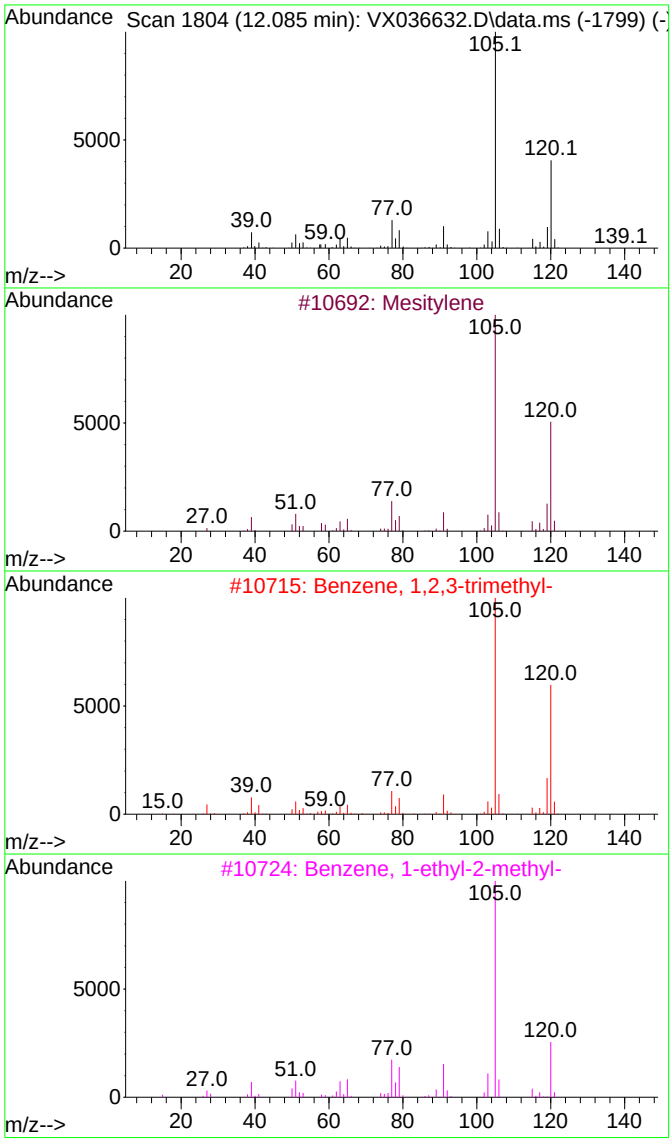
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

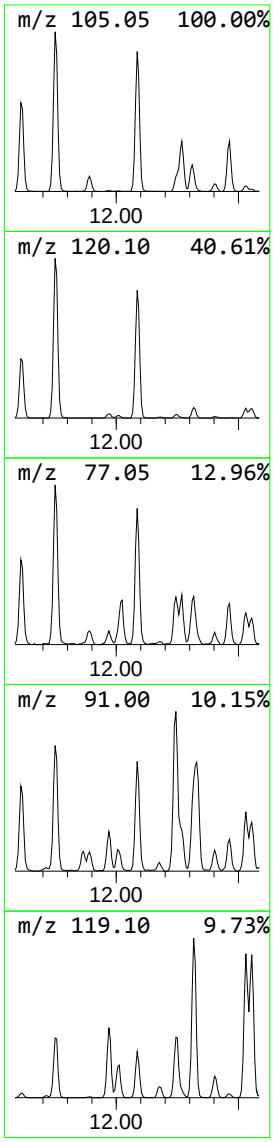
****

Peak Number 2 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	22.86 ug/l	340979	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-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

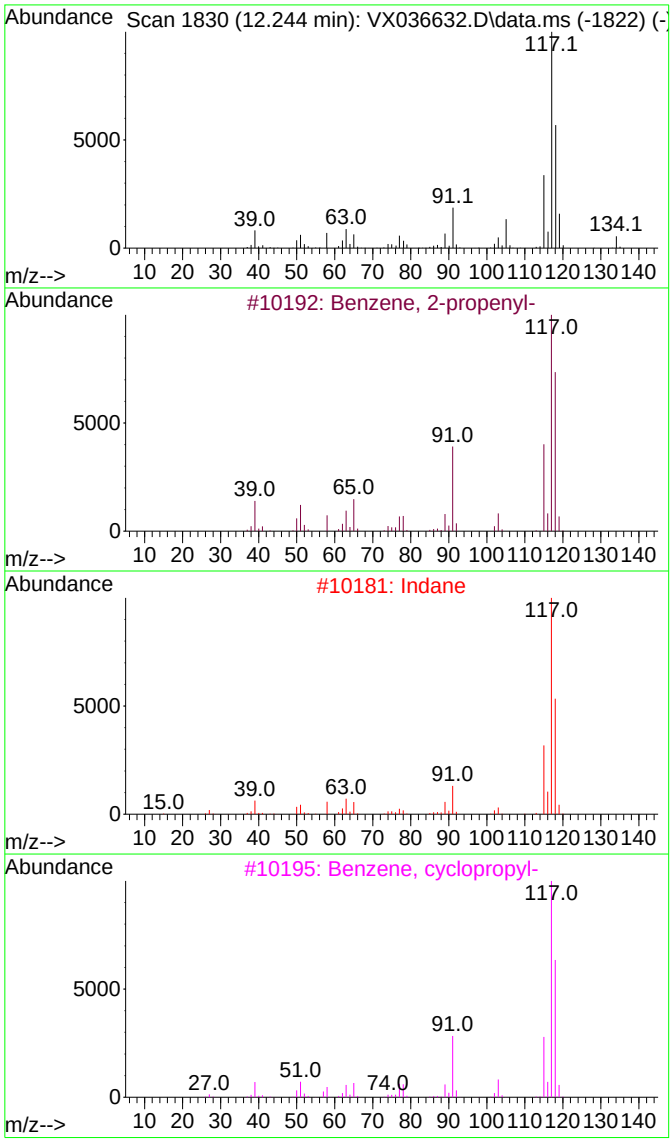
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

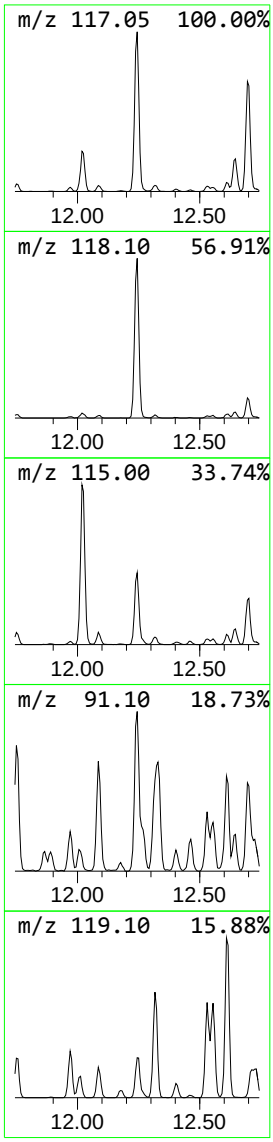
****

Peak Number 3 Benzene, 2-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	27.92 ug/l	416321	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
2	Indane	118	C9H10	000496-11-7	76
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	59
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	55





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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

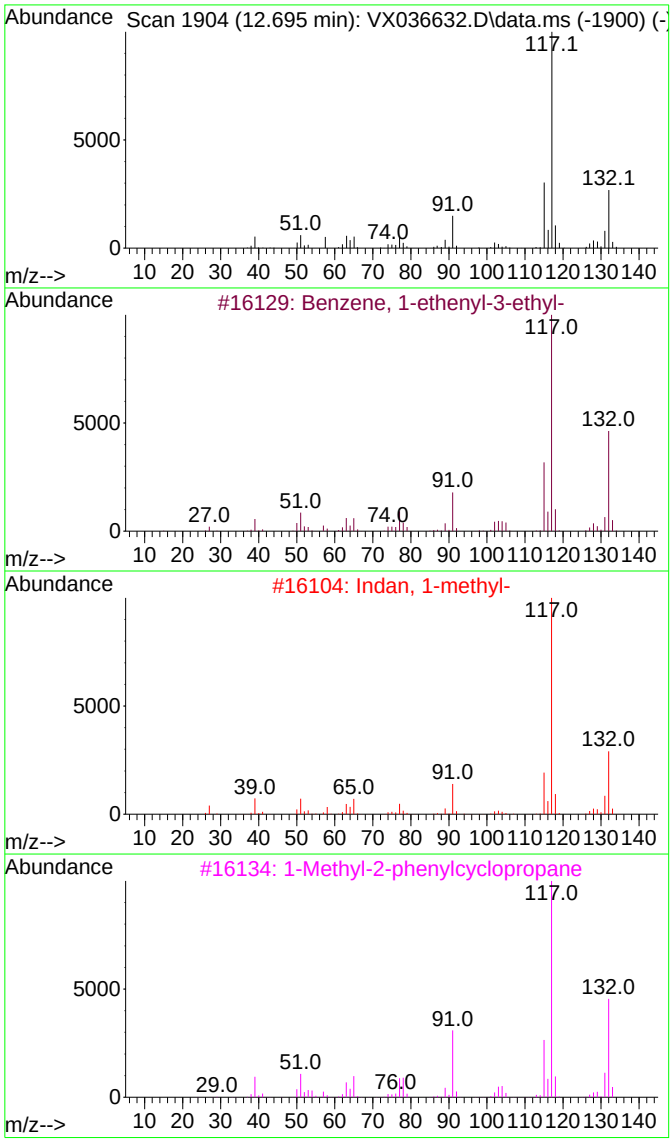
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

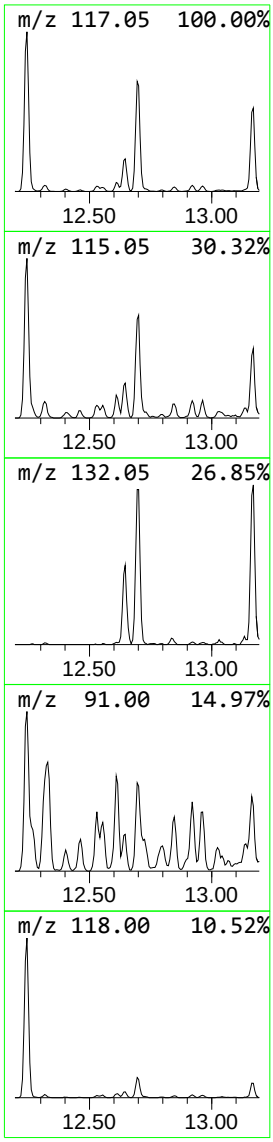
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	16.39 ug/l	244358	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	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83





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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

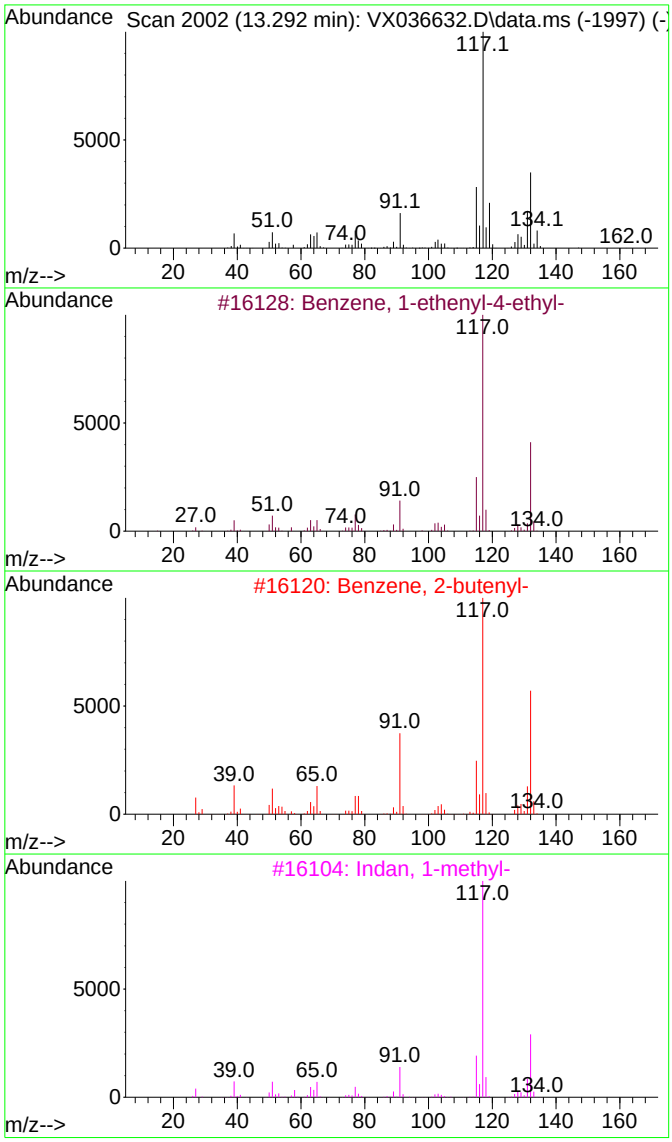
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

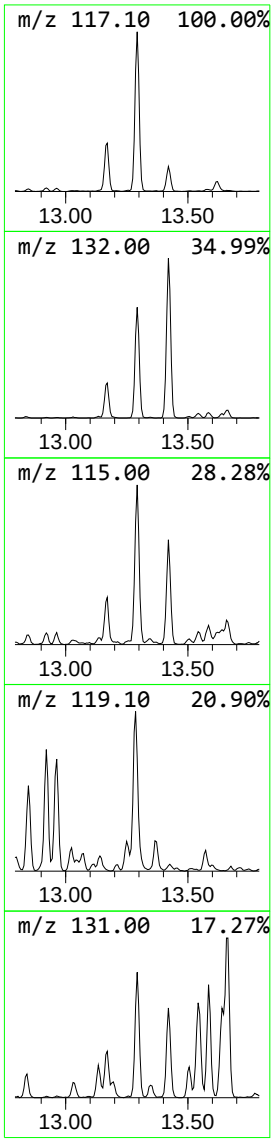
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	44.11 ug/l	657838	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	89
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	81





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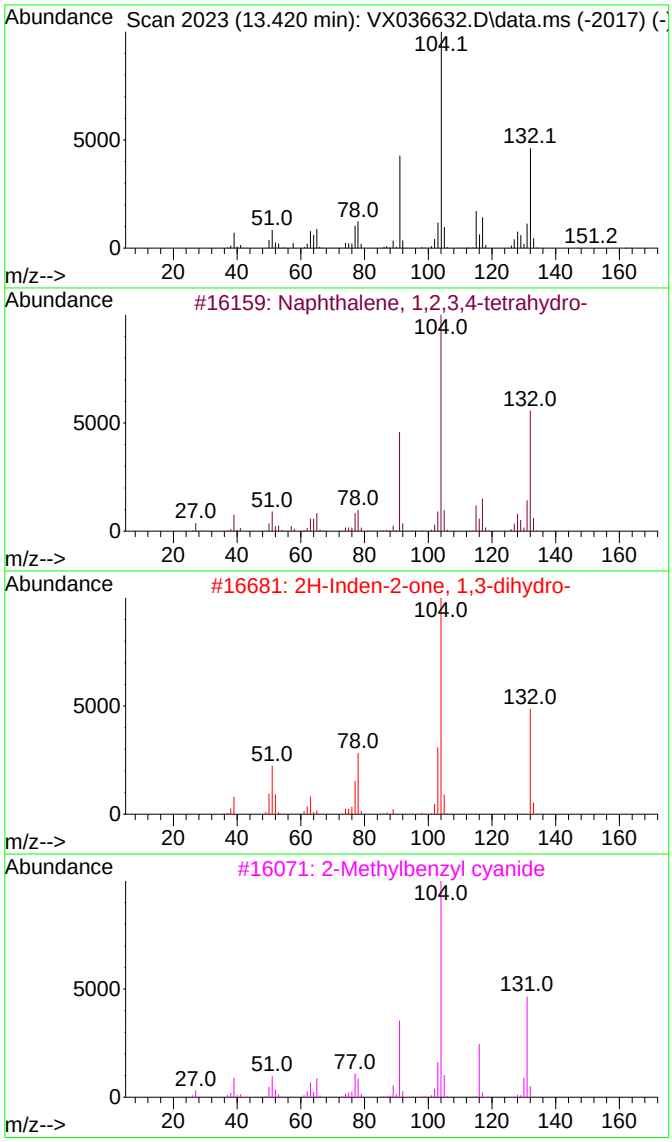
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

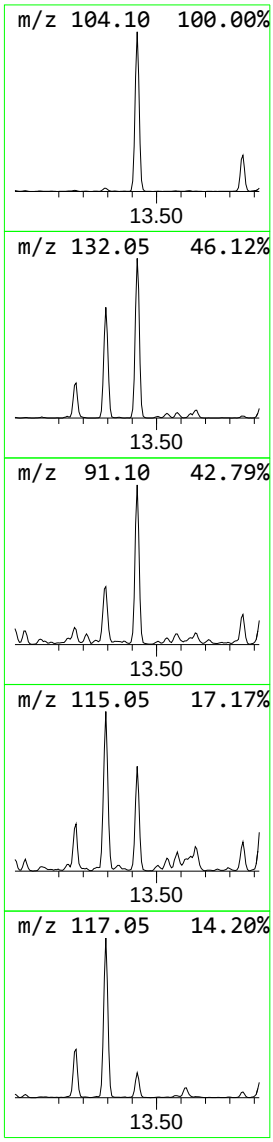
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	46.69 ug/l	696338	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
3	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46
4	Pentafluoropropionic acid, 2-phe...	268	C11H9F5O2	081089-48-7	38
5	3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	38





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Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

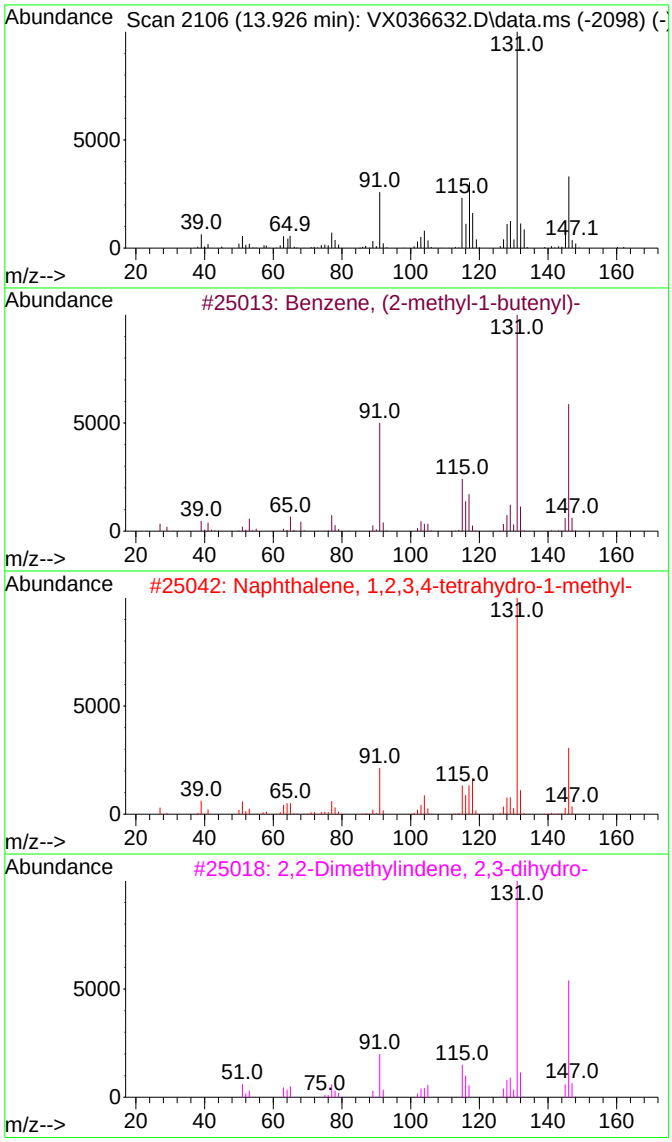
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

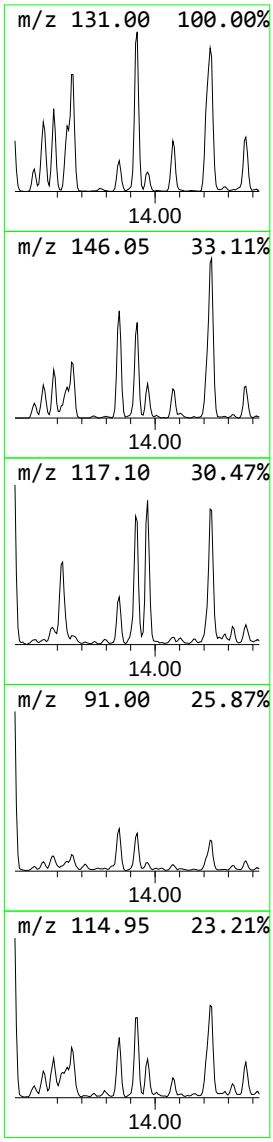
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	15.05 ug/l	224509	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	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	90
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

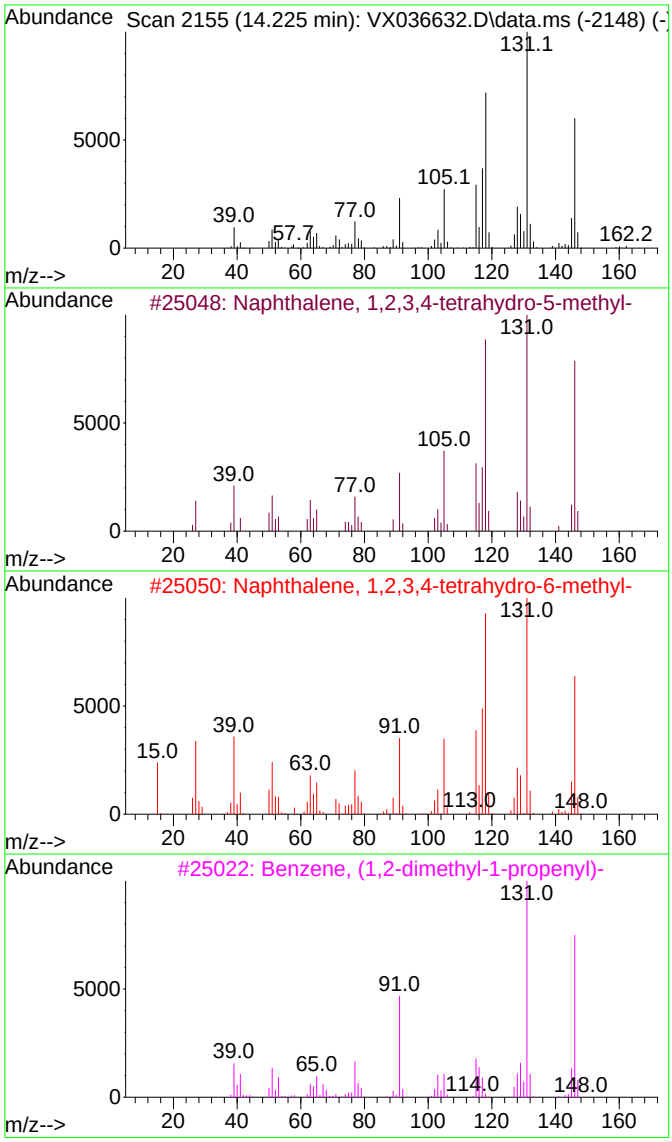
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

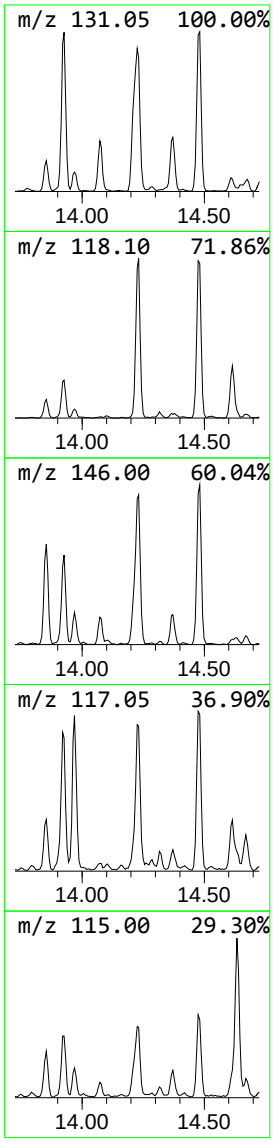
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	23.80 ug/l	354881	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

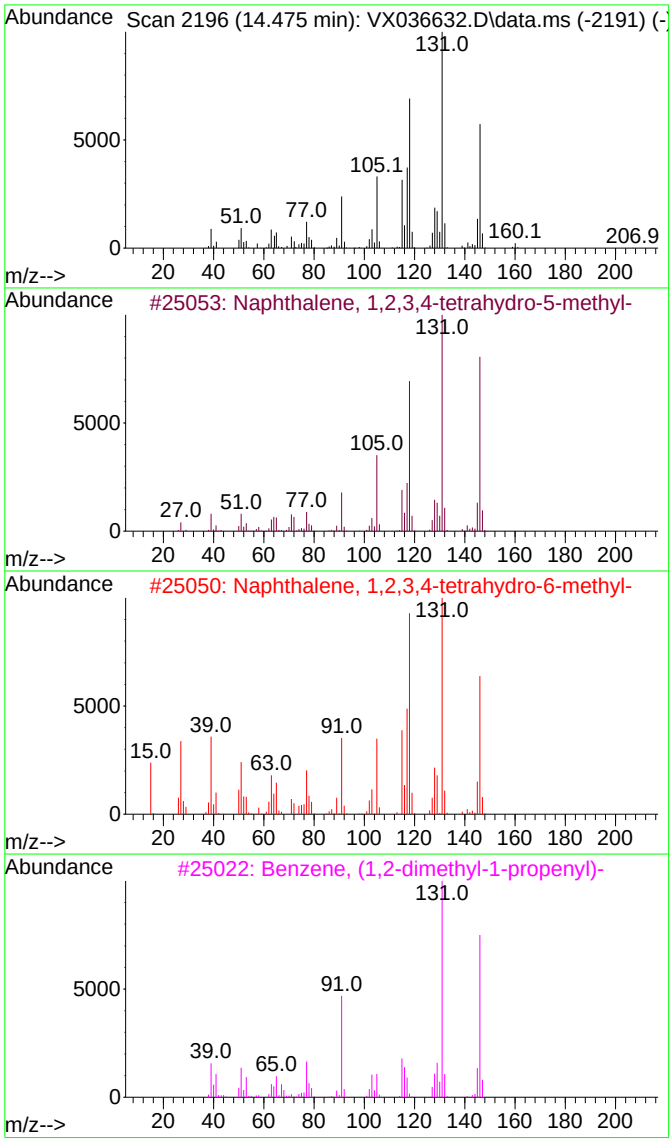
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

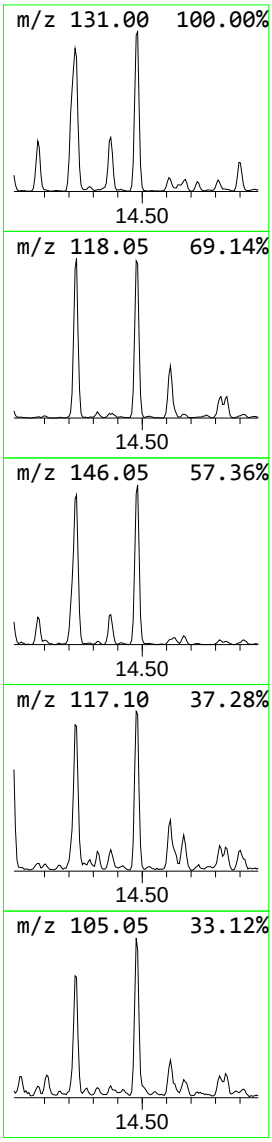
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.475	20.90 ug/l	311677	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

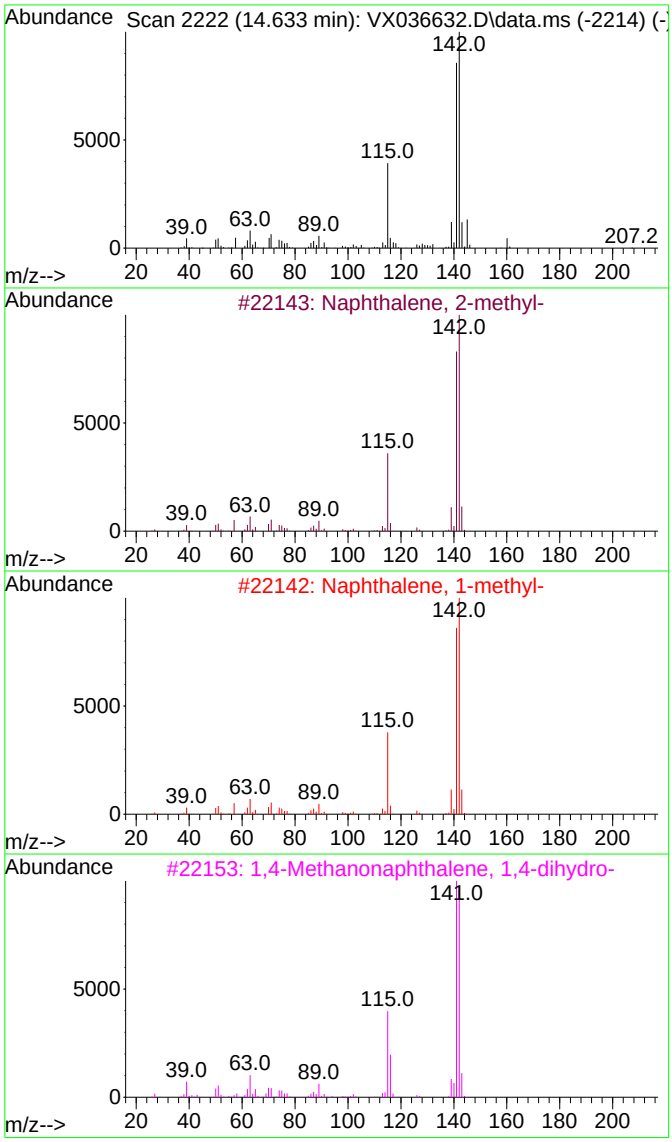
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

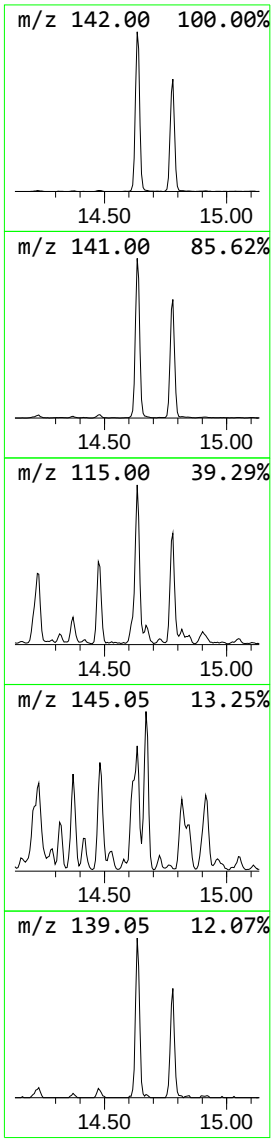
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	24.53 ug/l	365799	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	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	90
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
Data File : VX036632.D
Acq On : 24 Jul 2023 16:09
Operator : JC/MD
Sample : 03722-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
32399

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.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
Benzene, 1-ethy...	11.378	22.2	ug/l	331656	4	12.024	745665	50.0
Mesitylene	12.085	22.9	ug/l	340979	4	12.024	745665	50.0
Benzene, 2-prop...	12.244	27.9	ug/l	416321	4	12.024	745665	50.0
Benzene, 1-ethe...	12.695	16.4	ug/l	244358	4	12.024	745665	50.0
Benzene, 1-ethe...	13.292	44.1	ug/l	657838	4	12.024	745665	50.0
Naphthalene, 1,...	13.420	46.7	ug/l	696338	4	12.024	745665	50.0
Benzene, (2-met...	13.926	15.1	ug/l	224509	4	12.024	745665	50.0
Naphthalene, 1,...	14.225	23.8	ug/l	354881	4	12.024	745665	50.0
Naphthalene, 1,...	14.475	20.9	ug/l	311677	4	12.024	745665	50.0
Naphthalene, 2-...	14.634	24.5	ug/l	365799	4	12.024	745665	50.0