

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
 Data File : VY016168.D
 Acq On : 01 Nov 2023 15:03
 Operator : SY/MD
 Sample : 05102-05
 Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 67S-SB07-1418

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_Y\methods\82Y103123S.M

Title : SW846 8260

Signal : TIC: VY016168.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.167	1030	1041	1050	rBV	644889	1481768	1.32%	0.252%
2	9.831	1308	1314	1327	rVB	1152704	2180283	1.95%	0.371%
3	10.264	1376	1385	1398	rVV3	32901126	111981224	100.00%	19.050%
4	10.380	1398	1404	1412	rVB	763300	1287802	1.15%	0.219%
5	11.611	1597	1606	1615	rBV2	34346499	84252272	75.24%	14.333%
6	11.721	1615	1624	1634	rBV2	36889466	98731278	88.17%	16.796%
7	12.044	1665	1677	1688	rBV	26870597	53179382	47.49%	9.047%
8	12.343	1720	1726	1730	rBV	906795	1486010	1.33%	0.253%
9	12.684	1772	1782	1786	rBV	724722	1304934	1.17%	0.222%
10	12.757	1786	1794	1797	rBV2	38177374	91822218	82.00%	15.620%
11	12.788	1797	1799	1811	rVB	30682922	47873944	42.75%	8.144%
12	12.983	1823	1831	1838	rBV	1228653	2191177	1.96%	0.373%
13	13.129	1849	1855	1860	rBV	5094309	7525676	6.72%	1.280%
14	13.178	1860	1863	1867	rVB	1623915	2152933	1.92%	0.366%
15	13.233	1867	1872	1873	rBV	2153756	2909774	2.60%	0.495%
16	13.257	1873	1876	1883	rVB	5759938	8305514	7.42%	1.413%
17	13.465	1903	1910	1913	rBV	1947760	3199131	2.86%	0.544%
18	13.507	1913	1917	1924	rVB	7746349	10605970	9.47%	1.804%
19	13.599	1924	1932	1934	rBV4	741947	1298031	1.16%	0.221%
20	13.635	1934	1938	1941	rVV2	2333672	3745312	3.34%	0.637%
21	13.672	1941	1944	1953	rVV2	2508404	4717665	4.21%	0.803%
22	13.757	1953	1958	1963	rVV3	834426	1377836	1.23%	0.234%
23	13.830	1963	1970	1974	rVV2	784284	1359018	1.21%	0.231%
24	13.891	1974	1980	1983	rVV3	1797146	3575439	3.19%	0.608%
25	13.922	1983	1985	1990	rVV	1911773	2843284	2.54%	0.484%
26	13.983	1990	1995	1998	rVV	3334722	5015113	4.48%	0.853%
27	14.019	1998	2001	2007	rVV2	791487	1279314	1.14%	0.218%
28	14.086	2007	2012	2017	rVB2	1088824	1704339	1.52%	0.290%
29	14.221	2029	2034	2039	rBV	860483	1327663	1.19%	0.226%
30	14.306	2039	2048	2051	rVV	2741411	4751248	4.24%	0.808%
31	14.343	2051	2054	2058	rVV	3954276	5686846	5.08%	0.967%
32	14.391	2058	2062	2065	rVV	1202627	1883351	1.68%	0.320%
33	14.574	2088	2092	2096	rVV	1280800	2009497	1.79%	0.342%
34	14.617	2096	2099	2104	rVB	1014433	1477742	1.32%	0.251%
35	14.684	2104	2110	2116	rBV2	2603815	5806064	5.18%	0.988%
36	14.952	2151	2154	2166	rVB2	698672	1519971	1.36%	0.259%

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Instrument :
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ClientSampleId :
67S-SB07-1418

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_Y\methods\82Y103123S.M
Title : SW846 8260

37	15.245	2197	2202	2208	rVV	2301237	3983603	3.56%	0.678%
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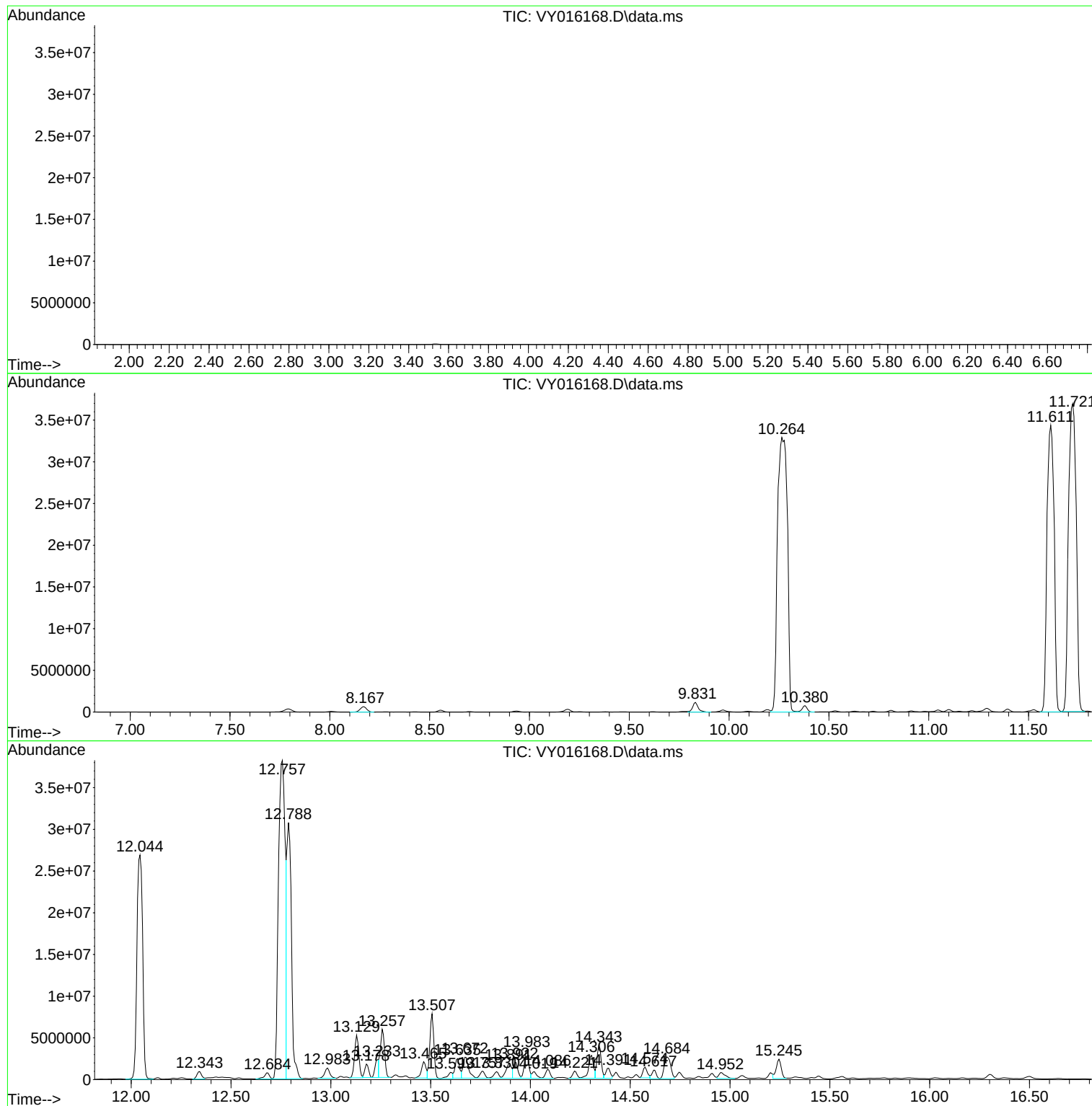
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.M
Quant Title : SW846 8260

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



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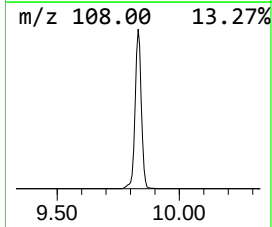
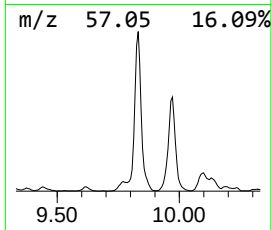
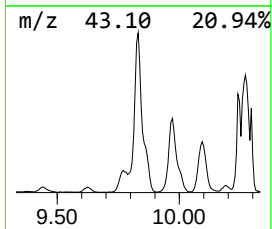
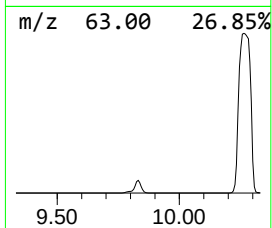
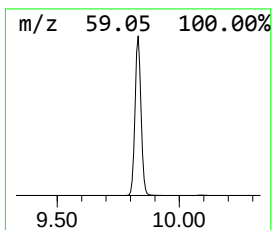
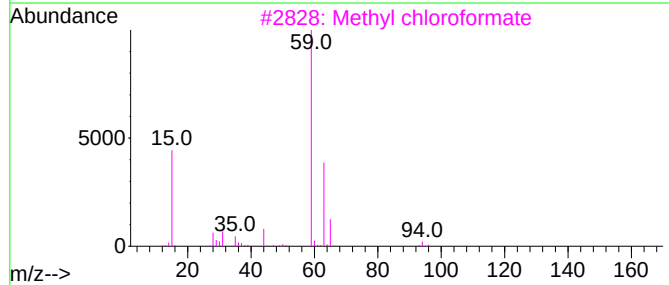
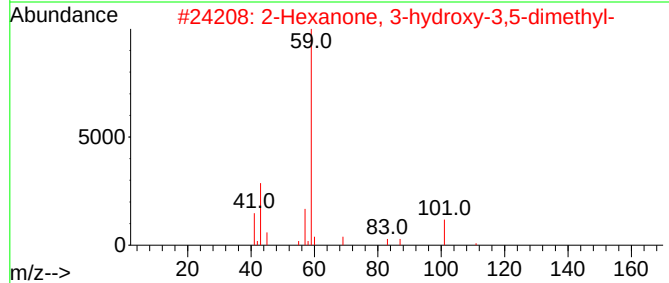
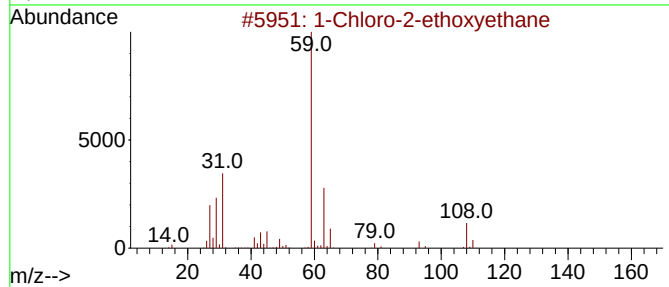
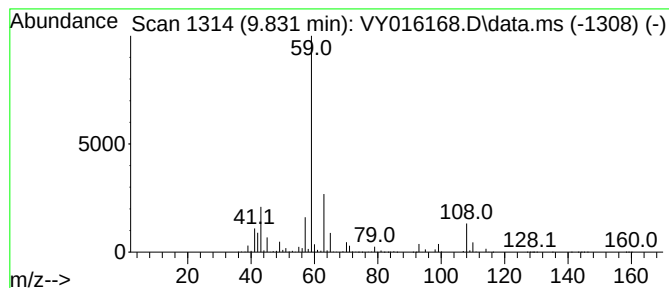
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Chloro-2-ethoxyethane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.831	73.57 ug/l	2180280	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Chloro-2-ethoxyethane	108	C4H9ClO	000628-34-2	87
2			2-Hexanone, 3-hydroxy-3,5-dimethyl-	144	C8H16O2	006321-14-8	50
3			Methyl chloroformate	94	C2H3ClO2	000079-22-1	36
4			2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	33
5			Propanoic acid, 2-hydroxy-2-meth...	132	C6H12O3	000080-55-7	25



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Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

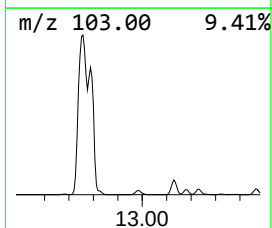
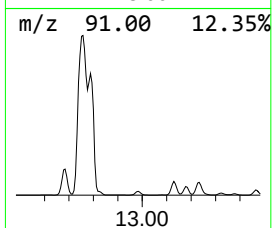
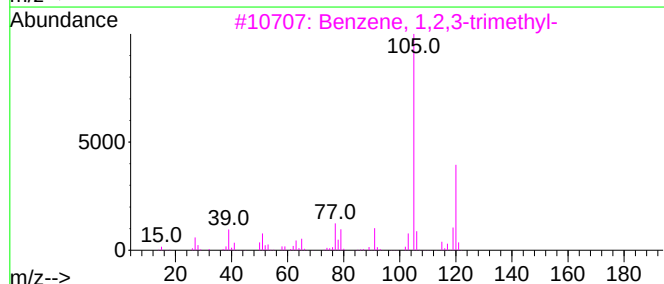
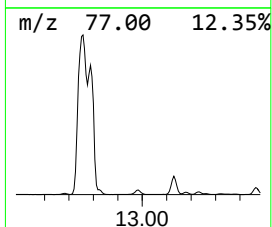
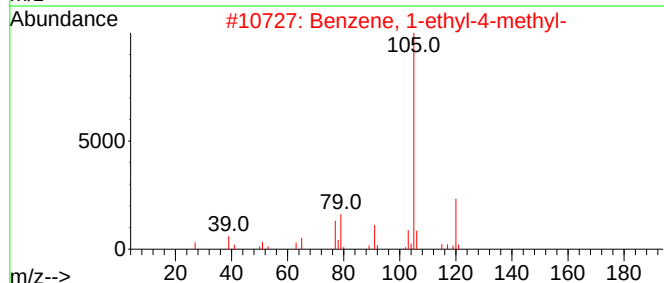
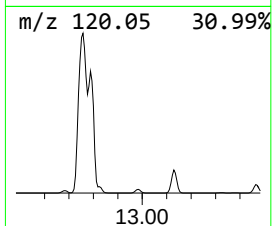
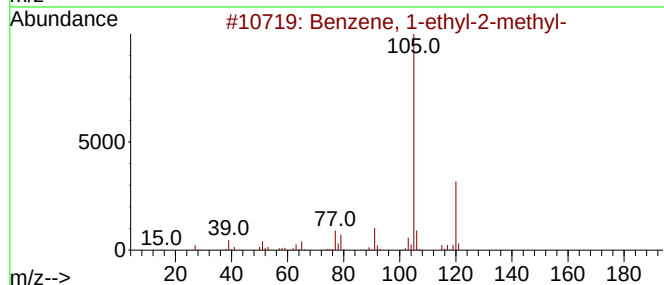
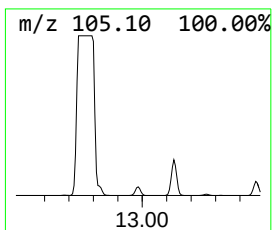
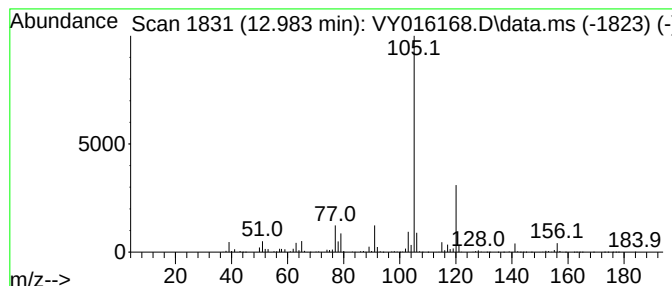
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.983	34.25 ug/l	2191180	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

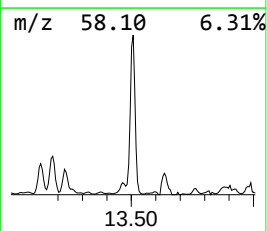
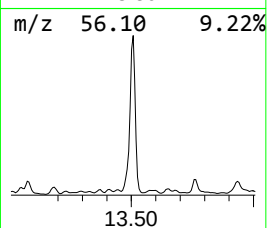
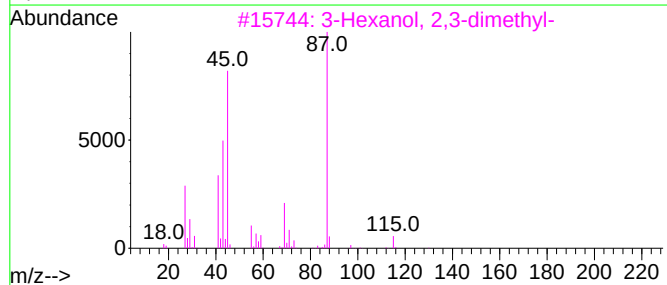
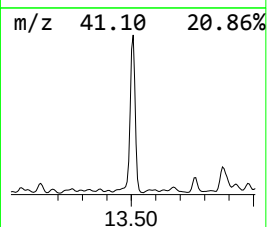
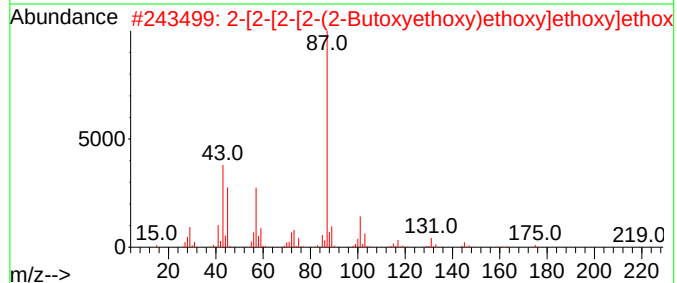
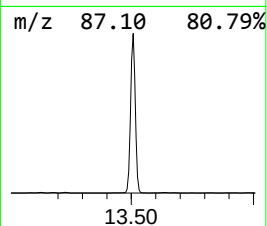
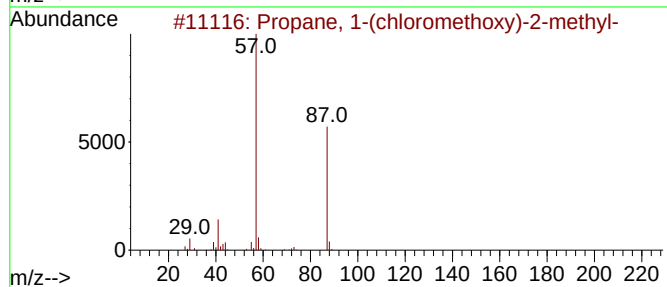
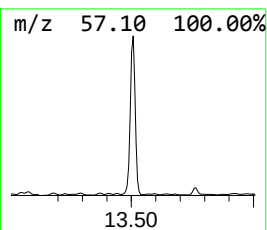
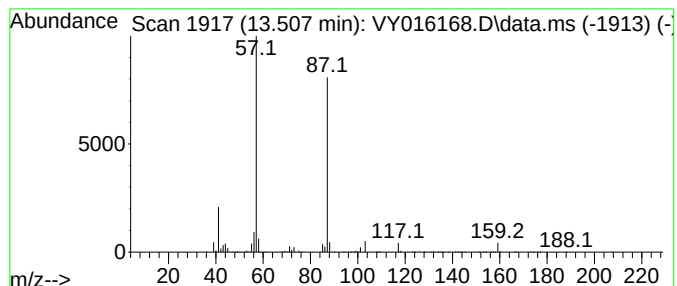
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.M
Quant Title : SW846 8260

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

Peak Number 3 Propane, 1-(chloromethoxy)-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.507	165.76 ug/l	10606000	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1-(chloromethoxy)-2-met...	122	C5H11ClO	034180-11-5	59
2			2-[2-[2-(2-Butoxyethoxy)ethox...	336	C16H32O7	1000351-94-0	56
3			3-Hexanol, 2,3-dimethyl-	130	C8H18O	004166-46-5	40
4			n-Butyl ether	130	C8H18O	000142-96-1	38
5			2-[2-[2-(2-Butoxyethoxy)ethoxy]e...	292	C14H28O6	1000351-93-9	38



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ClientSampleId :
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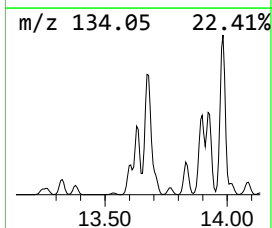
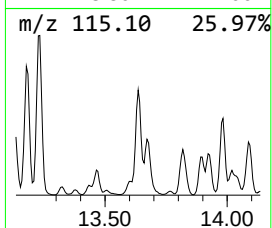
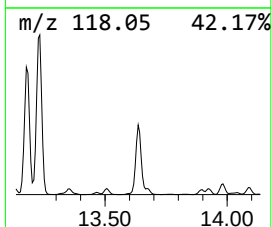
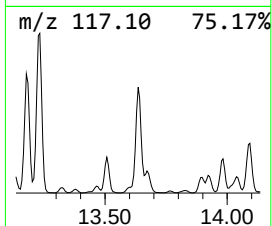
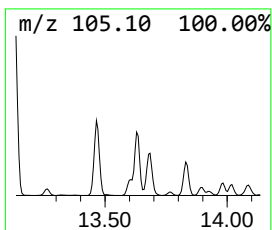
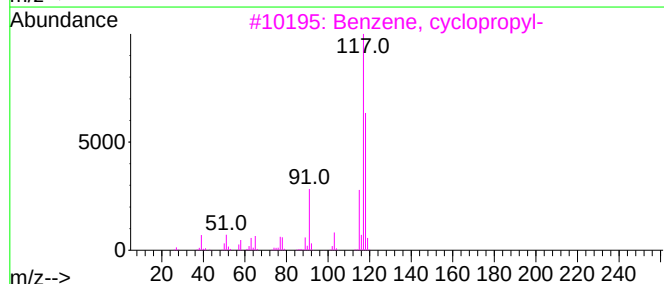
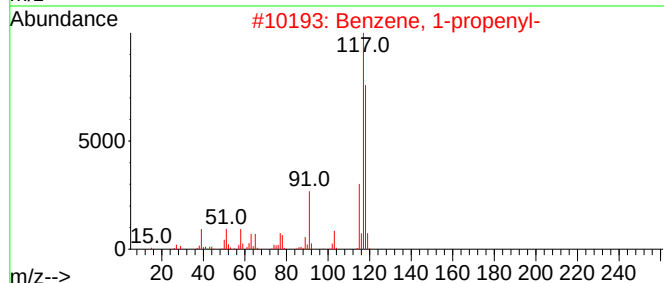
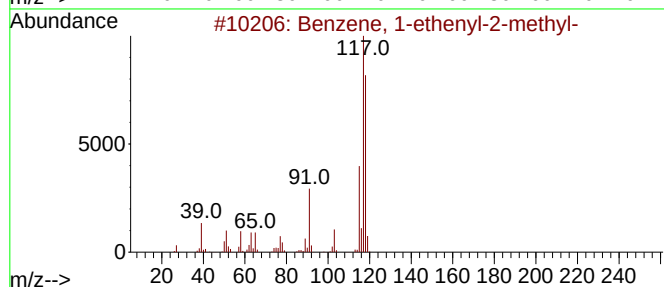
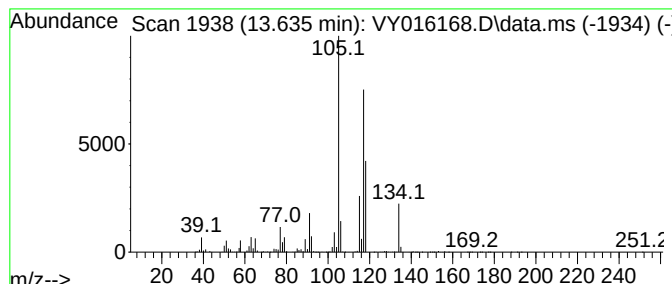
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.635	58.54 ug/l	3745310	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	58
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	58



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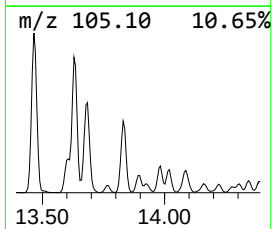
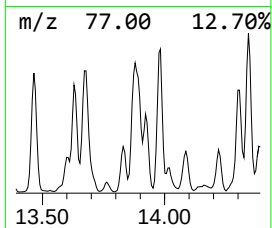
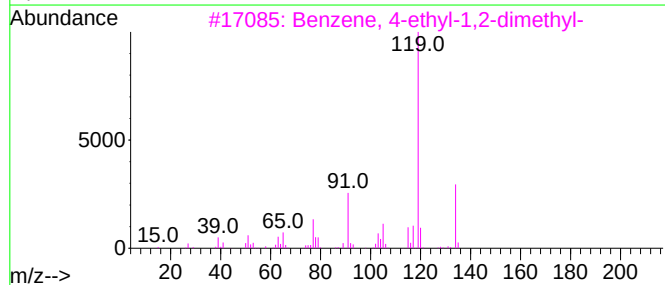
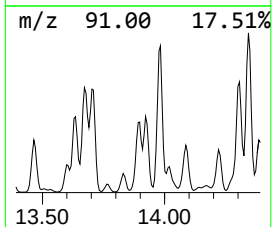
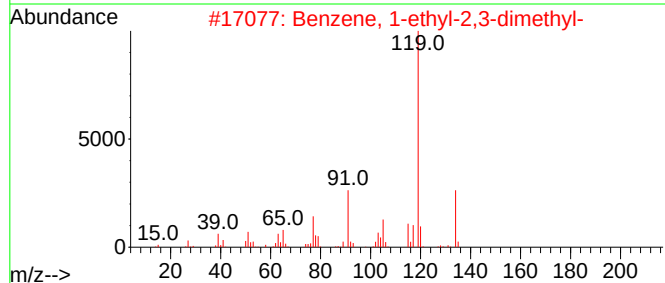
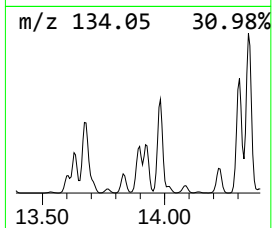
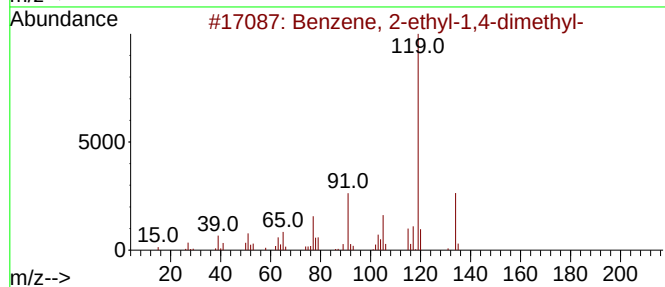
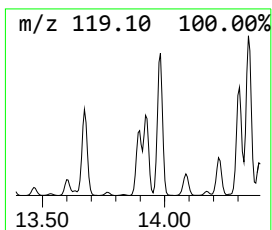
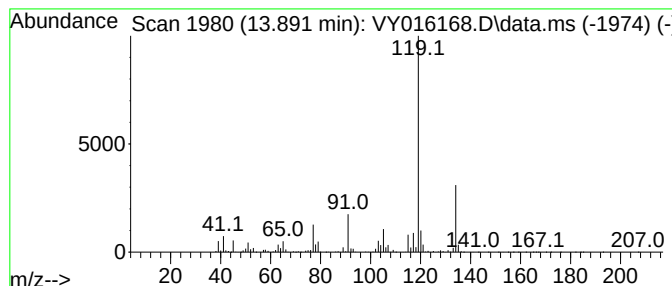
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.891	55.88 ug/l	3575440	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

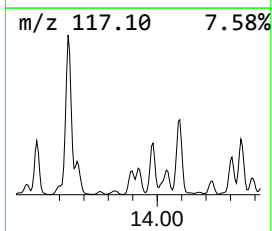
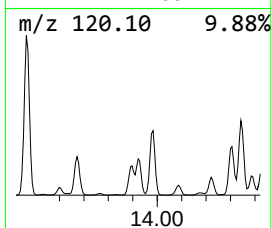
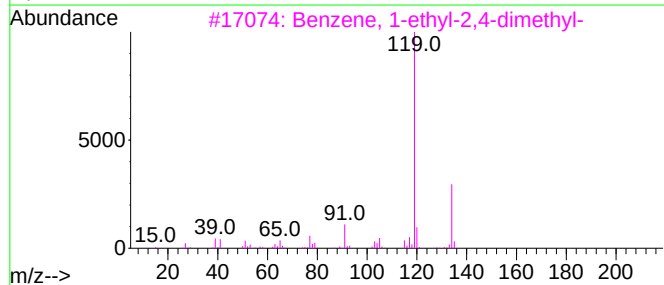
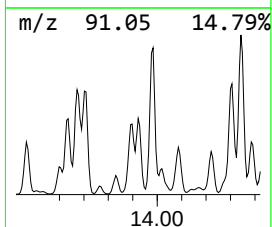
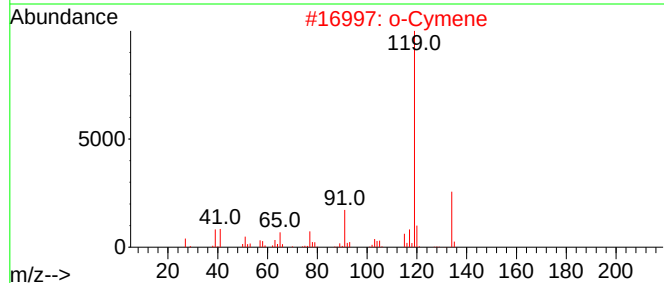
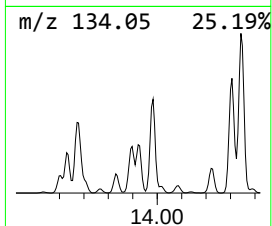
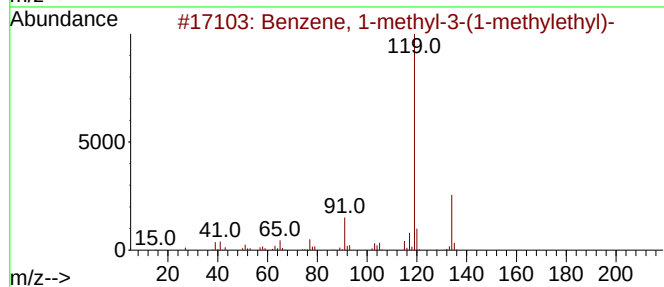
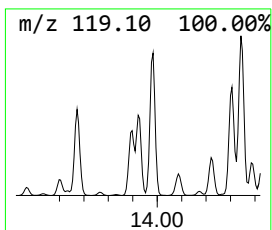
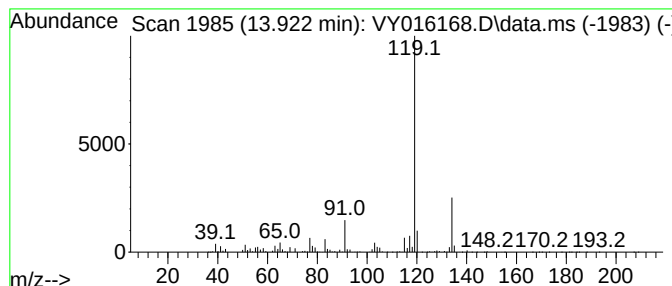
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-methyl-3-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.922	44.44 ug/l	2843280	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
2			o-Cymene	134	C10H14	000527-84-4	94
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			p-Cymene	134	C10H14	000099-87-6	94
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

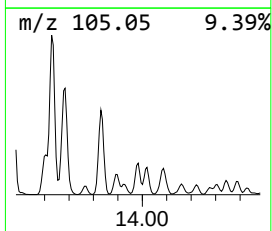
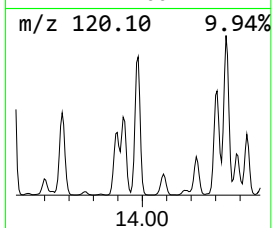
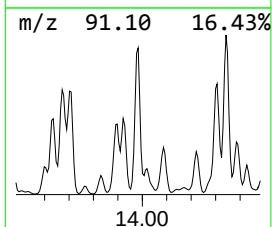
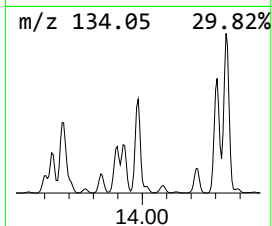
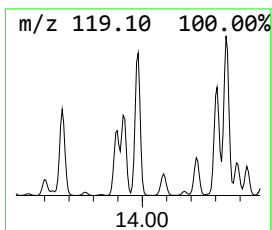
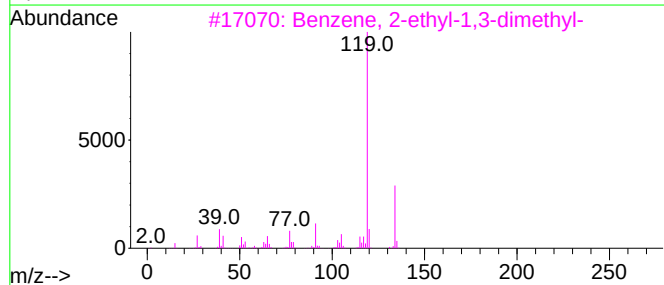
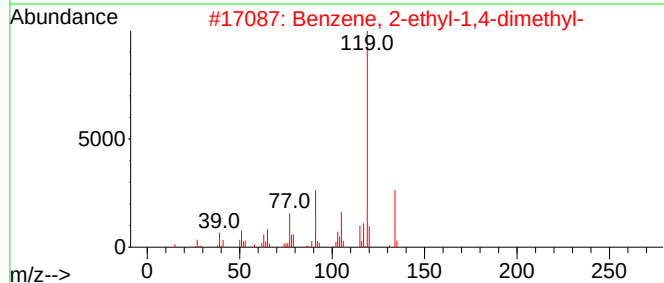
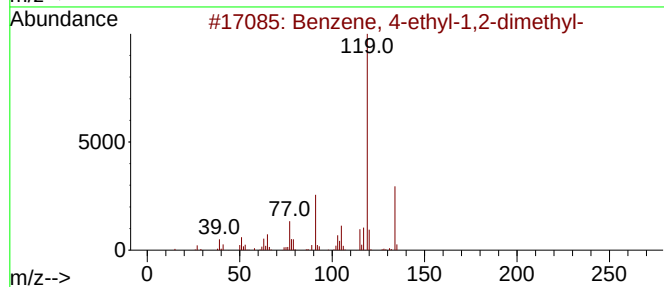
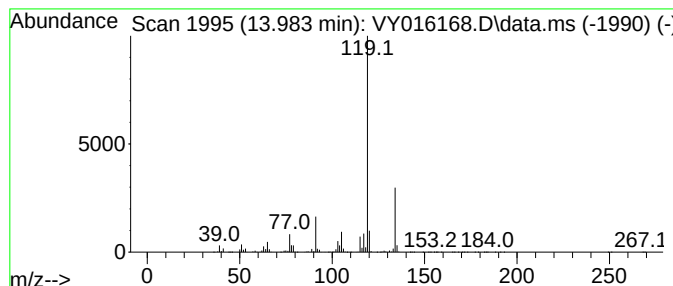
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.983	78.38 ug/l	5015110	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

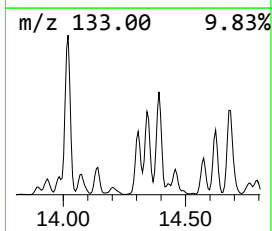
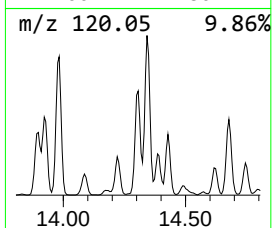
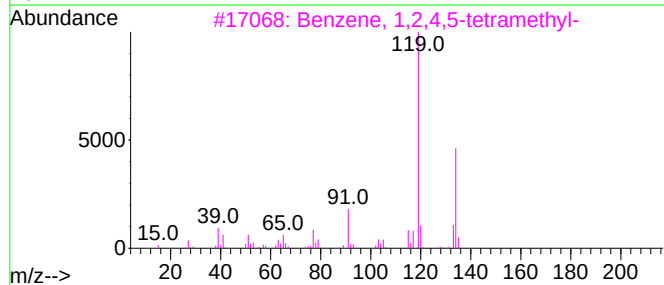
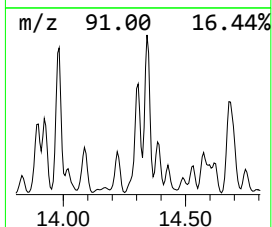
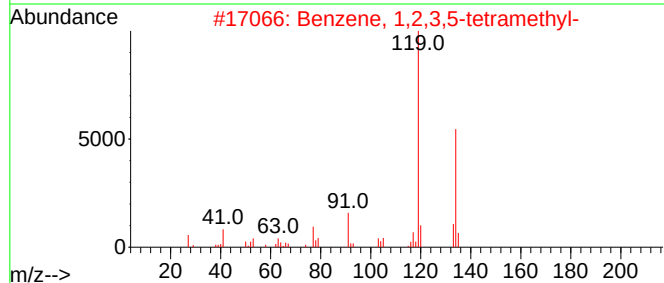
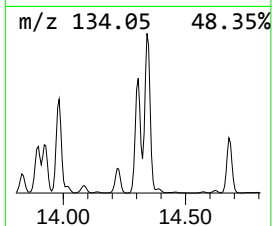
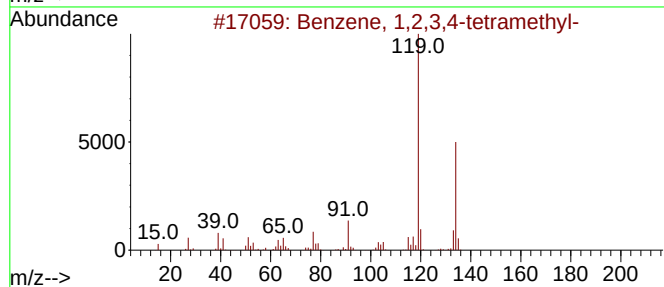
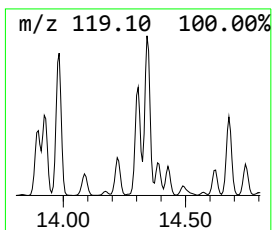
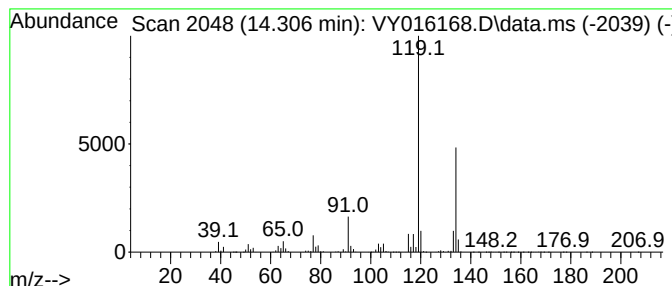
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.306	74.26 ug/l	4751250	1,4-Dichlorobenzene-d4	13.434

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, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

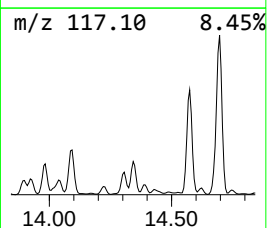
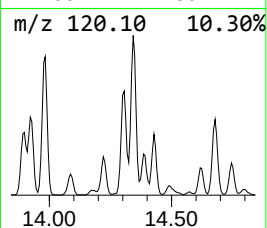
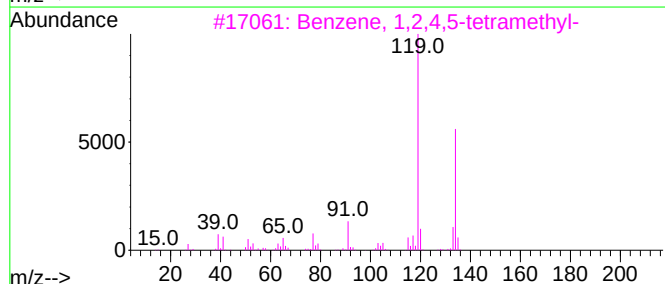
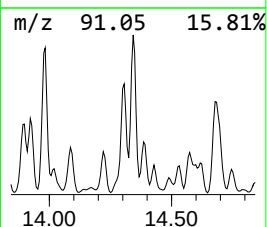
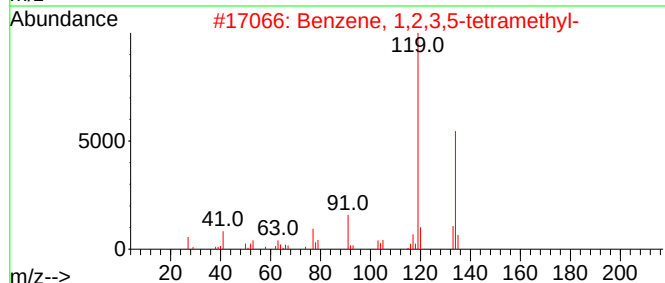
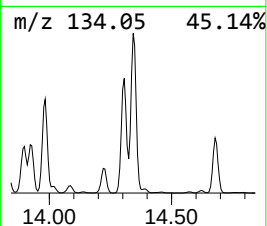
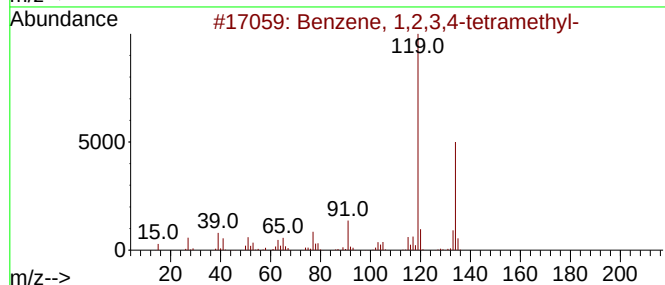
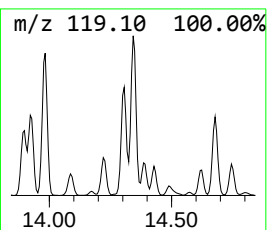
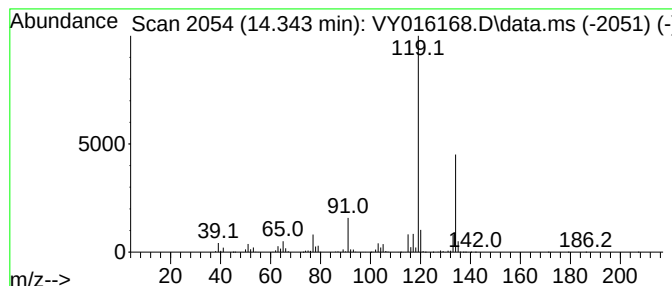
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.342	88.88 ug/l	5686850	1,4-Dichlorobenzene-d4	13.434

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	96
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

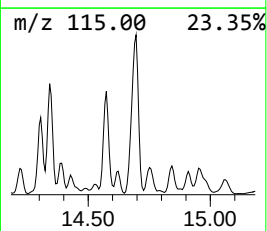
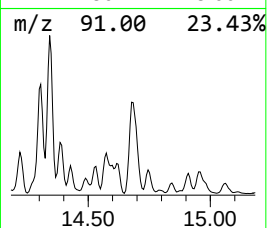
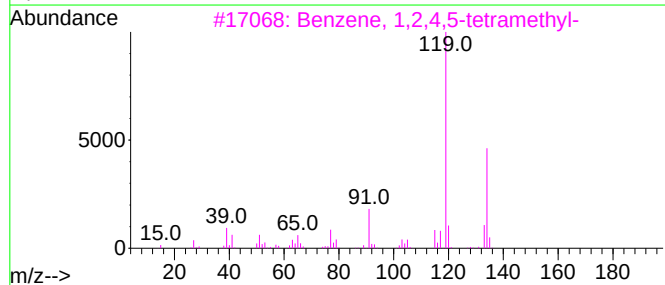
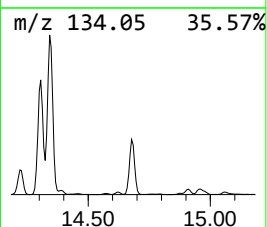
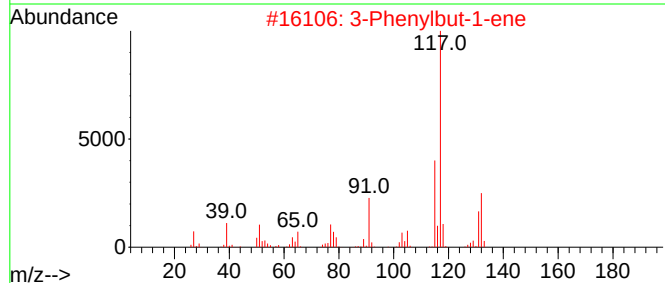
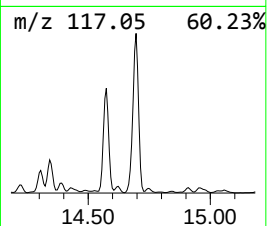
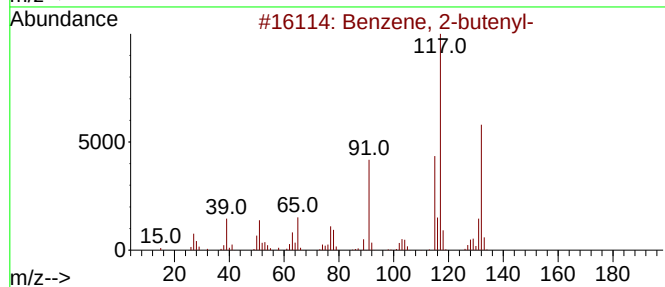
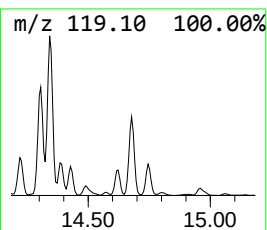
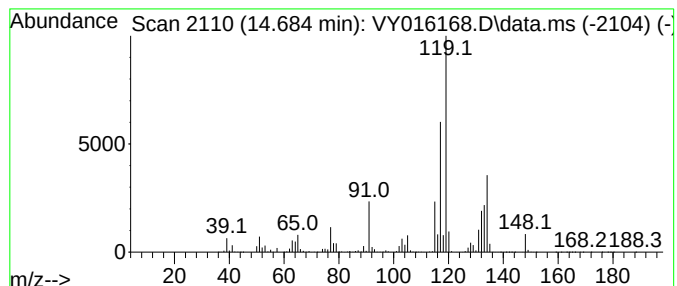
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	90.74 ug/l	5806060	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	80
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	55
4			o-Cymene	134	C10H14	000527-84-4	55
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110123\
Data File : VY016168.D
Acq On : 01 Nov 2023 15:03
Operator : SY/MD
Sample : 05102-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
67S-SB07-1418

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y103123S.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
1-Chloro-2-etho...	9.831	73.6	ug/l	2180280	2	8.697	1481770	50.0
Benzene, 1-ethy...	12.983	34.3	ug/l	2191180	4	13.434	3199130	50.0
Propane, 1-(chl...	13.507	165.8	ug/l	10606000	4	13.434	3199130	50.0
Benzene, 1-ethe...	13.635	58.5	ug/l	3745310	4	13.434	3199130	50.0
Benzene, 2-ethy...	13.891	55.9	ug/l	3575440	4	13.434	3199130	50.0
Benzene, 1-meth...	13.922	44.4	ug/l	2843280	4	13.434	3199130	50.0
Benzene, 4-ethy...	13.983	78.4	ug/l	5015110	4	13.434	3199130	50.0
Benzene, 1,2,3,...	14.306	74.3	ug/l	4751250	4	13.434	3199130	50.0
Benzene, 1,2,3,...	14.342	88.9	ug/l	5686850	4	13.434	3199130	50.0
Benzene, 2-bute...	14.684	90.7	ug/l	5806060	4	13.434	3199130	50.0