

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
 Data File : VX023003.D
 Acq On : 29 Jun 2021 15:08
 Operator : JC/MD
 Sample : M2875-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-14

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\82X062321W.M
 Title : SW846 8260

Signal : TIC: VX023003.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	27	36	50	rBV	136497	840200	2.28%	0.713%
2	5.513	715	726	739	rBV2	146373	604988	1.65%	0.514%
3	5.836	766	779	792	rBV	269009	824428	2.24%	0.700%
4	6.617	895	907	923	rBV	179685	432154	1.18%	0.367%
5	8.653	1235	1241	1249	rVB	244157	374553	1.02%	0.318%
6	10.963	1615	1620	1625	rBV	975672	1246188	3.39%	1.058%
7	11.305	1671	1676	1682	rBV	925128	1145276	3.11%	0.972%
8	11.402	1682	1692	1696	rVV	647142	796891	2.17%	0.677%
9	11.451	1696	1700	1708	rVB	414605	536121	1.46%	0.455%
10	11.616	1719	1727	1737	rBV	1922852	2380203	6.47%	2.021%
11	11.756	1740	1750	1755	rBV	6742707	8660766	23.55%	7.353%
12	12.018	1789	1793	1799	rVB2	428189	664588	1.81%	0.564%
13	12.091	1799	1805	1809	rBV	2507872	3022917	8.22%	2.567%
14	12.250	1824	1831	1838	rBV	5787057	7416329	20.17%	6.297%
15	12.323	1838	1843	1848	rBV2	809885	1177756	3.20%	1.000%
16	12.536	1873	1878	1879	rBV	438700	513950	1.40%	0.436%
17	12.555	1879	1881	1886	rVV	530032	655400	1.78%	0.556%
18	12.616	1886	1891	1894	rVV	1202827	1498571	4.08%	1.272%
19	12.701	1900	1905	1911	rBV	2472714	3331882	9.06%	2.829%
20	12.926	1934	1942	1945	rBV	464375	596851	1.62%	0.507%
21	12.969	1945	1949	1953	rVB2	932420	1213766	3.30%	1.031%
22	13.067	1960	1965	1968	rBV2	277387	393210	1.07%	0.334%
23	13.170	1977	1982	1987	rVB	1812231	2125657	5.78%	1.805%
24	13.292	1993	2002	2007	rBV2	1900204	2587038	7.04%	2.196%
25	13.347	2007	2011	2018	rVB	1753710	2102660	5.72%	1.785%
26	13.432	2019	2025	2030	rVB	6259401	8088216	22.00%	6.867%
27	13.506	2030	2037	2041	rBV	982686	1229074	3.34%	1.044%
28	13.548	2041	2044	2048	rVV	345794	455404	1.24%	0.387%
29	13.676	2061	2065	2070	rVB2	284300	416069	1.13%	0.353%
30	13.792	2074	2084	2089	rBV	21191011	36771216	100.00%	31.220%
31	14.225	2149	2155	2158	rBV2	530506	853603	2.32%	0.725%
32	14.323	2166	2171	2178	rVB	698263	1234201	3.36%	1.048%
33	14.524	2200	2204	2212	rVB	314823	488430	1.33%	0.415%
34	14.646	2218	2224	2233	rVB	8065921	10059430	27.36%	8.541%
35	14.786	2241	2247	2252	rBV	7237408	8846378	24.06%	7.511%
36	15.206	2308	2316	2321	rBV	333412	476840	1.30%	0.405%

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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

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Title : SW846 8260

37	15.377	2339	2344	2348	rBV	456643	647448	1.76%	0.550%
38	15.481	2354	2361	2366	rBV	363798	561214	1.53%	0.476%
39	15.615	2376	2383	2387	rBV	646272	1037231	2.82%	0.881%
40	15.834	2410	2419	2432	rVB2	172856	519976	1.41%	0.441%
41	16.328	2492	2500	2511	rBV	507804	954041	2.59%	0.810%

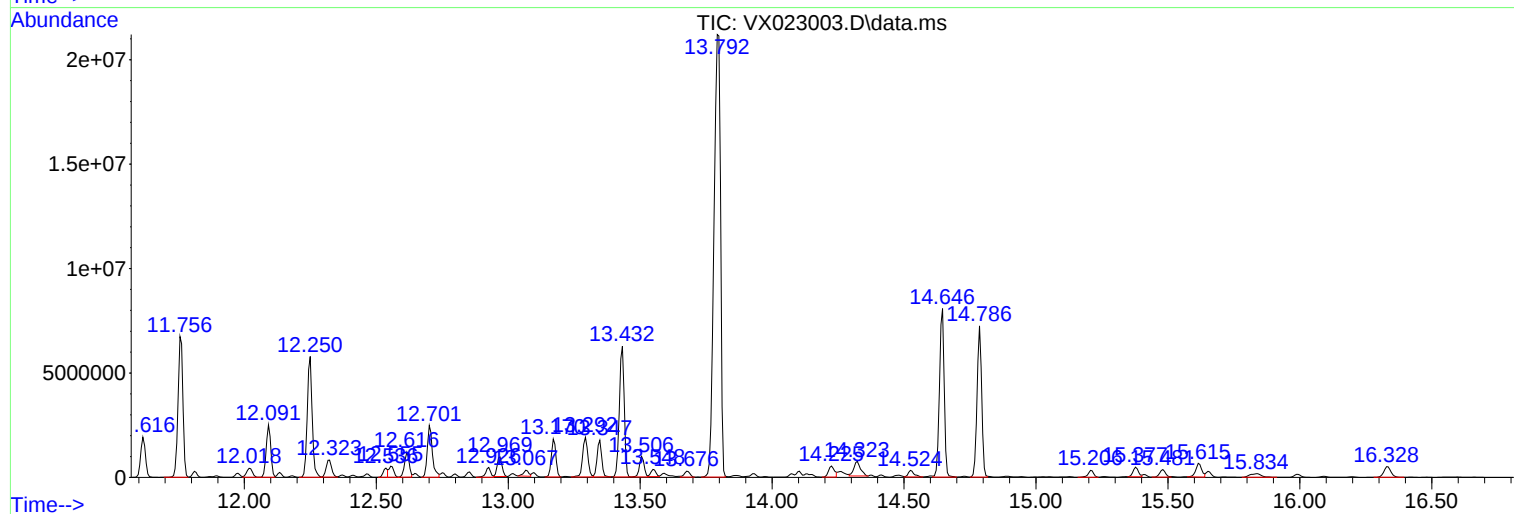
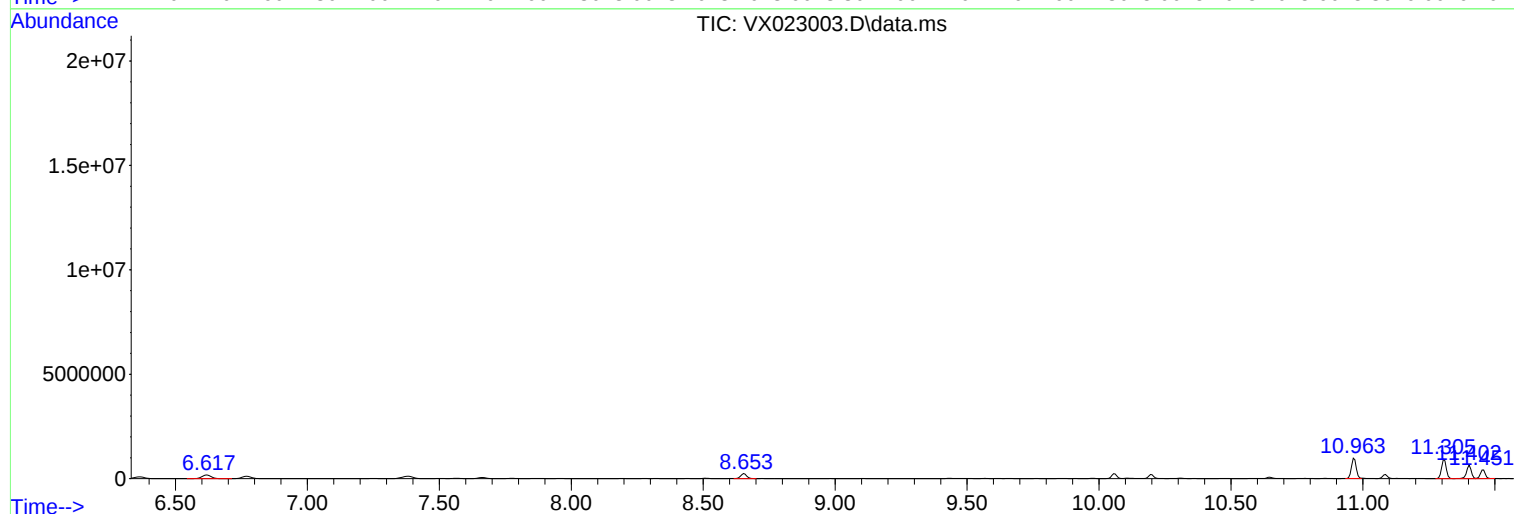
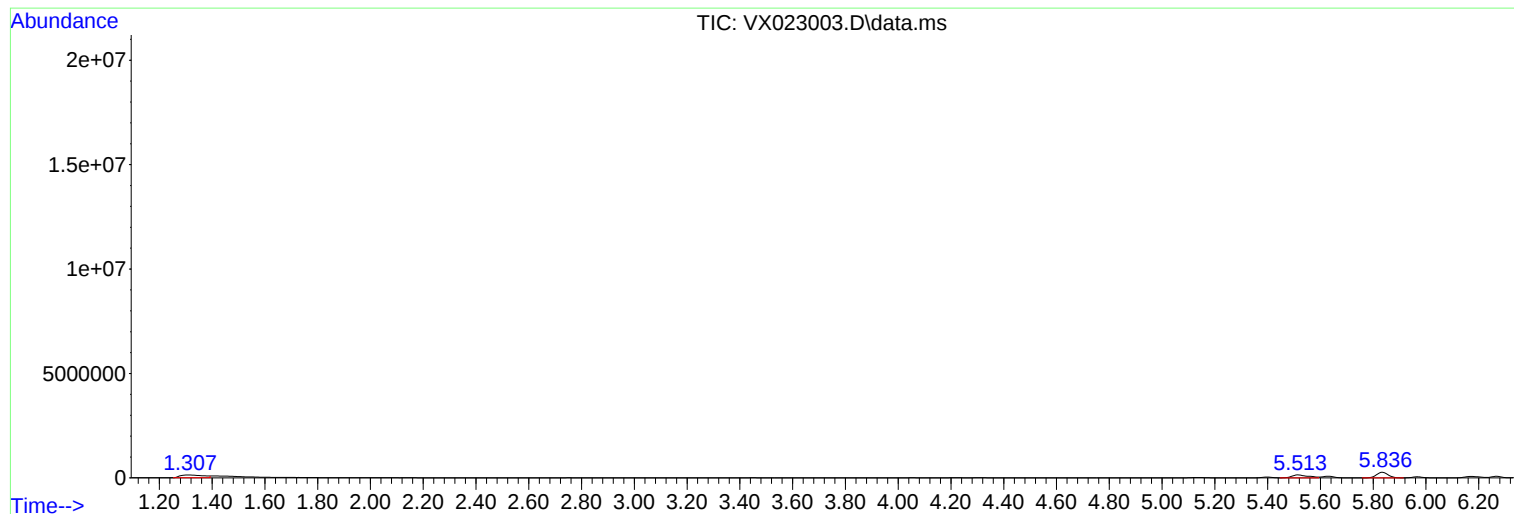
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
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Instrument :
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ClientSampleId :
MW-14

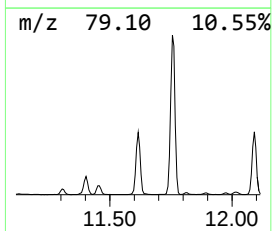
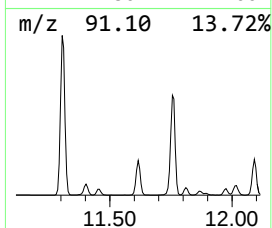
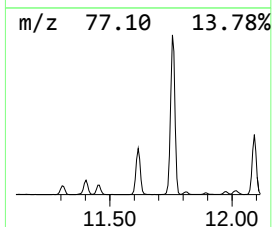
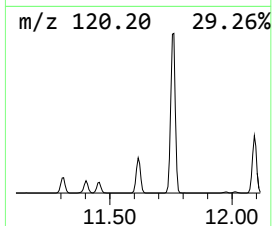
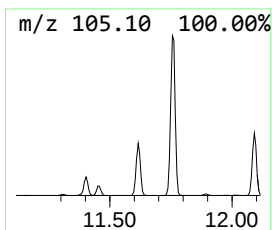
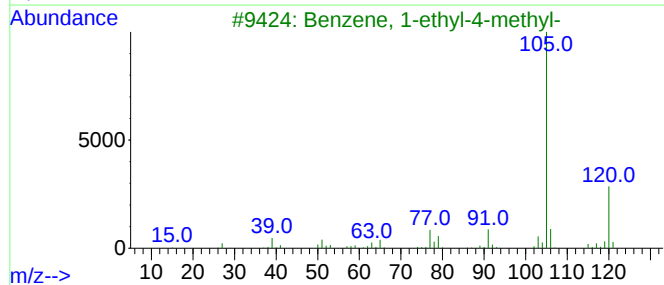
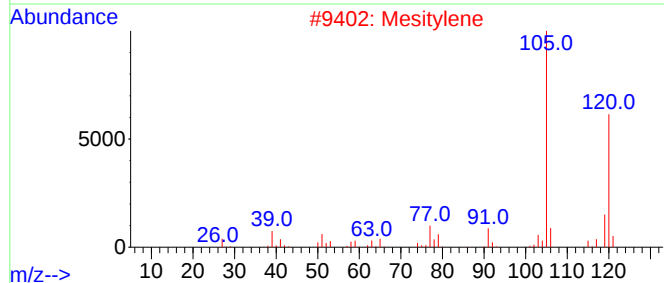
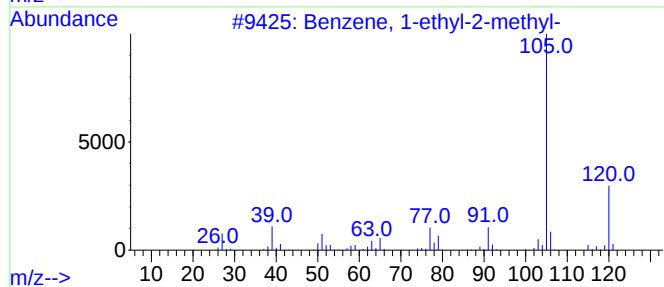
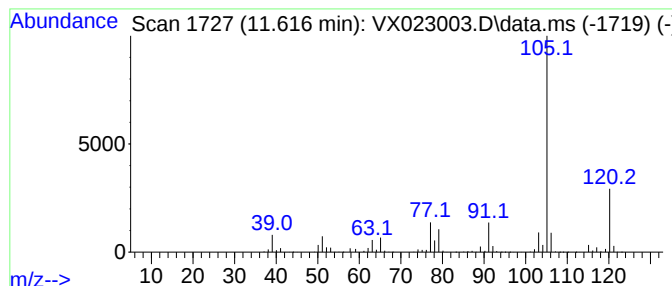
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	179.07 ug/l	2380200	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			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

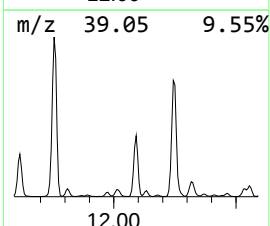
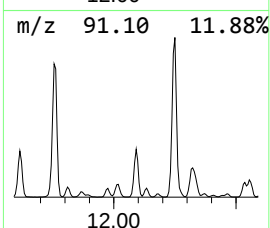
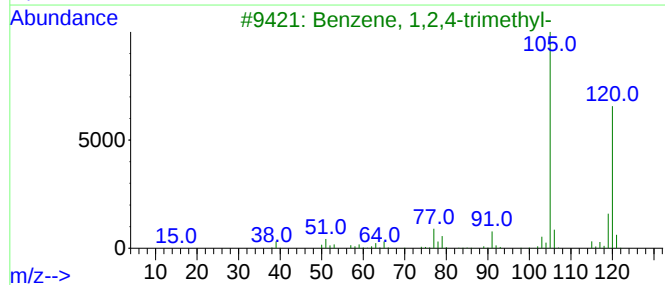
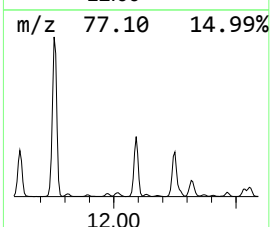
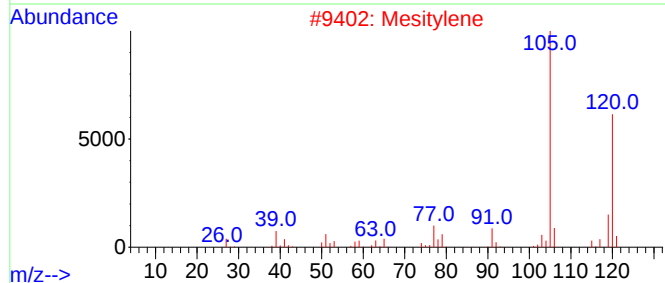
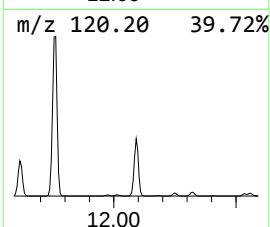
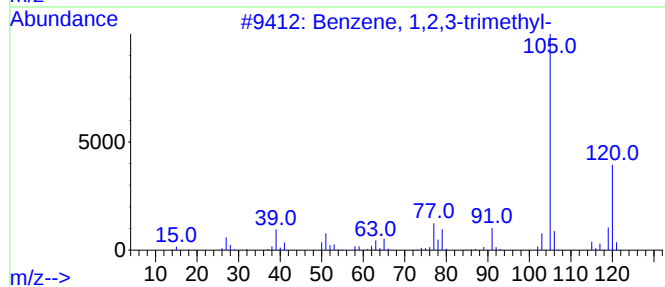
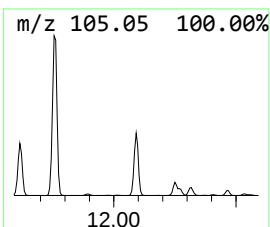
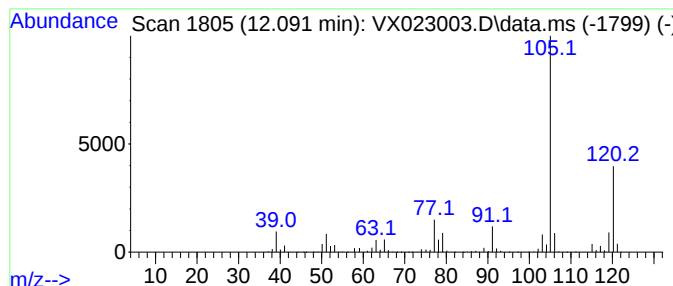
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	227.43 ug/l	3022920	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	94
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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

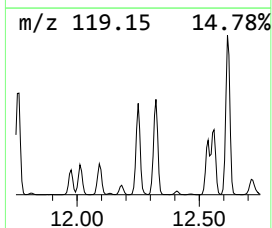
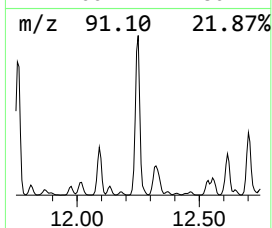
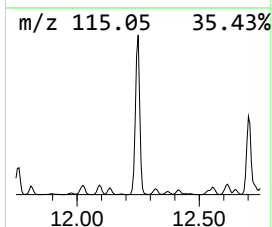
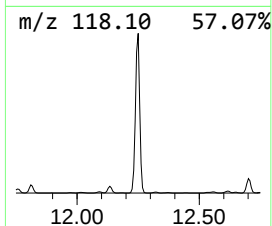
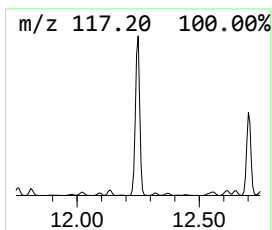
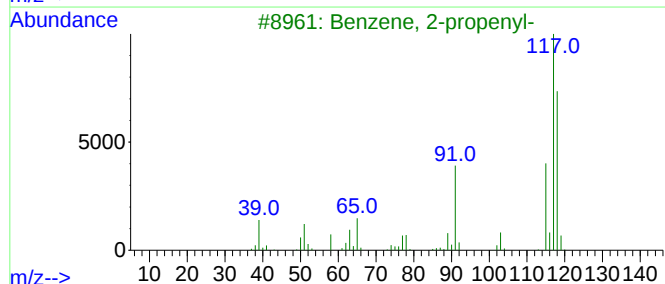
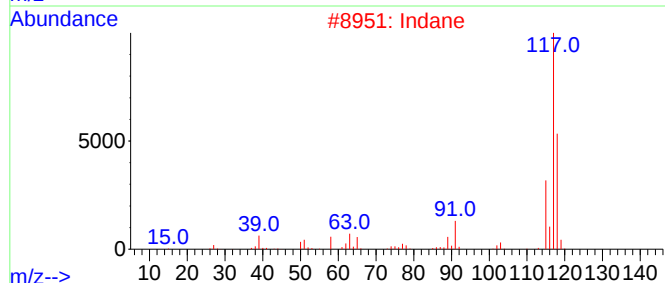
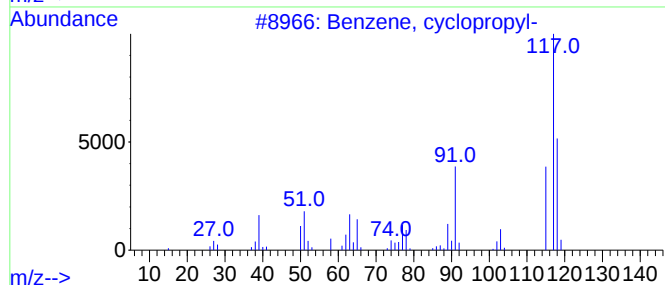
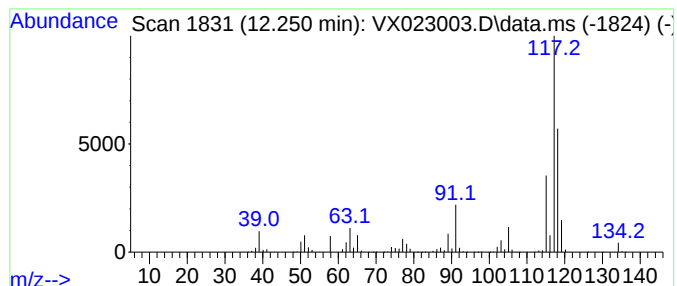
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	557.96 ug/l	7416330	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2			Indane	118	C9H10	000496-11-7	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4			Delta-cyclene	118	C9H10	007785-10-6	53
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	50



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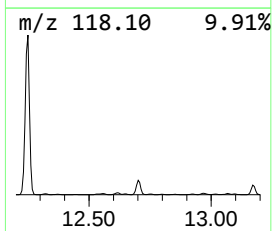
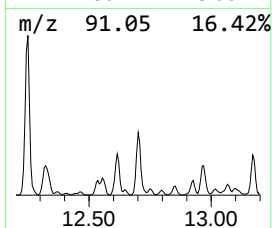
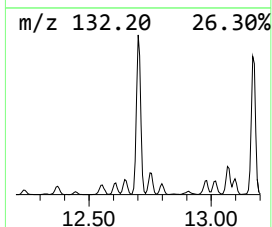
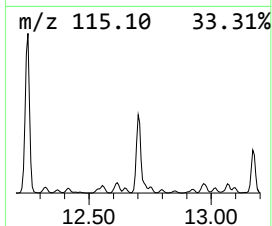
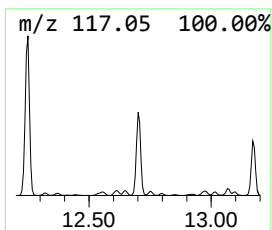
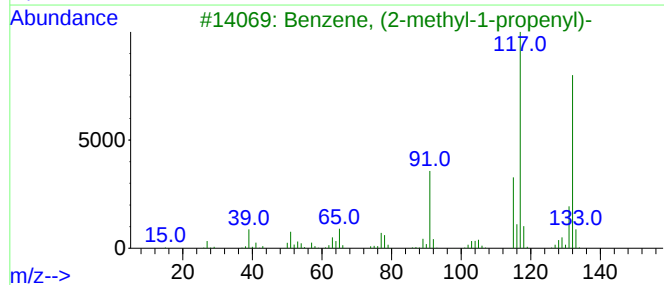
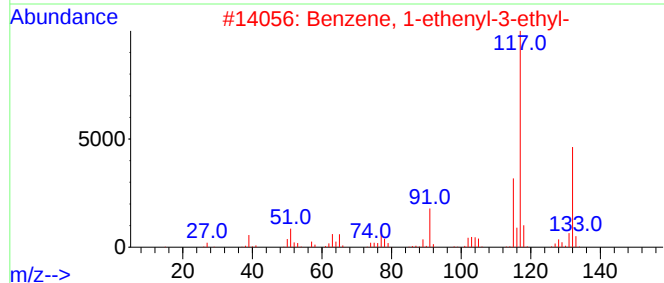
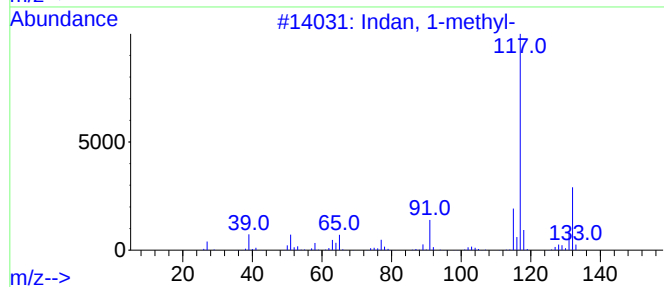
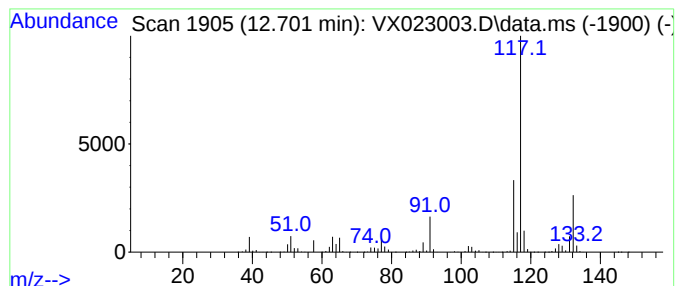
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	250.67 ug/l	3331880	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	91
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
4			Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83



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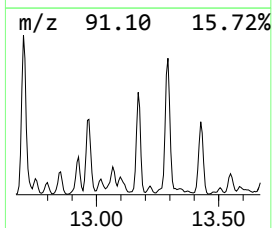
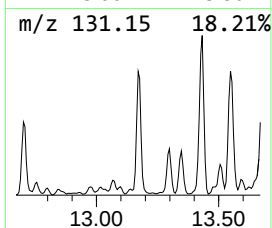
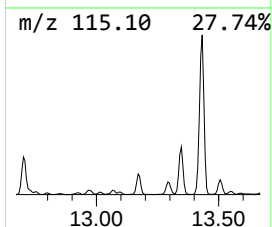
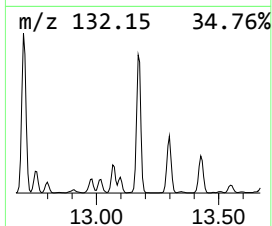
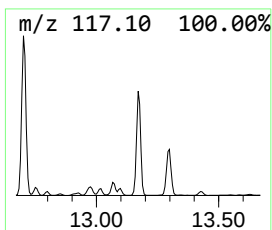
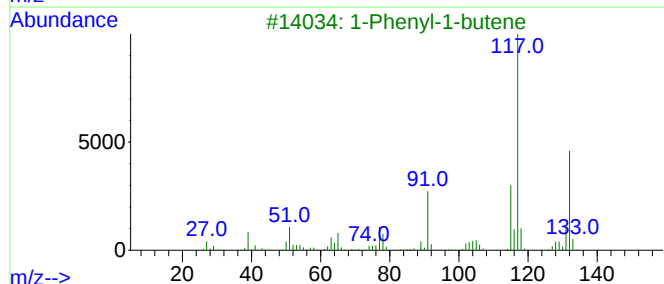
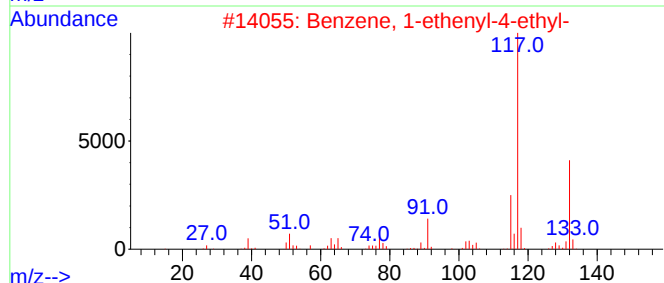
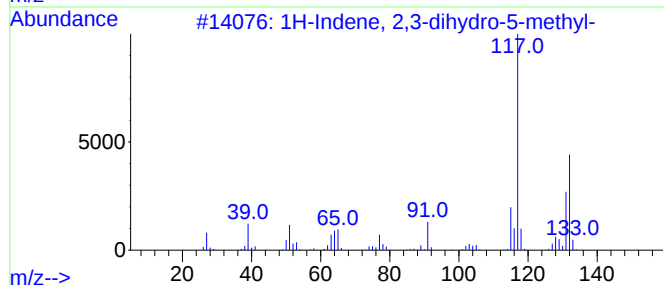
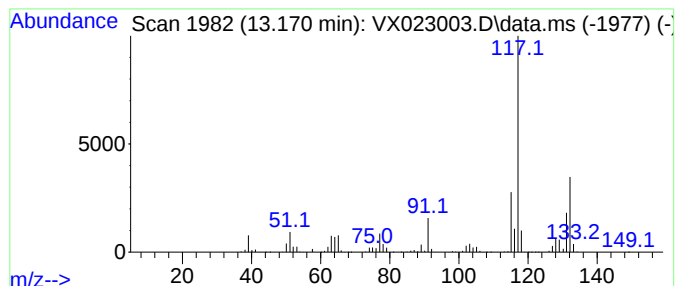
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	159.92 ug/l	2125660	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	94
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023003.D
Acq On : 29 Jun 2021 15:08
Operator : JC/MD
Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

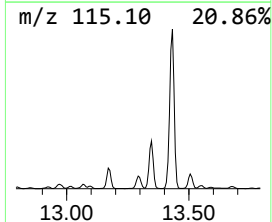
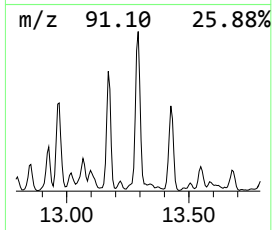
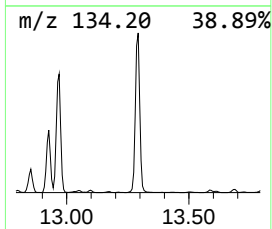
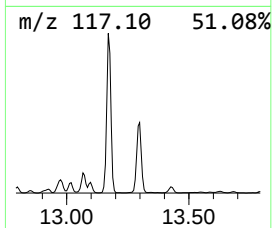
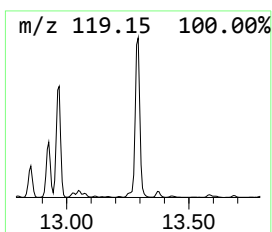
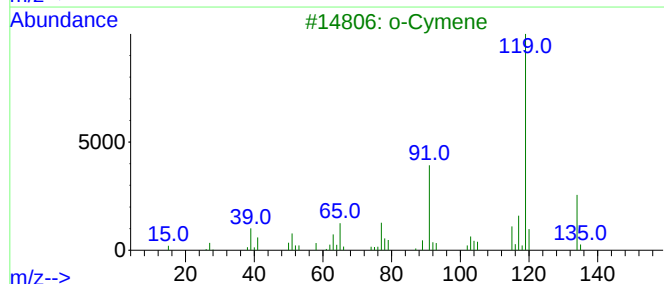
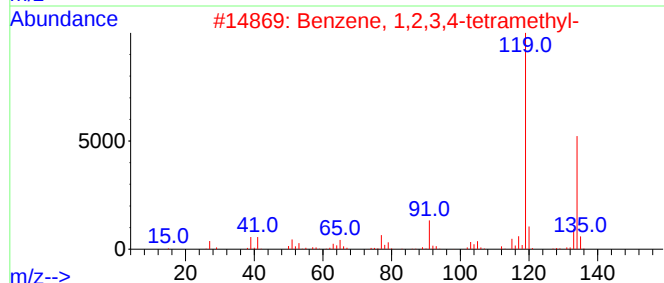
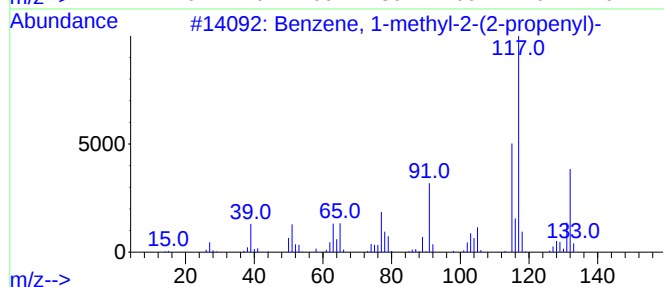
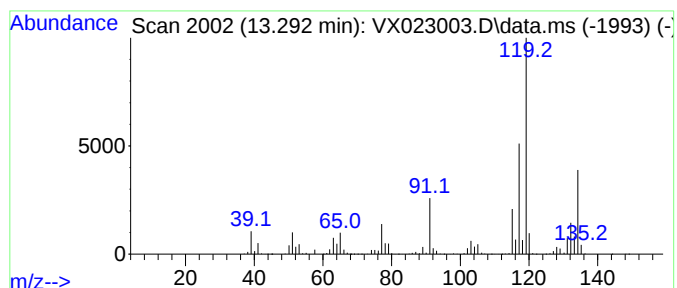
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-methyl-2-(2-prop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	194.63 ug/l	2587040	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
3			o-Cymene	134	C10H14	000527-84-4	70
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5			p-Cymene	134	C10H14	000099-87-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023003.D
Acq On : 29 Jun 2021 15:08
Operator : JC/MD
Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

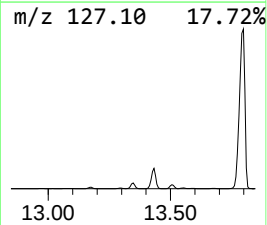
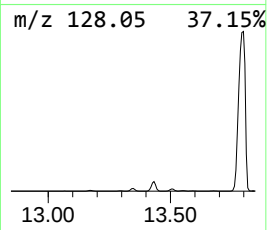
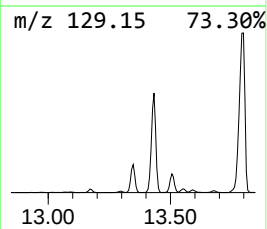
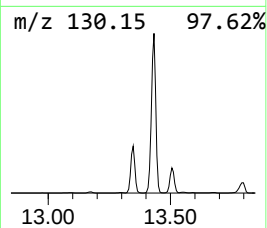
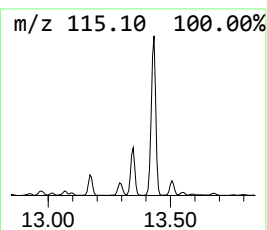
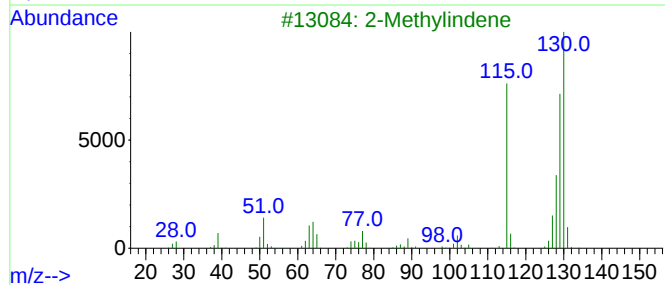
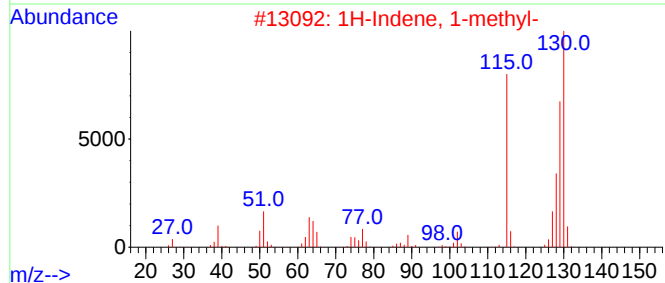
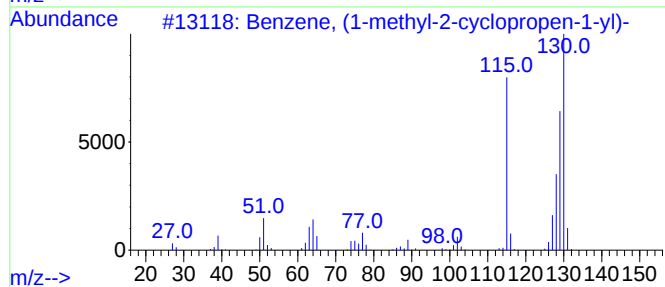
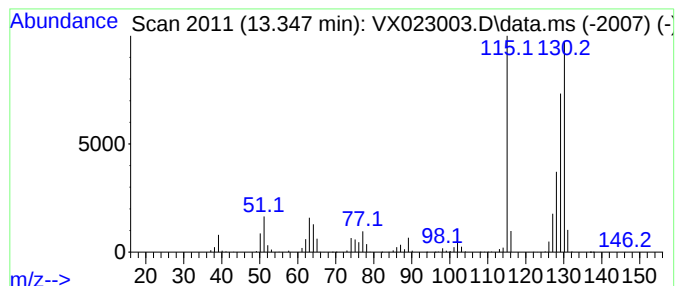
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.347	158.19 ug/l	2102660	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
2			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
3			2-Methylindene	130	C10H10	002177-47-1	96
4			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	96
5			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023003.D
Acq On : 29 Jun 2021 15:08
Operator : JC/MD
Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

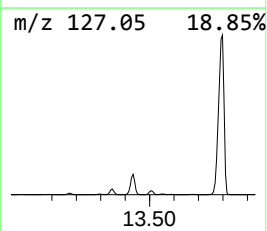
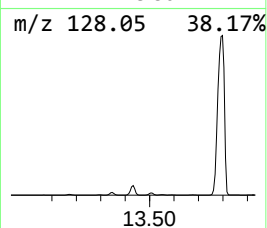
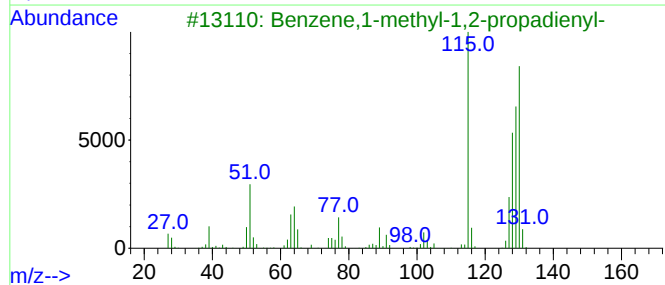
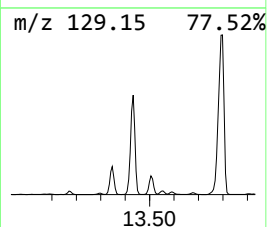
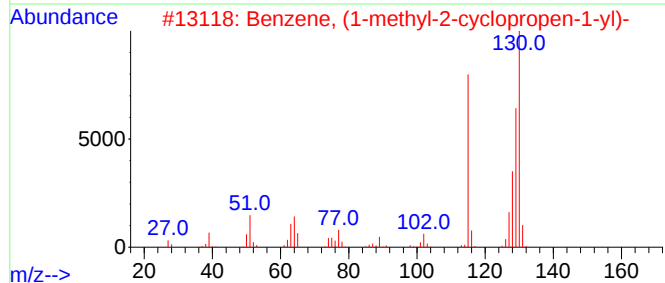
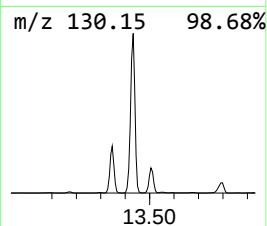
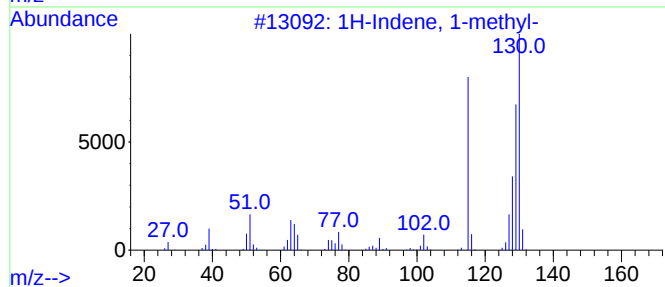
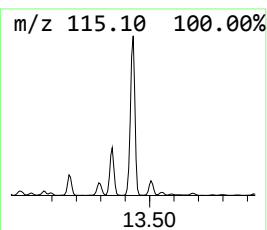
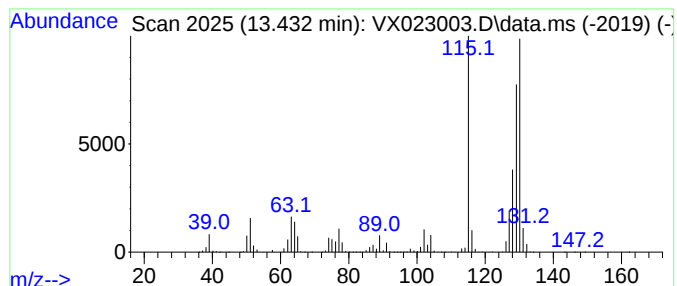
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.432	608.51 ug/l	8088220	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
3			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	96
4			2-Methylindene	130	C10H10	002177-47-1	96
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023003.D
Acq On : 29 Jun 2021 15:08
Operator : JC/MD
Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

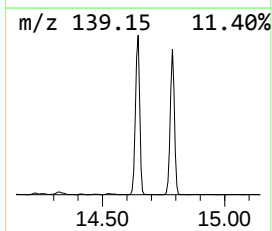
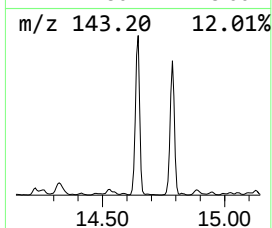
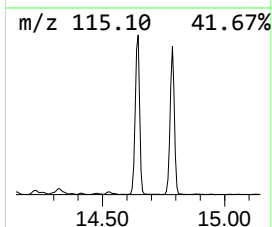
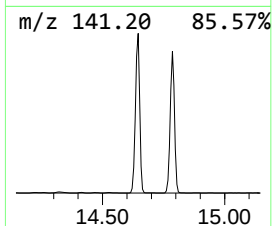
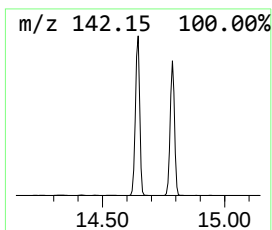
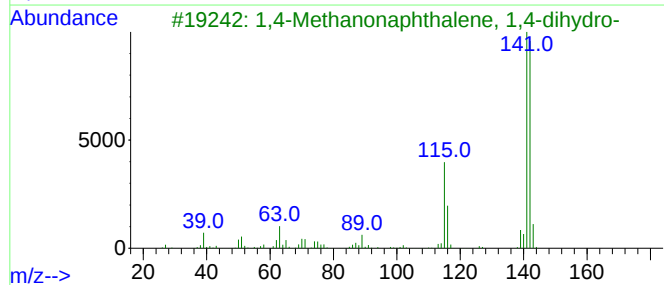
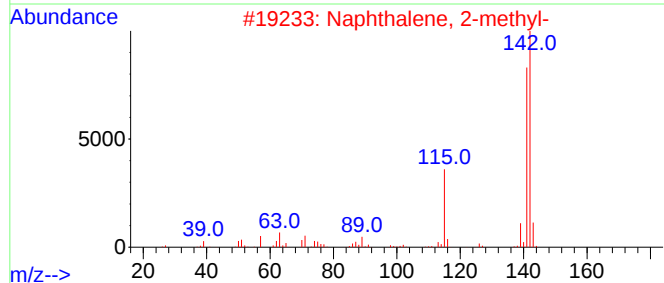
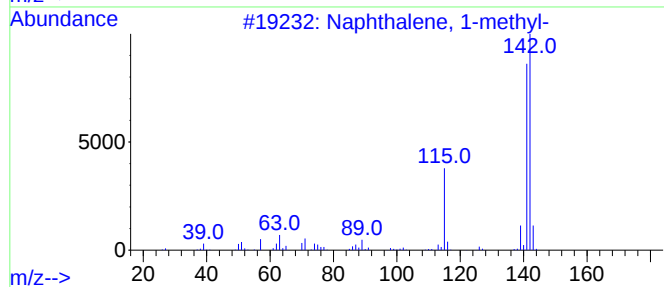
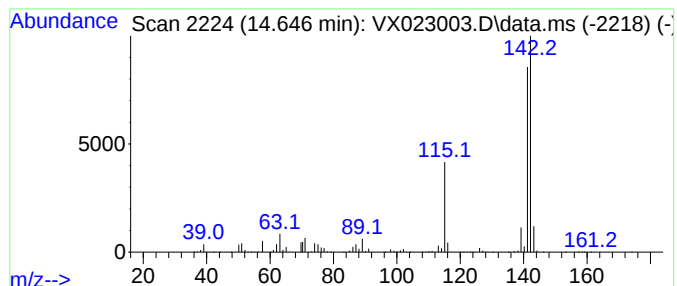
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.646	756.82 ug/l	10059400	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			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023003.D
Acq On : 29 Jun 2021 15:08
Operator : JC/MD
Sample : M2875-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

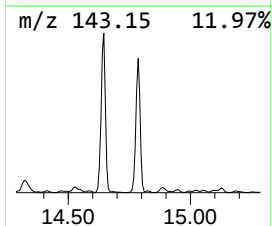
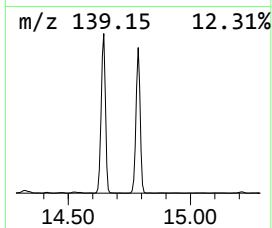
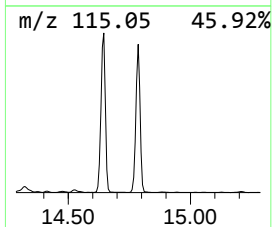
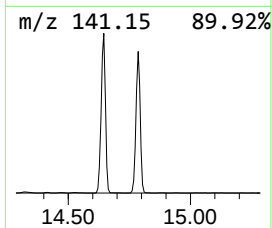
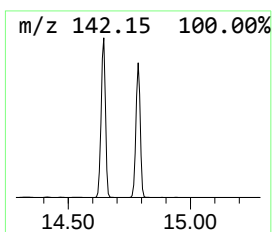
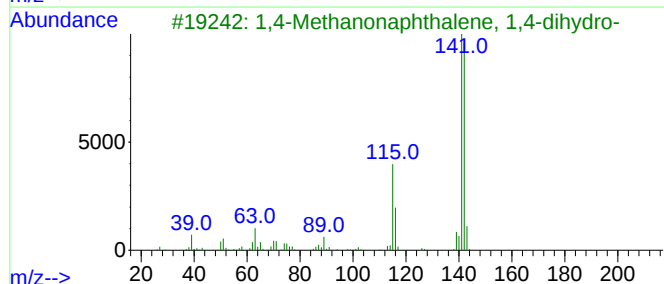
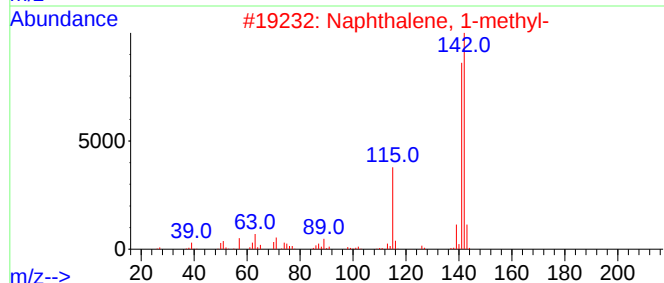
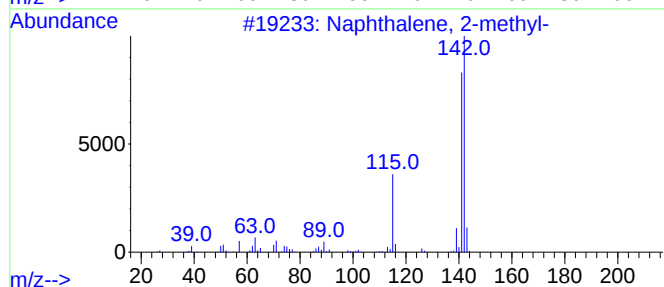
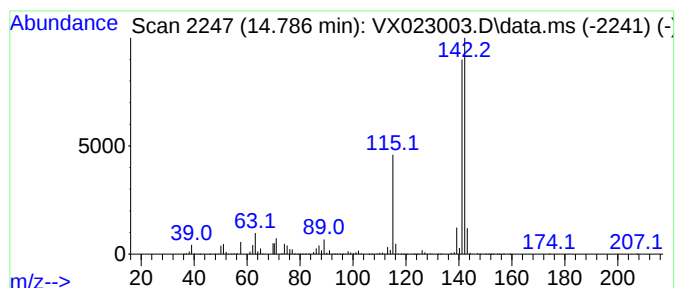
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.786	665.55 ug/l	8846380	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
 Data File : VX023003.D
 Acq On : 29 Jun 2021 15:08
 Operator : JC/MD
 Sample : M2875-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-14

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	11.616	179.1	ug/l	2380200	4	12.024	664588	50.0
Benzene, 1,2,3-...	12.091	227.4	ug/l	3022920	4	12.024	664588	50.0
Benzene, cyclop...	12.250	558.0	ug/l	7416330	4	12.024	664588	50.0
Indan, 1-methyl-	12.701	250.7	ug/l	3331880	4	12.024	664588	50.0
1H-Indene, 2,3-...	13.170	159.9	ug/l	2125660	4	12.024	664588	50.0
Benzene, 1-meth...	13.292	194.6	ug/l	2587040	4	12.024	664588	50.0
Benzene, (1-met...	13.347	158.2	ug/l	2102660	4	12.024	664588	50.0
1H-Indene, 1-me...	13.432	608.5	ug/l	8088220	4	12.024	664588	50.0
Naphthalene, 1-...	14.646	756.8	ug/l	10059400	4	12.024	664588	50.0
1,4-Methanonaph...	14.786	665.5	ug/l	8846380	4	12.024	664588	50.0