

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
 Data File : VX042115.D
 Acq On : 01 Jul 2024 15:32
 Operator : JC/MD
 Sample : P3092-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW3-20240626

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\82X061824W.M
 Title : SW846 8260

Signal : TIC: VX042115.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.240	21	25	34	rBV	56434	67389	4.96%	0.493%
2	4.300	518	527	539	rVB3	5617	16232	1.19%	0.119%
3	5.093	644	657	668	rBV	15787	49018	3.61%	0.359%
4	5.385	694	705	717	rBV2	77951	235235	17.32%	1.722%
5	5.550	722	732	754	rVB	131294	392802	28.92%	2.876%
6	5.958	789	799	817	rBV	74462	202363	14.90%	1.481%
7	6.245	841	846	856	rVB3	8624	25039	1.84%	0.183%
8	6.757	920	930	955	rBV	201194	498700	36.71%	3.651%
9	7.373	1021	1031	1043	rBV6	15827	44784	3.30%	0.328%
10	7.982	1122	1131	1139	rVB4	6091	15025	1.11%	0.110%
11	8.104	1145	1151	1155	rBV3	7996	17395	1.28%	0.127%
12	8.501	1211	1216	1224	rVB3	9562	18614	1.37%	0.136%
13	8.647	1234	1240	1247	rVV	431993	670699	49.38%	4.910%
14	8.720	1247	1252	1258	rVB	148125	230293	16.95%	1.686%
15	10.055	1465	1471	1477	rBV	452470	623998	45.94%	4.568%
16	10.195	1488	1494	1504	rBV	679483	905863	66.69%	6.632%
17	10.299	1504	1511	1521	rBV	1013540	1358361	100.00%	9.944%
18	10.640	1561	1567	1579	rVB	586505	742665	54.67%	5.437%
19	10.964	1613	1620	1627	rBV	43051	56779	4.18%	0.416%
20	11.079	1634	1639	1651	rBV	405745	520100	38.29%	3.808%
21	11.305	1670	1676	1683	rBV	81858	100960	7.43%	0.739%
22	11.378	1683	1688	1690	rBV	153617	200708	14.78%	1.469%
23	11.451	1696	1700	1709	rVB	163996	195059	14.36%	1.428%
24	11.610	1721	1726	1739	rVB	492153	604271	44.49%	4.424%
25	11.750	1744	1749	1756	rVV	1007639	1185900	87.30%	8.682%
26	12.018	1788	1793	1799	rVV	496537	646580	47.60%	4.734%
27	12.085	1799	1804	1810	rVB	569876	677420	49.87%	4.959%
28	12.244	1824	1830	1838	rBV	403942	533864	39.30%	3.908%
29	12.317	1838	1842	1852	rVB2	76552	108618	8.00%	0.795%
30	12.408	1852	1857	1862	rBV3	53663	78382	5.77%	0.574%
31	12.463	1862	1866	1872	rVB	79920	103352	7.61%	0.757%
32	12.530	1872	1877	1879	rBV	71303	92644	6.82%	0.678%
33	12.555	1879	1881	1886	rVB	77245	80502	5.93%	0.589%
34	12.610	1886	1890	1894	rBV	84954	111125	8.18%	0.814%
35	12.646	1894	1896	1900	rVB	26644	27589	2.03%	0.202%
36	12.695	1900	1904	1912	rBV2	172277	264821	19.50%	1.939%

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Integrator: RTE

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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37	12.847	1924	1929	1934	rVB2	59738	74913	5.51%	0.548%
38	12.921	1934	1941	1944	rBV	138955	169182	12.45%	1.239%
39	12.963	1944	1948	1954	rVB	147897	179501	13.21%	1.314%
40	13.030	1954	1959	1964	rBV3	16168	30000	2.21%	0.220%
41	13.134	1967	1976	1978	rBV4	16357	27340	2.01%	0.200%
42	13.170	1978	1982	1987	rVB	59448	72070	5.31%	0.528%
43	13.292	1997	2002	2007	rBV2	248673	324721	23.91%	2.377%
44	13.341	2007	2010	2013	rVB3	12201	15540	1.14%	0.114%
45	13.420	2019	2023	2031	rVB2	34512	50986	3.75%	0.373%
46	13.500	2031	2036	2039	rBV2	16781	25433	1.87%	0.186%
47	13.542	2039	2043	2046	rVV	46100	64415	4.74%	0.472%
48	13.579	2046	2049	2055	rVV3	56692	108793	8.01%	0.796%
49	13.640	2056	2059	2060	rVV	29597	35564	2.62%	0.260%
50	13.664	2060	2063	2069	rVB2	50794	70914	5.22%	0.519%
51	13.774	2075	2081	2091	rBV	335954	430158	31.67%	3.149%
52	13.865	2091	2096	2102	rVB	24271	36537	2.69%	0.267%
53	13.926	2102	2106	2110	rBV2	13995	19437	1.43%	0.142%
54	13.969	2110	2113	2117	rVB	18583	20169	1.48%	0.148%
55	14.073	2126	2130	2135	rBV	24552	31931	2.35%	0.234%
56	14.213	2148	2153	2159	rBV	28327	41650	3.07%	0.305%
57	14.323	2167	2171	2175	rVV	13712	17635	1.30%	0.129%
58	14.371	2175	2179	2184	rVB2	18071	21854	1.61%	0.160%
59	14.640	2219	2223	2227	rVB	23632	30248	2.23%	0.221%
60	14.780	2241	2246	2255	rBV	40635	57491	4.23%	0.421%

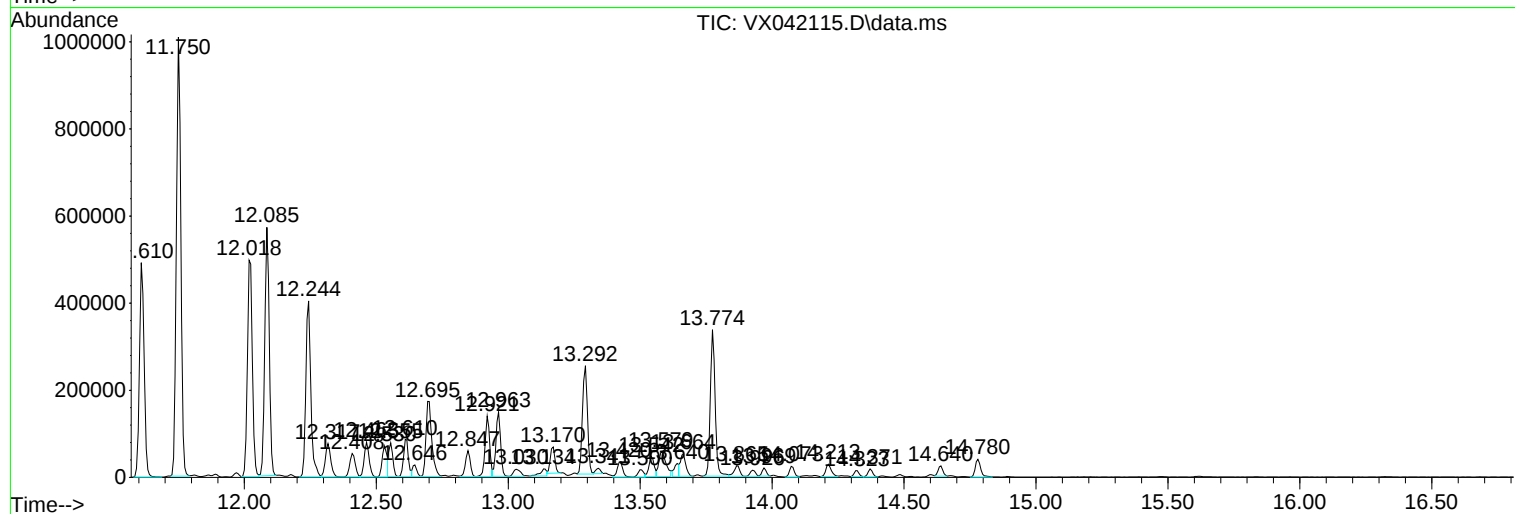
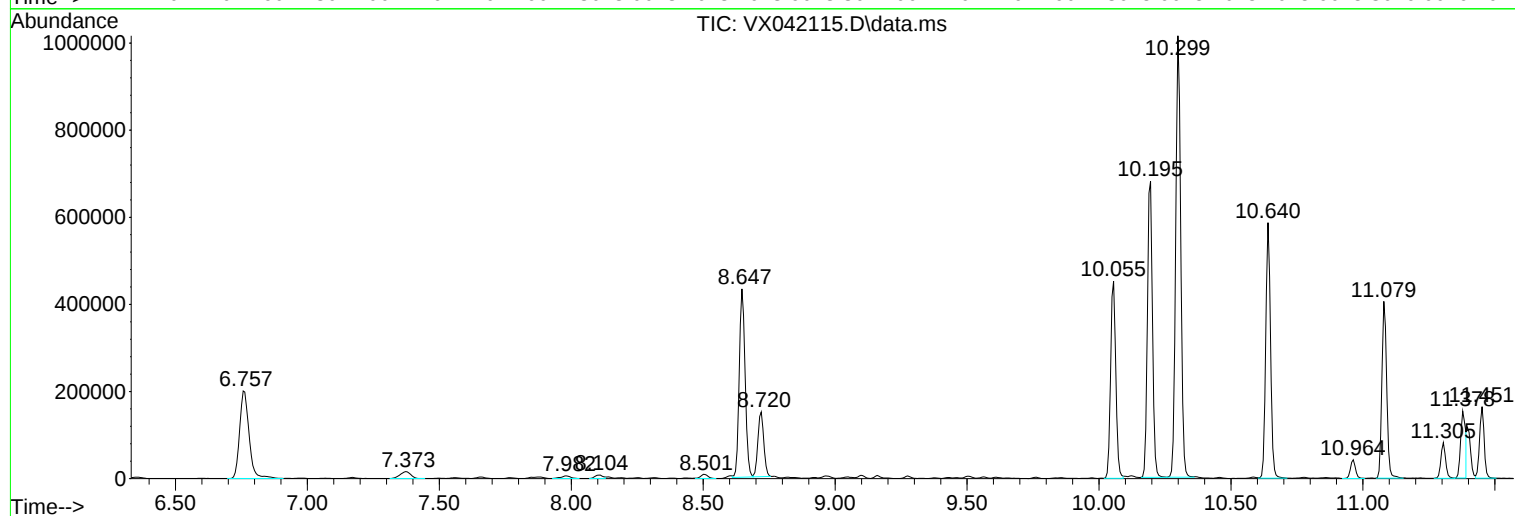
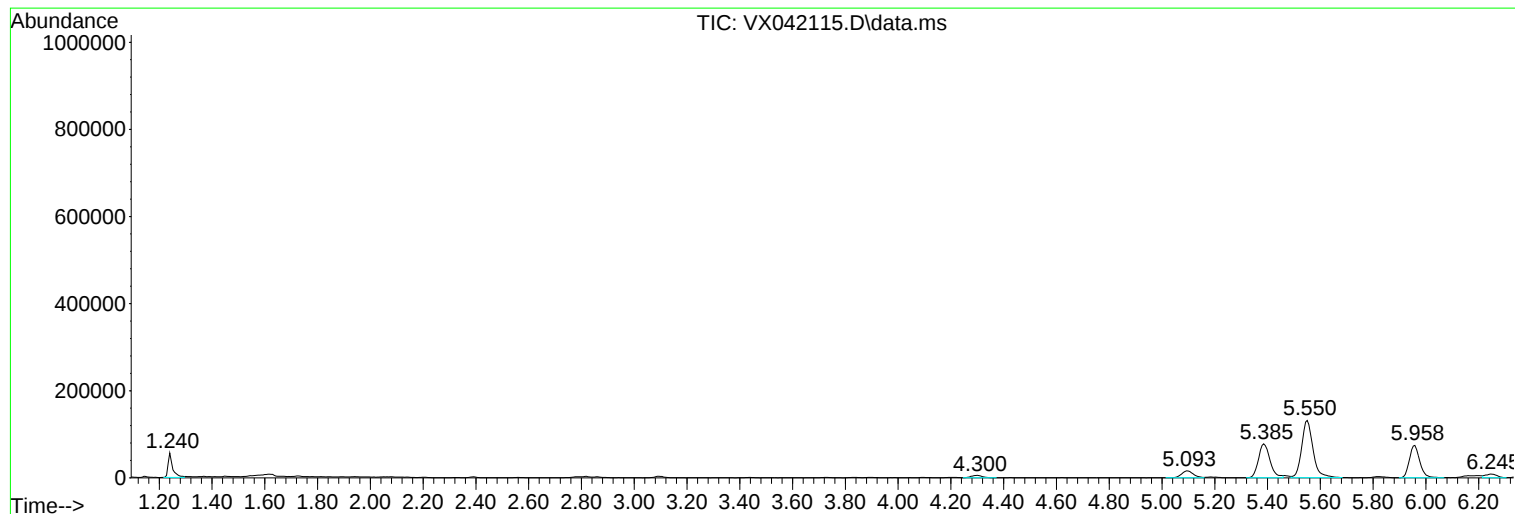
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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



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MSVOA_X
ClientSampleId :
MW3-20240626

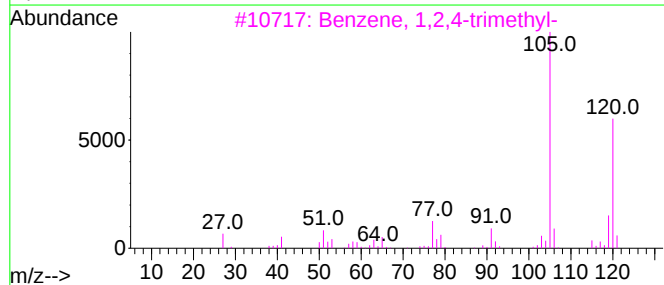
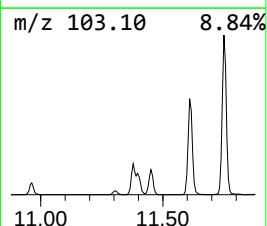
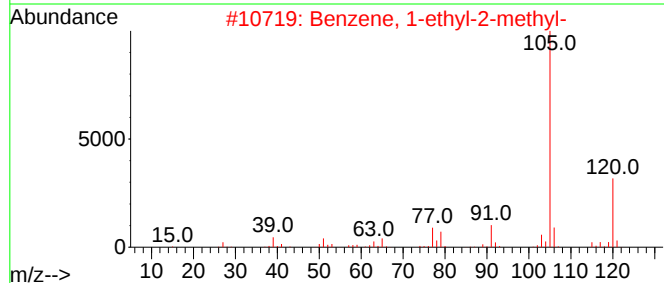
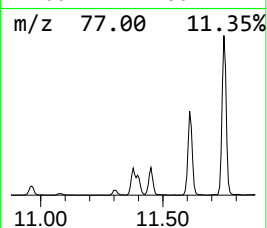
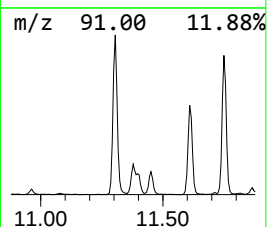
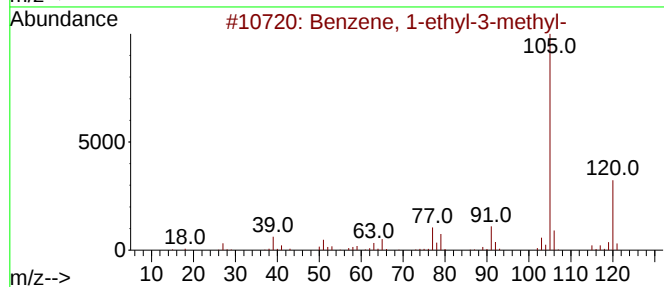
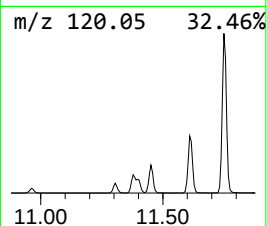
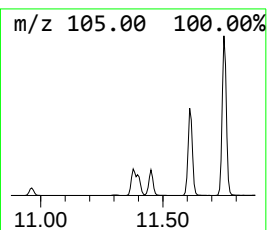
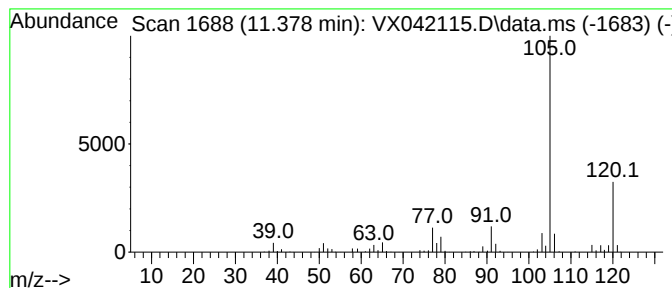
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	15.52 ug/l	200708	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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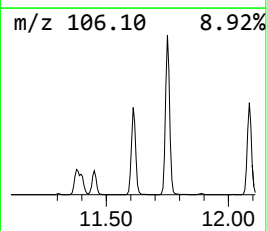
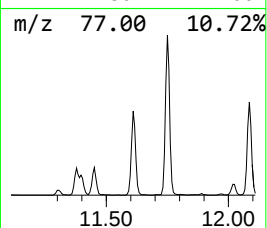
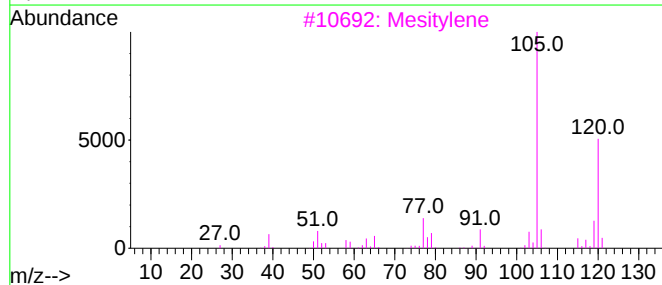
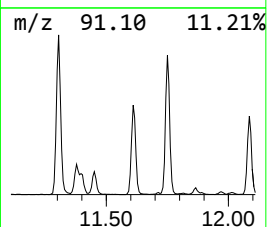
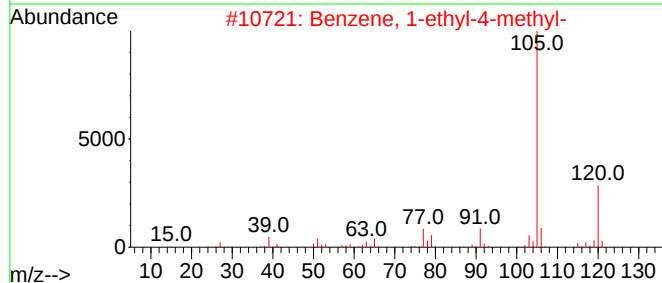
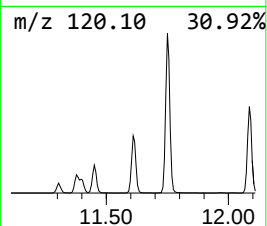
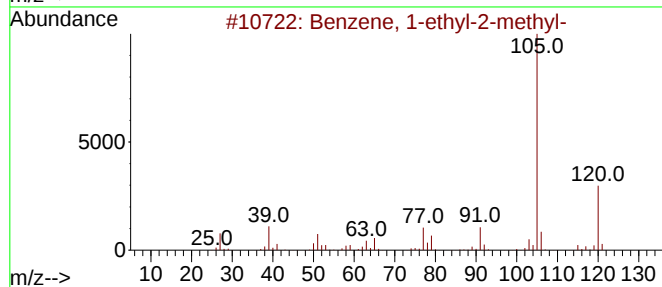
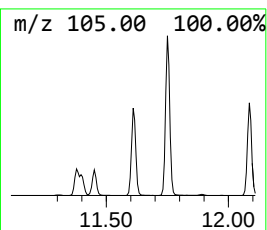
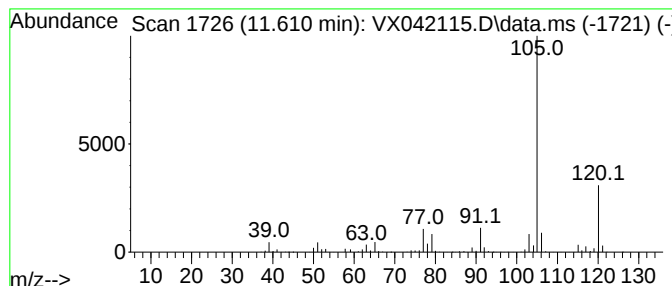
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	46.73 ug/l	604271	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	90
3			Mesitylene	120	C9H12	000108-67-8	90
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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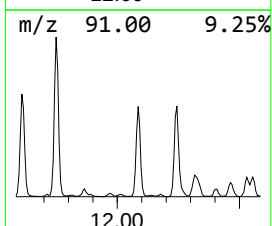
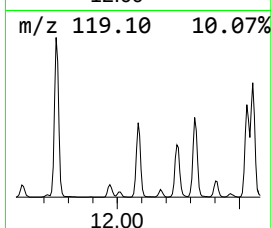
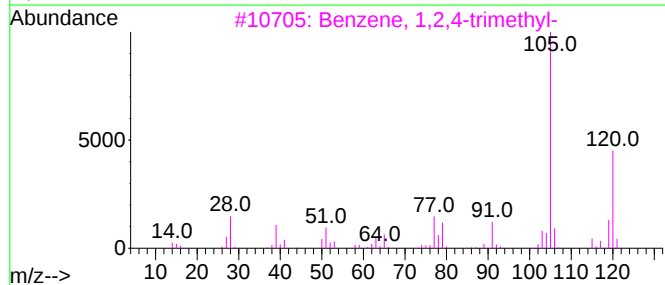
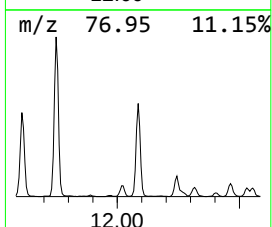
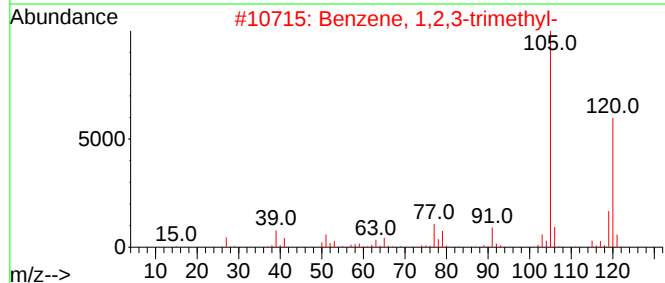
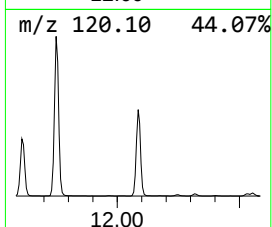
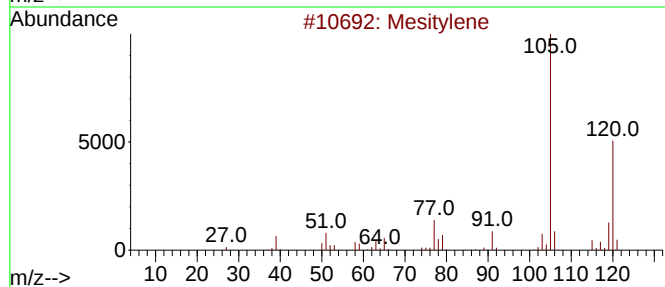
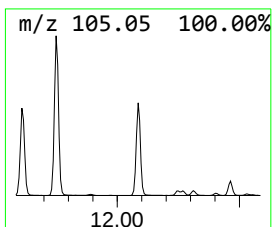
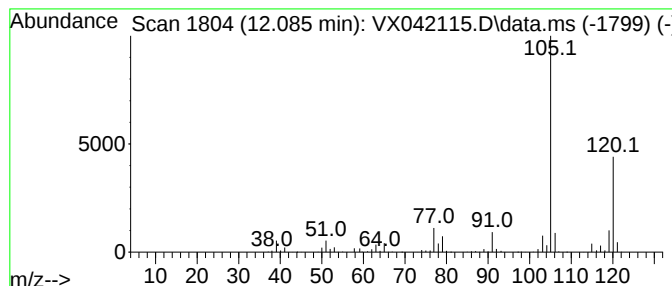
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	52.38 ug/l	677420	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	91



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Instrument :
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ClientSampleId :
MW3-20240626

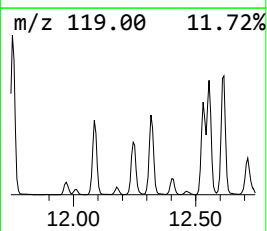
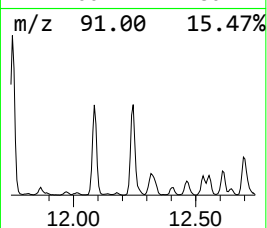
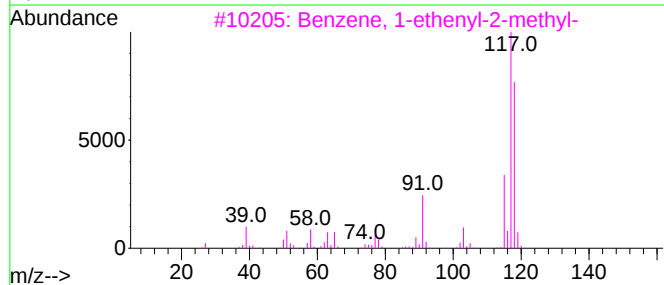
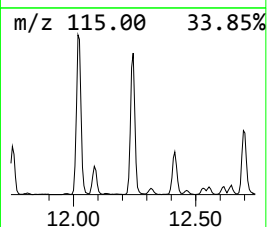
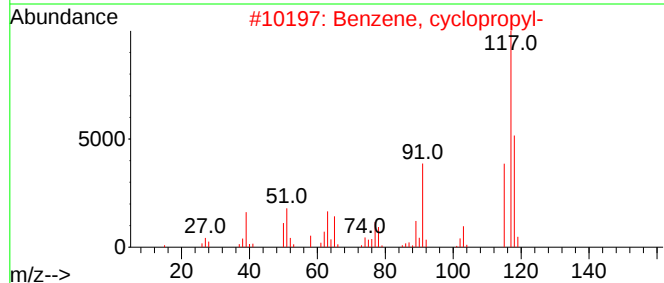
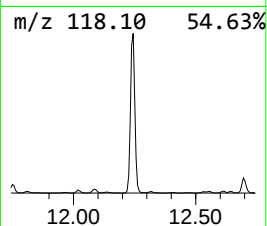
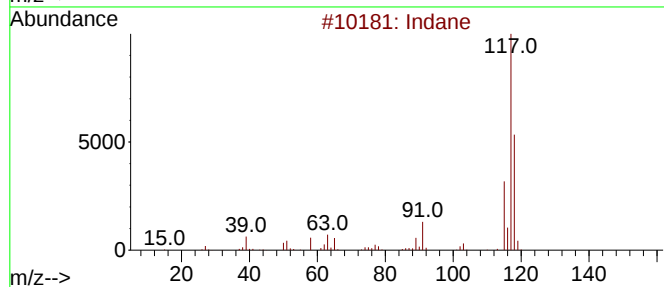
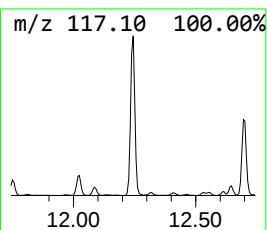
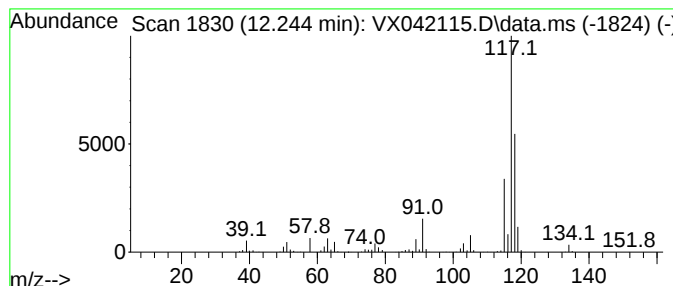
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	41.28 ug/l	533864	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	81
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VO070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

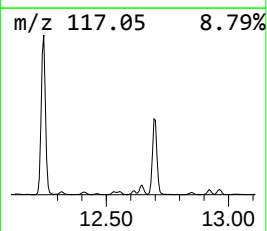
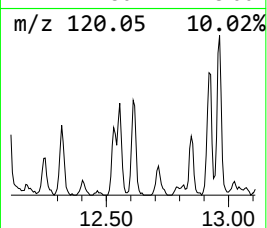
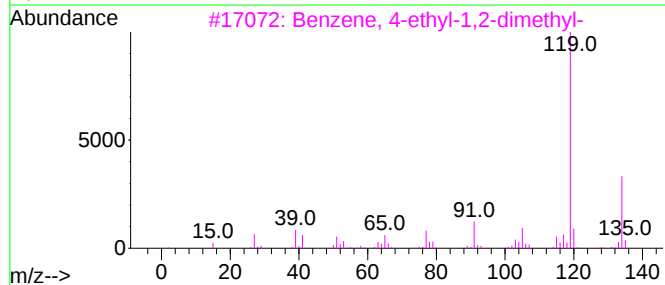
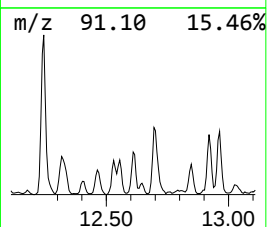
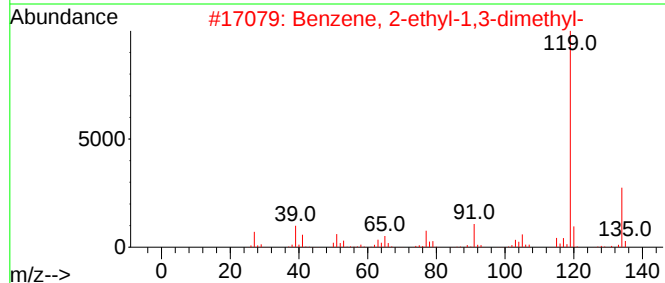
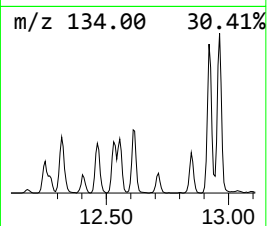
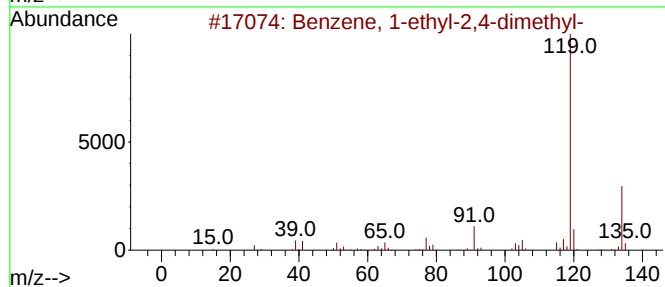
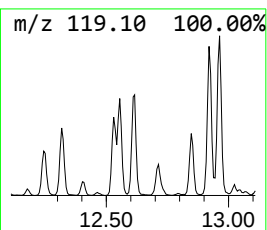
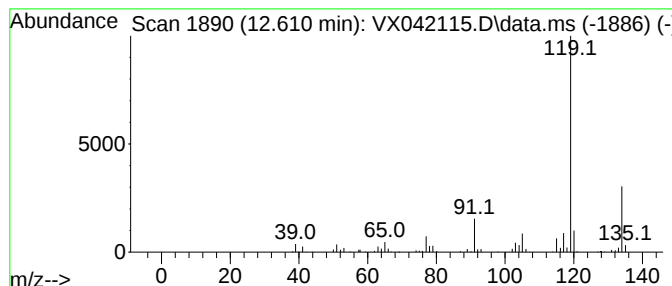
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	8.59 ug/l	111125	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-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

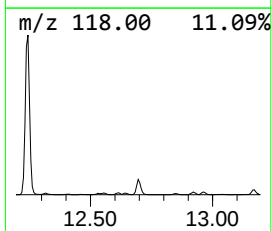
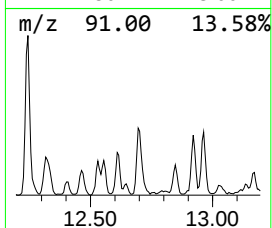
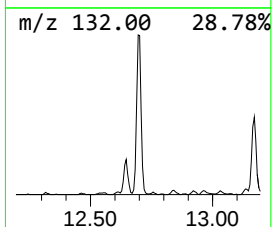
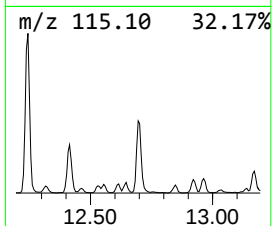
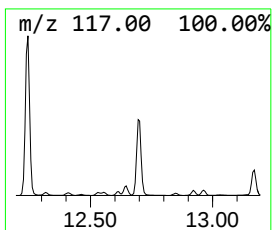
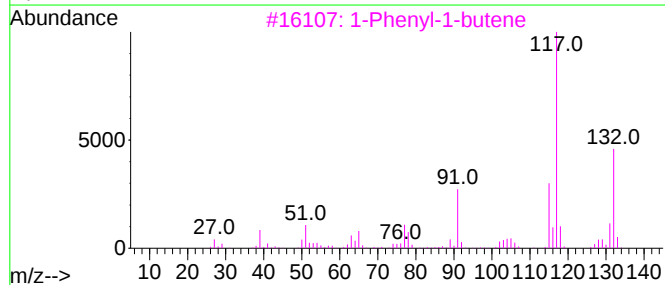
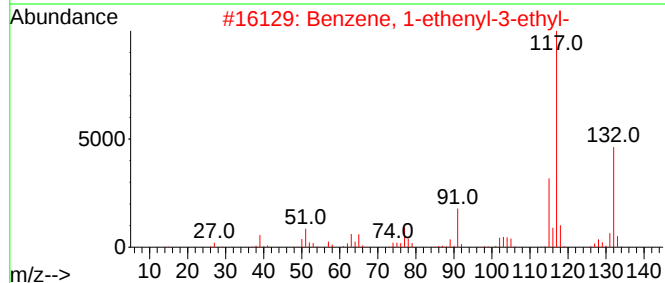
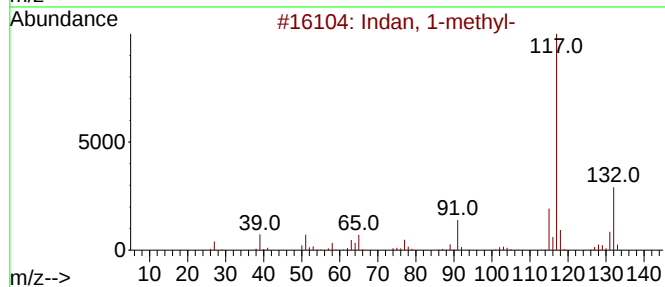
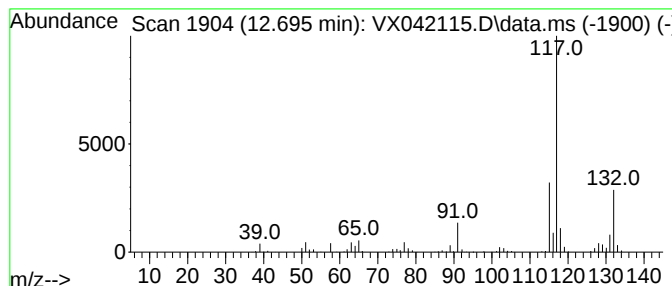
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	20.48 ug/l	264821	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

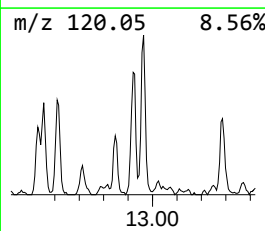
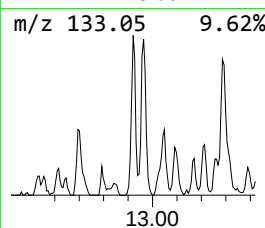
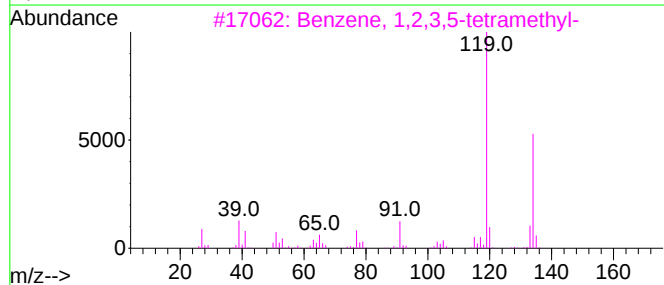
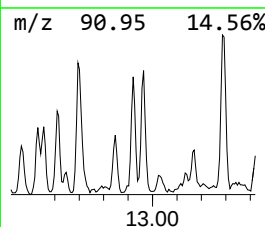
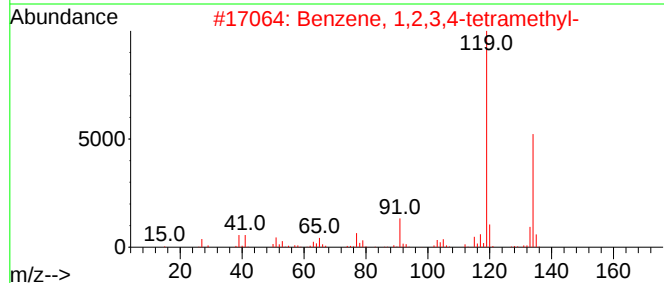
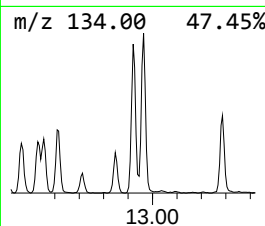
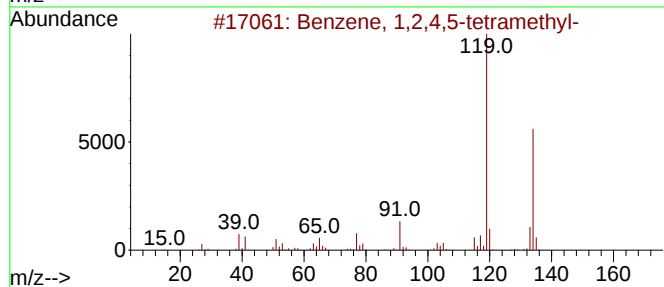
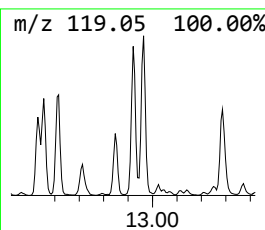
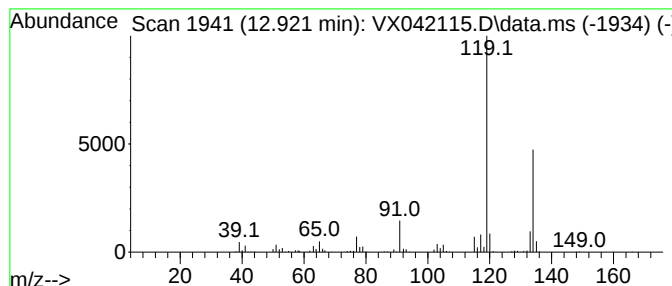
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.921	13.08 ug/l	169182	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5			o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

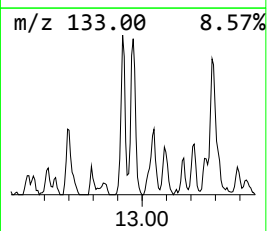
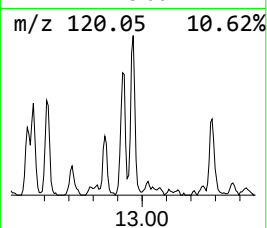
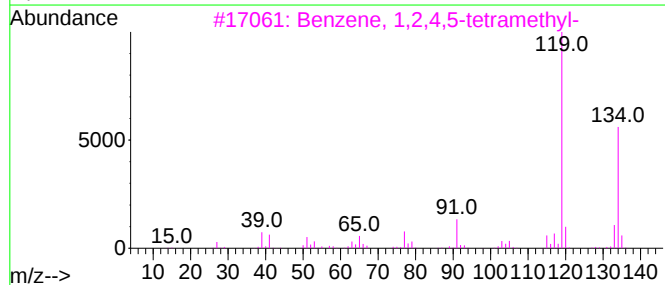
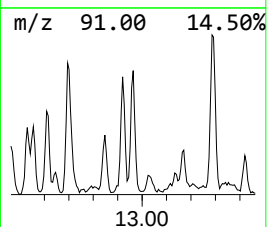
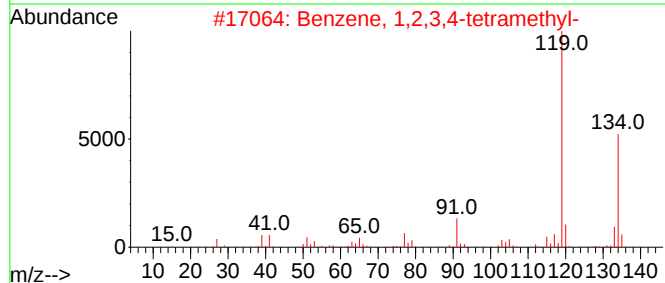
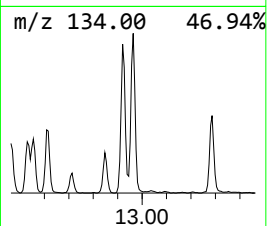
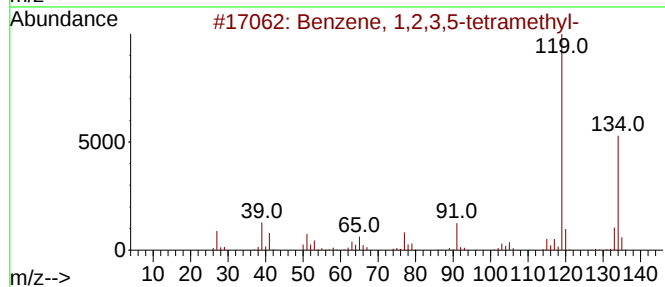
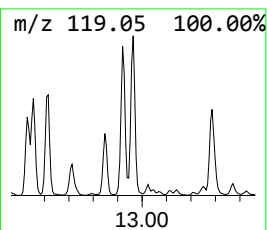
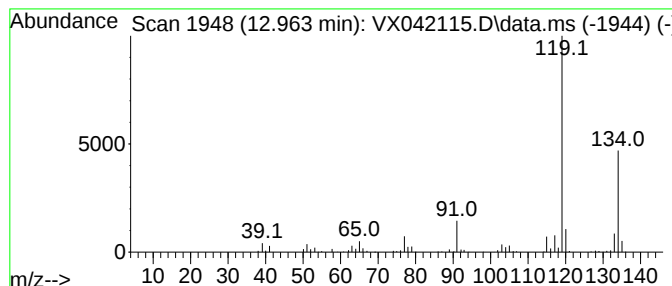
Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.963	13.88 ug/l	179501	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

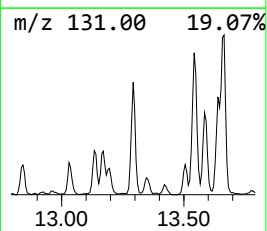
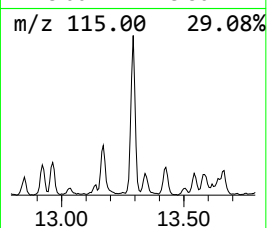
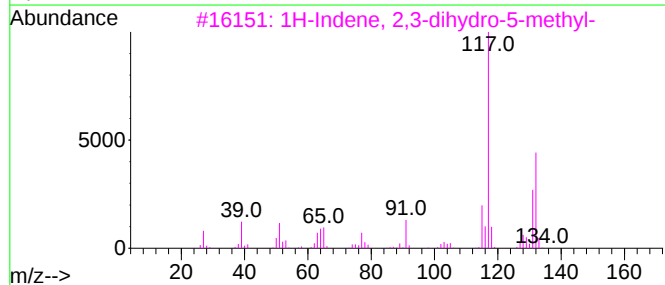
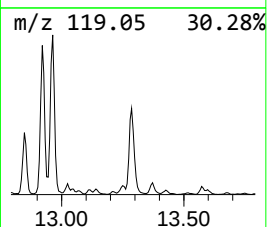
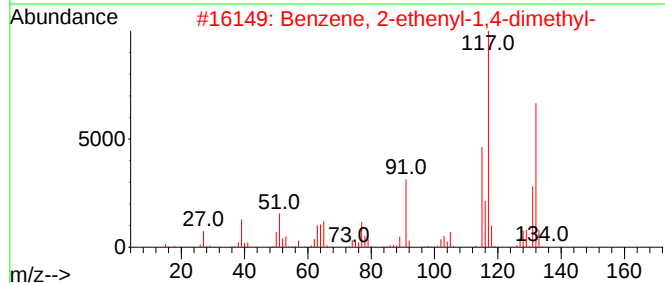
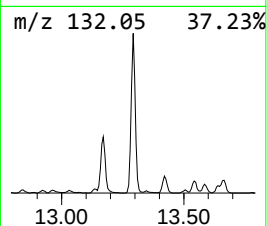
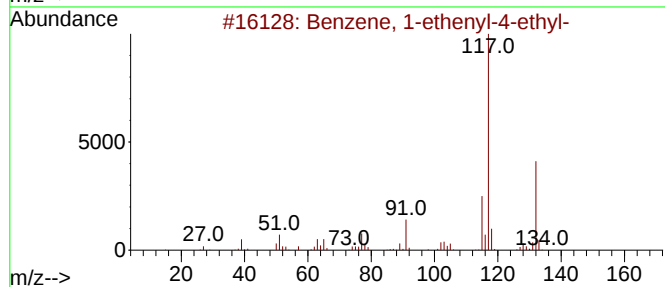
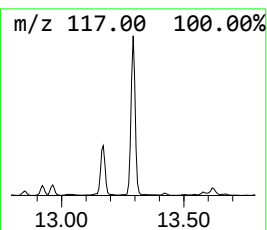
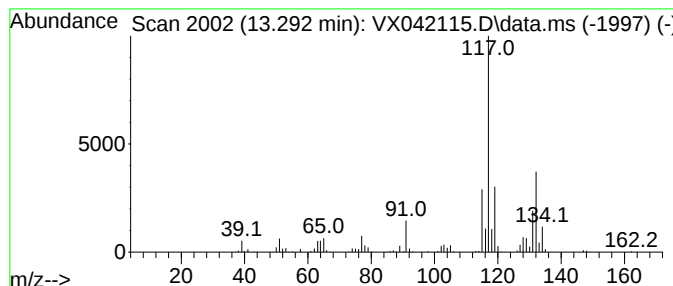
Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.292	25.11 ug/l	324721	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	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042115.D
Acq On : 01 Jul 2024 15:32
Operator : JC/MD
Sample : P3092-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW3-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.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	15.5	ug/l	200708	4	12.024	646580	50.0
Benzene, 1-ethy...	11.610	46.7	ug/l	604271	4	12.024	646580	50.0
Benzene, 1-ethy...	12.085	52.4	ug/l	677420	4	12.024	646580	50.0
Indane	12.244	41.3	ug/l	533864	4	12.024	646580	50.0
Benzene, 1-ethy...	12.610	8.6	ug/l	111125	4	12.024	646580	50.0
Indan, 1-methyl-	12.695	20.5	ug/l	264821	4	12.024	646580	50.0
Benzene, 1,2,4,...	12.921	13.1	ug/l	169182	4	12.024	646580	50.0
Benzene, 1,2,3,...	12.963	13.9	ug/l	179501	4	12.024	646580	50.0
Benzene, 2-ethe...	13.292	25.1	ug/l	324721	4	12.024	646580	50.0