

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
 Data File : VX031541.D
 Acq On : 22 Sep 2022 01:11
 Operator : JC/MD
 Sample : N4614-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-43

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\82X091922W.M
 Title : SW846 8260

Signal : TIC: VX031541.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.306	518	528	540	rVB	299109	847158	2.59%	0.464%
2	5.477	712	720	728	rBV5	99924	334815	1.02%	0.183%
3	6.043	805	813	823	rVB	336366	855810	2.62%	0.468%
4	6.763	924	931	939	rBV	134594	335248	1.03%	0.183%
5	7.379	1018	1032	1042	rBV4	235511	647819	1.98%	0.355%
6	8.647	1235	1240	1247	rVB	479833	764304	2.34%	0.418%
7	10.055	1461	1471	1477	rBV	299164	447903	1.37%	0.245%
8	10.207	1489	1496	1500	rBV	23732420	32692956	100.00%	17.892%
9	10.311	1506	1513	1518	rVB	13195390	16790690	51.36%	9.189%
10	10.646	1562	1568	1577	rVB	3965029	4629986	14.16%	2.534%
11	10.963	1614	1620	1628	rBV	1757344	2092443	6.40%	1.145%
12	11.085	1635	1640	1645	rBV	516176	655080	2.00%	0.359%
13	11.305	1671	1676	1683	rVB	5416810	6344154	19.41%	3.472%
14	11.402	1683	1692	1696	rVV	5718974	10013317	30.63%	5.480%
15	11.451	1696	1700	1707	rVB	2244677	2740399	8.38%	1.500%
16	11.616	1721	1727	1736	rBV	6880384	8230945	25.18%	4.505%
17	11.756	1744	1750	1755	rBV	18358498	23623084	72.26%	12.928%
18	12.024	1789	1794	1799	rVB2	642230	865824	2.65%	0.474%
19	12.091	1799	1805	1810	rBV	9350977	11083734	33.90%	6.066%
20	12.244	1825	1830	1838	rBV	9037054	12444045	38.06%	6.810%
21	12.317	1838	1842	1849	rVV3	1395075	2212467	6.77%	1.211%
22	12.408	1852	1857	1862	rBV2	389009	521620	1.60%	0.285%
23	12.463	1862	1866	1872	rVB	703026	825639	2.53%	0.452%
24	12.530	1872	1877	1880	rBV	1274719	1802942	5.51%	0.987%
25	12.555	1880	1881	1886	rVB	912406	764599	2.34%	0.418%
26	12.615	1886	1891	1894	rBV	3149336	3670327	11.23%	2.009%
27	12.646	1894	1896	1900	rVB	541868	562308	1.72%	0.308%
28	12.701	1900	1905	1913	rBV2	1655569	2326531	7.12%	1.273%
29	12.847	1925	1929	1934	rVB	707775	871486	2.67%	0.477%
30	12.920	1934	1941	1945	rBV	1859621	2355477	7.20%	1.289%
31	12.963	1945	1948	1954	rVB	1845061	2141973	6.55%	1.172%
32	13.170	1978	1982	1987	rVB	1633588	1888778	5.78%	1.034%
33	13.292	1998	2002	2008	rVB2	3595430	4747081	14.52%	2.598%
34	13.426	2019	2024	2032	rVB	477241	617922	1.89%	0.338%
35	13.585	2046	2050	2056	rVB3	307844	497876	1.52%	0.272%
36	13.664	2061	2063	2069	rVB2	281468	394389	1.21%	0.216%

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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

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Title : SW846 8260

37	13.780	2076	2082	2087	rBV	11212614	14223592	43.51%	7.784%
38	13.871	2092	2097	2101	rVB	308323	439503	1.34%	0.241%
39	14.640	2218	2223	2229	rVB	2081732	2513602	7.69%	1.376%
40	14.780	2241	2246	2255	rVB	1980942	2488432	7.61%	1.362%
41	15.615	2373	2383	2386	rBV	268920	419564	1.28%	0.230%

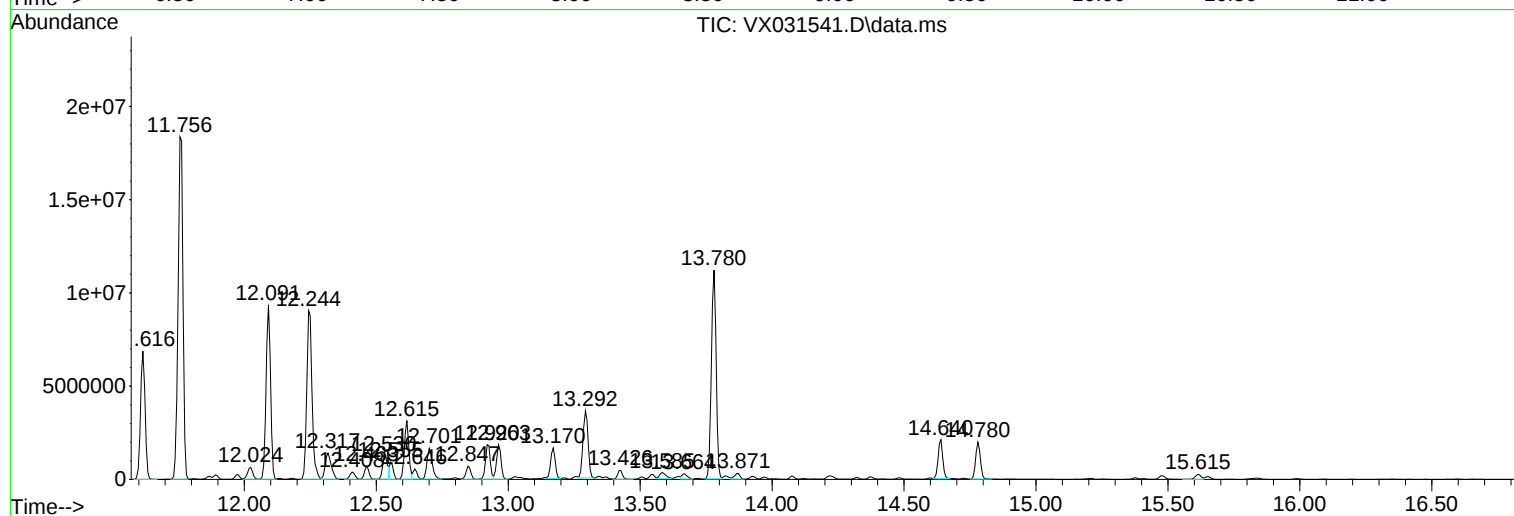
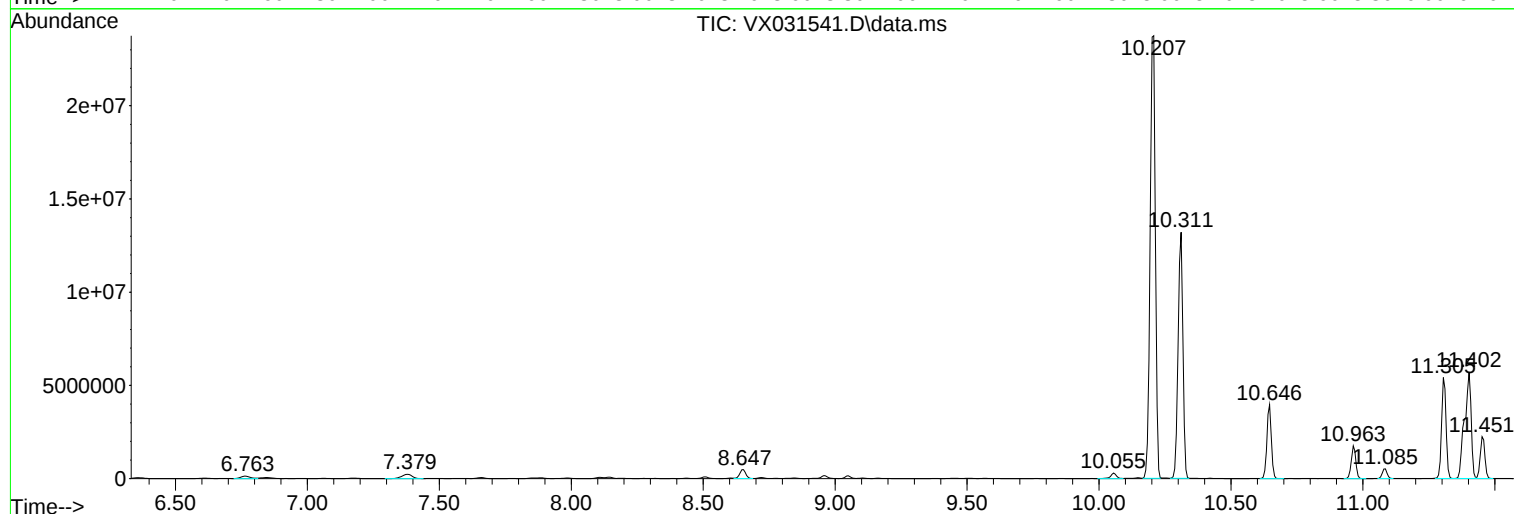
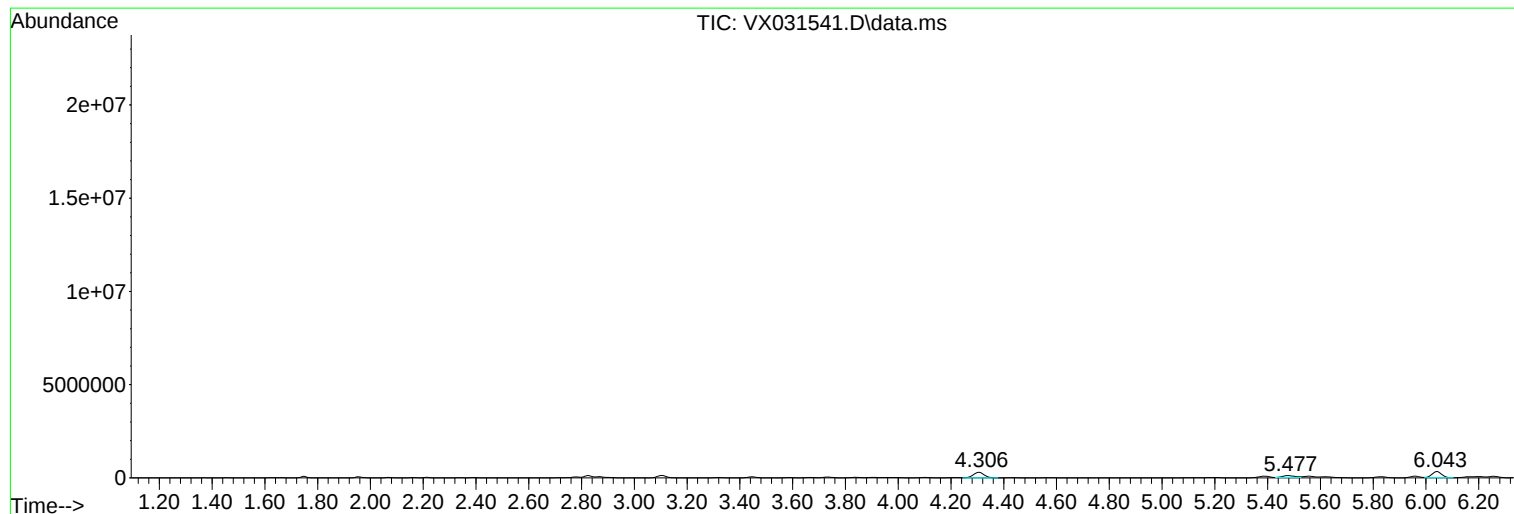
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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



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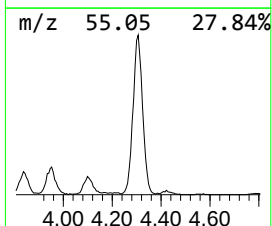
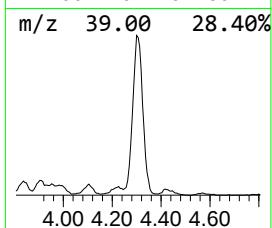
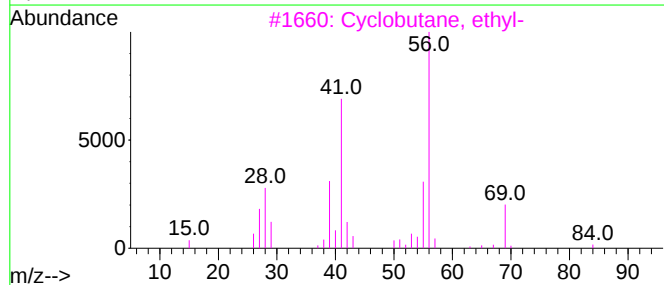
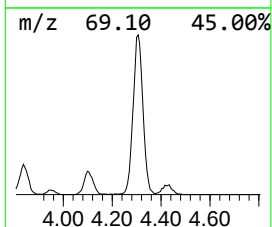
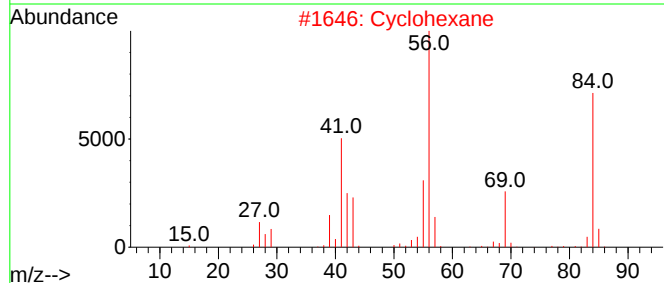
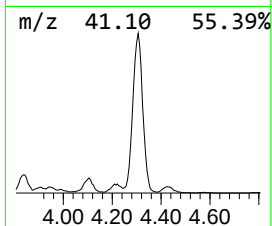
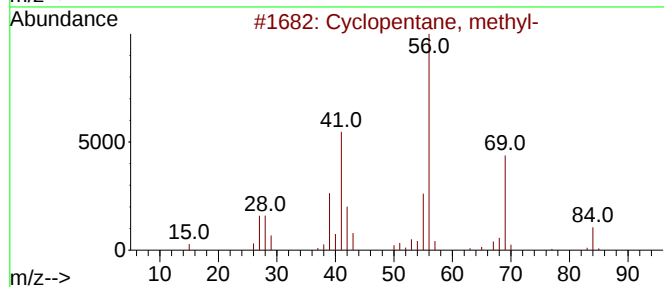
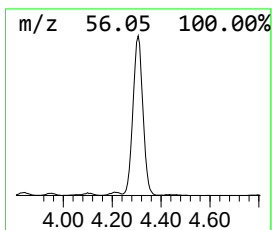
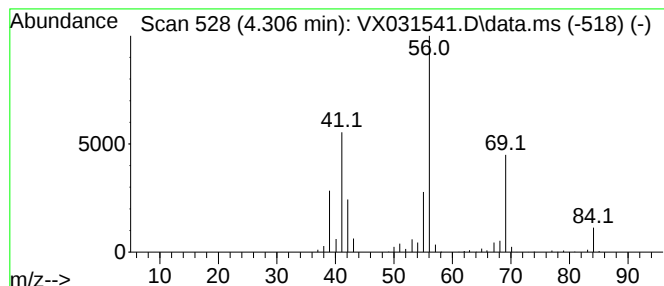
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	126.51 ug/l	847158	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclohexane	84	C6H12	000110-82-7	78
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
5			Cyclobutane	56	C4H8	000287-23-0	53



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ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

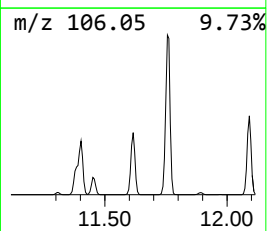
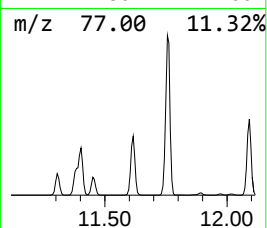
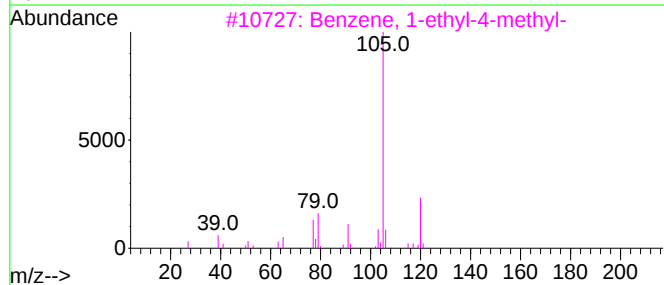
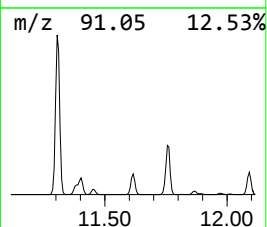
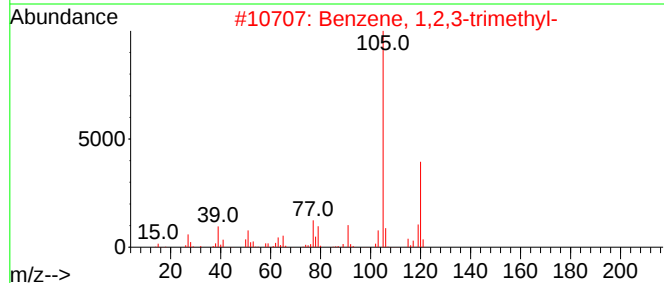
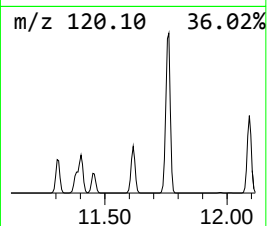
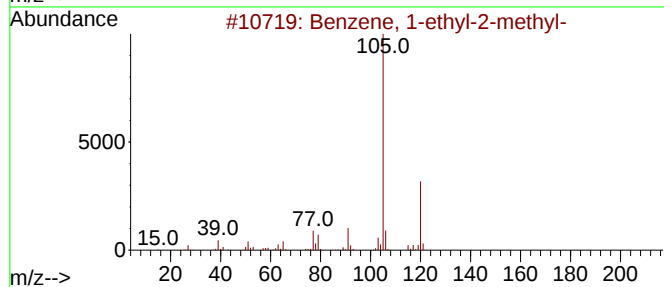
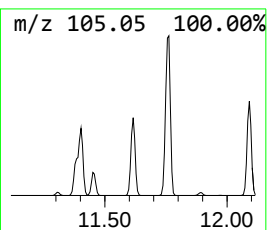
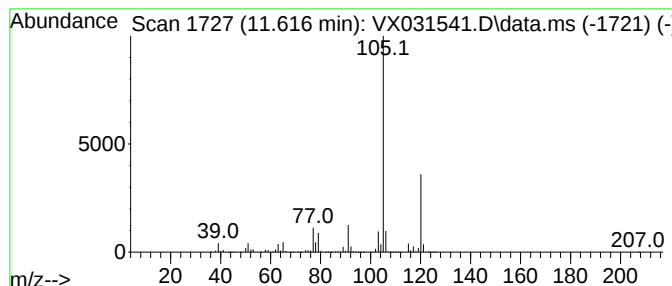
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	475.32 ug/l	8230950	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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Instrument :
MSVOA_X
ClientSampleId :
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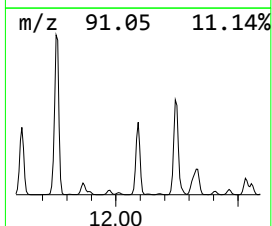
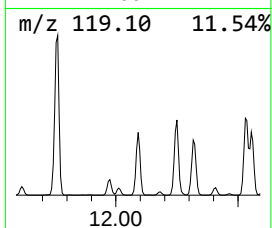
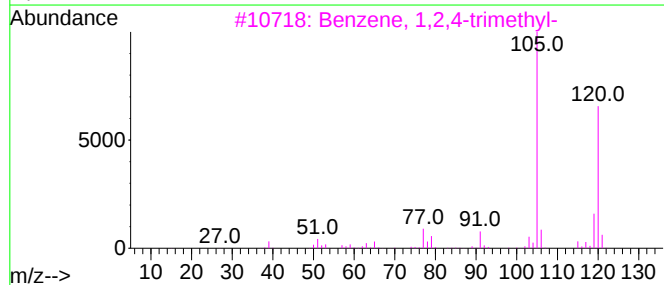
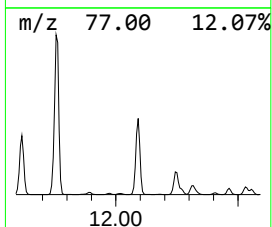
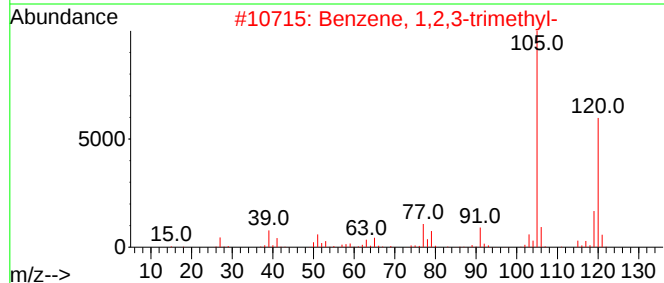
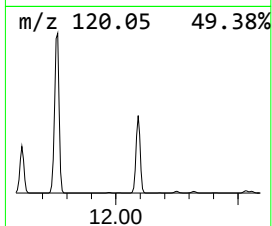
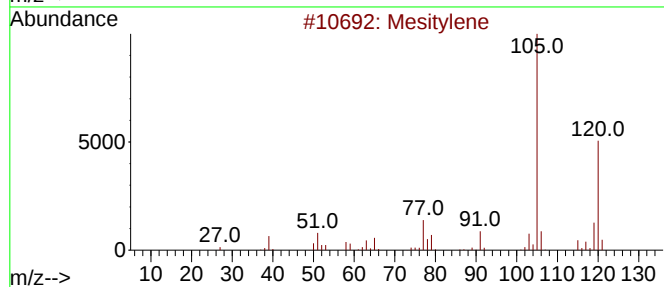
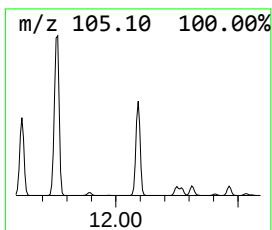
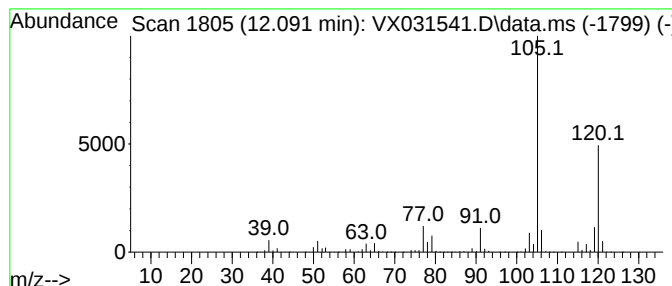
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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.
12.091	640.07 ug/l	11083700	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	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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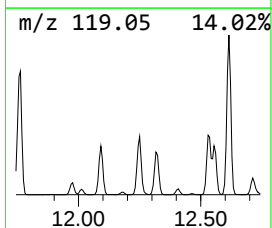
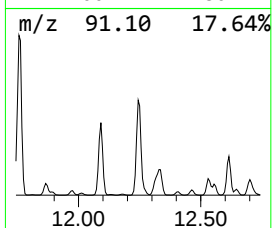
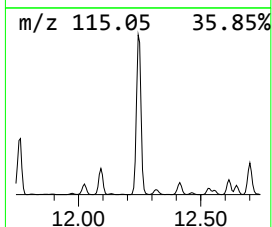
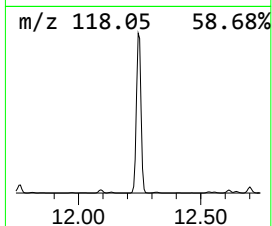
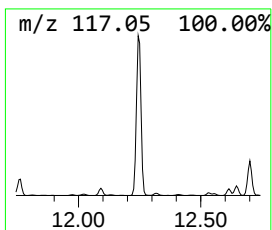
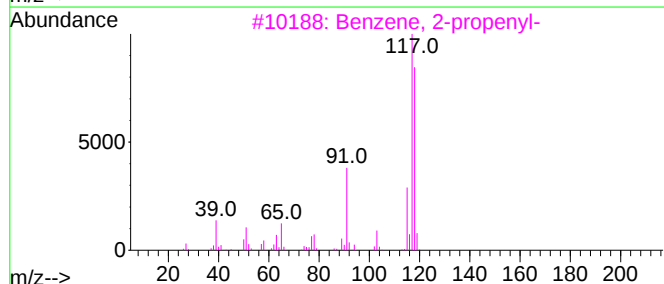
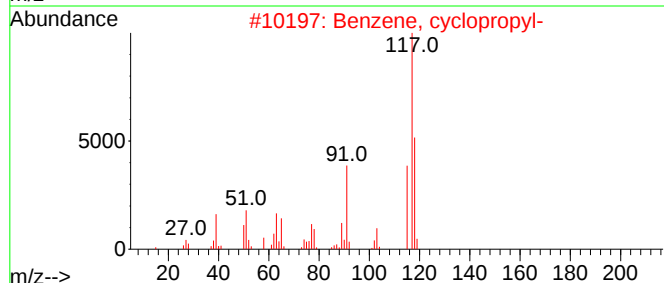
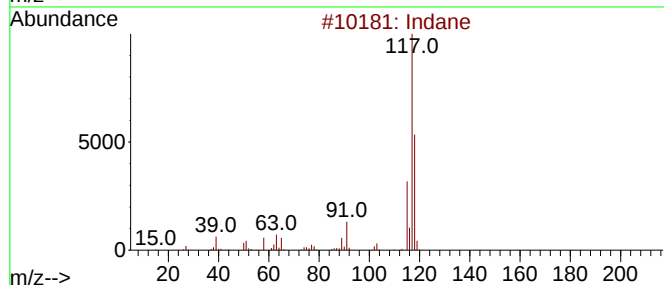
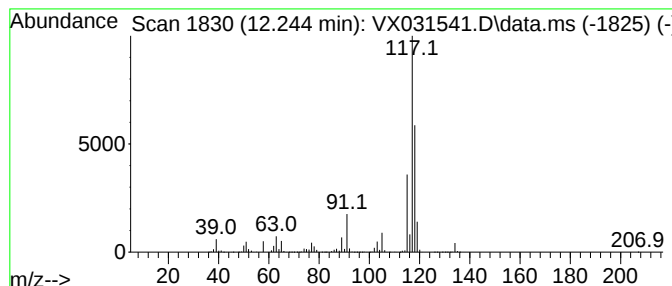
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	718.62 ug/l	12444000	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, 2-propenyl-	118	C9H10	000300-57-2	74
4			Benzeneethanol, .beta.-ethenyl-....	162	C11H14O	040596-14-3	50
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	49



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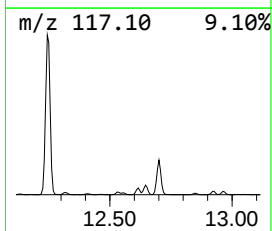
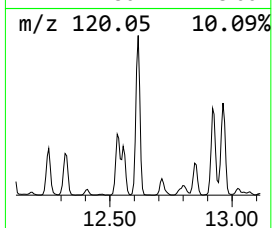
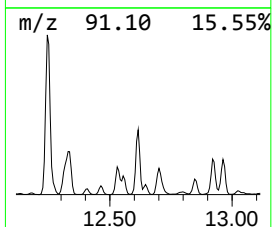
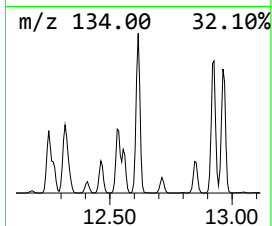
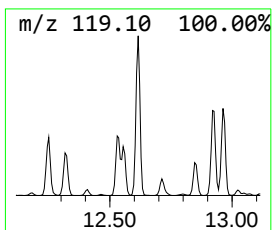
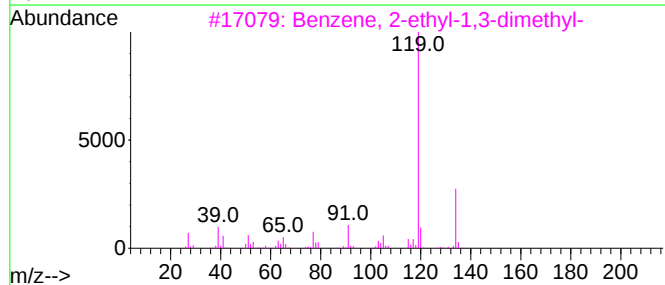
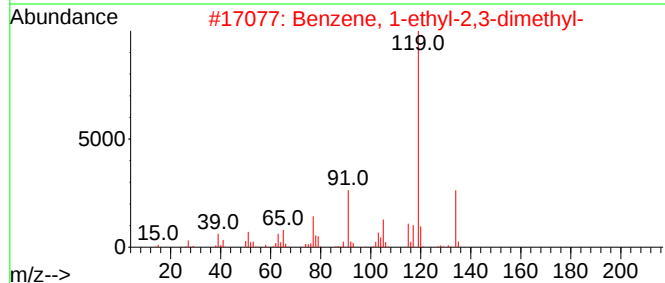
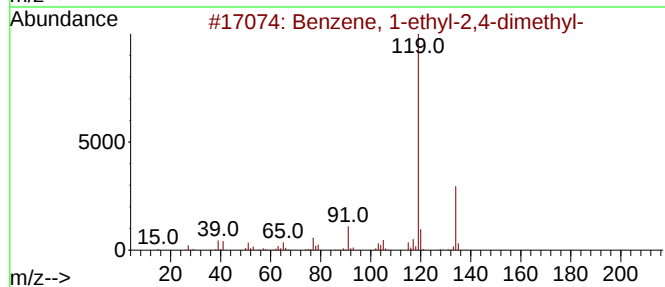
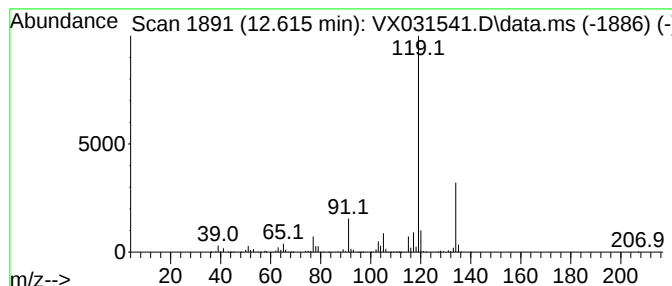
Instrument :
MSVOA_X
ClientSampleId :
MW-43

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.616	211.96 ug/l	3670330	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, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	95
4	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	95
5	o-Cymene		134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

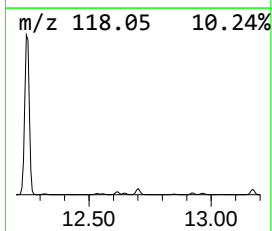
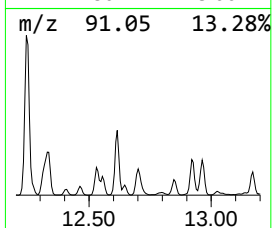
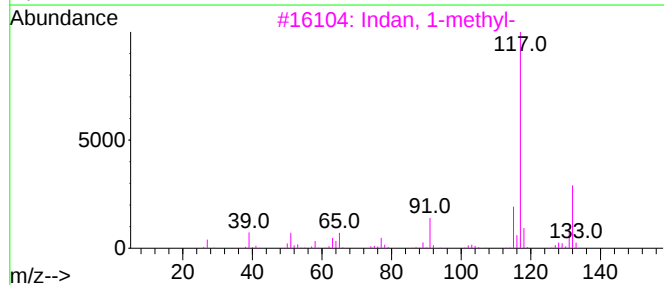
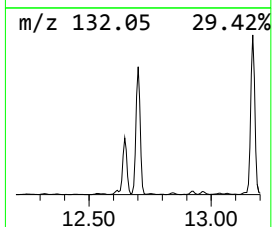
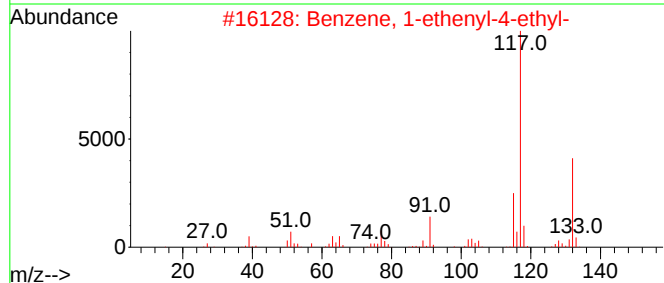
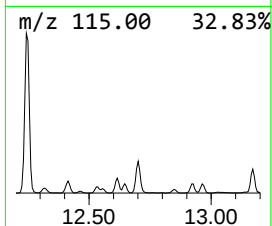
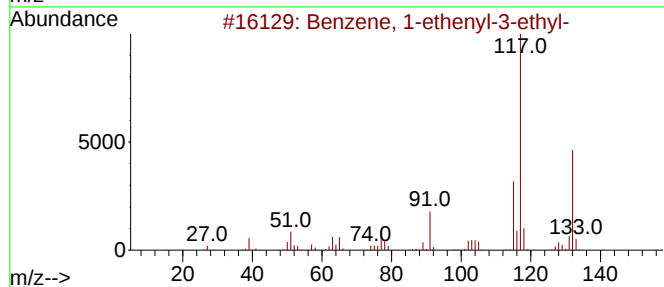
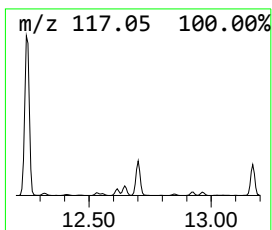
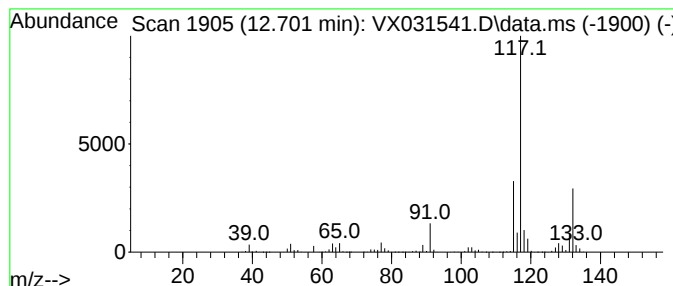
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	134.35 ug/l	2326530	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, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

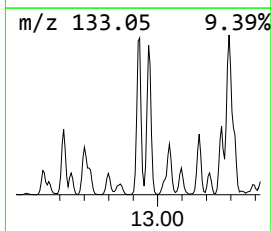
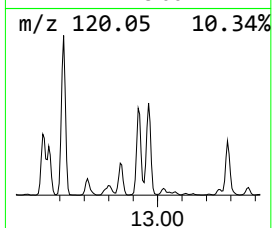
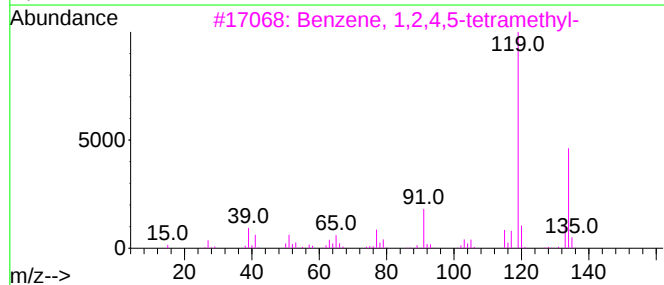
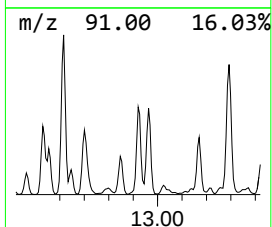
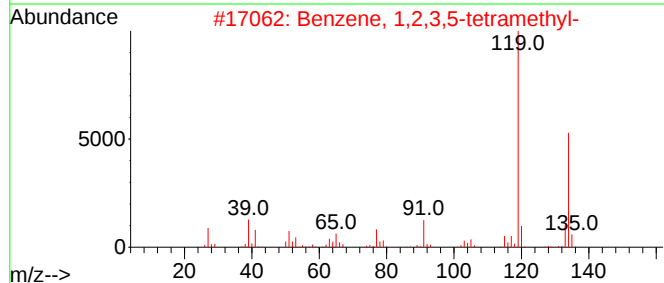
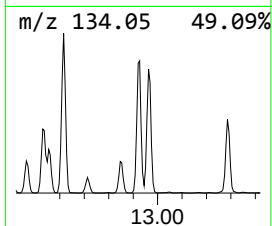
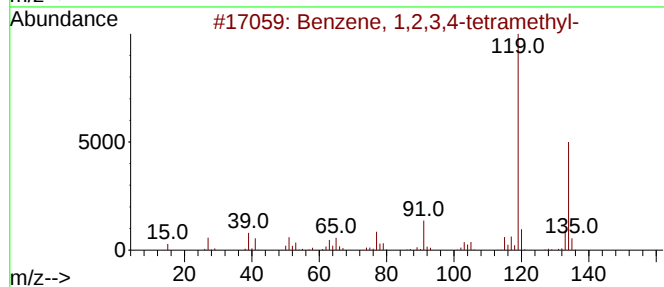
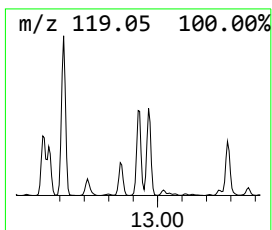
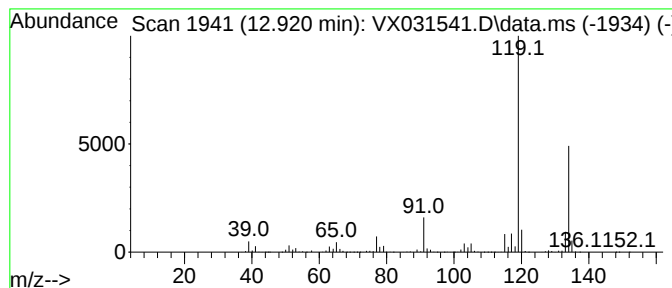
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	136.03 ug/l	2355480	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

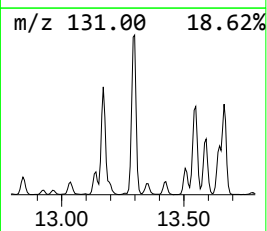
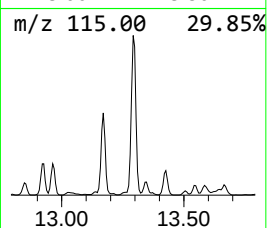
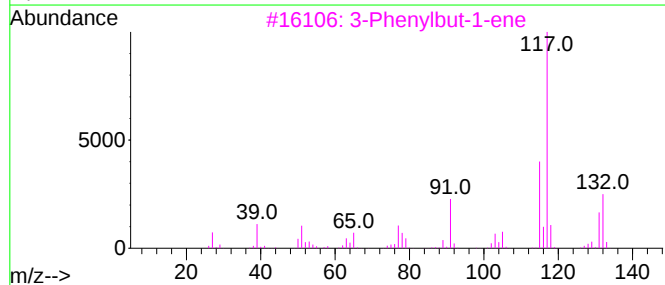
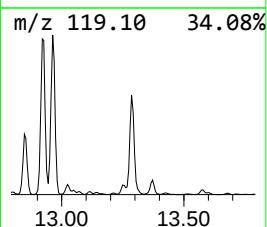
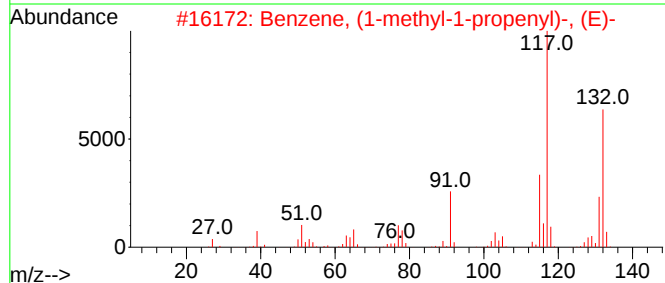
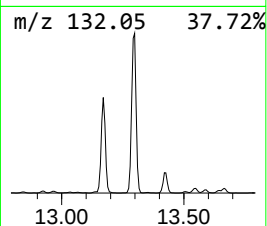
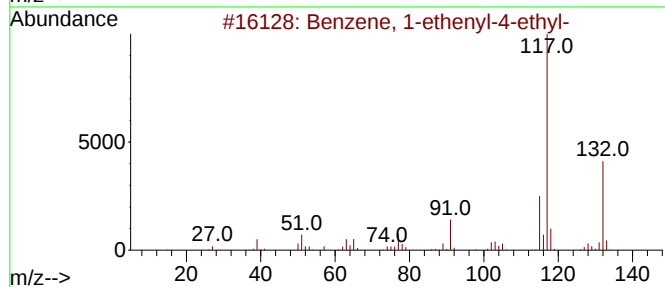
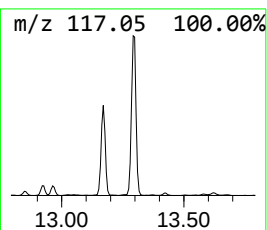
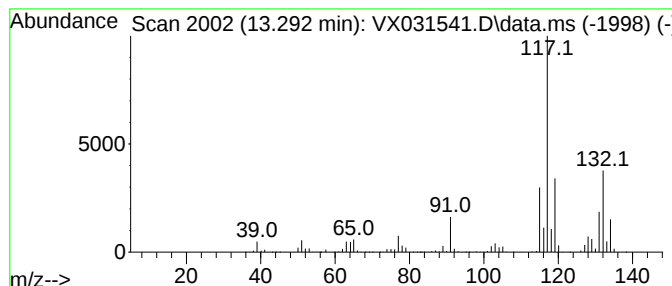
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	274.14 ug/l	4747080	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, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

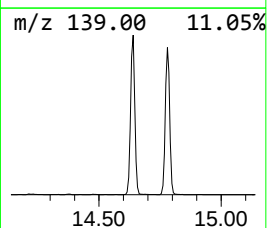
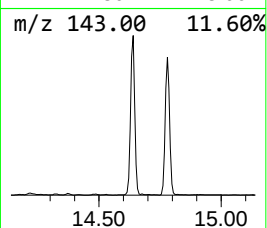
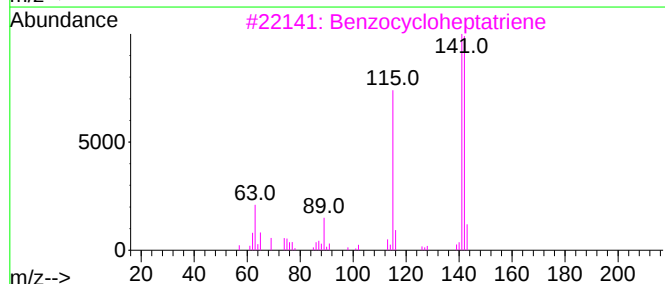
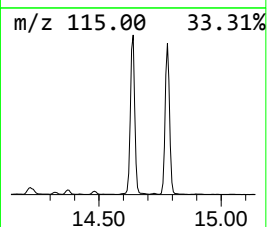
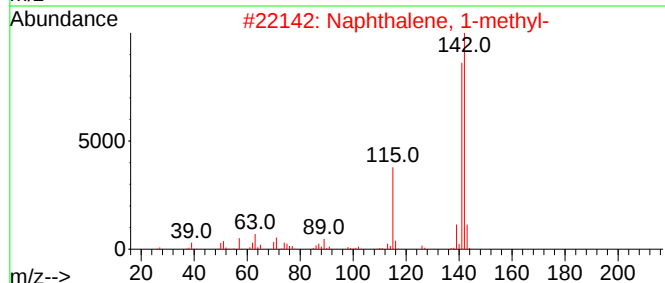
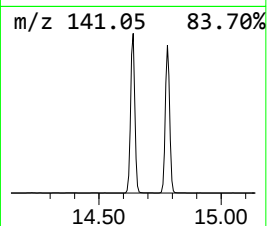
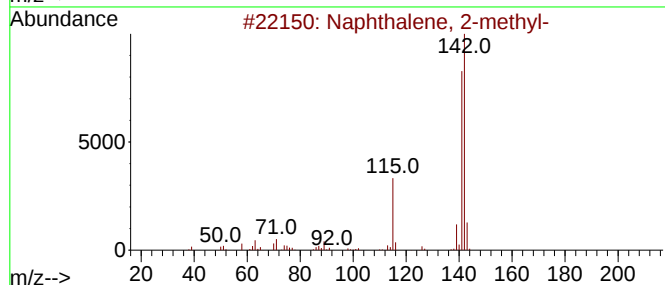
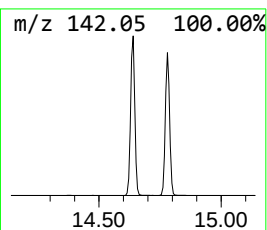
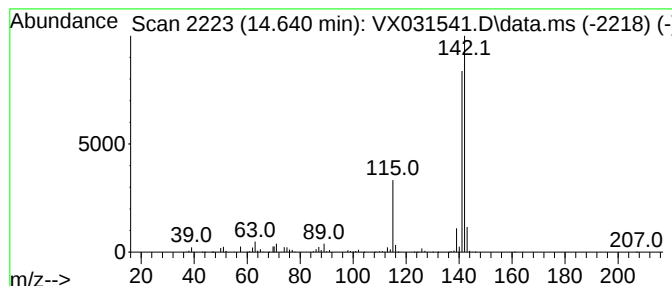
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	145.16 ug/l	2513600	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			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

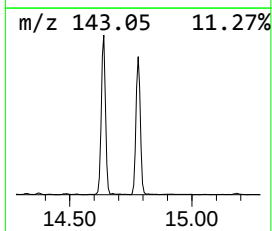
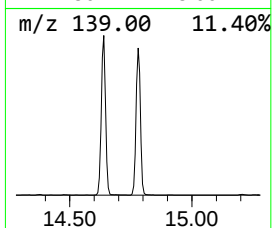
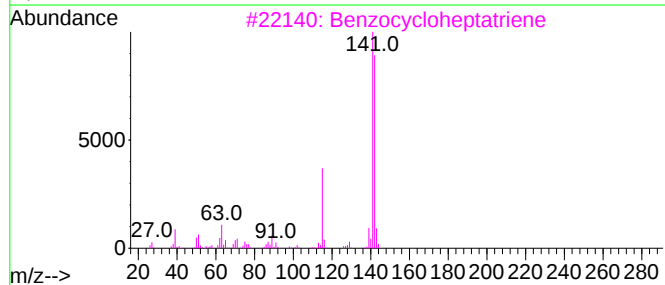
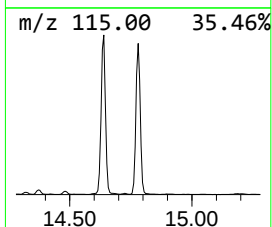
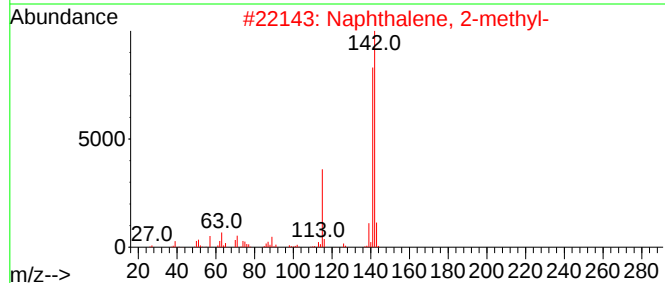
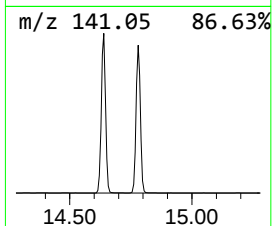
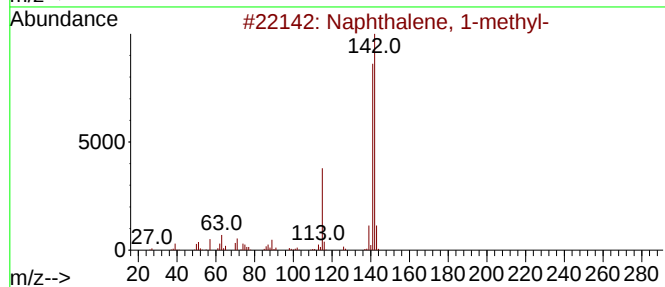
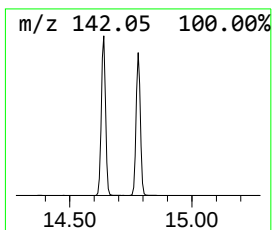
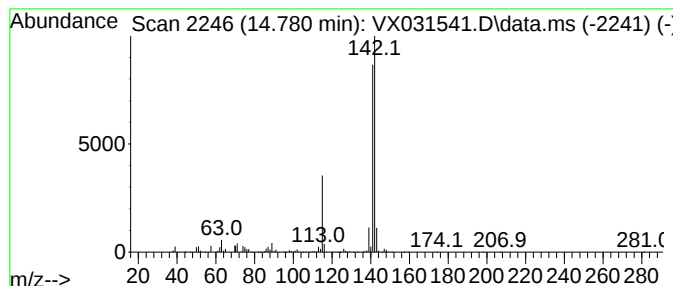
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	143.70 ug/l	2488430	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031541.D
Acq On : 22 Sep 2022 01:11
Operator : JC/MD
Sample : N4614-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-43

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.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
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Benzene, 1-ethy...	11.616	475.3	ug/l	8230950	4	12.024	865824	50.0
Benzene, 1,2,3-...	12.091	640.1	ug/l	11083700	4	12.024	865824	50.0
Indane	12.244	718.6	ug/l	12444000	4	12.024	865824	50.0
Benzene, 1-ethy...	12.616	212.0	ug/l	3670330	4	12.024	865824	50.0
Benzene, 1-ethe...	12.701	134.3	ug/l	2326530	4	12.024	865824	50.0
Benzene, 1,2,3,...	12.920	136.0	ug/l	2355480	4	12.024	865824	50.0
Benzene, 1-ethe...	13.292	274.1	ug/l	4747080	4	12.024	865824	50.0
Naphthalene, 2-...	14.640	145.2	ug/l	2513600	4	12.024	865824	50.0
Naphthalene, 1-...	14.780	143.7	ug/l	2488430	4	12.024	865824	50.0