

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
 Data File : VX029960.D
 Acq On : 07 Jul 2022 17:14
 Operator : JC/MD
 Sample : N3617-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-05

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\82X062822W.M
 Title : SW846 8260

Signal : TIC: VX029960.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	20	25	34	rBV	17187	31499	3.23%	0.349%
2	2.538	231	238	249	rBV	22765	40500	4.15%	0.448%
3	5.391	694	706	718	rBV2	109323	316597	32.47%	3.504%
4	5.562	722	734	746	rVB	187523	544364	55.84%	6.025%
5	5.958	788	799	814	rBV	120460	320859	32.91%	3.551%
6	6.763	922	931	945	rBV	307623	742826	76.20%	8.222%
7	8.653	1234	1241	1251	rVB	631912	974895	100.00%	10.790%
8	10.055	1466	1471	1478	rBV	614653	834291	85.58%	9.234%
9	10.287	1501	1509	1518	rVB7	4930	12296	1.26%	0.136%
10	10.963	1616	1620	1626	rBV2	7647	9774	1.00%	0.108%
11	11.085	1634	1640	1646	rBV	465055	594250	60.96%	6.577%
12	11.616	1723	1727	1731	rBV2	10057	13971	1.43%	0.155%
13	11.713	1737	1743	1747	rBV4	6722	9832	1.01%	0.109%
14	11.866	1762	1768	1770	rBV4	7062	13113	1.35%	0.145%
15	11.890	1770	1772	1780	rVB	17715	22576	2.32%	0.250%
16	12.024	1789	1794	1803	rBV	623348	772190	79.21%	8.547%
17	12.177	1815	1819	1824	rVB	22665	28442	2.92%	0.315%
18	12.244	1824	1830	1837	rBV2	363099	469407	48.15%	5.195%
19	12.317	1837	1842	1849	rVB	41379	56891	5.84%	0.630%
20	12.408	1852	1857	1864	rBV6	7370	13295	1.36%	0.147%
21	12.646	1891	1896	1901	rVV	123058	154877	15.89%	1.714%
22	12.701	1901	1905	1912	rVV2	112628	168316	17.27%	1.863%
23	12.762	1912	1915	1919	rVV2	10978	14961	1.53%	0.166%
24	12.841	1919	1928	1934	rVB	34561	57763	5.93%	0.639%
25	12.920	1934	1941	1946	rBV6	16765	36961	3.79%	0.409%
26	13.030	1954	1959	1965	rVB3	39788	61281	6.29%	0.678%
27	13.140	1965	1977	1980	rBV	42433	66791	6.85%	0.739%
28	13.195	1980	1986	1992	rVB2	28590	49456	5.07%	0.547%
29	13.292	1998	2002	2007	rBV	98057	121764	12.49%	1.348%
30	13.347	2007	2011	2015	rVB2	29462	37311	3.83%	0.413%
31	13.396	2015	2019	2020	rBV2	14146	16025	1.64%	0.177%
32	13.426	2020	2024	2032	rVB3	34461	54159	5.56%	0.599%
33	13.506	2032	2037	2039	rBV4	10764	16772	1.72%	0.186%
34	13.542	2039	2043	2047	rVV	28132	38751	3.97%	0.429%
35	13.585	2047	2050	2053	rVV	43531	55781	5.72%	0.617%
36	13.640	2053	2059	2060	rVV2	69344	116667	11.97%	1.291%

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Integrator: RTE

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Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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37	13.664	2060	2063	2073	rVB	142131	199391	20.45%	2.207%
38	13.756	2074	2078	2080	rVV	15984	19666	2.02%	0.218%
39	13.798	2081	2085	2090	rVV4	13045	25603	2.63%	0.283%
40	13.853	2090	2094	2099	rVB	110347	130495	13.39%	1.444%
41	13.926	2099	2106	2110	rBV	164493	208915	21.43%	2.312%
42	13.969	2110	2113	2117	rVV2	30999	41368	4.24%	0.458%
43	14.012	2117	2120	2122	rVB	13196	14515	1.49%	0.161%
44	14.073	2126	2130	2133	rBV	47064	61829	6.34%	0.684%
45	14.103	2133	2135	2141	rVB4	13525	16445	1.69%	0.182%
46	14.213	2148	2153	2163	rVV2	57313	105282	10.80%	1.165%
47	14.286	2163	2165	2167	rVV3	12603	13008	1.33%	0.144%
48	14.323	2167	2171	2175	rVV2	48080	61225	6.28%	0.678%
49	14.377	2175	2180	2184	rVB2	81829	113619	11.65%	1.258%
50	14.420	2184	2187	2192	rVB2	19029	26040	2.67%	0.288%
51	14.481	2192	2197	2202	rBV2	117174	170997	17.54%	1.893%
52	14.530	2202	2205	2210	rVB4	12487	19605	2.01%	0.217%
53	14.615	2210	2219	2222	rBV4	58807	130232	13.36%	1.441%
54	14.676	2225	2229	2233	rVB	52696	73718	7.56%	0.816%
55	14.725	2233	2237	2242	rBV2	24272	36424	3.74%	0.403%
56	14.792	2242	2248	2250	rBV3	37005	62615	6.42%	0.693%
57	14.816	2250	2252	2255	rVV3	30497	40829	4.19%	0.452%
58	14.847	2255	2257	2261	rVB3	29349	32616	3.35%	0.361%
59	14.902	2261	2266	2272	rBV4	29888	66065	6.78%	0.731%
60	15.048	2286	2290	2294	rVB4	14621	22331	2.29%	0.247%
61	15.182	2306	2312	2320	rBV3	58504	109756	11.26%	1.215%
62	15.255	2320	2324	2331	rVV6	20526	37587	3.86%	0.416%
63	15.322	2331	2335	2339	rVV5	7097	11684	1.20%	0.129%
64	15.377	2339	2344	2347	rVV4	16950	26500	2.72%	0.293%
65	15.481	2358	2361	2364	rBV4	10126	12723	1.31%	0.141%
66	15.524	2364	2368	2373	rVB9	14512	26881	2.76%	0.298%
67	15.603	2377	2381	2387	rVB7	9207	18866	1.94%	0.209%
68	15.658	2387	2390	2397	rVB9	11189	21189	2.17%	0.235%
69	15.725	2397	2401	2402	rBV3	10904	12018	1.23%	0.133%
70	15.810	2411	2415	2418	rBV2	23640	35438	3.64%	0.392%
71	15.847	2418	2421	2426	rVB5	18835	30226	3.10%	0.335%
72	15.987	2439	2444	2448	rBV2	31481	51708	5.30%	0.572%
73	16.328	2495	2500	2508	rVB5	13421	27941	2.87%	0.309%
74	16.426	2510	2516	2521	rVB7	5940	10211	1.05%	0.113%
75	16.530	2529	2533	2540	rBV9	8429	20012	2.05%	0.221%

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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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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Title : SW846 8260

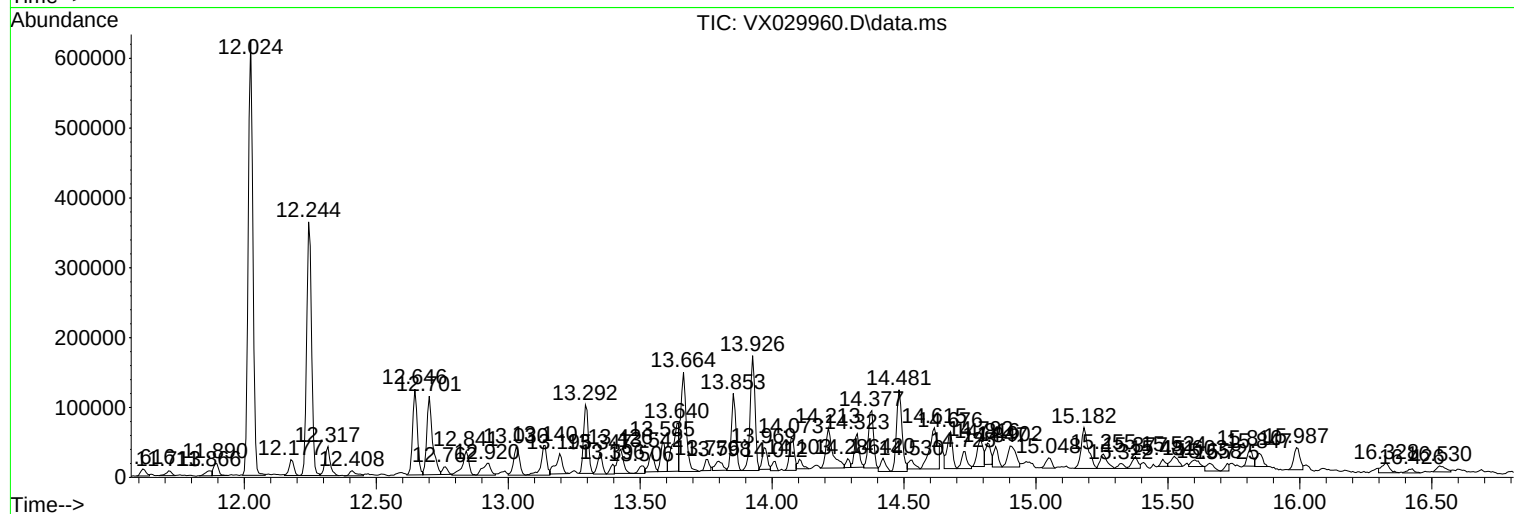
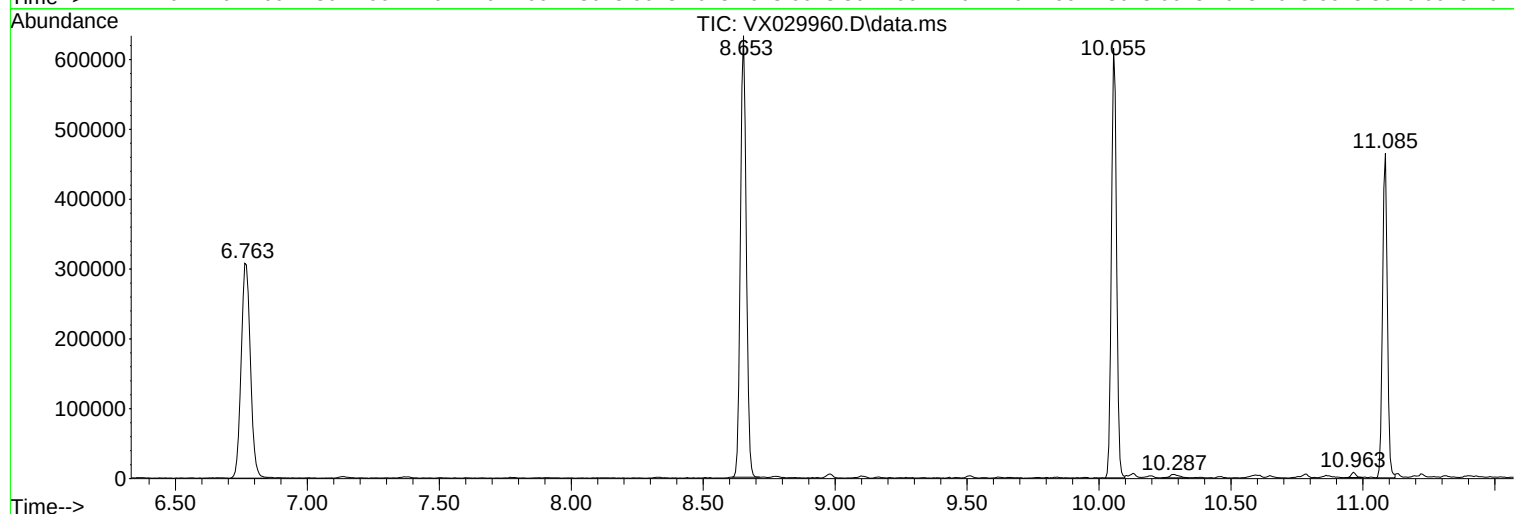
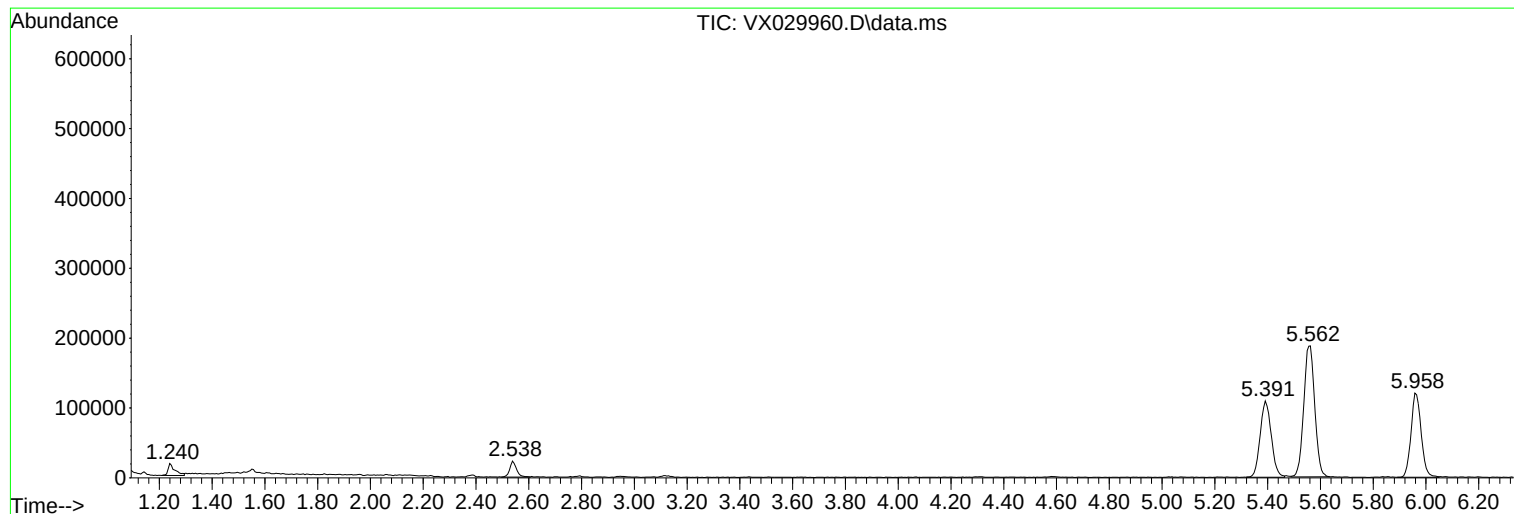
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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-05

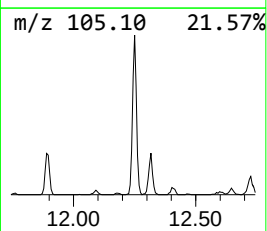
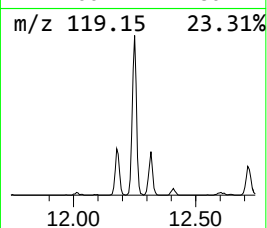
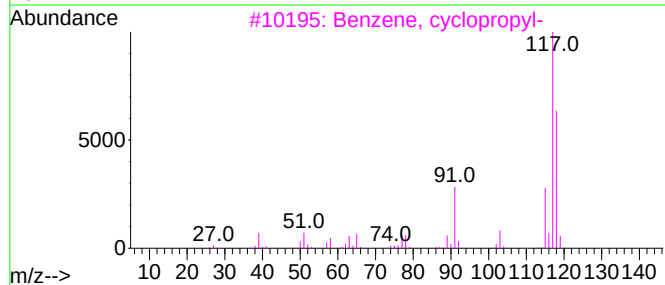
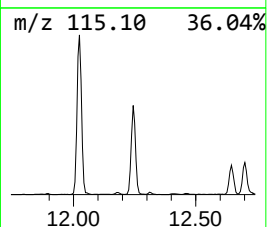
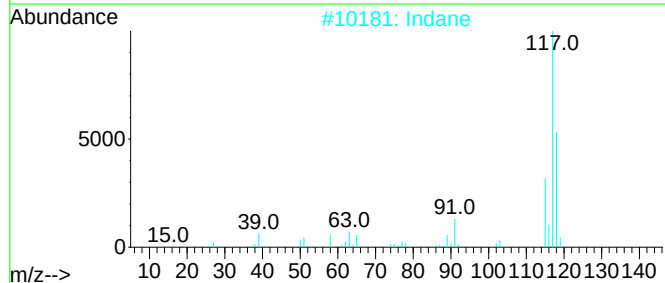
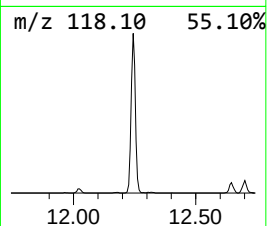
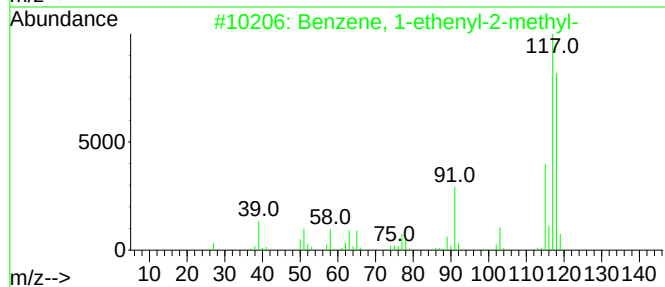
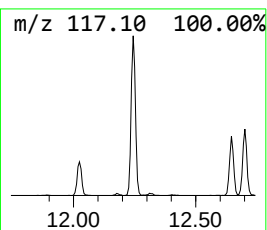
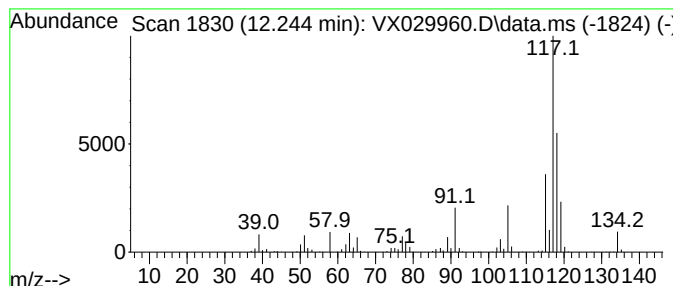
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	30.39 ug/l	469407	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			Indane	118	C9H10	000496-11-7	70
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64



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Instrument :
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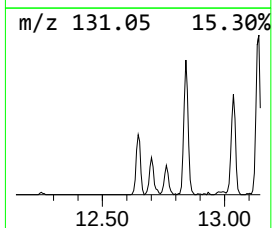
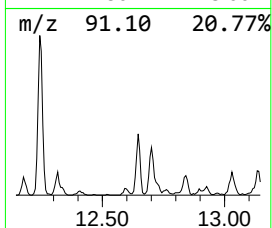
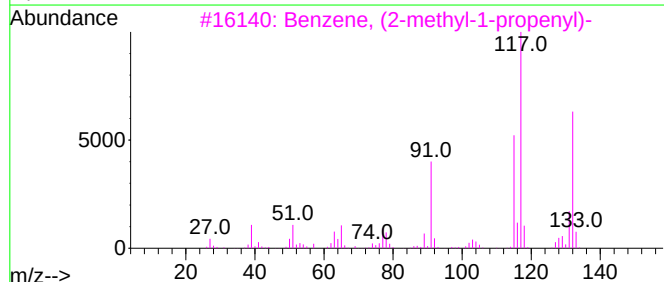
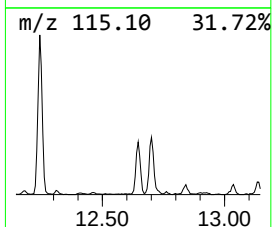
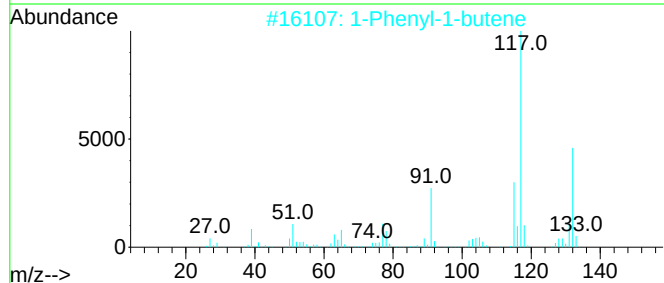
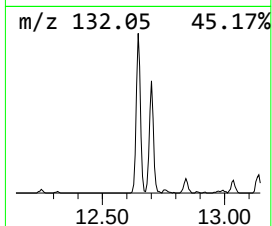
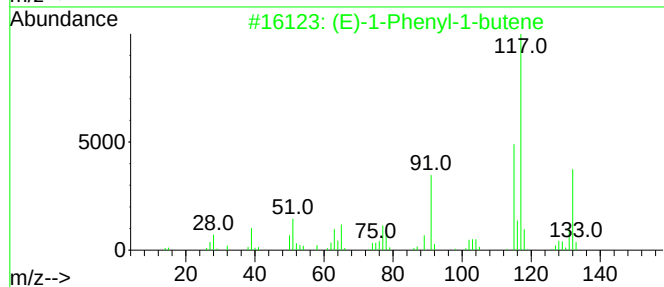
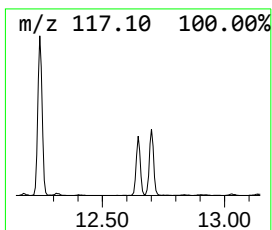
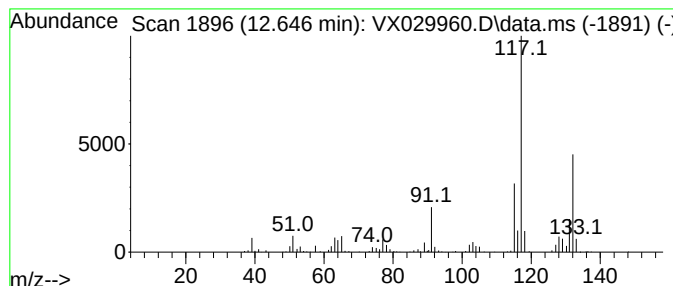
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Quant Title : SW846 8260

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

Peak Number 2 (E)-1-Phenyl-1-butene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	10.03 ug/l	154877	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	95
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	94
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	93
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	93
5			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92



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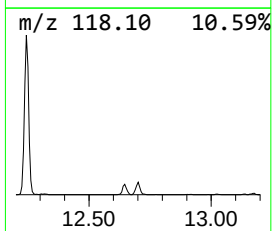
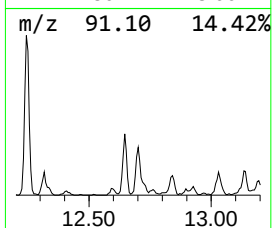
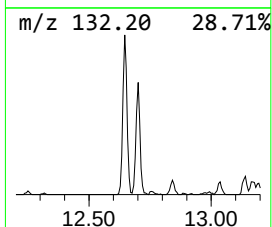
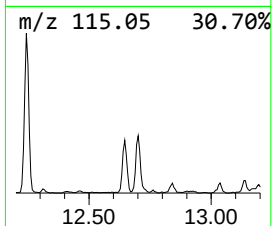
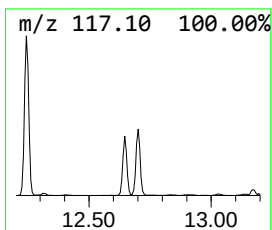
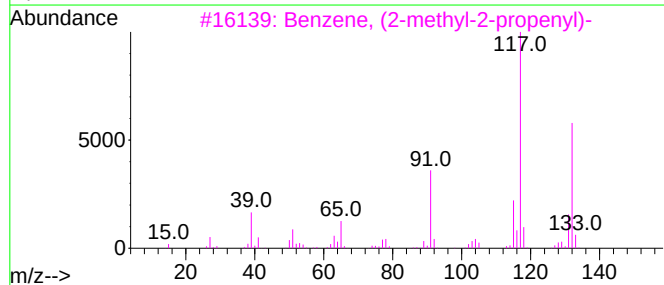
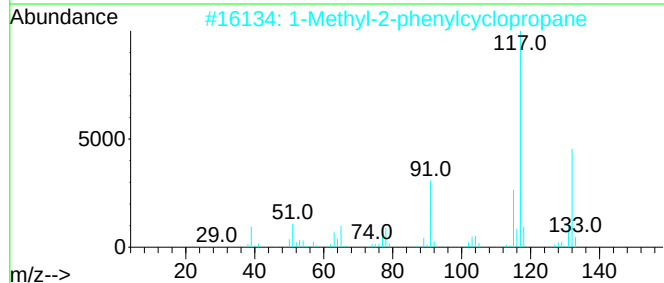
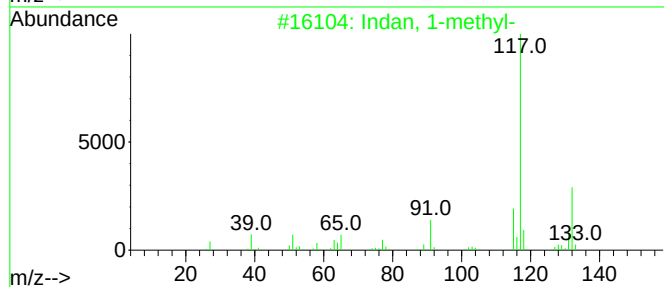
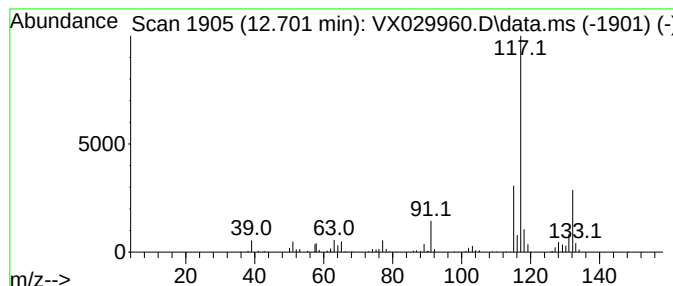
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Quant Title : SW846 8260

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

Peak Number 3 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	10.90 ug/l	168316	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			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
3			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	90
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



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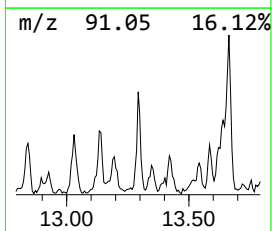
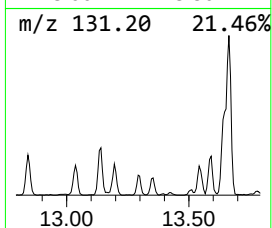
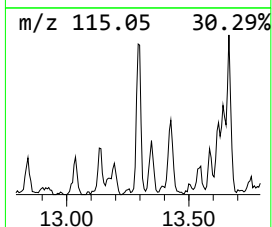
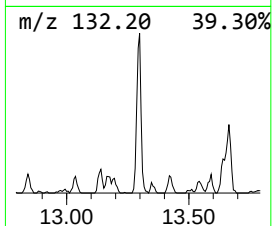
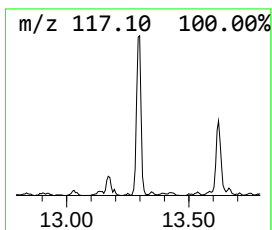
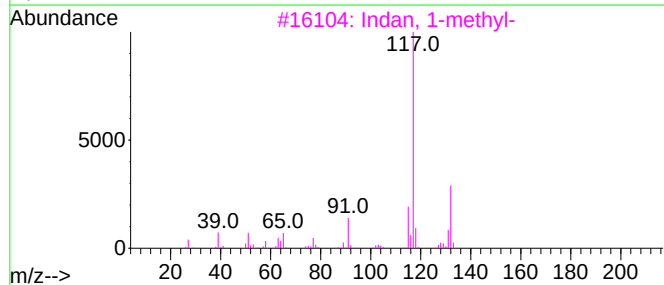
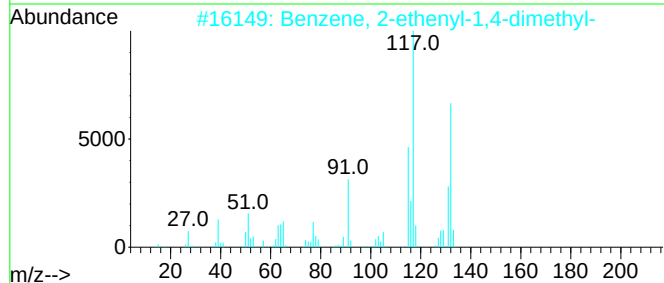
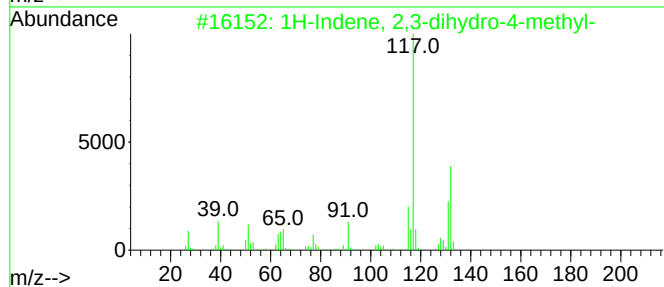
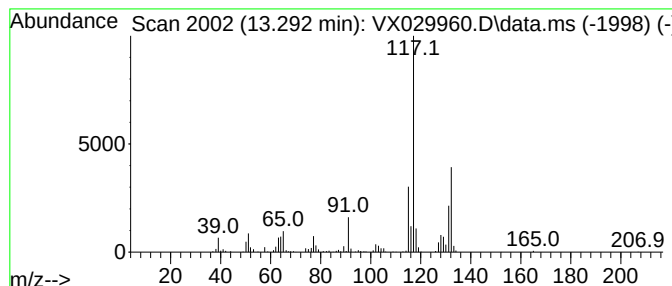
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	7.88 ug/l	121764	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	87
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
3		Indan, 1-methyl-	132	C10H12	000767-58-8	81
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	80
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-05

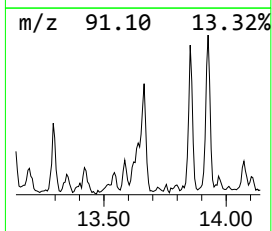
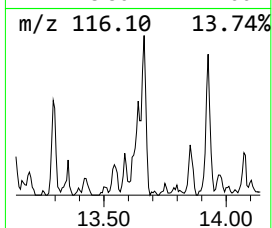
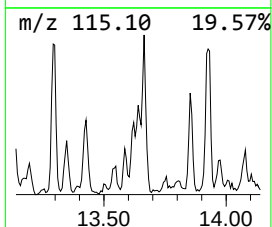
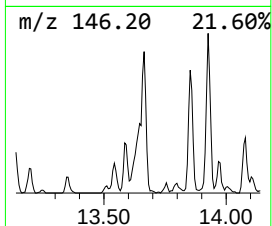
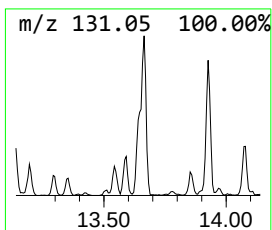
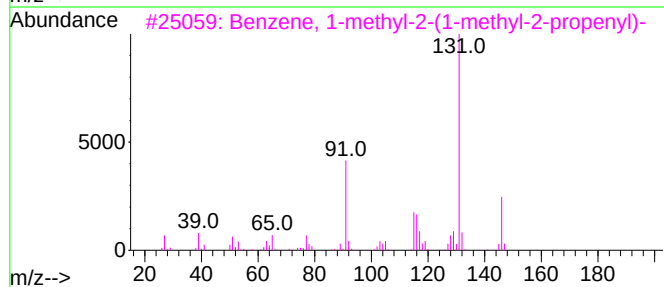
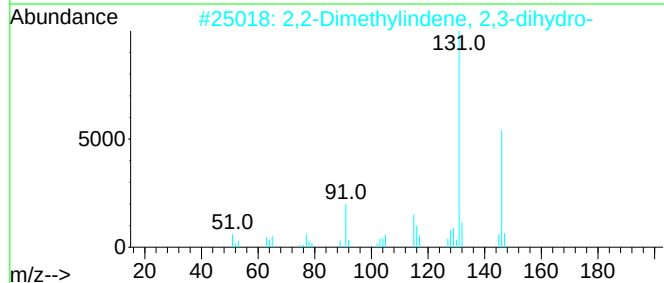
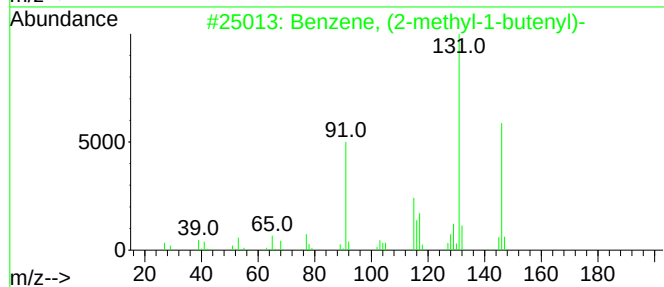
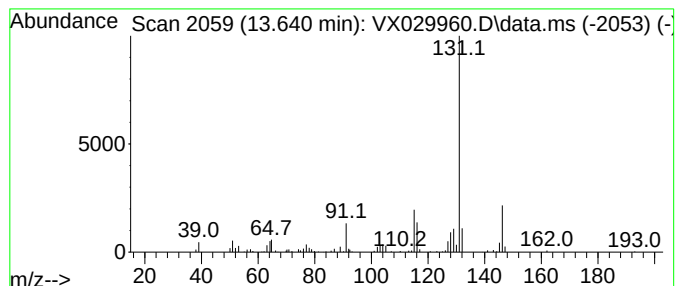
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	7.55 ug/l	116667	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			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
3			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-05

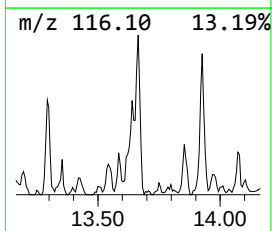
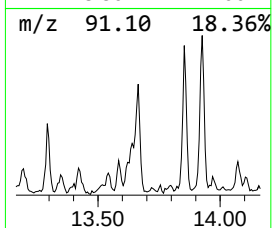
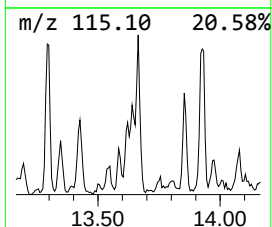
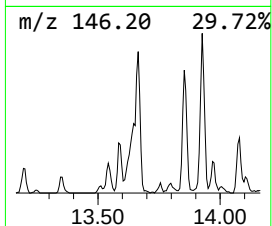
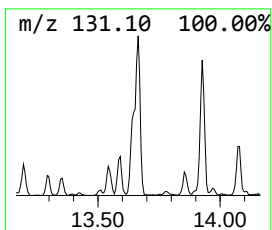
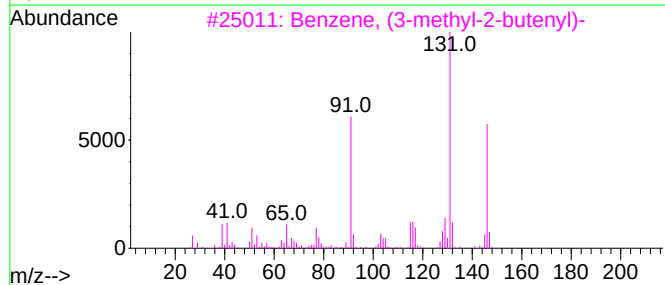
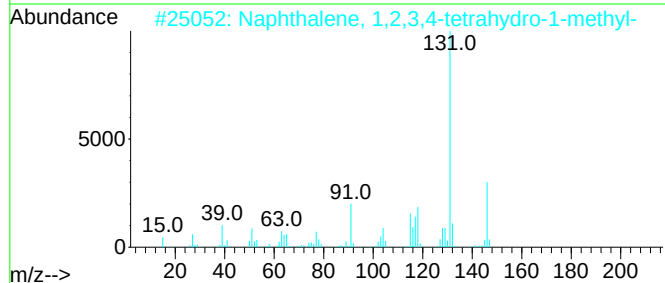
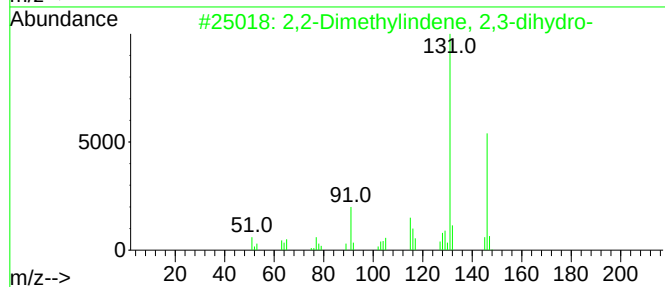
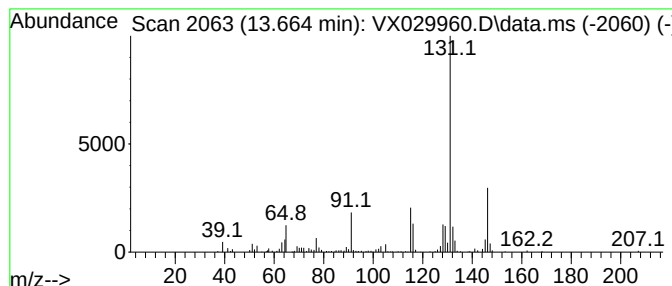
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Quant Title : SW846 8260

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

Peak Number 6 2,2-Dimethylindene, 2,3-dih... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	12.91 ug/l	199391	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

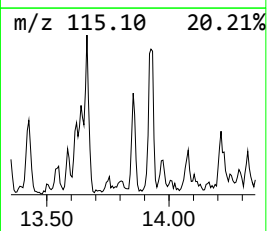
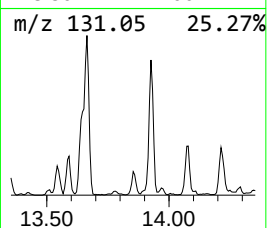
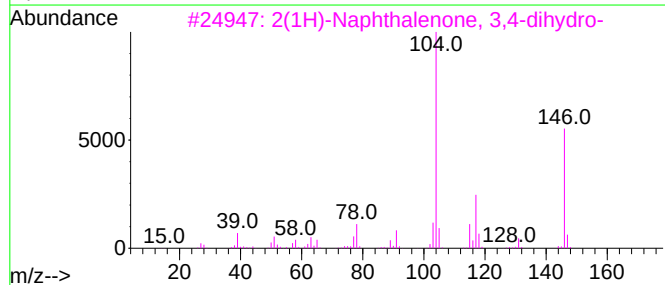
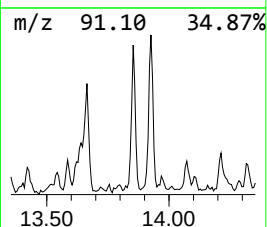
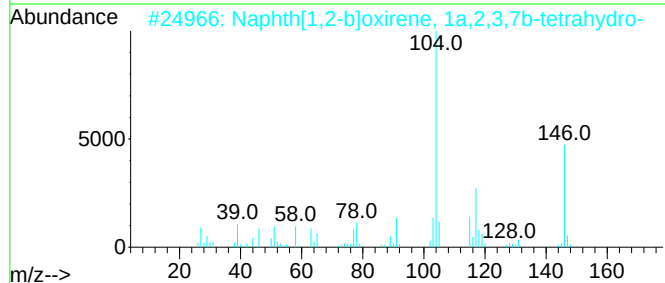
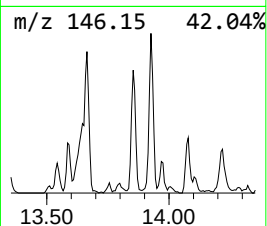
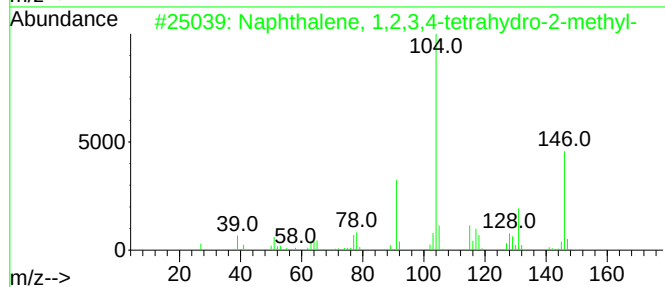
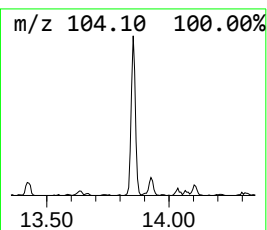
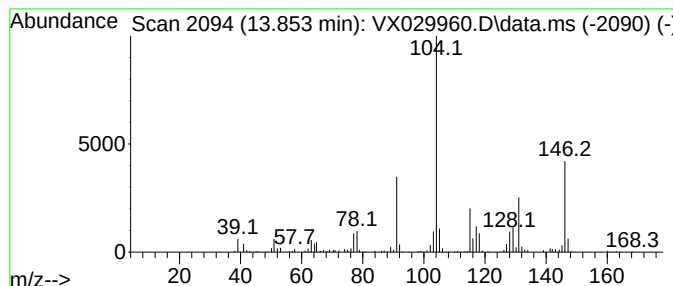
Instrument :
MSVOA_X
ClientSampleId :
MW-05

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.853	8.45 ug/l	130495	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	93
2	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	58
4	7-Azabicyclo[4.2.2]deca-2,4,9-tr...	147	C9H9NO	017198-06-0	35
5	2-Phenylethylamine, N,N-di(trifl...	313	C12H9F6NO2	1000463-67-7	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

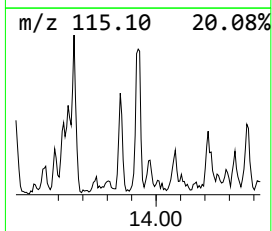
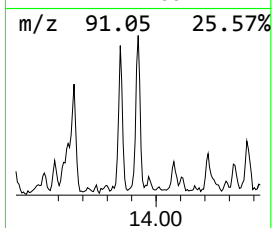
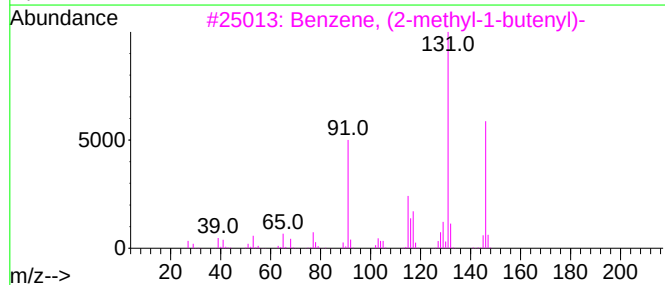
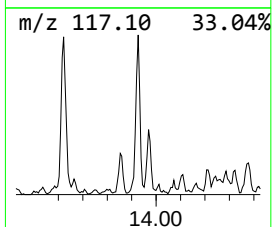
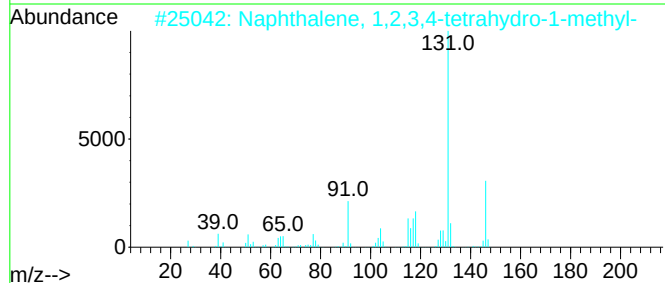
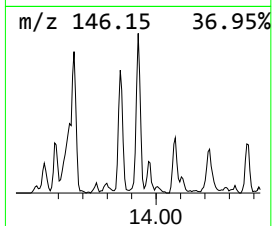
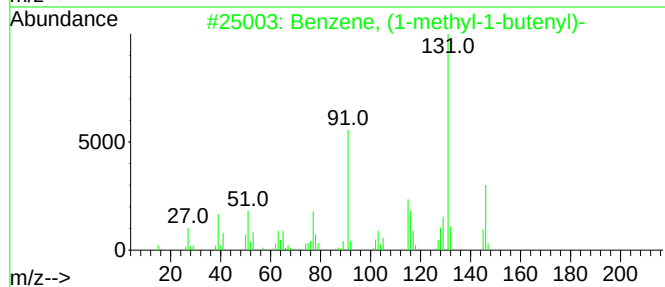
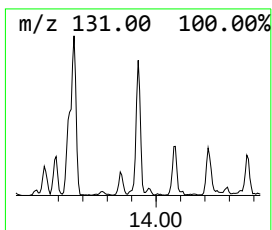
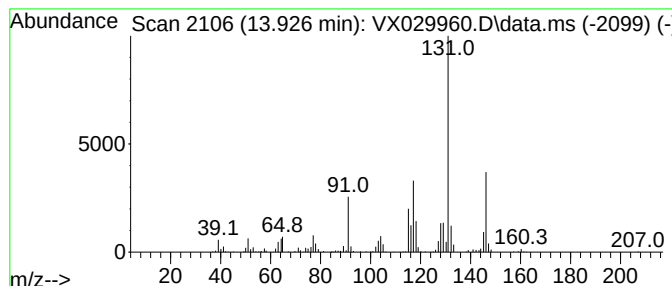
Instrument :
MSVOA_X
ClientSampleId :
MW-05

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-butenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.926	13.53 ug/l	208915	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-		146	C11H14	053172-84-2	93
2	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001559-81-5	93
3	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	93
4	1H-Indene, 2,3-dihydro-1,6-dimet...		146	C11H14	017059-48-2	93
5	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-05

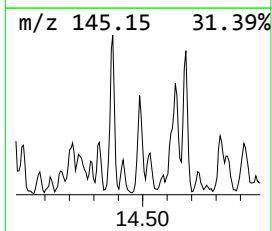
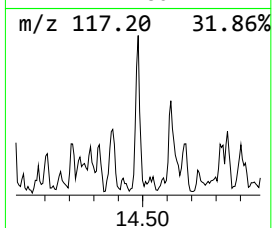
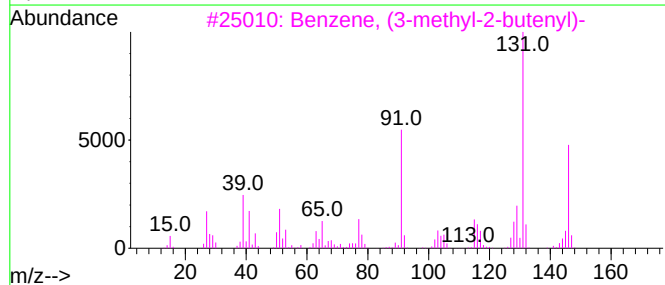
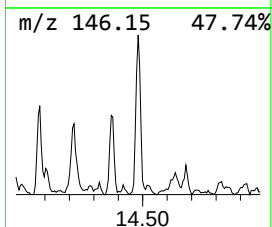
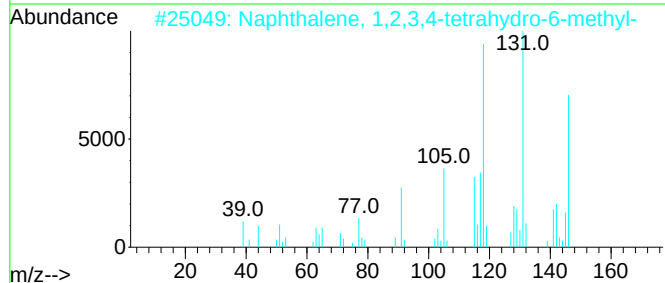
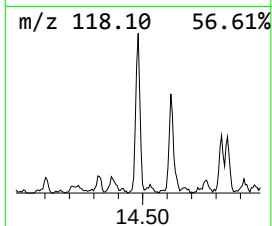
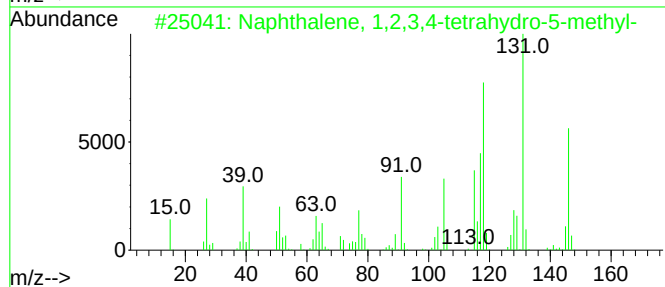
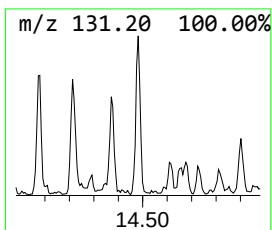
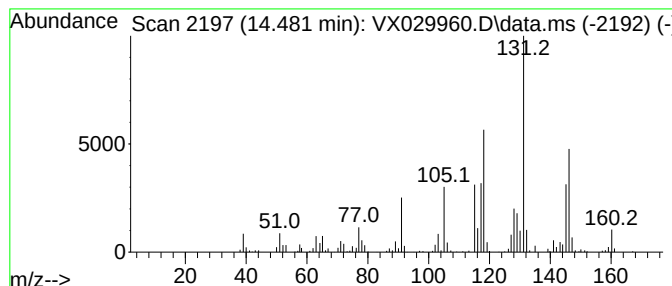
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	11.07 ug/l	170997	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	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
4			2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	58
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

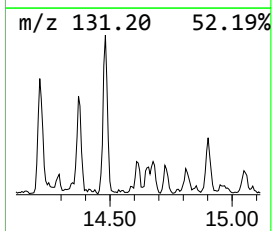
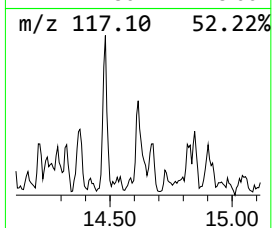
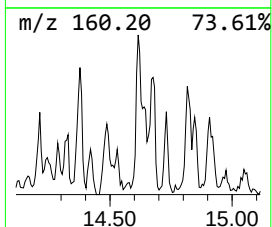
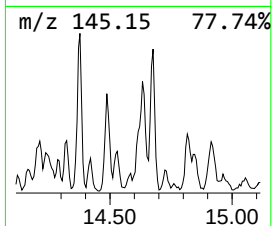
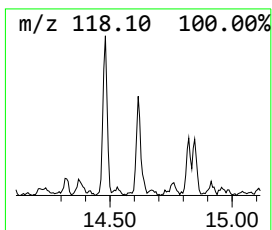
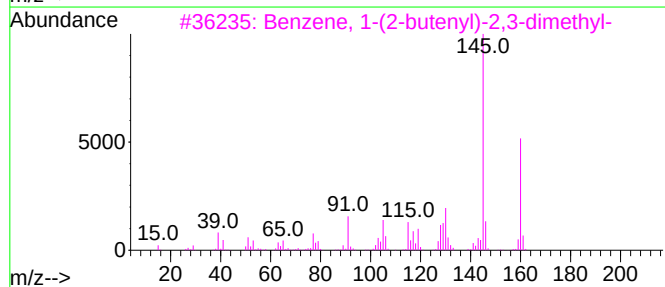
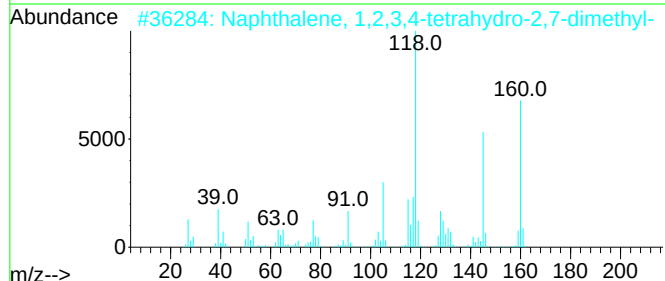
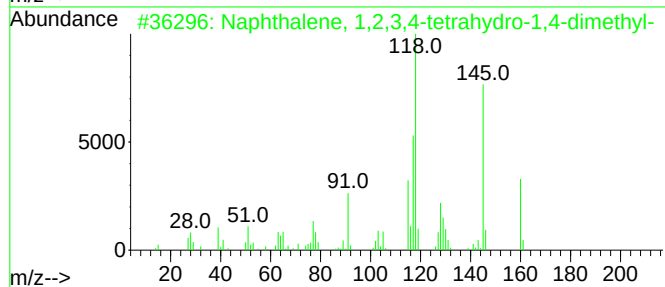
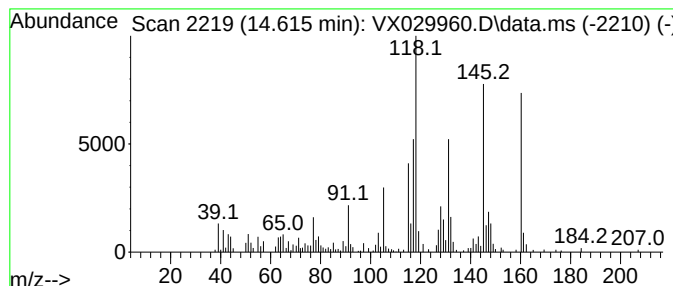
Instrument :
MSVOA_X
ClientSampleId :
MW-05

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
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.	
14.615	8.43 ug/l	130232	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	89
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	64
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	62
5	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070722\
Data File : VX029960.D
Acq On : 07 Jul 2022 17:14
Operator : JC/MD
Sample : N3617-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-05

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.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-ethe...	12.244	30.4	ug/l	469407	4	12.024	772190	50.0
(E)-1-Phenyl-1-...	12.646	10.0	ug/l	154877	4	12.024	772190	50.0
Indan, 1-methyl-	12.701	10.9	ug/l	168316	4	12.024	772190	50.0
1H-Indene, 2,3-...	13.292	7.9	ug/l	121764	4	12.024	772190	50.0
Benzene, (2-met...	13.640	7.5	ug/l	116667	4	12.024	772190	50.0
2,2-Dimethylind...	13.664	12.9	ug/l	199391	4	12.024	772190	50.0
Naphthalene, 1,...	13.853	8.4	ug/l	130495	4	12.024	772190	50.0
Benzene, (1-met...	13.926	13.5	ug/l	208915	4	12.024	772190	50.0
Naphthalene, 1,...	14.481	11.1	ug/l	170997	4	12.024	772190	50.0
Naphthalene, 1,...	14.615	8.4	ug/l	130232	4	12.024	772190	50.0