

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
 Data File : VV026799.D
 Acq On : 15 Jul 2022 18:37
 Operator : SY/MD
 Sample : N3710-17
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 H0B09

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

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_V\Method\SFAMVTR071122WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VV026799.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.105	17	20	24	rVV	34188	26497	2.71%	0.201%
2	1.134	25	29	38	rVB	62613	79563	8.13%	0.603%
3	1.208	48	52	69	rVV2	30606	37440	3.83%	0.284%
4	1.301	72	81	96	rVB2	111493	169917	17.36%	1.288%
5	1.465	125	132	147	rVB	206409	228941	23.39%	1.735%
6	1.561	156	162	175	rVB	93925	113251	11.57%	0.858%
7	1.947	273	282	299	rVB	648536	885296	90.45%	6.710%
8	2.101	320	330	342	rBV	154619	216927	22.16%	1.644%
9	2.175	346	353	356	rBV	16951	21902	2.24%	0.166%
10	2.288	379	388	402	rVB	17142	27380	2.80%	0.208%
11	2.690	498	513	515	rBV2	6600	12293	1.26%	0.093%
12	2.761	531	535	566	rVB5	12105	33094	3.38%	0.251%
13	2.953	585	595	610	rVB7	6833	14978	1.53%	0.114%
14	3.285	684	698	714	rBV2	14994	37239	3.80%	0.282%
15	3.915	872	894	950	rBV4	195302	975776	99.70%	7.395%
16	4.346	1011	1028	1080	rBV4	261867	874622	89.36%	6.629%
17	5.043	1227	1245	1255	rBV2	282683	697196	71.24%	5.284%
18	5.294	1312	1323	1357	rBV2	126956	332613	33.98%	2.521%
19	5.613	1411	1422	1459	rBV	271834	570374	58.28%	4.323%
20	6.066	1550	1563	1593	rBV	178795	382506	39.08%	2.899%
21	6.989	1839	1850	1870	rBV	61671	121450	12.41%	0.920%
22	7.313	1940	1951	1966	rBV	240627	424312	43.35%	3.216%
23	7.391	1968	1975	1997	rVB	27166	59402	6.07%	0.450%
24	7.625	2038	2048	2066	rBV2	36645	69793	7.13%	0.529%
25	7.937	2132	2145	2154	rBV4	14556	28063	2.87%	0.213%
26	8.088	2178	2192	2225	rBV2	480795	978719	100.00%	7.418%
27	8.230	2225	2236	2252	rVV	42771	103553	10.58%	0.785%
28	8.725	2376	2390	2401	rBV2	13355	26028	2.66%	0.197%
29	8.854	2418	2430	2433	rBV	435928	660885	67.53%	5.009%
30	8.879	2434	2438	2461	rVB	584047	885416	90.47%	6.711%
31	9.018	2473	2481	2484	rBV2	15087	20120	2.06%	0.152%
32	9.140	2509	2519	2534	rVB9	22532	53127	5.43%	0.403%
33	9.545	2637	2645	2663	rBV3	17317	33753	3.45%	0.256%
34	9.706	2681	2695	2703	rBV3	8589	16945	1.73%	0.128%
35	9.931	2753	2765	2780	rBV	59134	94147	9.62%	0.714%
36	10.181	2835	2843	2844	rBV6	11730	12534	1.28%	0.095%

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37	10.217	2844	2854	2871	rVB	166456	276408	28.24%	2.095%
38	10.352	2886	2896	2904	rBV2	44063	75935	7.76%	0.576%
39	10.452	2916	2927	2938	rVB4	58048	95219	9.73%	0.722%
40	10.545	2948	2956	2958	rBV3	14019	18012	1.84%	0.137%
41	10.693	2986	3002	3007	rBV4	15144	30356	3.10%	0.230%
42	10.735	3007	3015	3027	rVB3	41938	65652	6.71%	0.498%
43	10.915	3063	3071	3077	rBV	39647	59337	6.06%	0.450%
44	10.950	3077	3082	3090	rVB	29021	38473	3.93%	0.292%
45	11.085	3112	3124	3139	rVB2	357507	571988	58.44%	4.335%
46	11.191	3151	3157	3164	rVB4	15589	21478	2.19%	0.163%
47	11.246	3164	3174	3194	rBV2	469678	849064	86.75%	6.435%
48	11.336	3194	3202	3213	rVB	36055	56644	5.79%	0.429%
49	11.529	3252	3262	3276	rBV3	119716	223207	22.81%	1.692%
50	11.622	3281	3291	3322	rVB	351920	612717	62.60%	4.644%
51	11.744	3322	3329	3337	rBV5	7892	13922	1.42%	0.106%
52	11.924	3372	3385	3398	rVB3	12655	32655	3.34%	0.247%
53	12.014	3404	3413	3429	rBV2	46491	82562	8.44%	0.626%
54	12.114	3437	3444	3465	rVB2	24761	54399	5.56%	0.412%
55	12.236	3474	3482	3491	rVB7	12133	20323	2.08%	0.154%
56	12.316	3496	3507	3520	rBV3	27259	53232	5.44%	0.403%
57	12.413	3530	3537	3546	rVV	23619	32734	3.34%	0.248%
58	12.464	3546	3553	3560	rVB5	11332	17145	1.75%	0.130%
59	12.664	3603	3615	3625	rBV9	23486	45624	4.66%	0.346%
60	12.725	3626	3634	3647	rVB6	12207	24458	2.50%	0.185%
61	12.886	3669	3684	3694	rBV4	92878	167637	17.13%	1.271%
62	13.050	3727	3735	3745	rVB7	12805	23572	2.41%	0.179%
63	13.107	3747	3753	3762	rVV10	10278	17748	1.81%	0.135%
64	13.162	3762	3770	3780	rVB10	7645	13024	1.33%	0.099%
65	13.265	3796	3802	3813	rBV10	9368	17629	1.80%	0.134%
66	13.503	3866	3876	3886	rBV	66234	104629	10.69%	0.793%
67	14.631	4216	4227	4240	rBV2	33970	53782	5.50%	0.408%
68	14.760	4257	4267	4275	rBV7	9937	15722	1.61%	0.119%
69	14.815	4275	4284	4294	rVV2	23742	36660	3.75%	0.278%
70	15.802	4581	4591	4601	rBV8	8418	16456	1.68%	0.125%
71	16.307	4740	4748	4758	rBV4	19432	29745	3.04%	0.225%

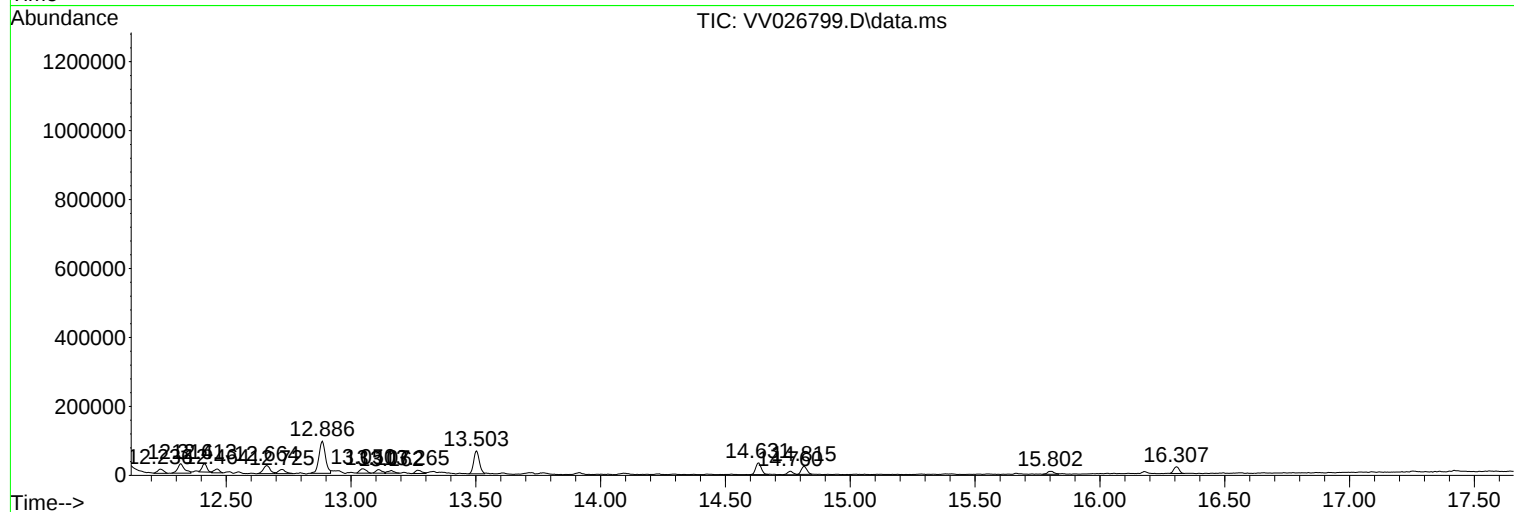
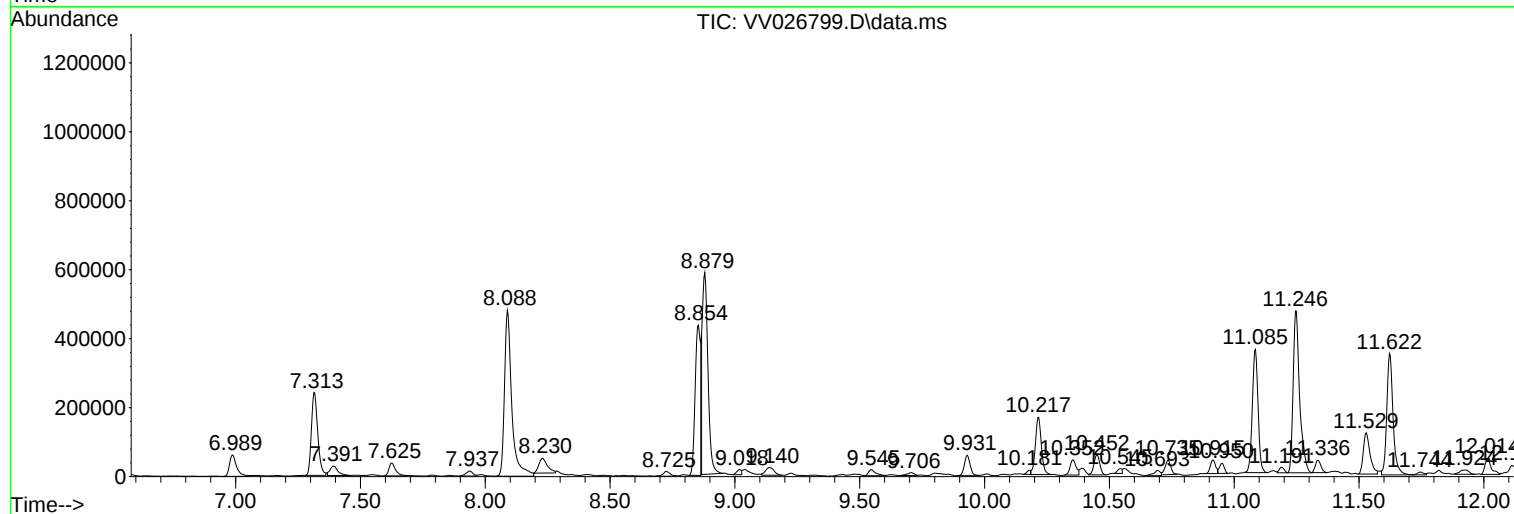
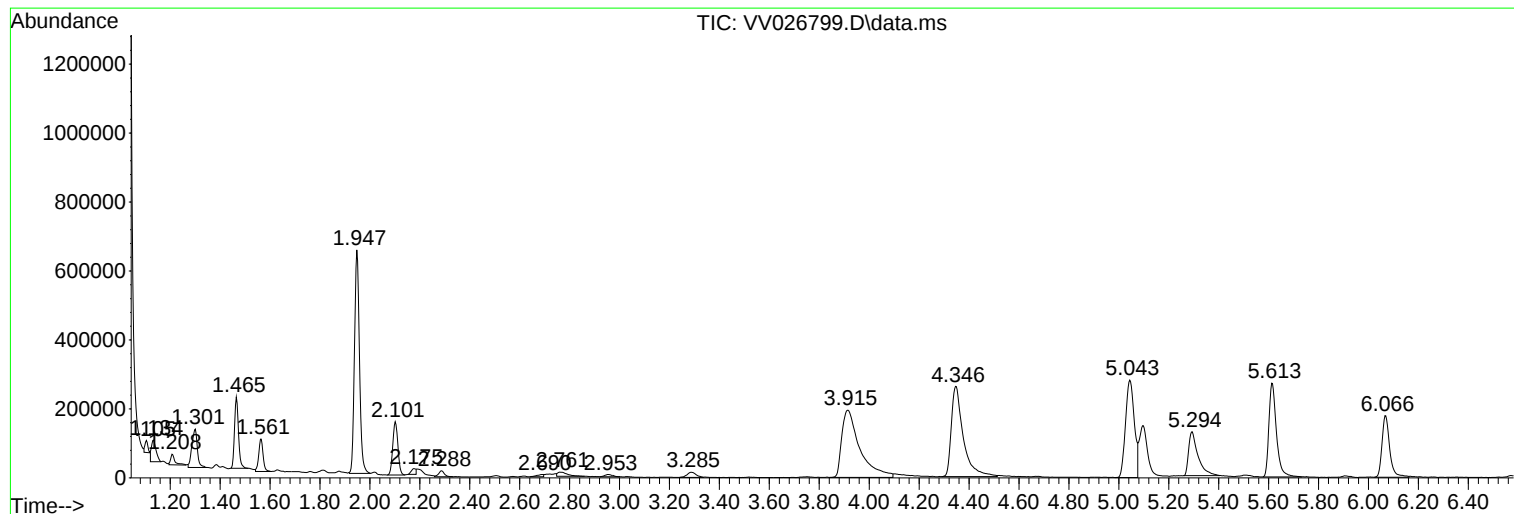
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Instrument :
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ClientSampleId :
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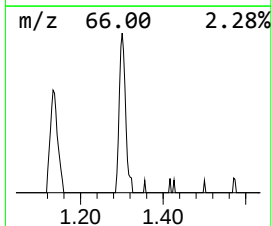
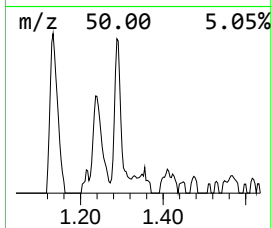
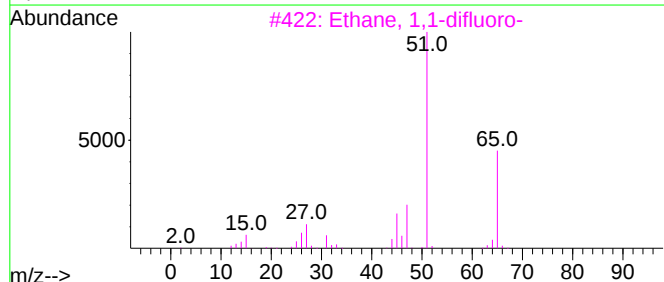
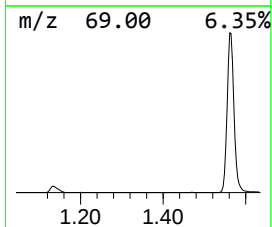
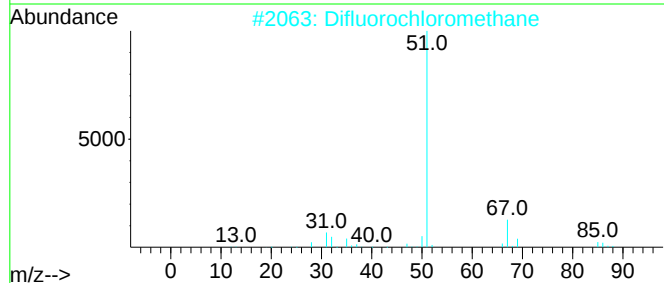
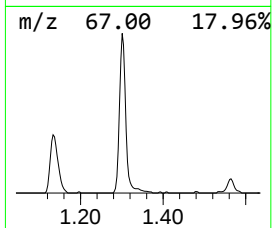
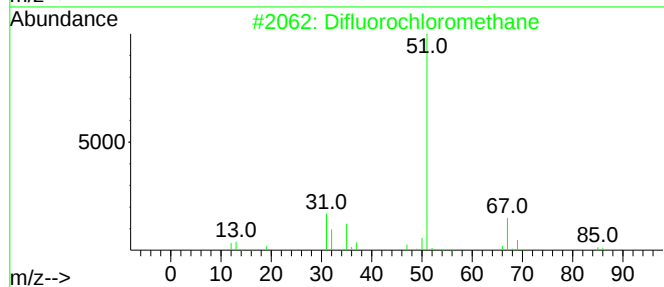
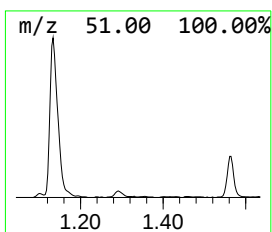
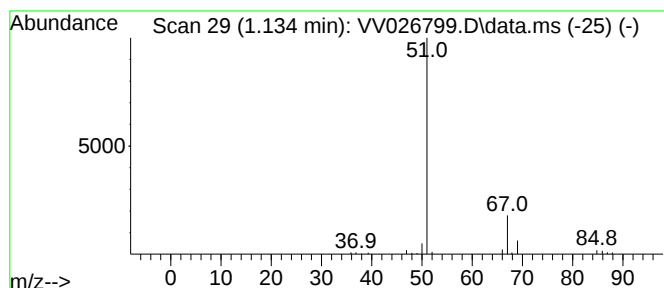
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Difluorochloromethane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.134	0.70 ug/L	79563	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Difluorochloromethane	86	CHClF2	000075-45-6	90
2			Difluorochloromethane	86	CHClF2	000075-45-6	11
3			Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	7
4			Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	5
5			Propiolonitrile	51	C3HN	001070-71-9	3



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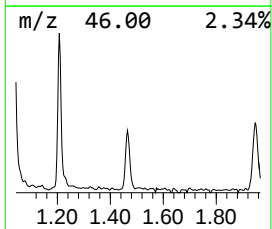
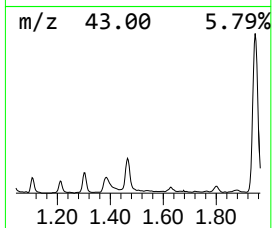
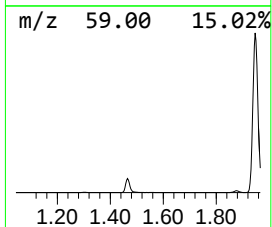
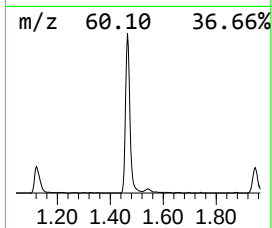
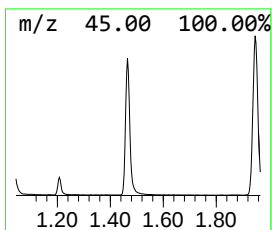
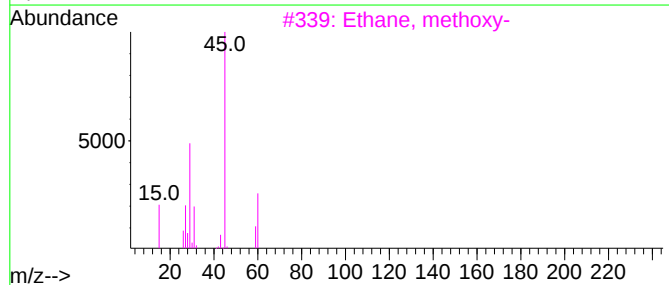
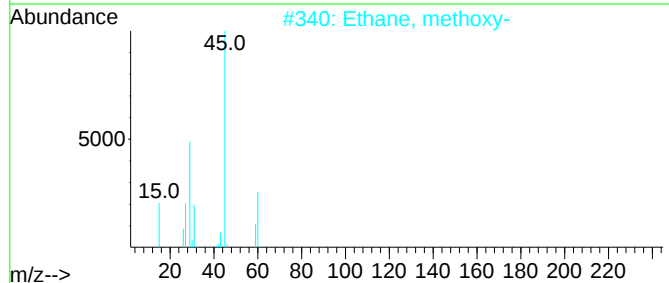
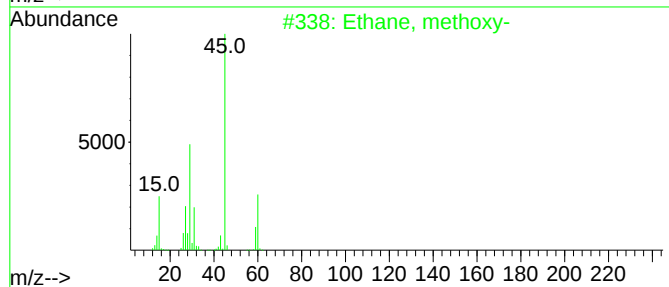
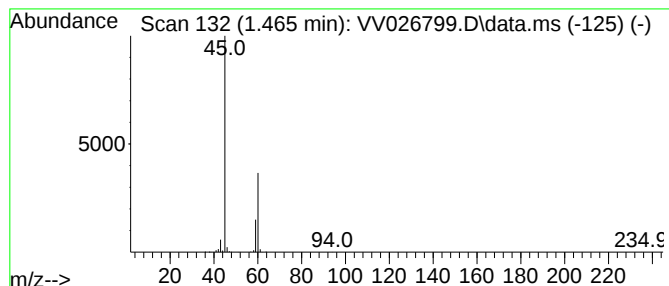
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Ethane, methoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.465	2.01 ug/L	228941	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, methoxy-	60	C3H8O	000540-67-0	78
2			Ethane, methoxy-	60	C3H8O	000540-67-0	72
3			Ethane, methoxy-	60	C3H8O	000540-67-0	72
4			Ethane, methoxy-	60	C3H8O	000540-67-0	56
5			Hydrazine, ethyl-	60	C2H8N2	000624-80-6	7



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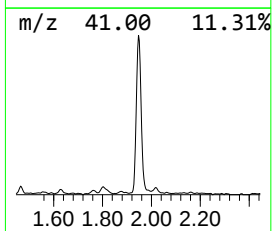
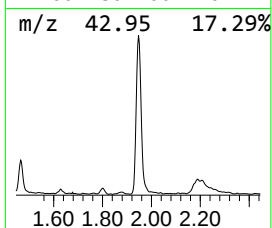
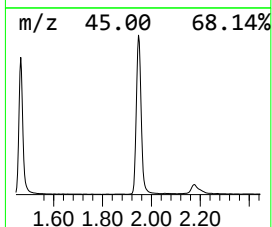
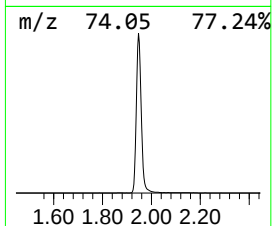
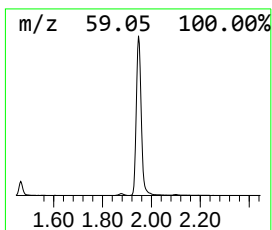
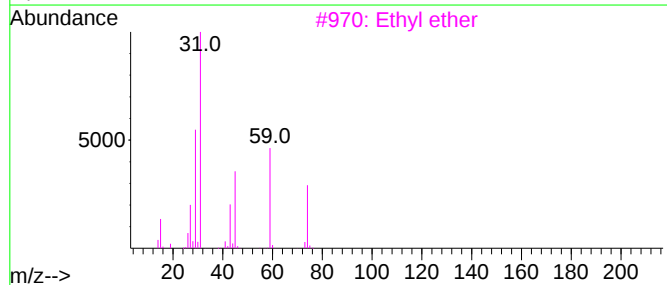
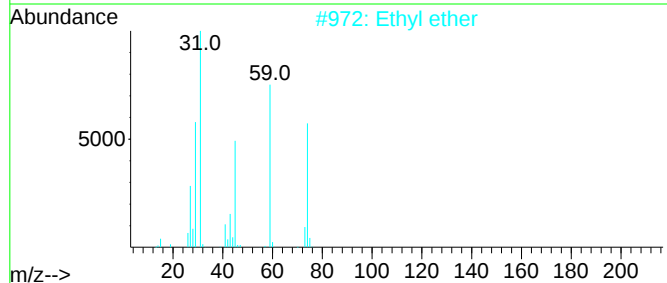
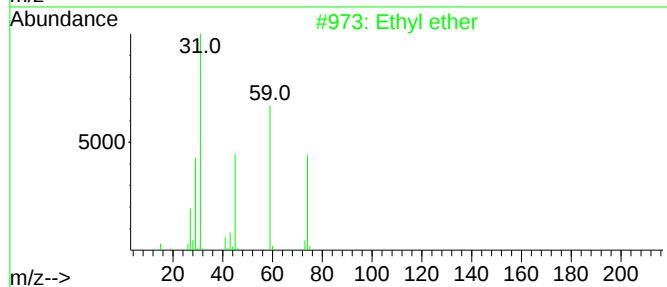
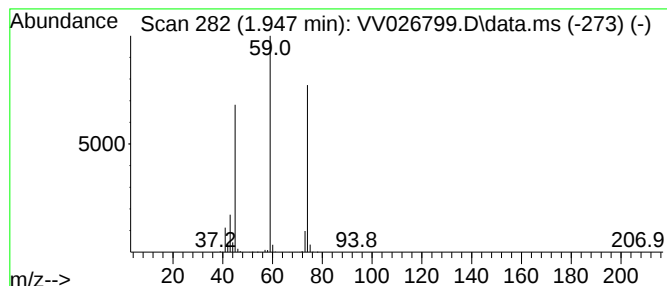
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Ethyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.947	7.76 ug/L	885296	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether			74	C4H10O	000060-29-7	91
2	Ethyl ether			74	C4H10O	000060-29-7	90
3	Ethyl ether			74	C4H10O	000060-29-7	83
4	Ethane, 1,2-diethoxy-			118	C6H14O2	000629-14-1	74
5	Ethyl ether			74	C4H10O	000060-29-7	72



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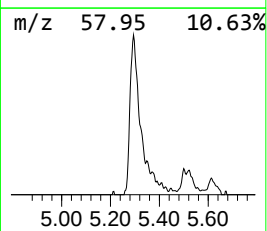
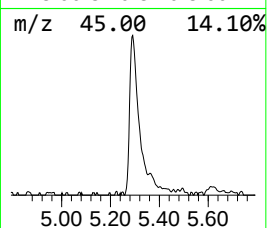
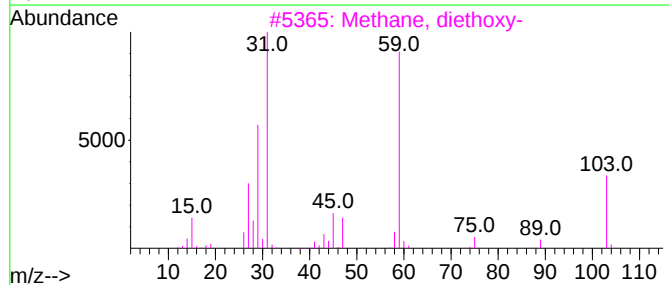
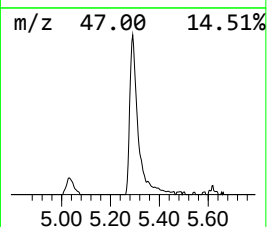
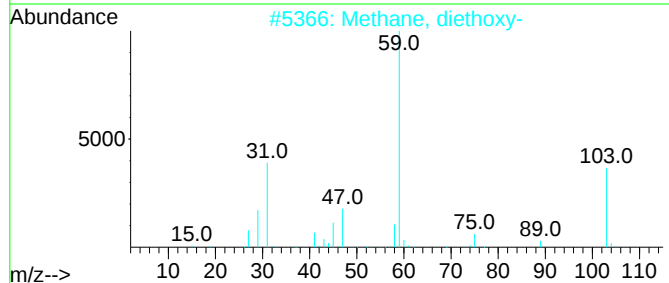
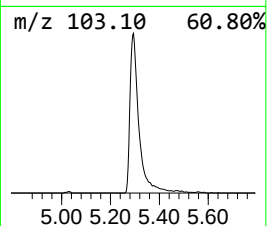
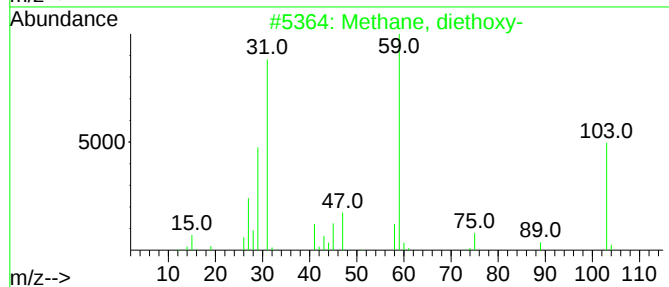
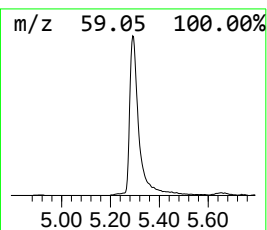
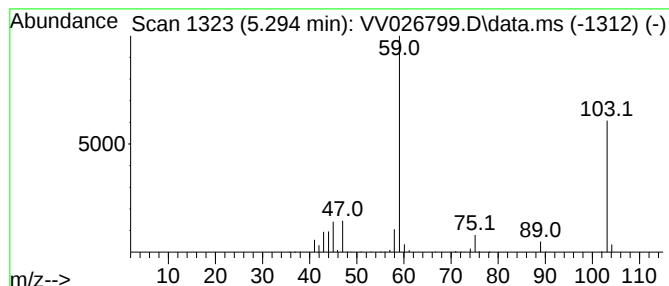
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Methane, diethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.294	2.92 ug/L	332613	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methane, diethoxy-	104	C5H12O2	000462-95-3	91
2			Methane, diethoxy-	104	C5H12O2	000462-95-3	78
3			Methane, diethoxy-	104	C5H12O2	000462-95-3	59
4			Propane, 1,2-dimethoxy-	104	C5H12O2	007778-85-0	47
5			Tetraglyme	222	C10H22O5	000143-24-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
H0B09

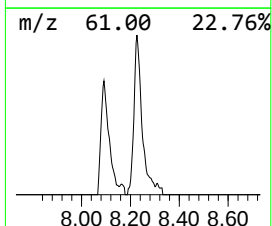
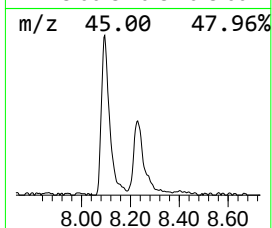
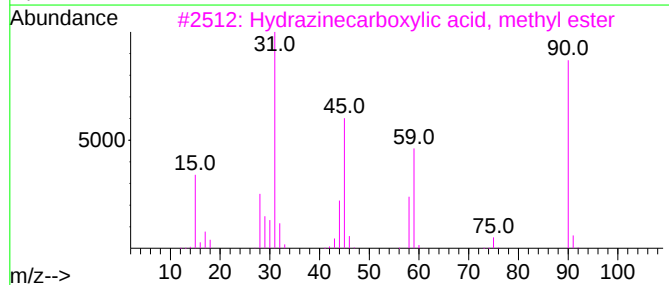
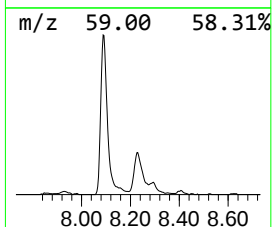
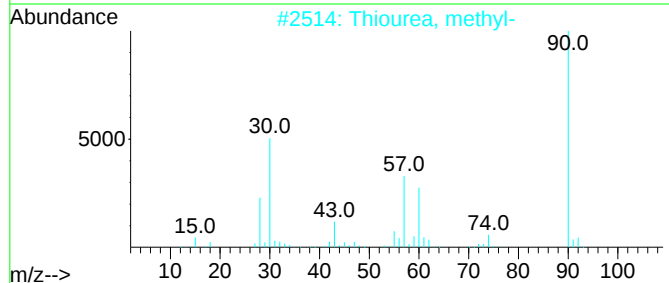
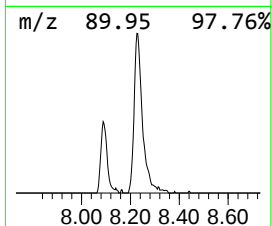
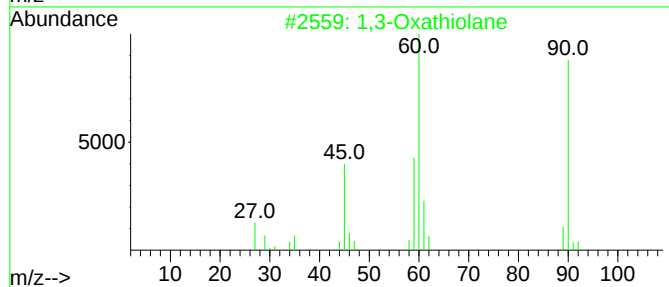
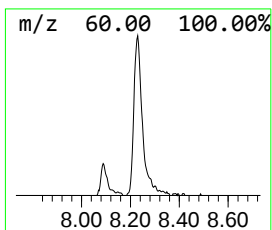
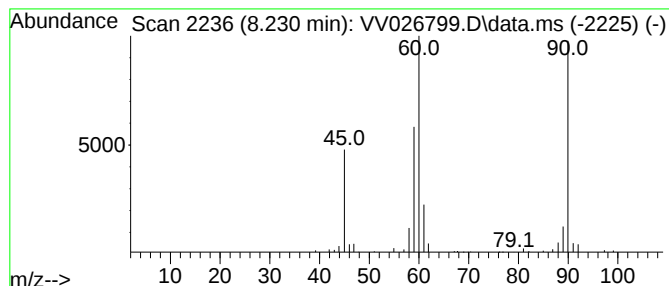
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.230	0.78 ug/L	103553	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Oxathiolane	90	C3H6OS	002094-97-5	91
2			Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
3			Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	45
4			1,3-Oxathiolane	90	C3H6OS	002094-97-5	38
5			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
H0B09

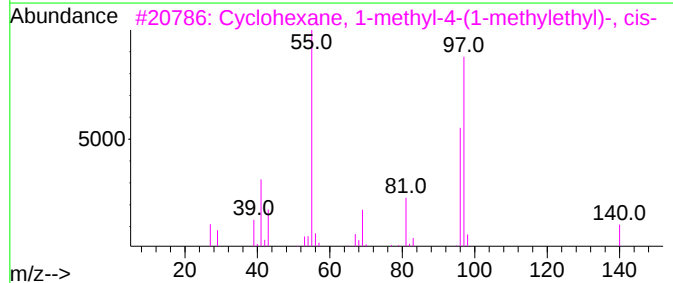
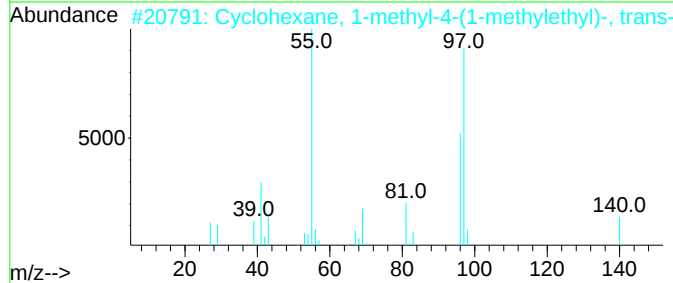
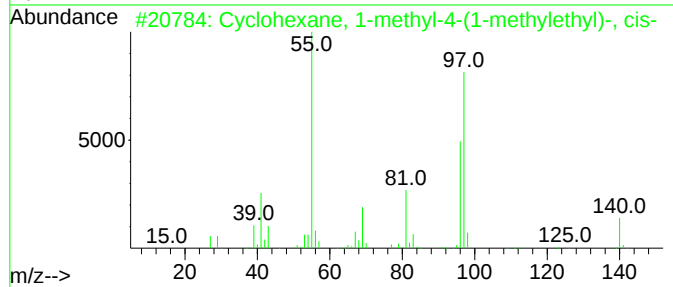
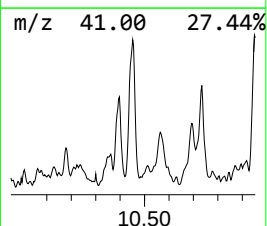
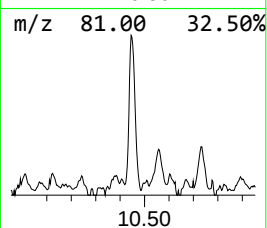
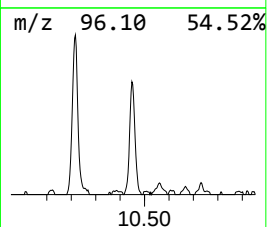
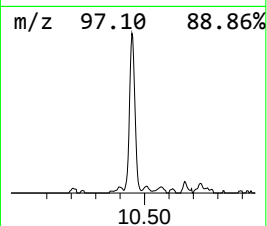
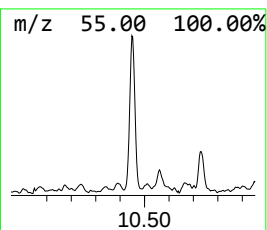
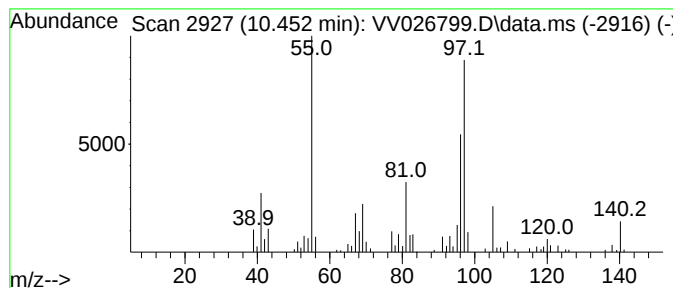
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Cyclic10.452 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.452	0.56 ug/L	95219	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	93
2			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	81
3			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	81
4			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	81
5			1-Methyl-4-(1-methylethyl)-cyclo...	140	C10H20	000099-82-1	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
H0B09

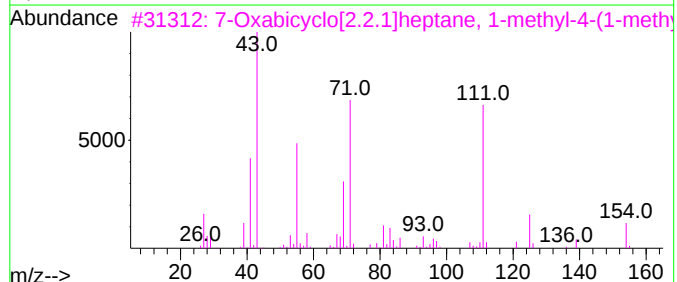
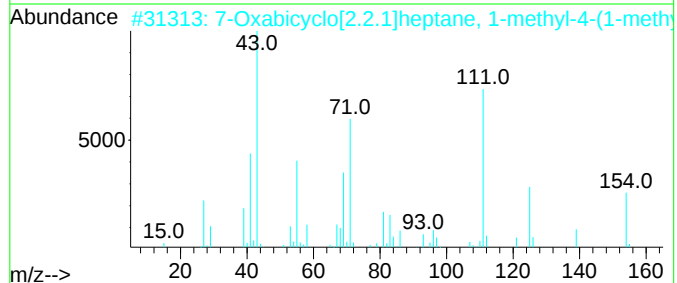
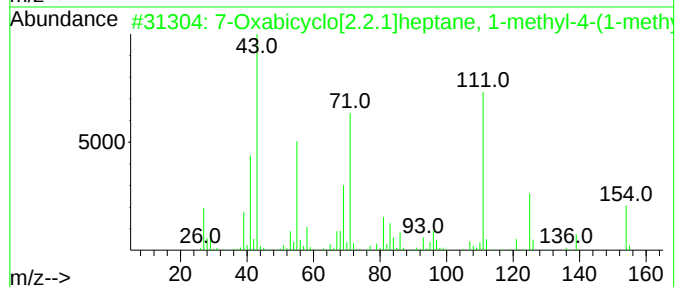
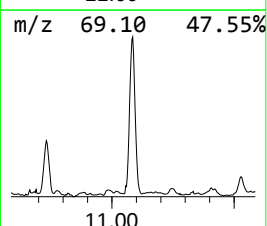
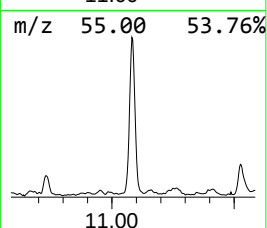
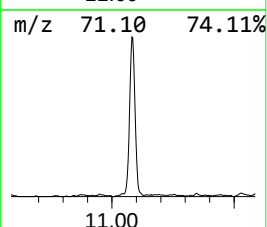
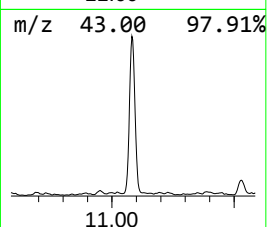
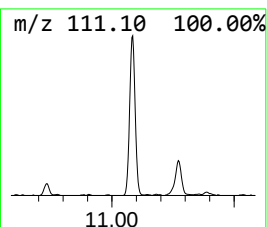
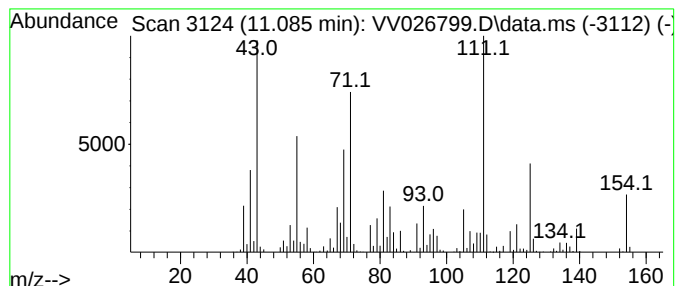
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 7-Oxabicyclo[2.2.1]heptane,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.085	3.37 ug/L	571988	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	93
2			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	93
3			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	76
4			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	55
5			1-Propanone, 2-methyl-1-[2-(1-me...	154	C10H18O	056259-15-5	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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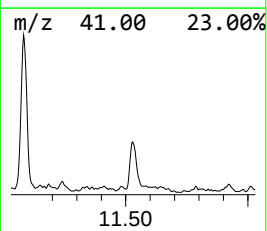
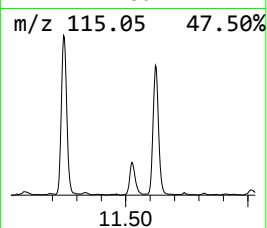
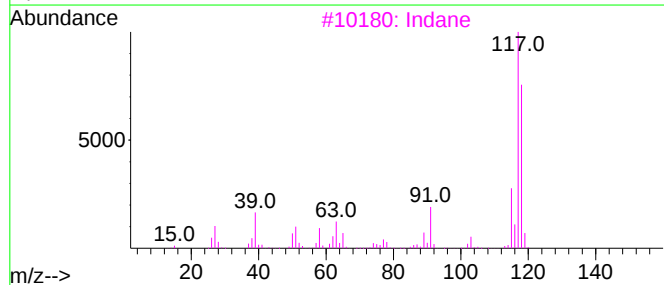
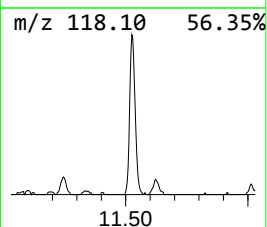
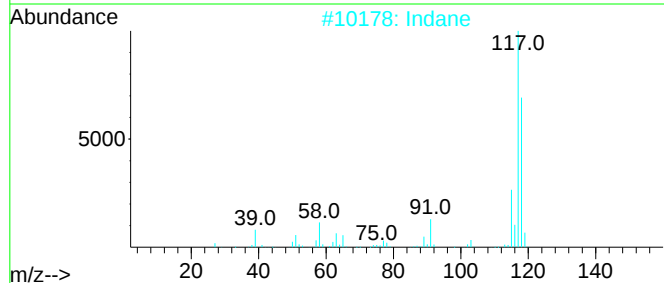
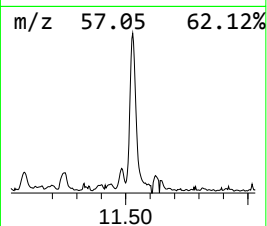
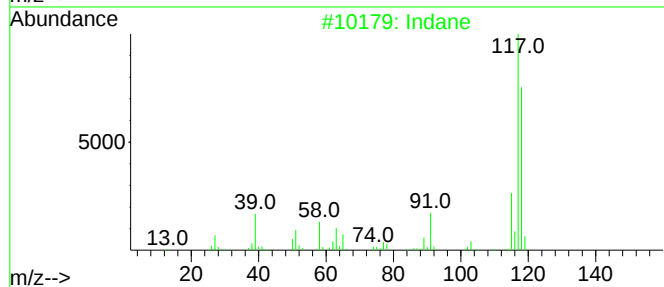
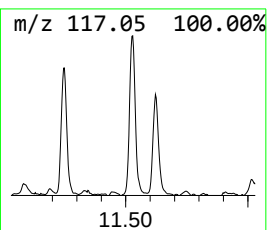
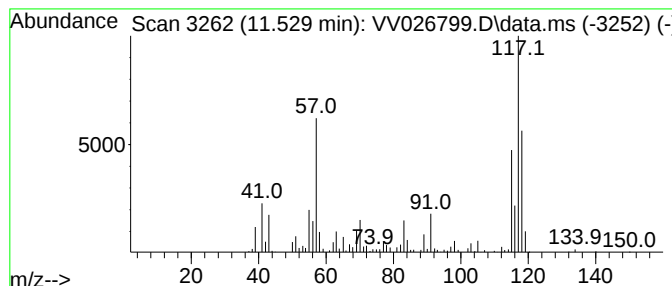
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	1.31 ug/L	223207	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	95
2	Indane			118	C9H10	000496-11-7	93
3	Indane			118	C9H10	000496-11-7	92
4	Indane			118	C9H10	000496-11-7	70
5	Benzene, 2-propenyl-			118	C9H10	000300-57-2	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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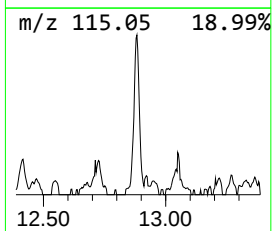
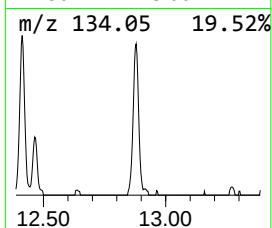
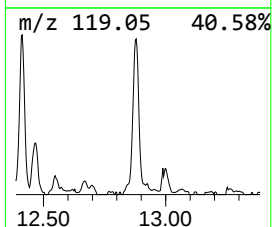
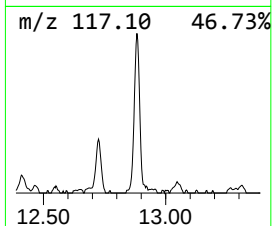
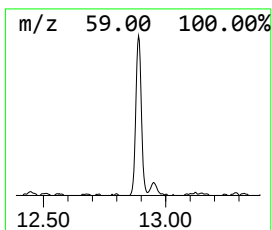
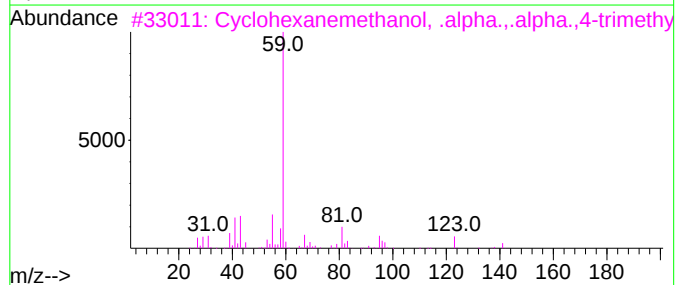
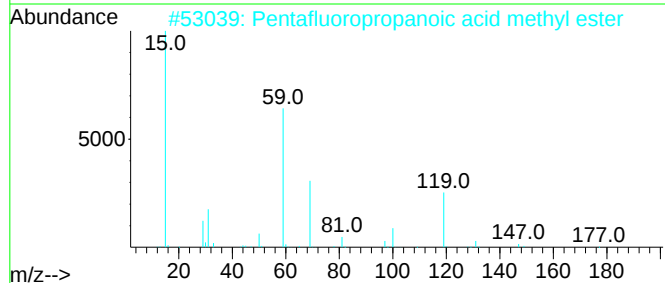
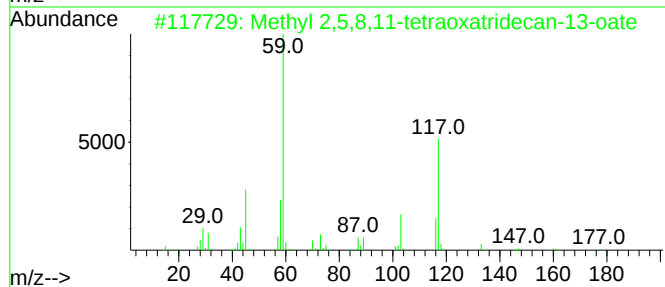
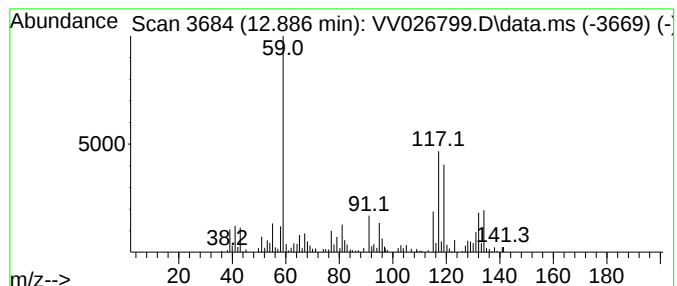
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	0.99 ug/L	167637	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	25
2			Pentafluoropropanoic acid methyl...	178	C4H3F5O2	000378-75-6	25
3			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	22
4			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	22
5			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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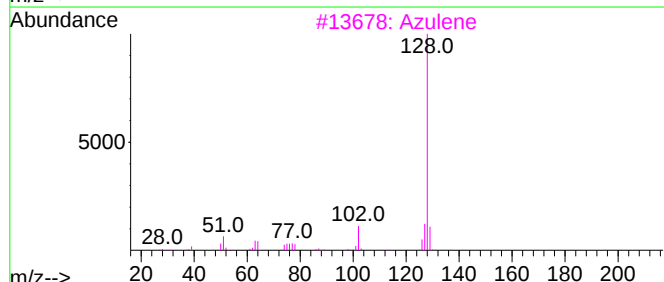
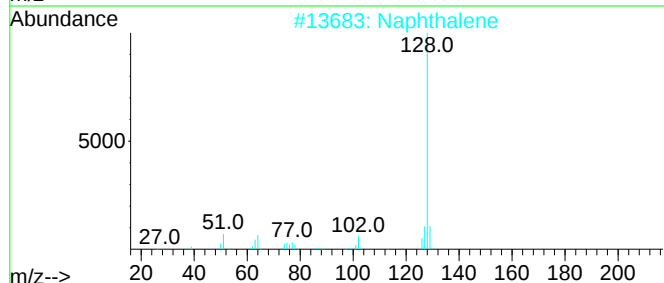
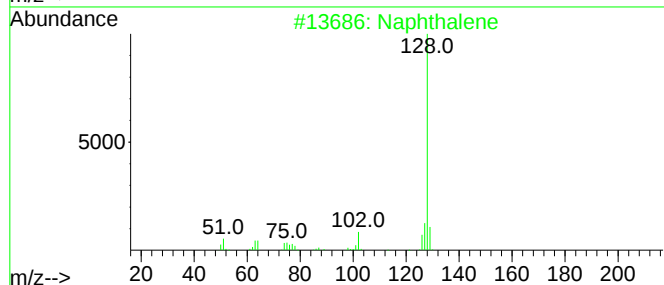
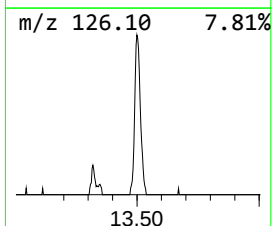
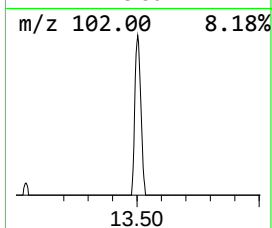
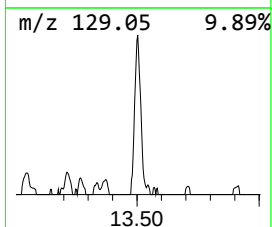
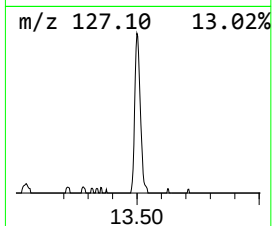
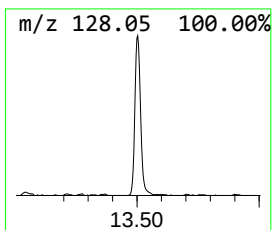
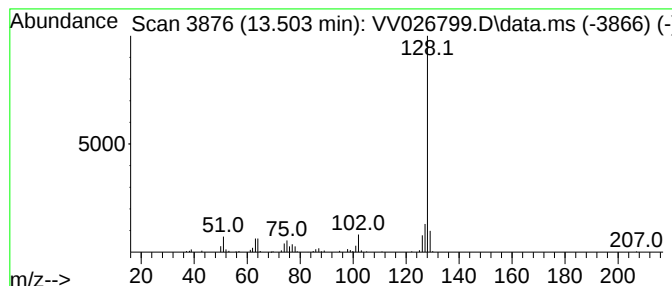
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Azulene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.503	0.62 ug/L	104629	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Azulene	128	C10H8	000275-51-4	93
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV071522\
Data File : VV026799.D
Acq On : 15 Jul 2022 18:37
Operator : SY/MD
Sample : N3710-17
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
H0B09

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Difluorochlorom...	1.134	0.7	ug/L	79563	1	5.616	570374	5.0
Ethane, methoxy-	1.465	2.0	ug/L	228941	1	5.616	570374	5.0
Ethyl ether	1.947	7.8	ug/L	885296	1	5.616	570374	5.0
Methane, diethoxy-	5.294	2.9	ug/L	332613	1	5.616	570374	5.0
1,3-Oxathiolane	8.230	0.8	ug/L	103553	2	8.850	660885	5.0
(DEL) Alkane: C...	10.452	0.6	ug/L	95219	3	11.249	849064	5.0
7-Oxabicyclo[2....	11.085	3.4	ug/L	571988	3	11.249	849064	5.0
Indane	11.529	1.3	ug/L	223207	3	11.249	849064	5.0
unknown-01	12.886	1.0	ug/L	167637	3	11.249	849064	5.0
Azulene	13.503	0.6	ug/L	104629	3	11.249	849064	5.0