

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
 Data File : VX035635.D
 Acq On : 12 May 2023 14:07
 Operator : JC/MD
 Sample : 02702-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 50033

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\82X051123W.M
 Title : SW846 8260

Signal : TIC: VX035635.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.368	41	46	53	rVV	169792	185266	3.90%	0.341%
2	1.752	104	109	117	rVB	61468	82938	1.74%	0.153%
3	1.953	138	142	153	rVB2	26057	48722	1.02%	0.090%
4	2.215	181	185	195	rVB	32763	54662	1.15%	0.101%
5	3.111	320	332	345	rBV2	59948	141164	2.97%	0.260%
6	4.300	514	527	540	rVB3	17282	51685	1.09%	0.095%
7	5.385	694	705	714	rBV2	129356	378460	7.96%	0.696%
8	5.550	723	732	751	rVB	222987	641361	13.48%	1.179%
9	5.958	788	799	804	rBV	155755	408724	8.59%	0.752%
10	6.038	804	812	826	rVB	601983	1592647	33.49%	2.928%
11	6.757	919	930	948	rBV	358448	876763	18.43%	1.612%
12	7.379	1021	1032	1042	rBV2	36256	84472	1.78%	0.155%
13	8.647	1234	1240	1246	rBV	744285	1144207	24.06%	2.104%
14	8.720	1246	1252	1259	rVB	3090046	4756130	100.00%	8.745%
15	10.055	1465	1471	1479	rBV	769418	1038899	21.84%	1.910%
16	10.195	1488	1494	1499	rBV	352321	463081	9.74%	0.851%
17	10.305	1503	1512	1518	rVV	544088	750693	15.78%	1.380%
18	10.640	1561	1567	1579	rVB	3595842	4664365	98.07%	8.577%
19	10.963	1614	1620	1626	rBV	38934	52009	1.09%	0.096%
20	11.079	1634	1639	1644	rBV	678453	846035	17.79%	1.556%
21	11.378	1683	1688	1690	rBV	796556	1069009	22.48%	1.966%
22	11.451	1696	1700	1709	rVB	1016900	1213852	25.52%	2.232%
23	11.610	1721	1726	1734	rVB	1284299	1623678	34.14%	2.986%
24	11.750	1744	1749	1755	rVB	140417	171156	3.60%	0.315%
25	11.890	1769	1772	1777	rVB	53768	65083	1.37%	0.120%
26	11.969	1777	1785	1789	rBV	219901	271842	5.72%	0.500%
27	12.018	1789	1793	1799	rVB2	841708	1126868	23.69%	2.072%
28	12.085	1799	1804	1810	rBV	1901378	2241273	47.12%	4.121%
29	12.244	1823	1830	1833	rBV	1416696	1931195	40.60%	3.551%
30	12.317	1838	1842	1850	rVB2	793714	1104200	23.22%	2.030%
31	12.402	1850	1856	1861	rBV2	133648	184680	3.88%	0.340%
32	12.463	1861	1866	1872	rVB	399305	481780	10.13%	0.886%
33	12.530	1872	1877	1879	rBV	276265	354244	7.45%	0.651%
34	12.555	1879	1881	1885	rVB	325393	324227	6.82%	0.596%
35	12.610	1885	1890	1894	rBV	782319	1015697	21.36%	1.868%
36	12.646	1894	1896	1900	rVB	268704	271057	5.70%	0.498%

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Integrator: RTE

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37	12.701	1900	1905	1913	rBV2	664485	1103421	23.20%	2.029%
38	12.798	1917	1921	1924	rBV2	62464	75525	1.59%	0.139%
39	12.847	1924	1929	1934	rVB2	344491	431881	9.08%	0.794%
40	12.920	1934	1941	1944	rBV	545543	680591	14.31%	1.251%
41	12.963	1944	1948	1954	rVB	778261	925946	19.47%	1.703%
42	13.024	1954	1958	1960	rBV	124894	174121	3.66%	0.320%
43	13.042	1960	1961	1964	rVV2	94341	105222	2.21%	0.193%
44	13.067	1964	1965	1968	rVB	67567	55960	1.18%	0.103%
45	13.140	1973	1977	1978	rVV2	128607	161670	3.40%	0.297%
46	13.164	1978	1981	1986	rVV	725734	974087	20.48%	1.791%
47	13.213	1986	1989	1992	rVV	118530	141338	2.97%	0.260%
48	13.256	1992	1996	1998	rVV2	152461	234933	4.94%	0.432%
49	13.292	1998	2002	2007	rVV2	2292031	3049347	64.11%	5.607%
50	13.335	2007	2009	2012	rVV3	103273	153330	3.22%	0.282%
51	13.372	2012	2015	2018	rVV	104524	132365	2.78%	0.243%
52	13.420	2018	2023	2031	rVB	1831656	2231571	46.92%	4.103%
53	13.506	2032	2037	2040	rBV2	170583	234986	4.94%	0.432%
54	13.542	2040	2043	2046	rVV	305760	405921	8.53%	0.746%
55	13.585	2046	2050	2054	rVV3	454746	759612	15.97%	1.397%
56	13.640	2054	2059	2060	rVV2	213650	345048	7.25%	0.634%
57	13.664	2060	2063	2069	rVB2	419717	604917	12.72%	1.112%
58	13.774	2074	2081	2088	rBV	276826	413037	8.68%	0.759%
59	13.853	2089	2094	2099	rVB	539819	677223	14.24%	1.245%
60	13.926	2099	2106	2110	rBV2	549698	767556	16.14%	1.411%
61	13.969	2110	2113	2117	rVB	208174	234406	4.93%	0.431%
62	14.073	2126	2130	2135	rBV	421084	514595	10.82%	0.946%
63	14.231	2149	2156	2163	rVV	1113824	1897009	39.89%	3.488%
64	14.323	2167	2171	2175	rBV	140416	183022	3.85%	0.337%
65	14.371	2175	2179	2183	rVB	392003	481112	10.12%	0.885%
66	14.420	2183	2187	2192	rVB2	56951	79473	1.67%	0.146%
67	14.481	2192	2197	2202	rBV2	854341	1102406	23.18%	2.027%
68	14.615	2214	2219	2220	rBV	311027	346912	7.29%	0.638%
69	14.634	2220	2222	2226	rVV	384212	507297	10.67%	0.933%
70	14.670	2226	2228	2233	rVB	183071	213287	4.48%	0.392%
71	14.725	2233	2237	2240	rBV2	71007	91106	1.92%	0.168%
72	14.780	2240	2246	2249	rVV	218323	289203	6.08%	0.532%
73	14.816	2249	2252	2255	rVV2	128106	192308	4.04%	0.354%
74	14.847	2255	2257	2261	rVB2	100386	104337	2.19%	0.192%
75	14.902	2261	2266	2272	rBV2	123958	210870	4.43%	0.388%
76	15.048	2283	2290	2295	rBV2	50205	78911	1.66%	0.145%

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Title : SW846 8260

77	15.182	2307	2312	2320	rVV2	204707	397833	8.36%	0.732%
78	15.255	2320	2324	2331	rVB5	32645	64318	1.35%	0.118%
79	15.371	2339	2343	2352	rVB	50556	90719	1.91%	0.167%
80	15.475	2352	2360	2365	rVV3	91806	177114	3.72%	0.326%
81	15.542	2365	2371	2378	rVV4	89774	201322	4.23%	0.370%
82	15.609	2378	2382	2386	rVV	83642	144421	3.04%	0.266%
83	15.652	2386	2389	2397	rVB	48803	76395	1.61%	0.140%
84	15.841	2416	2420	2434	rVB7	19475	54902	1.15%	0.101%
85	16.475	2517	2524	2534	rBV2	49066	99798	2.10%	0.184%

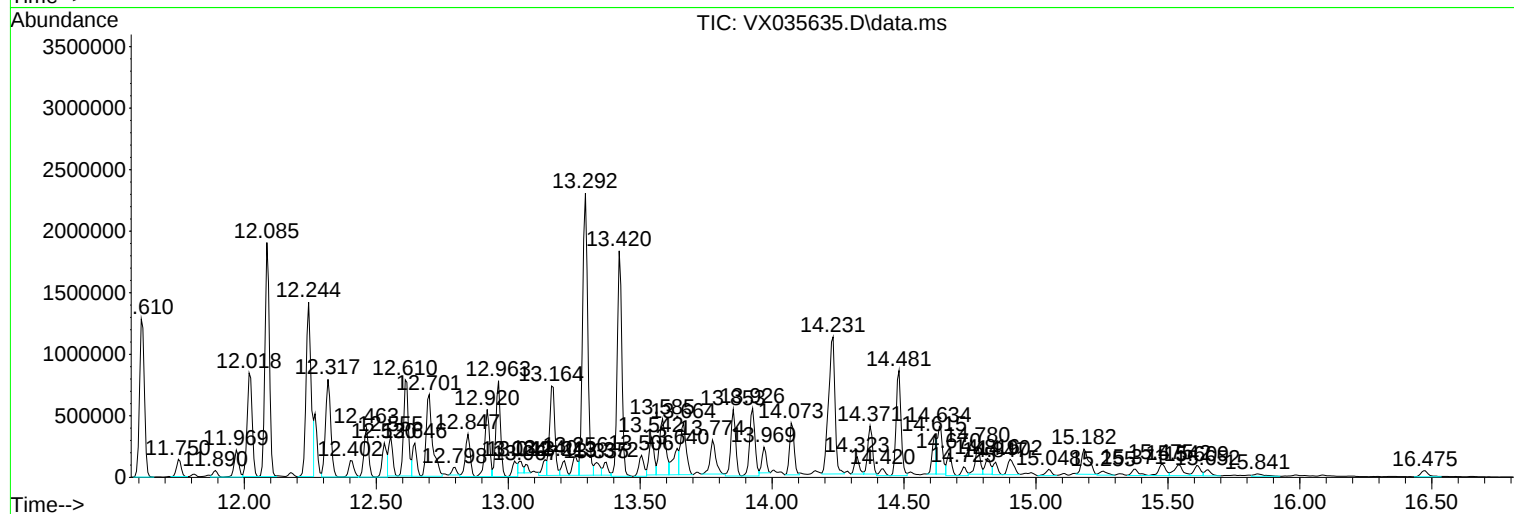
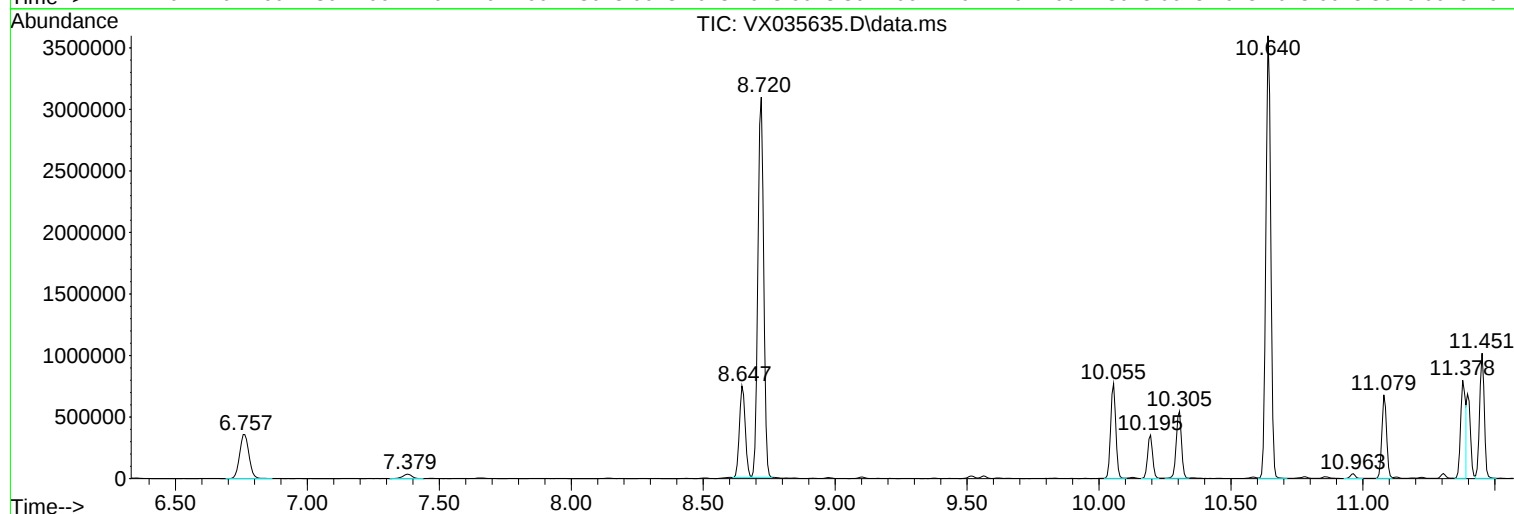
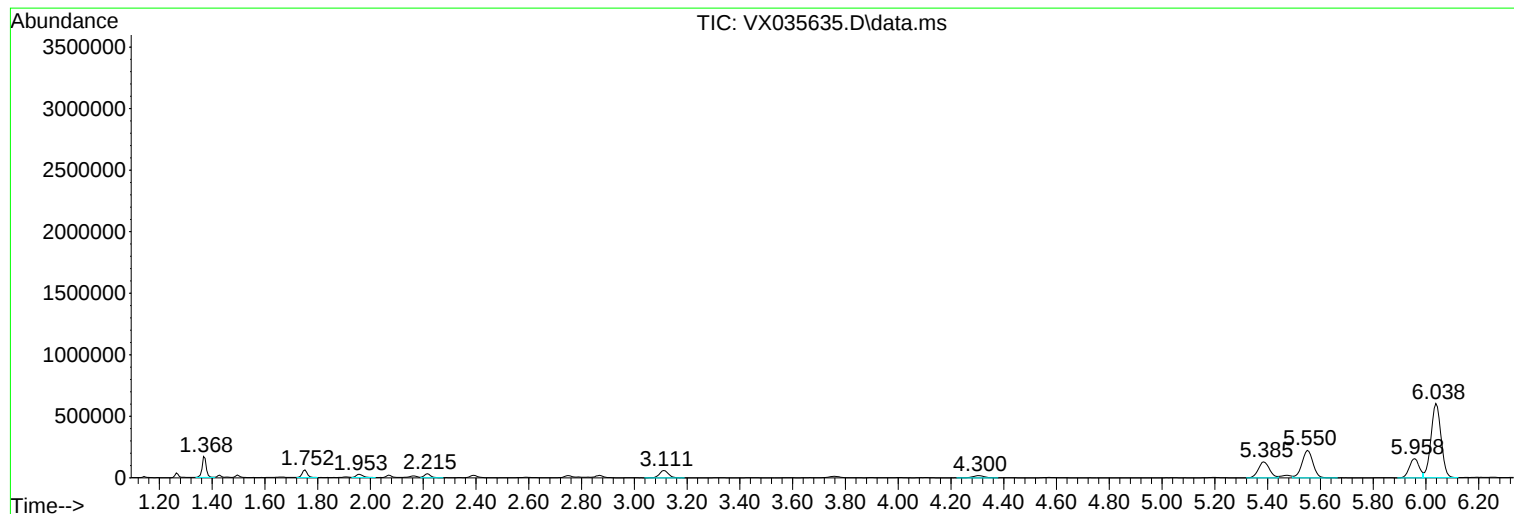
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Instrument :
MSVOA_X
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
Quant Title : SW846 8260

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



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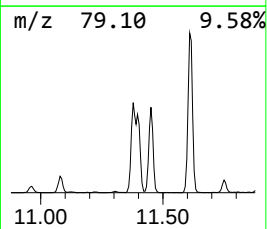
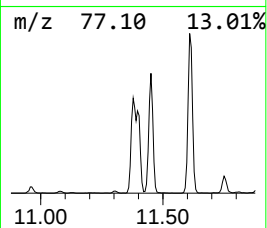
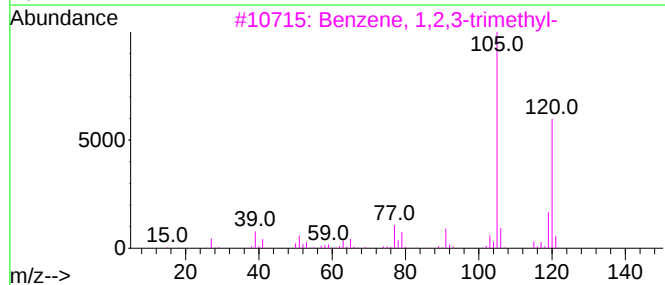
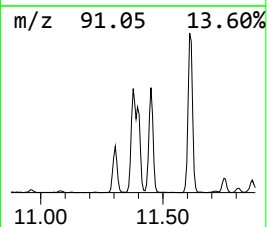
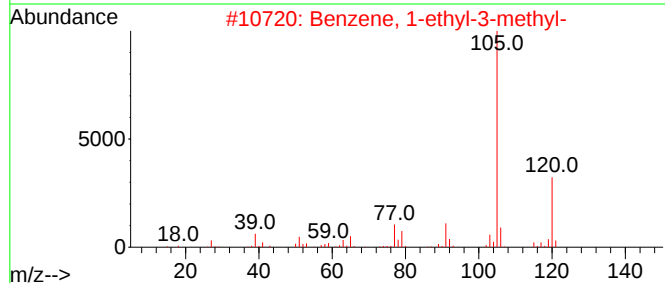
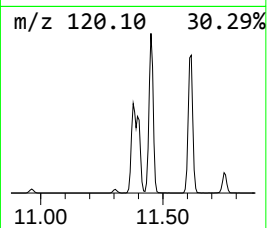
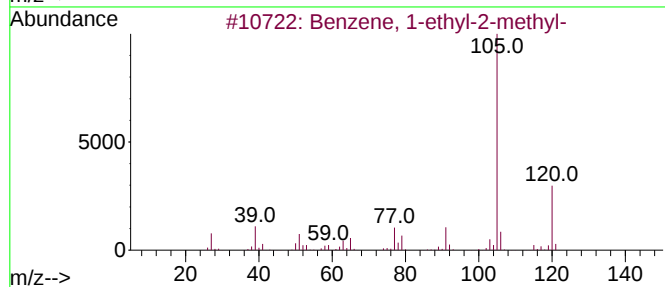
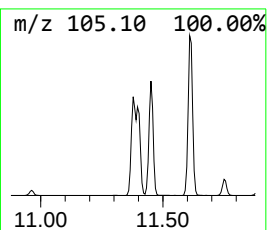
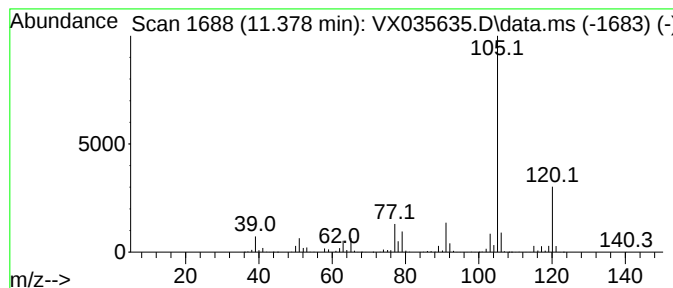
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	47.43 ug/l	1069010	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	94
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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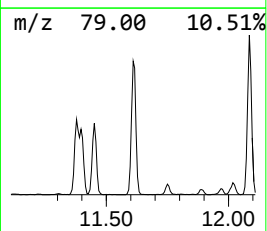
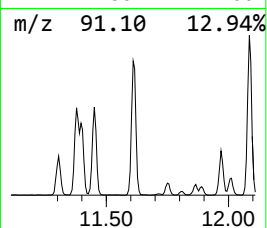
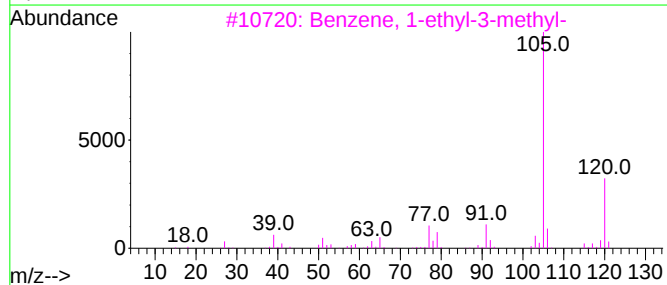
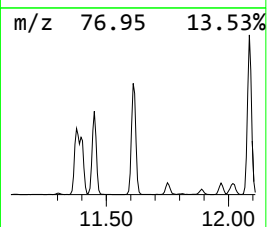
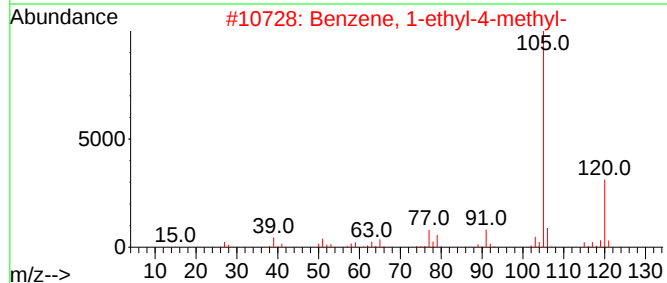
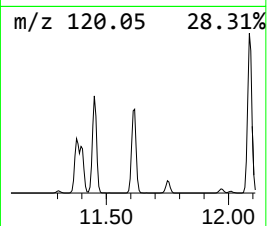
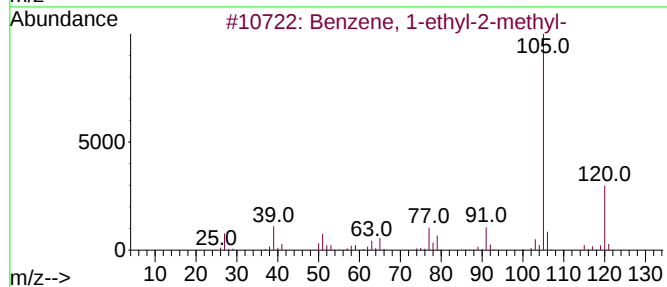
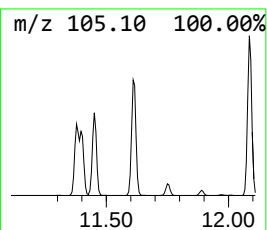
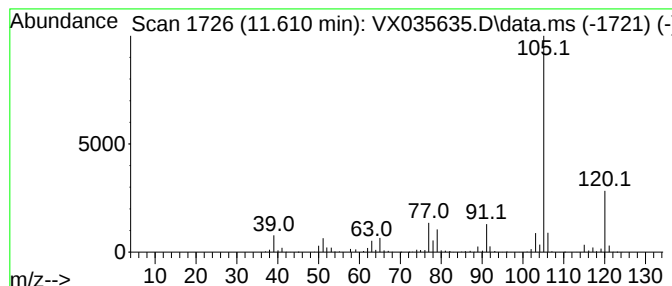
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	72.04 ug/l	1623680	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	91
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



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Instrument :
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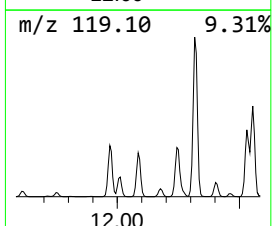
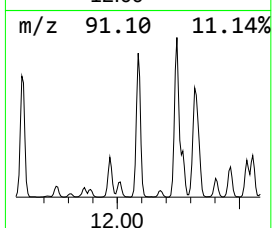
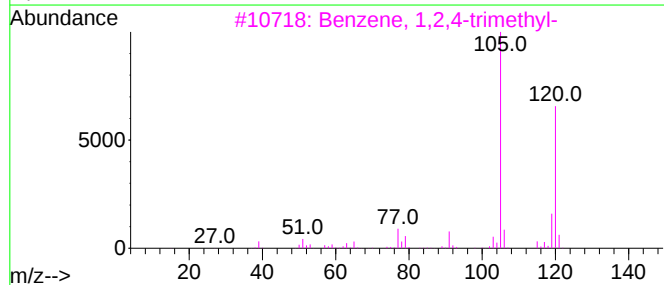
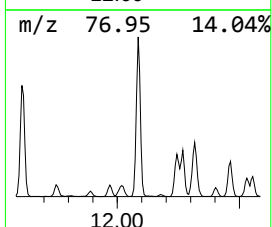
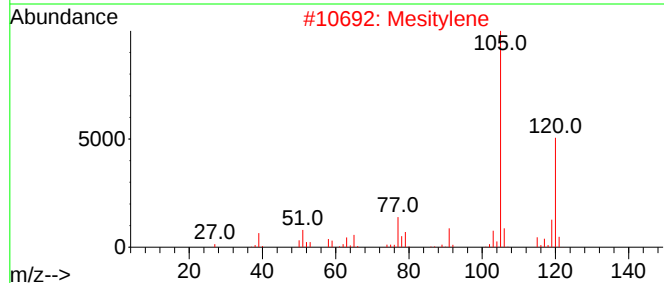
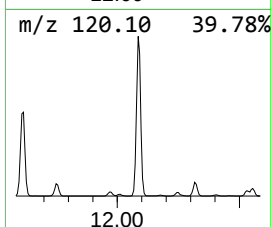
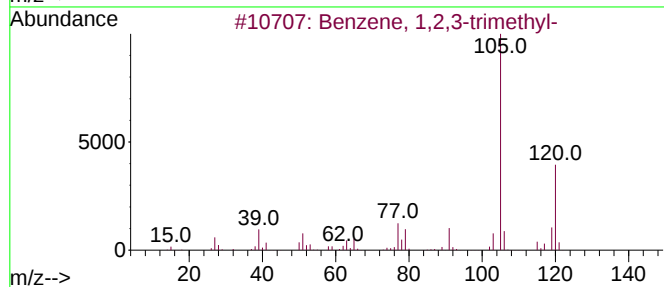
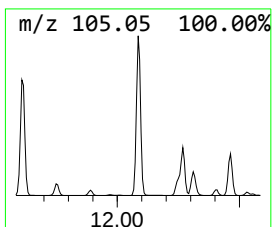
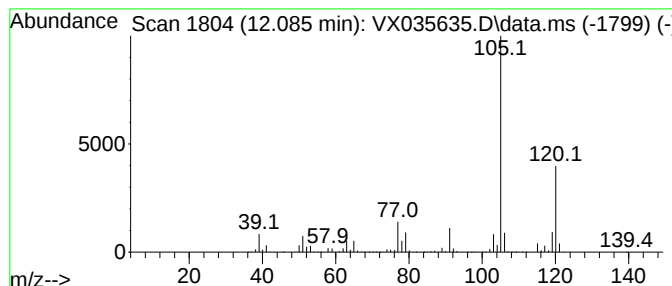
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	99.45 ug/l	2241270	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
50033

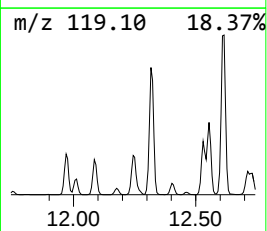
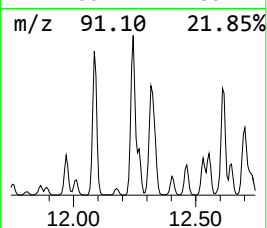
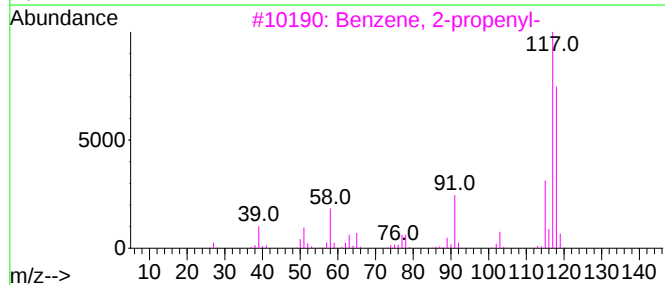
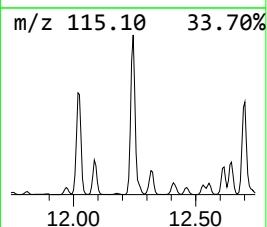
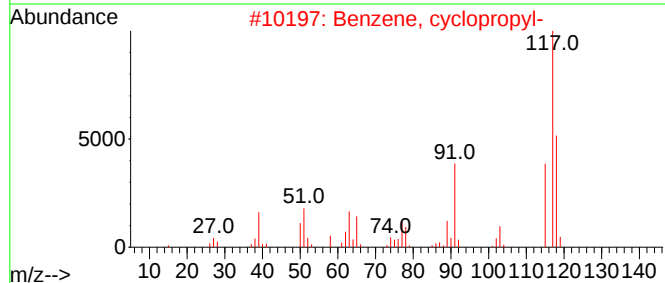
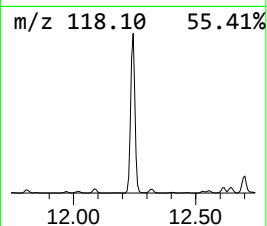
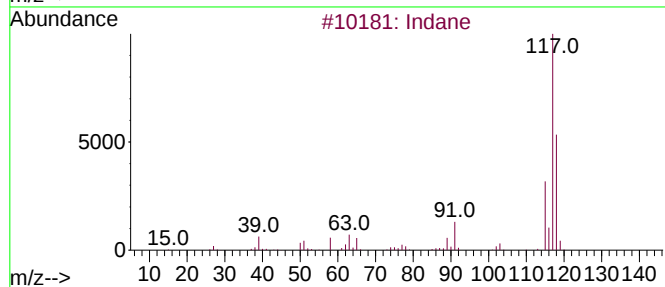
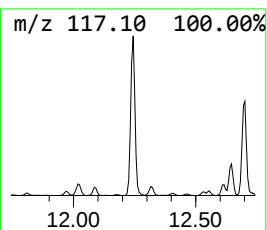
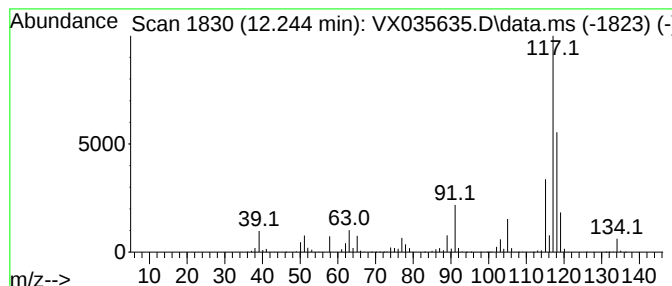
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	85.69 ug/l	1931200	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50033

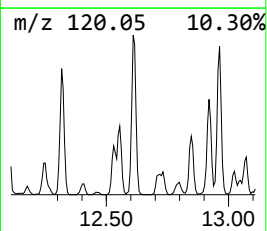
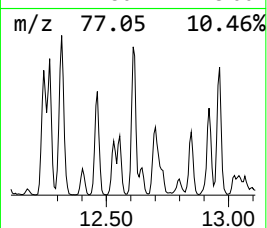
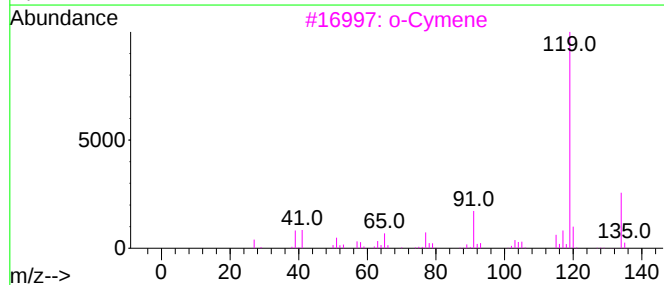
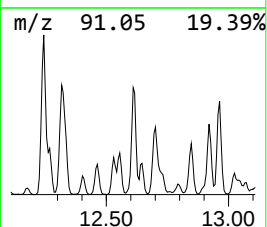
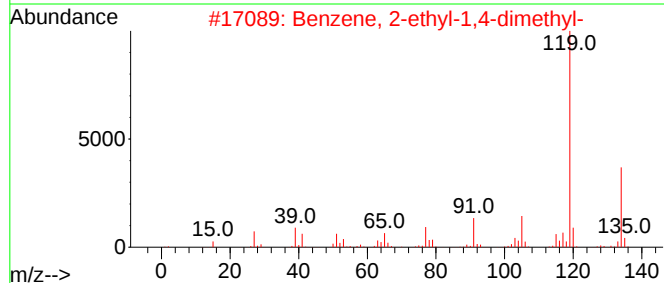
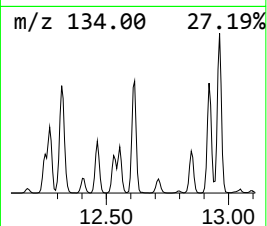
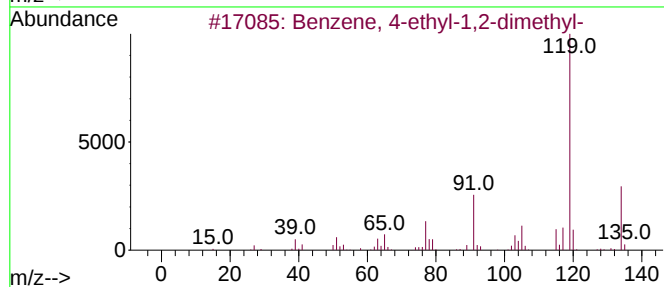
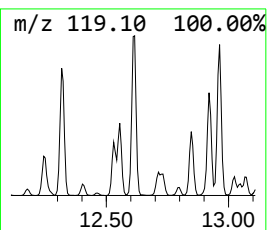
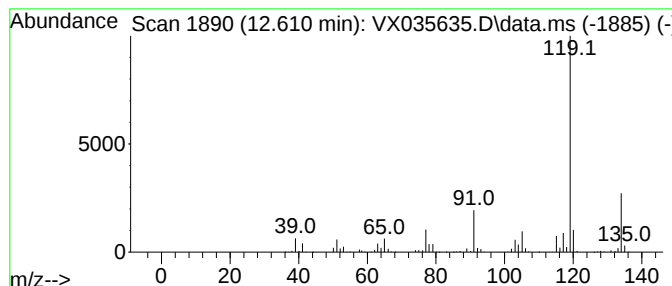
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	45.07 ug/l	1015700	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50033

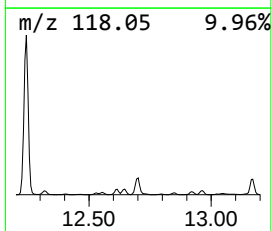
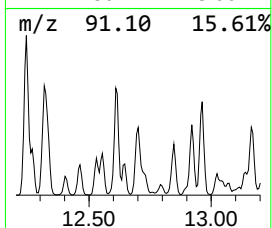
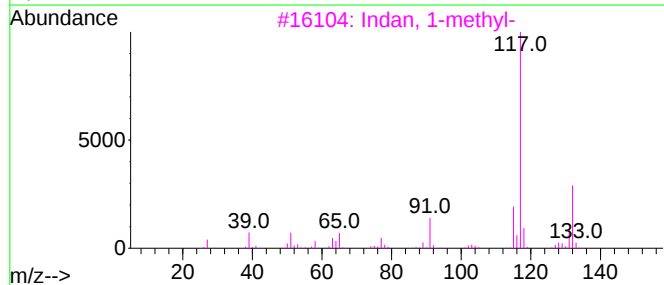
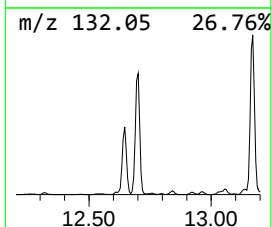
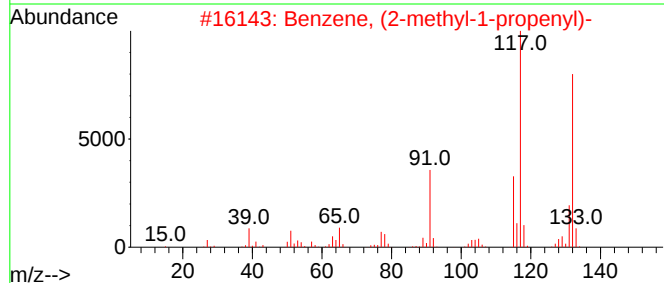
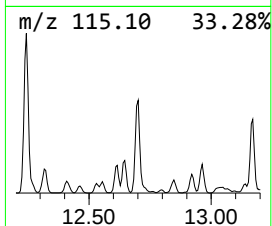
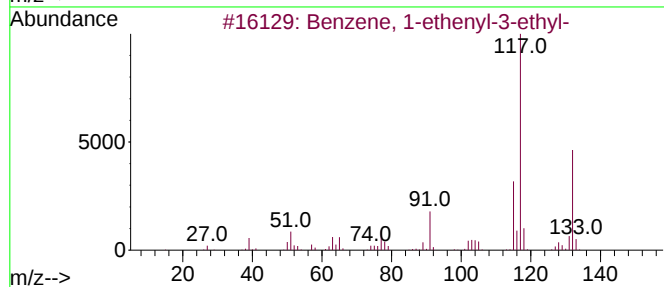
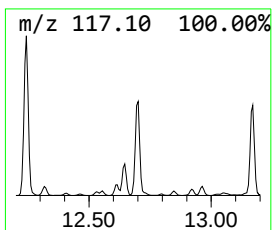
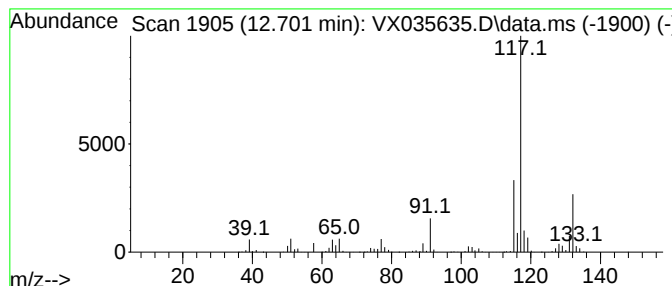
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	48.96 ug/l	1103420	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			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50033

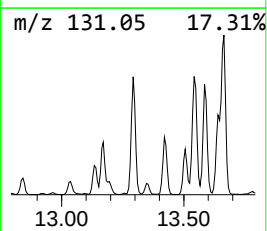
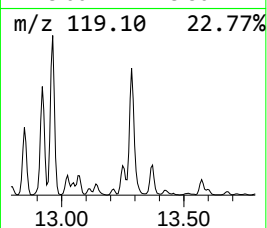
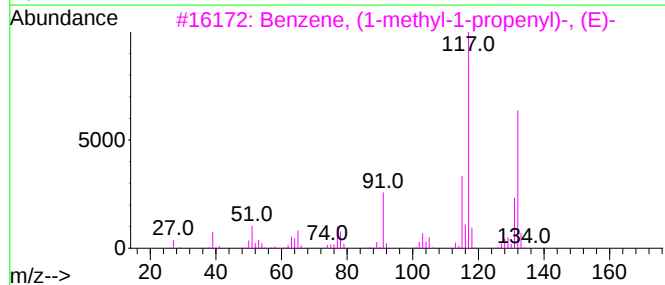
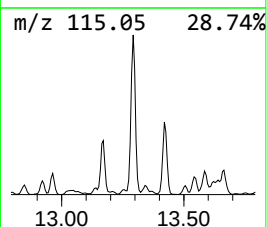
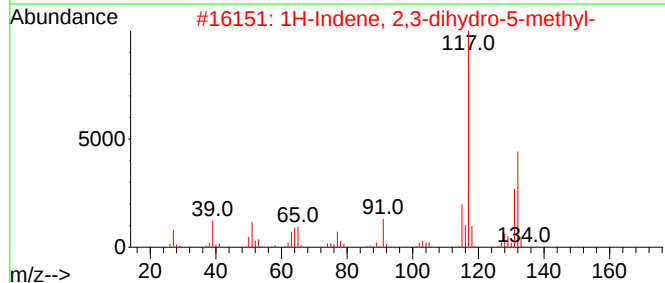
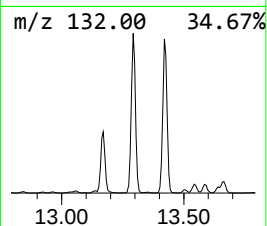
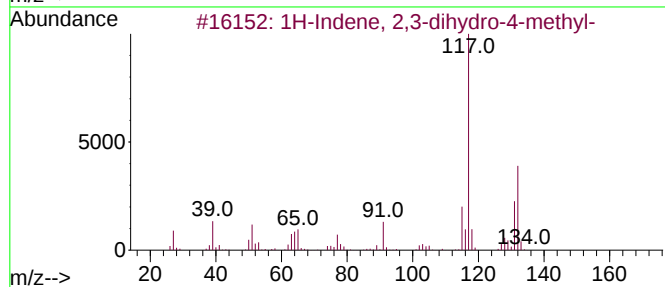
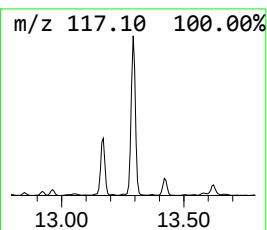
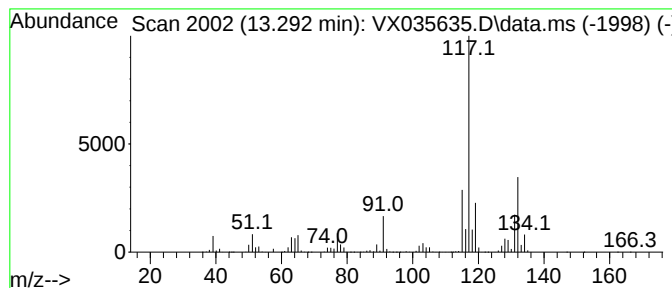
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	135.30 ug/l	3049350	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	93
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50033

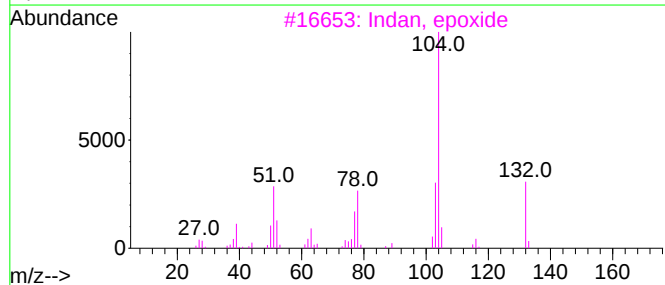
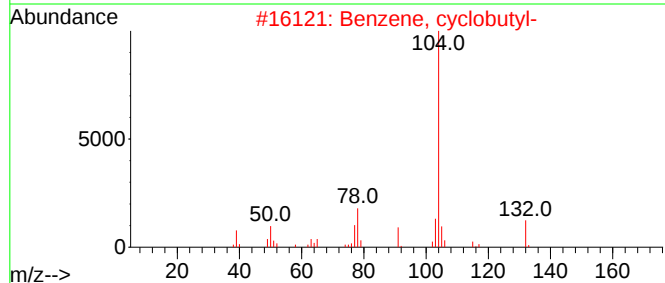
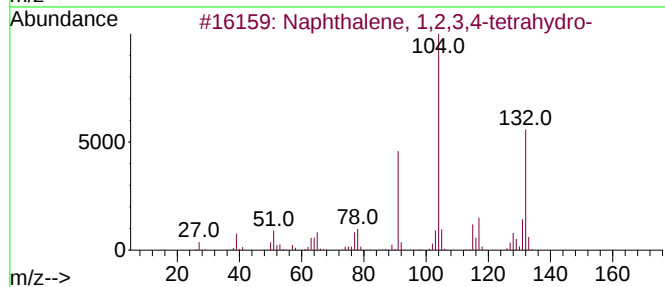
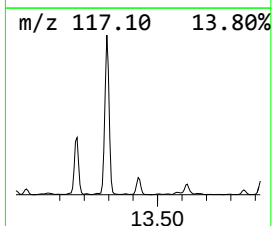
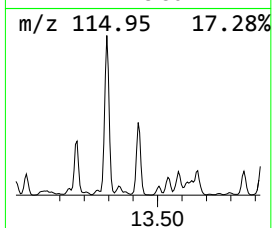
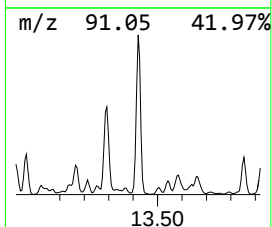
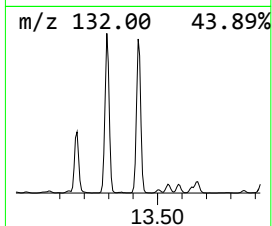
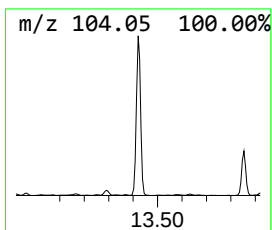
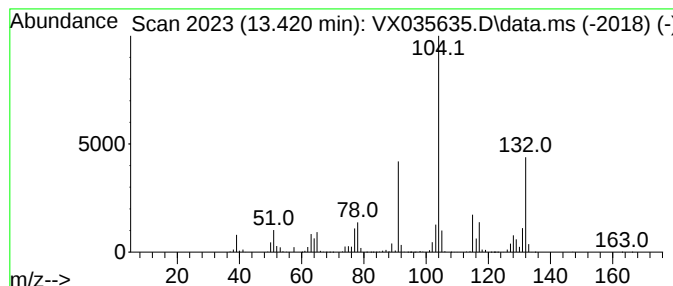
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	99.02 ug/l	2231570	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	97
2			Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
3			Indan, epoxide	132	C9H8O	000768-22-9	58
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	52
5			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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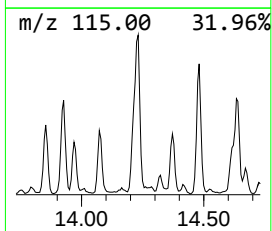
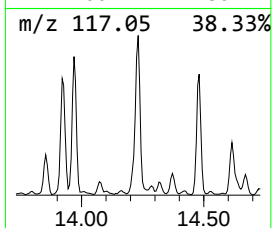
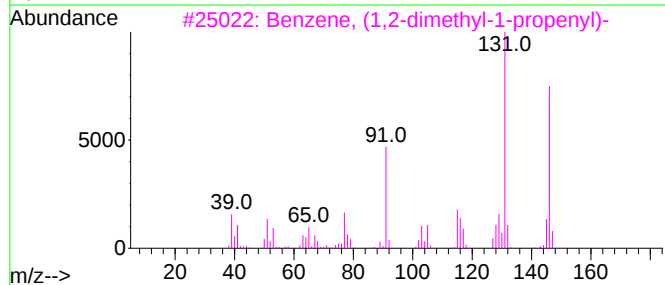
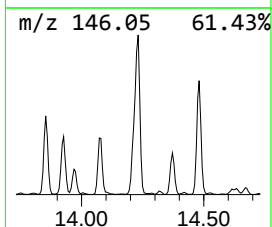
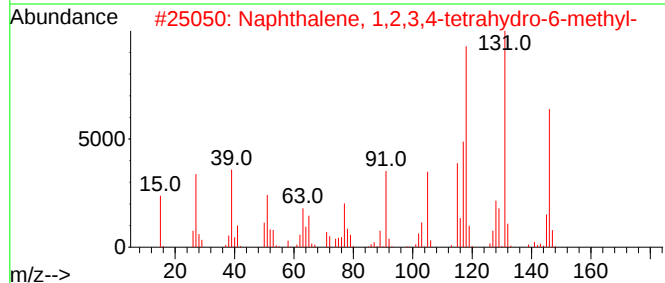
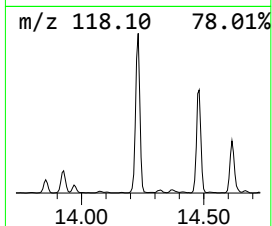
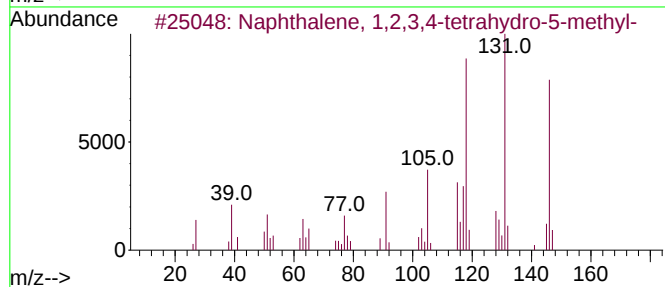
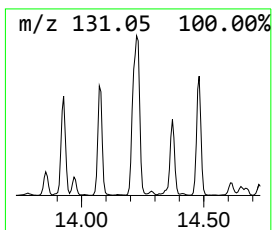
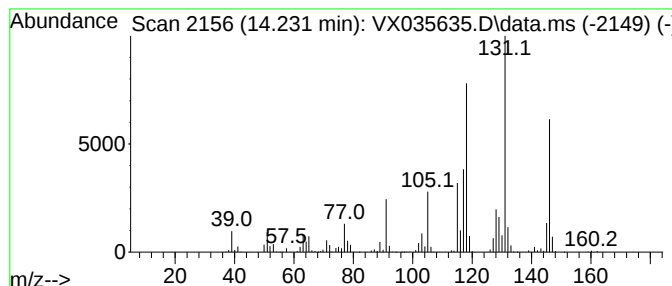
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.231	84.17 ug/l	1897010	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	87
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	60
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
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Operator : JC/MD
Sample : 02702-02
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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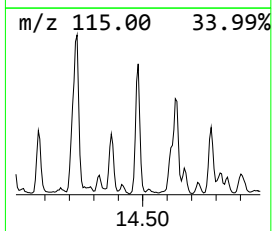
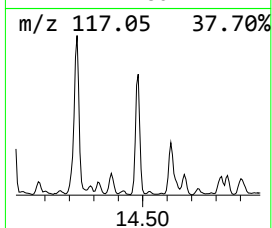
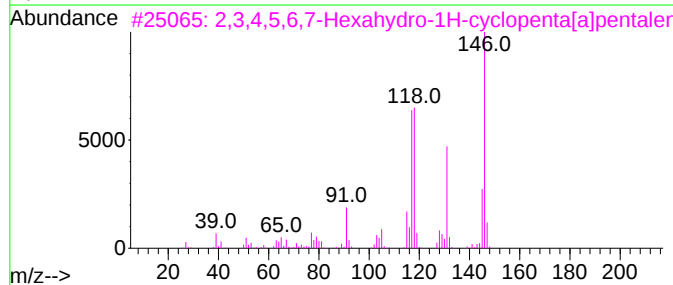
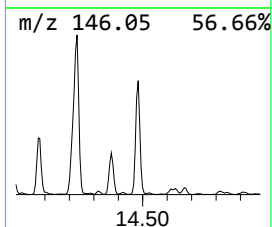
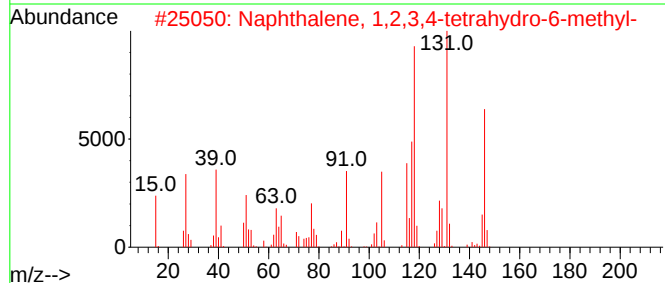
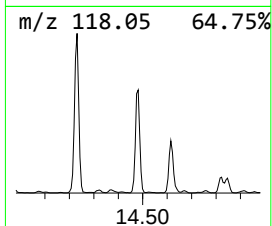
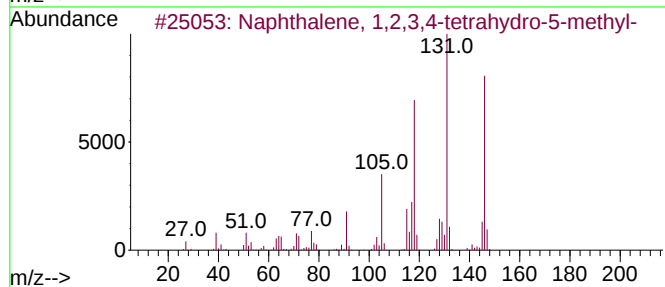
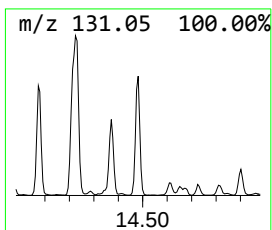
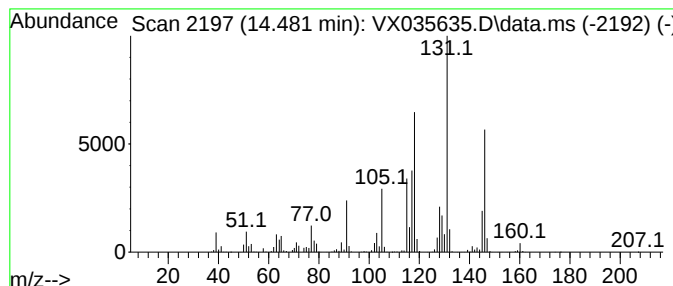
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	48.91 ug/l	1102410	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	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
3			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	76
4			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	72
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051223\
Data File : VX035635.D
Acq On : 12 May 2023 14:07
Operator : JC/MD
Sample : 02702-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50033

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	11.378	47.4	ug/l	1069010	4	12.024	1126870	50.0
Benzene, 1-ethy...	11.610	72.0	ug/l	1623680	4	12.024	1126870	50.0
Benzene, 1,2,3-...	12.085	99.5	ug/l	2241270	4	12.024	1126870	50.0
Indane	12.244	85.7	ug/l	1931200	4	12.024	1126870	50.0
Benzene, 4-ethy...	12.610	45.1	ug/l	1015700	4	12.024	1126870	50.0
Benzene, 1-ethe...	12.701	49.0	ug/l	1103420	4	12.024	1126870	50.0
1H-Indene, 2,3-...	13.292	135.3	ug/l	3049350	4	12.024	1126870	50.0
Naphthalene, 1,...	13.420	99.0	ug/l	2231570	4	12.024	1126870	50.0
Naphthalene, 1,...	14.231	84.2	ug/l	1897010	4	12.024	1126870	50.0
Naphthalene, 1,...	14.481	48.9	ug/l	1102410	4	12.024	1126870	50.0