

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
 Data File : VN075800.D
 Acq On : 15 Dec 2022 05:15
 Operator : JC\MD
 Sample : N5990-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-43

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Title : SW846 8260

Signal : TIC: VN075800.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.336	887	898	908	rBV	760185	2180274	2.31%	0.372%
2	8.236	1044	1051	1062	rVB3	549251	1792397	1.90%	0.306%
3	8.612	1103	1115	1126	rBV3	757287	2159598	2.29%	0.369%
4	8.741	1126	1137	1143	rBV4	316415	1091835	1.16%	0.187%
5	9.106	1189	1199	1206	rBV3	909547	2073017	2.20%	0.354%
6	9.606	1268	1284	1299	rBV2	534902	1695029	1.80%	0.290%
7	10.571	1440	1448	1454	rBV	1014951	1885539	2.00%	0.322%
8	11.865	1661	1668	1676	rBV	889109	1684470	1.79%	0.288%
9	11.971	1677	1686	1695	rVV2	56208511	94292451	100.00%	16.107%
10	12.076	1695	1704	1715	rVB2	46369717	85625203	90.81%	14.627%
11	12.400	1748	1759	1775	rBV	24216617	39139377	41.51%	6.686%
12	12.700	1799	1810	1817	rBV	3023700	4928837	5.23%	0.842%
13	12.853	1827	1836	1845	rVB	706485	1239453	1.31%	0.212%
14	13.041	1860	1868	1873	rBV	8498754	13775454	14.61%	2.353%
15	13.100	1873	1878	1881	rVV	12929762	21285833	22.57%	3.636%
16	13.135	1881	1884	1888	rVV	12721941	19422898	20.60%	3.318%
17	13.176	1888	1891	1899	rVB	10324346	14905569	15.81%	2.546%
18	13.341	1911	1919	1930	rBV	17582062	27697279	29.37%	4.731%
19	13.488	1935	1944	1957	rVV2	55349711	88248033	93.59%	15.075%
20	13.823	1990	2001	2017	rVB	20688500	33802340	35.85%	5.774%
21	13.953	2017	2023	2025	rBV	2089217	3086536	3.27%	0.527%
22	14.000	2025	2031	2046	rVV2	15680211	33315794	35.33%	5.691%
23	14.182	2056	2062	2068	rVV2	1718890	2959539	3.14%	0.506%
24	14.247	2068	2073	2076	rVV	2817260	4749898	5.04%	0.811%
25	14.276	2076	2078	2083	rVV	2081982	2716991	2.88%	0.464%
26	14.335	2083	2088	2094	rVV	5502754	8559367	9.08%	1.462%
27	14.400	2094	2099	2102	rVV	828425	1345972	1.43%	0.230%
28	14.447	2102	2107	2116	rVB2	3068621	5271481	5.59%	0.900%
29	14.576	2123	2129	2135	rBV	1353117	2111482	2.24%	0.361%
30	14.659	2135	2143	2146	rBV	3311157	5596724	5.94%	0.956%
31	14.694	2146	2149	2155	rVB	3892107	5763957	6.11%	0.985%
32	14.929	2184	2189	2195	rBV	3168978	5145881	5.46%	0.879%
33	15.053	2201	2210	2217	rBV2	5791141	12732525	13.50%	2.175%
34	15.117	2217	2221	2229	rVB2	633564	1233054	1.31%	0.211%
35	15.212	2230	2237	2245	rBV3	1175418	2184754	2.32%	0.373%
36	15.323	2245	2256	2261	rBV3	783706	2074071	2.20%	0.354%

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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_N\methods\82N121422W.M
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37	15.441	2268	2276	2285	rVB3	633539	1627098	1.73%	0.278%
38	15.647	2302	2311	2319	rBV	13683707	26004490	27.58%	4.442%

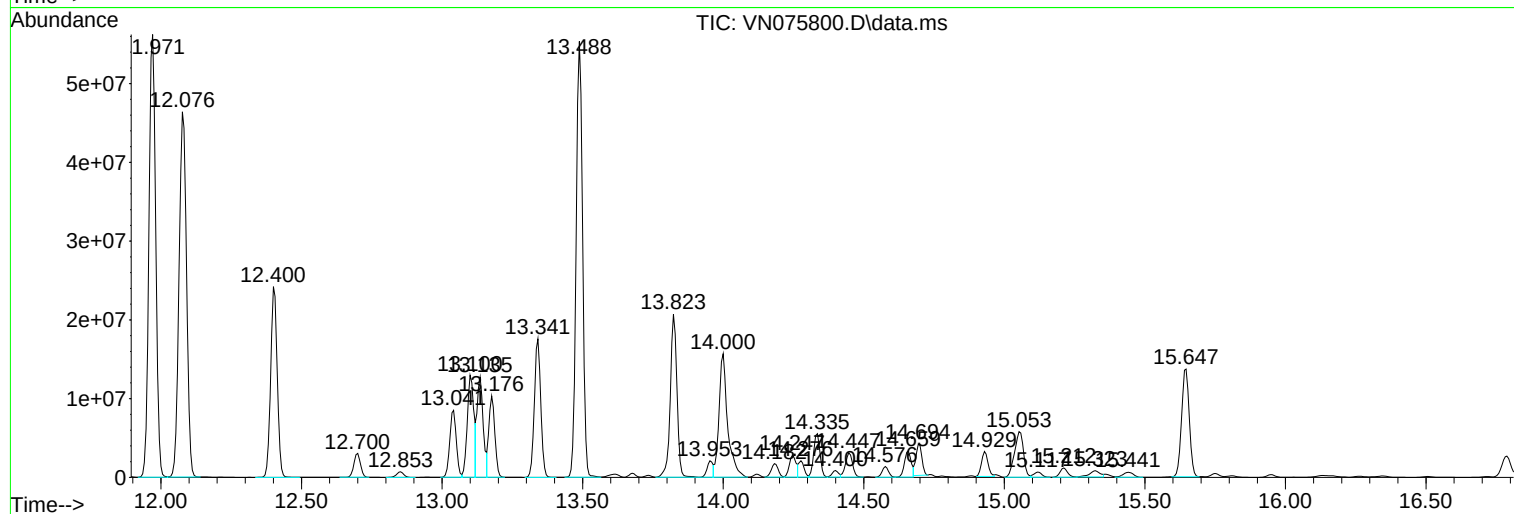
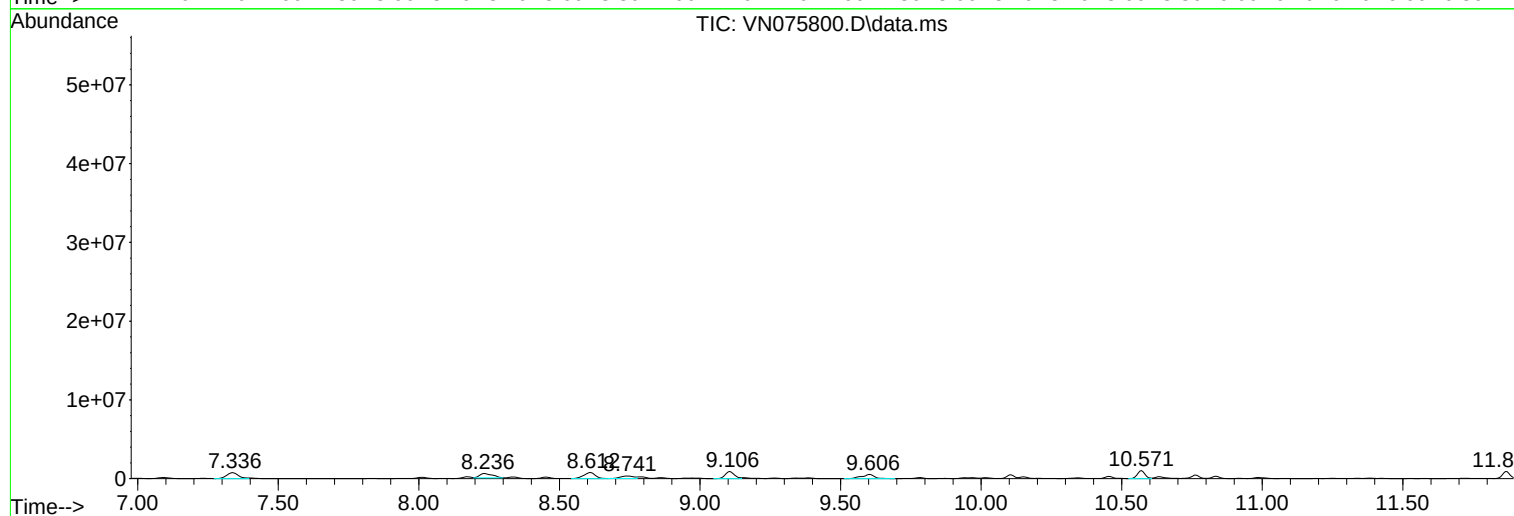
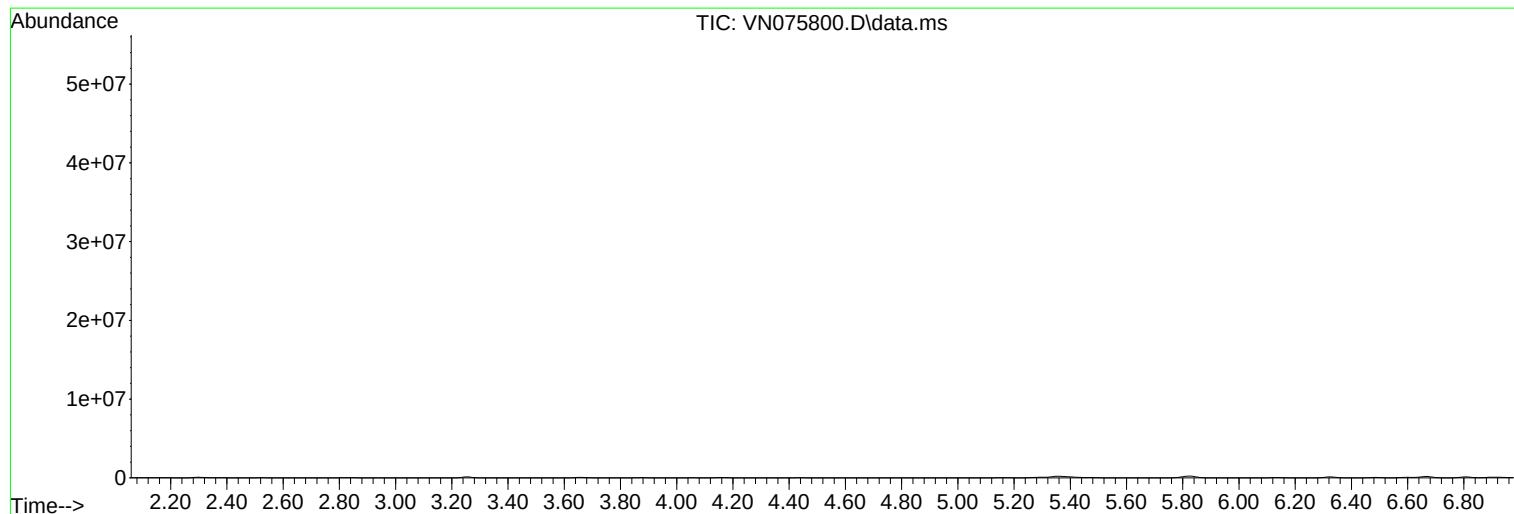
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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



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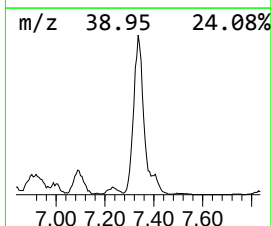
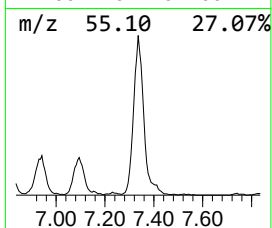
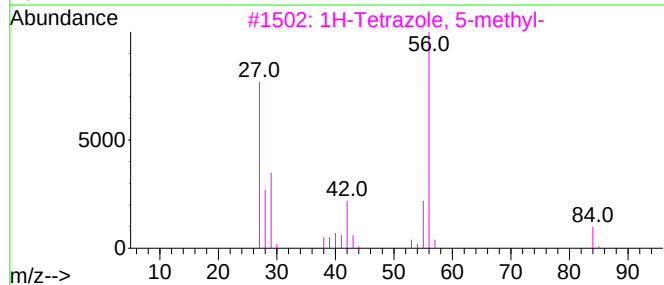
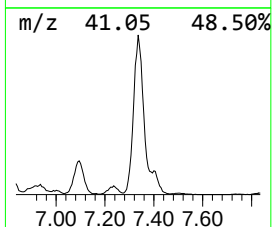
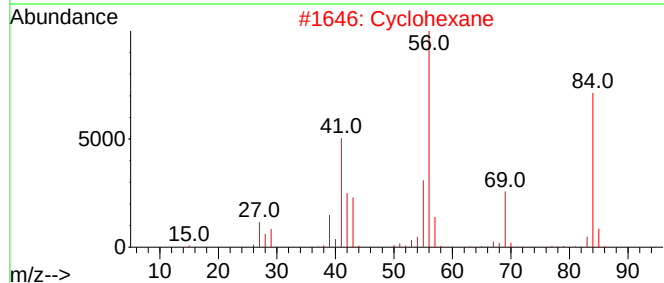
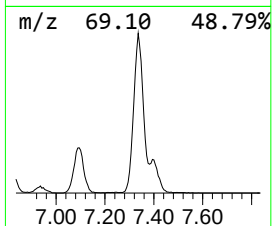
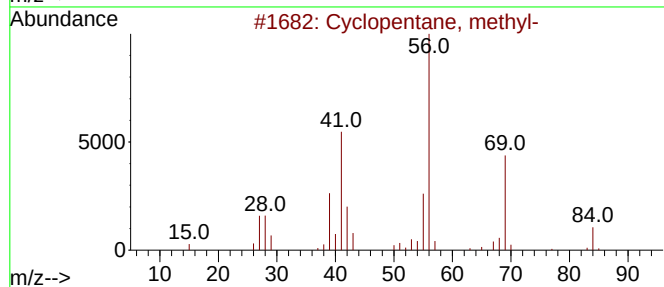
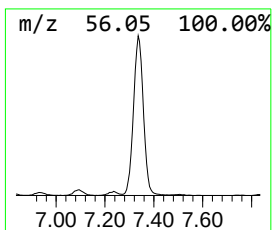
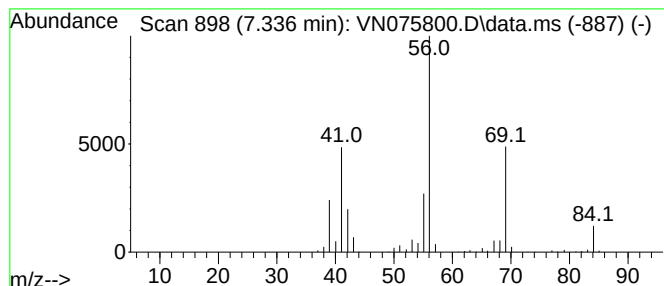
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.336	60.82 ug/l	2180270	Pentafluorobenzene	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclohexane	84	C6H12	000110-82-7	78
3			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72



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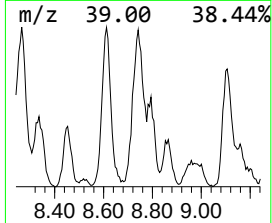
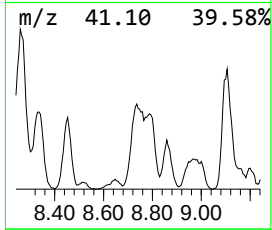
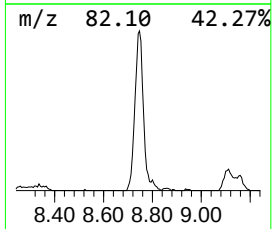
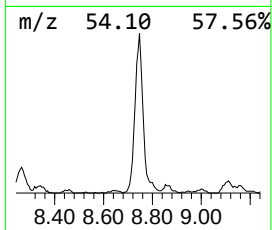
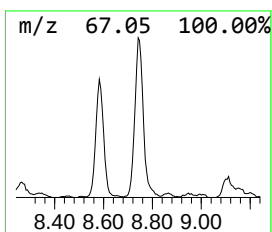
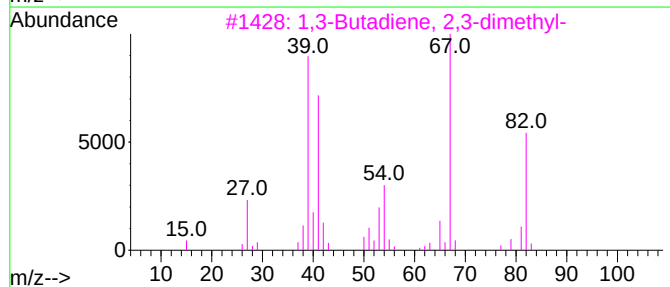
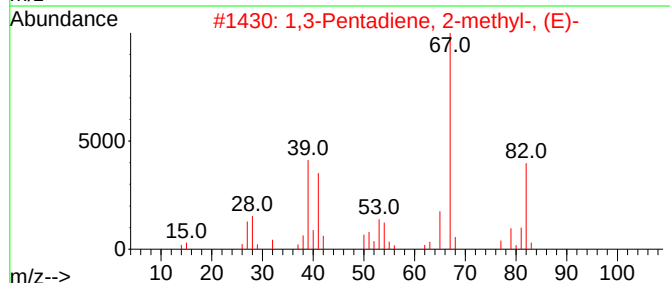
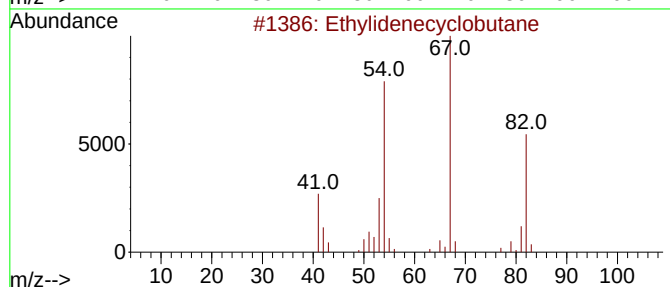
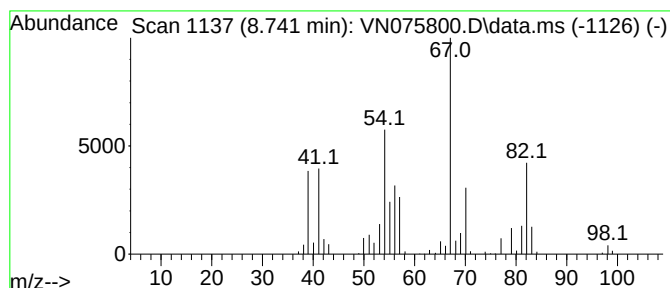
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Ethylidenecyclobutane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.741	26.33 ug/l	1091840	1,4-Difluorobenzene	9.106

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethylidenecyclobutane	82	C6H10	001528-21-8	76
2		1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	64
3		1,3-Butadiene, 2,3-dimethyl-	82	C6H10	000513-81-5	62
4		Cyclobutane, ethenyl-	82	C6H10	002597-49-1	60
5		Cyclohexene	82	C6H10	000110-83-8	60



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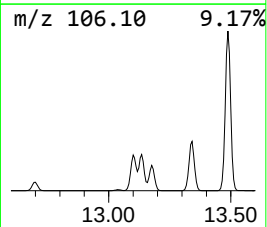
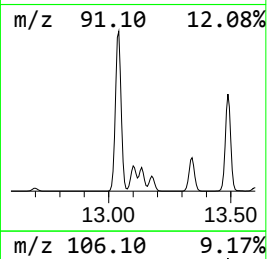
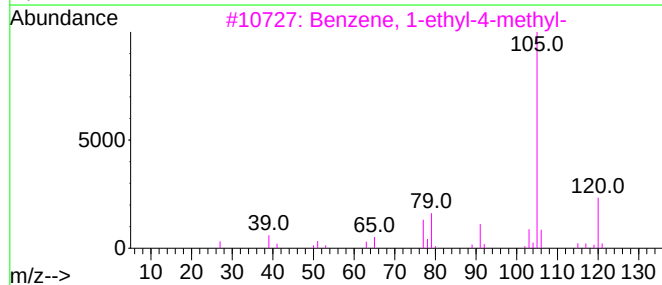
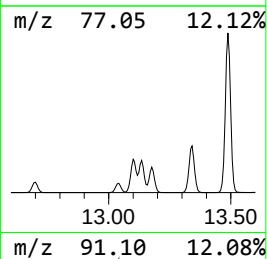
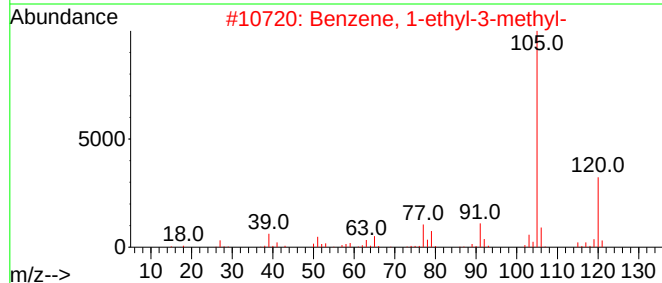
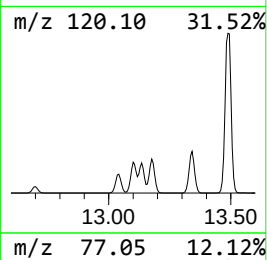
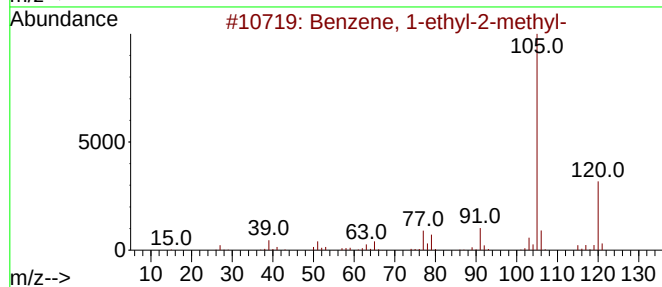
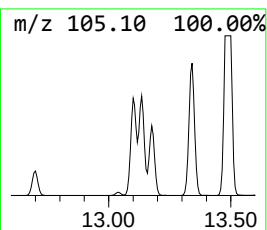
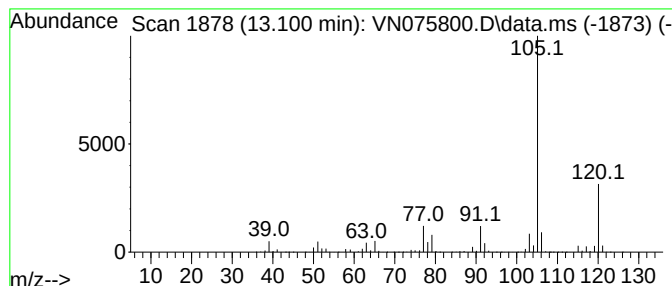
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	31.49 ug/l	21285800	1,4-Dichlorobenzene-d4	13.794

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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ClientSampleId :
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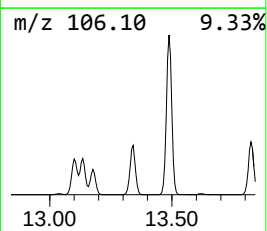
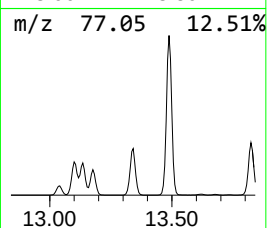
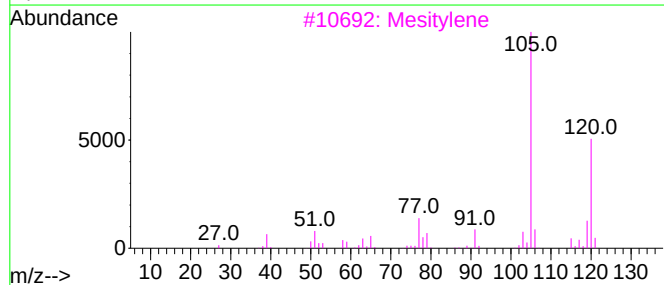
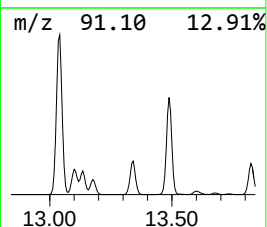
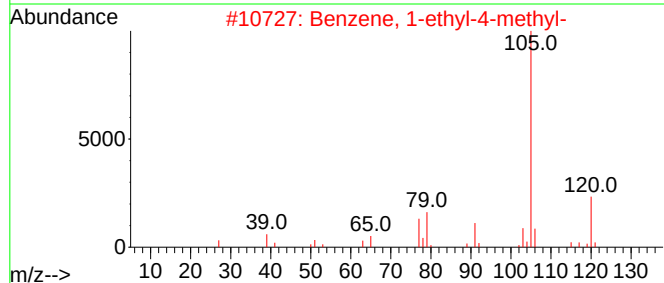
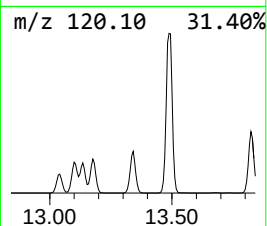
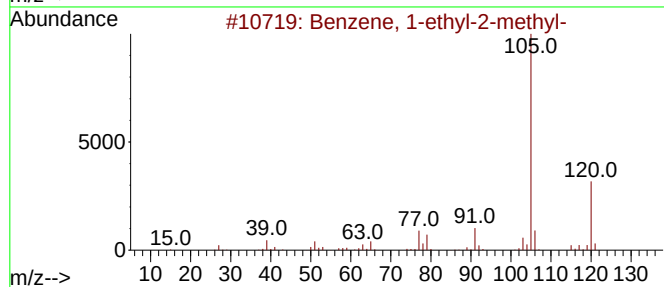
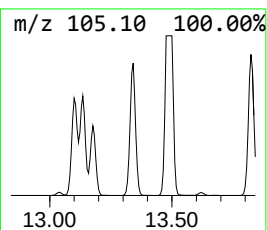
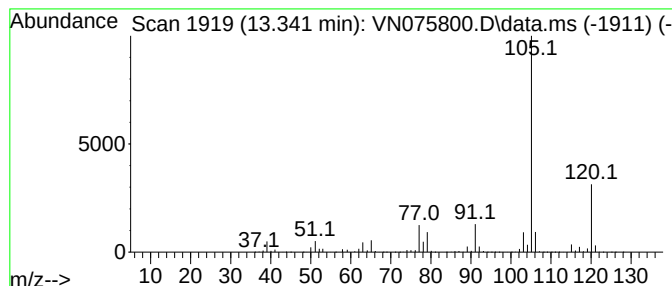
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.341	40.97 ug/l	27697300	1,4-Dichlorobenzene-d4	13.794

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	93
3			Mesitylene	120	C9H12	000108-67-8	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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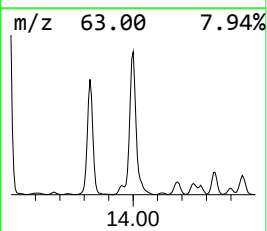
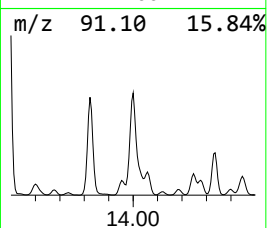
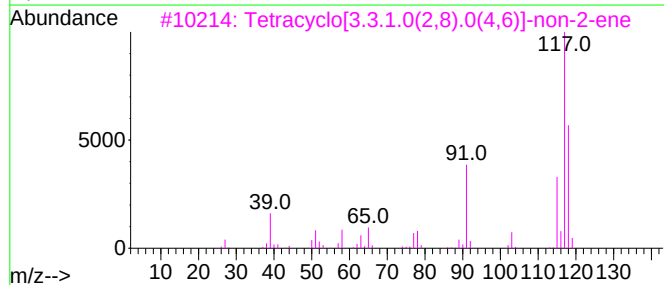
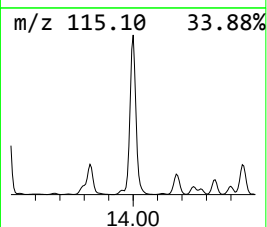
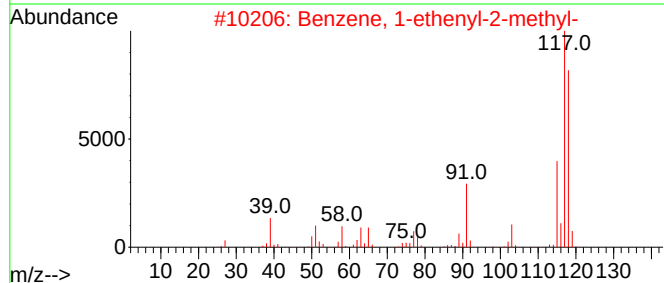
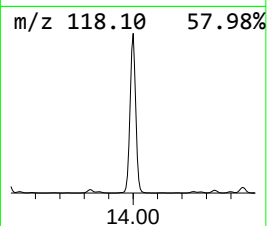
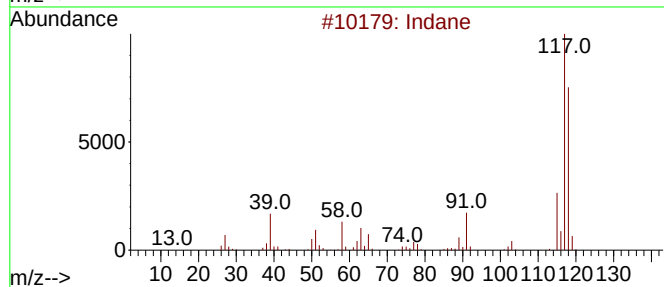
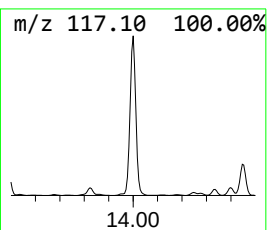
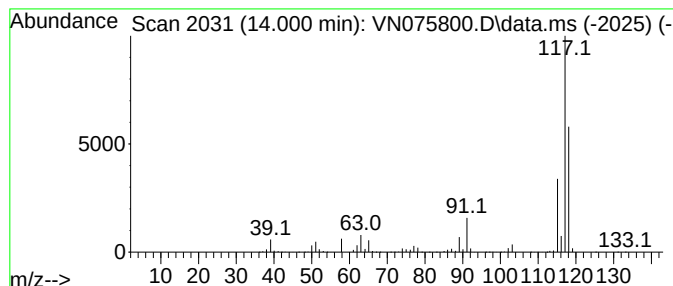
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.000	49.28 ug/l	33315800	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

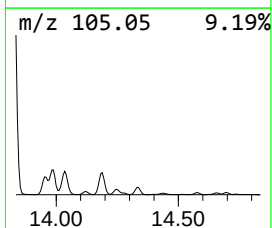
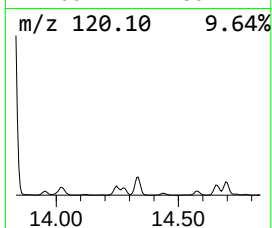
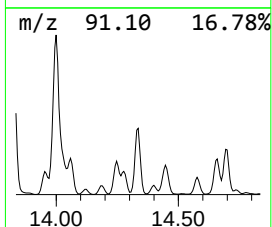
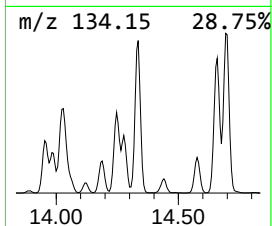
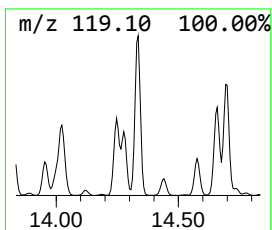
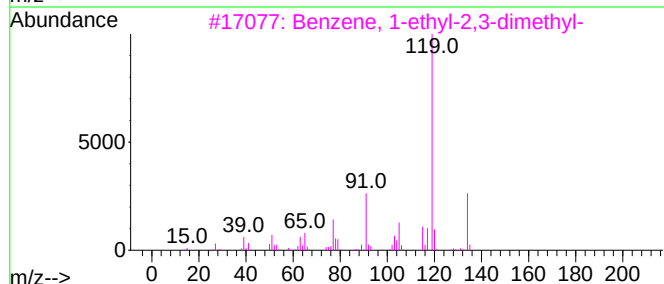
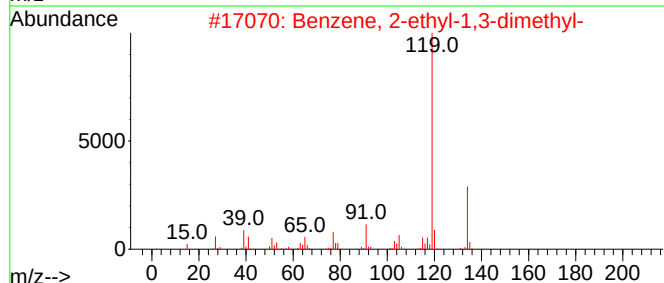
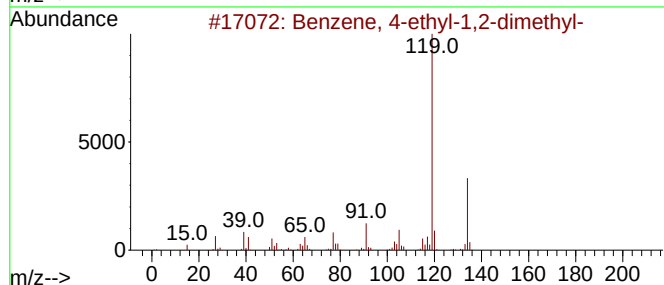
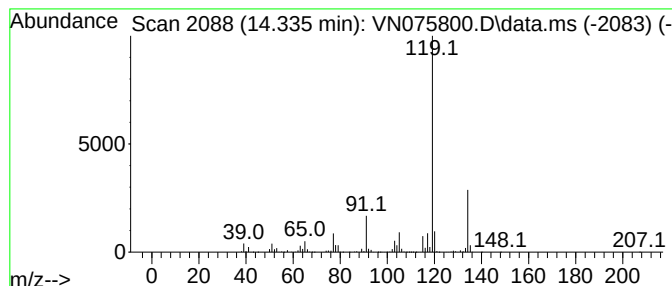
Instrument :
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ClientSampleId :
MW-43

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.335	12.66 ug/l	8559370	1,4-Dichlorobenzene-d4		13.794	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-		134	C10H14	000934-80-5	95
2	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	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_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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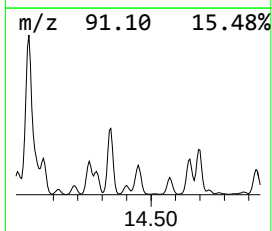
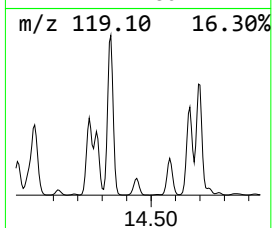
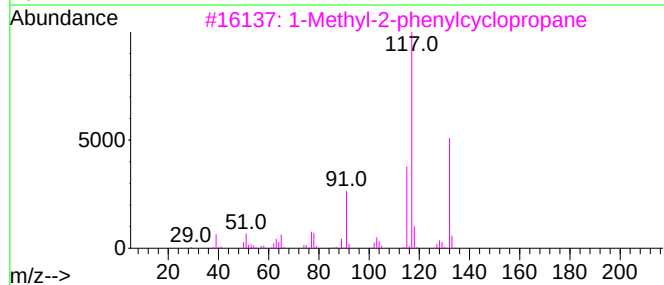
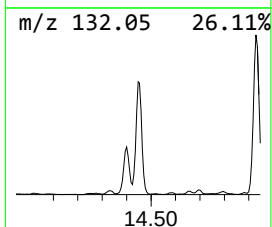
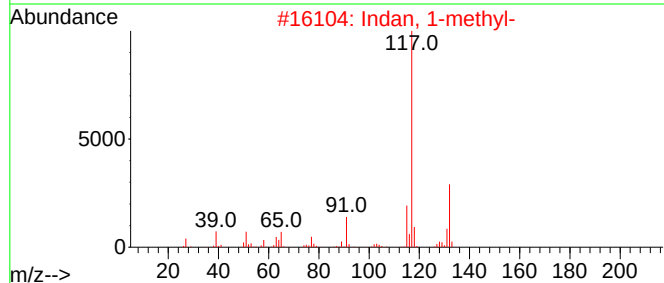
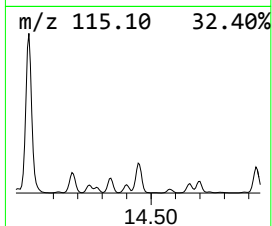
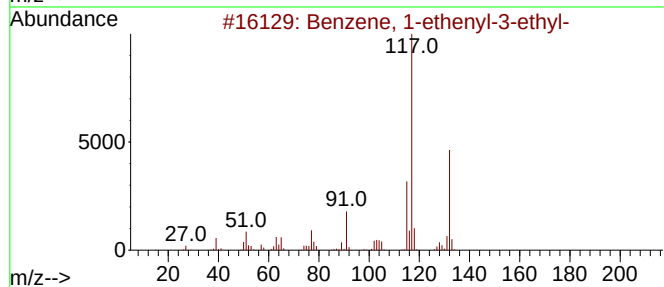
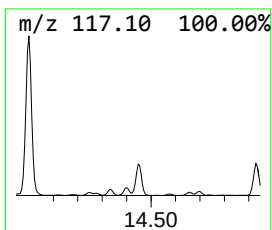
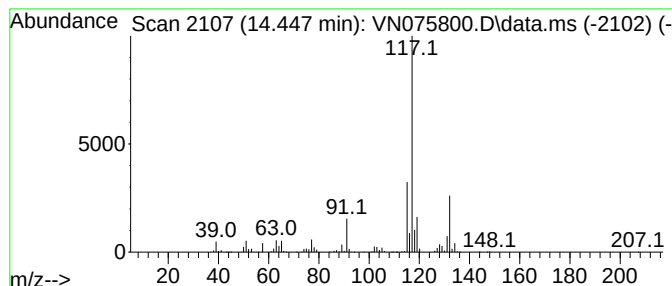
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	7.80 ug/l	5271480	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	83
2			Indan, 1-methyl-	132	C10H12	000767-58-8	81
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	74
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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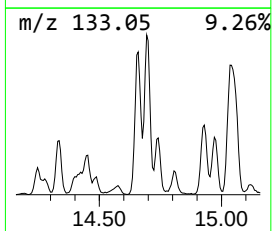
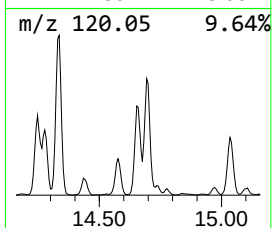
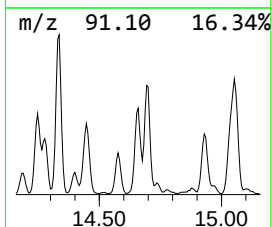
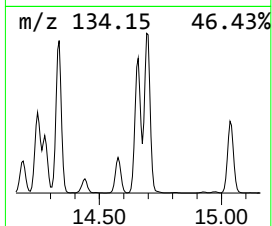
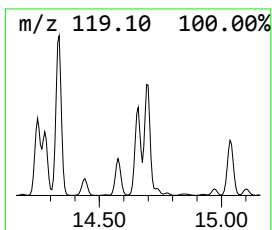
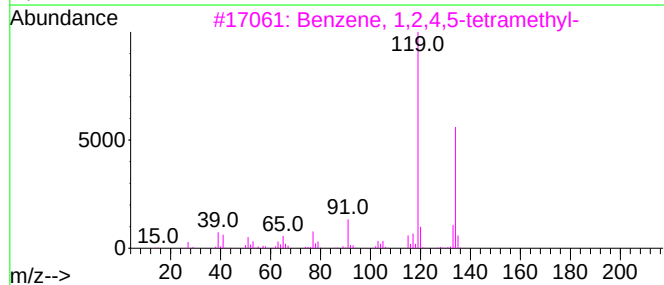
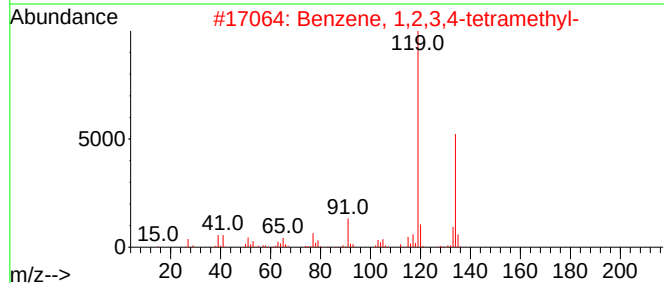
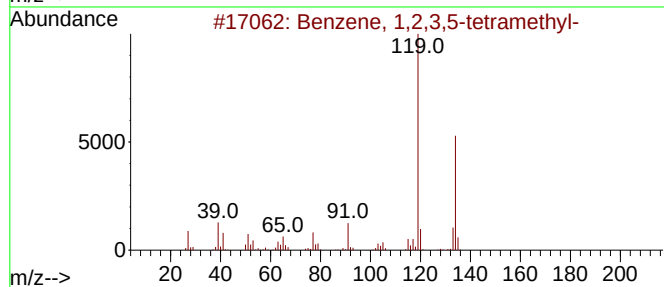
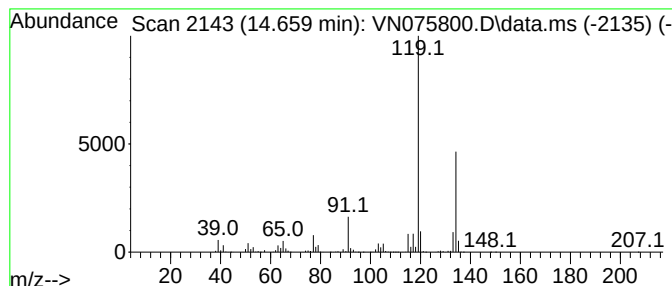
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.659	8.28 ug/l	5596720	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
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_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-43

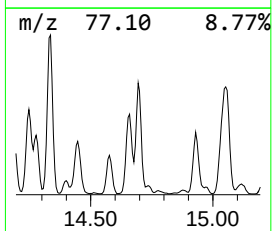
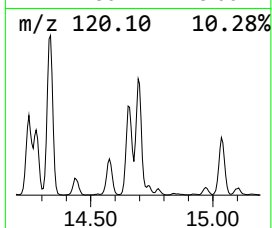
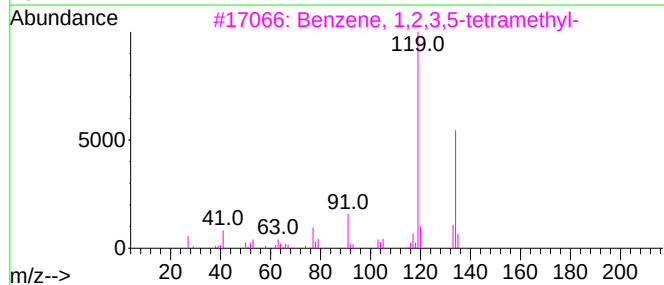
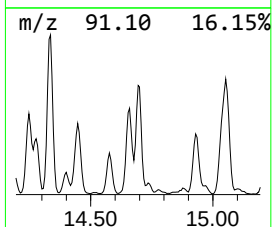
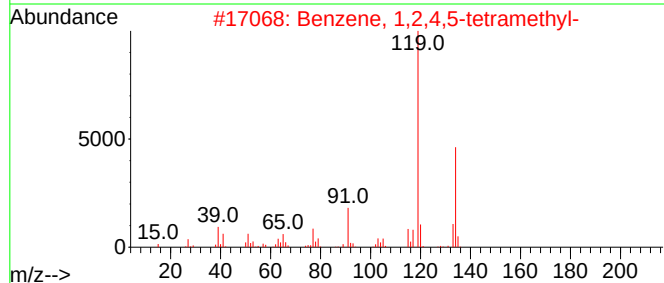
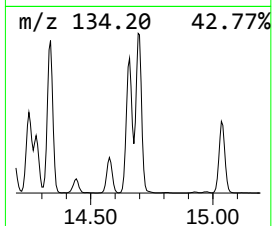
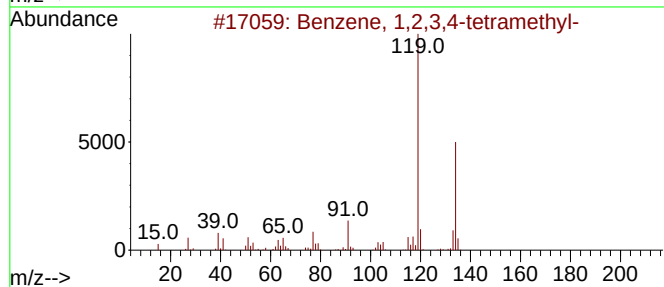
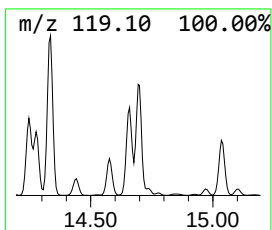
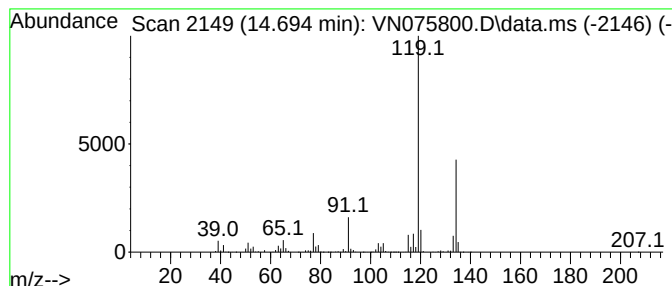
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.694	8.53 ug/l	5763960	1,4-Dichlorobenzene-d4	13.794

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, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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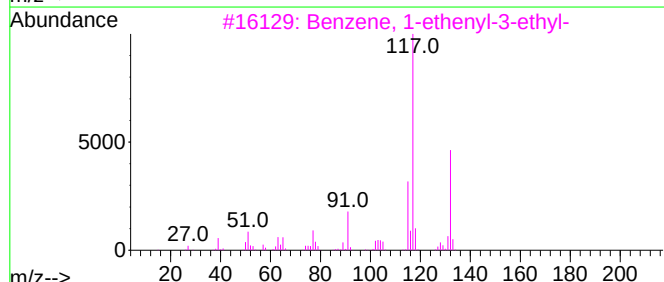
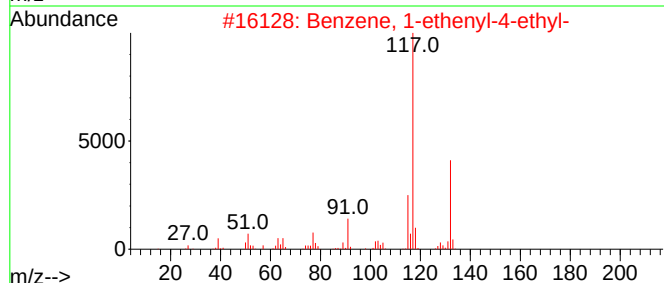
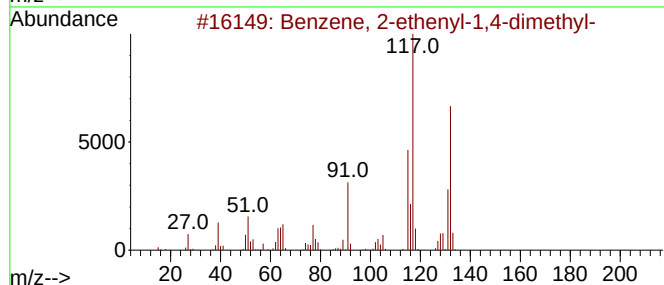
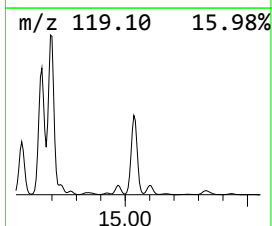
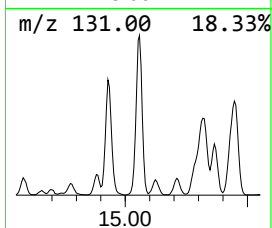
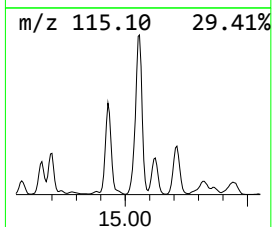
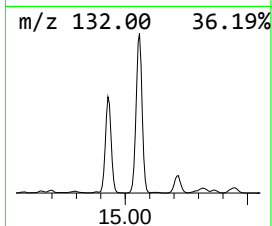
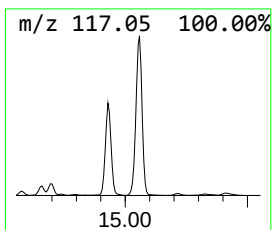
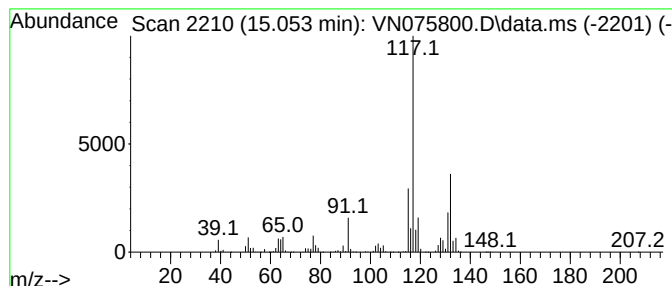
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	18.83 ug/l	12732500	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
4			Indan, 1-methyl-	132	C10H12	000767-58-8	90
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075800.D
Acq On : 15 Dec 2022 05:15
Operator : JC\MD
Sample : N5990-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Ethylidenecyclo...	8.741	26.3	ug/l	1091840	2	9.106	2073020	50.0
Benzene, 1-ethy...	13.100	31.5	ug/l	21285800	4	13.794	33802300	50.0
Benzene, 1-ethy...	13.341	41.0	ug/l	27697300	4	13.794	33802300	50.0
Indane	14.000	49.3	ug/l	33315800	4	13.794	33802300	50.0
Benzene, 4-ethy...	14.335	12.7	ug/l	8559370	4	13.794	33802300	50.0
Benzene, 1-ethe...	14.447	7.8	ug/l	5271480	4	13.794	33802300	50.0
Benzene, 1,2,3,...	14.659	8.3	ug/l	5596720	4	13.794	33802300	50.0
Benzene, 1,2,3,...	14.694	8.5	ug/l	5763960	4	13.794	33802300	50.0
Benzene, 2-ethe...	15.053	18.8	ug/l	12732500	4	13.794	33802300	50.0