

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
 Data File : VN072609.D
 Acq On : 25 May 2022 00:32
 Operator : JC\MD
 Sample : N2968-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-38

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\82N052422W.M
 Title : SW846 8260

Signal : TIC: VN072609.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.434	65	76	86	rBV	226561	383440	3.01%	0.413%
2	3.116	182	192	213	rVB	259218	615143	4.83%	0.662%
3	3.510	249	259	274	rVB4	105242	344213	2.70%	0.371%
4	3.710	282	293	307	rBV	129789	321302	2.52%	0.346%
5	3.887	310	323	331	rVV	74454	199614	1.57%	0.215%
6	3.987	331	340	356	rVB2	158716	435320	3.41%	0.469%
7	4.957	490	505	516	rBV2	92274	324467	2.55%	0.349%
8	5.181	530	543	554	rVV3	136529	528740	4.15%	0.569%
9	5.351	563	572	594	rVB2	115159	406515	3.19%	0.438%
10	5.610	602	616	635	rBV4	101554	366126	2.87%	0.394%
11	6.463	755	761	774	rVV3	75663	248390	1.95%	0.267%
12	6.598	774	784	792	rVV2	46554	143706	1.13%	0.155%
13	6.692	792	800	811	rVV3	52351	216080	1.69%	0.233%
14	6.904	826	836	848	rVB3	48526	143975	1.13%	0.155%
15	7.145	864	877	886	rBV	274604	818409	6.42%	0.881%
16	8.016	1015	1025	1030	rBV	97793	239017	1.87%	0.257%
17	8.086	1030	1037	1046	rVV2	352033	931683	7.31%	1.003%
18	8.457	1088	1100	1110	rBV	4938087	11203107	87.88%	12.061%
19	8.580	1116	1121	1129	rVB2	83258	177739	1.39%	0.191%
20	8.963	1175	1186	1199	rBV	334589	761618	5.97%	0.820%
21	9.457	1257	1270	1277	rBV3	141638	391570	3.07%	0.422%
22	9.969	1350	1357	1361	rBV	123700	237527	1.86%	0.256%
23	10.439	1428	1437	1442	rBV	472141	887334	6.96%	0.955%
24	10.498	1442	1447	1454	rVV	441511	764475	6.00%	0.823%
25	10.633	1461	1470	1477	rVV	77879	147464	1.16%	0.159%
26	11.739	1651	1658	1667	rBV	542068	951849	7.47%	1.025%
27	11.839	1667	1675	1685	rVV	7413533	12019803	94.28%	12.940%
28	11.951	1685	1694	1705	rVB	7053450	12748773	100.00%	13.725%
29	12.274	1741	1749	1761	rBV	4076470	6528191	51.21%	7.028%
30	12.574	1793	1800	1809	rVB	584993	957198	7.51%	1.030%
31	12.727	1819	1826	1835	rBV	607843	1021157	8.01%	1.099%
32	12.915	1848	1858	1863	rBV	1413829	2232373	17.51%	2.403%
33	12.980	1863	1869	1871	rVV	1484603	2476789	19.43%	2.666%
34	13.010	1871	1874	1878	rVV	1564237	2426064	19.03%	2.612%
35	13.051	1878	1881	1889	rVB	1129241	1748776	13.72%	1.883%
36	13.215	1901	1909	1916	rBV	1466061	2424582	19.02%	2.610%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
 Data File : VN072609.D
 Acq On : 25 May 2022 00:32
 Operator : JC\MD
 Sample : N2968-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-38

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\82N052422W.M
 Title : SW846 8260

37	13.363	1926	1934	1948	rVB	6381704	9883242	77.52%	10.640%
38	13.698	1980	1991	2001	rBV2	1602886	3443200	27.01%	3.707%
39	13.833	2006	2014	2015	rVV2	194137	306774	2.41%	0.330%
40	13.868	2015	2020	2037	rVB	1750083	3496760	27.43%	3.764%
41	14.057	2046	2052	2058	rVB2	179310	336364	2.64%	0.362%
42	14.121	2058	2063	2066	rBV	217687	355064	2.79%	0.382%
43	14.151	2066	2068	2073	rVV	171188	249176	1.95%	0.268%
44	14.210	2073	2078	2084	rVV	453508	711198	5.58%	0.766%
45	14.274	2084	2089	2092	rVV	113319	195342	1.53%	0.210%
46	14.321	2092	2097	2104	rVB2	381628	648649	5.09%	0.698%
47	14.451	2114	2119	2125	rVB2	115311	182410	1.43%	0.196%
48	14.533	2125	2133	2136	rBV	269376	430048	3.37%	0.463%
49	14.574	2136	2140	2145	rVB	303197	450072	3.53%	0.485%
50	14.804	2174	2179	2184	rBV	293727	459149	3.60%	0.494%
51	14.927	2191	2200	2206	rBV3	564055	1250651	9.81%	1.346%
52	15.080	2218	2226	2233	rVV	171992	308343	2.42%	0.332%
53	15.186	2233	2244	2248	rVV4	96894	249377	1.96%	0.268%
54	15.304	2257	2264	2271	rVB3	90635	224758	1.76%	0.242%
55	15.498	2290	2297	2305	rBV	1129045	2052918	16.10%	2.210%
56	16.609	2478	2486	2500	rVB	389954	883166	6.93%	0.951%

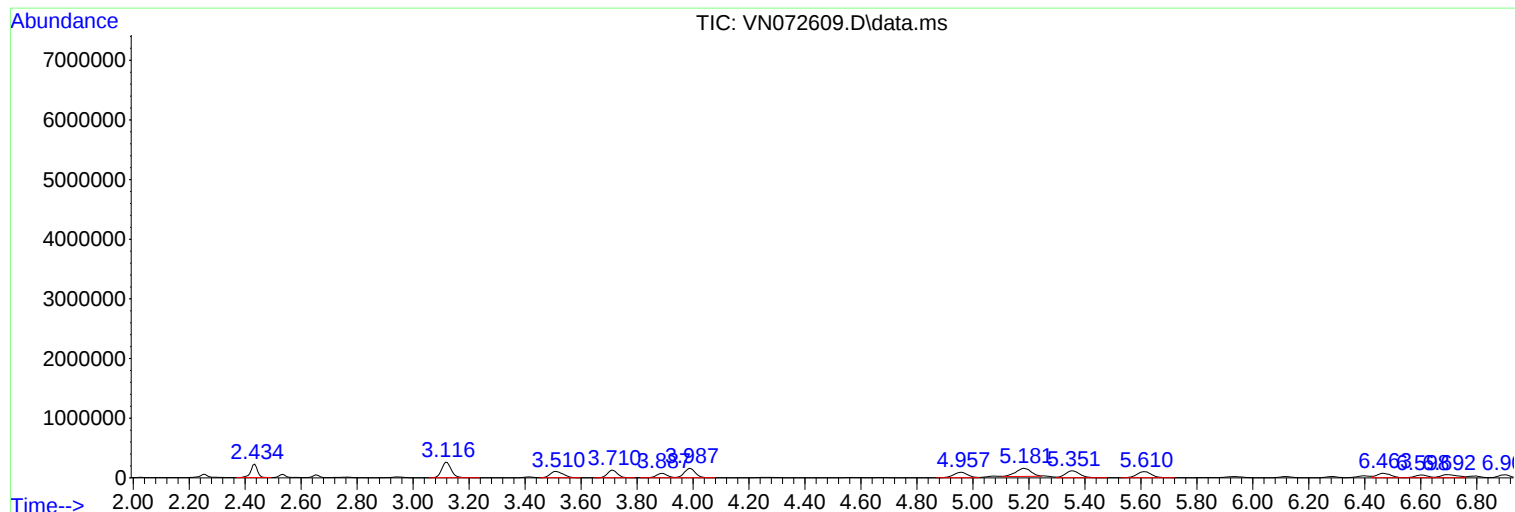
Sum of corrected areas: 92889190

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

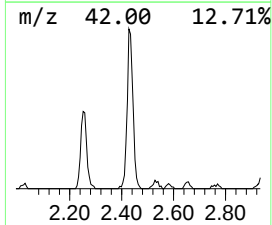
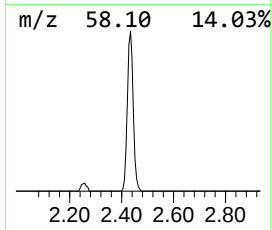
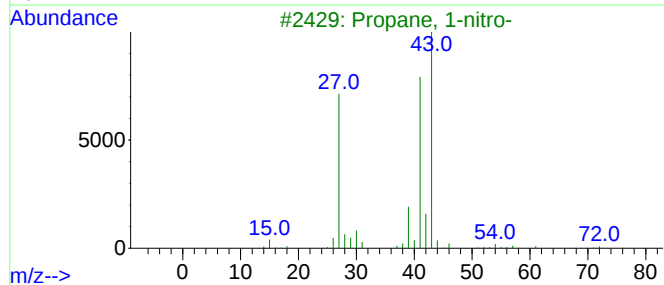
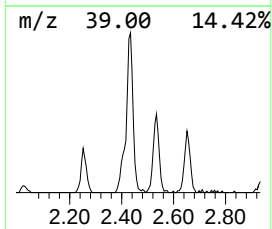
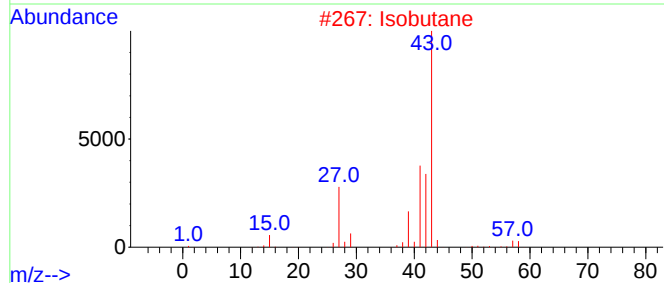
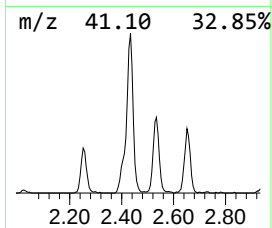
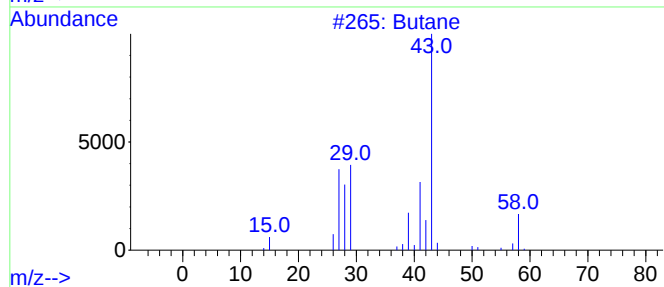
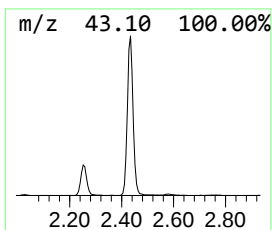
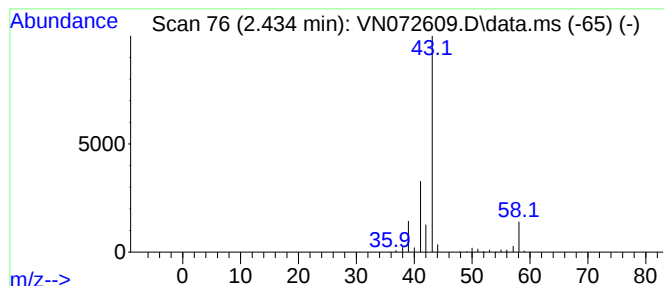
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.434	20.58 ug/l	383440	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane			58	C4H10	000106-97-8	80
2	Isobutane			58	C4H10	000075-28-5	9
3	Propane, 1-nitro-			89	C3H7NO2	000108-03-2	9
4	Isopropylsulfonyl chloride			142	C3H7ClO2S	010147-37-2	9
5	Acetone			58	C3H6O	000067-64-1	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

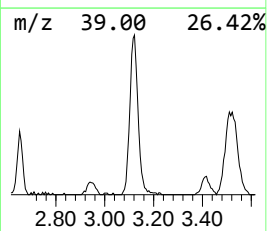
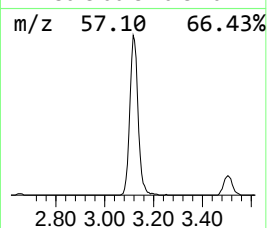
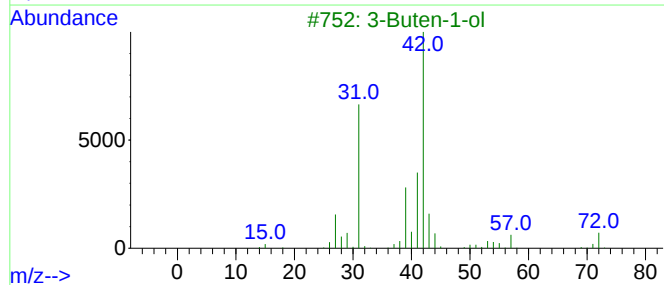
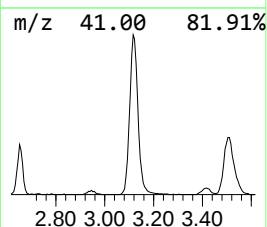
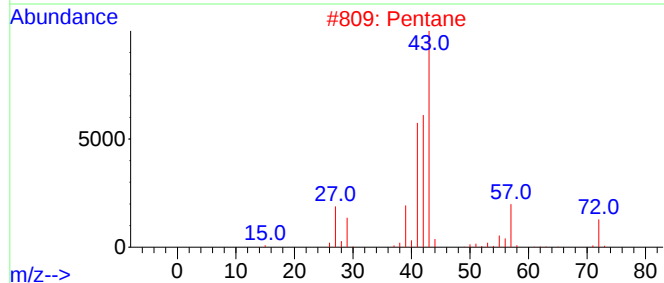
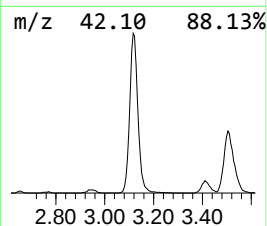
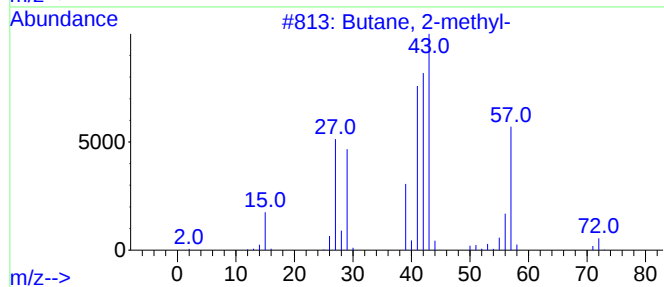
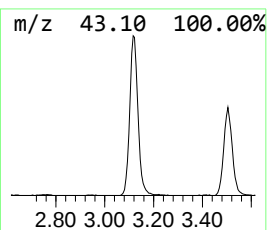
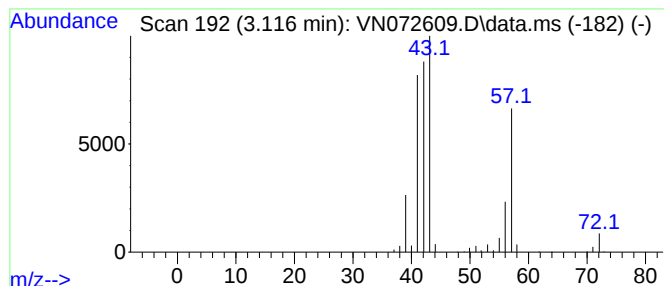
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.116	33.01 ug/l	615143	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	90
2			Pentane	72	C5H12	000109-66-0	50
3			3-Buten-1-ol	72	C4H8O	000627-27-0	28
4			1-Butene	56	C4H8	000106-98-9	10
5			Butane, 2-nitro-	103	C4H9NO2	000600-24-8	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

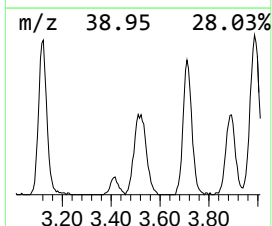
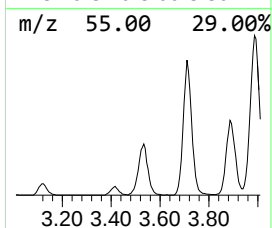
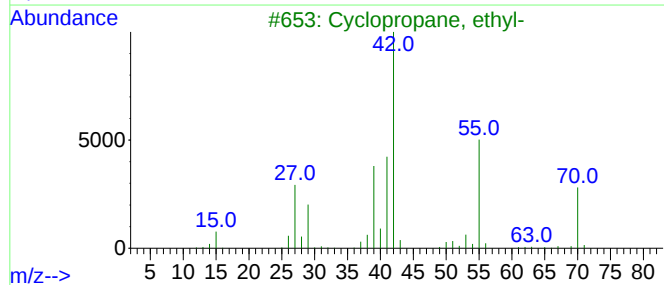
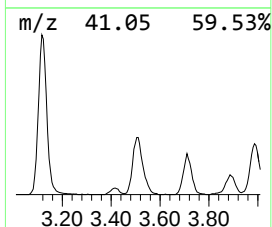
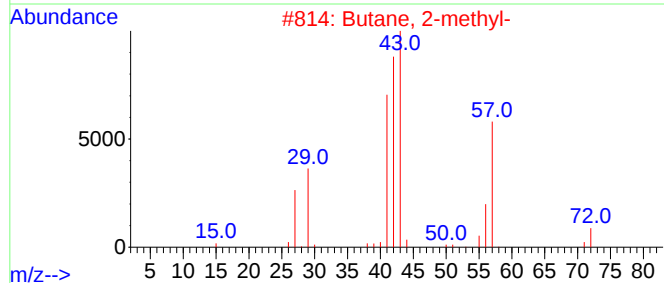
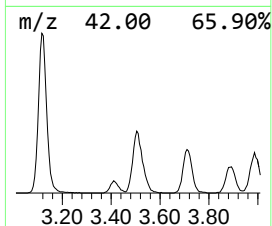
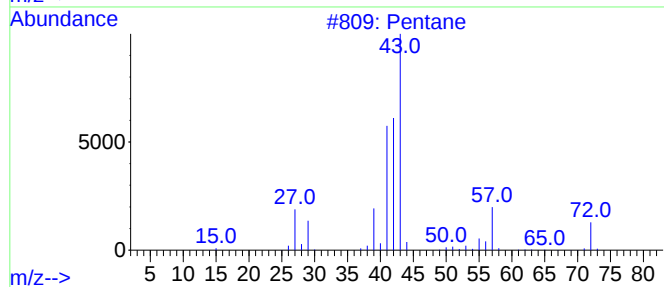
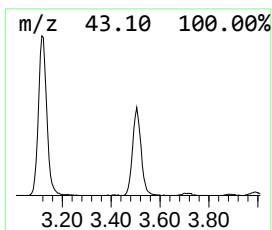
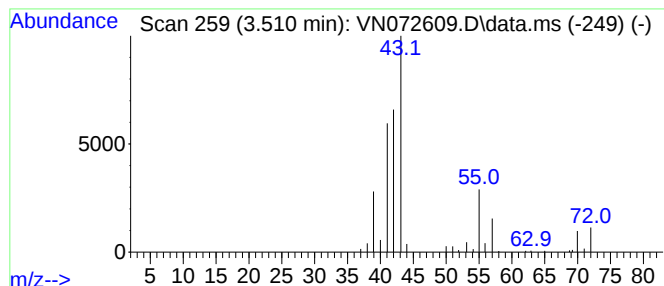
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.510	18.47 ug/l	344213	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	86
2			Butane, 2-methyl-	72	C5H12	000078-78-4	45
3			Cyclopropane, ethyl-	70	C5H10	001191-96-4	43
4			Oxirane, ethyl-	72	C4H8O	000106-88-7	37
5			1-Pentene	70	C5H10	000109-67-1	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

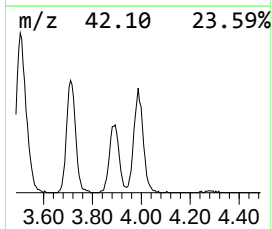
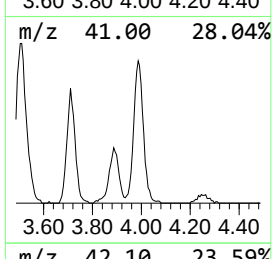
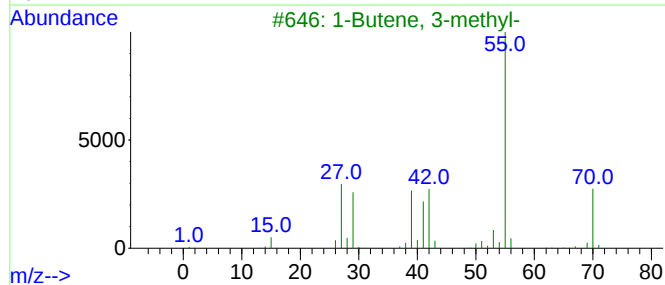
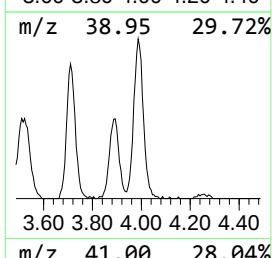
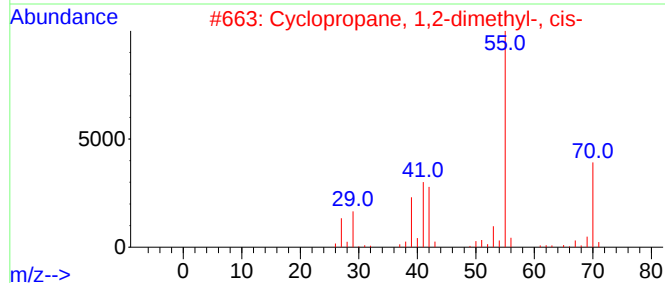
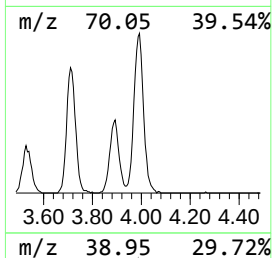
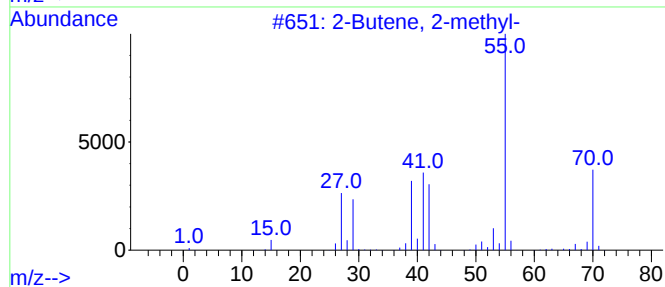
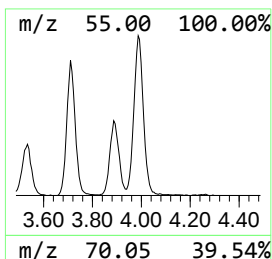
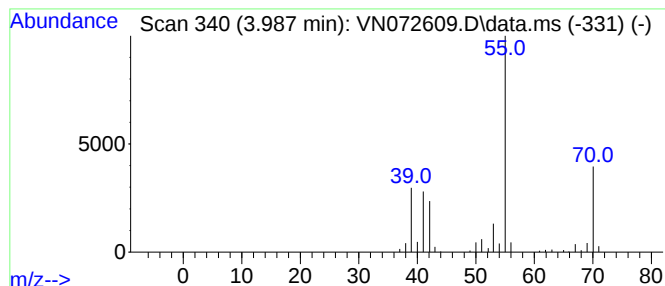
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 4 2-Butene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.987	23.36 ug/l	435320	Pentafluorobenzene	8.081

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Butene, 2-methyl-	70	C5H10	000513-35-9	91
2		Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
3		1-Butene, 3-methyl-	70	C5H10	000563-45-1	86
4		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	83
5		Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

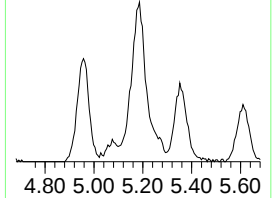
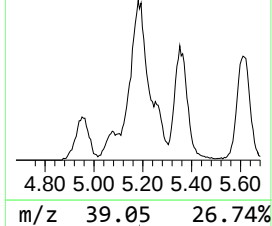
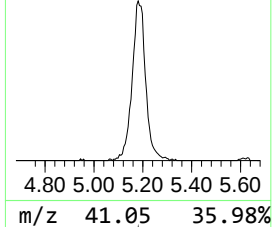
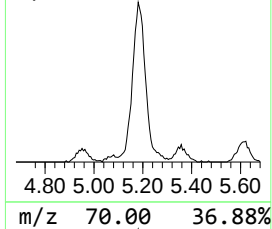
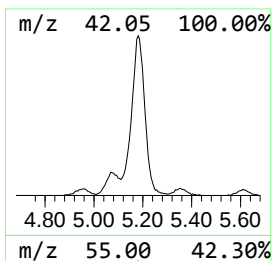
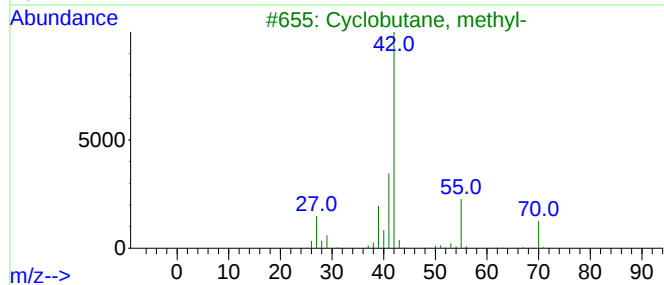
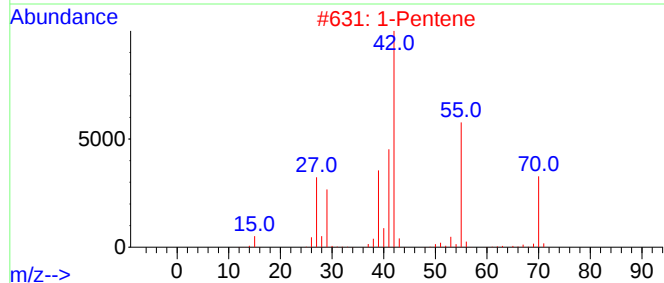
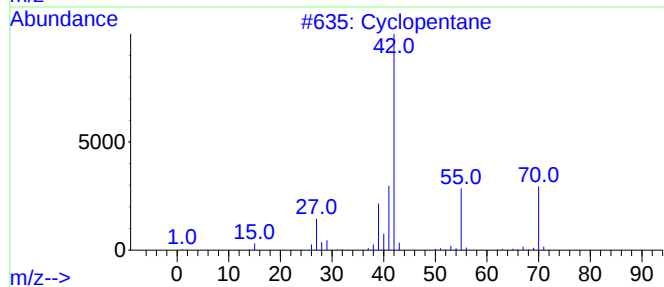
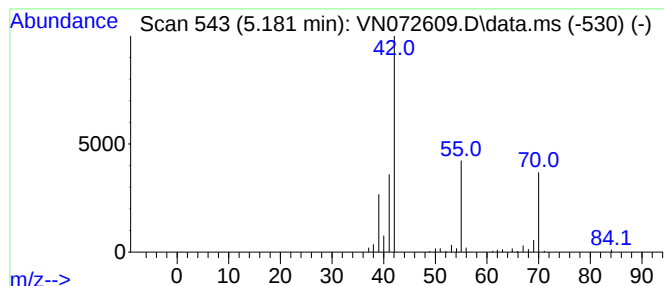
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.181	28.38 ug/l	528740	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	80
2			1-Pentene	70	C5H10	000109-67-1	78
3			Cyclobutane, methyl-	70	C5H10	000598-61-8	72
4			Cyclopropane, ethyl-	70	C5H10	001191-96-4	38
5			Cyclobutanone	70	C4H6O	001191-95-3	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

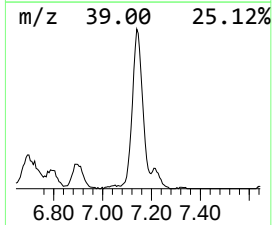
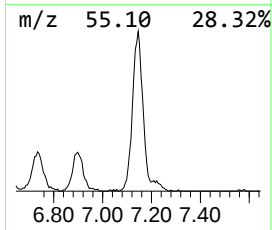
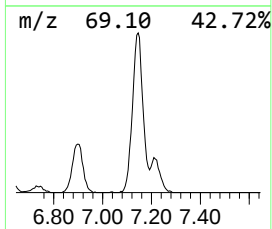
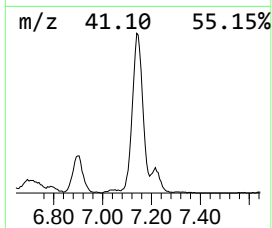
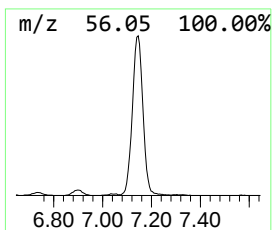
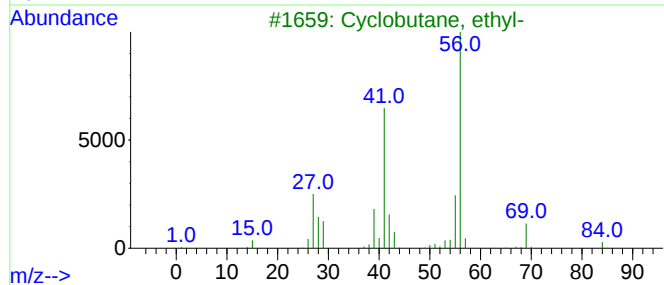
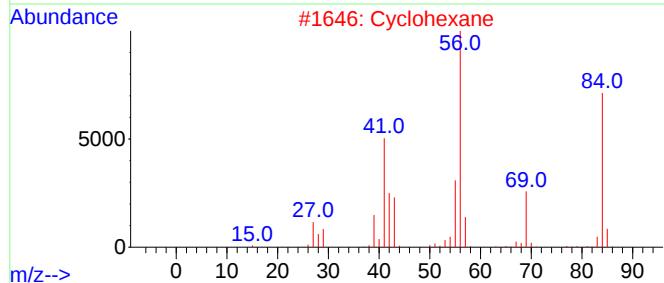
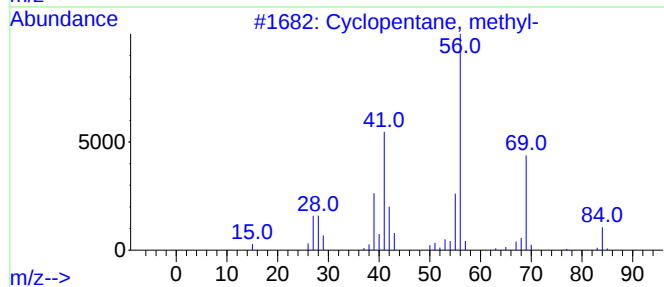
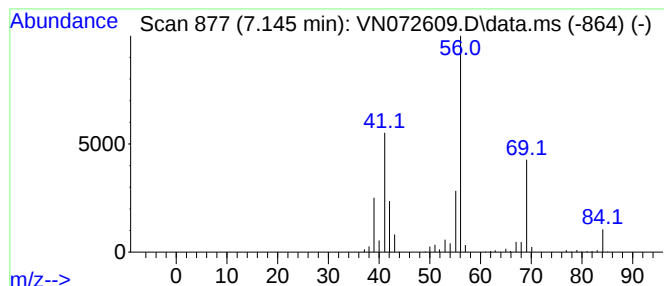
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.145	43.92 ug/l	818409	Pentafluorobenzene	8.081

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	83
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5			Cyclopropane, propyl-	84	C6H12	002415-72-7	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

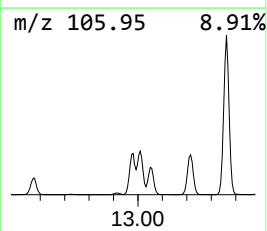
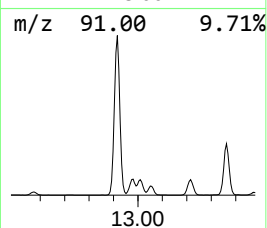
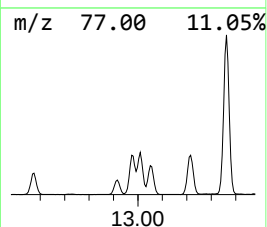
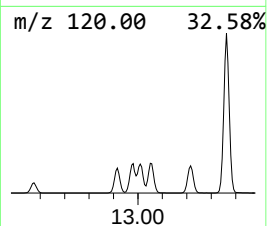
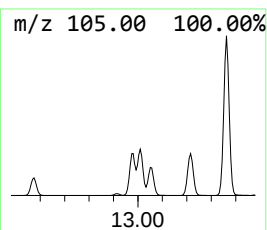
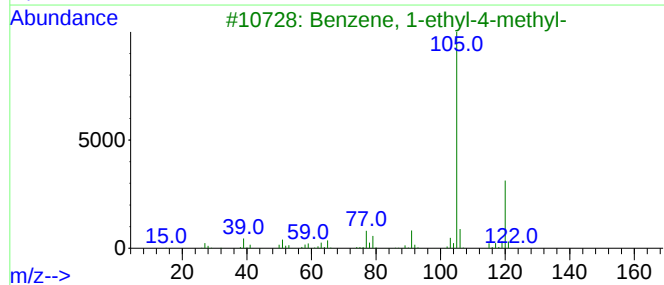
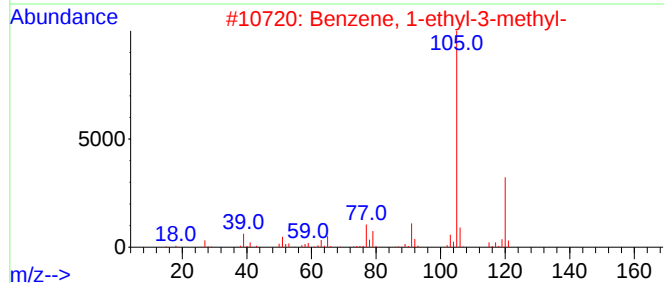
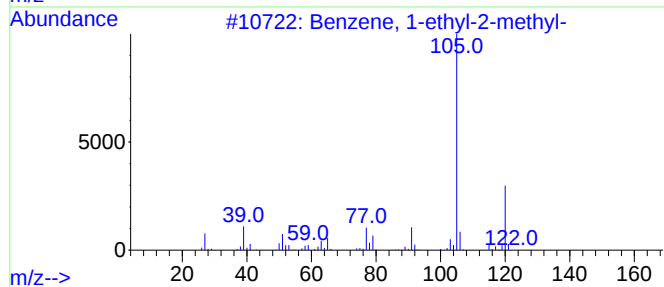
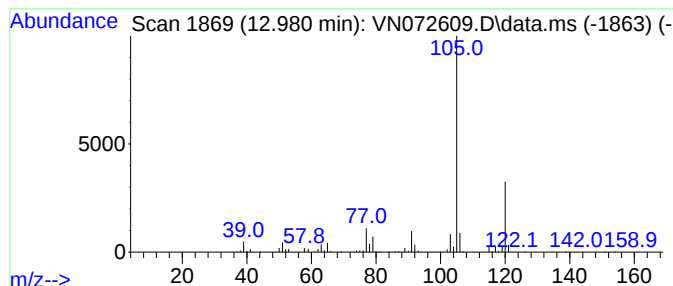
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	35.97 ug/l	2476790	1,4-Dichlorobenzene-d4	13.668

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			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_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

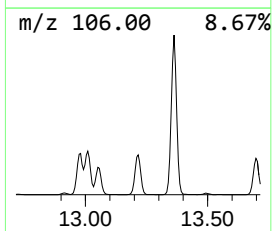
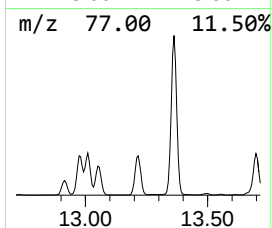
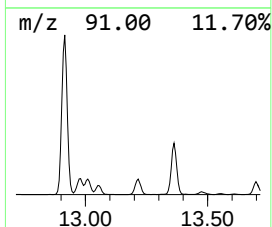
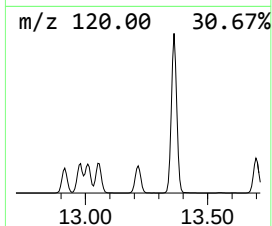
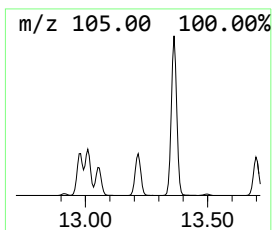
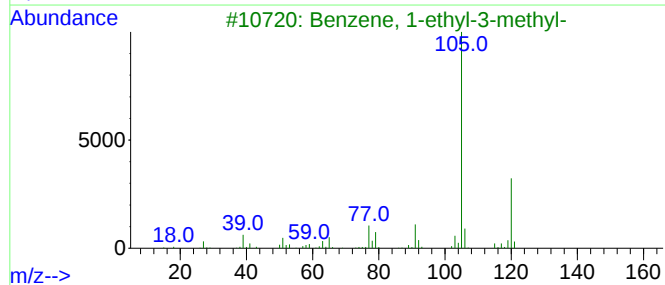
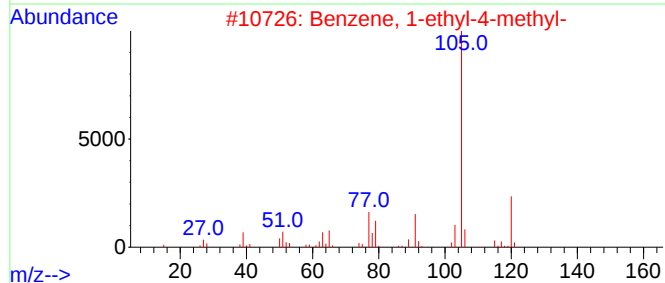
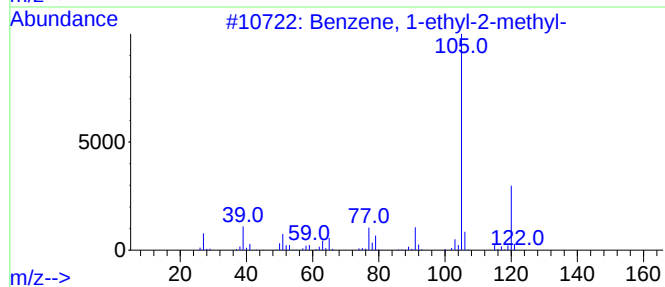
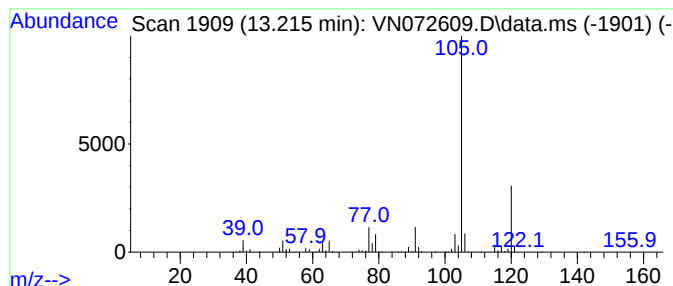
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	35.21 ug/l	2424580	1,4-Dichlorobenzene-d4	13.668

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-ethyl-3-methyl-	120	C9H12	000620-14-4	90
4			Mesitylene	120	C9H12	000108-67-8	90
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

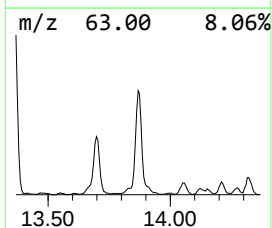
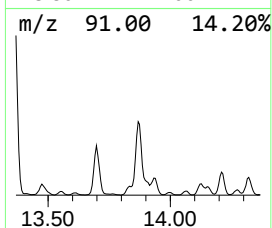
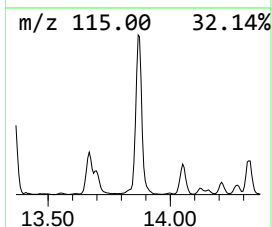
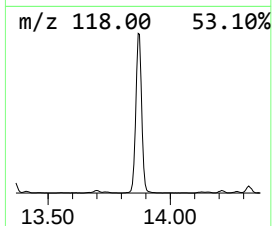
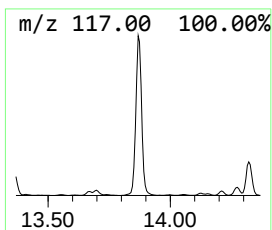
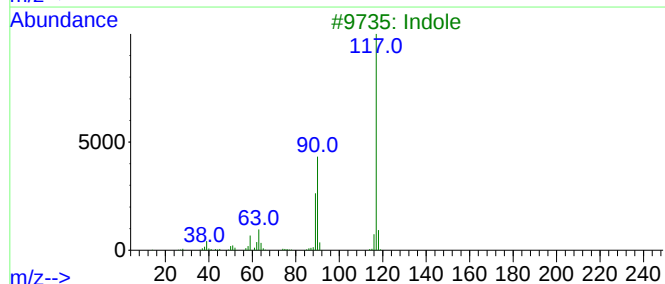
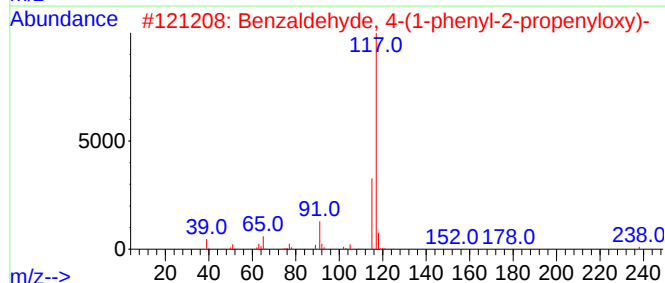
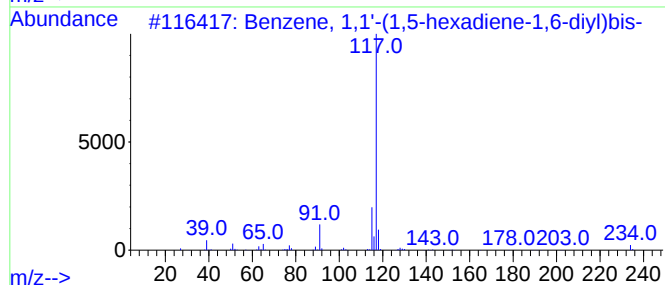
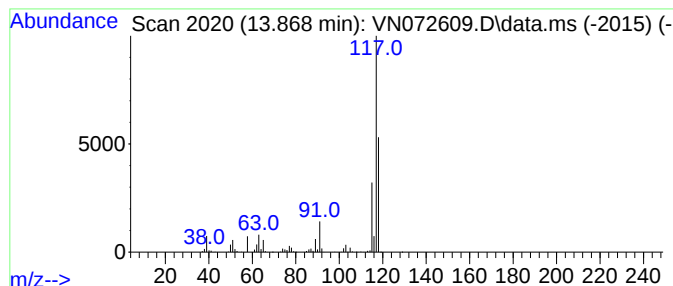
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.868	50.78 ug/l	3496760	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	50
3			Indole	117	C8H7N	000120-72-9	49
4			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
5			4-Ethynylaniline	117	C8H7N	014235-81-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

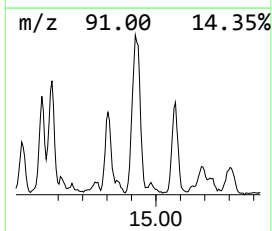
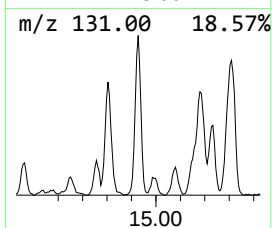
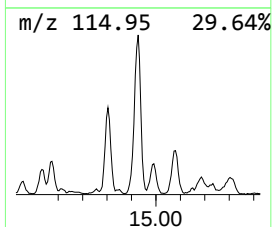
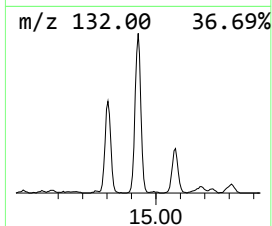
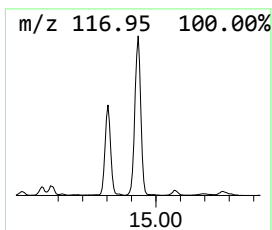
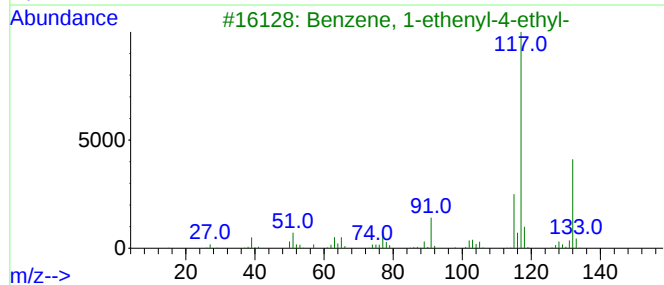
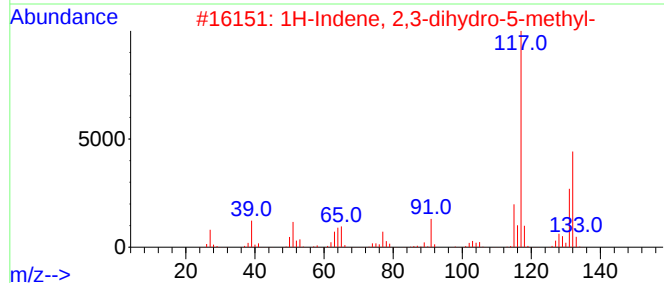
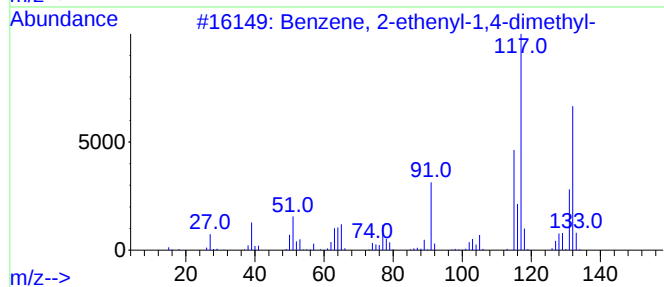
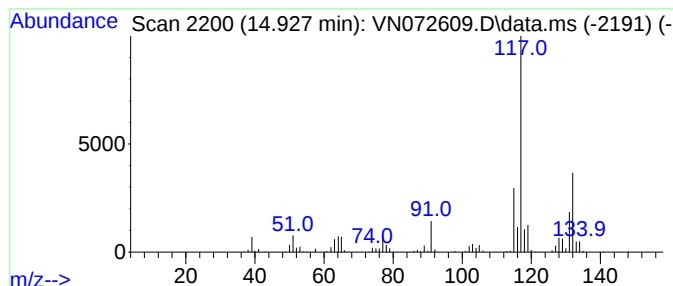
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	18.16 ug/l	1250650	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052422\
Data File : VN072609.D
Acq On : 25 May 2022 00:32
Operator : JC\MD
Sample : N2968-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-38

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N052422W.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
Butane	2.434	20.6	ug/l	383440	1	8.081	931683	50.0
Butane, 2-methyl-	3.116	33.0	ug/l	615143	1	8.081	931683	50.0
Pentane	3.510	18.5	ug/l	344213	1	8.081	931683	50.0
2-Butene, 2-met...	3.987	23.4	ug/l	435320	1	8.081	931683	50.0
Cyclopentane	5.181	28.4	ug/l	528740	1	8.081	931683	50.0
Cyclopentane, m...	7.145	43.9	ug/l	818409	1	8.081	931683	50.0
Benzene, 1-ethy...	12.980	36.0	ug/l	2476790	4	13.668	3443200	50.0
Benzene, 1-ethy...	13.216	35.2	ug/l	2424580	4	13.668	3443200	50.0
Benzene, 1,1'-(...	13.868	50.8	ug/l	3496760	4	13.668	3443200	50.0
Benzene, 2-ethe...	14.927	18.2	ug/l	1250650	4	13.668	3443200	50.0