

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120921\
 Data File : VN069914.D
 Acq On : 9 Dec 2021 16:21
 Operator : JC/MD
 Sample : M4940-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-24

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\82N120221W.M
 Title : SW846 8260

Signal : TIC: VN069914.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.234	80	91	103	rBV	220952	520355	6.12%	0.968%
2	2.446	162	170	182	rVB	123961	187557	2.21%	0.349%
3	3.143	414	430	454	rVB	502120	1141620	13.43%	2.123%
4	4.256	825	845	881	rVB5	41912	178452	2.10%	0.332%
5	5.082	1120	1153	1171	rBV4	143992	575835	6.77%	1.071%
6	5.374	1237	1262	1292	rBV3	136497	510891	6.01%	0.950%
7	5.618	1328	1353	1382	rVB4	139350	480034	5.65%	0.893%
8	6.498	1660	1681	1704	rVB2	41286	128804	1.51%	0.240%
9	7.147	1900	1923	1943	rBV3	211488	627926	7.39%	1.168%
10	8.021	2226	2249	2260	rBV	964477	2295715	27.00%	4.269%
11	8.086	2261	2273	2298	rVB	1724784	4026804	47.36%	7.489%
12	8.437	2384	2404	2427	rBV	1099521	2562705	30.14%	4.766%
13	8.536	2428	2441	2460	rVV3	130009	310145	3.65%	0.577%
14	8.968	2582	2602	2642	rBV	2701186	5742776	67.55%	10.680%
15	9.971	2960	2976	2990	rBV2	146796	290102	3.41%	0.540%
16	10.202	3042	3062	3067	rBV2	94626	175659	2.07%	0.327%
17	10.441	3132	3151	3181	rBV	4420996	8237657	96.89%	15.320%
18	10.550	3182	3192	3213	rVV2	127281	291350	3.43%	0.542%
19	10.642	3213	3226	3240	rVV3	85061	161693	1.90%	0.301%
20	10.717	3241	3254	3271	rVB2	67090	122855	1.45%	0.228%
21	11.130	3399	3408	3422	rVB	88521	157497	1.85%	0.293%
22	11.623	3575	3592	3618	rBV7	51563	157922	1.86%	0.294%
23	11.744	3618	3637	3671	rBV	4684846	8501915	100.00%	15.811%
24	12.731	3988	4005	4037	rBV	3577782	6398713	75.26%	11.900%
25	13.675	4340	4357	4383	rBV	3899290	6886510	81.00%	12.807%
26	13.876	4411	4432	4463	rVB	1693881	2967851	34.91%	5.519%
27	14.329	4590	4601	4616	rVB2	76238	132307	1.56%	0.246%

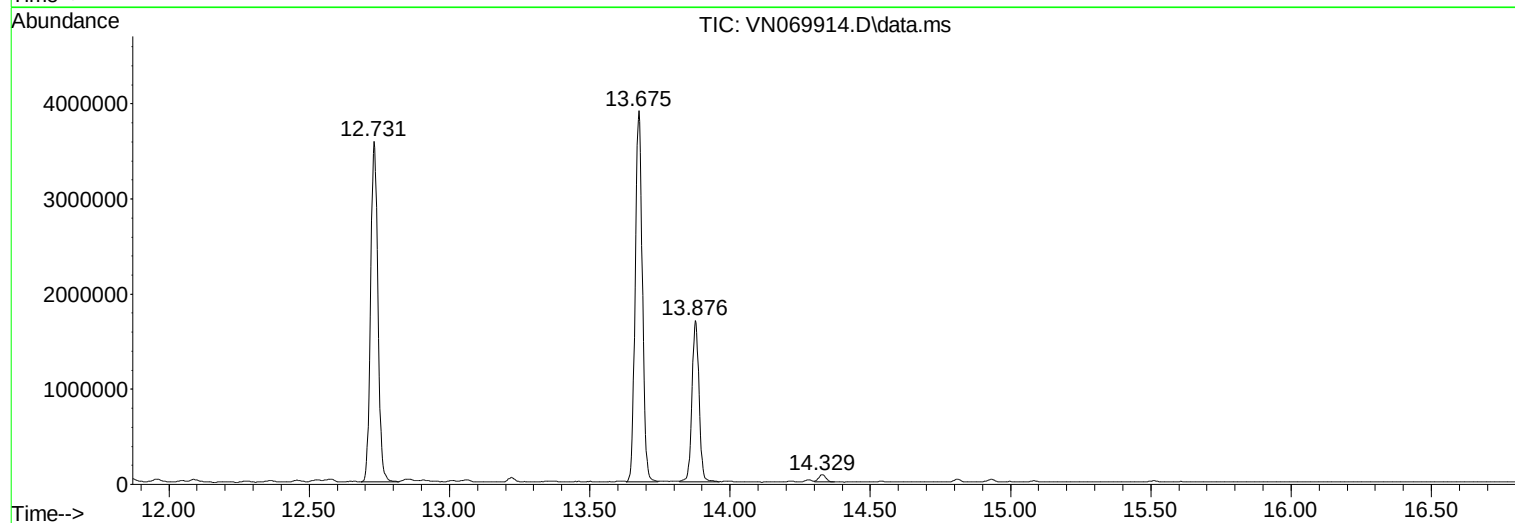
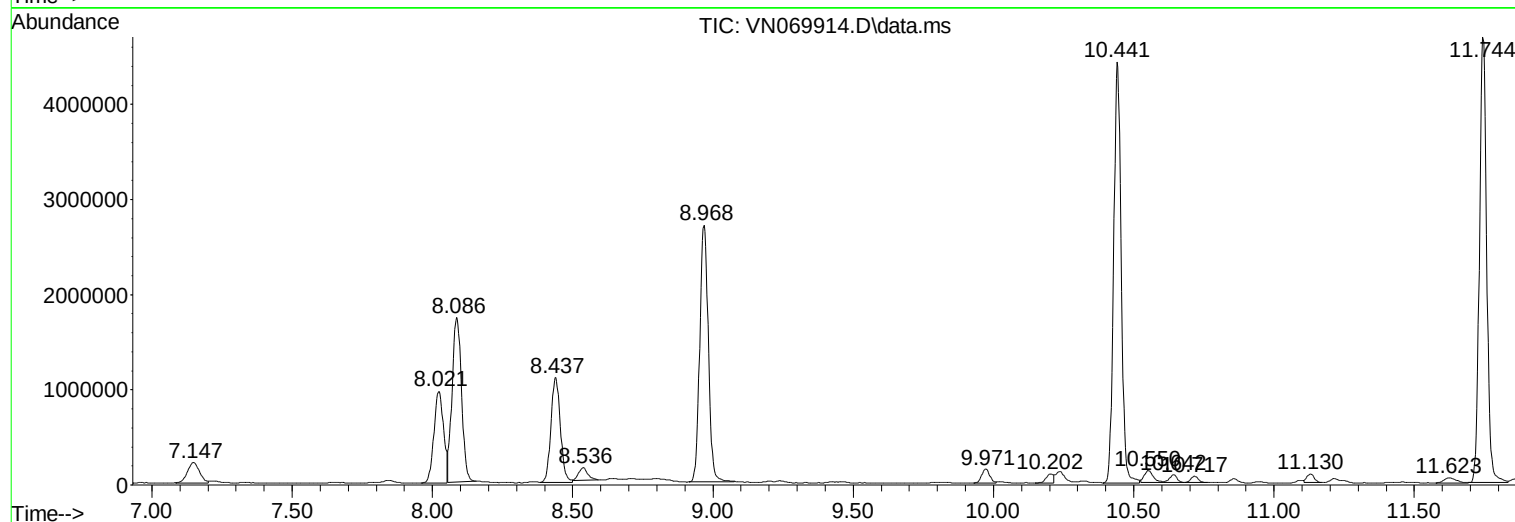
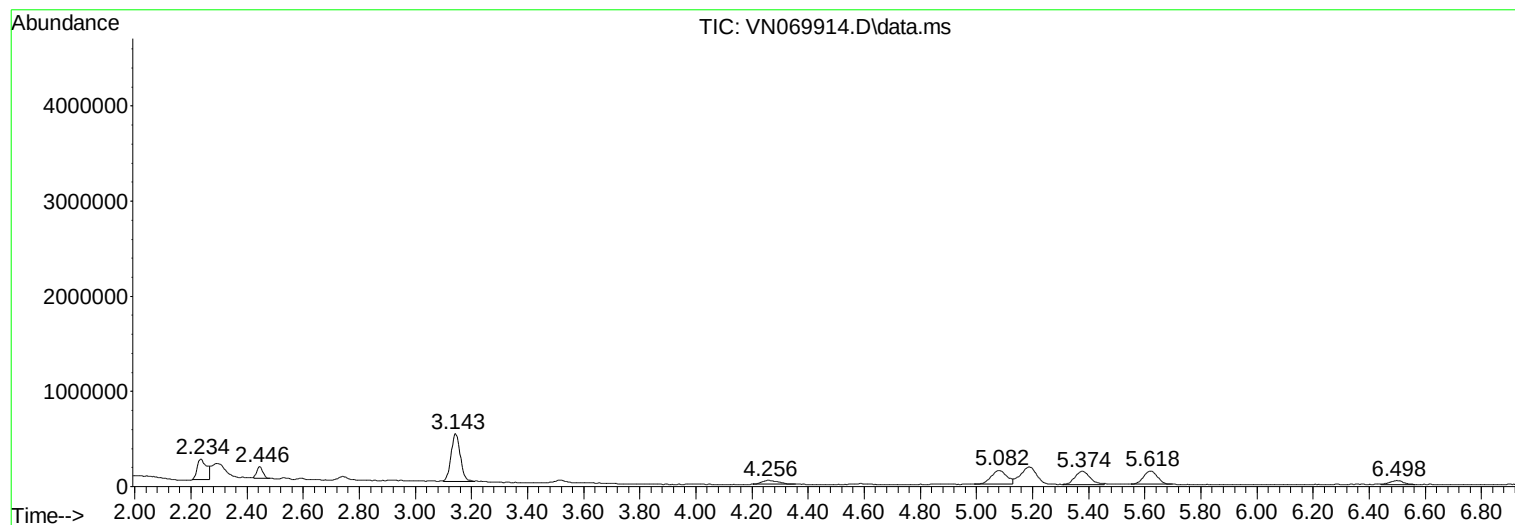
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Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

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



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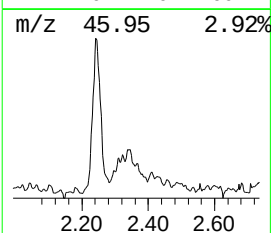
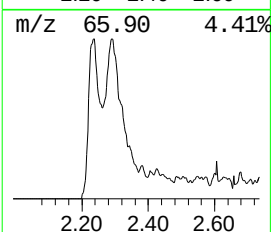
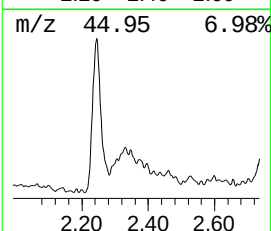
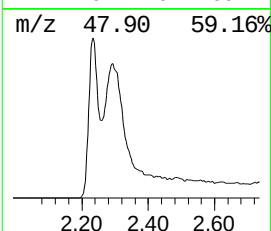
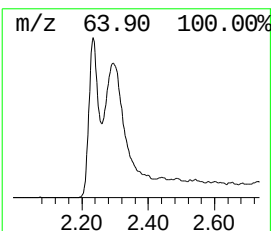
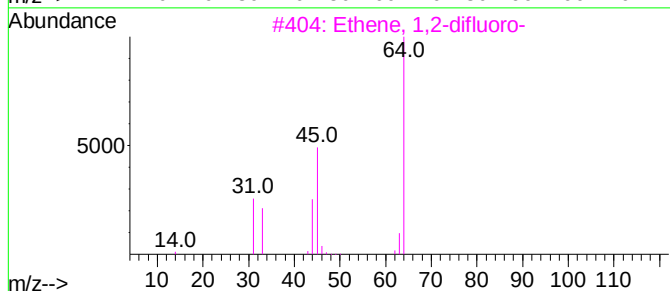
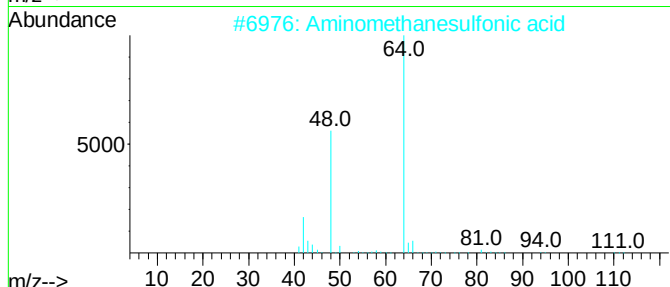
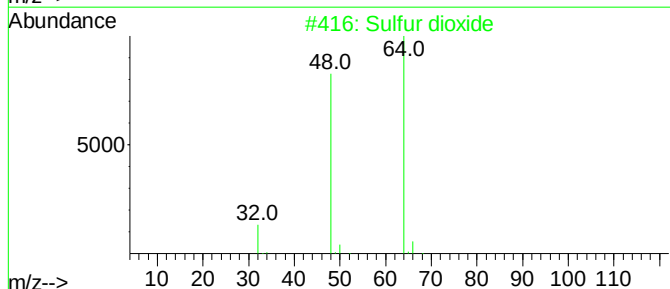
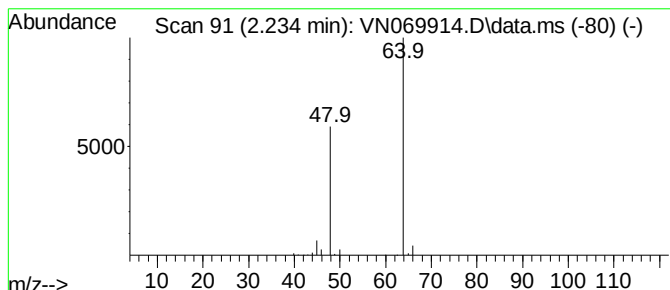
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Quant Title : SW846 8260

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

Peak Number 1 Sulfur dioxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.234	6.46 ug/l	520355	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
3	Ethene, 1,2-difluoro-	64	C2H2F2	001691-13-0	5
4	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	3



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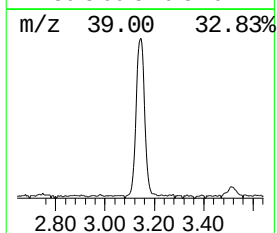
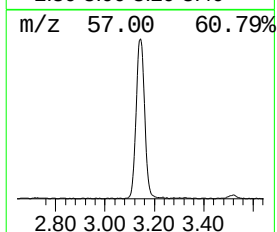
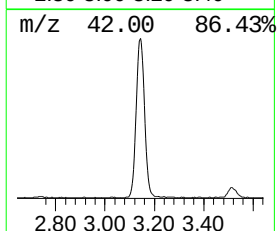
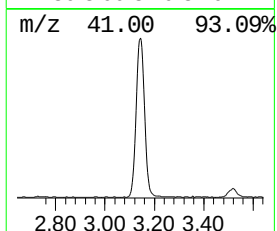
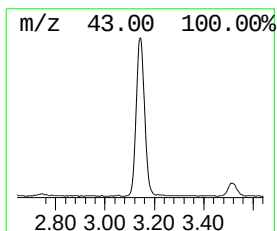
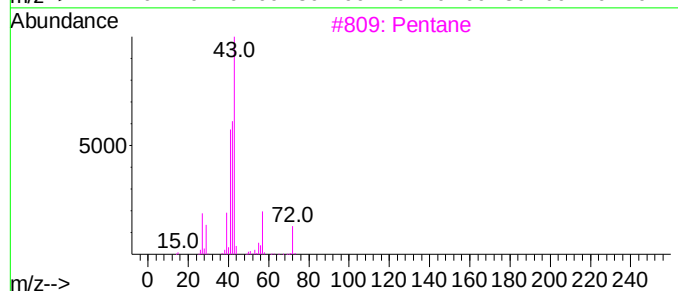
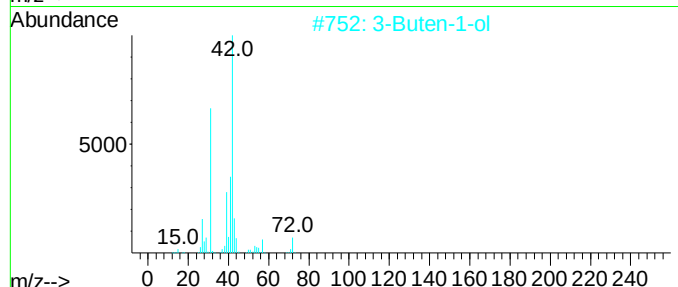
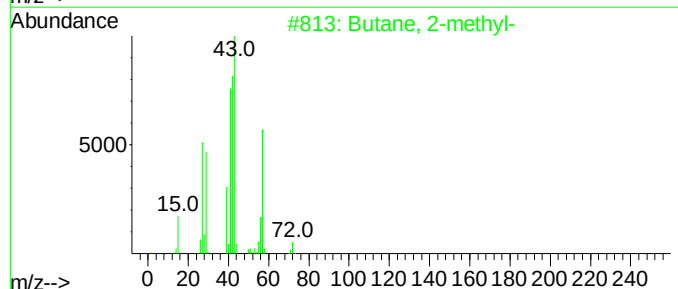
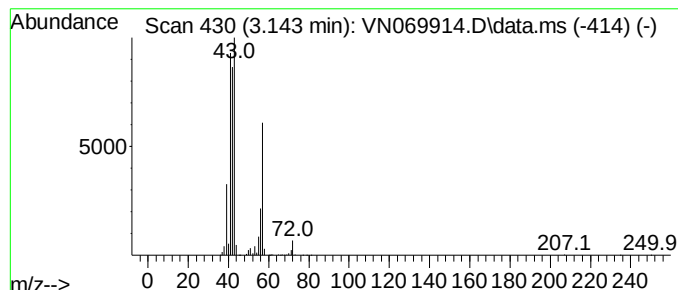
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Quant Title : SW846 8260

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

Peak Number 2 Butane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.143	14.18 ug/l	1141620	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	87
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	42
4	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	38
5	2,5-Furandione, dihydro-3-methyl-	114	C5H6O3	004100-80-5	23



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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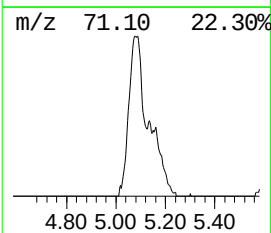
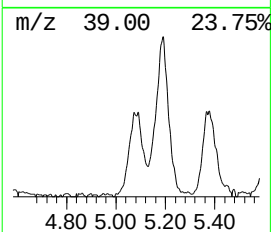
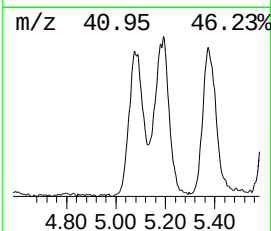
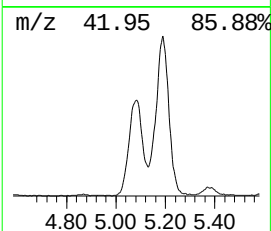
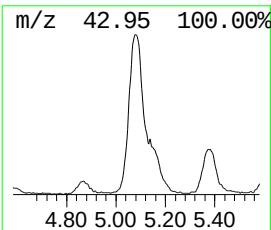
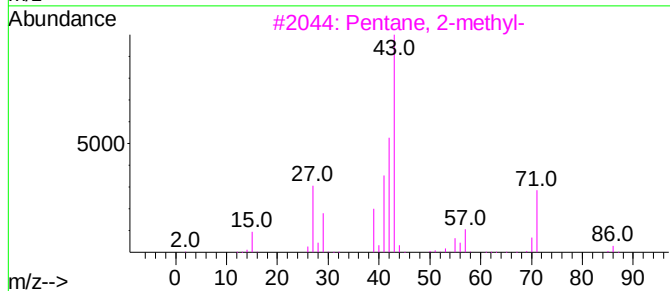
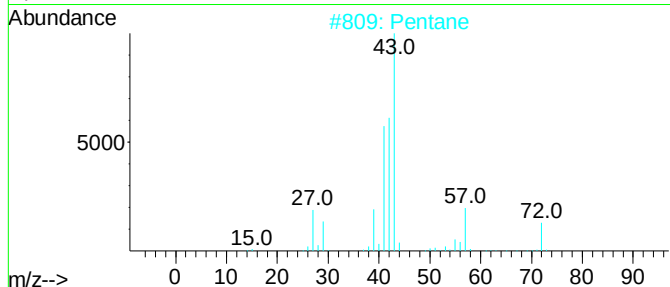
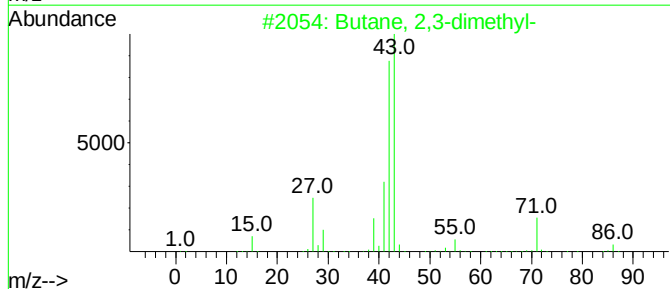
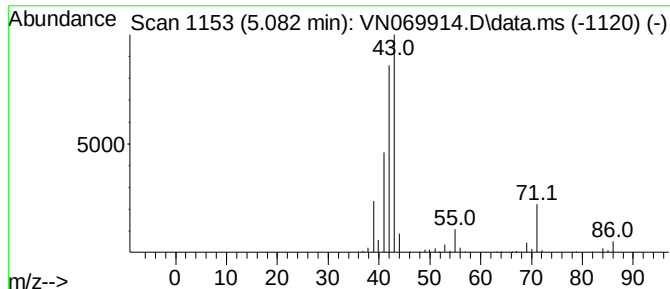
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.082	7.15 ug/l	575835	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane	72	C5H12	000109-66-0	36
3	Pentane, 2-methyl-	86	C6H14	000107-83-5	17
4	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9
5	1-Hexene, 4,5-dimethyl-	112	C8H16	016106-59-5	9



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Sample : M4940-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 3

Instrument :
MSVOA_N
ClientSampleId :
MW-24

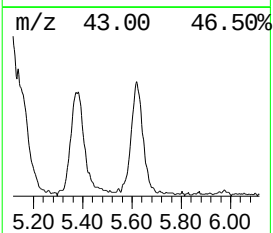
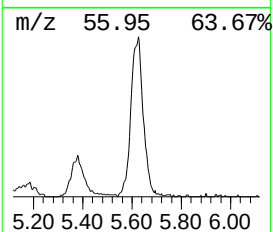
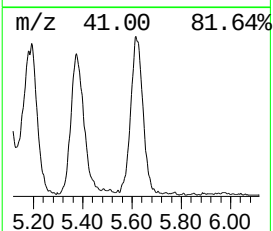
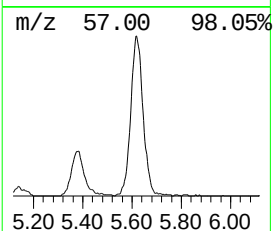
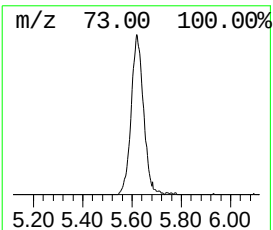
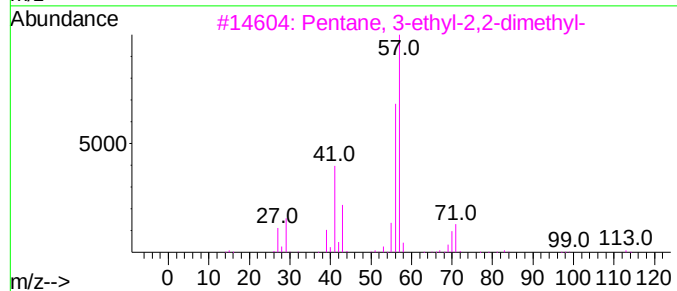
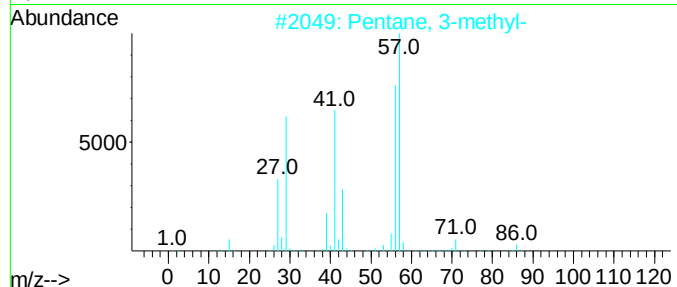
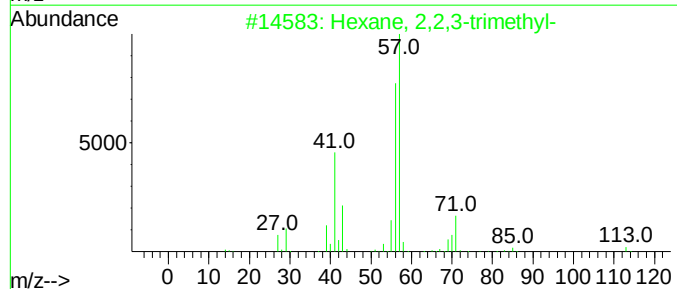
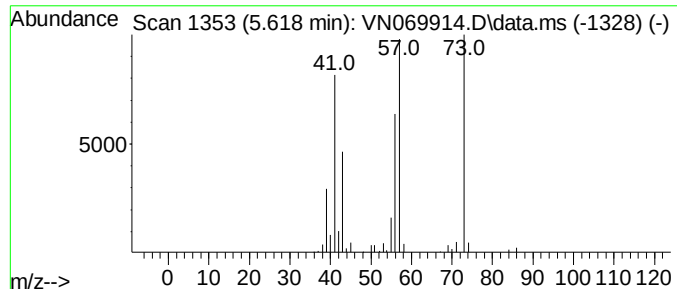
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

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

Peak Number 4 Hexane, 2,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.618	5.96 ug/l	480034	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	59
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	50
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	42
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	38
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	38



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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

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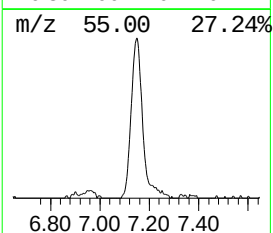
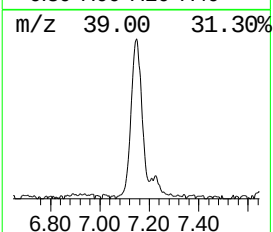
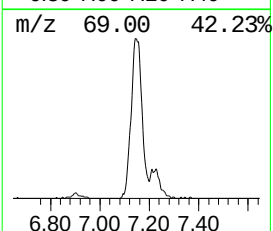
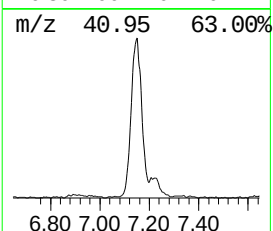
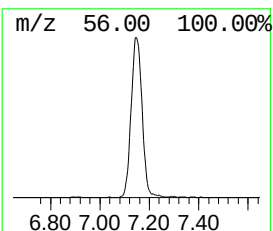
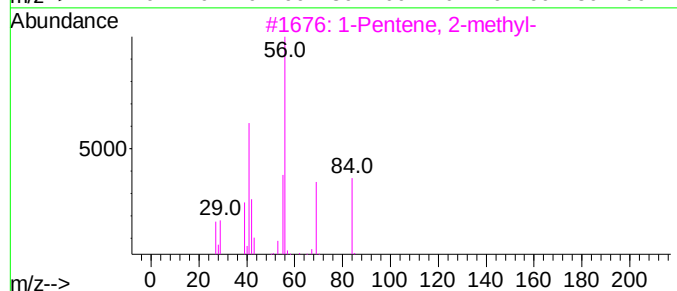
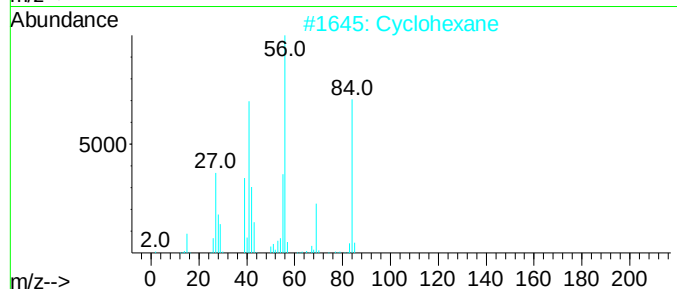
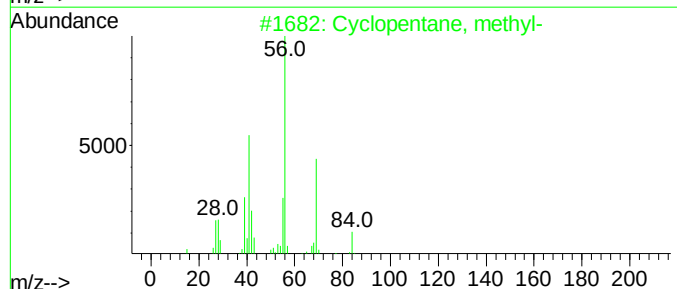
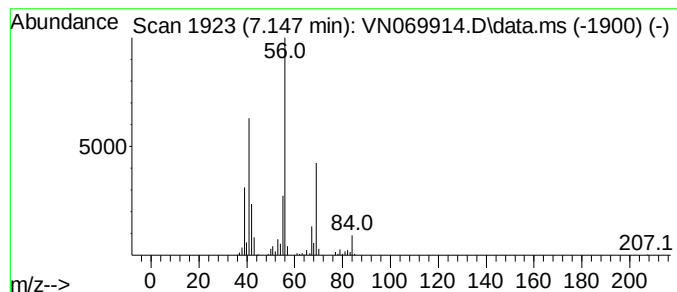
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.147	7.80 ug/l	627926	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclohexane	84	C6H12	000110-82-7	80
3	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59



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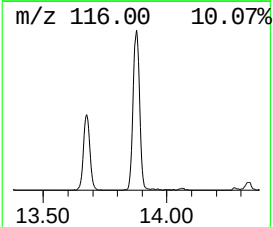
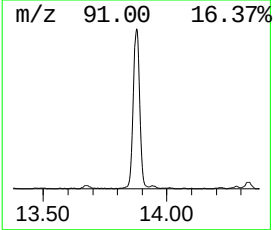
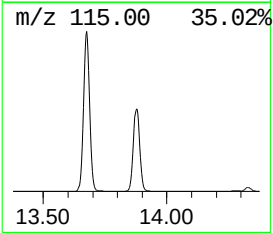
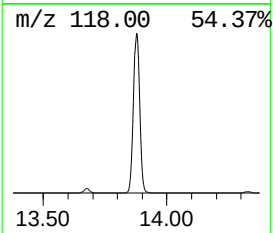
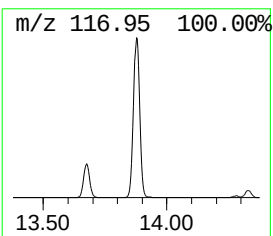
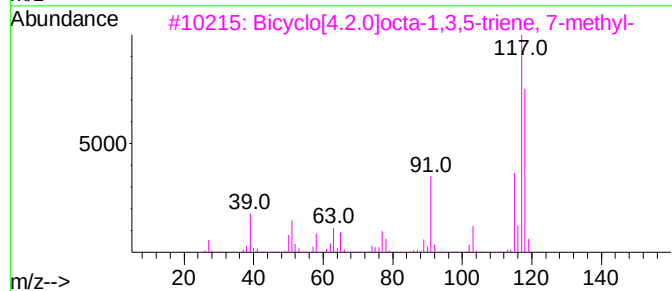
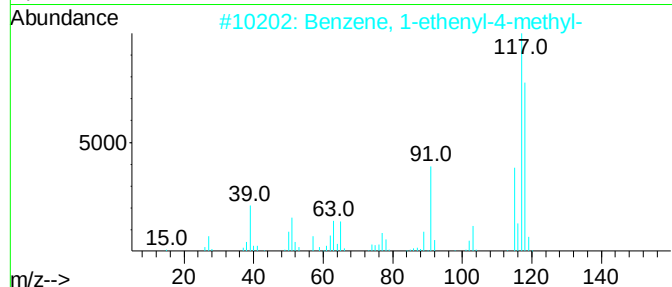
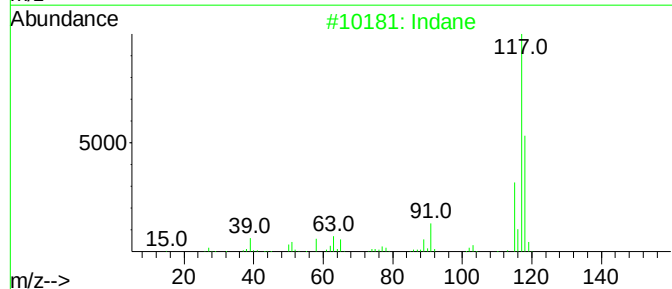
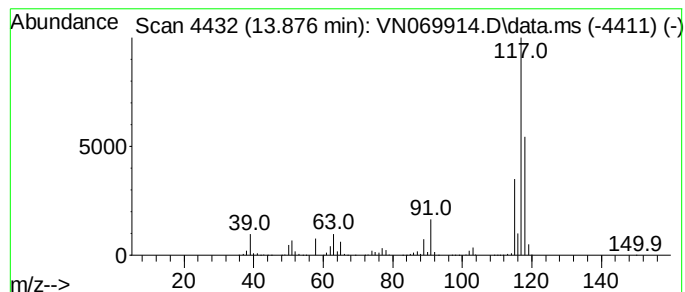
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.876	21.55 ug/l	2967850	1,4-Dichlorobenzene-d4	13.675

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95	
2	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83	
3	Bicyclo[4.2.0]octa-1,3,5-triene, ...	118	C9H10	055337-80-9	74	
4	Deltacyclene	118	C9H10	007785-10-6	68	
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	68	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120921\
Data File : VN069914.D
Acq On : 9 Dec 2021 16:21
Operator : JC/MD
Sample : M4940-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-24

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.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
Sulfur dioxide	2.234	6.5	ug/l	520355	1	8.086	4026800	50.0
Butane, 2-methyl-	3.143	14.2	ug/l	1141620	1	8.086	4026800	50.0
Butane, 2,3-dim...	5.082	7.2	ug/l	575835	1	8.086	4026800	50.0
Hexane, 2,2,3-t...	5.618	6.0	ug/l	480034	1	8.086	4026800	50.0
Cyclopentane, m...	7.147	7.8	ug/l	627926	1	8.086	4026800	50.0
Indane	13.876	21.6	ug/l	2967850	4	13.675	6886510	50.0