

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
 Data File : VV023347.D
 Acq On : 10 Nov 2021 21:09
 Operator : SY/MD
 Sample : M4542-03
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGGE9

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: VV023347.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.304	73	82	93	rVB	66998	70375	5.84%	0.811%
2	1.568	156	164	177	rBV	44341	52543	4.36%	0.605%
3	2.108	321	332	346	rBV	95378	140711	11.68%	1.621%
4	3.191	656	669	687	rBV2	11257	25551	2.12%	0.294%
5	3.908	877	892	921	rBV3	65714	229487	19.05%	2.643%
6	4.352	1014	1030	1062	rBV3	71431	198061	16.44%	2.281%
7	4.612	1094	1111	1137	rVB3	22583	58656	4.87%	0.676%
8	5.050	1229	1247	1287	rBV2	153080	393522	32.66%	4.533%
9	5.619	1412	1424	1448	rBV	145064	298676	24.79%	3.440%
10	5.915	1499	1516	1541	rBV	225669	450712	37.41%	5.192%
11	6.072	1552	1565	1580	rBV	85680	176684	14.66%	2.035%
12	6.992	1841	1851	1870	rBV3	41982	77063	6.40%	0.888%
13	7.320	1941	1953	1965	rBV	188173	320168	26.57%	3.688%
14	7.391	1965	1975	1993	rVB	213166	369972	30.71%	4.262%
15	7.628	2040	2049	2065	rBV2	24796	45932	3.81%	0.529%
16	7.982	2150	2159	2169	rVB4	12719	21610	1.79%	0.249%
17	8.091	2182	2193	2227	rBV	178923	329573	27.35%	3.796%
18	8.857	2419	2431	2435	rBV	240736	394428	32.74%	4.543%
19	8.882	2436	2439	2461	rVB	266433	364302	30.24%	4.196%
20	9.017	2474	2481	2491	rVB	9384	12814	1.06%	0.148%
21	9.143	2509	2520	2533	rVB3	13770	26680	2.21%	0.307%
22	9.477	2608	2624	2637	rBV3	4730	13776	1.14%	0.159%
23	9.548	2637	2646	2666	rVB	33976	57792	4.80%	0.666%
24	9.709	2677	2696	2705	rBV9	8809	19231	1.60%	0.222%
25	9.805	2715	2726	2740	rBV9	7304	18192	1.51%	0.210%
26	9.934	2756	2766	2778	rBV2	49174	78805	6.54%	0.908%
27	10.217	2844	2854	2875	rVB	63304	114353	9.49%	1.317%
28	10.355	2886	2897	2917	rBV	136244	222121	18.44%	2.559%
29	10.451	2917	2927	2943	rBV	57205	126359	10.49%	1.456%
30	10.738	3003	3016	3036	rBV	255753	396704	32.93%	4.570%
31	10.914	3060	3071	3092	rVB	817894	1204813	100.00%	13.878%
32	11.059	3096	3116	3120	rBV2	28717	51154	4.25%	0.589%
33	11.091	3120	3126	3137	rVB	40010	61515	5.11%	0.709%
34	11.152	3137	3145	3151	rBV6	9806	15217	1.26%	0.175%
35	11.194	3152	3158	3165	rVV4	14865	20986	1.74%	0.242%
36	11.252	3165	3176	3192	rVV	271179	469039	38.93%	5.403%

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37	11.336	3192	3202	3223	rVB	246276	377627	31.34%	4.350%
38	11.455	3229	3239	3250	rVB	11246	17952	1.49%	0.207%
39	11.532	3251	3263	3283	rBV3	164229	393167	32.63%	4.529%
40	11.628	3283	3293	3312	rVB4	241481	475224	39.44%	5.474%
41	11.747	3321	3330	3340	rBV3	11794	18082	1.50%	0.208%
42	11.821	3344	3353	3363	rVB	24207	35850	2.98%	0.413%
43	11.914	3371	3382	3387	rBV	34394	53376	4.43%	0.615%
44	11.943	3387	3391	3401	rVB	24323	30033	2.49%	0.346%
45	12.017	3401	3414	3420	rBV	38637	61612	5.11%	0.710%
46	12.117	3435	3445	3456	rBV3	34412	70521	5.85%	0.812%
47	12.246	3475	3485	3496	rBV10	7184	17111	1.42%	0.197%
48	12.310	3496	3505	3510	rBV6	9253	17011	1.41%	0.196%
49	12.345	3510	3516	3524	rVB3	13057	17219	1.43%	0.198%
50	12.416	3532	3538	3546	rVB	15848	21805	1.81%	0.251%
51	12.467	3546	3554	3568	rVB2	18816	30699	2.55%	0.354%
52	12.554	3573	3581	3590	rBV4	8625	12638	1.05%	0.146%
53	12.882	3675	3683	3695	rVB3	24081	38307	3.18%	0.441%
54	12.943	3695	3702	3715	rVB9	7817	12622	1.05%	0.145%
55	13.049	3726	3735	3746	rVB7	7573	13106	1.09%	0.151%
56	13.522	3868	3882	3896	rBV4	10392	25283	2.10%	0.291%
57	17.101	4988	4995	5005	rVB	9586	14558	1.21%	0.168%

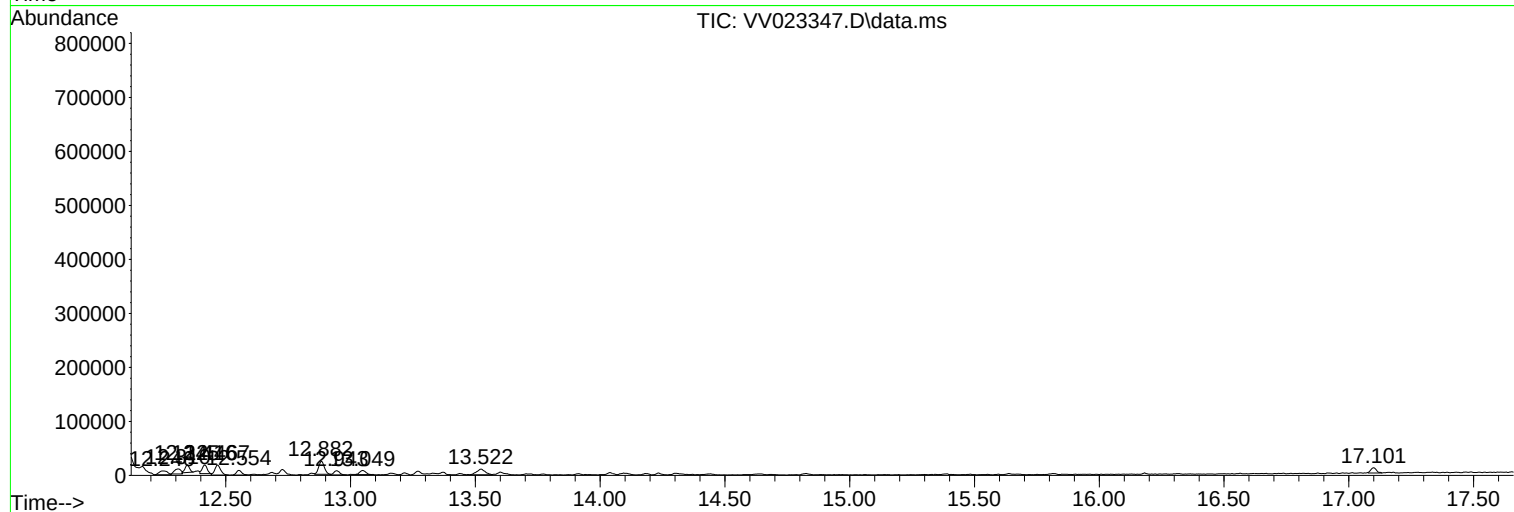
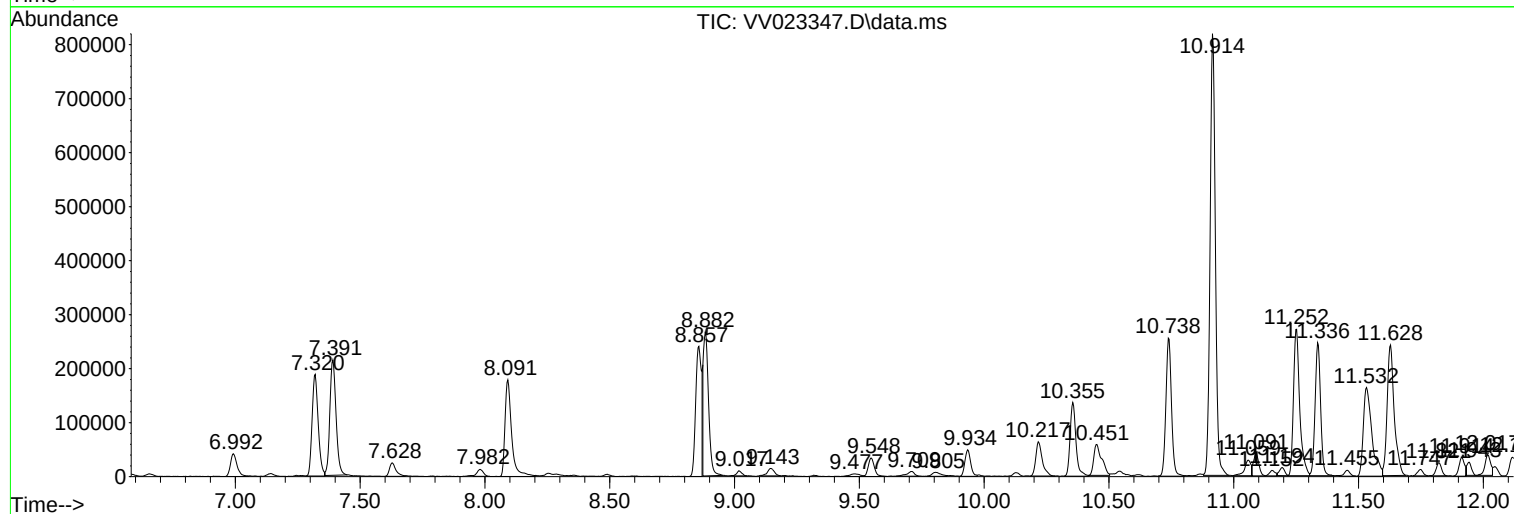
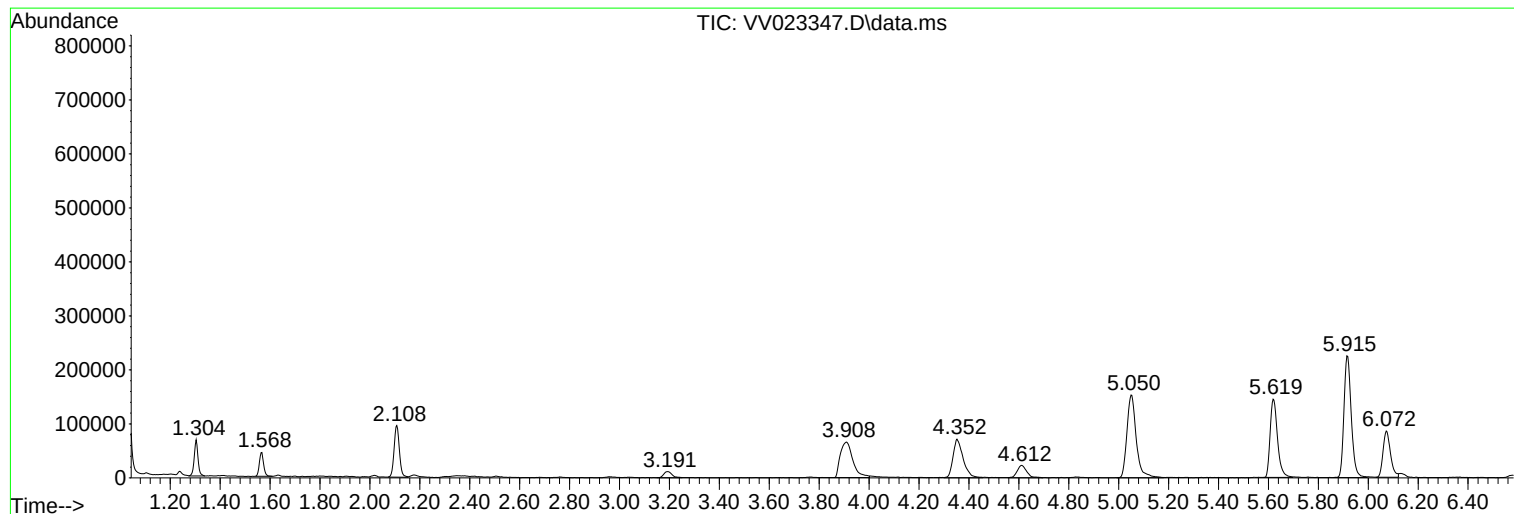
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Quant Title : TRACE VOA SFAM1.0

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



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Instrument :
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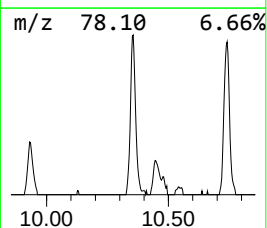
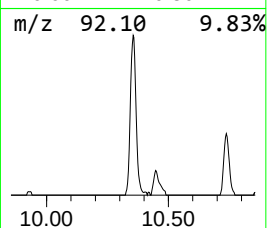
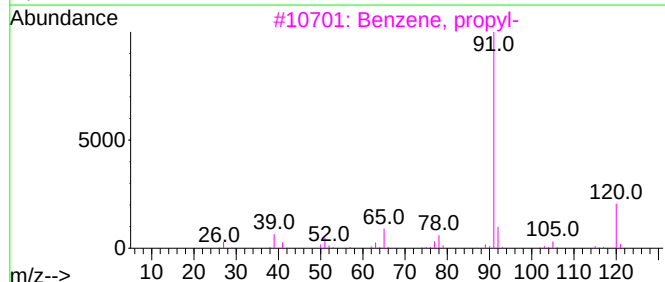
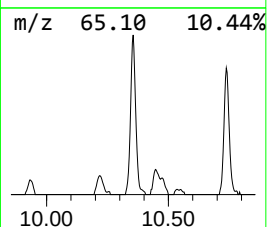
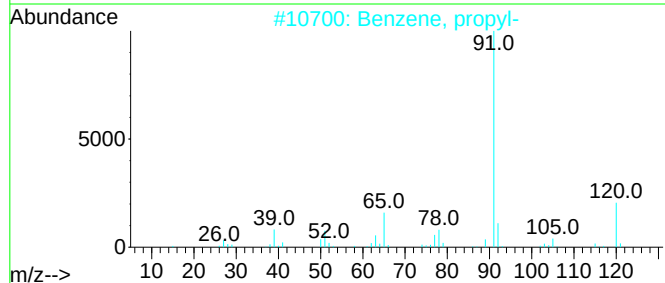
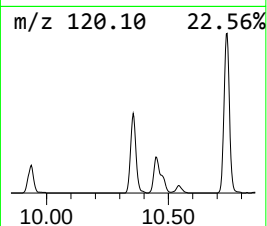
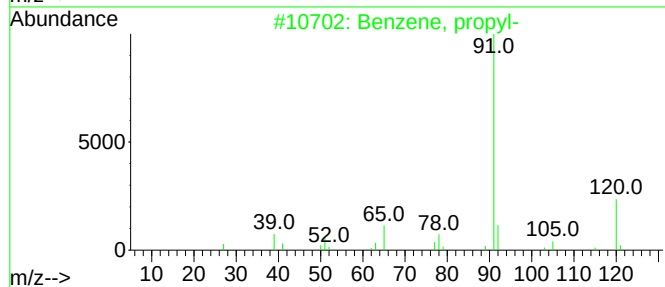
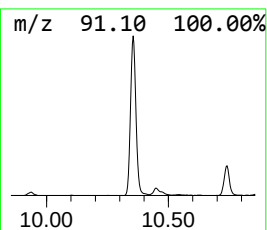
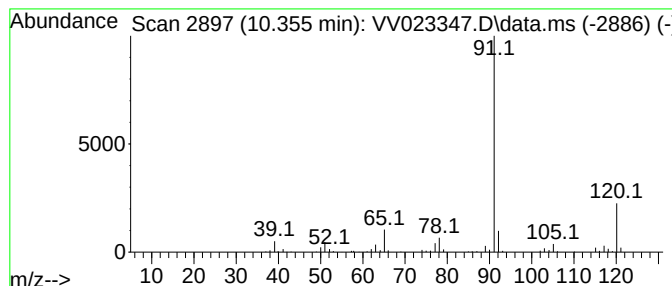
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Quant Title : TRACE VOA SFAM1.0

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	2.37 ug/L	222121	1,4-Dichlorobenzene-d4	11.252

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	80
5			1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	72



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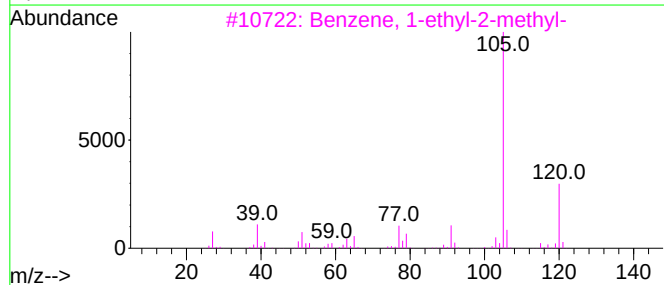
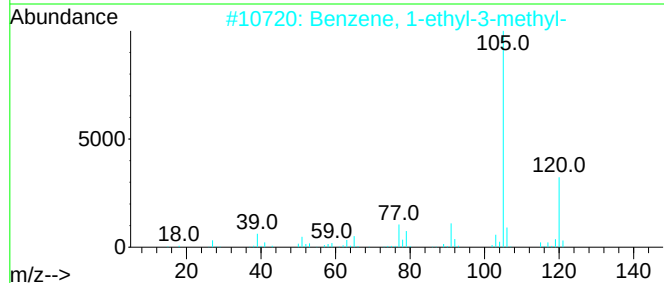
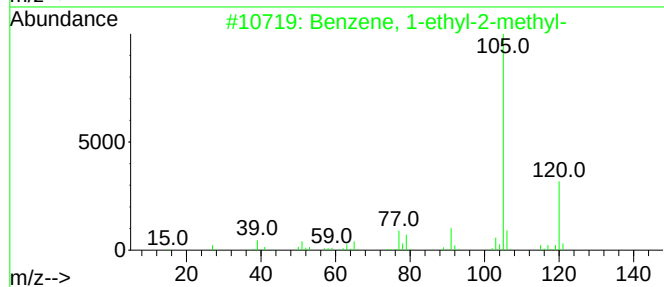
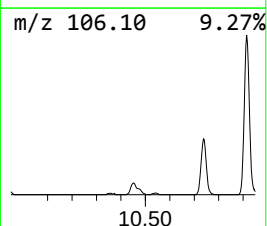
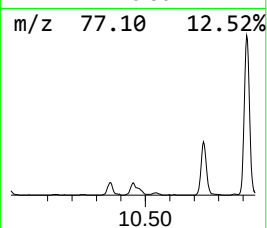
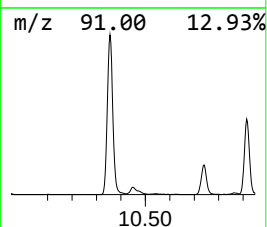
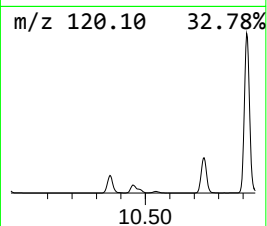
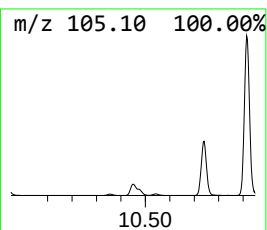
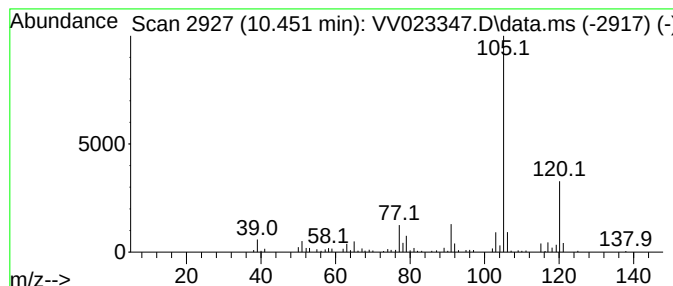
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.451	1.35 ug/L	126359	1,4-Dichlorobenzene-d4	11.252

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



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Instrument :
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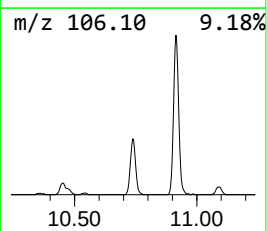
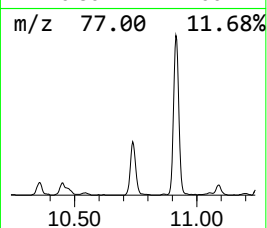
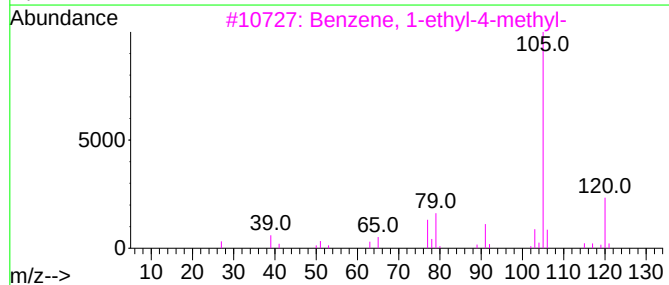
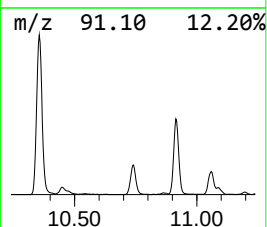
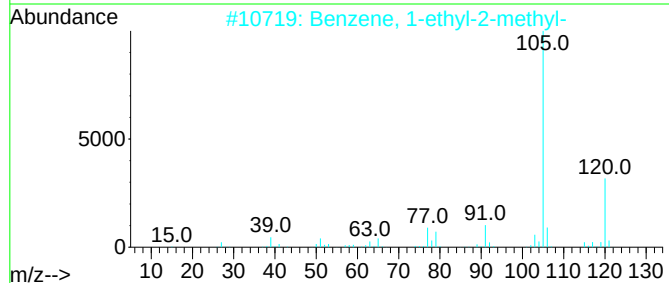
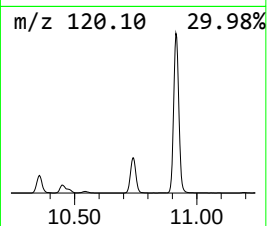
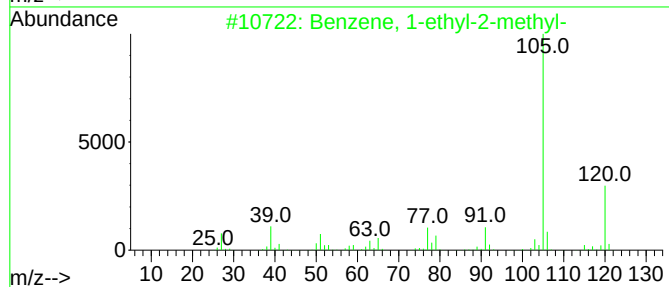
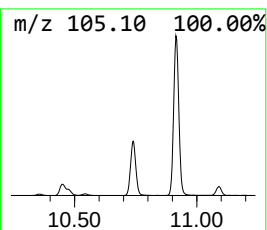
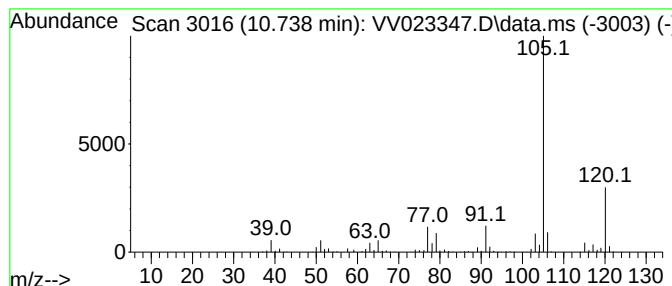
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.738	4.23 ug/L	396704	1,4-Dichlorobenzene-d4	11.252

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	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93



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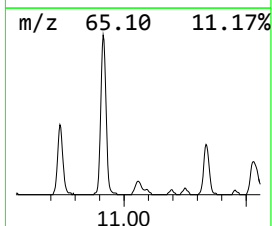
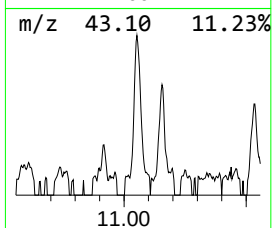
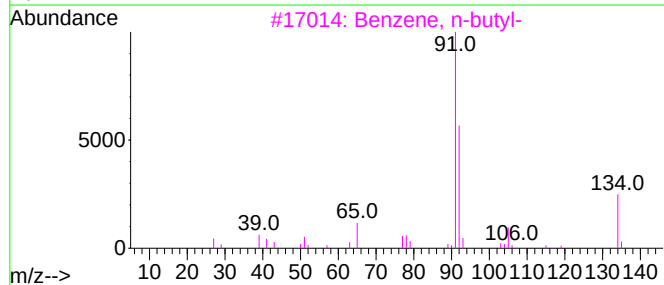
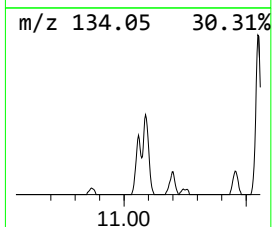
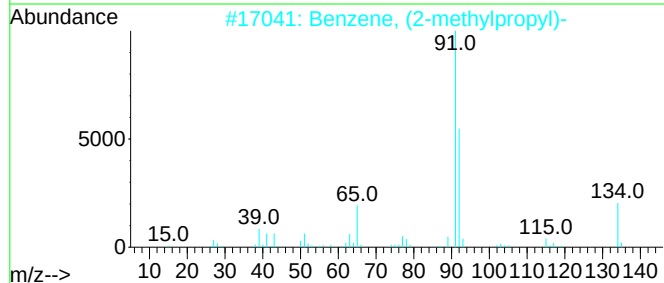
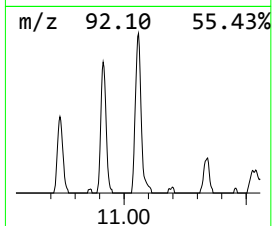
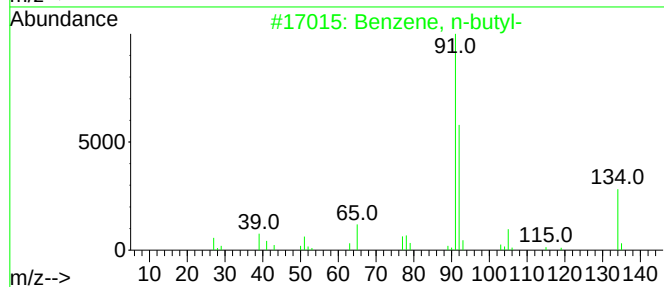
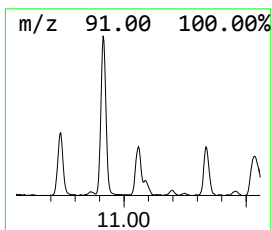
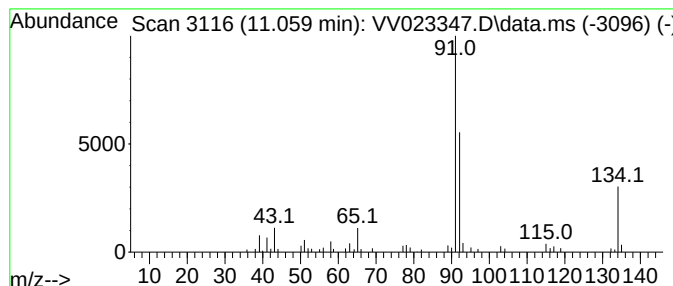
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, n-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.059	0.55 ug/L	51154	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, n-butyl-	134	C10H14	000104-51-8	90
2			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3			Benzene, n-butyl-	134	C10H14	000104-51-8	90
4			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90



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BGGE9

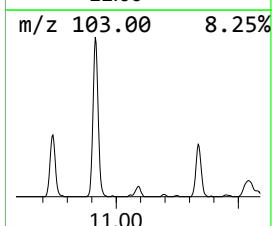
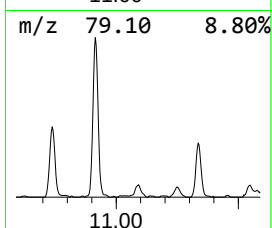
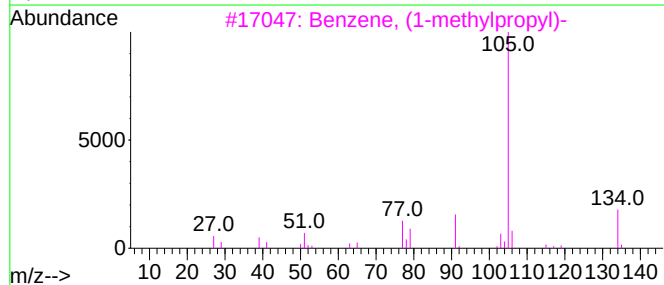
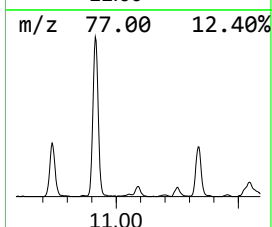
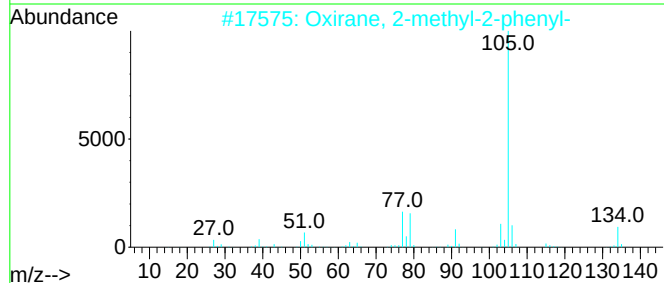
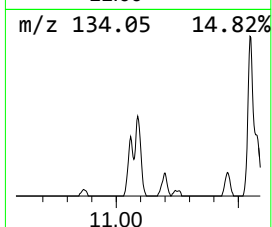
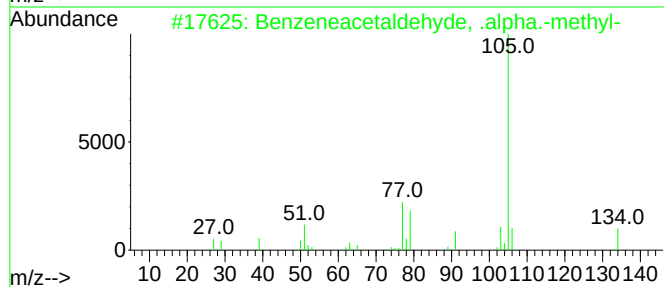
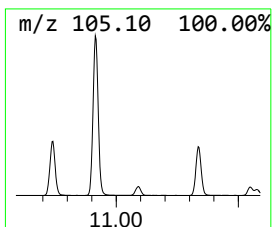
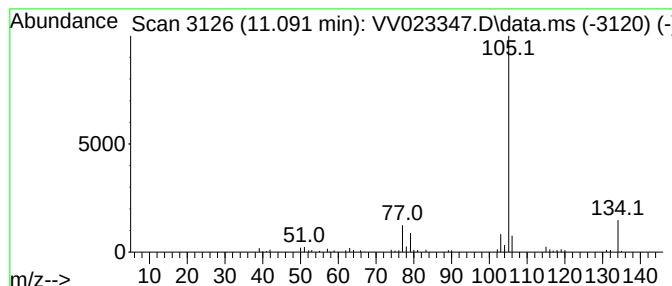
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzeneacetaldehyde, .alpha... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.091	0.66 ug/L	61515	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	80
2			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	80
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72
4			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	64
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE9

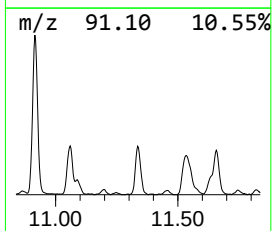
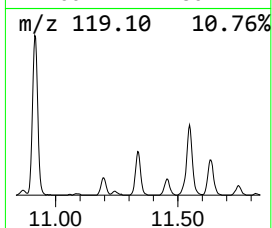
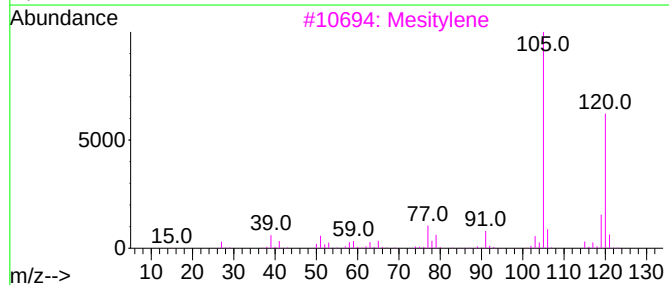
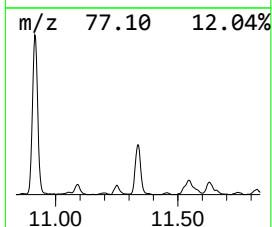
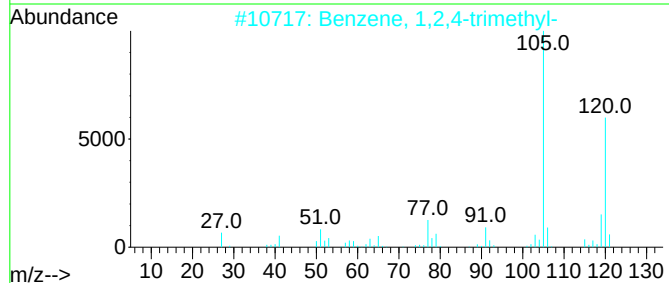
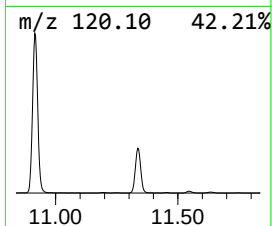
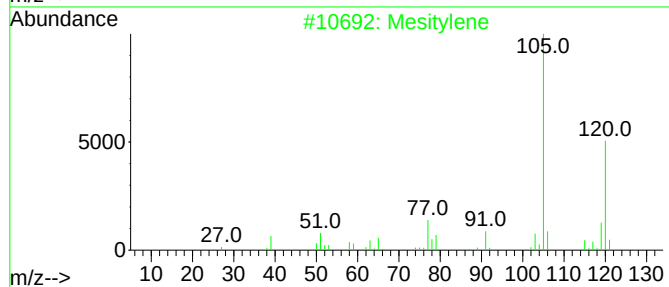
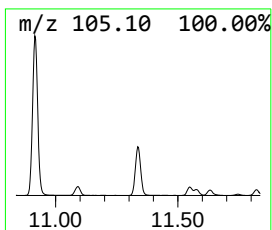
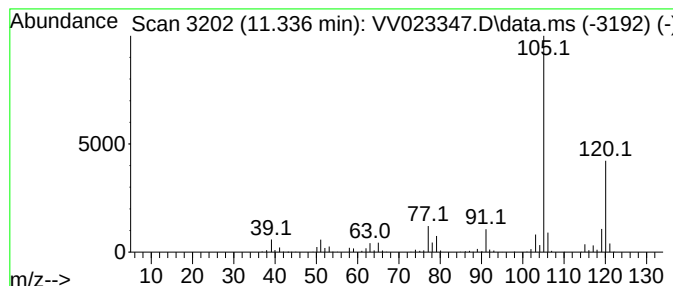
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	4.03 ug/L	377627	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3			Mesitylene	120	C9H12	000108-67-8	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

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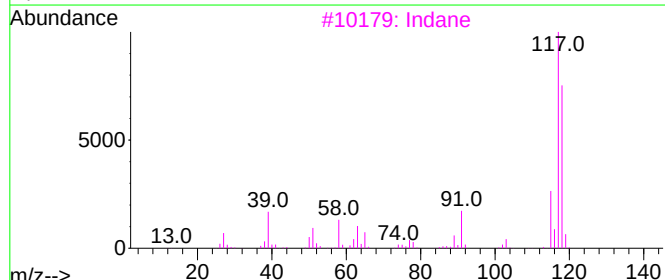
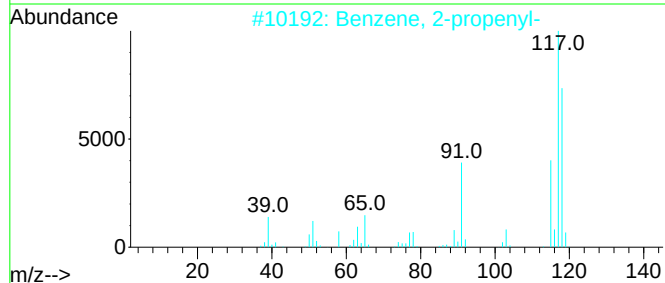
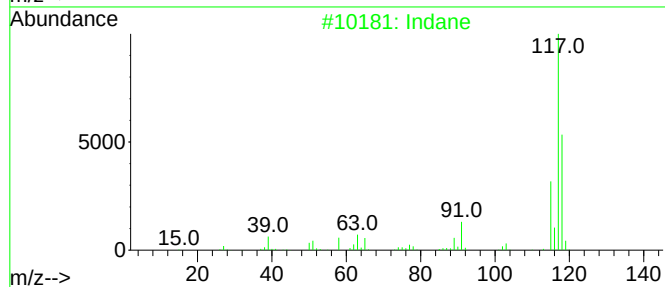
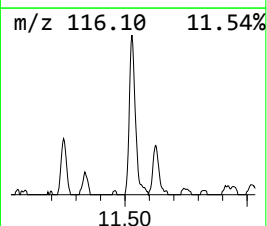
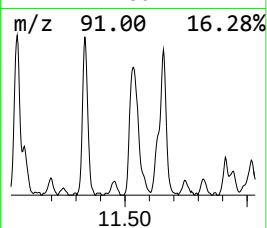
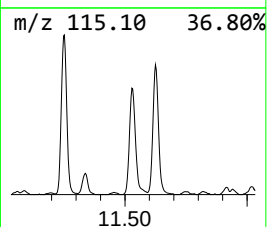
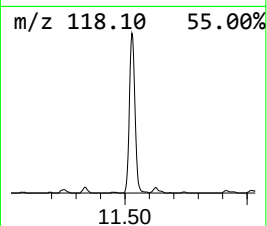
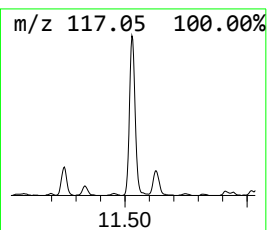
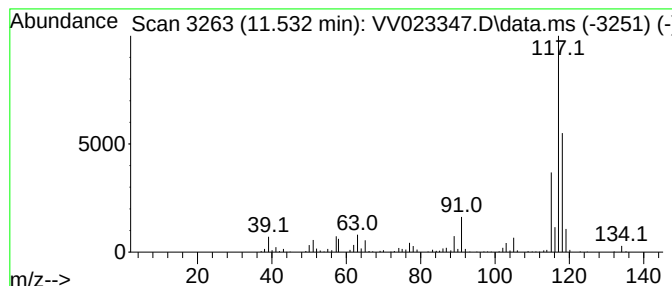
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.532	4.19 ug/L	393167	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3			Indane	118	C9H10	000496-11-7	81
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

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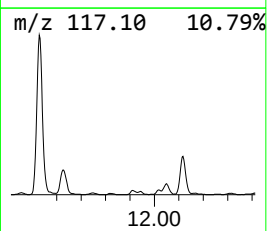
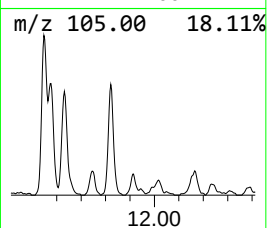
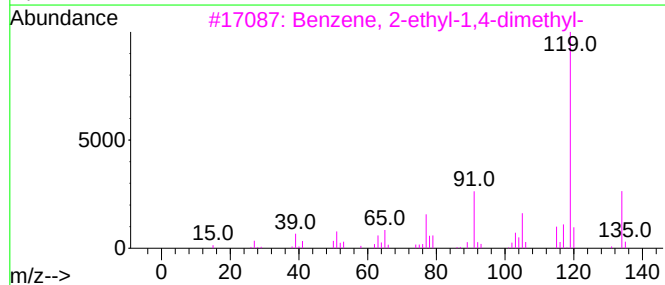
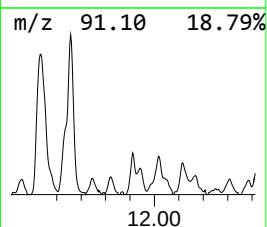
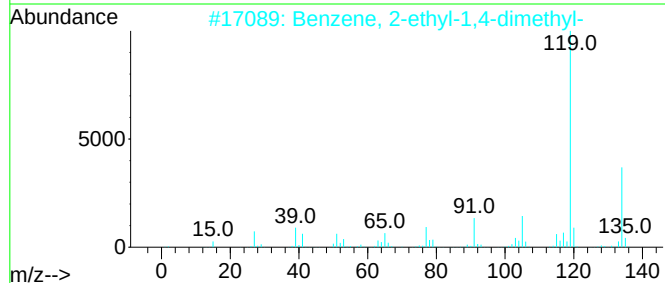
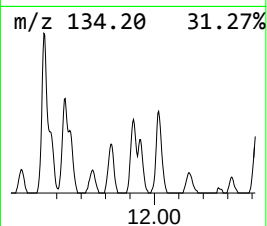
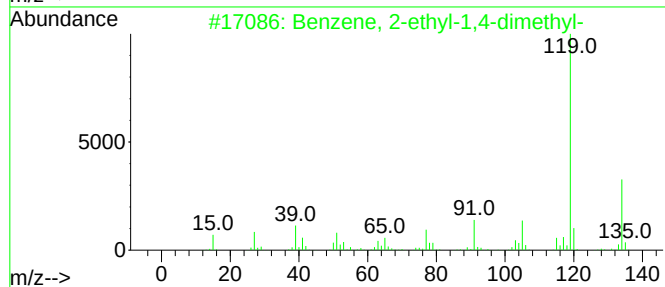
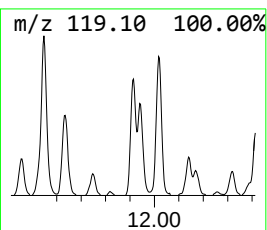
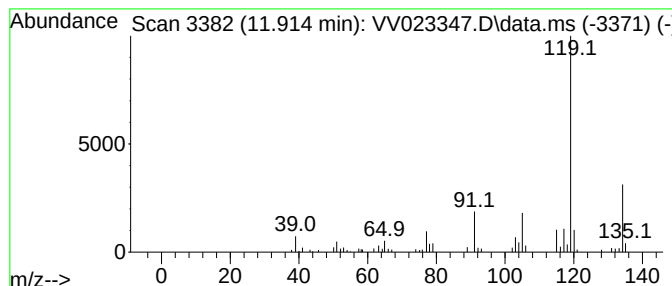
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.914	0.57 ug/L	53376	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
4			p-Cymene	134	C10H14	000099-87-6	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE9

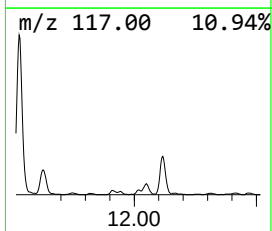
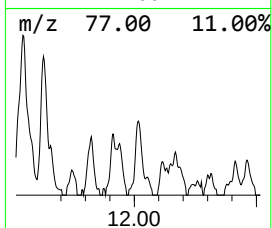
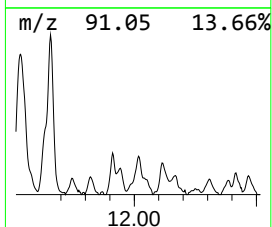
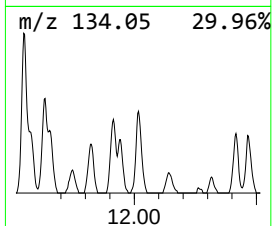
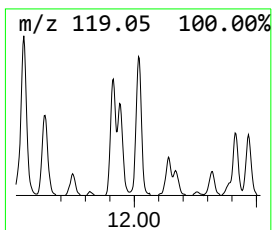
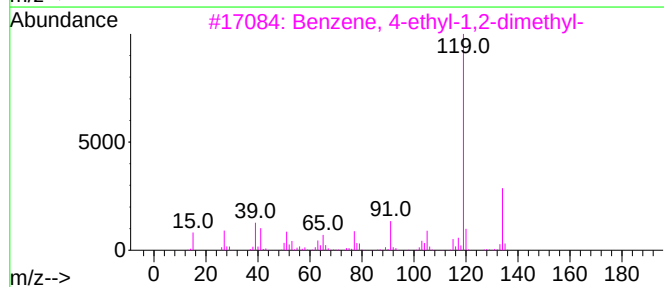
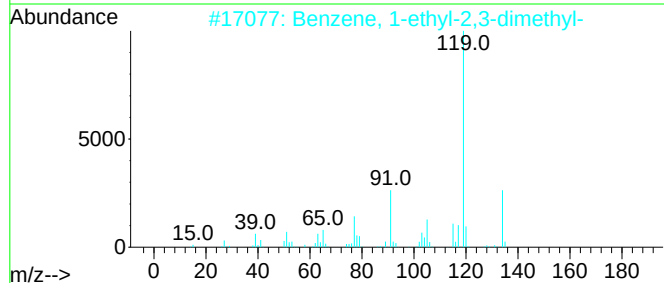
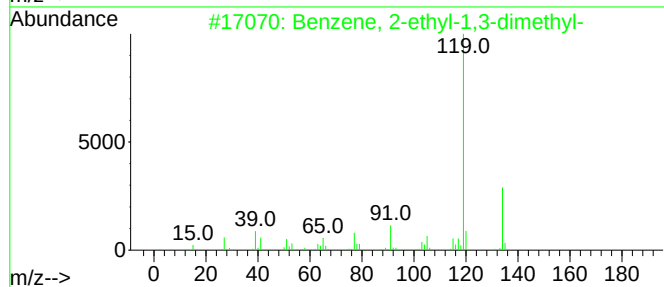
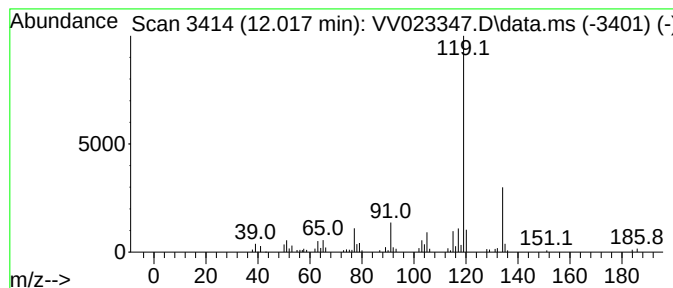
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.017	0.66 ug/L	61612	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

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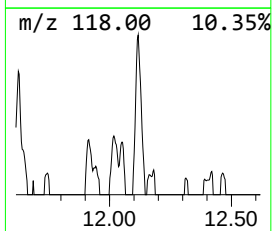
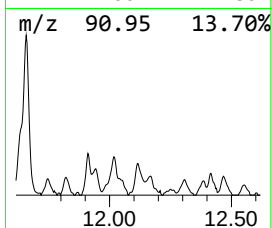
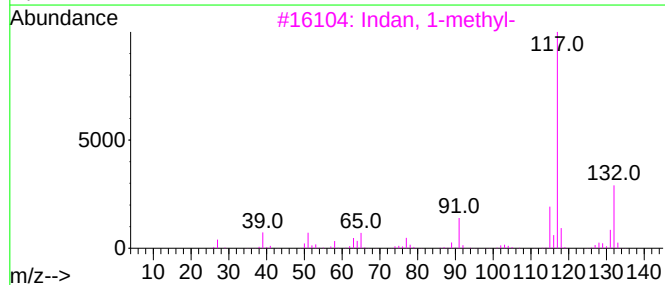
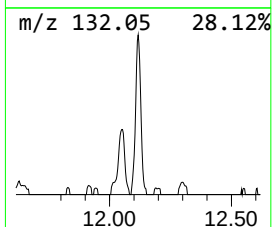
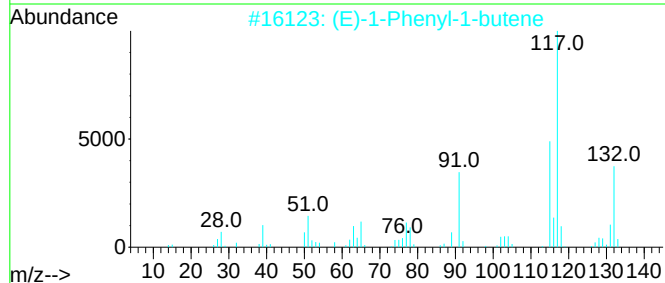
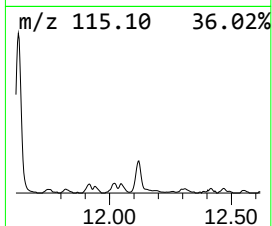
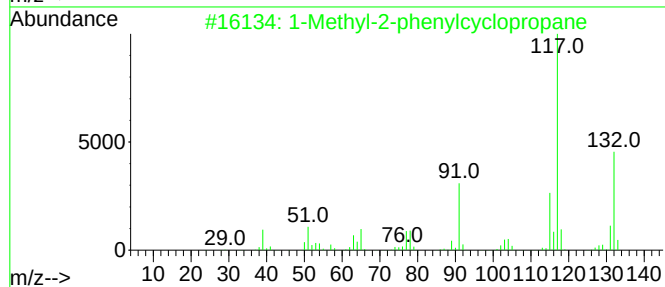
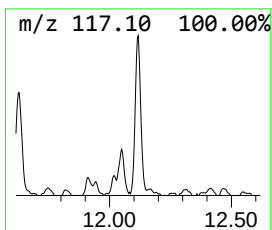
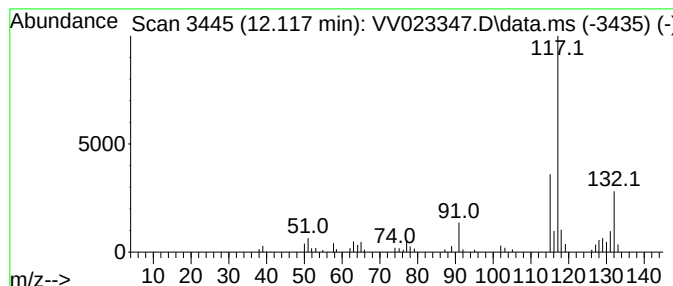
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 1-Methyl-2-phenylcyclopropane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	0.75 ug/L	70521	1,4-Dichlorobenzene-d4	11.252

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
2		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110921\
Data File : VV023347.D
Acq On : 10 Nov 2021 21:09
Operator : SY/MD
Sample : M4542-03
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGGE9

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
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Benzene, 1-ethy...	10.451	1.4	ug/L	126359	3	11.252	469039	5.0
Benzene, 1-ethy...	10.738	4.2	ug/L	396704	3	11.252	469039	5.0
Benzene, n-butyl-	11.059	0.6	ug/L	51154	3	11.252	469039	5.0
Benzeneacetalde...	11.091	0.7	ug/L	61515	3	11.252	469039	5.0
Benzene, 1,2,3-...	11.336	4.0	ug/L	377627	3	11.252	469039	5.0
Indane	11.532	4.2	ug/L	393167	3	11.252	469039	5.0
Benzene, 2-ethy...	11.914	0.6	ug/L	53376	3	11.252	469039	5.0
Benzene, 2-ethy...	12.017	0.7	ug/L	61612	3	11.252	469039	5.0
1-Methyl-2-phen...	12.117	0.8	ug/L	70521	3	11.252	469039	5.0