

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
 Data File : VV023345.D
 Acq On : 10 Nov 2021 20:24
 Operator : SY/MD
 Sample : M4542-01
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGGE7

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\SFAMVTR110421WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV023345.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	75	83	96	rVB2	53720	58635	3.09%	0.543%
2	1.568	156	164	174	rBV	42716	48132	2.54%	0.446%
3	2.108	321	332	346	rBV	77967	115962	6.11%	1.073%
4	3.195	657	670	684	rBV3	11151	23166	1.22%	0.214%
5	3.902	875	890	921	rBV3	70944	270486	14.25%	2.504%
6	4.352	1008	1030	1057	rBV4	72963	212040	11.17%	1.963%
7	4.616	1095	1112	1128	rBV5	22389	57775	3.04%	0.535%
8	5.050	1231	1247	1277	rBV4	147399	385853	20.32%	3.571%
9	5.619	1412	1424	1447	rBV	153113	314637	16.57%	2.912%
10	5.915	1503	1516	1552	rVB	290404	575905	30.34%	5.331%
11	6.072	1552	1565	1578	rBV	90584	186466	9.82%	1.726%
12	6.989	1839	1850	1870	rBV	43281	80022	4.22%	0.741%
13	7.320	1940	1953	1966	rBV	164569	290876	15.32%	2.692%
14	7.391	1966	1975	1992	rVB	92471	161827	8.52%	1.498%
15	7.628	2037	2049	2075	rBV2	25490	50403	2.65%	0.467%
16	7.982	2150	2159	2170	rVB2	14364	24598	1.30%	0.228%
17	8.088	2182	2192	2225	rBV	231226	409143	21.55%	3.787%
18	8.853	2419	2430	2433	rBV	263392	359156	18.92%	3.324%
19	8.882	2434	2439	2465	rVB	367257	586971	30.92%	5.433%
20	9.477	2604	2624	2636	rBV	8813	25690	1.35%	0.238%
21	9.545	2637	2645	2663	rVB	37043	61622	3.25%	0.570%
22	9.706	2674	2695	2708	rVB6	15726	35532	1.87%	0.329%
23	9.799	2708	2724	2735	rBV10	12152	25293	1.33%	0.234%
24	9.934	2752	2766	2776	rBV	72713	114087	6.01%	1.056%
25	10.127	2816	2826	2836	rBV4	11756	19179	1.01%	0.178%
26	10.217	2843	2854	2874	rBV	74987	144920	7.63%	1.341%
27	10.355	2885	2897	2916	rBV	205420	336163	17.71%	3.111%
28	10.448	2916	2926	2943	rBV	88382	190792	10.05%	1.766%
29	10.545	2950	2956	2971	rVB2	10520	20622	1.09%	0.191%
30	10.738	3002	3016	3034	rBV	396254	605372	31.89%	5.603%
31	10.914	3059	3071	3095	rVB	1284424	1898482	100.00%	17.572%
32	11.056	3095	3115	3120	rVV	48024	84814	4.47%	0.785%
33	11.088	3120	3125	3138	rVB	64038	96482	5.08%	0.893%
34	11.194	3150	3158	3165	rVB2	19830	28373	1.49%	0.263%
35	11.249	3165	3175	3192	rVV2	294931	521735	27.48%	4.829%
36	11.336	3192	3202	3225	rVB	366323	560728	29.54%	5.190%

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37	11.455	3231	3239	3249	rVB2	16839	27743	1.46%	0.257%
38	11.532	3249	3263	3275	rBV3	237439	524482	27.63%	4.855%
39	11.625	3282	3292	3314	rVB6	287918	593134	31.24%	5.490%
40	11.744	3319	3329	3343	rVB3	18030	29610	1.56%	0.274%
41	11.821	3343	3353	3366	rBV	38099	57575	3.03%	0.533%
42	11.911	3372	3381	3386	rBV	51460	77519	4.08%	0.718%
43	11.940	3387	3390	3399	rVB	37752	46072	2.43%	0.426%
44	12.017	3399	3414	3420	rBV	63086	101701	5.36%	0.941%
45	12.117	3434	3445	3454	rBV2	53861	101540	5.35%	0.940%
46	12.249	3476	3486	3495	rBV7	13304	28565	1.50%	0.264%
47	12.304	3495	3503	3515	rVB6	15860	30944	1.63%	0.286%
48	12.413	3531	3537	3546	rVB2	23150	32490	1.71%	0.301%
49	12.467	3546	3554	3568	rVB3	28753	47970	2.53%	0.444%
50	12.551	3568	3580	3590	rBV5	13641	21577	1.14%	0.200%
51	12.725	3628	3634	3651	rVB3	14617	25780	1.36%	0.239%
52	12.885	3675	3684	3694	rVB2	33648	51288	2.70%	0.475%
53	12.943	3695	3702	3714	rVB8	14901	23982	1.26%	0.222%

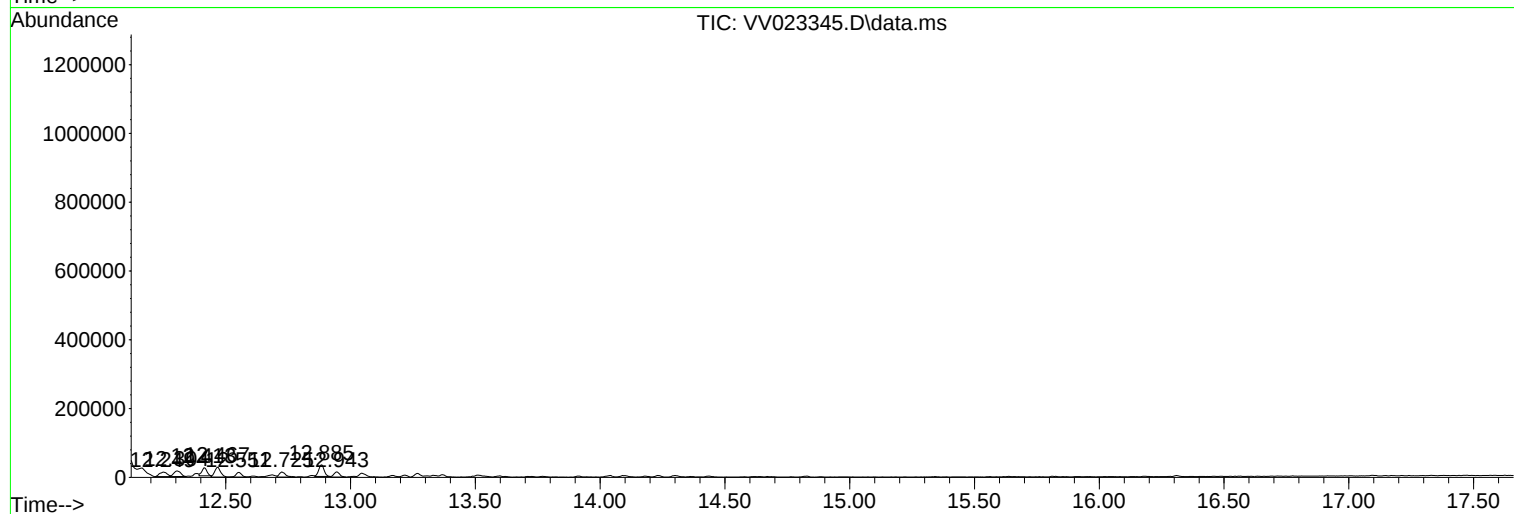
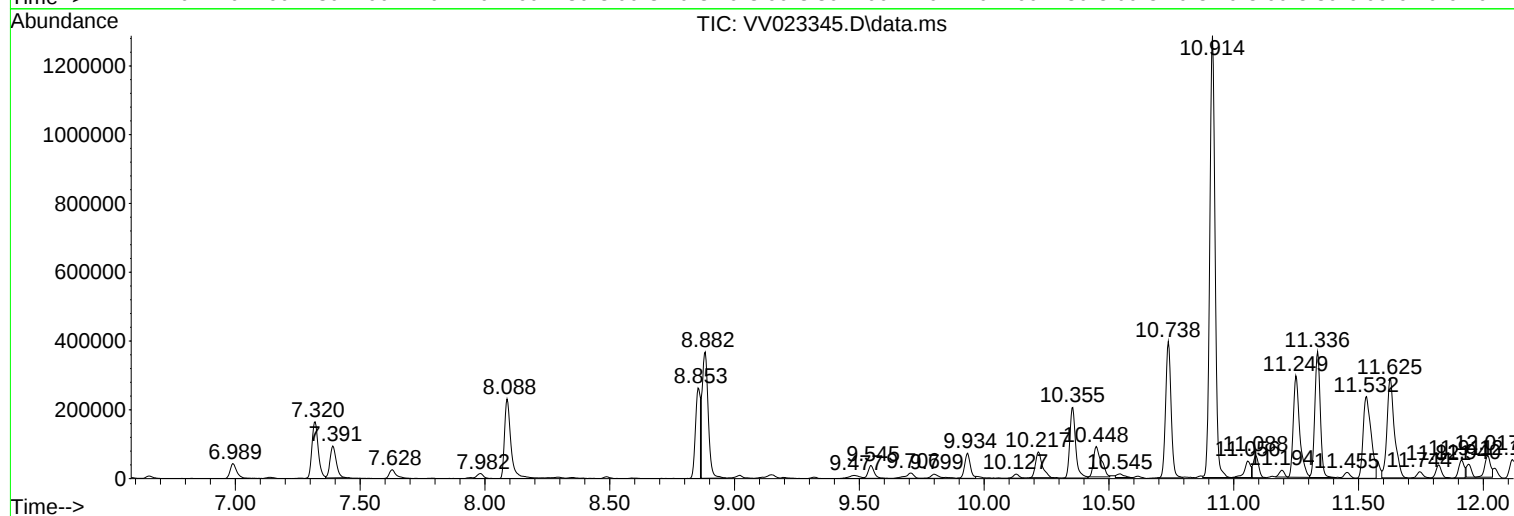
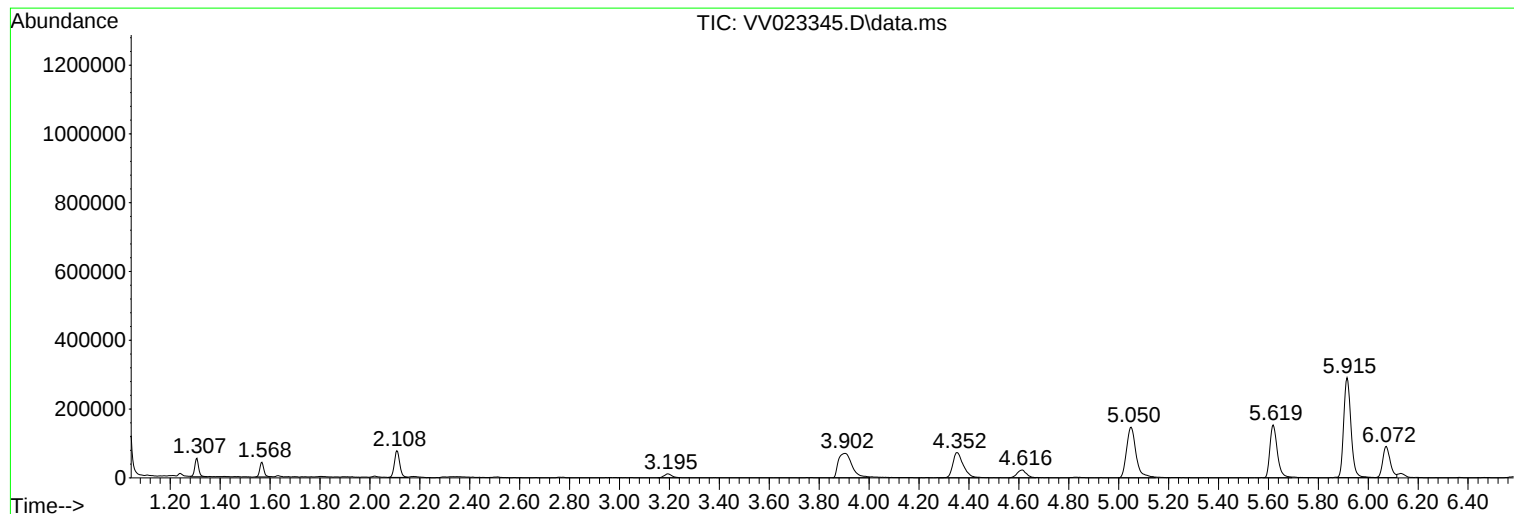
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Quant Title : TRACE VOA SFAM1.0

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



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ClientSampleId :
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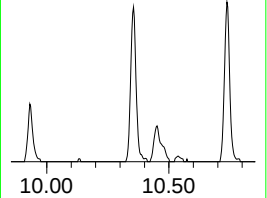
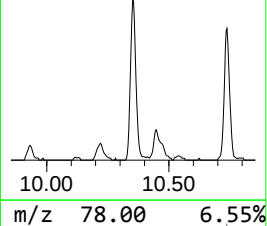
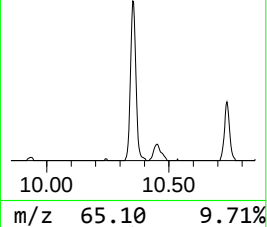
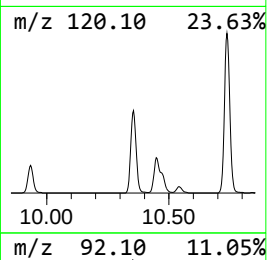
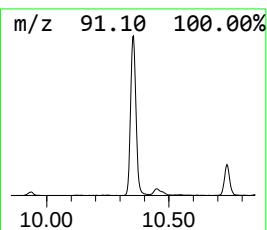
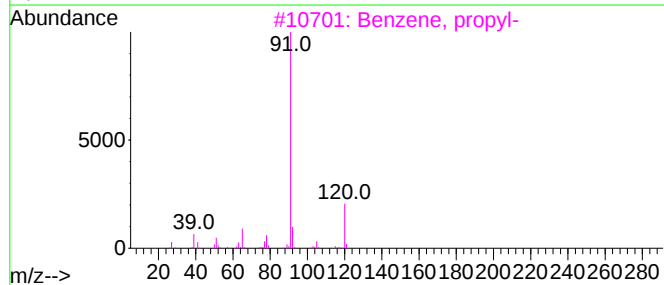
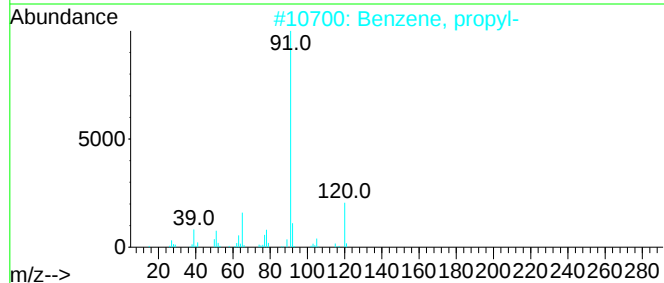
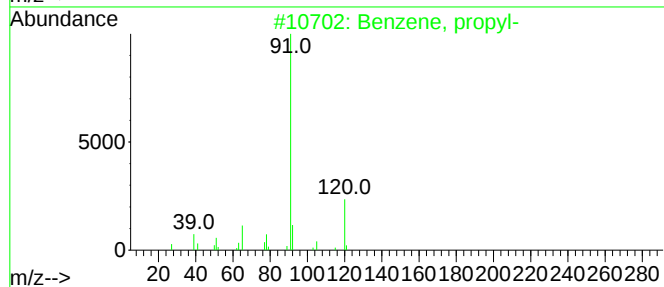
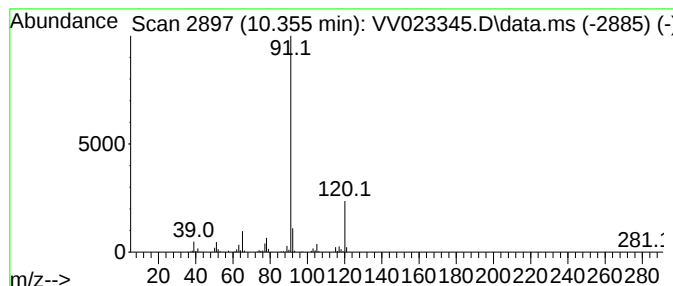
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Benzene, propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	3.22 ug/L	336163	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	87
5			1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	72



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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE7

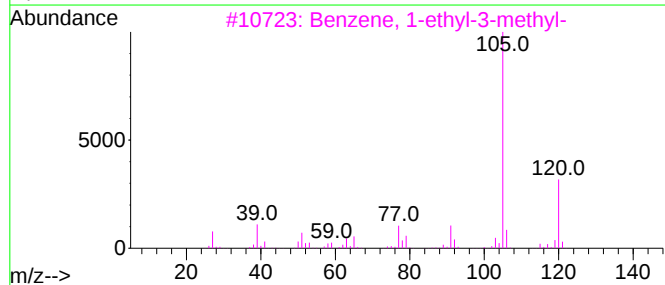
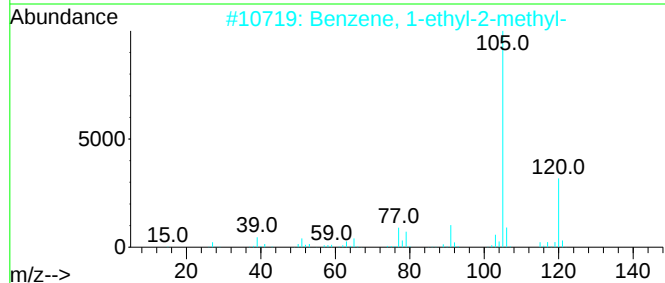
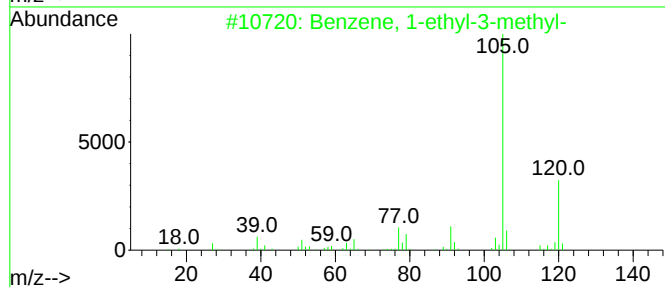
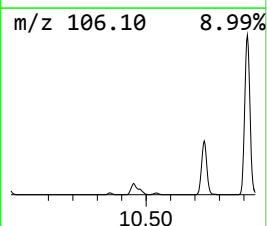
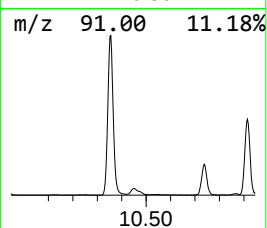
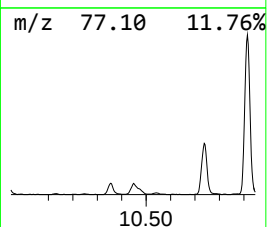
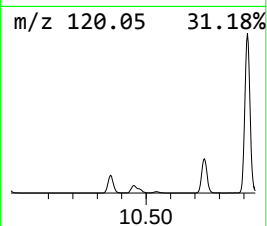
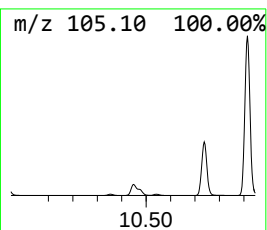
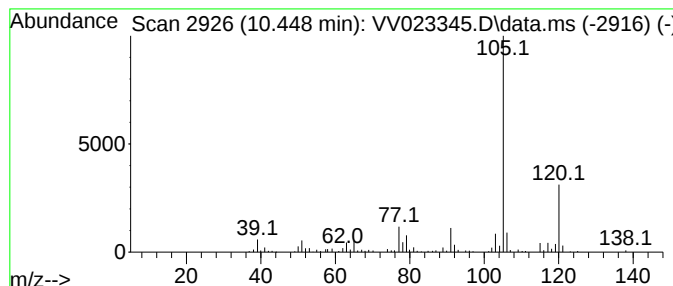
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	1.83 ug/L	190792	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

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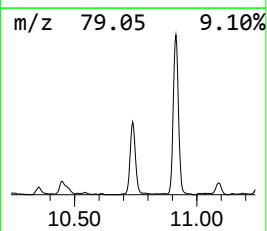
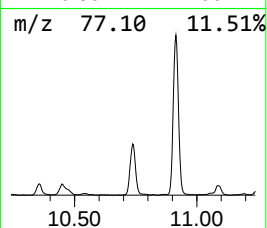
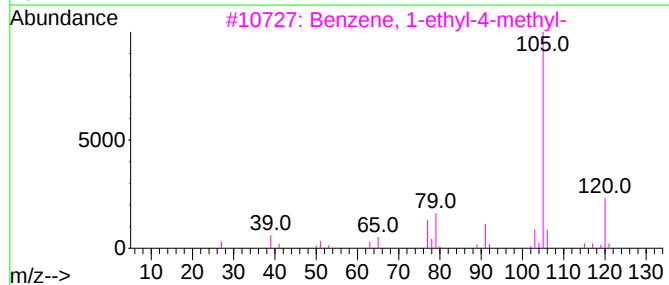
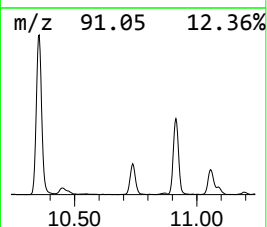
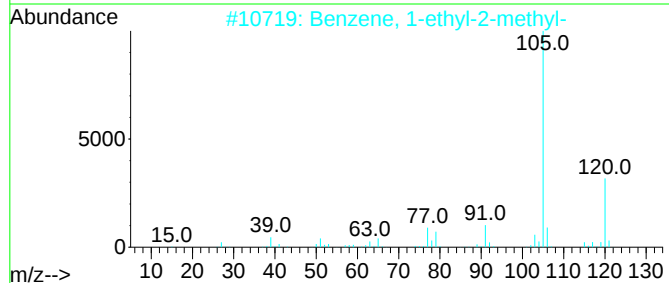
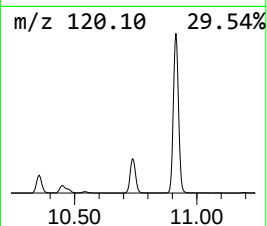
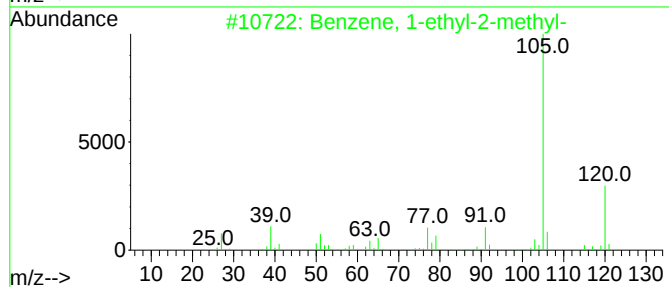
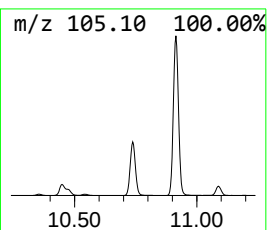
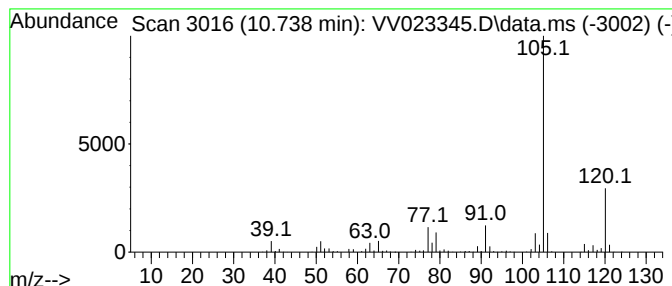
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.738	5.80 ug/L	605372	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5			Mesitylene	120	C9H12	000108-67-8	91



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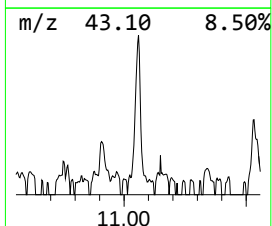
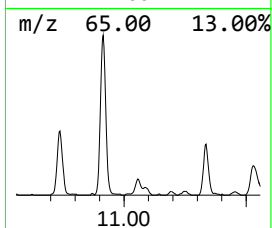
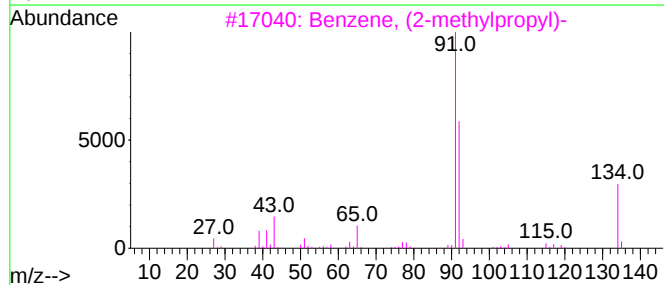
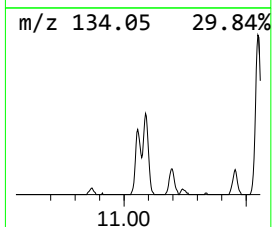
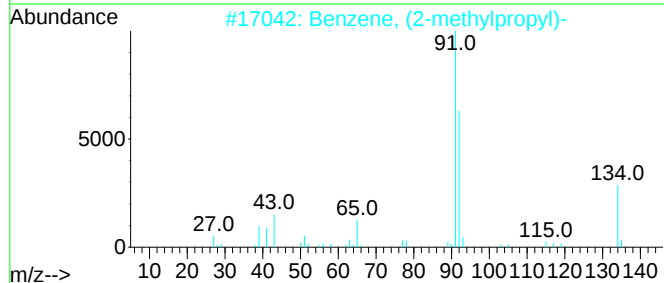
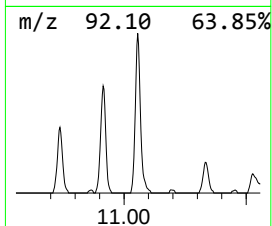
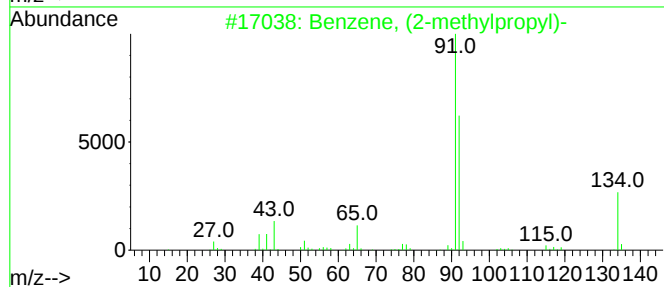
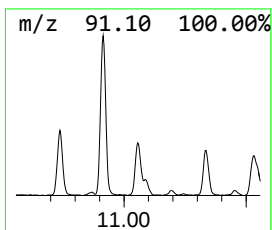
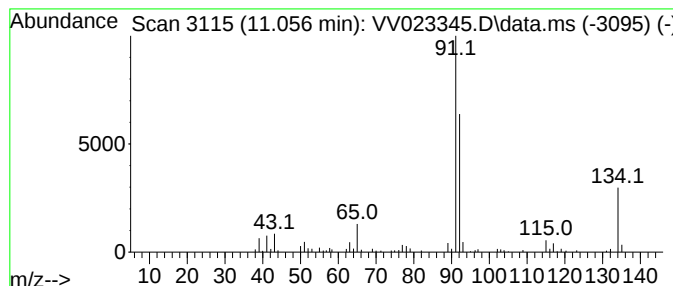
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, (2-methylpropyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.056	0.81 ug/L	84814	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87



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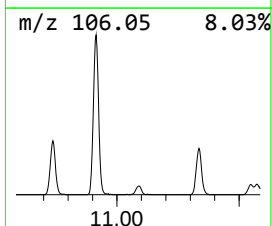
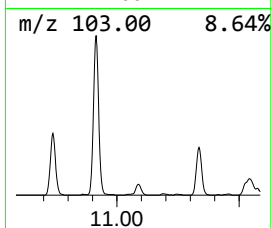
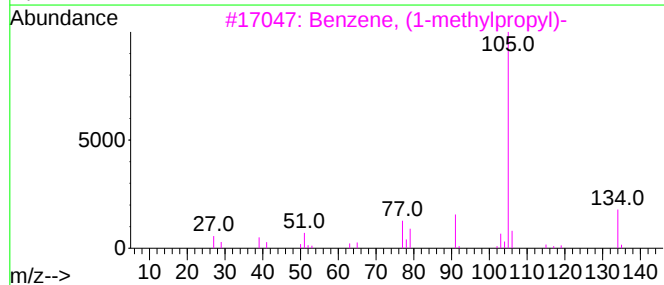
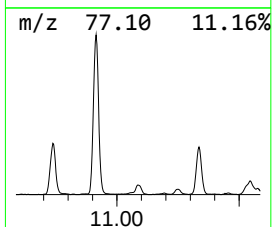
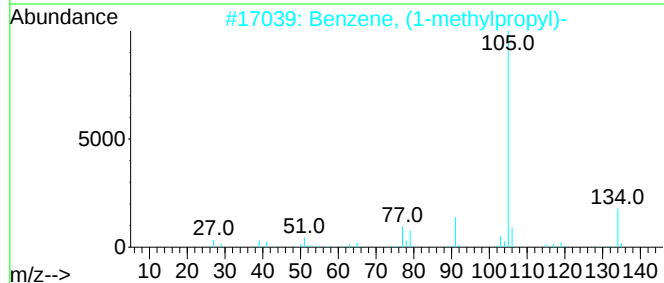
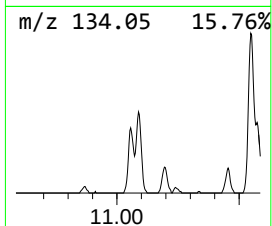
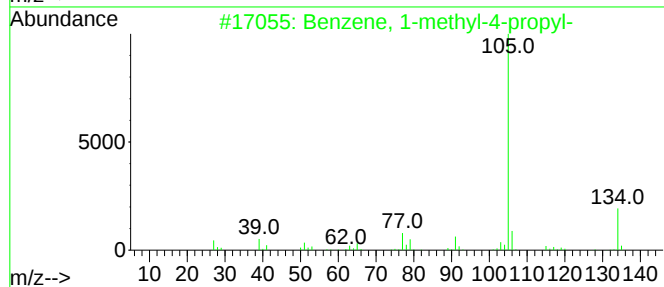
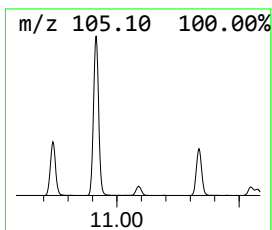
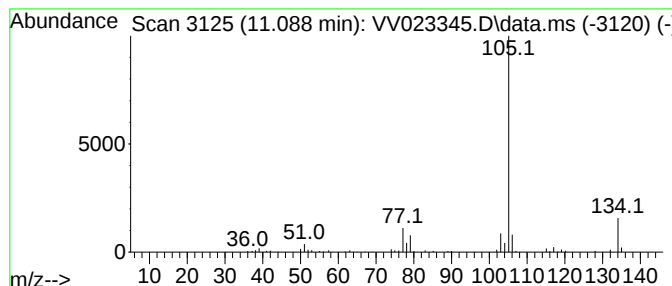
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, 1-methyl-4-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.088	0.92 ug/L	96482	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	86
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE7

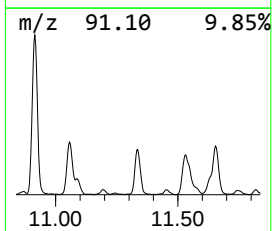
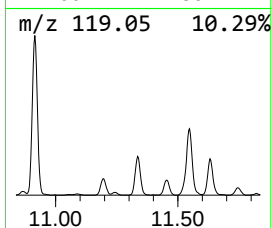
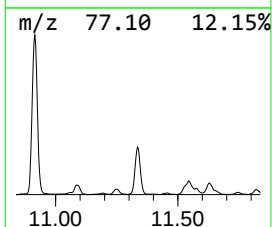
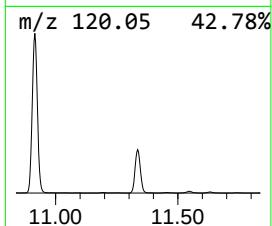
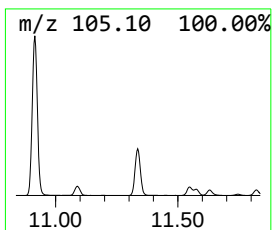
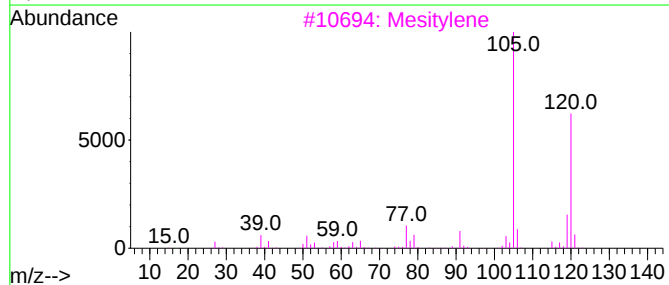
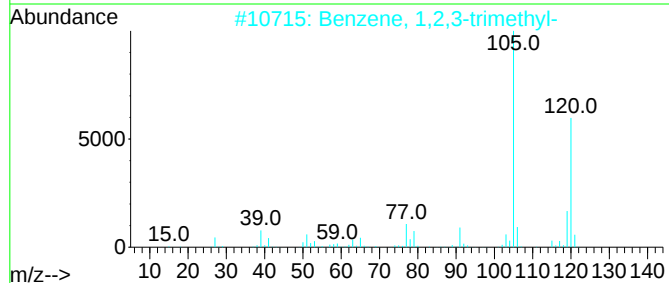
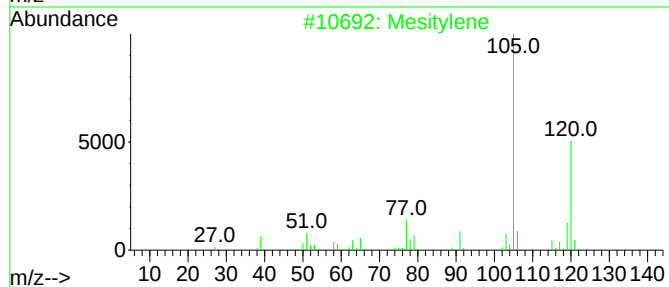
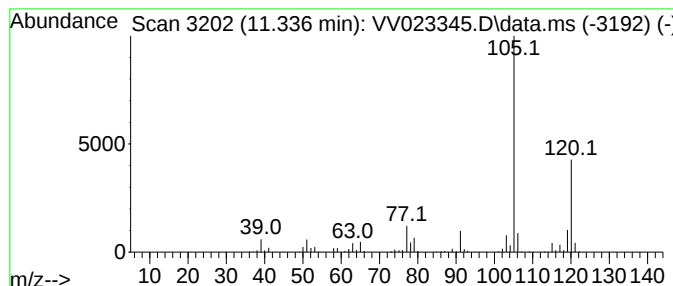
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	5.37 ug/L	560728	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Mesitylene	120	C9H12	000108-67-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE7

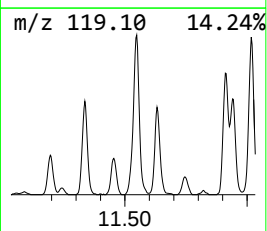
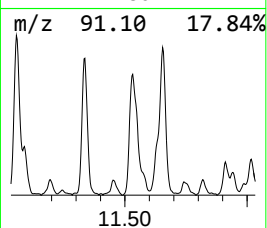
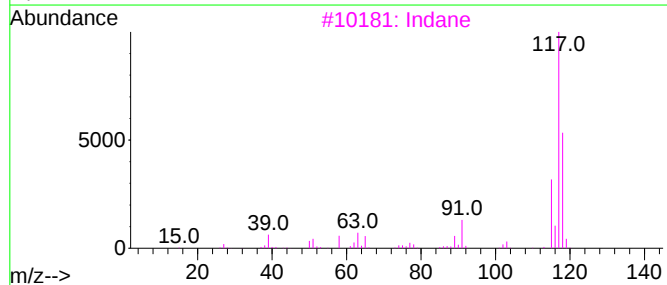
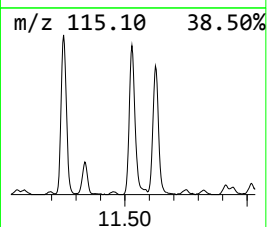
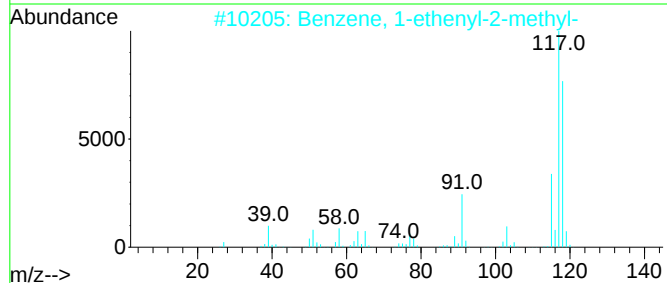
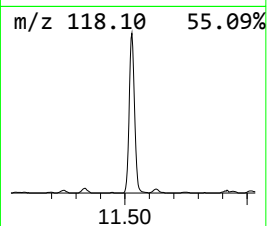
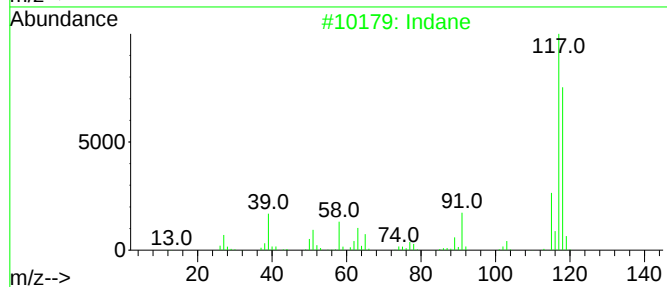
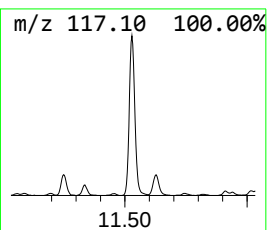
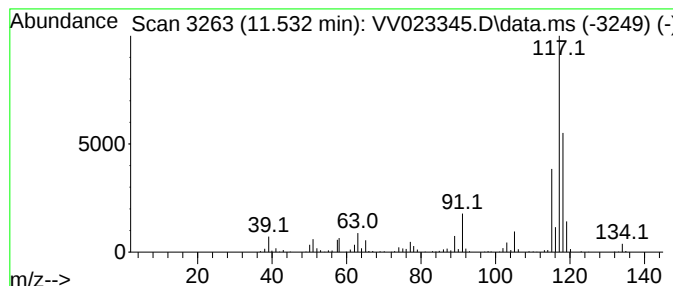
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.532	5.03 ug/L	524482	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	70
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70
3			Indane	118	C9H10	000496-11-7	70
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE7

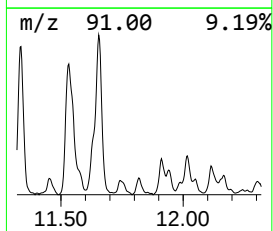
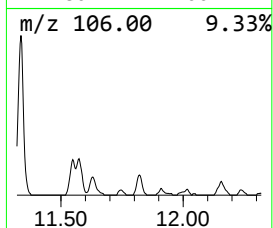
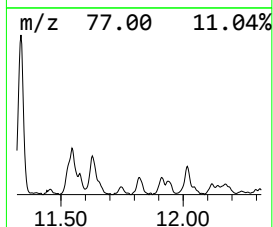
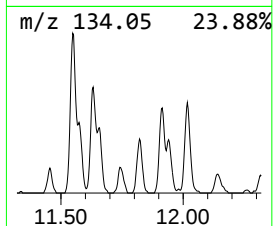
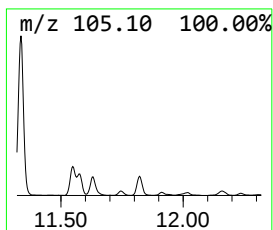
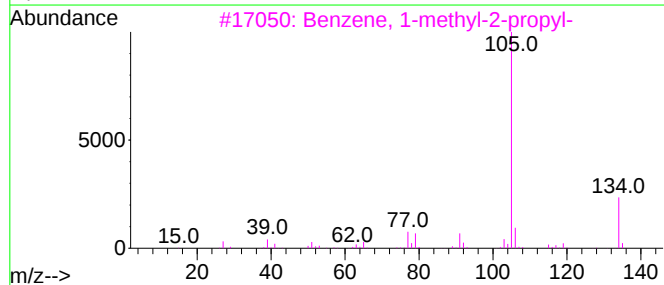
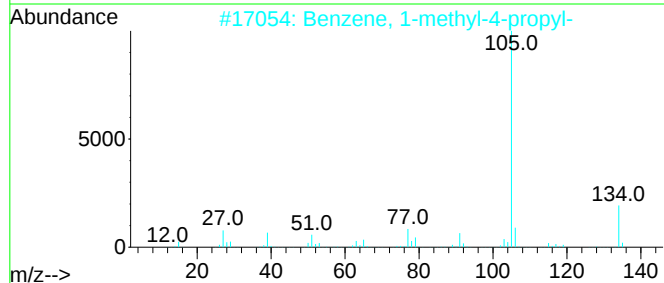
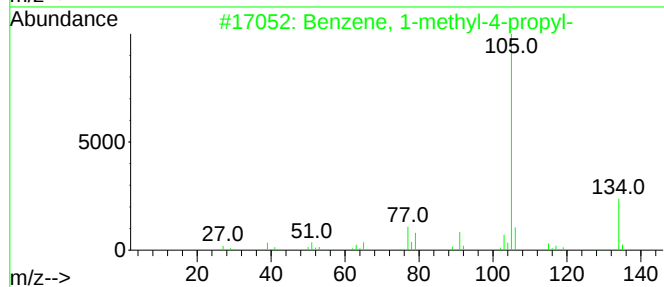
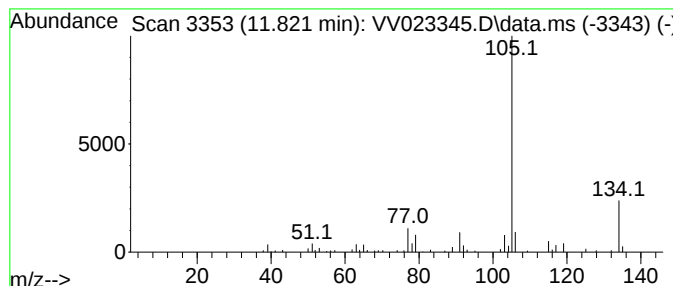
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, 1-methyl-2-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.821	0.55 ug/L	57575	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

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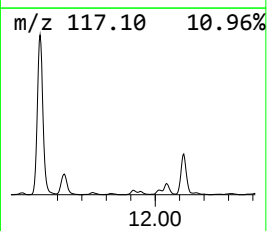
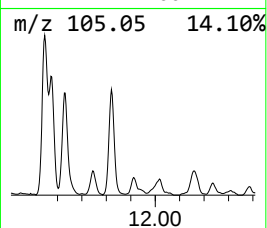
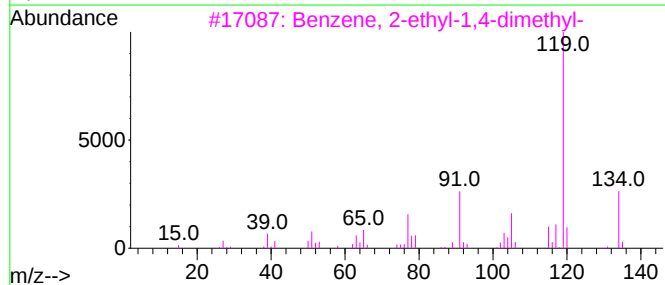
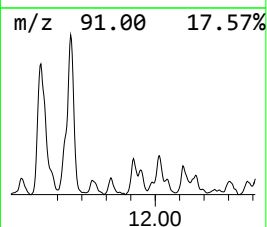
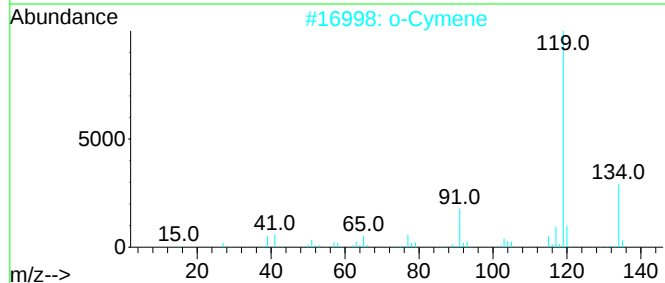
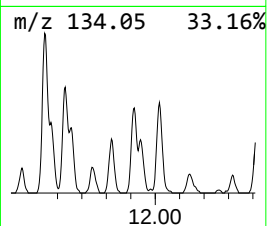
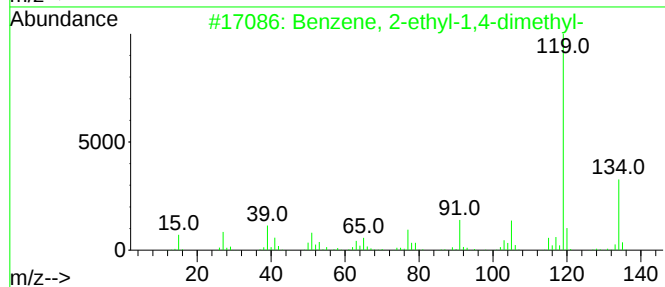
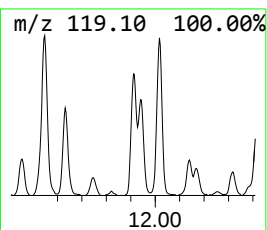
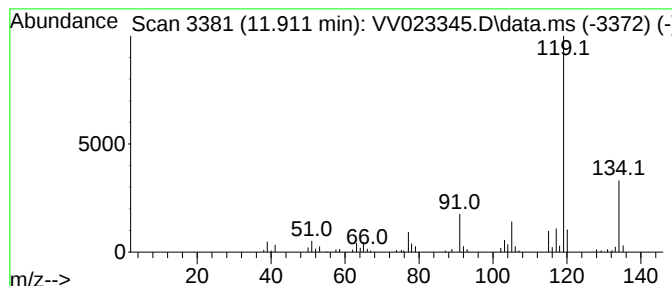
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	0.74 ug/L	77519	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

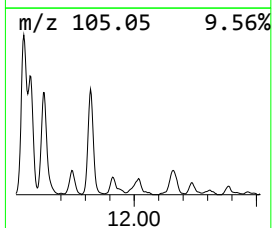
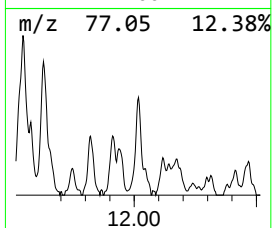
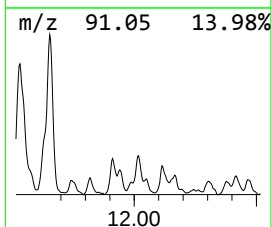
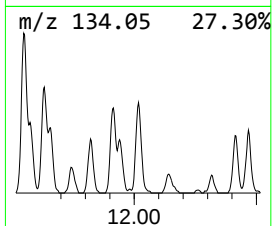
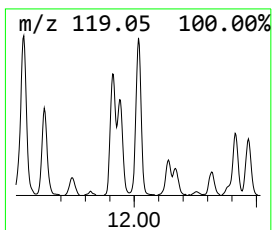
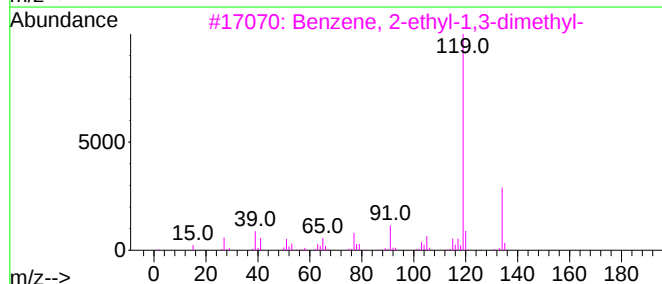
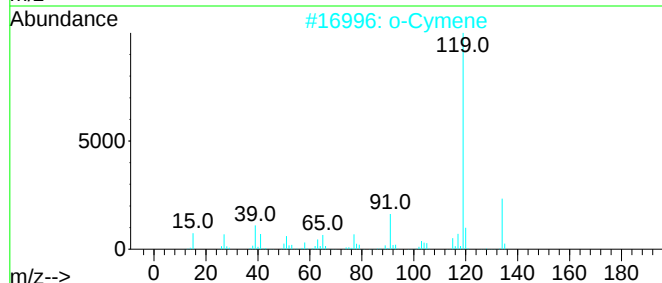
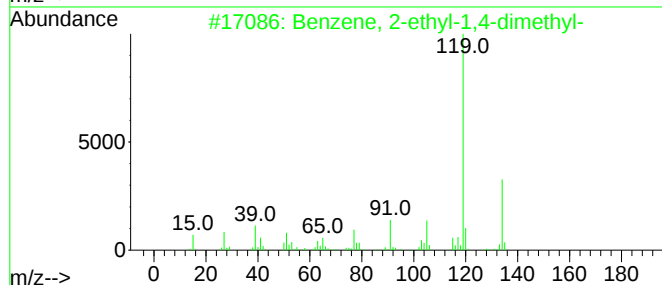
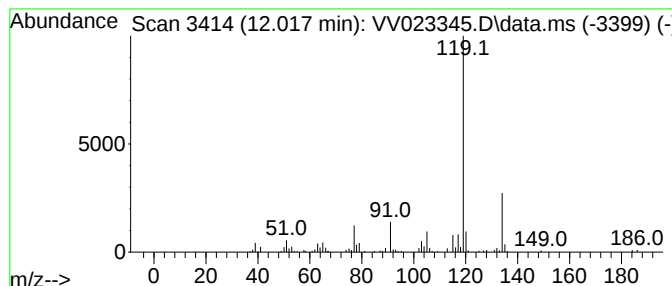
Instrument :
MSVOA_V
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.017	0.97 ug/L	101701	1,4-Dichlorobenzene-d4			11.249
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-		134	C10H14	001758-88-9	95
2	o-Cymene		134	C10H14	000527-84-4	94
3	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	94
4	Benzene, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	94
5	Benzene, 4-ethyl-1,2-dimethyl-		134	C10H14	000934-80-5	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

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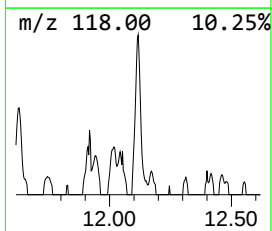
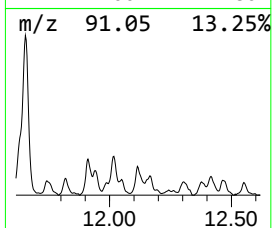
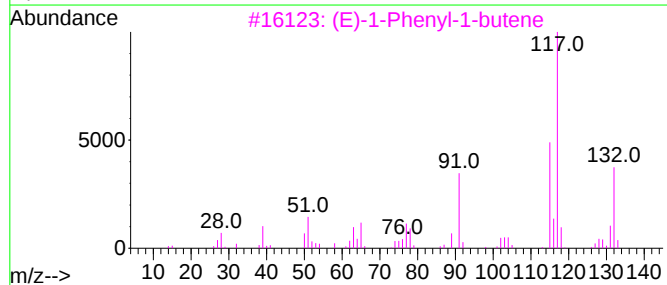
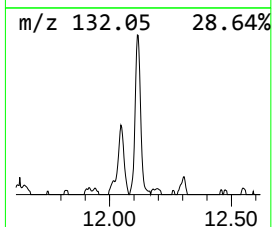
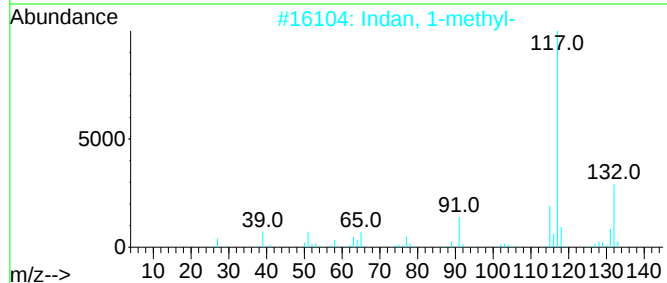
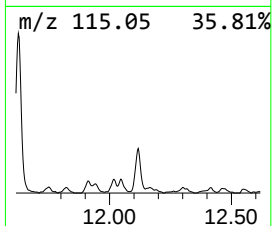
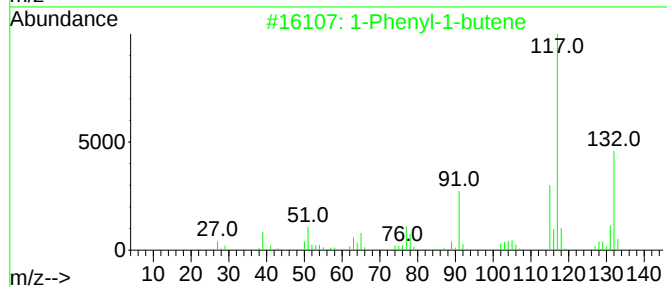
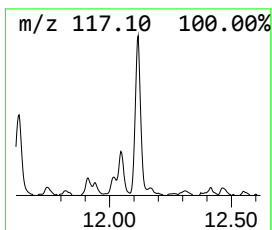
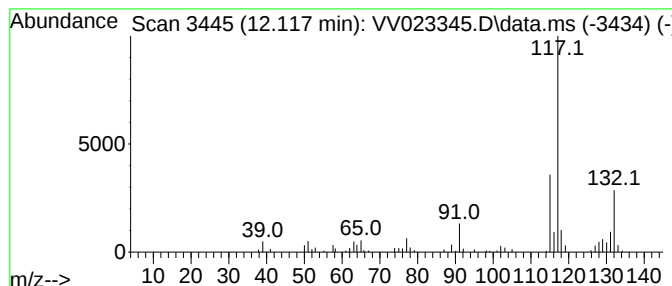
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	0.97 ug/L	101540	1,4-Dichlorobenzene-d4	11.249

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	87
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW110921\
Data File : VV023345.D
Acq On : 10 Nov 2021 20:24
Operator : SY/MD
Sample : M4542-01
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.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
Benzene, propyl-	10.355	3.2	ug/L	336163	3	11.249	521735	5.0
Benzene, 1-ethy...	10.448	1.8	ug/L	190792	3	11.249	521735	5.0
Benzene, 1-ethy...	10.738	5.8	ug/L	605372	3	11.249	521735	5.0
Benzene, (2-met...	11.056	0.8	ug/L	84814	3	11.249	521735	5.0
Benzene, 1-meth...	11.088	0.9	ug/L	96482	3	11.249	521735	5.0
Benzene, 1,2,3-...	11.336	5.4	ug/L	560728	3	11.249	521735	5.0
Indane	11.532	5.0	ug/L	524482	3	11.249	521735	5.0
Benzene, 1-meth...	11.821	0.6	ug/L	57575	3	11.249	521735	5.0
Benzene, 2-ethy...	11.911	0.7	ug/L	77519	3	11.249	521735	5.0
o-Cymene	12.017	1.0	ug/L	101701	3	11.249	521735	5.0
1-Phenyl-1-butene	12.117	1.0	ug/L	101540	3	11.249	521735	5.0