

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
 Data File : VV030362.D
 Acq On : 30 Mar 2023 16:19
 Operator : SY/MD
 Sample : 02092-21
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 CB994

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\SFAMVTR032323WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV030362.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.089	11	15	18	rBV	42163	38401	6.96%	0.513%
2	1.114	18	23	43	rVB	143419	188946	34.23%	2.524%
3	1.278	65	74	86	rVB2	75563	114401	20.73%	1.528%
4	1.388	102	108	118	rVB	54024	64922	11.76%	0.867%
5	1.536	147	154	170	rVB	62304	76907	13.93%	1.027%
6	1.770	224	227	237	rVB5	6376	8069	1.46%	0.108%
7	2.063	309	318	336	rVB	107993	155674	28.21%	2.079%
8	2.163	336	349	359	rBV2	5366	14526	2.63%	0.194%
9	2.243	370	374	383	rVB5	4586	6522	1.18%	0.087%
10	2.896	566	577	589	rBV4	10395	23634	4.28%	0.316%
11	3.220	663	678	696	rBV2	52820	126109	22.85%	1.684%
12	3.831	853	868	910	rBV3	44663	217702	39.45%	2.908%
13	4.265	987	1003	1017	rBV	92929	237957	43.12%	3.178%
14	4.593	1096	1105	1117	rVB	5431	12007	2.18%	0.160%
15	4.966	1203	1221	1247	rBV2	198691	551909	100.00%	7.372%
16	5.252	1299	1310	1315	rBV6	13675	26443	4.79%	0.353%
17	5.439	1357	1368	1387	rVB9	7126	18296	3.32%	0.244%
18	5.542	1387	1400	1424	rBV	152940	323483	58.61%	4.321%
19	5.998	1522	1542	1558	rBV	124998	272312	49.34%	3.637%
20	6.921	1817	1829	1850	rBV2	51702	103492	18.75%	1.382%
21	7.252	1920	1932	1947	rBV	214355	383594	69.50%	5.124%
22	7.326	1949	1955	1962	rVB3	8095	12751	2.31%	0.170%
23	7.561	2019	2028	2045	rBV2	32414	62394	11.31%	0.833%
24	7.883	2118	2128	2137	rBV6	6940	14258	2.58%	0.190%
25	8.037	2163	2176	2206	rBV	226590	494682	89.63%	6.607%
26	8.188	2215	2223	2243	rVB6	20325	48207	8.73%	0.644%
27	8.792	2399	2411	2442	rBV2	233897	482422	87.41%	6.444%
28	8.956	2453	2462	2471	rBV6	5372	8395	1.52%	0.112%
29	9.014	2472	2480	2482	rBV5	6457	8148	1.48%	0.109%
30	9.079	2492	2500	2513	rVB4	18772	36639	6.64%	0.489%
31	9.442	2607	2613	2619	rVB7	5232	6724	1.22%	0.090%
32	9.487	2620	2627	2642	rVB2	16394	29742	5.39%	0.397%
33	9.876	2738	2748	2762	rBV	43062	70243	12.73%	0.938%
34	9.953	2764	2772	2778	rVV	3966	5916	1.07%	0.079%
35	10.085	2803	2813	2818	rBV4	24748	40569	7.35%	0.542%
36	10.127	2818	2826	2830	rVV2	46202	75461	13.67%	1.008%

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37	10.162	2830	2837	2851	rVB	97475	158491	28.72%	2.117%
38	10.345	2884	2894	2904	rVV6	27872	67211	12.18%	0.898%
39	10.397	2905	2910	2921	rVB9	22398	37594	6.81%	0.502%
40	10.513	2933	2946	2966	rVB	153960	269580	48.85%	3.601%
41	10.680	2991	2998	3008	rBV	17283	28907	5.24%	0.386%
42	10.808	3031	3038	3044	rBV4	8612	11975	2.17%	0.160%
43	10.860	3046	3054	3062	rVB	13672	20672	3.75%	0.276%
44	10.998	3089	3097	3103	rBV6	6336	10211	1.85%	0.136%
45	11.034	3103	3108	3117	rVB4	15682	21609	3.92%	0.289%
46	11.133	3132	3139	3147	rVB6	7866	11593	2.10%	0.155%
47	11.194	3147	3158	3176	rVB	272298	444231	80.49%	5.933%
48	11.281	3178	3185	3195	rVB3	20352	30204	5.47%	0.403%
49	11.349	3195	3206	3214	rBV10	10196	19418	3.52%	0.259%
50	11.397	3214	3221	3230	rVV	27734	40848	7.40%	0.546%
51	11.474	3234	3245	3262	rVV	163196	288234	52.22%	3.850%
52	11.571	3264	3275	3293	rVB	278850	457763	82.94%	6.114%
53	11.693	3305	3313	3321	rBV5	10462	16350	2.96%	0.218%
54	11.963	3388	3397	3413	rBV3	35143	61597	11.16%	0.823%
55	12.059	3413	3427	3437	rBV2	27970	55461	10.05%	0.741%
56	12.204	3464	3472	3478	rBV9	4489	6983	1.27%	0.093%
57	12.329	3498	3511	3518	rBV	103545	174874	31.69%	2.336%
58	12.413	3529	3537	3547	rBV3	18029	28041	5.08%	0.375%
59	12.612	3588	3599	3607	rVV3	10093	18807	3.41%	0.251%
60	12.673	3607	3618	3626	rVB4	15690	24415	4.42%	0.326%
61	12.828	3652	3666	3679	rBV3	47021	82226	14.90%	1.098%
62	13.107	3746	3753	3764	rVB6	16358	28812	5.22%	0.385%
63	13.159	3766	3769	3778	rVB7	6333	8040	1.46%	0.107%
64	13.217	3778	3787	3797	rBV7	10666	20987	3.80%	0.280%
65	13.448	3846	3859	3869	rBV	45321	83573	15.14%	1.116%
66	13.567	3887	3896	3904	rBV	5779	11360	2.06%	0.152%
67	13.615	3904	3911	3916	rVV7	7746	13316	2.41%	0.178%
68	13.648	3916	3921	3933	rVV10	13201	26148	4.74%	0.349%
69	13.709	3933	3940	3951	rVB10	7809	16513	2.99%	0.221%
70	13.863	3978	3988	3994	rBV6	10221	18358	3.33%	0.245%
71	14.040	4033	4043	4051	rBV9	7877	17311	3.14%	0.231%
72	14.246	4095	4107	4114	rBV10	8572	18669	3.38%	0.249%
73	14.525	4182	4194	4200	rBV10	10746	17857	3.24%	0.239%
74	14.580	4200	4211	4220	rVV2	14099	27428	4.97%	0.366%
75	14.628	4222	4226	4236	rVB2	6191	8145	1.48%	0.109%
76	14.763	4258	4268	4279	rBV	44019	75953	13.76%	1.014%

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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

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Peak separation: 5

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Title : TRACE VOA SFAM1.0

77	15.612	4523	4532	4545	rBV2	17310	33391	6.05%	0.446%
78	15.760	4568	4578	4585	rBV	23190	37881	6.86%	0.506%
79	15.795	4585	4589	4596	rVB6	10267	13398	2.43%	0.179%
80	16.127	4686	4692	4701	rVB6	5934	9677	1.75%	0.129%
81	16.400	4767	4777	4781	rBV	75955	113857	20.63%	1.521%
82	16.432	4782	4787	4801	rVB	88996	137052	24.83%	1.831%

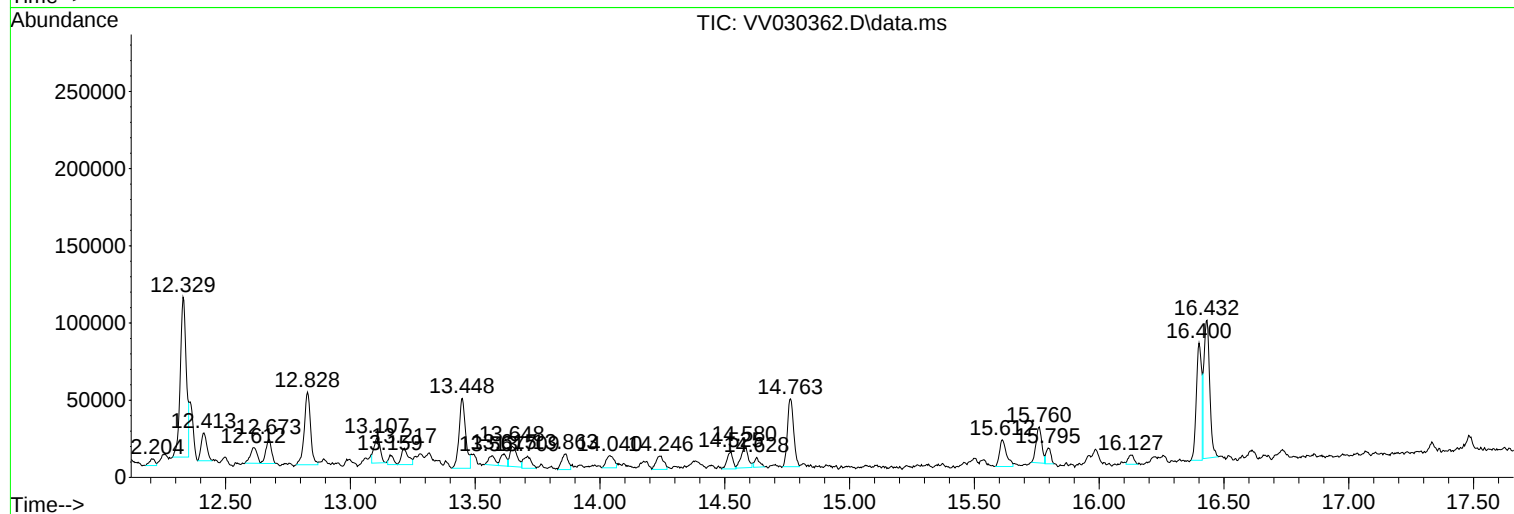
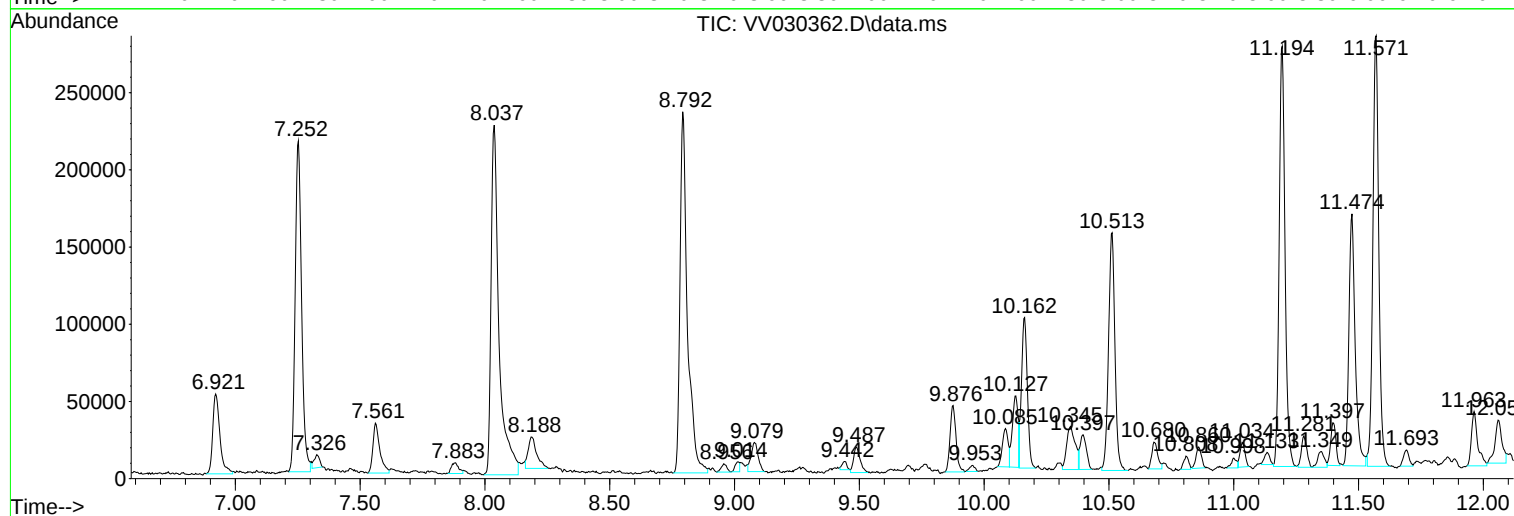
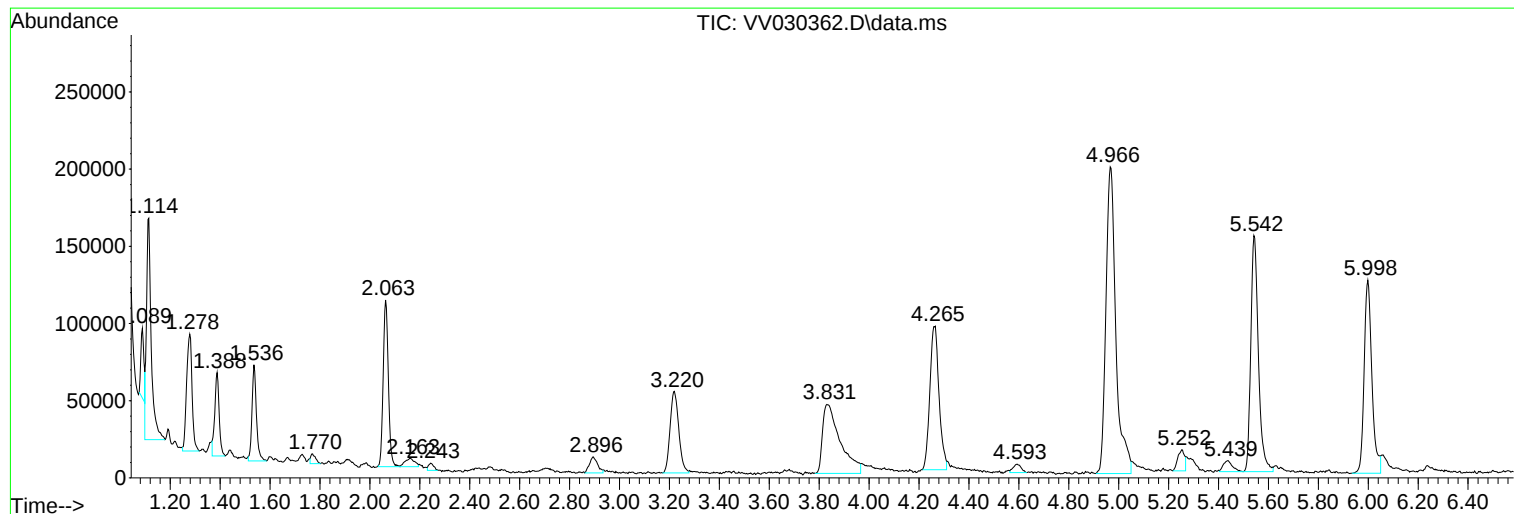
Sum of corrected areas: 7486878

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR032323WMA.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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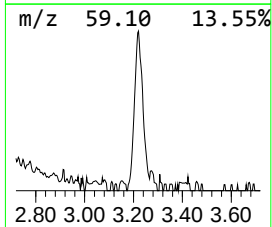
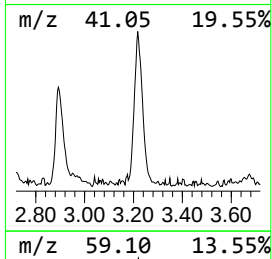
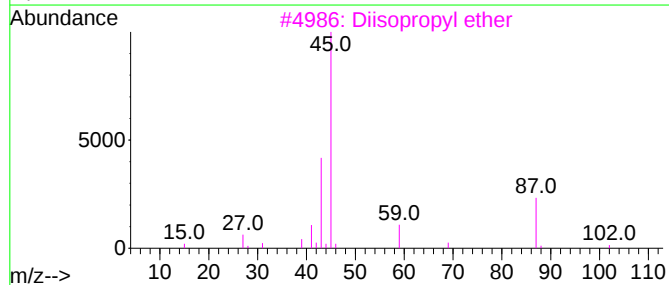
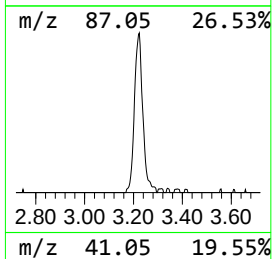
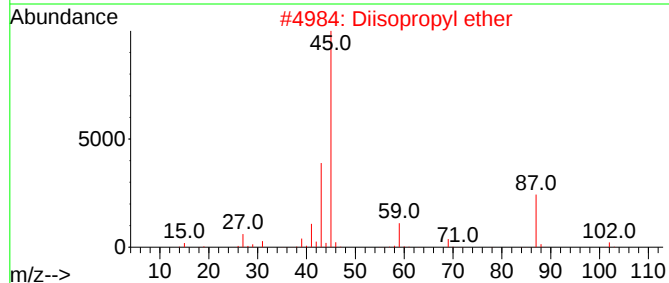
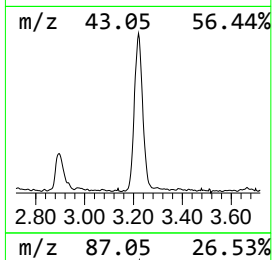
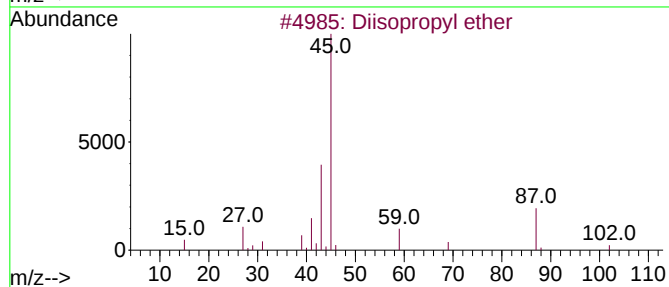
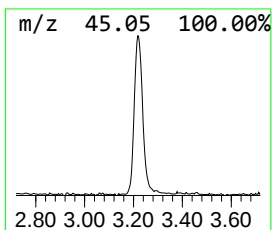
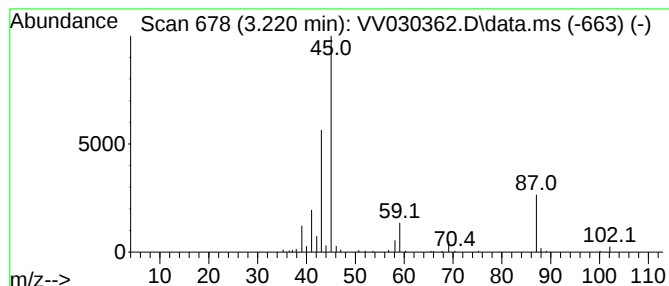
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR032323WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Diisopropyl ether Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.220	1.95 ug/L	126109	1,4-Difluorobenzene	5.542

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisopropyl ether	102	C6H14O	000108-20-3	90
2			Diisopropyl ether	102	C6H14O	000108-20-3	87
3			Diisopropyl ether	102	C6H14O	000108-20-3	86
4			1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	78
5			Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	72



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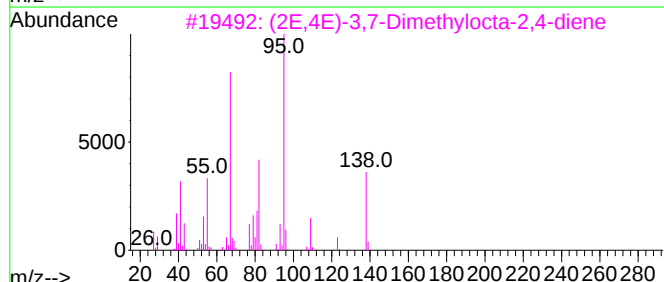
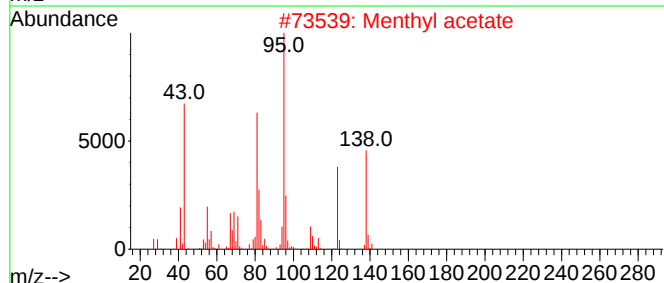
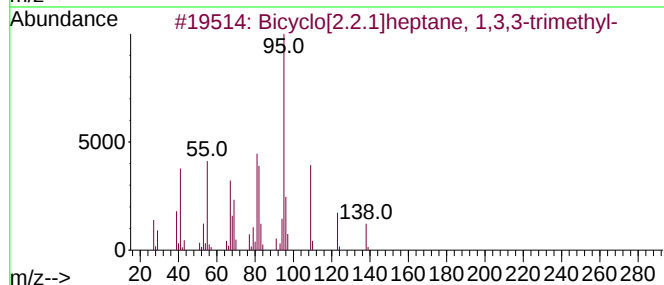
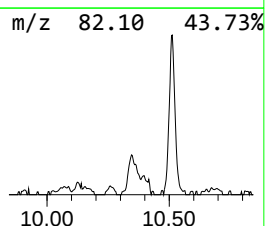
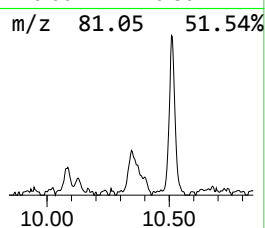
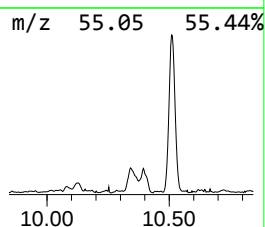
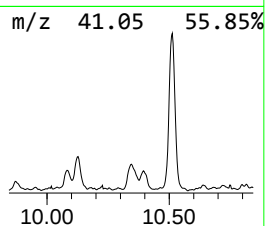
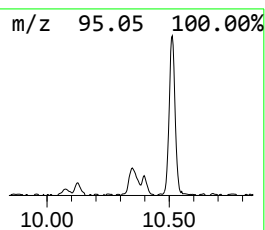
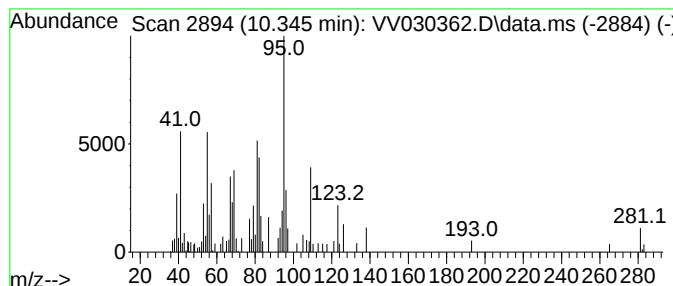
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR032323WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Bicyclo[2.2.1]heptane, 1,3,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.345	0.76 ug/L	67211	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	94
2			Menthyl acetate	198	C12H22O2	000089-48-5	47
3			(2E,4E)-3,7-Dimethylocta-2,4-diene	138	C10H18	006874-39-1	47
4			Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	002778-68-9	43
5			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	38



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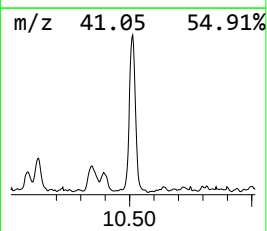
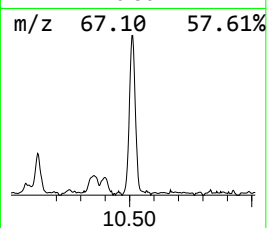
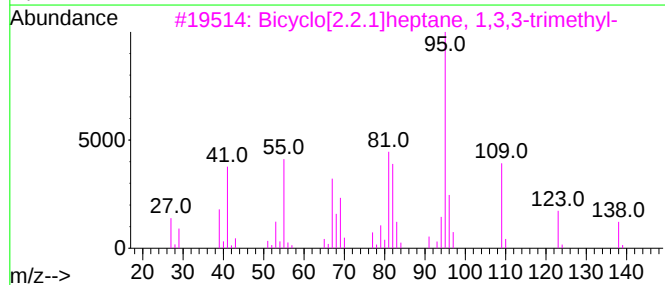
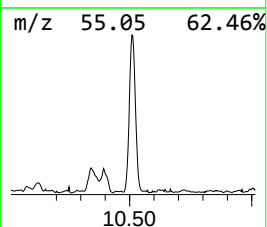
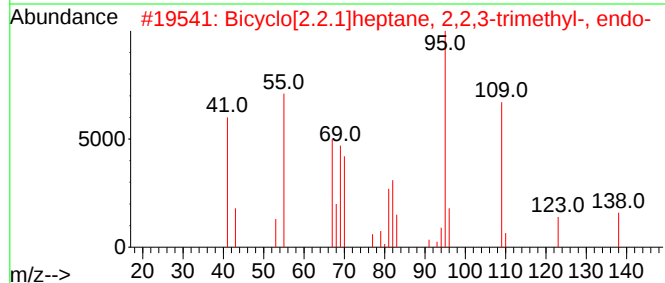
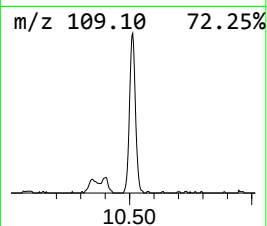
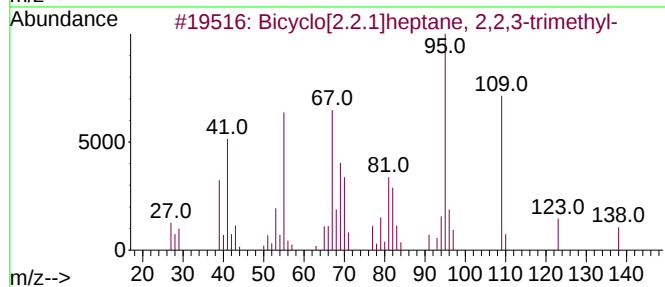
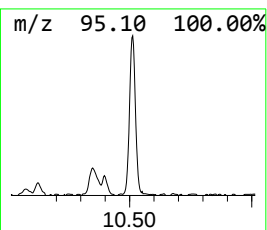
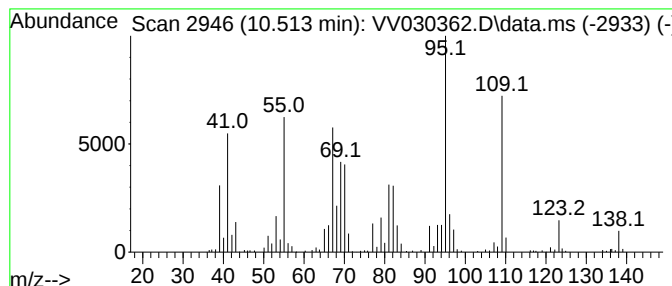
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR032323WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Bicyclo[2.2.1]heptane, 2,2,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.513	3.03 ug/L	269580	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	98
2			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	97
3			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	92
4			Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	87
5			(2E,4E)-3,7-Dimethylocta-2,4-diene	138	C10H18	006874-39-1	53



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ClientSampleId :
CB994

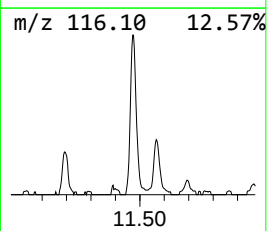
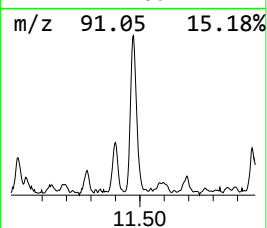
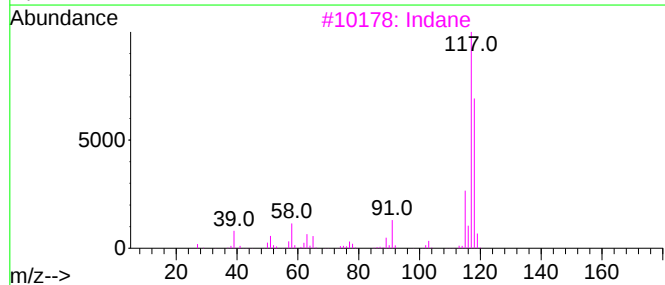
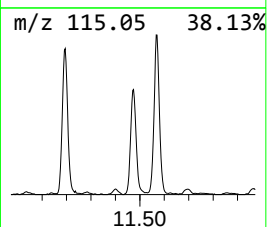
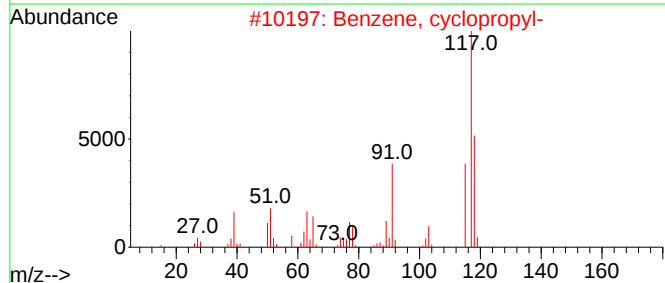
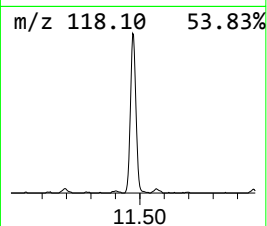
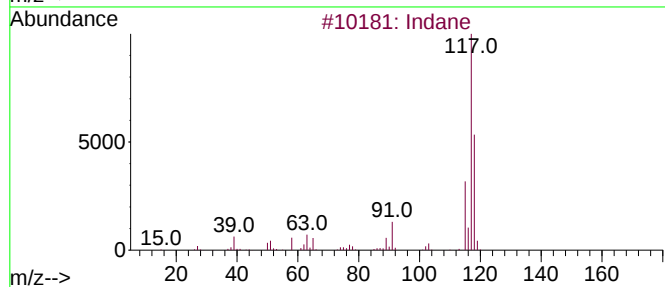
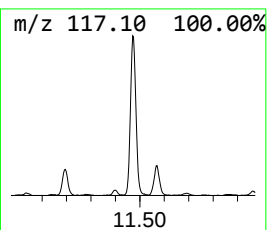
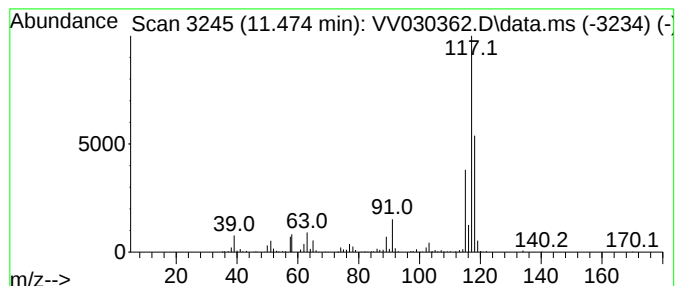
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Quant Title : TRACE VOA SFAM1.0

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.
11.474	3.24 ug/L	288234	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	95
2	Benzene, cyclopropyl-			118	C9H10	000873-49-4	94
3	Indane			118	C9H10	000496-11-7	93
4	Indane			118	C9H10	000496-11-7	90
5	Indane			118	C9H10	000496-11-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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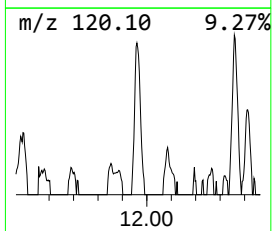
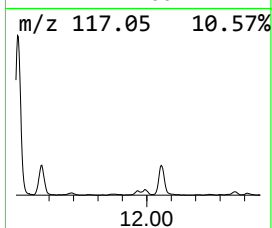
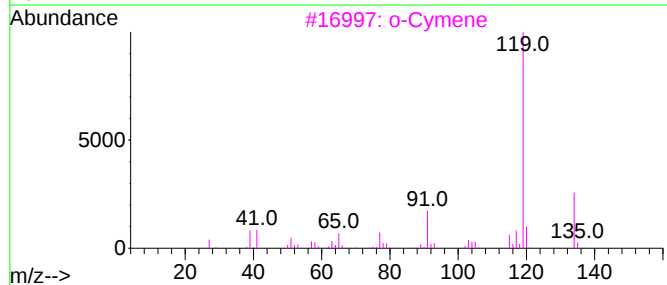
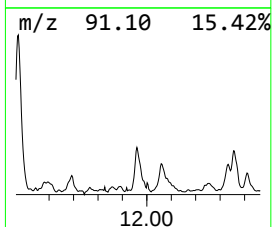
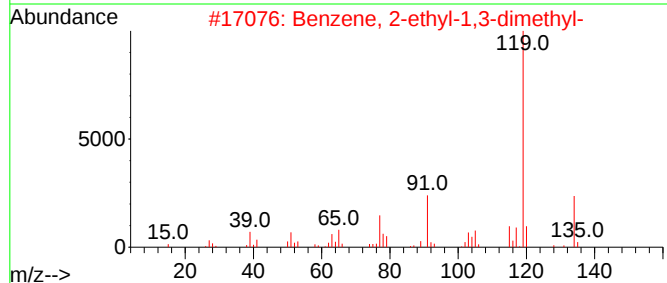
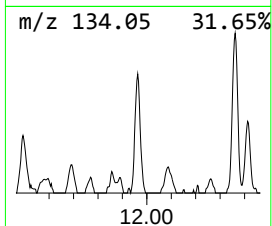
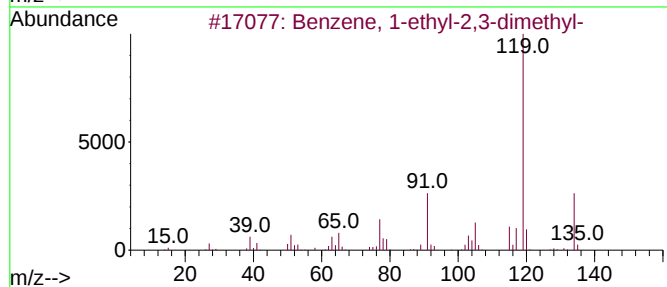
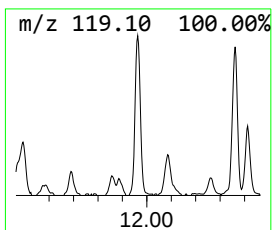
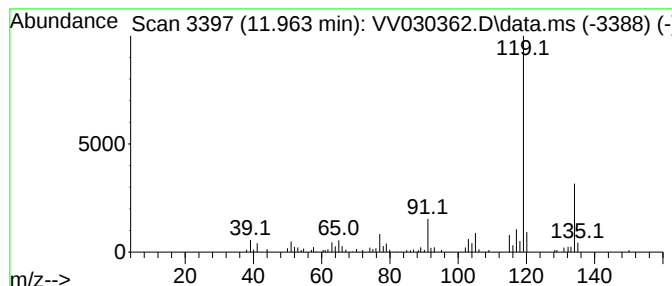
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	0.69 ug/L	61597	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			o-Cymene	134	C10H14	000527-84-4	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
CB994

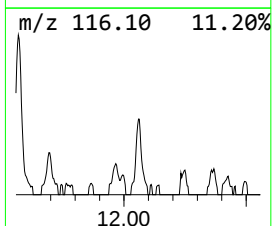
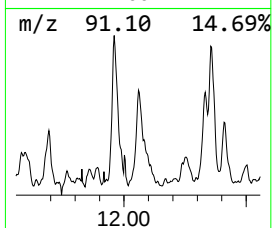
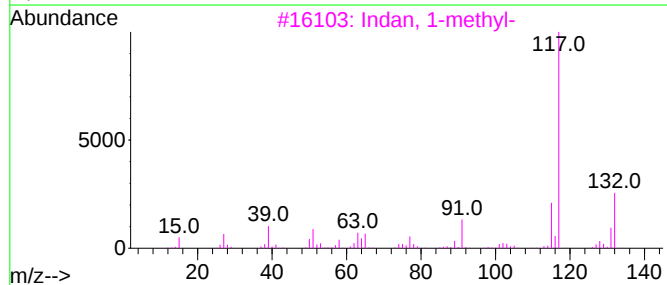
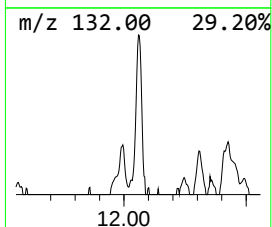
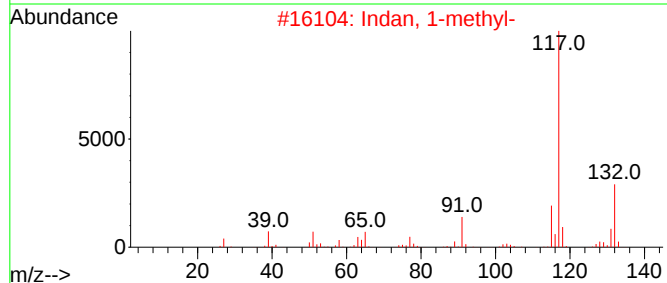
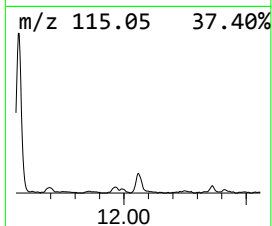
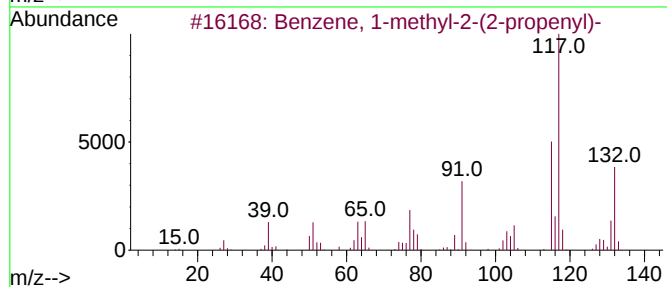
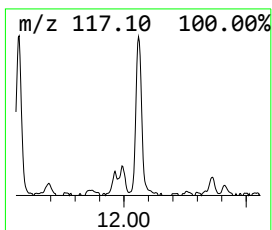
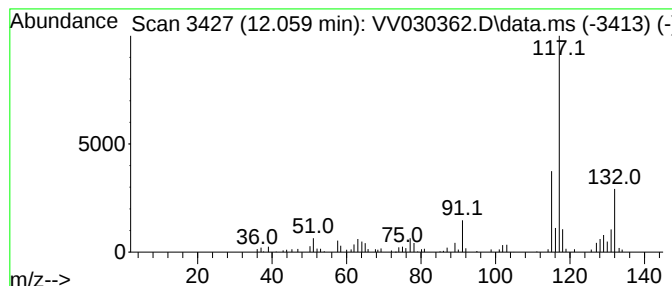
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Benzene, 1-methyl-2-(2-prop... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.059	0.62 ug/L	55461	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2			Indan, 1-methyl-	132	C10H12	000767-58-8	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	81
4			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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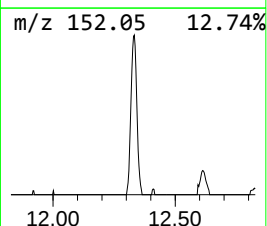
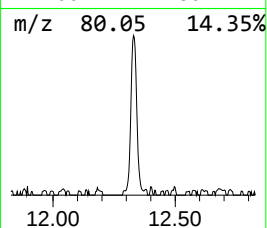
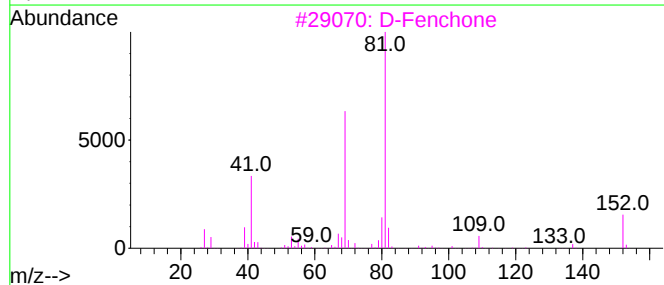
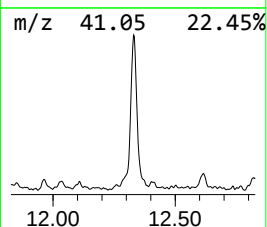
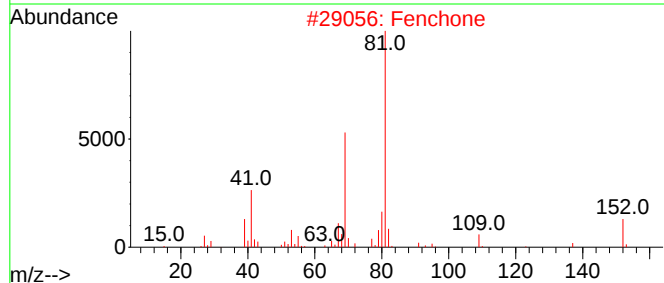
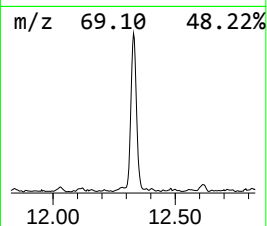
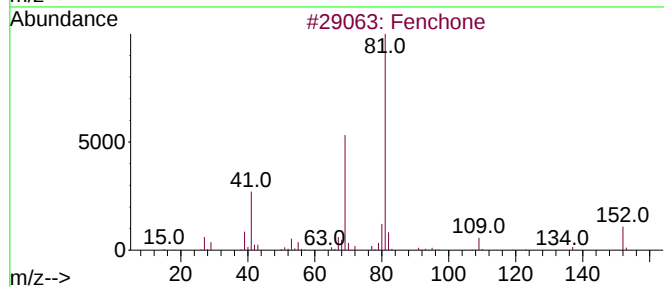
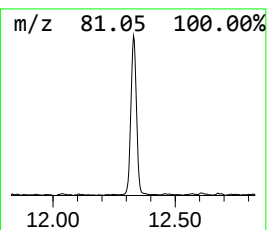
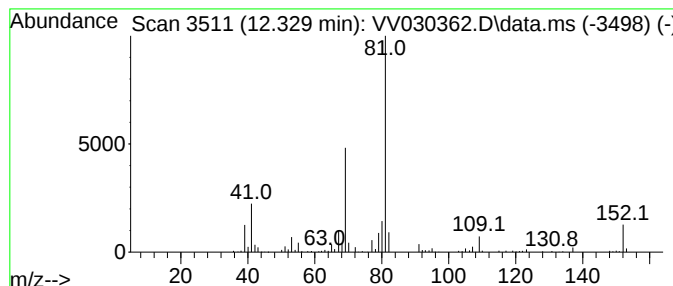
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Fenchone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.329	1.97 ug/L	174874	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchone	152	C10H16O	001195-79-5	95
2			Fenchone	152	C10H16O	001195-79-5	95
3			D-Fenchone	152	C10H16O	004695-62-9	94
4			L-Fenchone	152	C10H16O	007787-20-4	94
5			L-Fenchone	152	C10H16O	007787-20-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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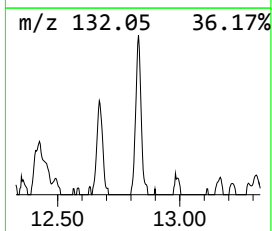
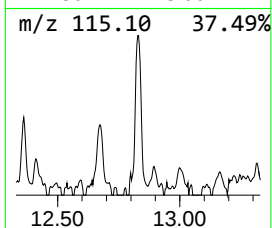
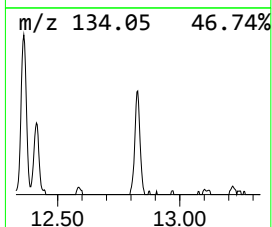
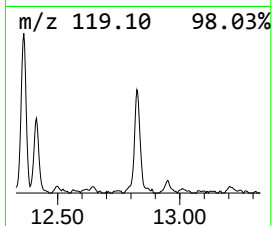
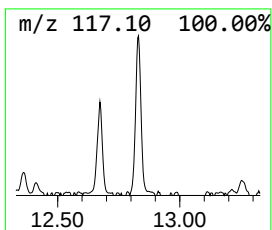
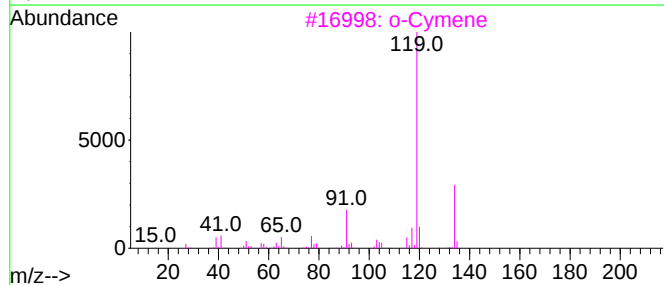
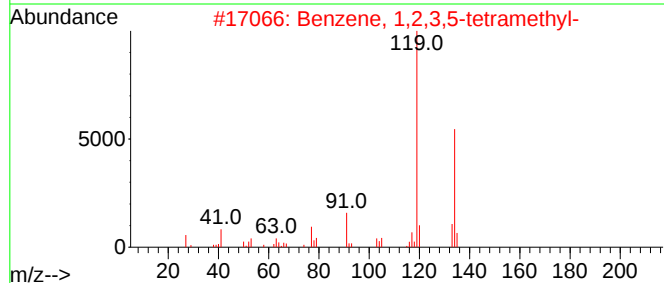
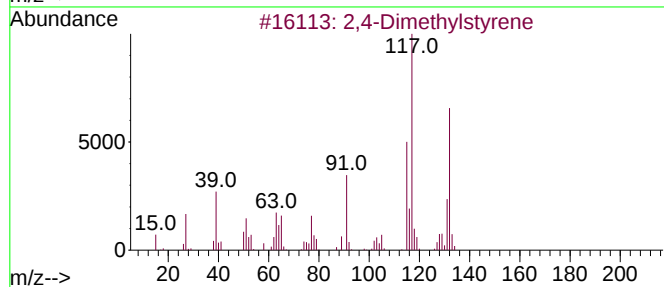
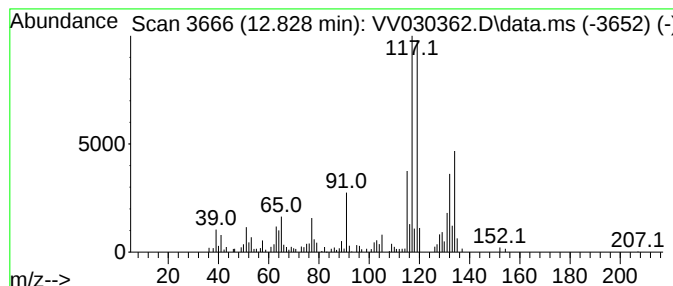
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 2,4-Dimethylstyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	0.93 ug/L	82226	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylstyrene	132	C10H12	002234-20-0	74
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50
3			o-Cymene	134	C10H14	000527-84-4	50
4			p-Cymene	134	C10H14	000099-87-6	50
5			o-Cymene	134	C10H14	000527-84-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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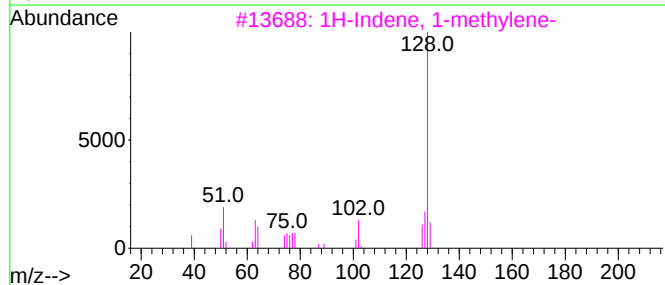
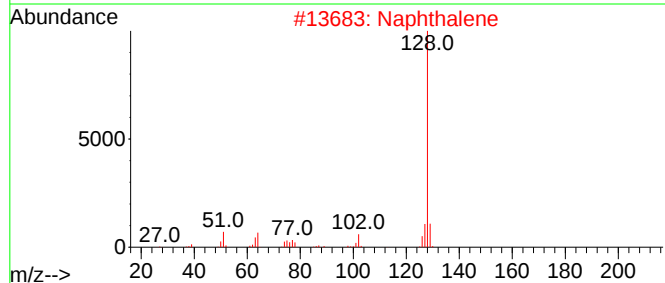
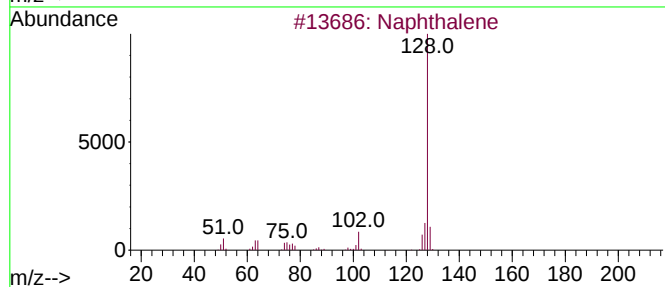
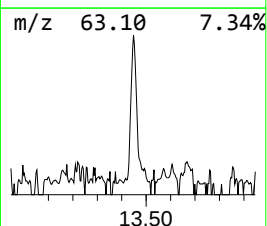
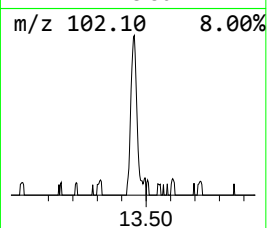
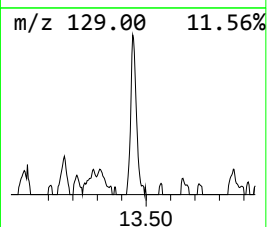
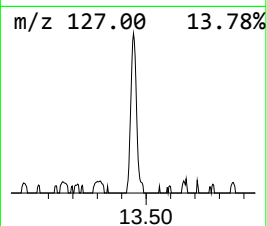
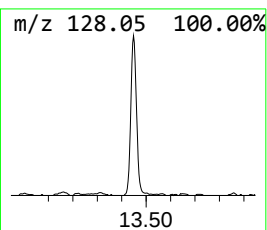
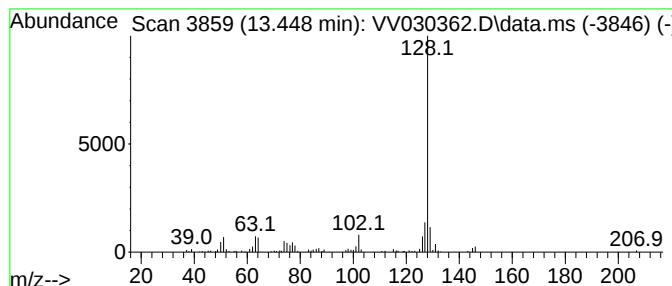
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Azulene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.448	0.94 ug/L	83573	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	94
2			Naphthalene	128	C10H8	000091-20-3	93
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	93
4			Azulene	128	C10H8	000275-51-4	90
5			Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

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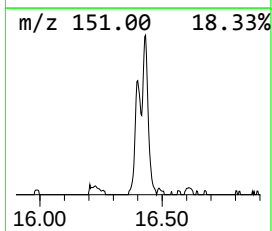
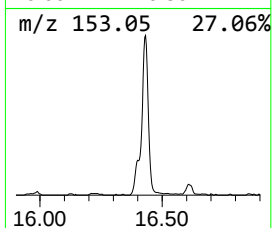
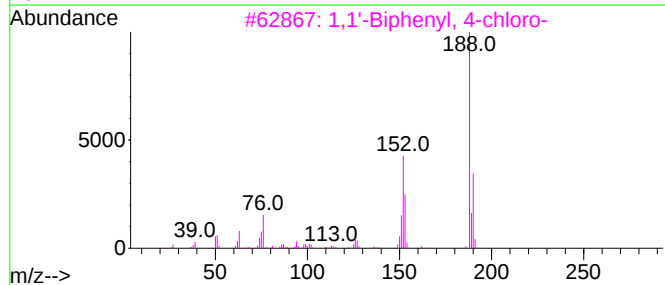
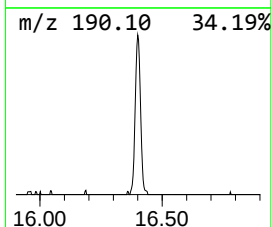
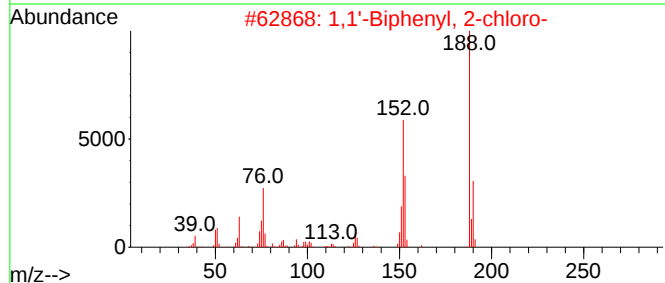
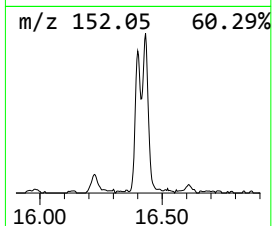
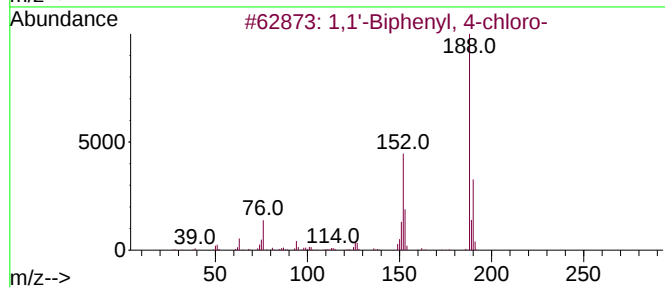
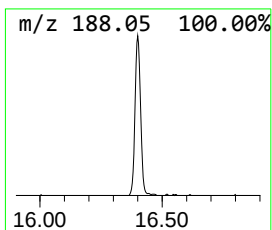
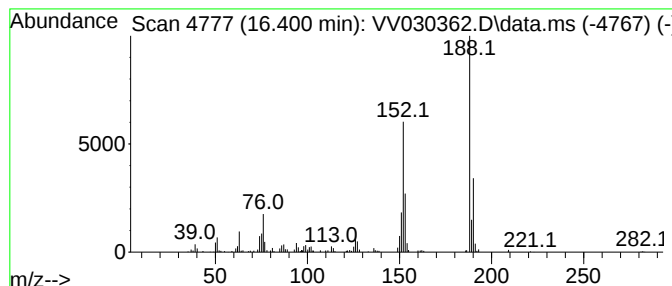
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 1,1'-Biphenyl, 4-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.400	1.28 ug/L	113857	1,4-Dichlorobenzene-d4	11.194

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	98
2		1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	97
3		1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	97
4		3-Chlorobiphenyl	188	C12H9Cl	002051-61-8	96
5		3-Chlorobiphenyl	188	C12H9Cl	002051-61-8	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
Data File : VV030362.D
Acq On : 30 Mar 2023 16:19
Operator : SY/MD
Sample : 02092-21
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
CB994

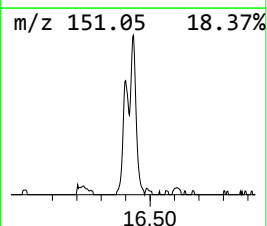
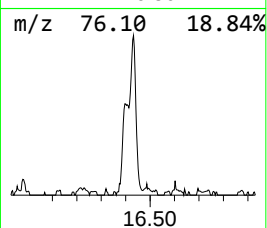
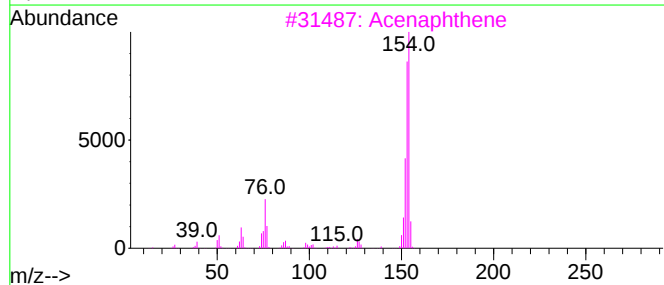
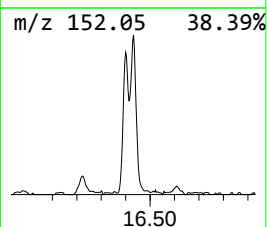
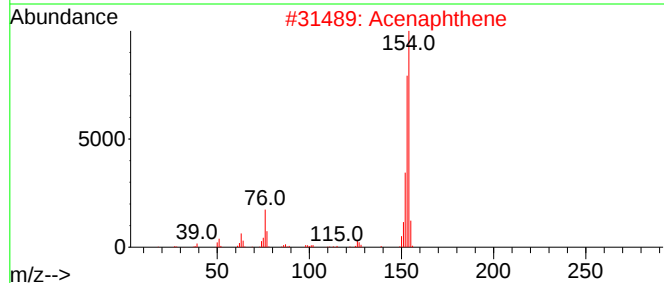
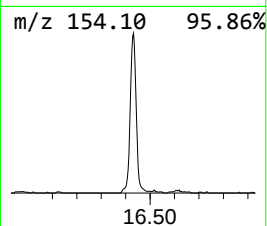
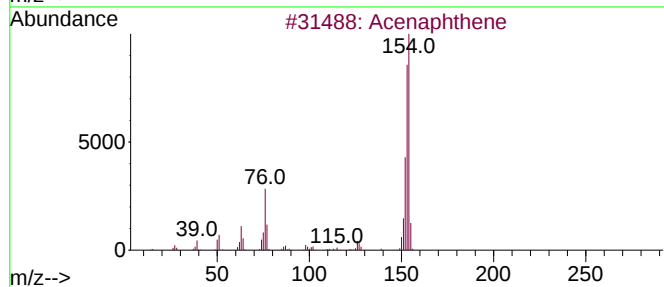
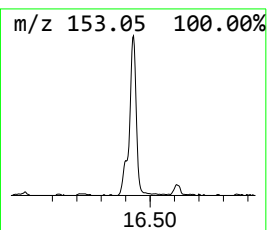
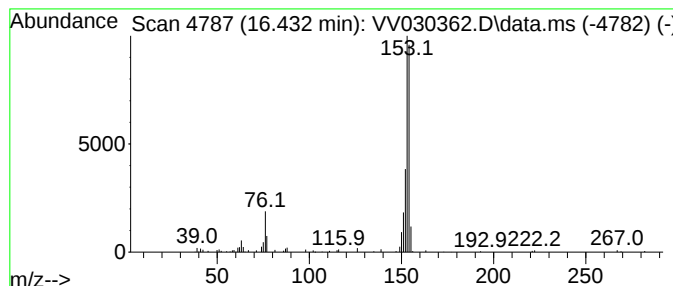
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Naphthalene, 2-ethenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.432	1.54 ug/L	137052	1,4-Dichlorobenzene-d4	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	74
2			Acenaphthene	154	C12H10	000083-32-9	74
3			Acenaphthene	154	C12H10	000083-32-9	72
4			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	59
5			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV033023\
 Data File : VV030362.D
 Acq On : 30 Mar 2023 16:19
 Operator : SY/MD
 Sample : 02092-21
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 CB994

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR032323WMA.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
Diisopropyl ether	3.220	2.0	ug/L	126109	1	5.542	323483	5.0
Bicyclo[2.2.1]h...	10.345	0.8	ug/L	67211	3	11.194	444231	5.0
Bicyclo[2.2.1]h...	10.513	3.0	ug/L	269580	3	11.194	444231	5.0
Indane	11.474	3.2	ug/L	288234	3	11.194	444231	5.0
Benzene, 1-ethy...	11.963	0.7	ug/L	61597	3	11.194	444231	5.0
Benzene, 1-meth...	12.059	0.6	ug/L	55461	3	11.194	444231	5.0
Fenchone	12.329	2.0	ug/L	174874	3	11.194	444231	5.0
2,4-Dimethylsty...	12.828	0.9	ug/L	82226	3	11.194	444231	5.0
Azulene	13.448	0.9	ug/L	83573	3	11.194	444231	5.0
1,1'-Biphenyl, ...	16.400	1.3	ug/L	113857	3	11.194	444231	5.0
Naphthalene, 2-...	16.432	1.5	ug/L	137052	3	11.194	444231	5.0