

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
 Data File : VX026660.D
 Acq On : 28 Jan 2022 18:40
 Operator : JC/MD
 Sample : N1297-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GES-1

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\82X011122W.M
 Title : SW846 8260

Signal : TIC: VX026660.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.752	104	109	126	rVB	134785	187564	2.32%	0.364%
2	2.221	181	186	195	rVB	76661	120322	1.49%	0.234%
3	2.873	289	293	308	rVB2	100560	217609	2.69%	0.423%
4	3.105	321	331	345	rVB	53872	122771	1.52%	0.238%
5	3.739	428	435	443	rVB	40510	94887	1.17%	0.184%
6	4.111	486	496	507	rVB	35370	90546	1.12%	0.176%
7	4.312	519	529	540	rVB	93621	264571	3.27%	0.514%
8	5.202	652	675	693	rVB2	65598	210541	2.60%	0.409%
9	5.391	693	706	714	rBV	77600	225803	2.79%	0.438%
10	5.483	714	721	726	rVV	42665	114317	1.41%	0.222%
11	5.556	726	733	742	rVB	116379	312388	3.86%	0.607%
12	5.958	790	799	807	rBV	84100	219046	2.71%	0.425%
13	6.044	807	813	825	rVB2	45262	116932	1.44%	0.227%
14	6.202	825	839	846	rBV2	42122	155739	1.92%	0.302%
15	6.769	922	932	941	rBV	203766	540351	6.67%	1.049%
16	7.385	1022	1033	1044	rBV4	45719	128598	1.59%	0.250%
17	8.147	1154	1158	1163	rVB	51454	81535	1.01%	0.158%
18	8.507	1209	1217	1225	rVB	67572	114268	1.41%	0.222%
19	8.653	1235	1241	1246	rBV	416408	650175	8.03%	1.262%
20	8.720	1246	1252	1264	rVB	2108561	3220261	39.77%	6.253%
21	10.055	1461	1471	1476	rBV	419832	569726	7.04%	1.106%
22	10.201	1489	1495	1500	rBV	6046531	8096779	100.00%	15.721%
23	10.305	1505	1512	1527	rVB	3883986	5260197	64.97%	10.214%
24	10.646	1562	1568	1578	rVB	3721449	4842294	59.81%	9.402%
25	10.963	1614	1620	1628	rVB	422382	525717	6.49%	1.021%
26	11.085	1634	1640	1645	rBV	309137	395890	4.89%	0.769%
27	11.305	1671	1676	1683	rVB	1111057	1322177	16.33%	2.567%
28	11.384	1683	1689	1690	rBV	1117712	1416478	17.49%	2.750%
29	11.402	1690	1692	1696	rVV	1116510	1180232	14.58%	2.292%
30	11.451	1696	1700	1707	rVB	1062057	1342248	16.58%	2.606%
31	11.616	1721	1727	1738	rBV	1513158	1843078	22.76%	3.579%
32	11.756	1744	1750	1756	rBV	4052076	4844624	59.83%	9.407%
33	12.024	1789	1794	1799	rVB	414176	525272	6.49%	1.020%
34	12.091	1799	1805	1809	rBV	906082	1129016	13.94%	2.192%
35	12.244	1825	1830	1838	rBV2	1902868	2557410	31.59%	4.966%
36	12.317	1838	1842	1852	rVB4	383225	560524	6.92%	1.088%

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Filtering: 5

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 >

Peak separation: 5

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37	12.414	1852	1858	1862	rBV2	211253	284349	3.51%	0.552%
38	12.463	1862	1866	1872	rVB	92134	112988	1.40%	0.219%
39	12.530	1872	1877	1880	rBV	292124	388807	4.80%	0.755%
40	12.616	1886	1891	1894	rBV	747755	863323	10.66%	1.676%
41	12.646	1894	1896	1900	rVB	112510	120385	1.49%	0.234%
42	12.701	1900	1905	1913	rBV	345320	473433	5.85%	0.919%
43	12.926	1934	1942	1945	rBV	332845	435443	5.38%	0.845%
44	12.963	1945	1948	1954	rVB	458302	539973	6.67%	1.048%
45	13.170	1978	1982	1987	rVB	349167	398210	4.92%	0.773%
46	13.292	1998	2002	2007	rVB2	718303	919260	11.35%	1.785%
47	13.426	2019	2024	2032	rVB	171906	219679	2.71%	0.427%
48	13.585	2046	2050	2056	rVB3	66282	112924	1.39%	0.219%
49	13.664	2061	2063	2069	rVB2	62175	94143	1.16%	0.183%
50	13.780	2076	2082	2090	rVV	1439517	1832419	22.63%	3.558%
51	13.871	2090	2097	2101	rVB	60531	82970	1.02%	0.161%
52	14.219	2149	2154	2159	rBV2	48756	82464	1.02%	0.160%
53	14.640	2218	2223	2233	rVB	491426	610718	7.54%	1.186%
54	14.780	2241	2246	2262	rVB	259061	330162	4.08%	0.641%

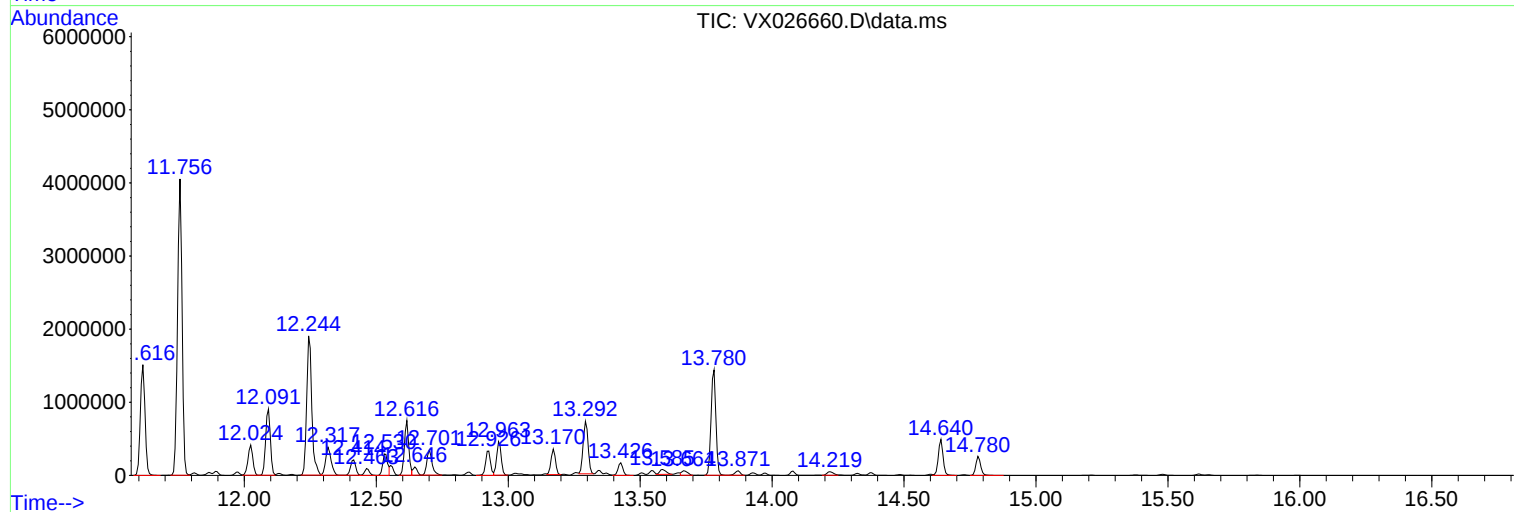
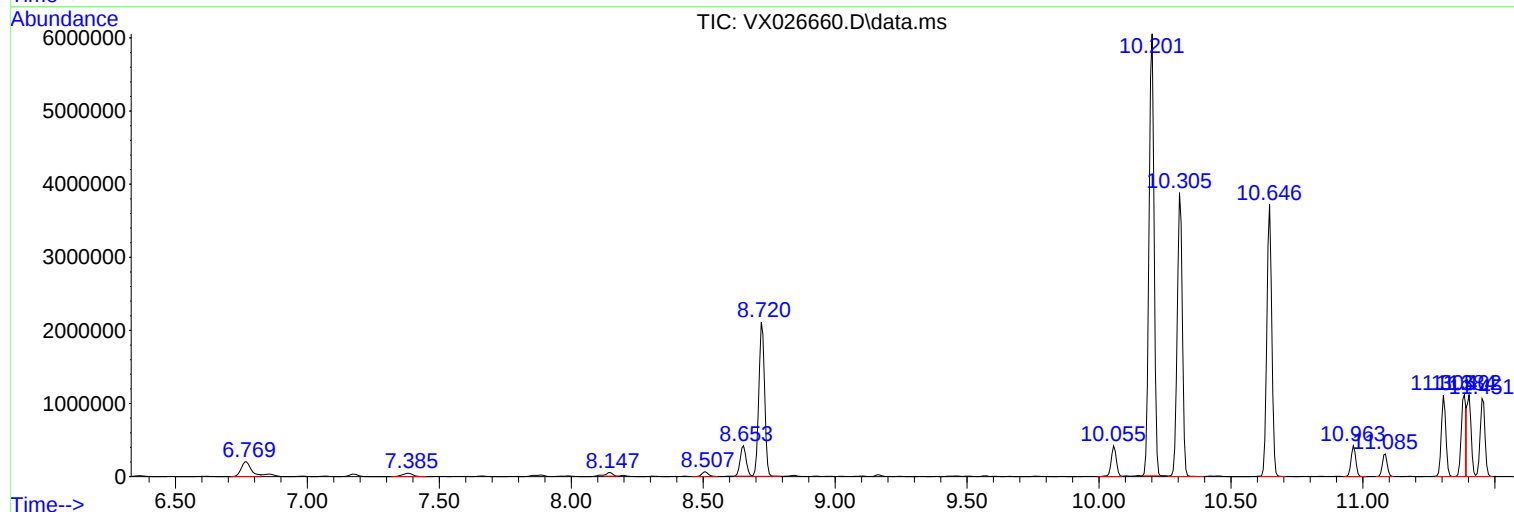
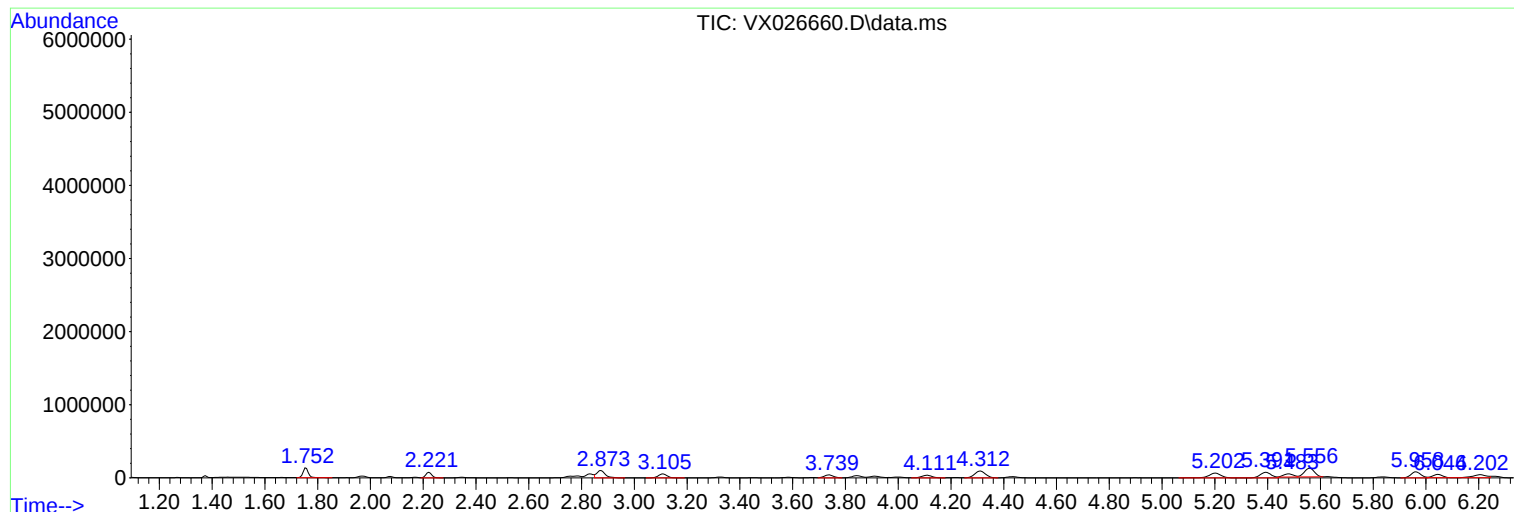
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Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampleId :
GES-1

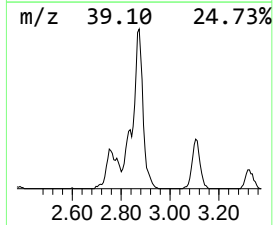
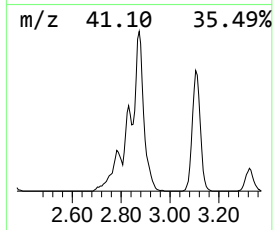
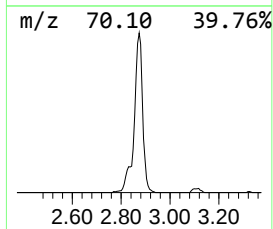
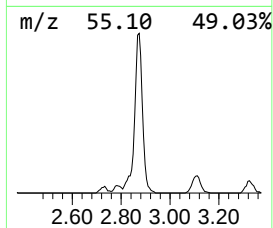
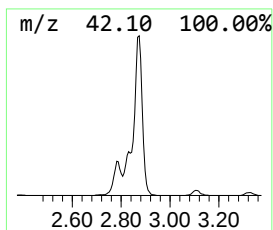
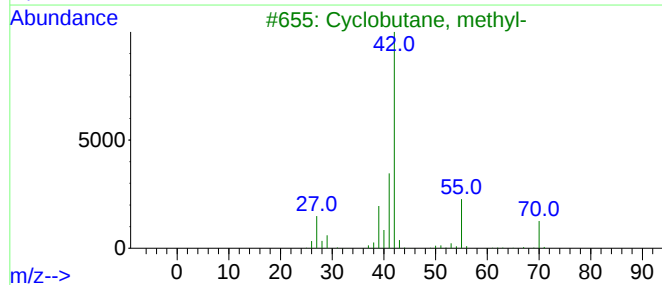
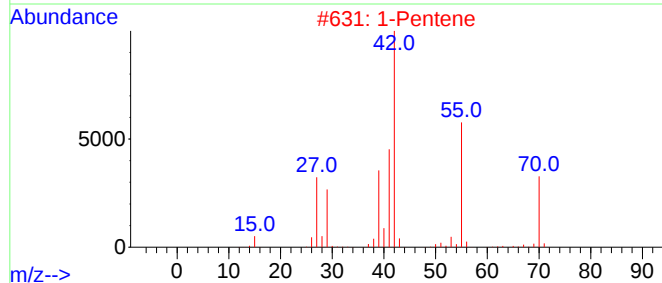
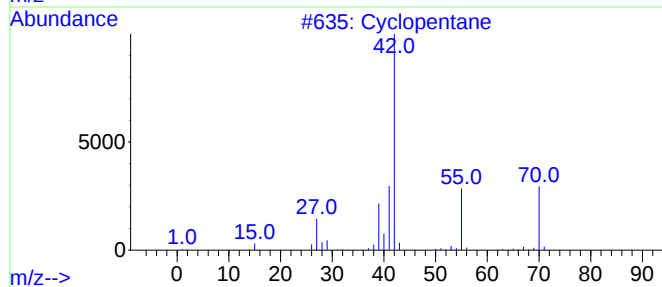
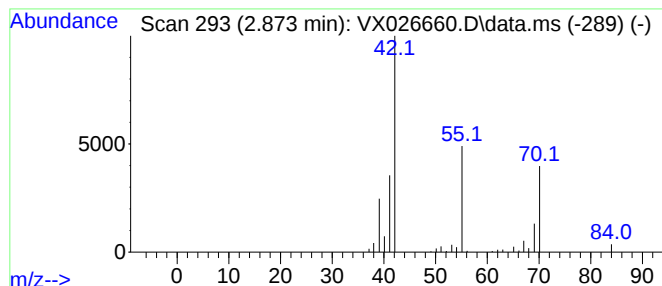
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.873	34.83 ug/l	217609	Pentafluorobenzene	5.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	64
2			1-Pentene	70	C5H10	000109-67-1	64
3			Cyclobutane, methyl-	70	C5H10	000598-61-8	64
4			3-Buten-1-ol	72	C4H8O	000627-27-0	37
5			1-Pentanol	88	C5H12O	000071-41-0	9



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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GES-1

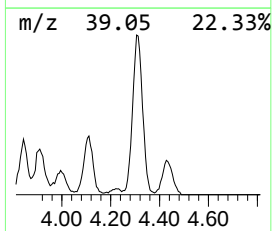
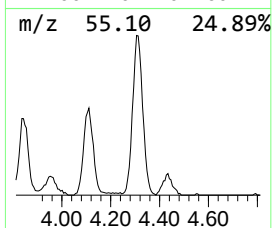
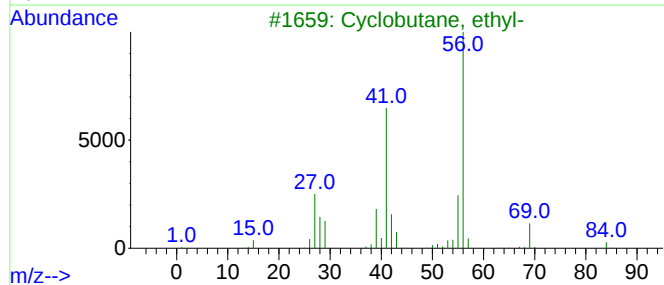
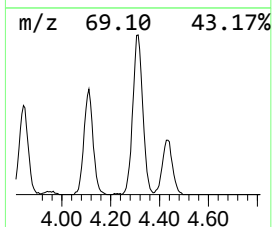
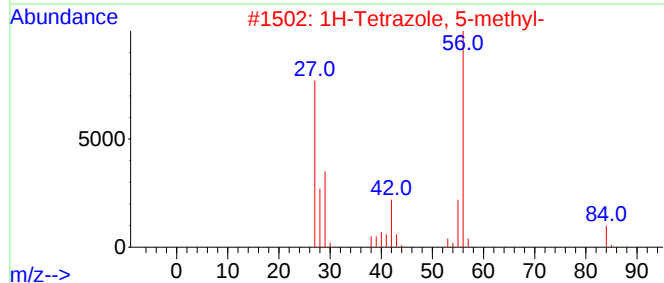
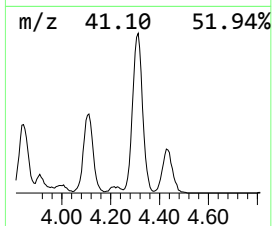
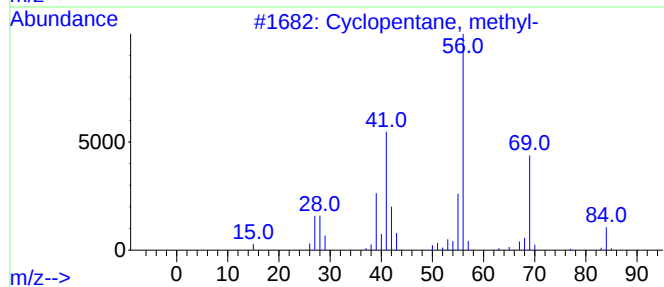
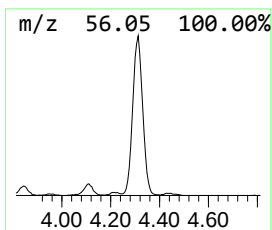
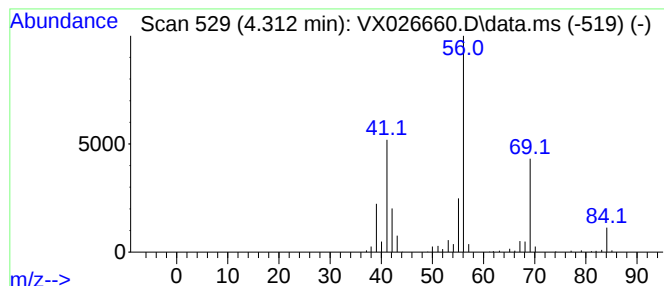
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.312	42.35 ug/l	264571	Pentafluorobenzene	5.562

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50
5		Cyclobutane	56	C4H8	000287-23-0	50



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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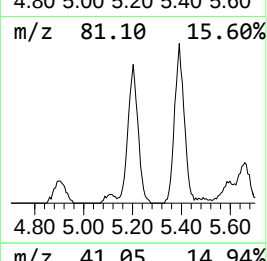
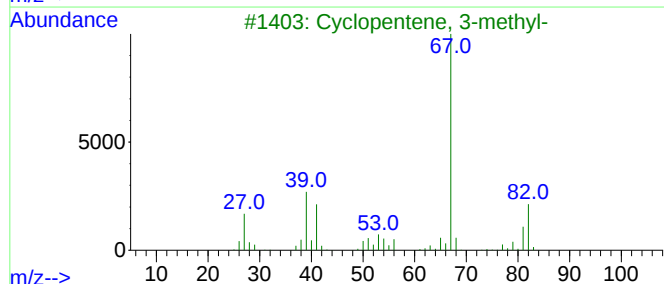
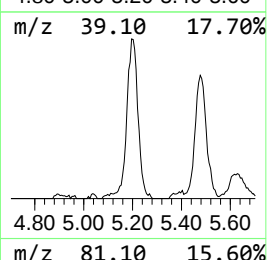
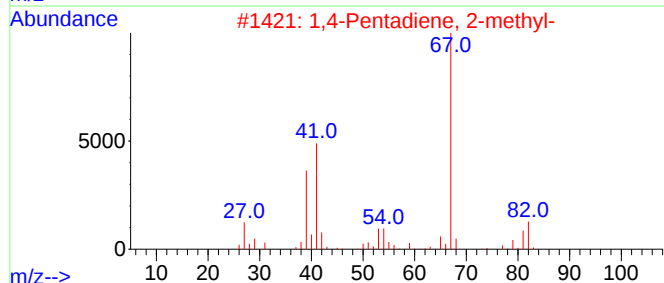
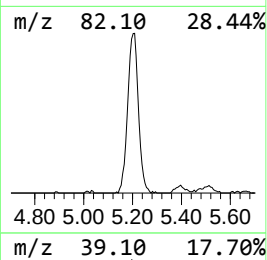
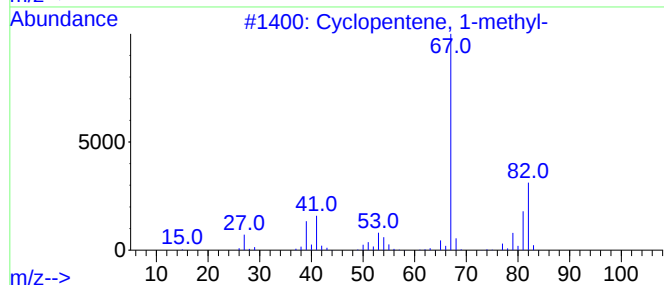
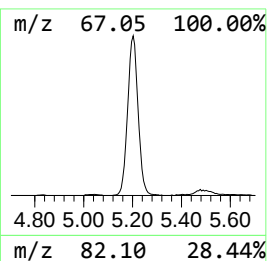
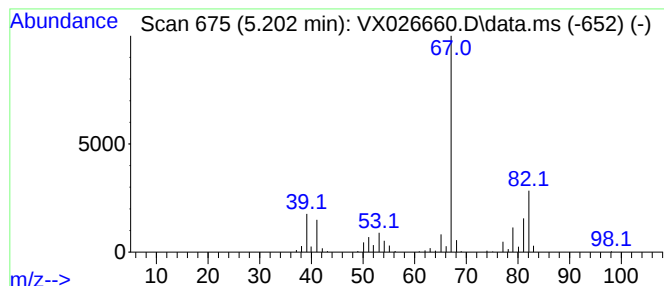
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.202	33.70 ug/l	210541	Pentafluorobenzene	5.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	90
2			1,4-Pentadiene, 2-methyl-	82	C6H10	000763-30-4	90
3			Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	86
4			Cyclopentane, methylene-	82	C6H10	001528-30-9	83
5			1,4-Hexadiene, (Z)-	82	C6H10	007318-67-4	80



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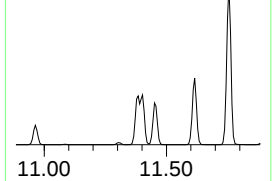
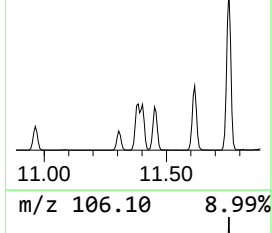
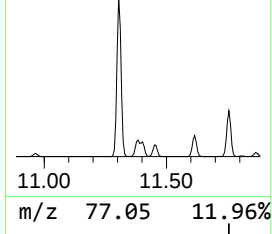
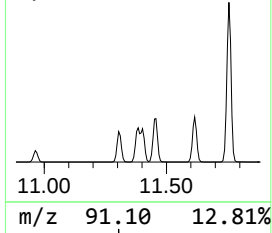
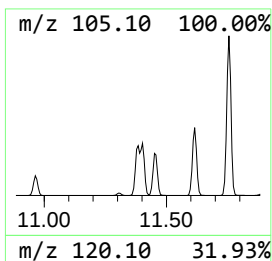
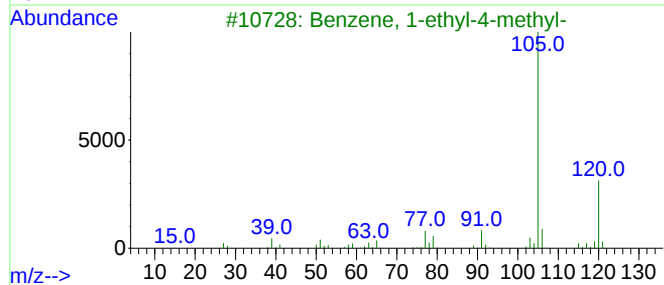
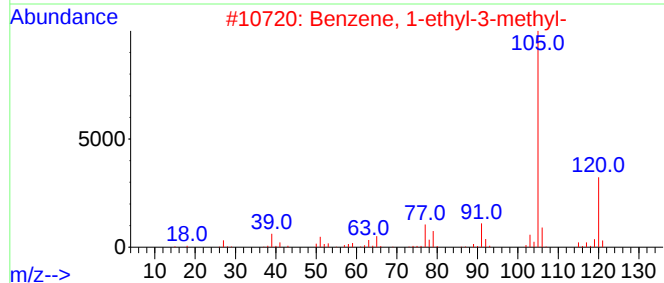
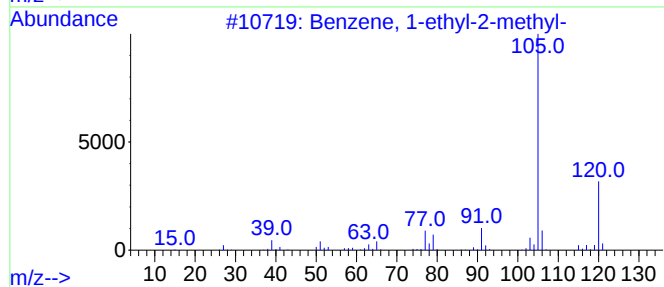
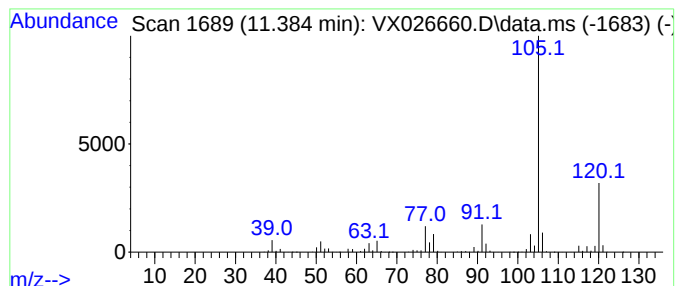
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	134.83 ug/l	1416480	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	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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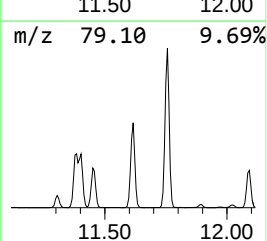
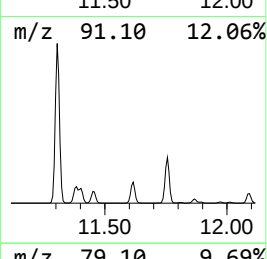
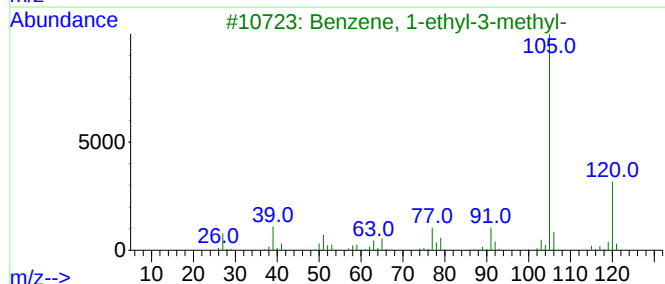
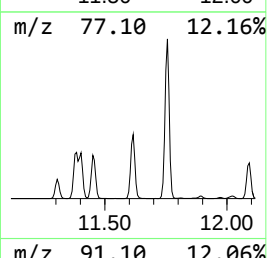
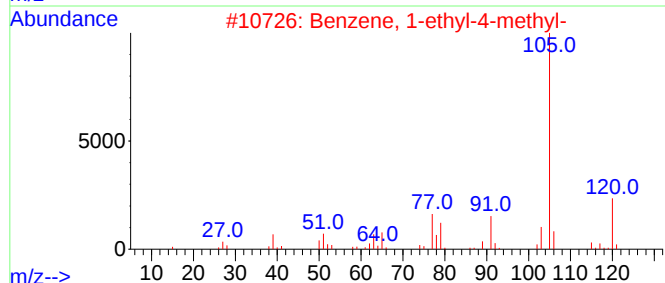
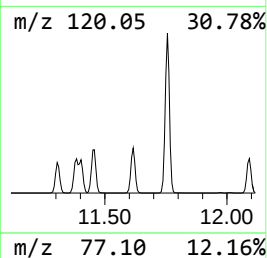
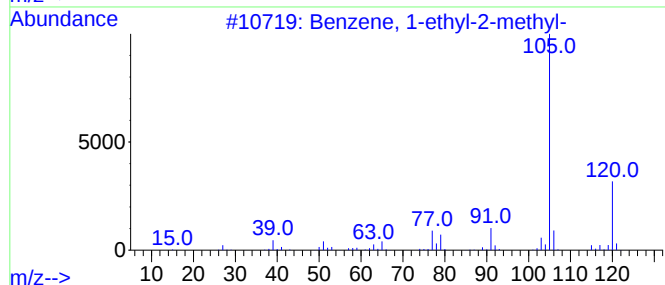
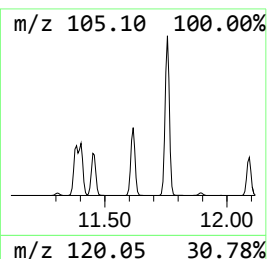
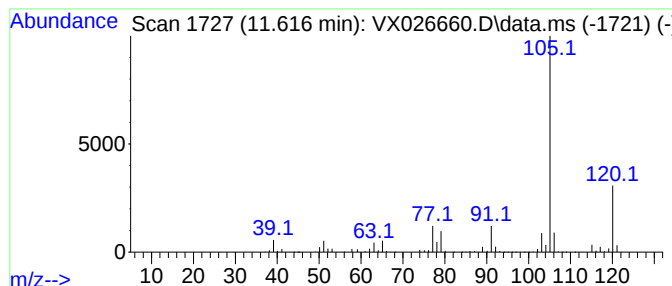
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	175.44 ug/l	1843080	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

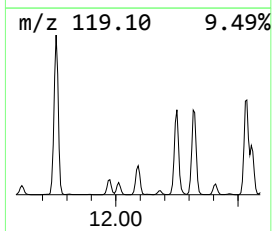
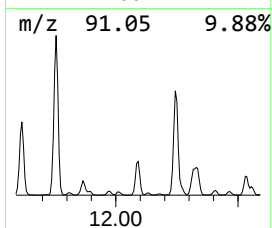
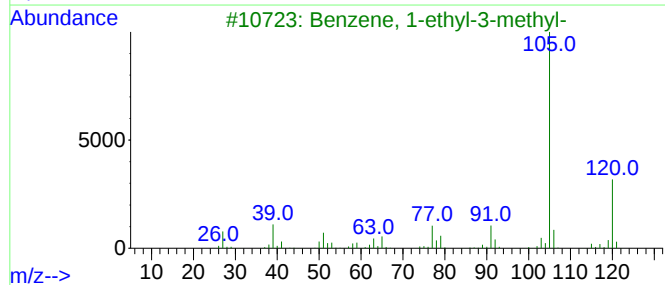
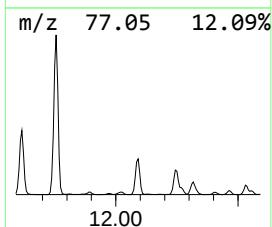
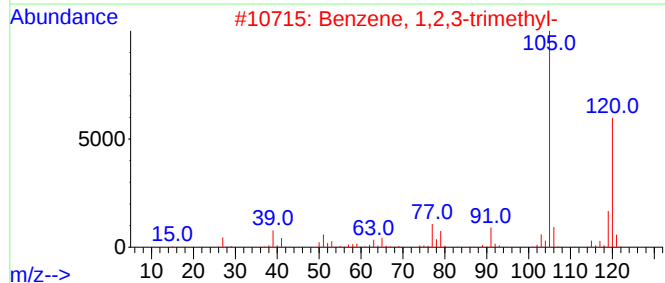
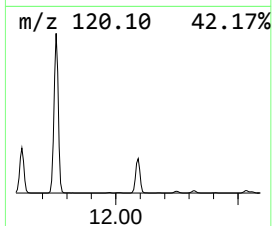
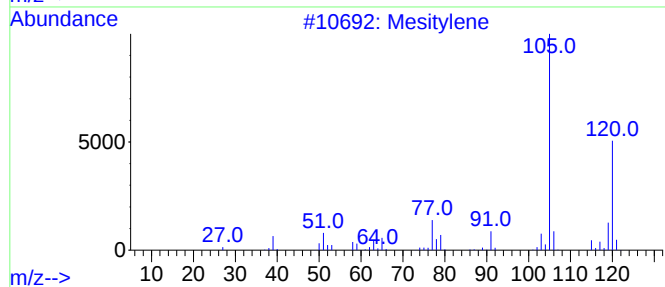
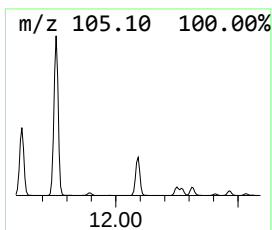
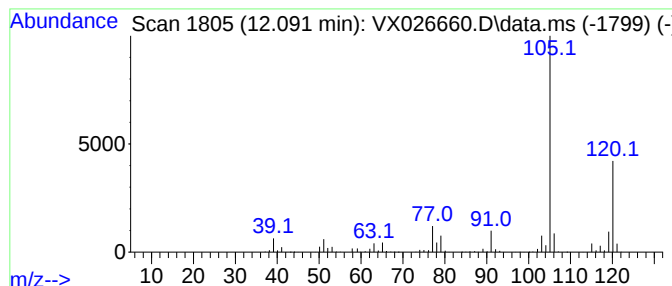
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	107.47 ug/l	1129020	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

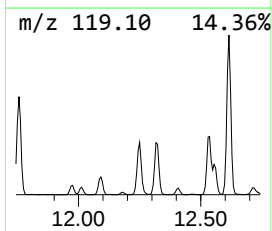
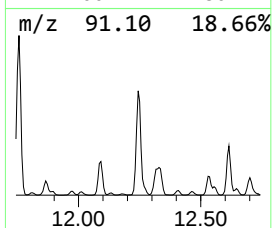
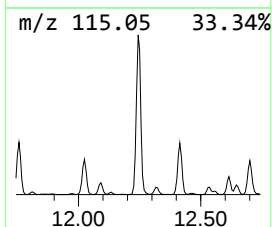
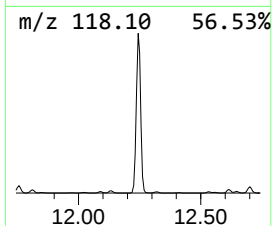
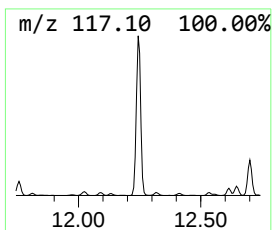
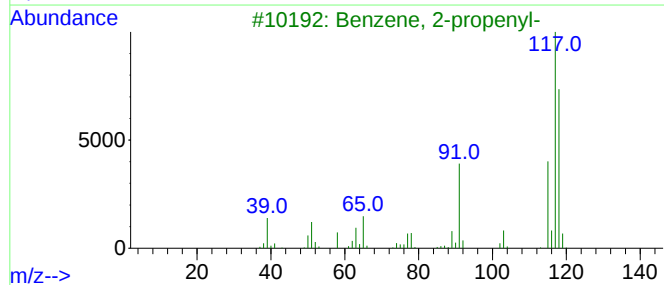
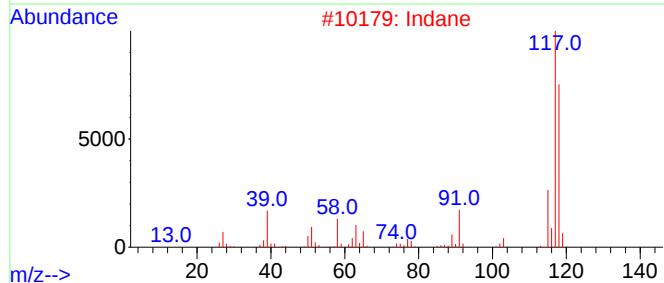
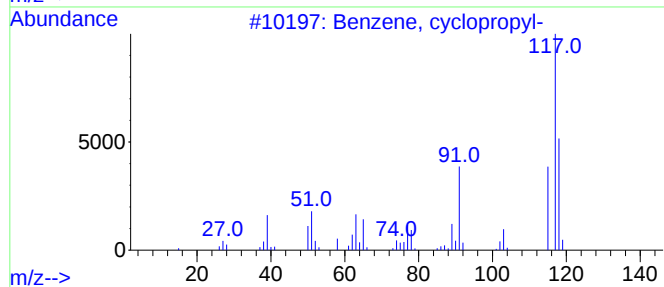
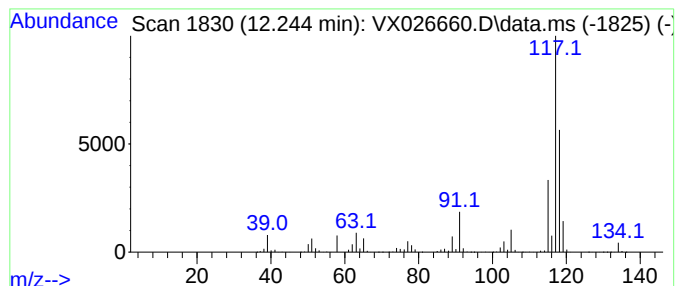
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	243.44 ug/l	2557410	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			Δ-tetralene	118	C9H10	007785-10-6	58
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

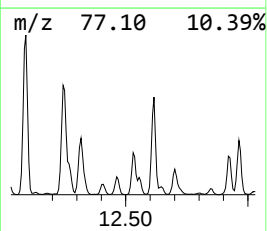
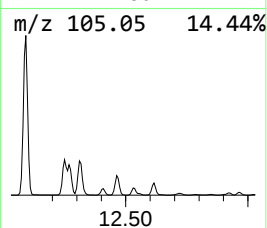
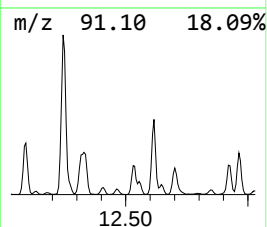
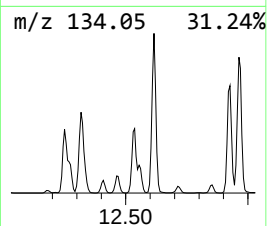
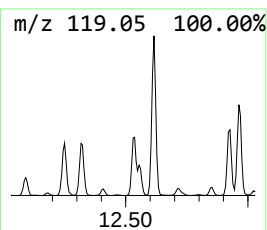
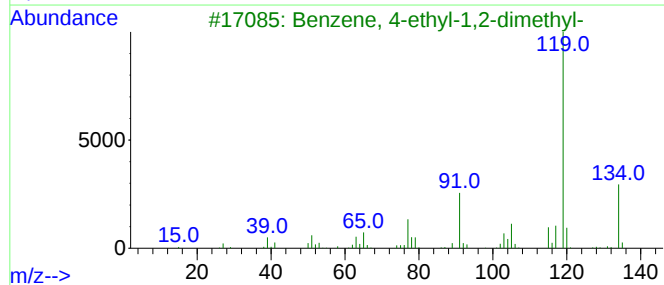
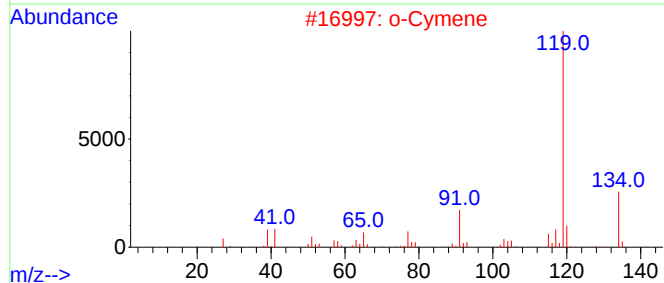
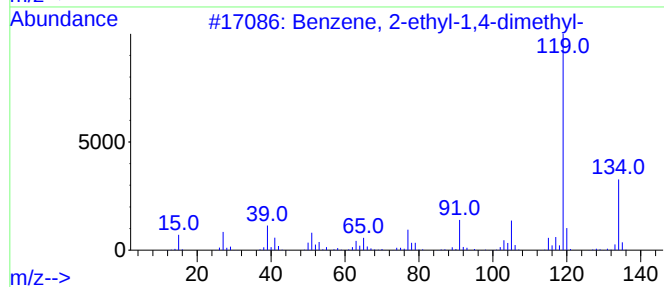
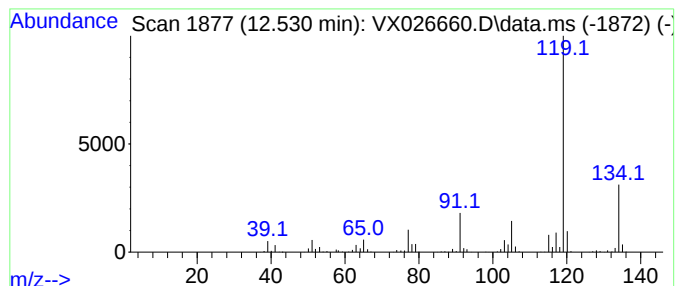
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.530	37.01 ug/l	388807	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

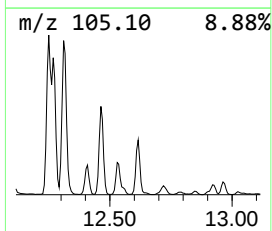
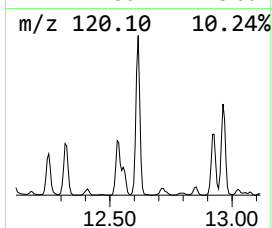
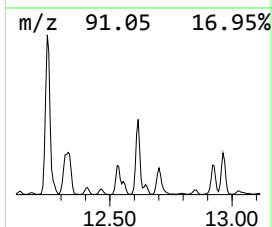
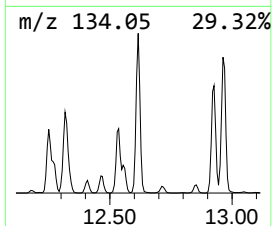
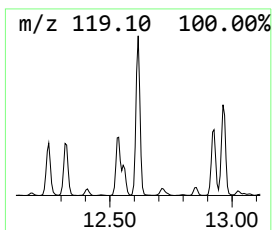
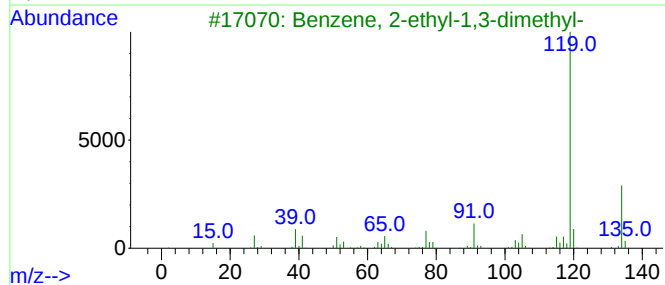
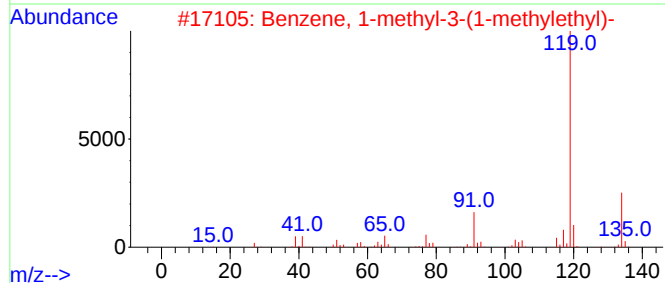
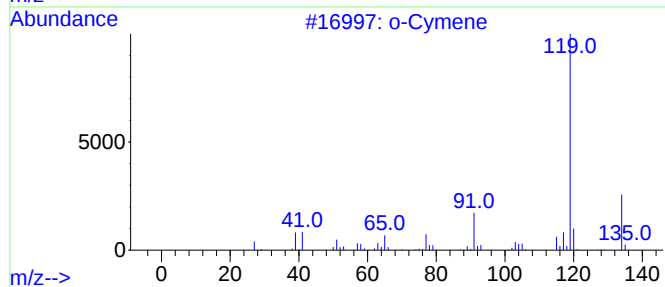
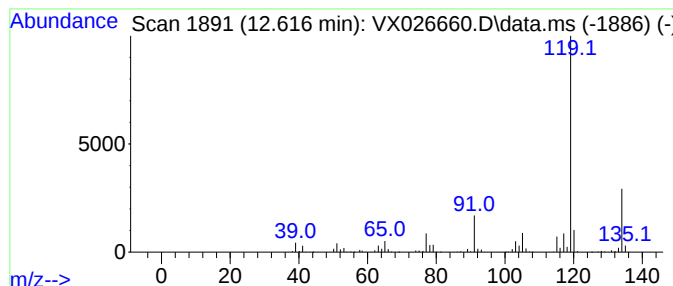
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Quant Title : SW846 8260

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

Peak Number 9 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	82.18 ug/l	863323	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

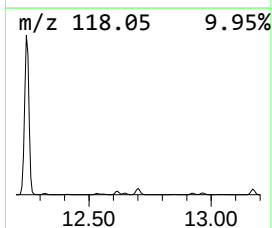
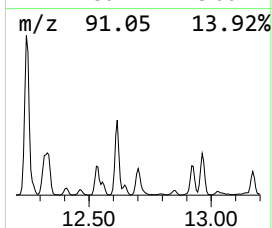
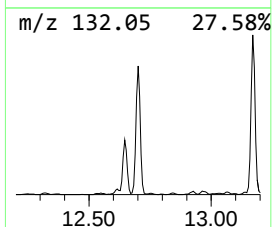
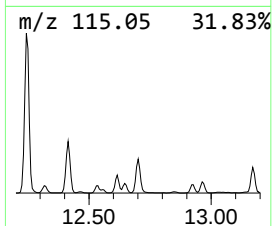
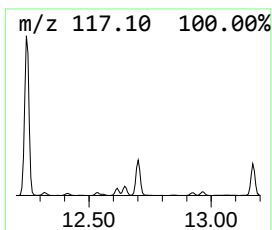
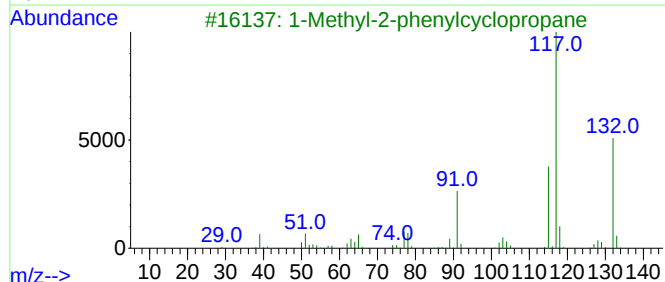
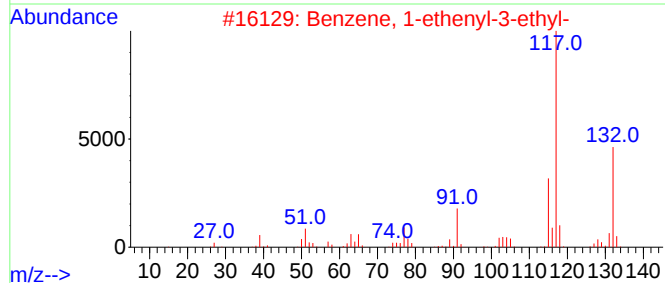
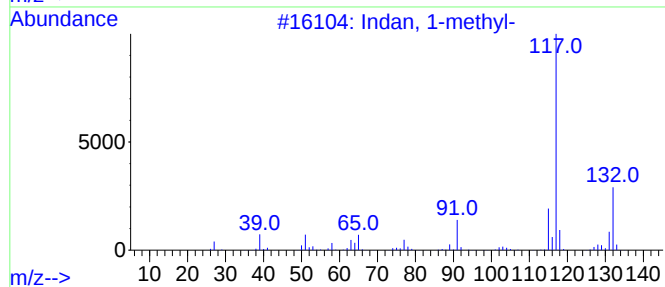
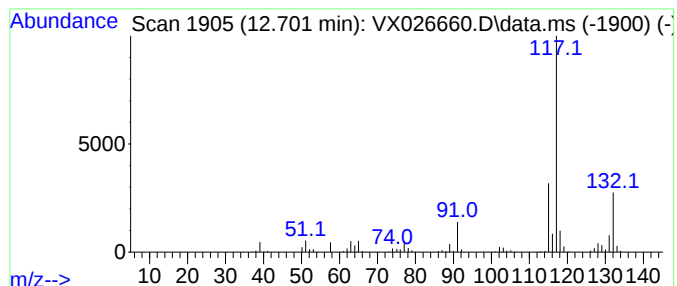
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Quant Title : SW846 8260

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

Peak Number 10 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	45.07 ug/l	473433	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	93
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

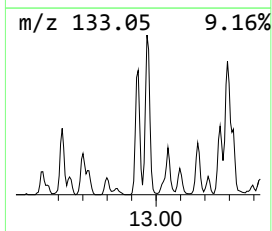
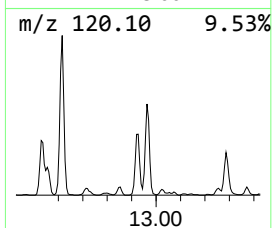
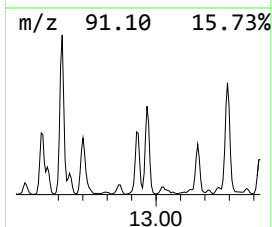
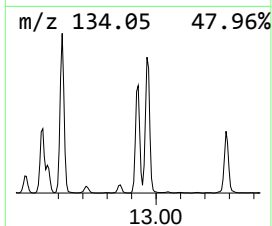
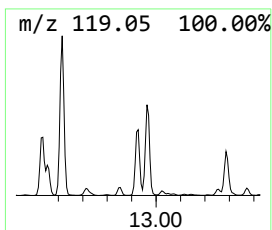
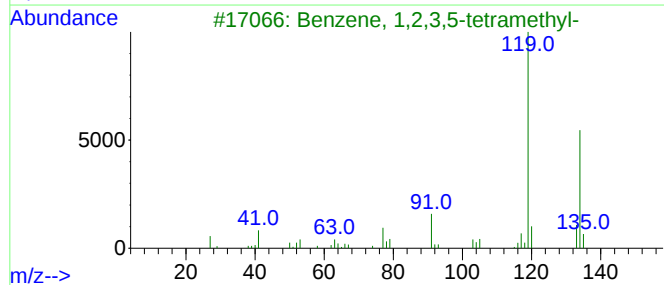
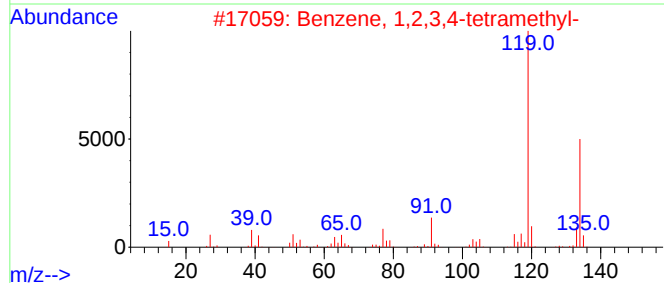
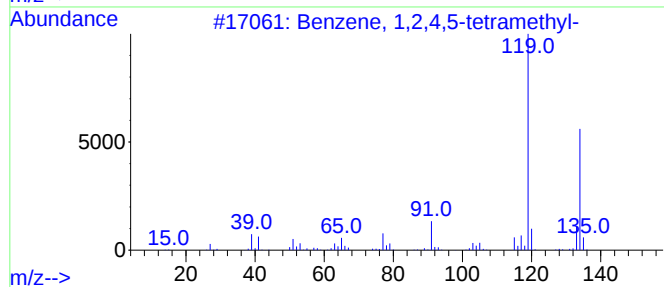
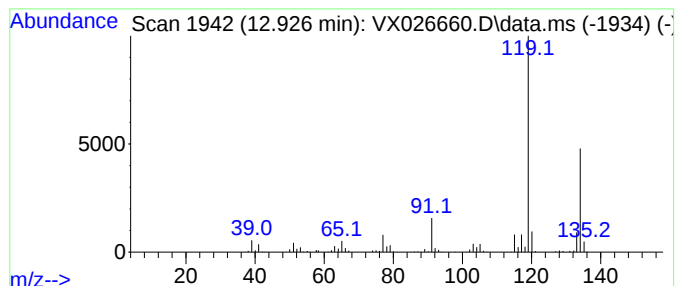
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Quant Title : SW846 8260

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

Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	41.45 ug/l	435443	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

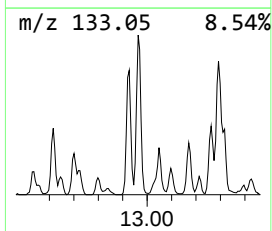
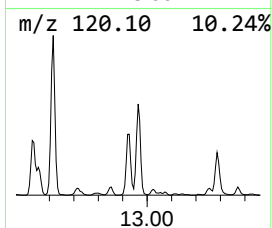
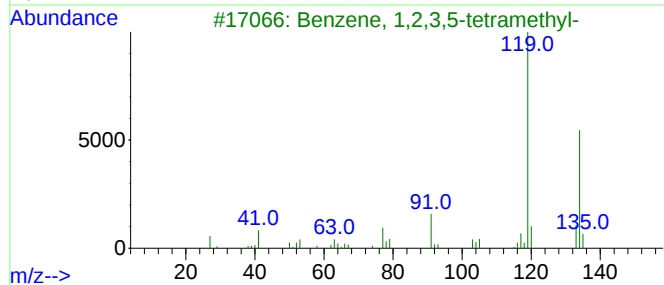
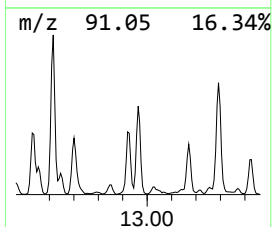
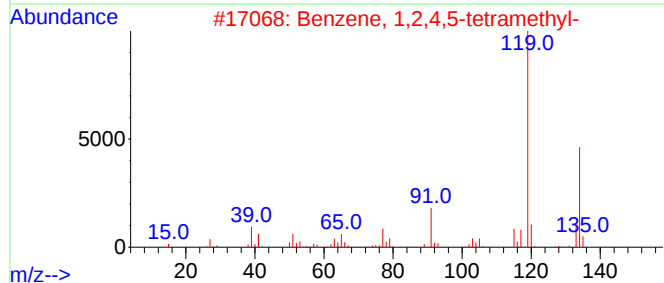
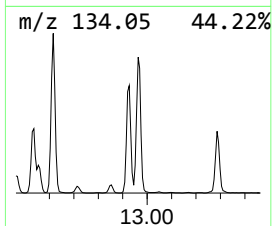
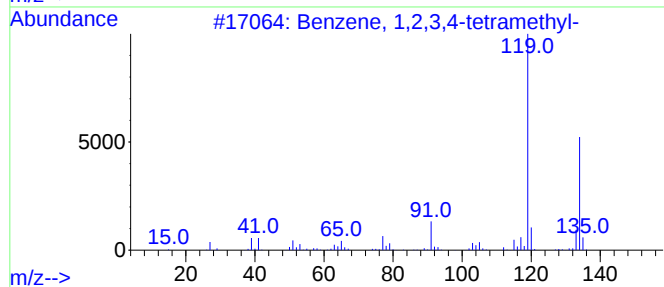
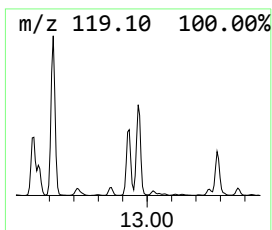
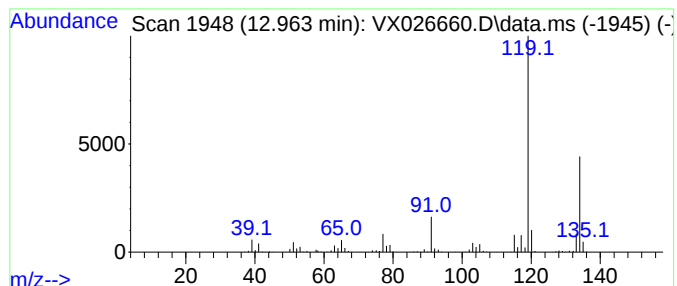
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 12 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	51.40 ug/l	539973	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, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

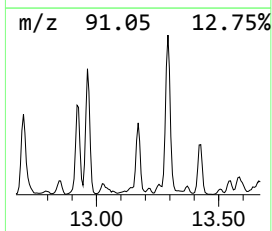
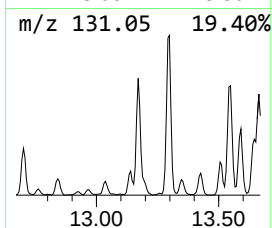
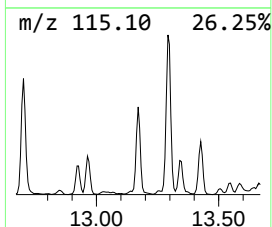
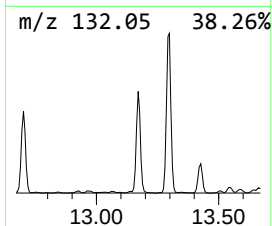
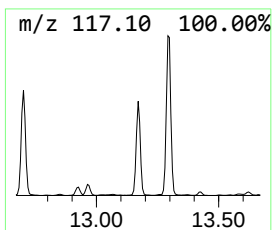
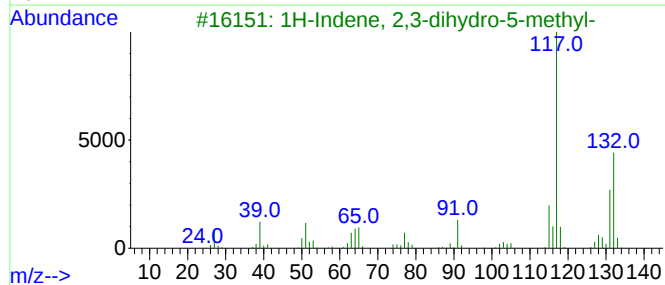
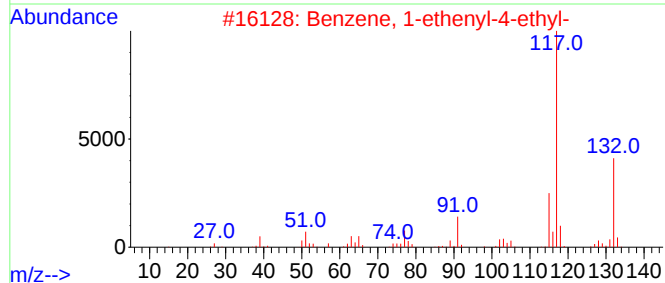
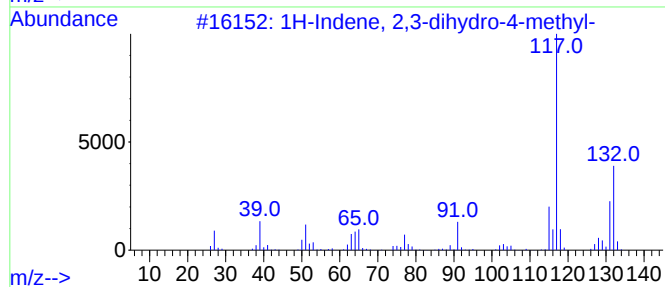
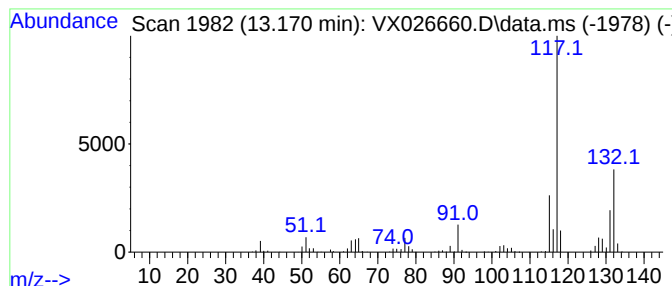
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Quant Title : SW846 8260

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

Peak Number 13 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	37.91 ug/l	398210	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	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

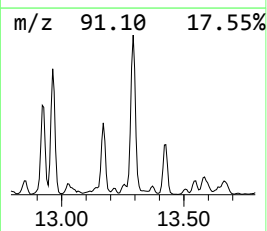
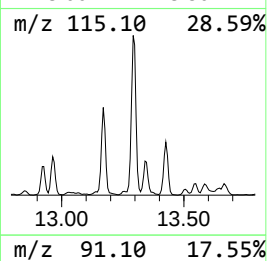
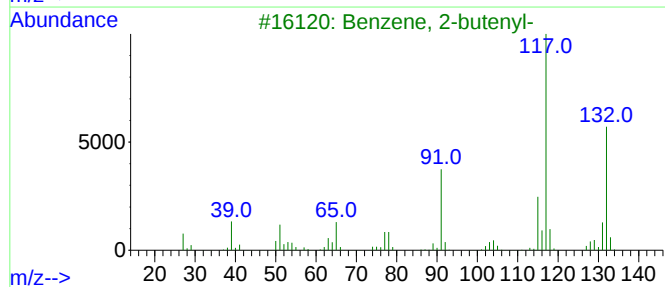
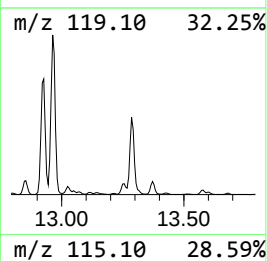
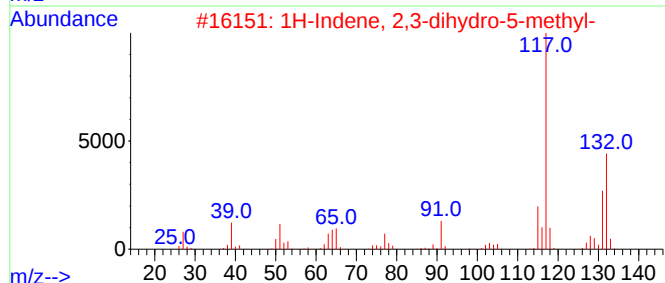
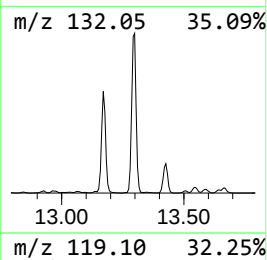
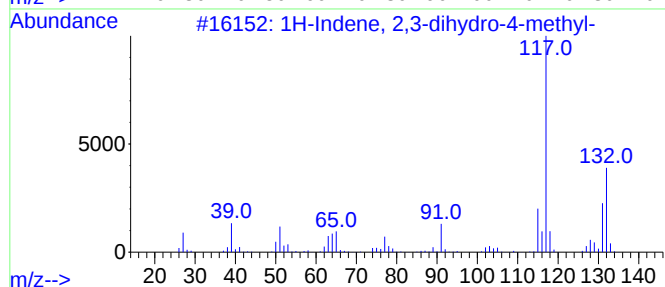
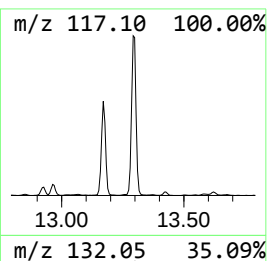
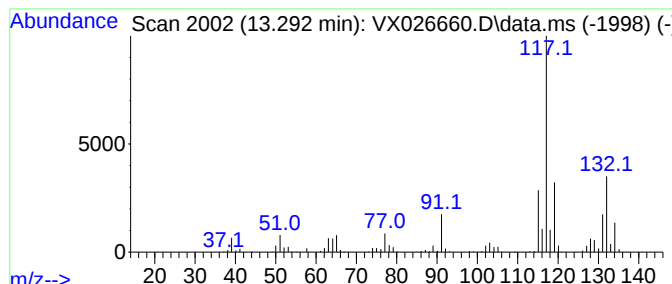
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 14 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	87.50 ug/l	919260	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	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

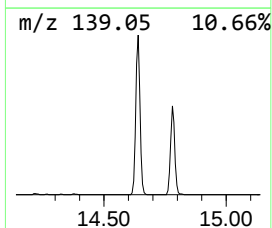
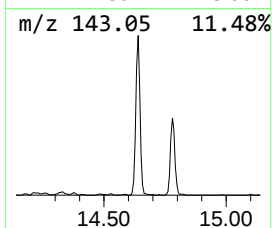
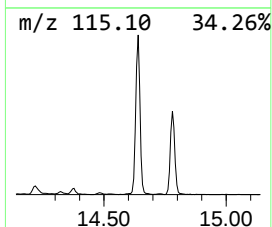
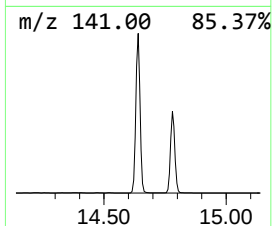
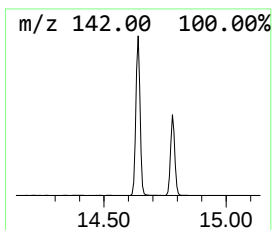
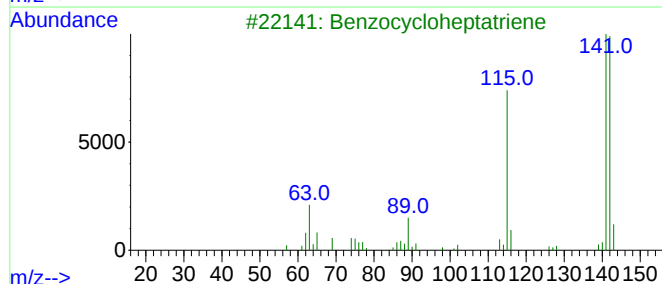
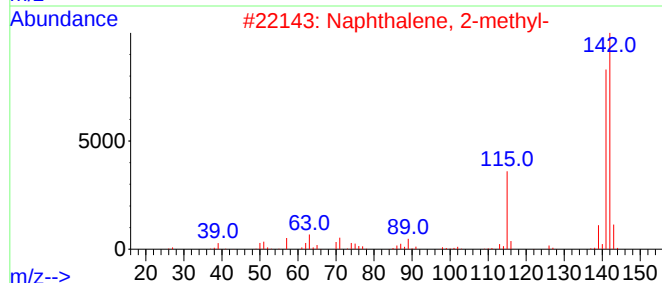
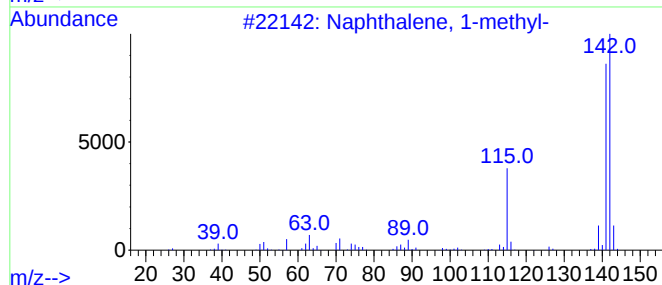
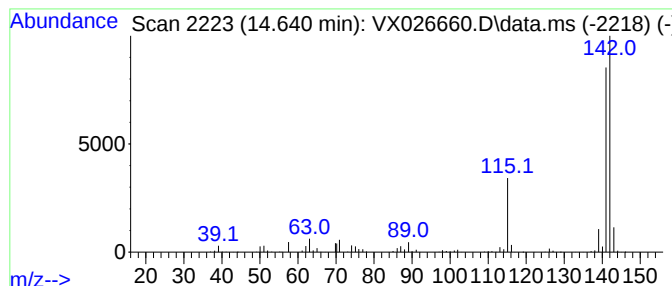
Instrument :
MSVOA_X
ClientSampleId :
GES-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260

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

Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.640	58.13 ug/l	610718	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	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012822\
Data File : VX026660.D
Acq On : 28 Jan 2022 18:40
Operator : JC/MD
Sample : N1297-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GES-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.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
Cyclopentane	2.873	34.8	ug/l	217609	1	5.562	312388	50.0
Cyclopentane, m...	4.312	42.4	ug/l	264571	1	5.562	312388	50.0
Cyclopentene, 1...	5.202	33.7	ug/l	210541	1	5.562	312388	50.0
Benzene, 1-ethy...	11.384	134.8	ug/l	1416480	4	12.024	525272	50.0
Benzene, 1-ethy...	11.616	175.4	ug/l	1843080	4	12.024	525272	50.0
Benzene, 1,2,3-...	12.091	107.5	ug/l	1129020	4	12.024	525272	50.0
Benzene, cyclop...	12.244	243.4	ug/l	2557410	4	12.024	525272	50.0
Benzene, 2-ethy...	12.530	37.0	ug/l	388807	4	12.024	525272	50.0
o-Cymene	12.616	82.2	ug/l	863323	4	12.024	525272	50.0
Indan, 1-methyl-	12.701	45.1	ug/l	473433	4	12.024	525272	50.0
Benzene, 1,2,4,...	12.926	41.5	ug/l	435443	4	12.024	525272	50.0
Benzene, 1,2,3,...	12.963	51.4	ug/l	539973	4	12.024	525272	50.0
1H-Indene, 2,3-...	13.170	37.9	ug/l	398210	4	12.024	525272	50.0
1H-Indene, 2,3-...	13.292	87.5	ug/l	919260	4	12.024	525272	50.0
Naphthalene, 1-...	14.640	58.1	ug/l	610718	4	12.024	525272	50.0