

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
 Data File : VX033053.D
 Acq On : 24 Nov 2022 10:39
 Operator : JC/MD
 Sample : N5747-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-28

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\82X112222W.M
 Title : SW846 8260

Signal : TIC: VX033053.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.264	25	29	42	rBV2	40080	66855	4.81%	0.890%
2	1.374	42	47	53	rVB	55032	60783	4.37%	0.809%
3	1.752	104	109	118	rVB	85136	113914	8.19%	1.516%
4	1.965	138	144	150	rVB3	17698	35954	2.59%	0.479%
5	2.215	181	185	196	rVB	28484	44539	3.20%	0.593%
6	2.757	261	274	276	rBV	10380	21958	1.58%	0.292%
7	2.867	287	292	301	rVB	27940	57398	4.13%	0.764%
8	2.989	301	312	326	rVV	140790	562355	40.45%	7.485%
9	3.111	326	332	345	rVB	65043	157877	11.36%	2.101%
10	3.751	427	437	446	rBV4	16433	50449	3.63%	0.671%
11	4.312	519	529	539	rVV3	20894	60110	4.32%	0.800%
12	5.385	695	705	714	rBV	63695	183387	13.19%	2.441%
13	5.550	724	732	742	rVB2	127377	352296	25.34%	4.689%
14	5.958	788	799	804	rBV	67720	176805	12.72%	2.353%
15	6.037	804	812	827	rVV	522125	1390188	100.00%	18.503%
16	6.251	831	847	858	rVV6	17088	72413	5.21%	0.964%
17	6.354	860	864	878	rVB8	4709	14677	1.06%	0.195%
18	6.763	921	931	942	rBV	205505	492425	35.42%	6.554%
19	7.366	1022	1030	1041	rBV7	4644	14299	1.03%	0.190%
20	8.110	1144	1152	1155	rBV3	8532	18555	1.33%	0.247%
21	8.141	1155	1157	1163	rVB	10408	15161	1.09%	0.202%
22	8.427	1198	1204	1211	rBV	10377	19328	1.39%	0.257%
23	8.506	1211	1217	1225	rVB4	9082	17530	1.26%	0.233%
24	8.647	1233	1240	1247	rBV	324400	503453	36.21%	6.701%
25	8.720	1247	1252	1260	rVB	45462	72774	5.23%	0.969%
26	8.958	1281	1291	1299	rVB3	17959	31443	2.26%	0.418%
27	9.049	1299	1306	1312	rBV2	20400	31886	2.29%	0.424%
28	9.457	1365	1373	1377	rBV2	22379	37296	2.68%	0.496%
29	10.055	1465	1471	1477	rBV	435471	574659	41.34%	7.648%
30	10.195	1489	1494	1501	rBV	80002	99508	7.16%	1.324%
31	10.299	1506	1511	1520	rVB	145861	205996	14.82%	2.742%
32	10.640	1562	1567	1577	rVB	69998	91741	6.60%	1.221%
33	10.963	1614	1620	1625	rBV	69298	86001	6.19%	1.145%
34	11.079	1634	1639	1647	rBV	274432	356280	25.63%	4.742%
35	11.305	1671	1676	1684	rVB	53948	65324	4.70%	0.869%
36	11.616	1722	1727	1736	rVB	18281	24400	1.76%	0.325%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
 Data File : VX033053.D
 Acq On : 24 Nov 2022 10:39
 Operator : JC/MD
 Sample : N5747-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-28

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\82X112222W.M
 Title : SW846 8260

37	11.750	1745	1749	1755	rVB	32257	40654	2.92%	0.541%
38	12.024	1789	1794	1800	rVV	437165	551865	39.70%	7.345%
39	12.085	1800	1804	1809	rVB	16718	21405	1.54%	0.285%
40	12.244	1824	1830	1837	rBV	280262	341492	24.56%	4.545%
41	12.311	1837	1841	1848	rVV2	12289	18587	1.34%	0.247%
42	12.463	1862	1866	1871	rVB	22410	27006	1.94%	0.359%
43	12.615	1884	1891	1893	rBV	20451	25157	1.81%	0.335%
44	12.646	1893	1896	1900	rVV	19146	22440	1.61%	0.299%
45	12.701	1900	1905	1913	rVV	65605	86632	6.23%	1.153%
46	12.920	1934	1941	1945	rVV	28694	36163	2.60%	0.481%
47	13.170	1979	1982	1991	rVB	17445	22393	1.61%	0.298%
48	13.292	1998	2002	2007	rVB2	33999	41880	3.01%	0.557%
49	13.579	2046	2049	2054	rVV3	8101	13950	1.00%	0.186%
50	13.664	2060	2063	2070	rVB3	11623	16015	1.15%	0.213%
51	13.774	2074	2081	2090	rBV	28226	39643	2.85%	0.528%
52	14.780	2241	2246	2254	rBV2	19904	28171	2.03%	0.375%

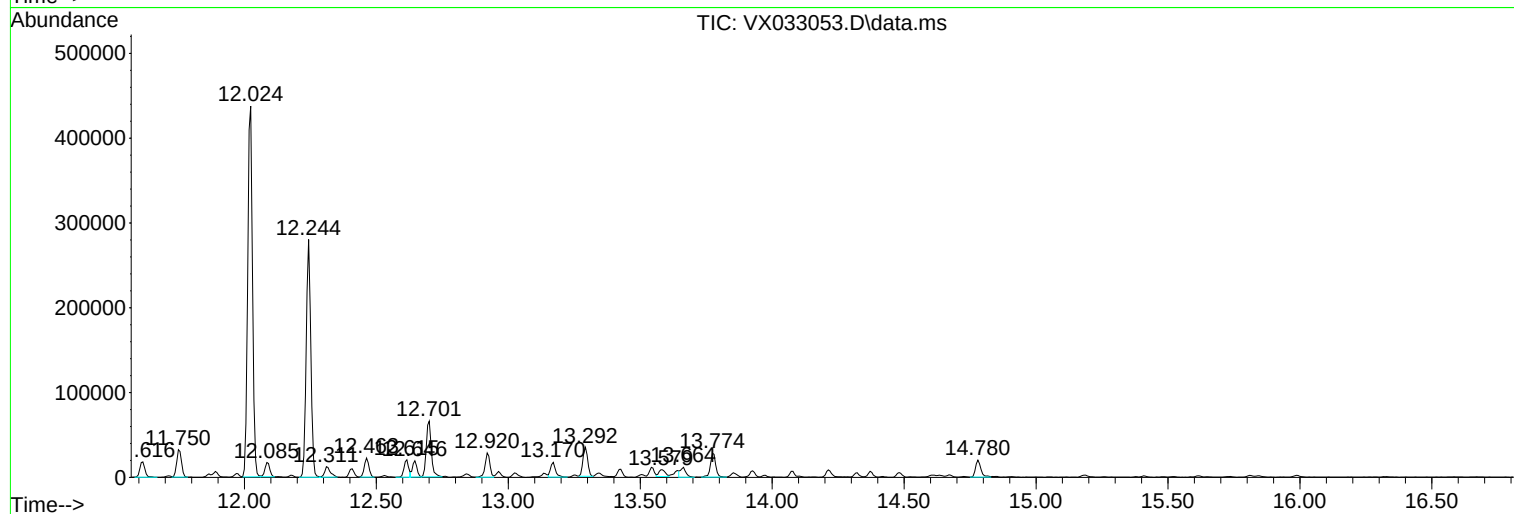
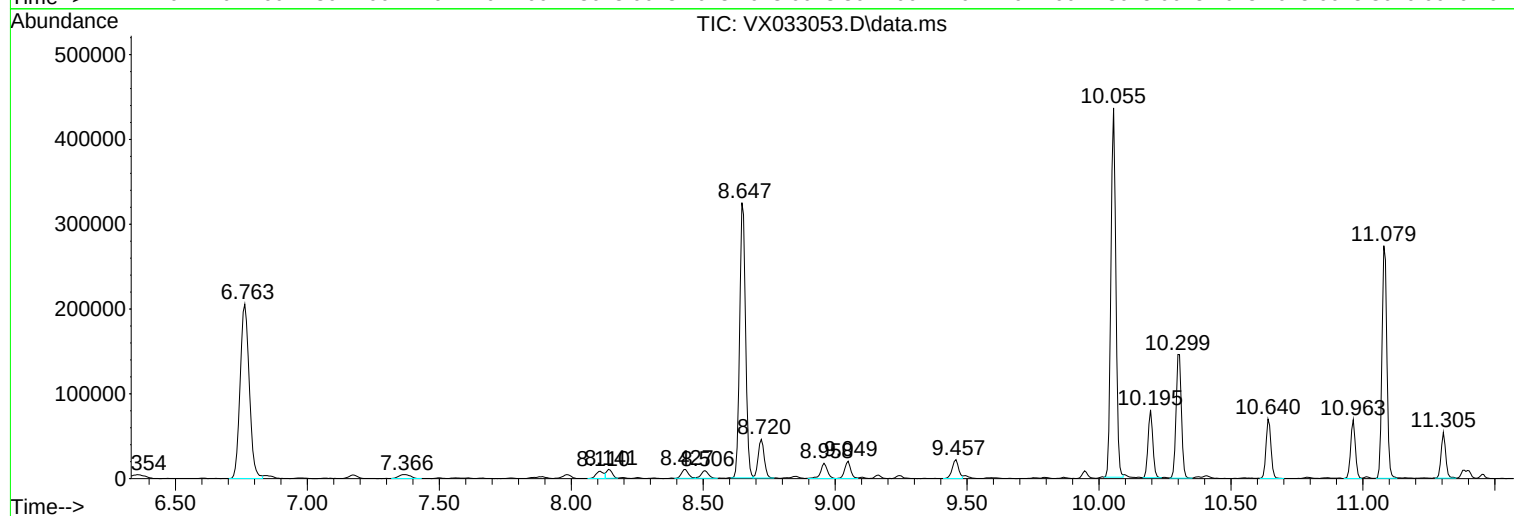
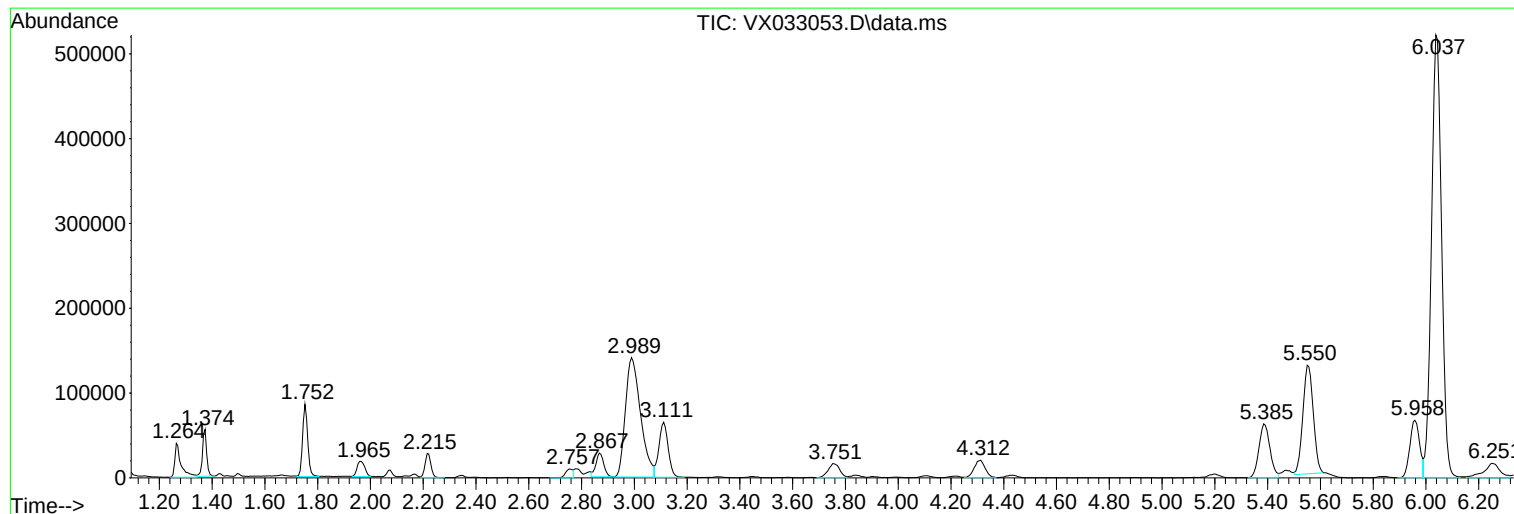
Sum of corrected areas: 7513470

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

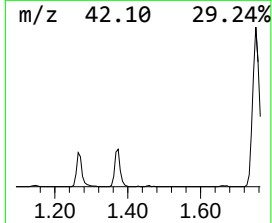
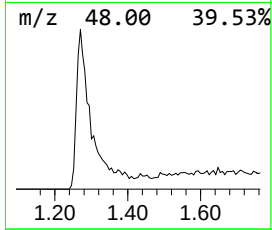
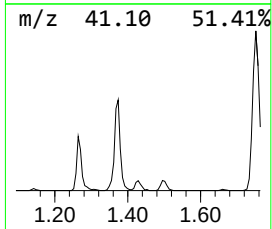
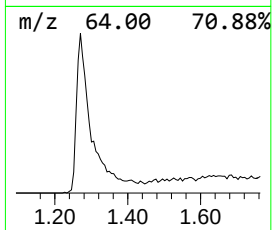
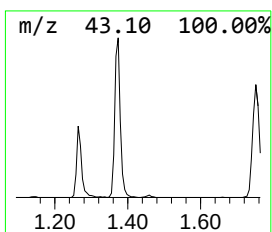
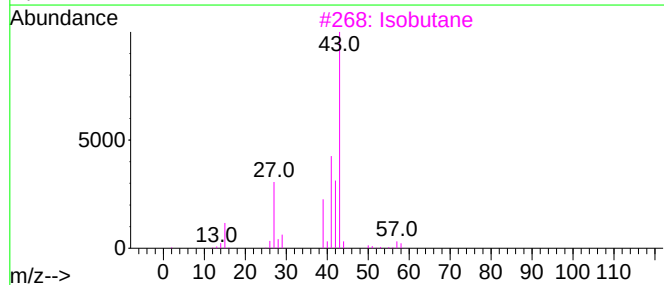
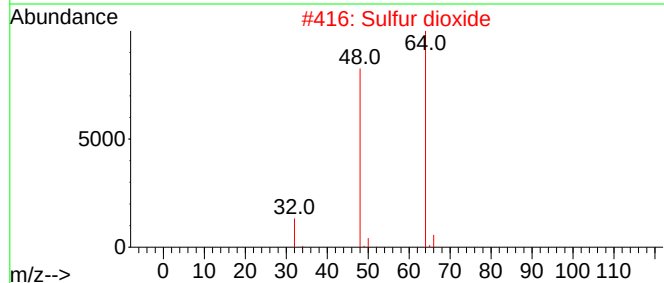
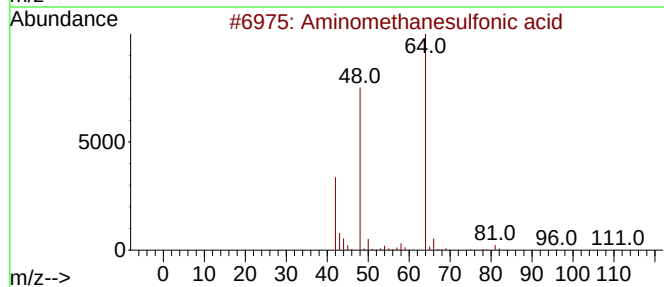
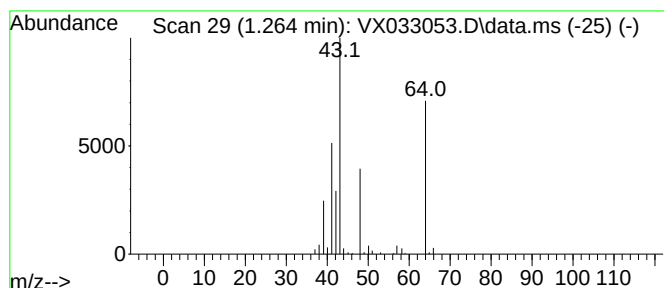
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 1 Aminomethanesulfonic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.264	9.49 ug/l	66855	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	56
2			Sulfur dioxide	64	O2S	007446-09-5	45
3			Isobutane	58	C4H10	000075-28-5	10
4			Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5			Hepta-4,6-diyne-2-ol	108	C7H8O	1000154-66-3	4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

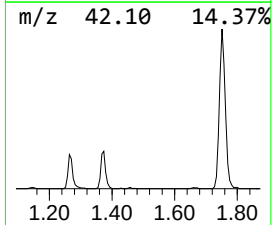
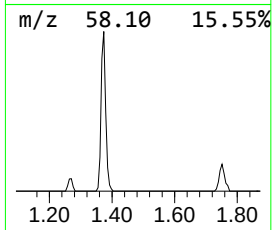
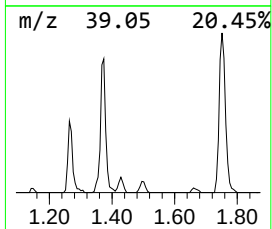
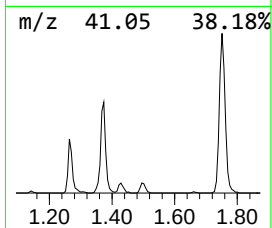
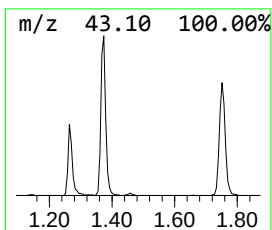
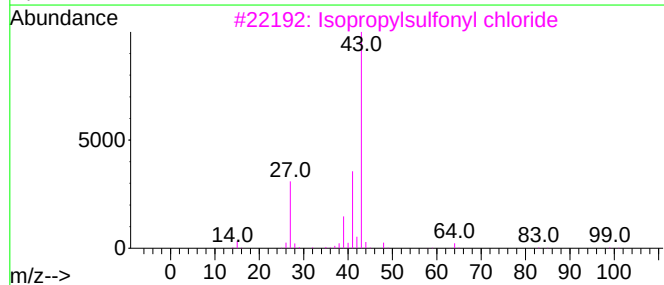
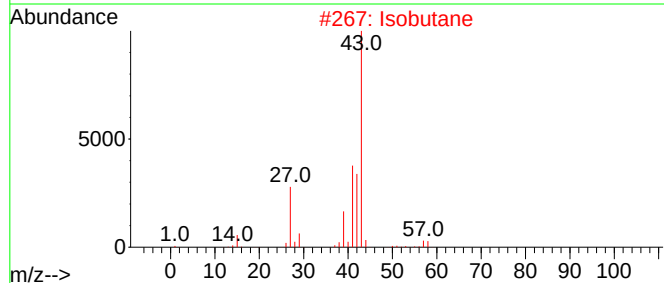
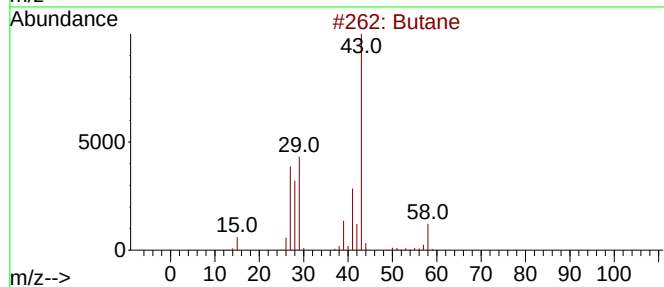
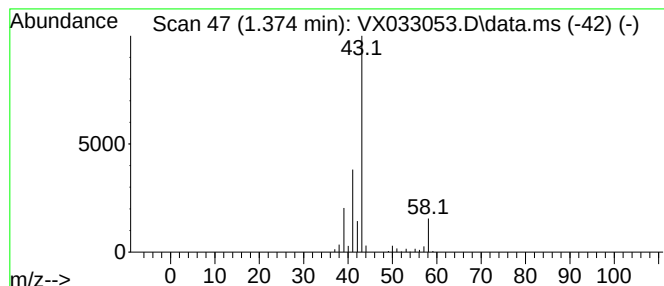
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.374	8.63 ug/l	60783	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane	58	C4H10	000106-97-8	72
2			Isobutane	58	C4H10	000075-28-5	9
3			Isopropylsulfonyl chloride	142	C3H7ClO2S	010147-37-2	9
4			Propane, 1-nitro-	89	C3H7NO2	000108-03-2	9
5			Acetone	58	C3H6O	000067-64-1	7



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

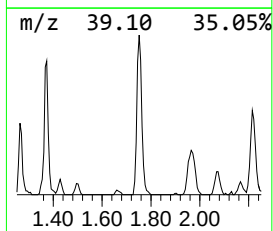
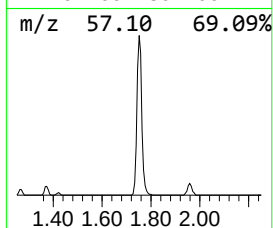
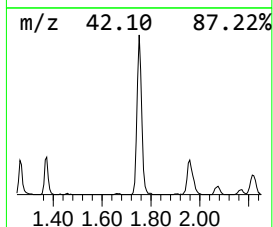
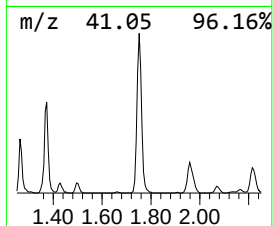
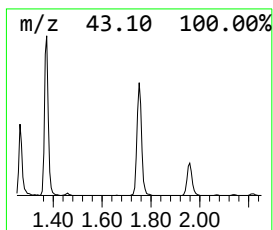
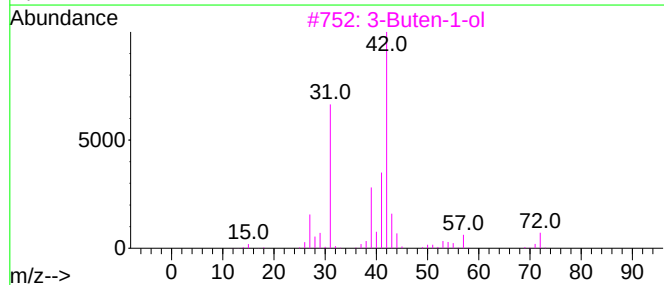
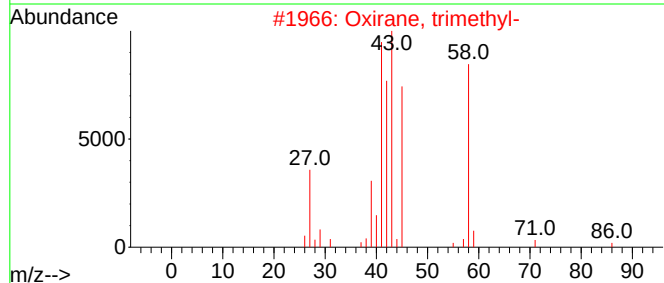
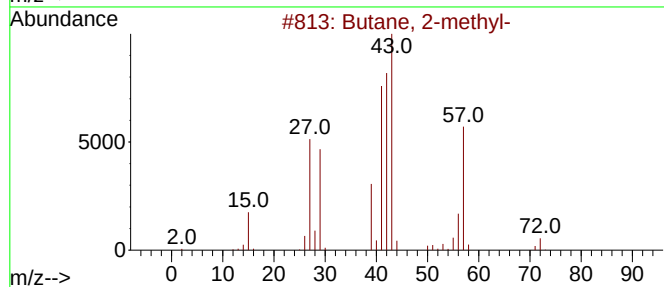
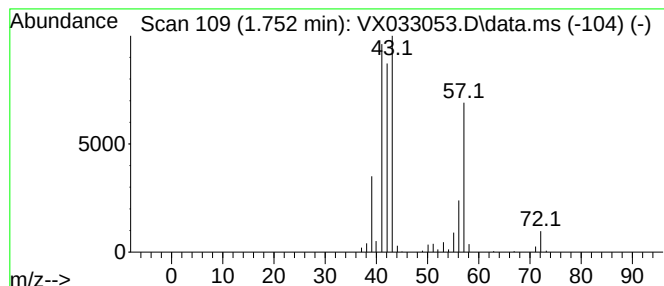
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	16.17 ug/l	113914	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	90
2			Oxirane, trimethyl-	86	C5H10O	005076-19-7	40
3			3-Buten-1-ol	72	C4H8O	000627-27-0	38
4			Pentane	72	C5H12	000109-66-0	33
5			2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	28



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

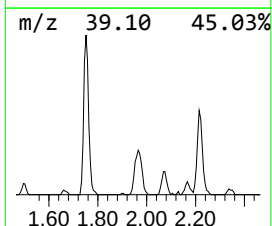
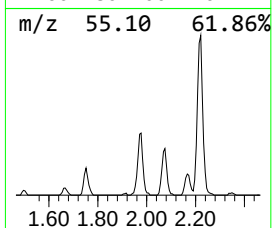
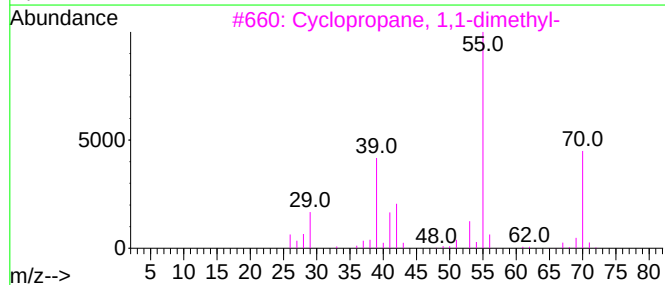
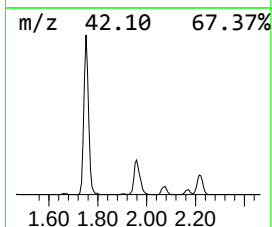
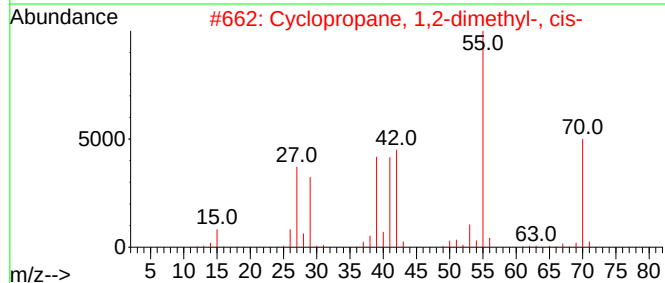
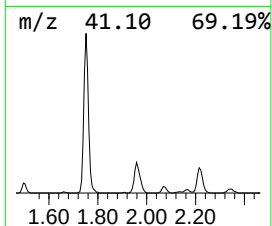
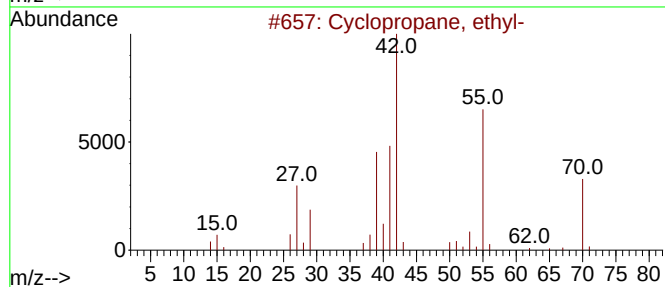
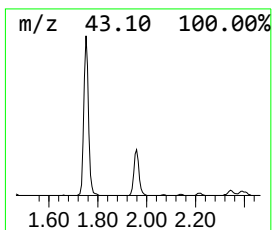
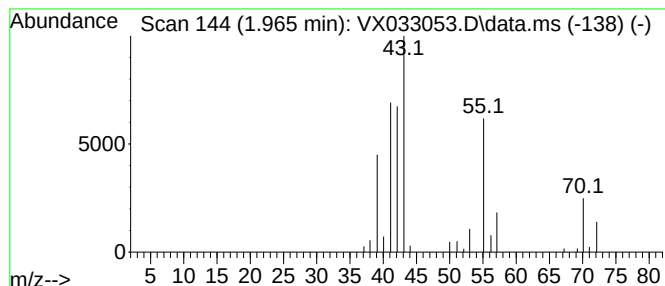
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 4 Cyclopropane, ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.965	5.10 ug/l	35954	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, ethyl-	70	C5H10	001191-96-4	62
2			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	58
3			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	50
4			Pentane	72	C5H12	000109-66-0	38
5			2-Pentene, (E)-	70	C5H10	000646-04-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

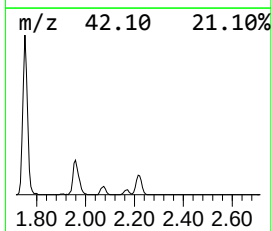
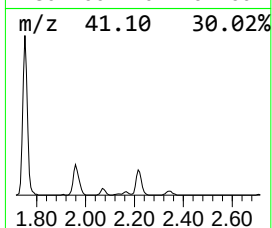
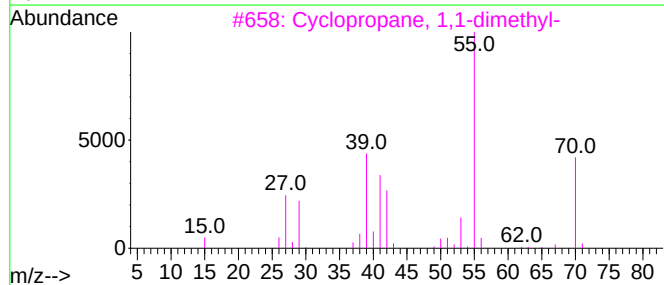
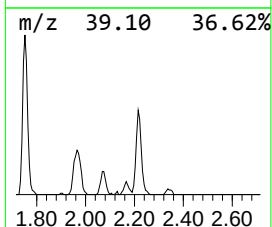
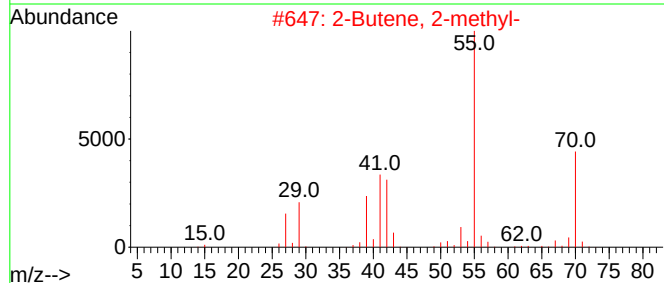
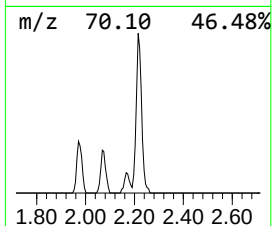
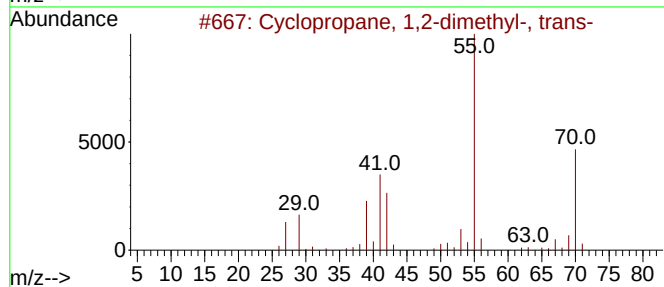
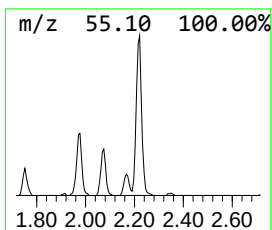
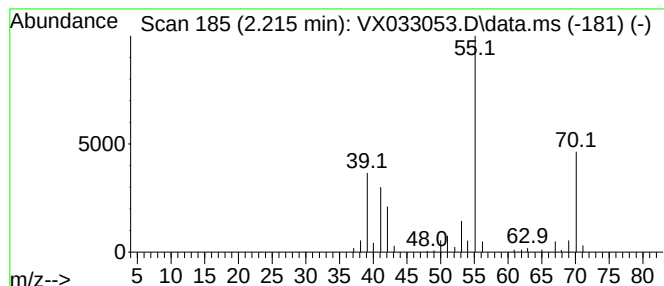
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 5 Cyclopropane, 1,2-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.215	6.32 ug/l	44539	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	91
2			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
3			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80
4			1-Butene, 3-methyl-	70	C5H10	000563-45-1	72
5			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

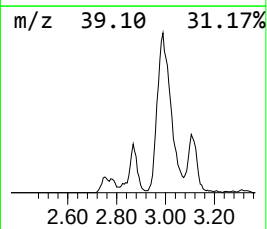
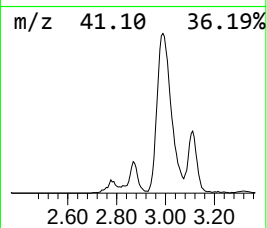
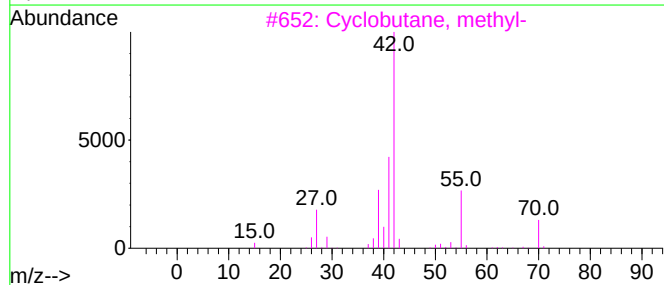
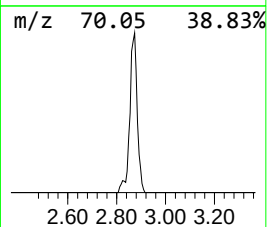
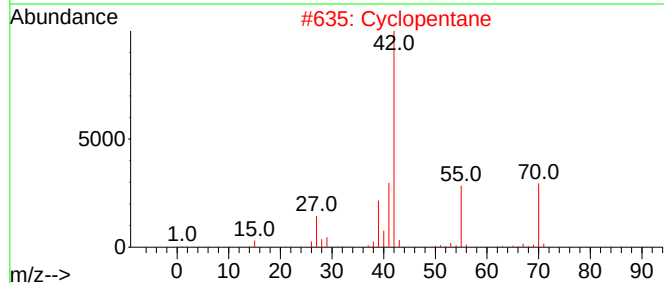
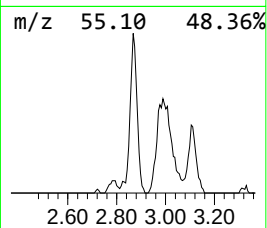
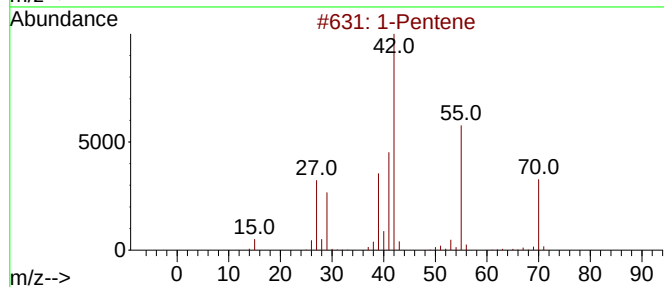
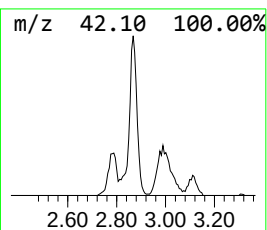
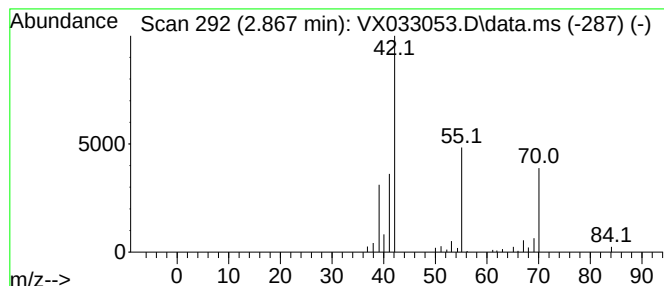
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 6 1-Pentene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.867	8.15 ug/l	57398	Pentafluorobenzene	5.550

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene	70	C5H10	000109-67-1	86
2		Cyclopentane	70	C5H10	000287-92-3	86
3		Cyclobutane, methyl-	70	C5H10	000598-61-8	78
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	53
5		1-Pentanol	88	C5H12O	000071-41-0	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

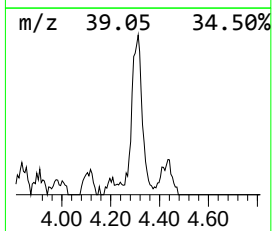
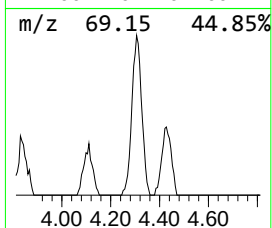
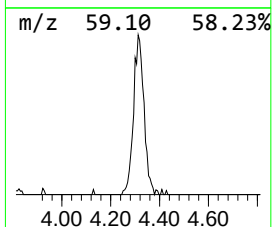
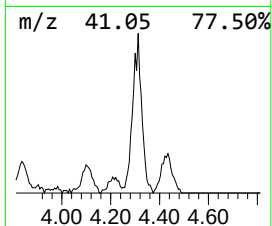
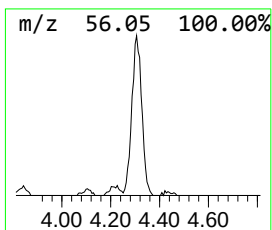
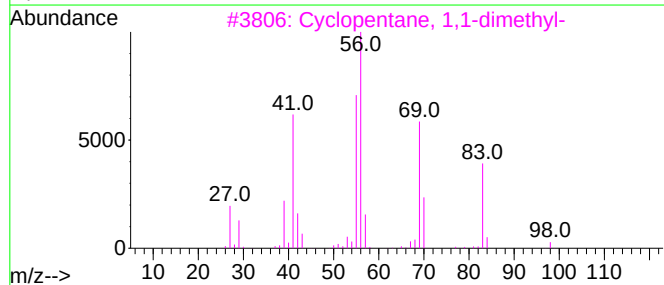
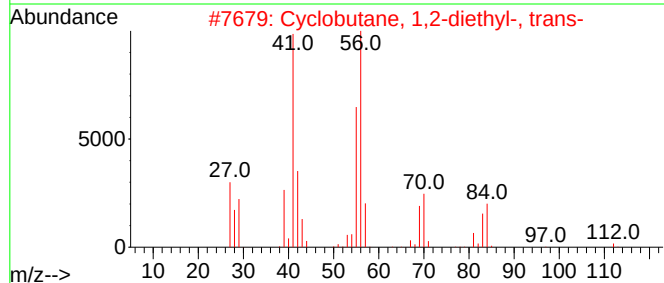
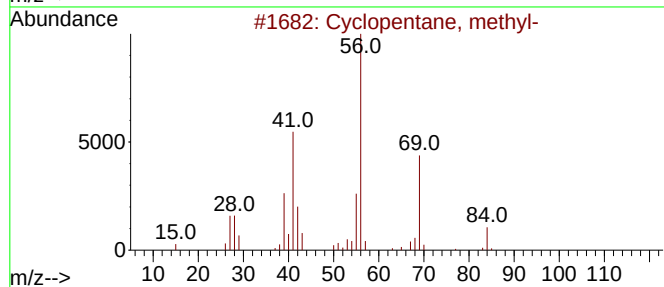
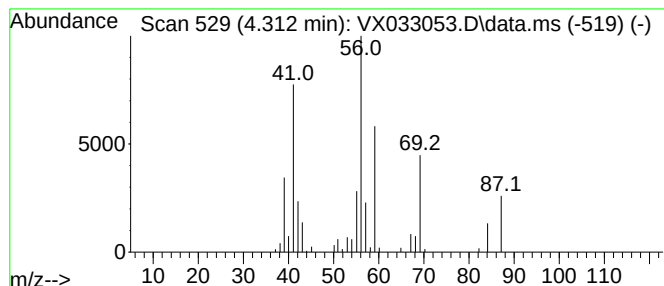
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.312	8.53 ug/l	60110	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	70
2			Cyclobutane, 1,2-diethyl-, trans-	112	C8H16	019341-98-1	50
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	47
4			Cyclohexane	84	C6H12	000110-82-7	47
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

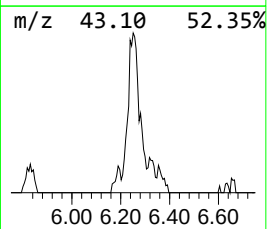
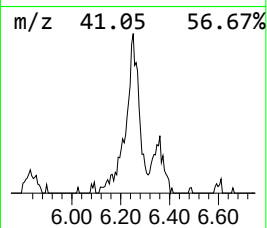
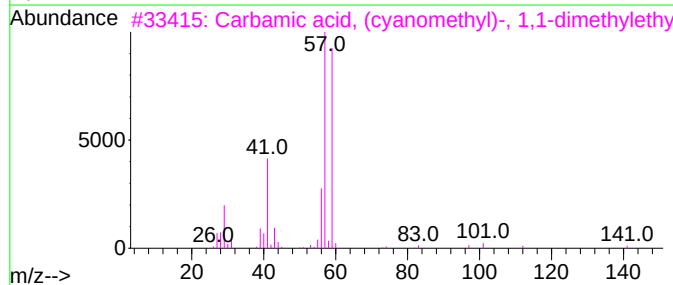
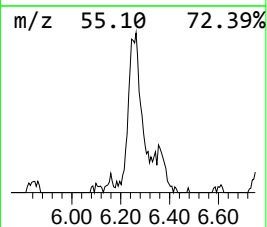
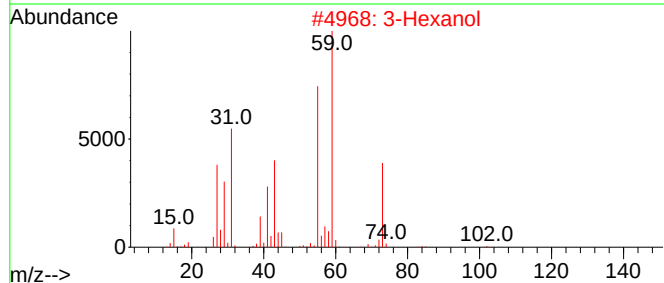
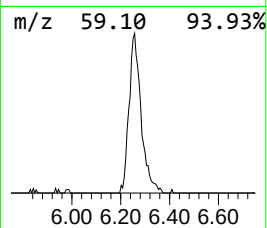
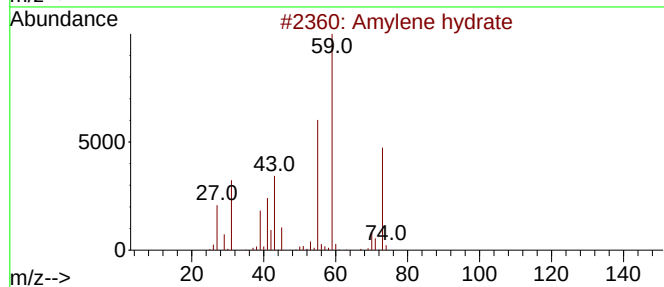
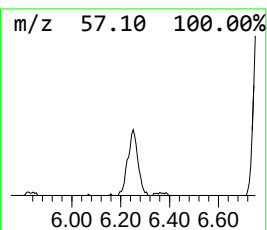
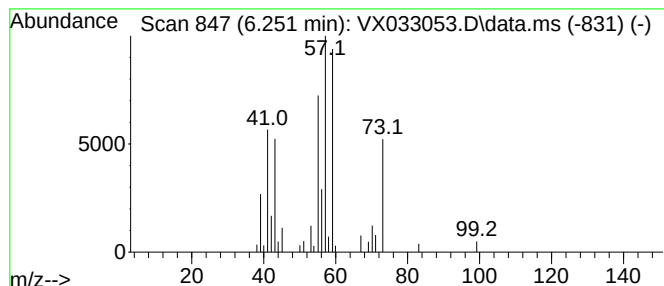
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 8 Amylene hydrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	7.35 ug/l	72413	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Amylene hydrate	88	C5H12O	000075-85-4	53
2			3-Hexanol	102	C6H14O	000623-37-0	50
3			Carbamic acid, (cyanomethyl)-, 1...	156	C7H12N2O2	085363-04-8	40
4			Formic acid, 1,1-dimethylethyl e...	102	C5H10O2	000762-75-4	40
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

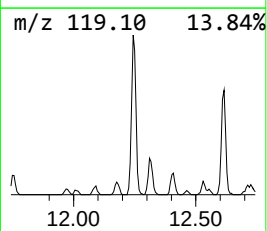
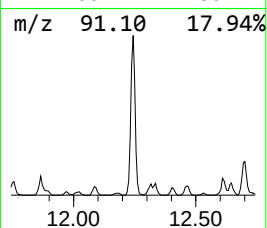
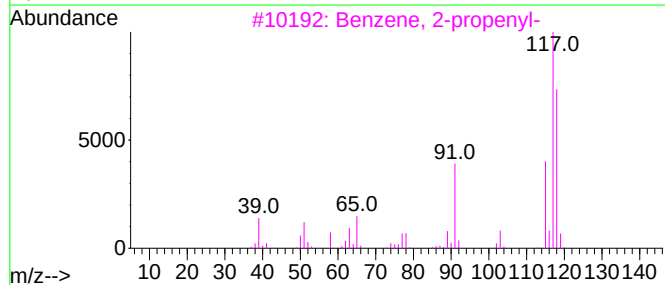
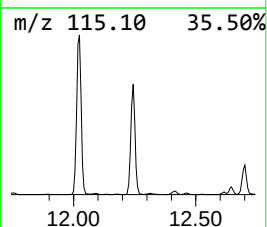
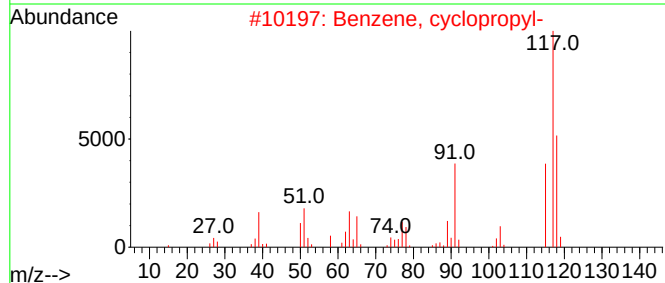
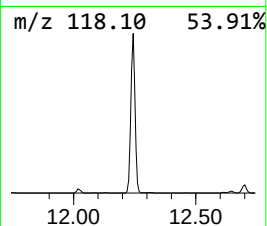
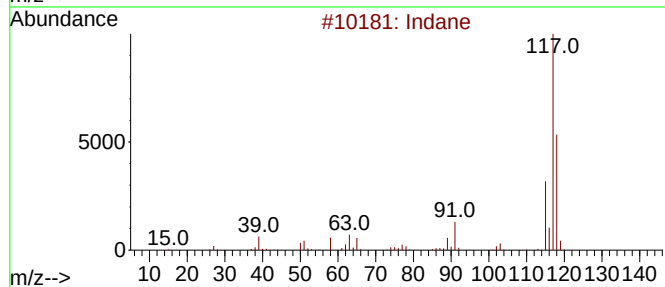
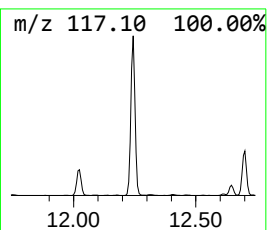
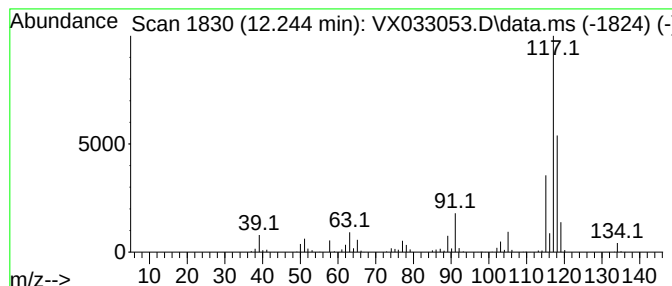
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 9 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	30.94 ug/l	341492	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Delatcyclene	118	C9H10	007785-10-6	62
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

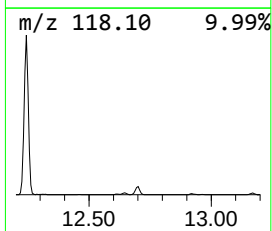
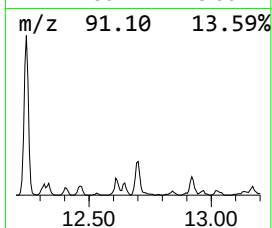
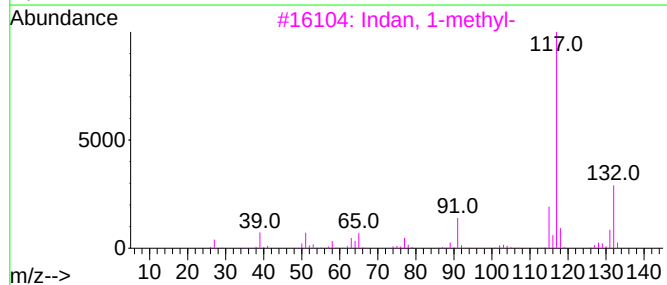
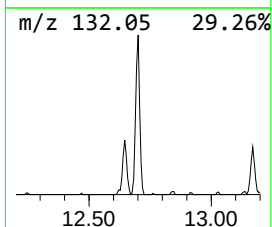
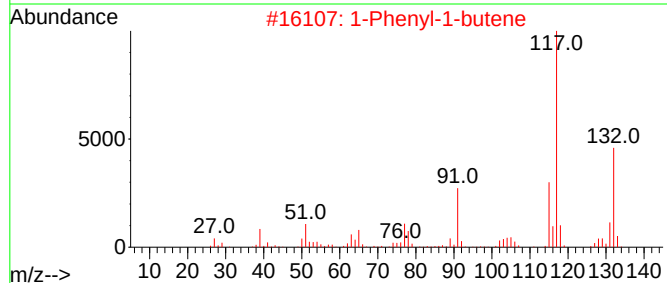
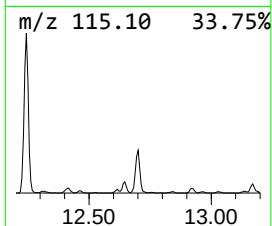
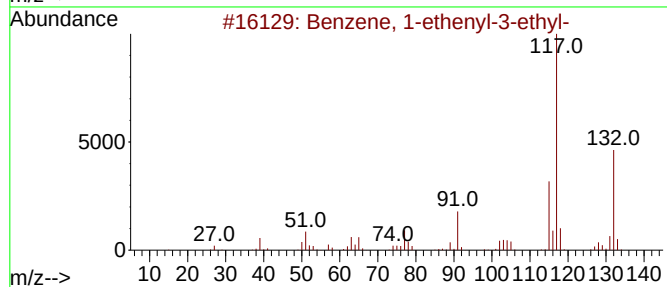
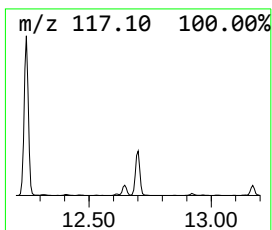
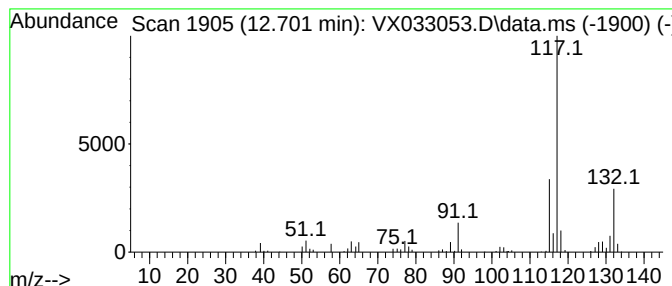
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	7.85 ug/l	86632	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
3			Indan, 1-methyl-	132	C10H12	000767-58-8	90
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
Data File : VX033053.D
Acq On : 24 Nov 2022 10:39
Operator : JC/MD
Sample : N5747-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-28

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.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
Aminomethanesul...	1.264	9.5	ug/l	66855	1	5.550	352296	50.0
Butane	1.374	8.6	ug/l	60783	1	5.550	352296	50.0
Butane, 2-methyl-	1.752	16.2	ug/l	113914	1	5.550	352296	50.0
Cyclopropane, e...	1.965	5.1	ug/l	35954	1	5.550	352296	50.0
Cyclopropane, 1...	2.215	6.3	ug/l	44539	1	5.550	352296	50.0
1-Pentene	2.867	8.2	ug/l	57398	1	5.550	352296	50.0
Cyclopentane, m...	4.312	8.5	ug/l	60110	1	5.550	352296	50.0
Amylene hydrate	6.251	7.3	ug/l	72413	2	6.763	492425	50.0
Indane	12.244	30.9	ug/l	341492	4	12.024	551865	50.0
Benzene, 1-ethe...	12.701	7.8	ug/l	86632	4	12.024	551865	50.0