

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
 Data File : VU059922.D
 Acq On : 12 Jul 2024 15:25
 Operator : MD/SY
 Sample : P3217-04
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 YDZC7

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_U\Method\SFAMUTR061724WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU059922.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.599	84	96	132	rBV	15300466	18300791	19.27%	3.791%
2	2.637	407	419	458	rBV	348431	1126864	1.19%	0.233%
3	4.657	1028	1047	1136	rVB3	41392631	94948711	100.00%	19.670%
4	5.766	1383	1392	1408	rVB	618875	1253288	1.32%	0.260%
5	7.968	2062	2077	2108	rVB3	42082715	73758706	77.68%	15.280%
6	9.438	2516	2534	2552	rBV	5793964	9945320	10.47%	2.060%
7	9.563	2561	2573	2591	rBV	4405591	6918443	7.29%	1.433%
8	9.685	2591	2611	2627	rBV	17495033	28254634	29.76%	5.853%
9	10.094	2708	2738	2757	rBV	9585685	15065906	15.87%	3.121%
10	10.476	2843	2857	2869	rBV	567977	956808	1.01%	0.198%
11	10.898	2968	2988	3005	rBV	1183729	2057815	2.17%	0.426%
12	10.991	3005	3017	3035	rVV	5240165	11411139	12.02%	2.364%
13	11.084	3035	3046	3059	rVV2	3965558	6907697	7.28%	1.431%
14	11.155	3059	3068	3083	rVB	2875887	4640478	4.89%	0.961%
15	11.283	3097	3108	3126	rVB	3468085	5385077	5.67%	1.116%
16	11.460	3151	3163	3182	rVB	12576080	18911533	19.92%	3.918%
17	11.631	3197	3216	3231	rBV4	3013273	5842142	6.15%	1.210%
18	11.714	3231	3242	3256	rBV2	1194205	2579407	2.72%	0.534%
19	11.788	3256	3265	3273	rBV2	1065130	1928054	2.03%	0.399%
20	11.888	3285	3296	3316	rVB	9757570	16483295	17.36%	3.415%
21	12.087	3345	3358	3366	rBV	3310729	5379599	5.67%	1.114%
22	12.203	3377	3394	3413	rVB2	8060815	14732516	15.52%	3.052%
23	12.460	3460	3474	3478	rBV2	898655	1533526	1.62%	0.318%
24	12.486	3479	3482	3497	rVV	1032422	1572107	1.66%	0.326%
25	12.563	3497	3506	3513	rVV	1057135	1574805	1.66%	0.326%
26	12.676	3531	3541	3561	rVB2	1356348	2682344	2.83%	0.556%
27	12.868	3590	3601	3611	rVB3	682300	1117718	1.18%	0.232%
28	12.939	3611	3623	3639	rBV2	5257106	9637035	10.15%	1.996%
29	13.020	3639	3648	3658	rVB	1335890	2032512	2.14%	0.421%
30	13.286	3723	3731	3744	rVB	811307	1207541	1.27%	0.250%
31	13.399	3755	3766	3772	rBV	2620162	4104330	4.32%	0.850%
32	13.444	3772	3780	3789	rVV3	3493606	5922047	6.24%	1.227%
33	13.512	3789	3801	3822	rVB	18009127	28204875	29.71%	5.843%
34	13.618	3823	3834	3846	rVB	2308083	3695050	3.89%	0.765%
35	13.730	3847	3869	3880	rBV2	6368458	11762669	12.39%	2.437%
36	13.833	3891	3901	3914	rBV4	737481	1326832	1.40%	0.275%

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Title : TRACE VOA SFAM1.0

37	14.077	3964	3977	3998	rVB	7075077	11382147	11.99%	2.358%
38	14.479	4092	4102	4120	rVB	656167	1223733	1.29%	0.254%
39	14.669	4148	4161	4183	rBV4	918337	2366140	2.49%	0.490%
40	14.868	4213	4223	4235	rVB2	776498	1246297	1.31%	0.258%
41	15.010	4256	4267	4278	rBV2	547316	960926	1.01%	0.199%
42	15.216	4317	4331	4344	rBV	14650276	22759364	23.97%	4.715%
43	15.402	4376	4389	4401	rBV	8747542	13682573	14.41%	2.835%
44	16.254	4642	4654	4680	rVB	1310597	2325711	2.45%	0.482%
45	16.399	4689	4699	4706	rBV	1399374	2255350	2.38%	0.467%
46	16.437	4706	4711	4727	rVB	880583	1351273	1.42%	0.280%

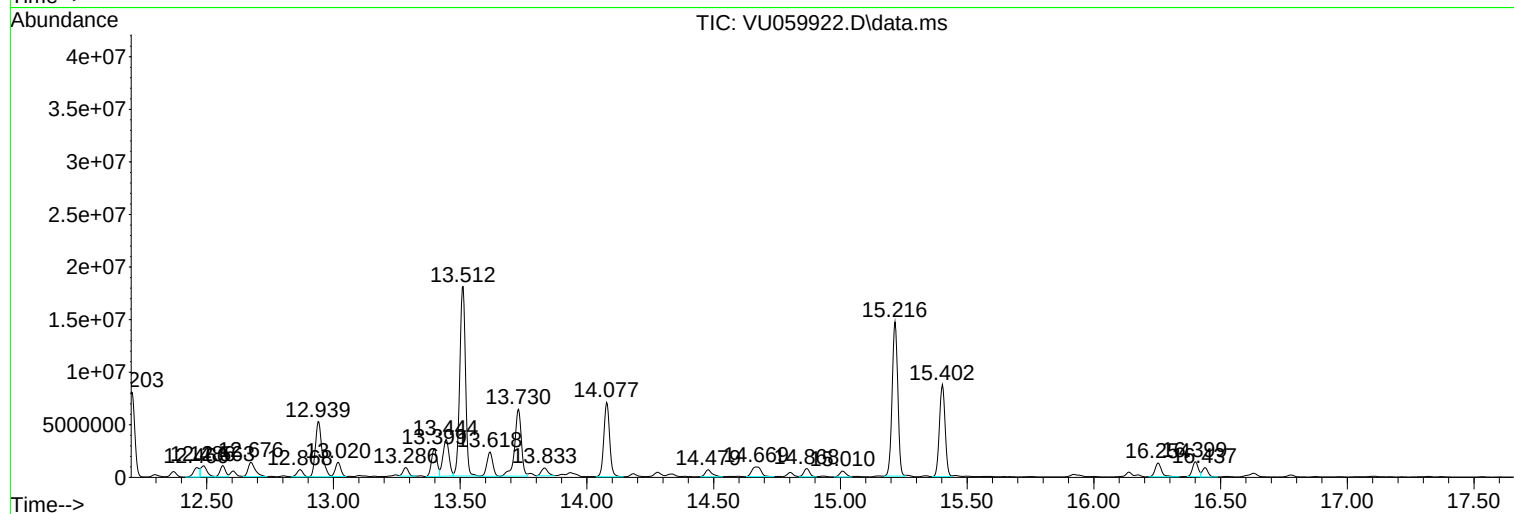
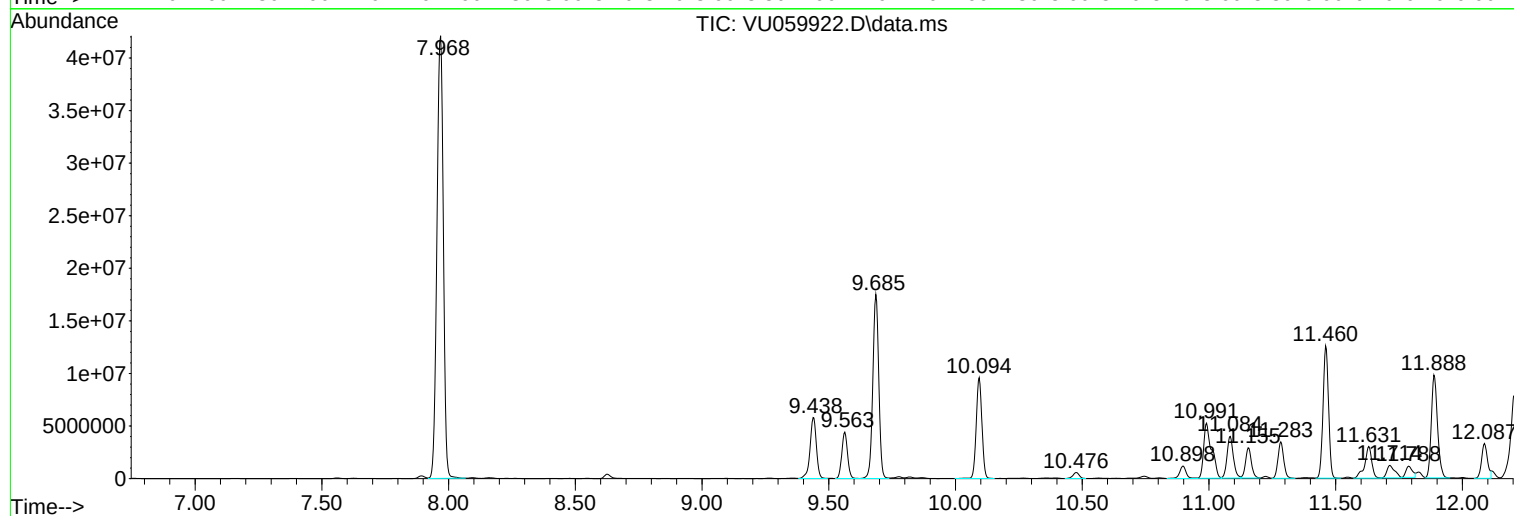
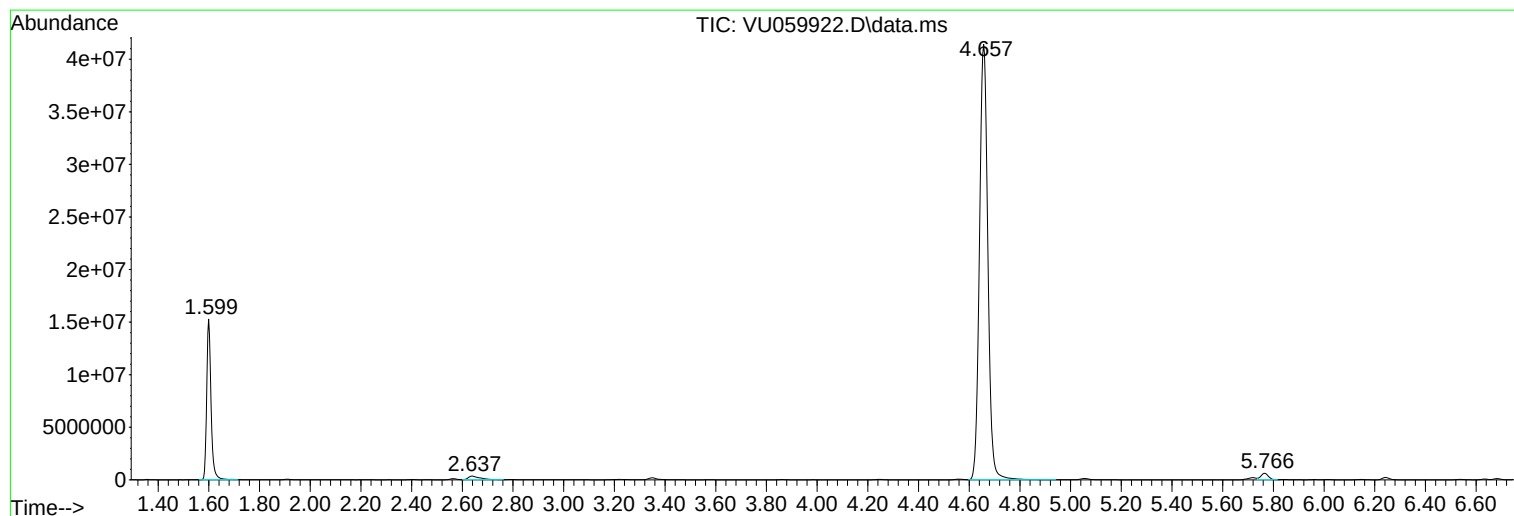
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Instrument :
MSVOA_U
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YDZC7

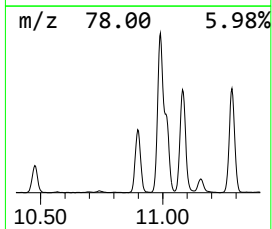
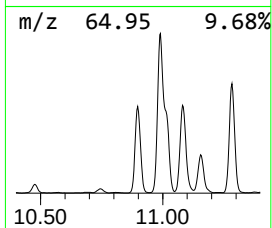
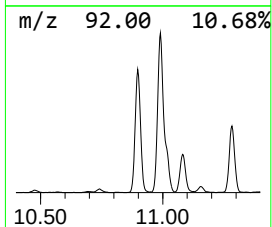
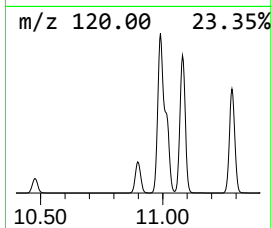
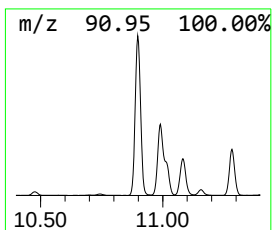
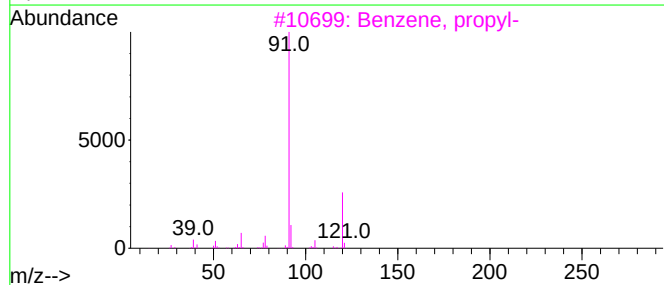
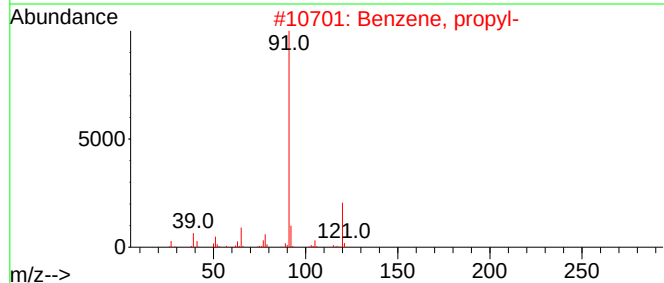
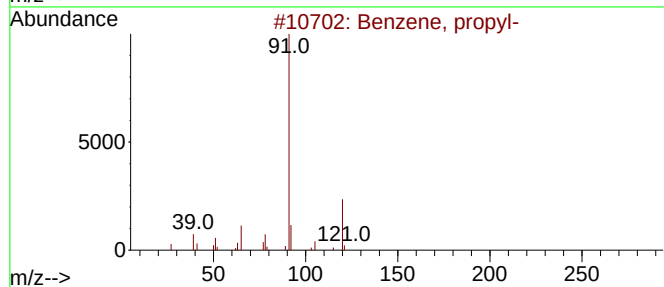
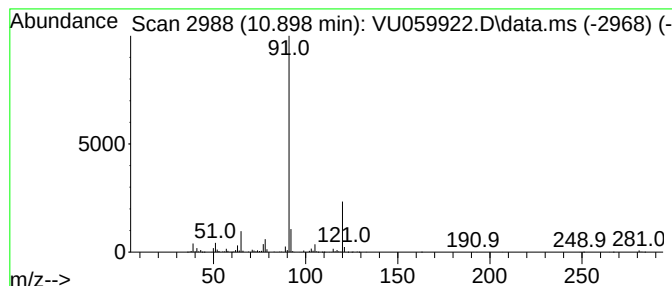
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Benzene, propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.898	5.34 ug/L	2057820	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
5			N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	64



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Instrument :
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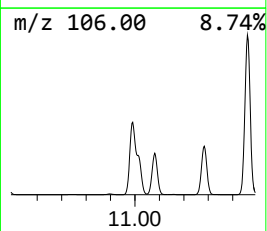
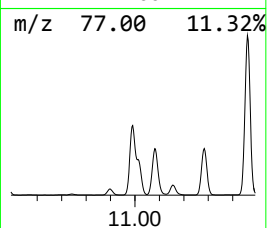
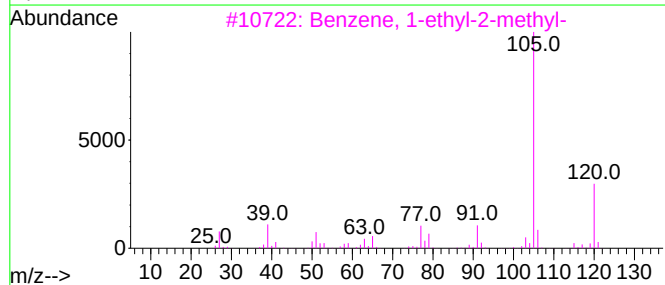
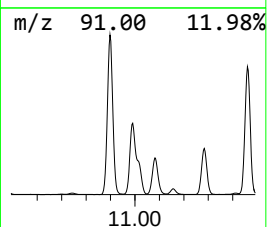
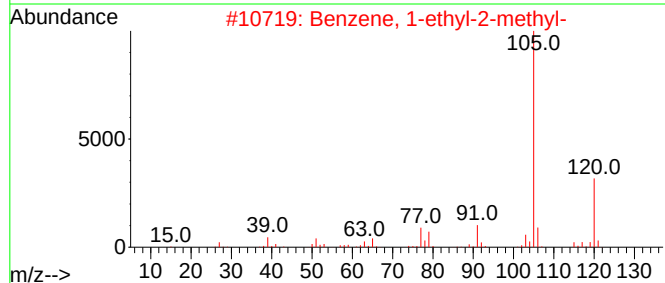
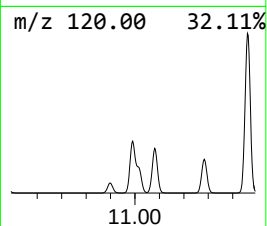
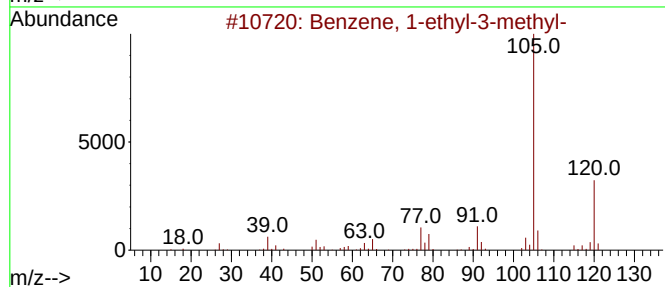
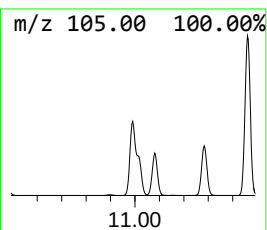
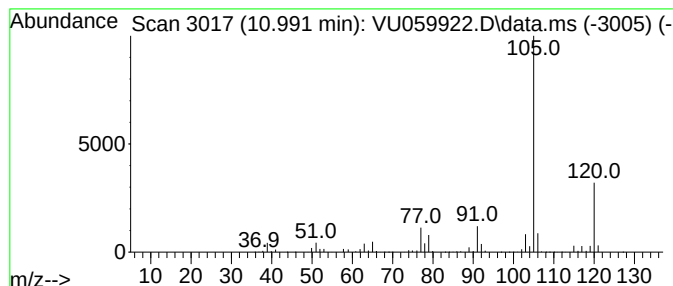
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.991	29.59 ug/L	11411100	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
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Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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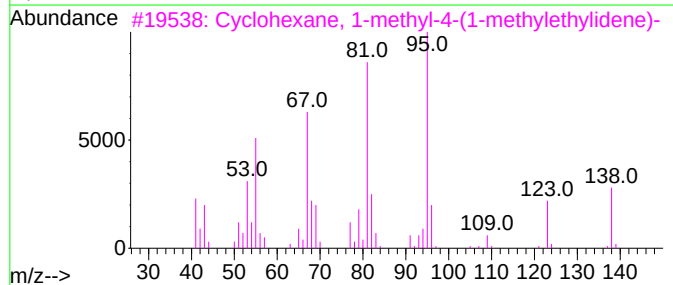
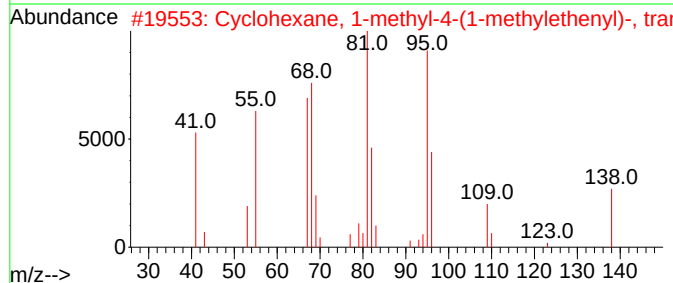
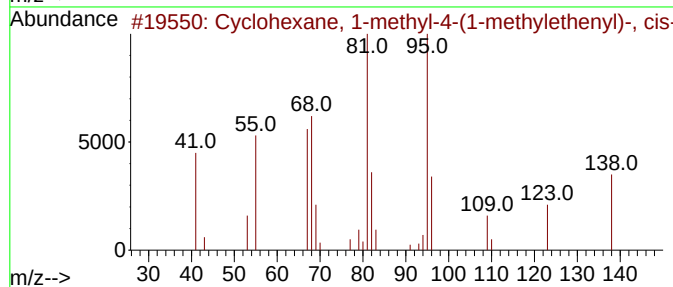
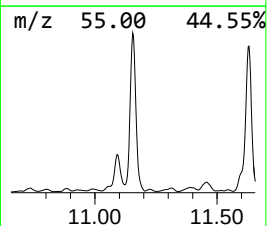
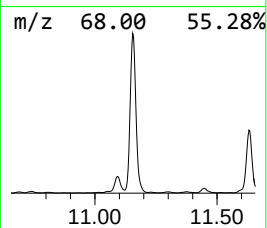
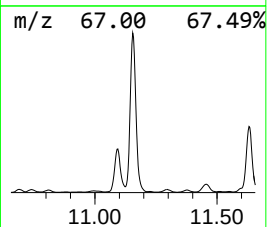
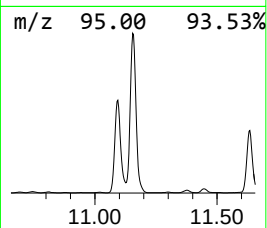
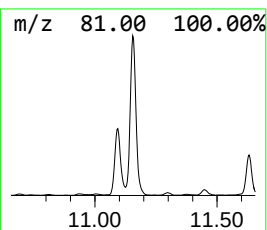
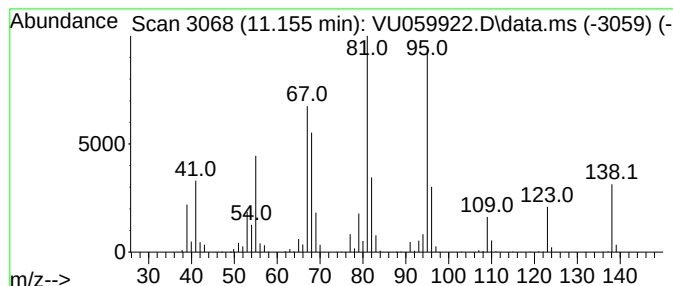
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Cyclohexane, 1-methyl-4-(1-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.155	12.03 ug/L	4640480	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001879-07-8	98
2			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-25-0	87
3			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-27-2	83
4			Cyclohexene, 4-methyl-1-(1-methy...	138	C10H18	000500-00-5	81
5			Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	018968-23-5	74



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Instrument :
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ClientSampleId :
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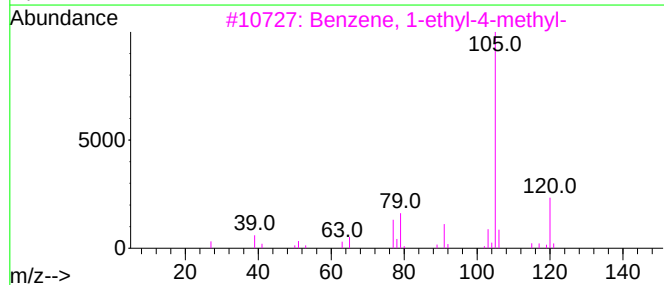
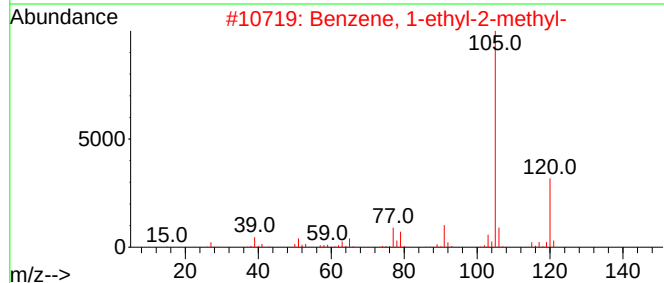
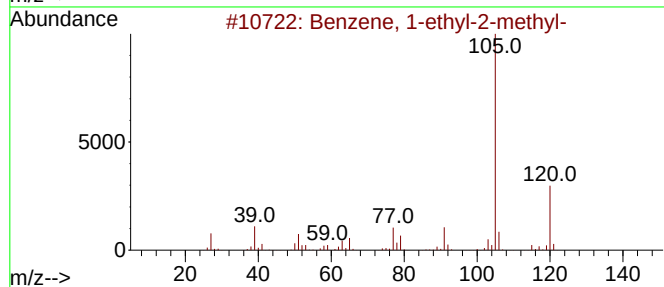
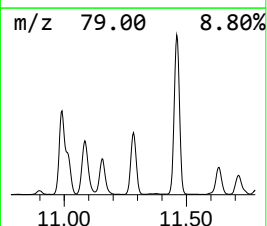
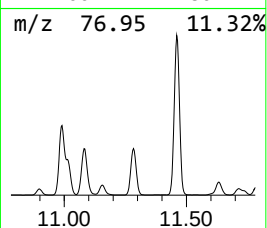
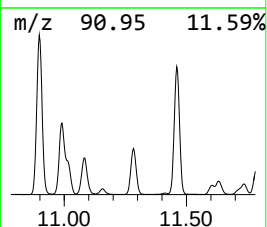
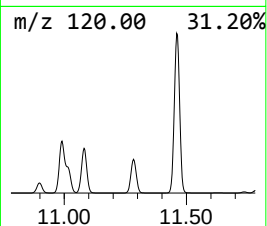
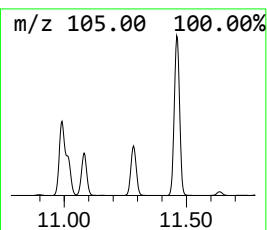
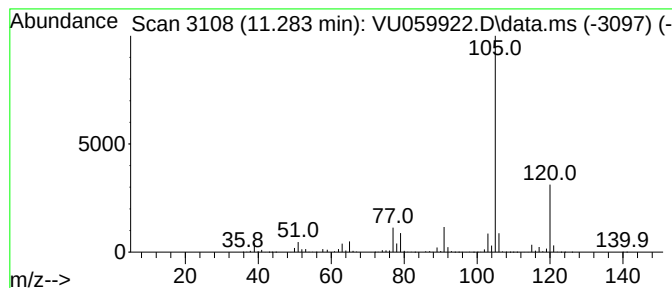
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.283	13.97 ug/L	5385080	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



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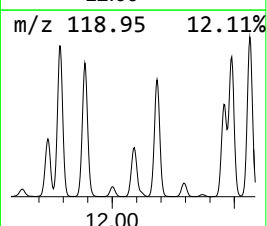
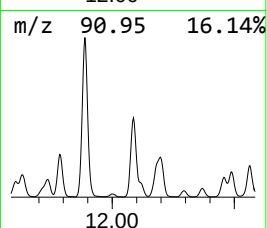
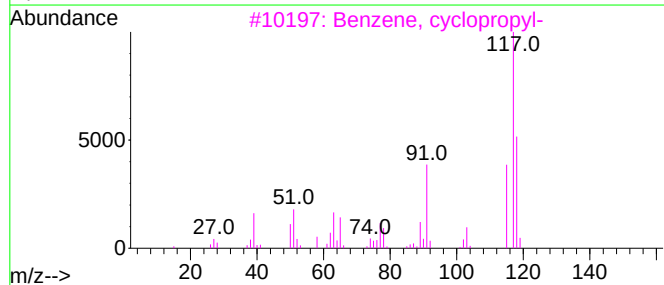
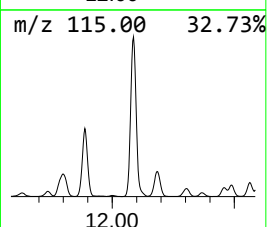
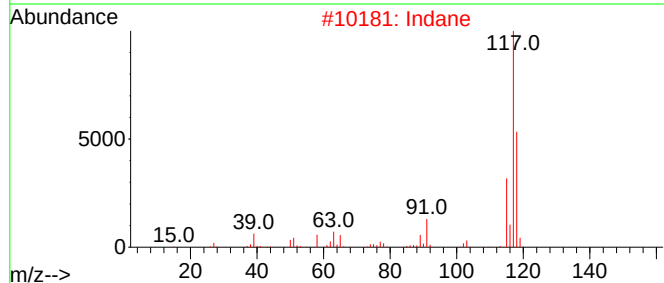
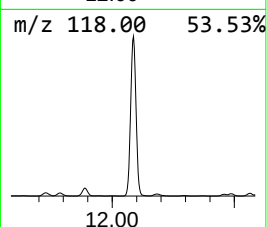
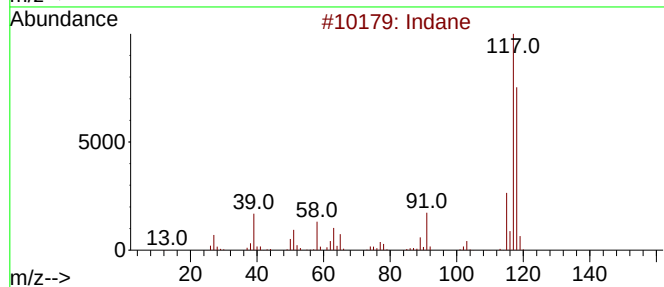
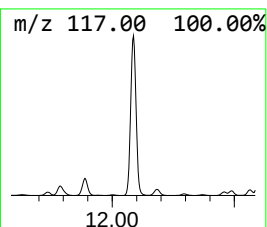
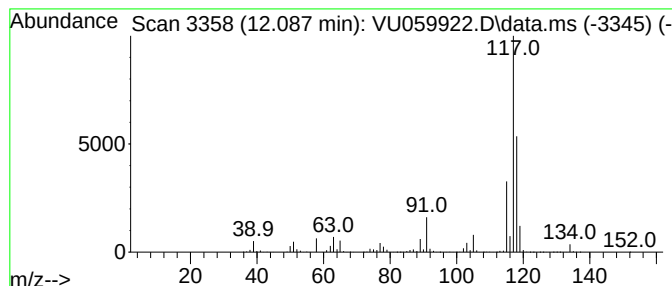
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.087	13.95 ug/L	5379600	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	81
2	Indane			118	C9H10	000496-11-7	81
3	Benzene, cyclopropyl-			118	C9H10	000873-49-4	81
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	74
5	Benzene, cyclopropyl-			118	C9H10	000873-49-4	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

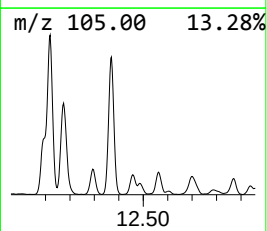
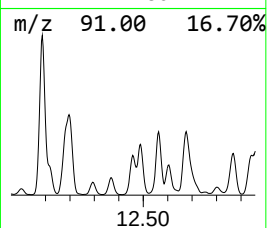
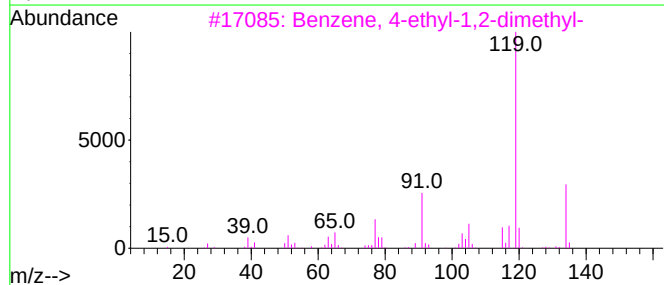
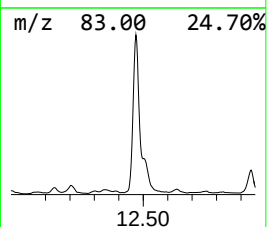
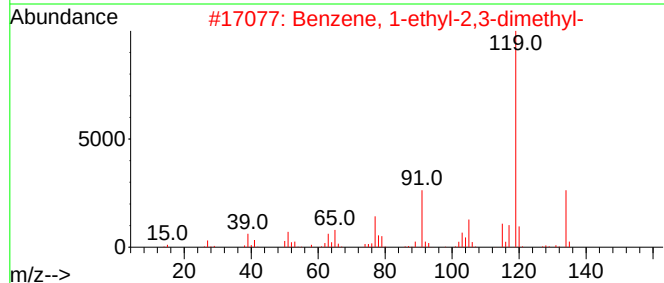
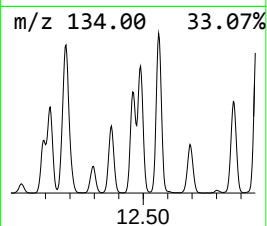
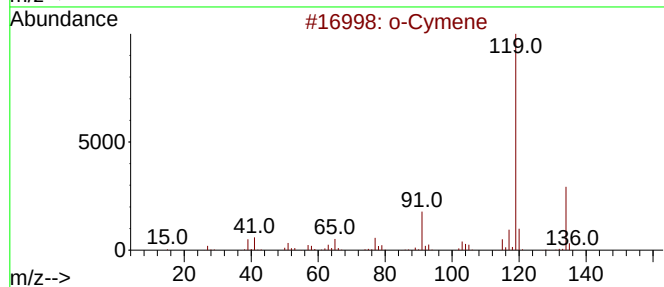
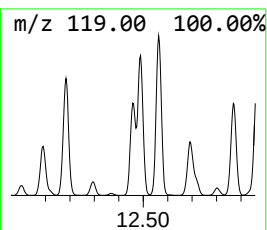
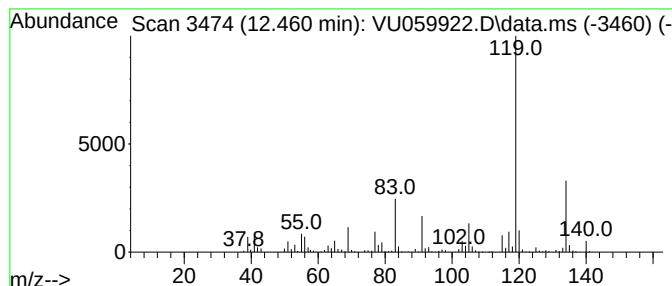
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 o-Cymene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.460	3.98 ug/L	1533530	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	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	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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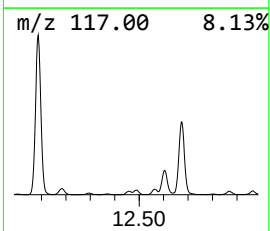
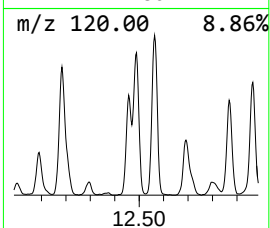
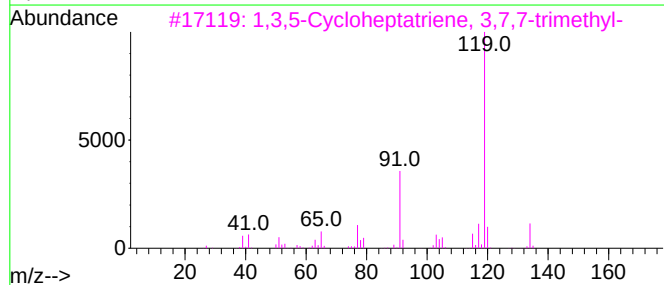
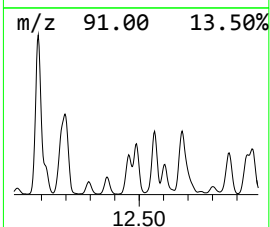
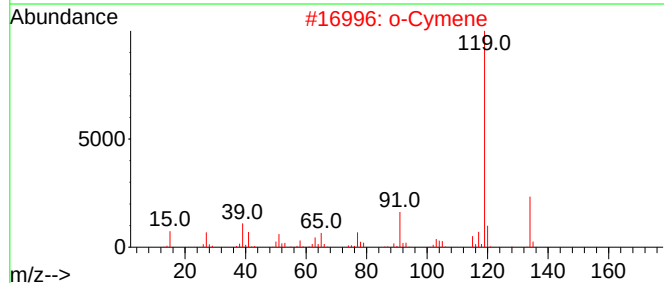
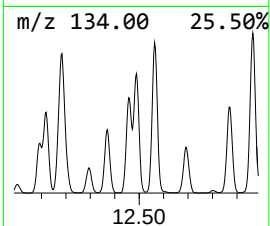
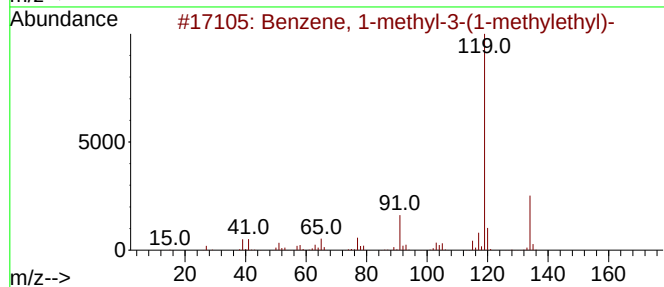
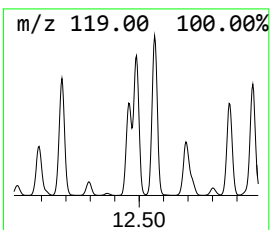
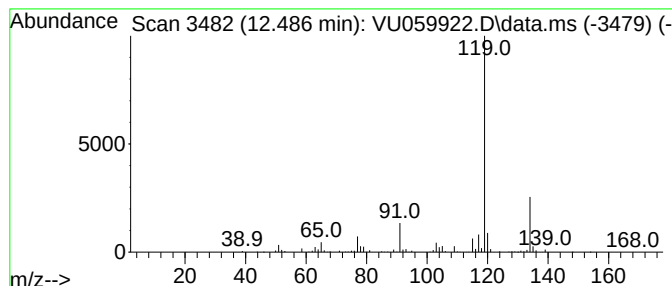
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, 1-methyl-3-(1-meth... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	4.08 ug/L	1572110	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	94
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

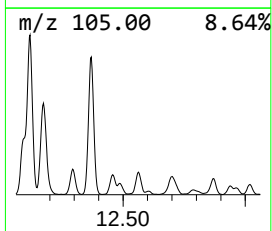
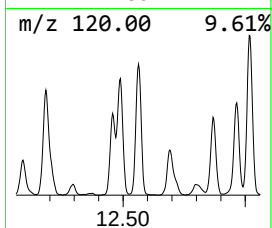
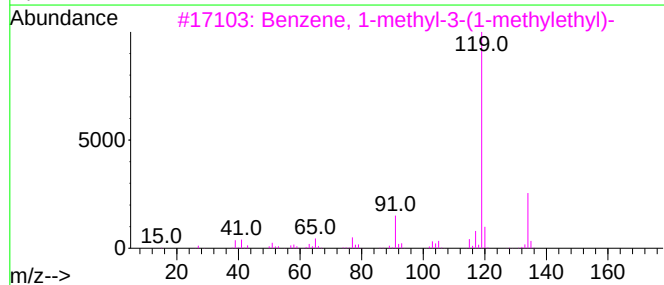
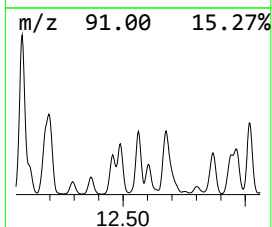
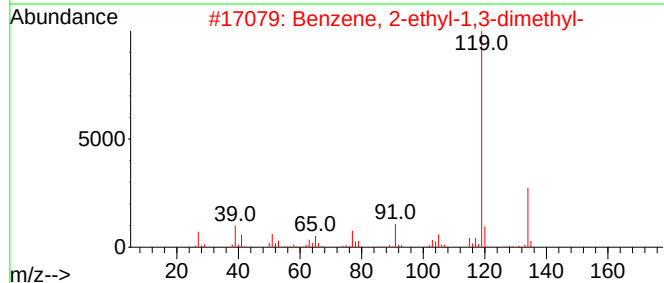
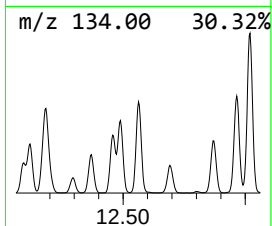
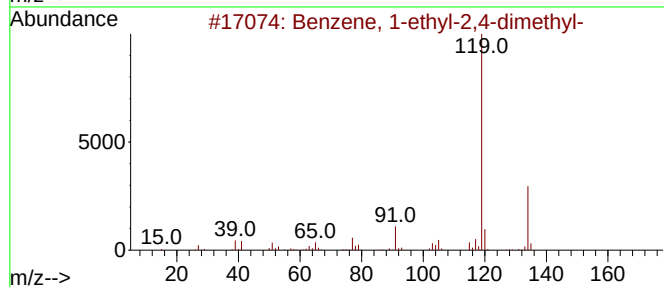
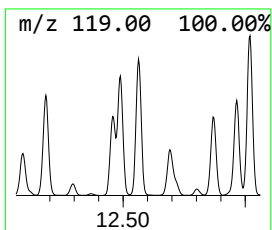
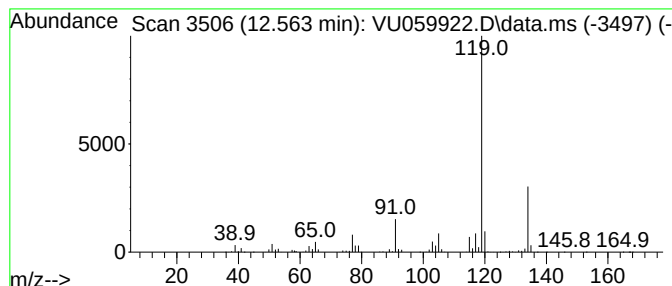
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.563	4.08 ug/L	1574810	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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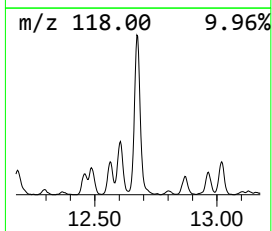
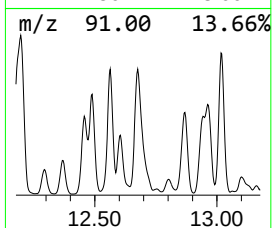
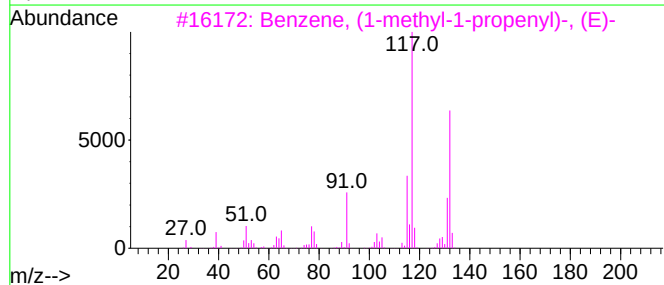
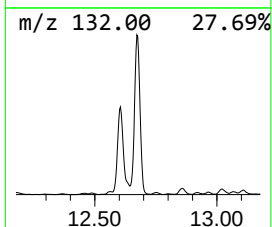
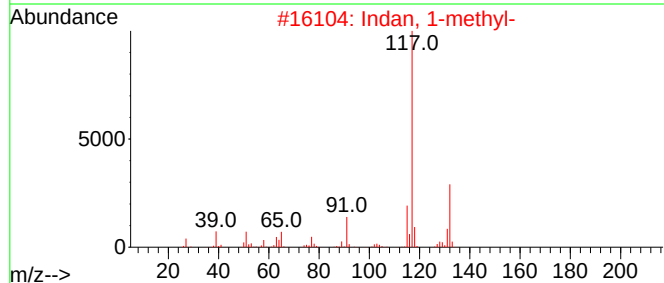
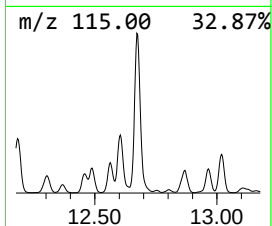
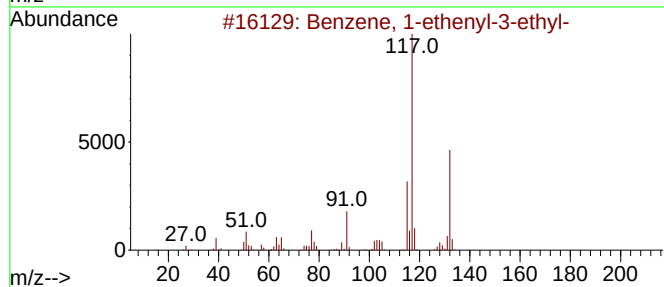
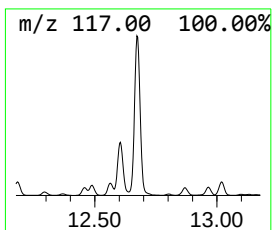
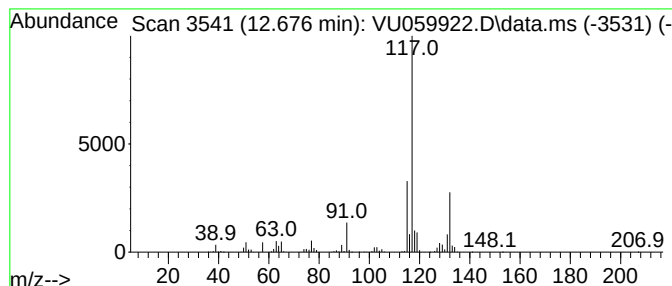
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.676	6.96 ug/L	2682340	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2			Indan, 1-methyl-	132	C10H12	000767-58-8	87
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

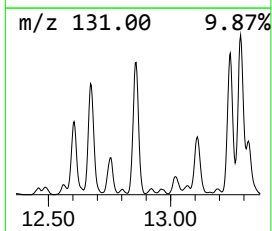
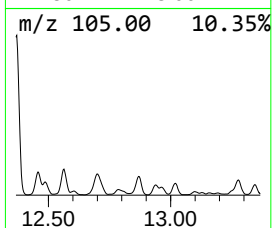
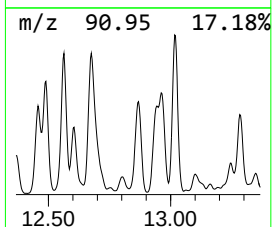
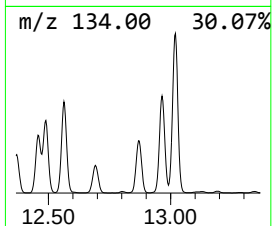
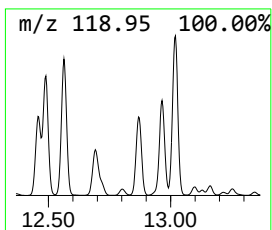
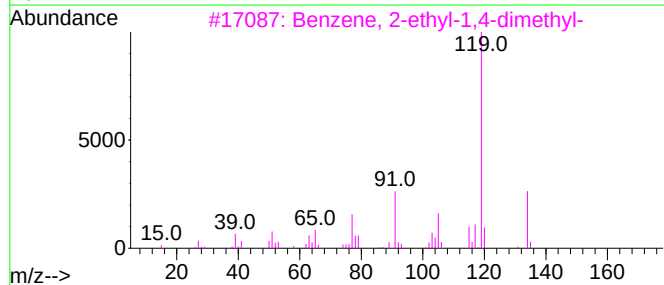
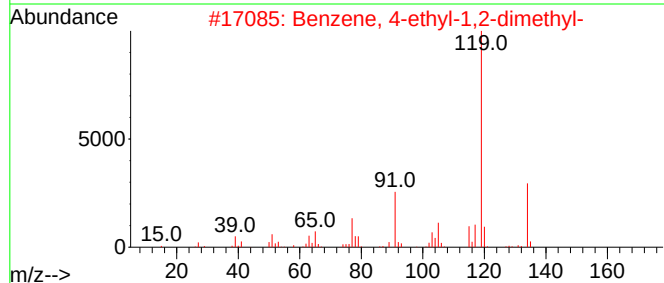
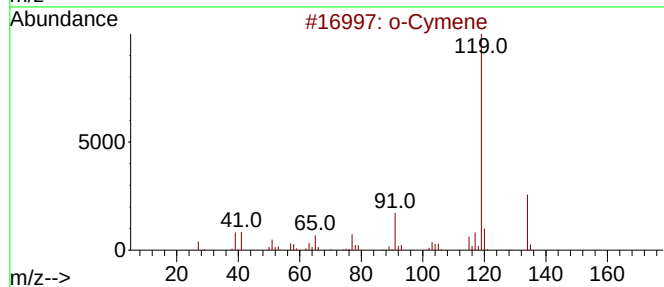
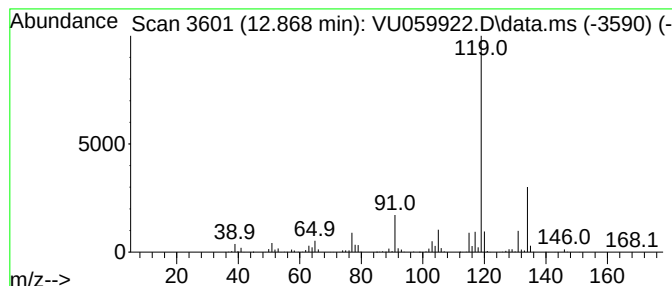
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.868	2.90 ug/L	1117720	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

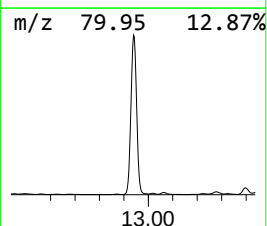
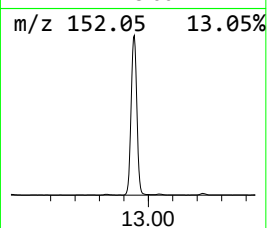
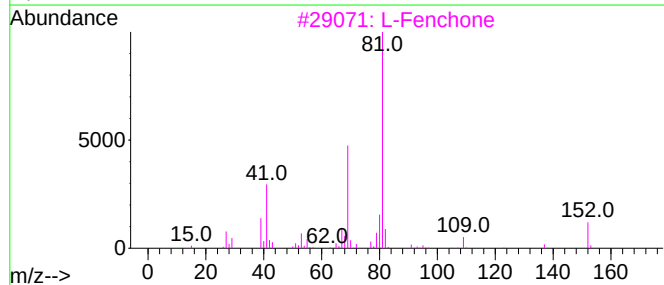
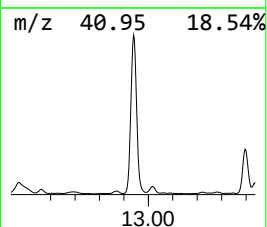
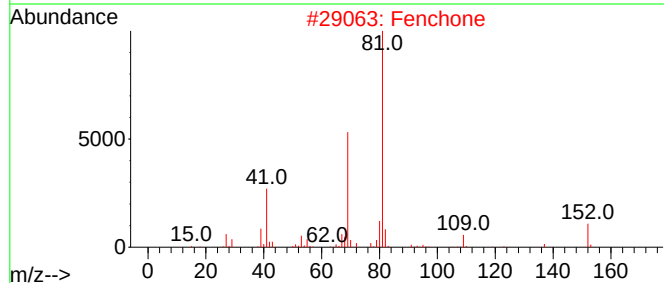
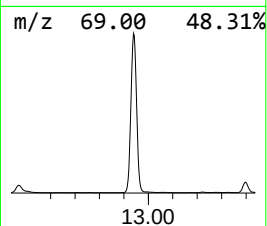
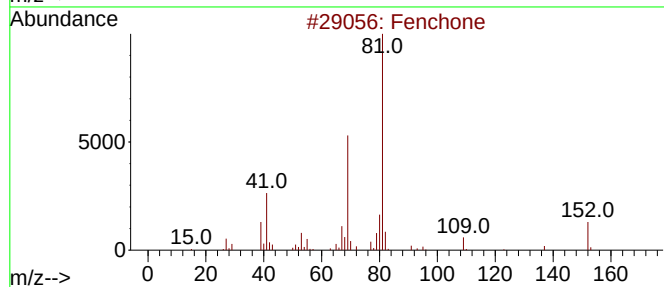
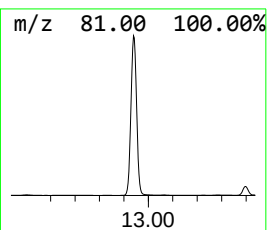
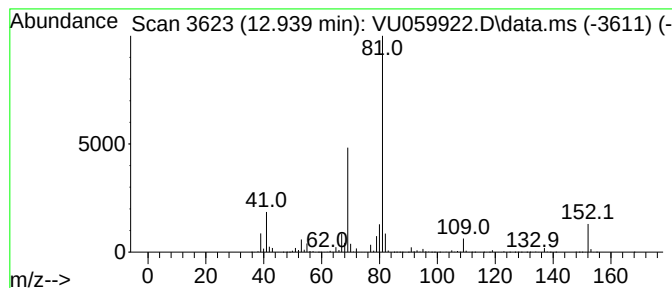
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Fenchone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.939	24.99 ug/L	9637040	1,4-Dichlorobenzene-d4	11.804

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			L-Fenchone	152	C10H16O	007787-20-4	95
4			D-Fenchone	152	C10H16O	004695-62-9	94
5			L-Fenchone	152	C10H16O	007787-20-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

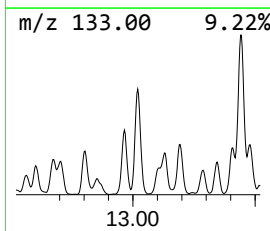
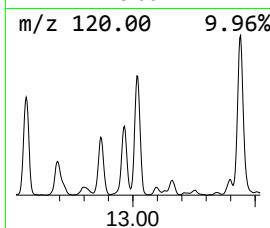
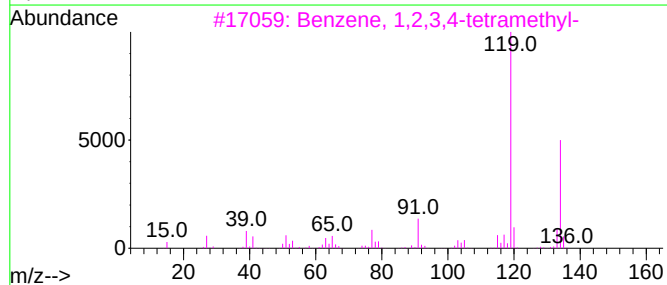
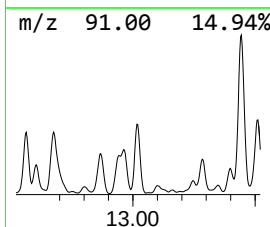
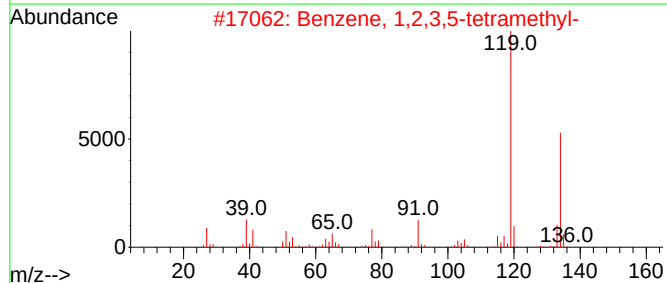
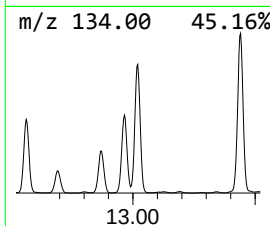
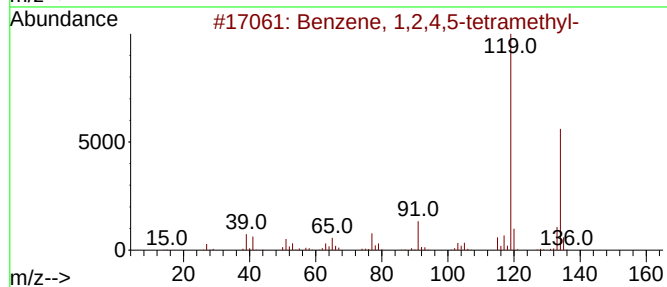
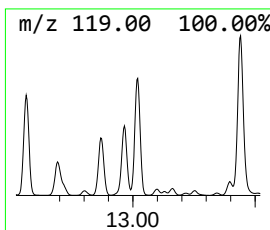
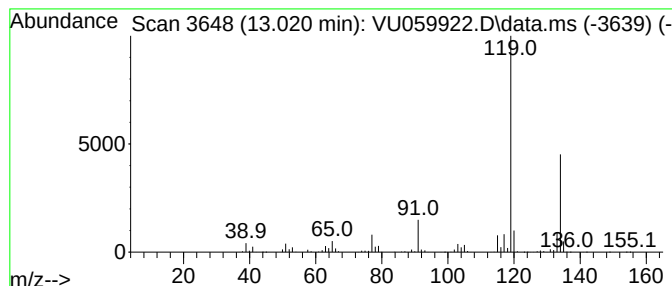
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.020	5.27 ug/L	2032510	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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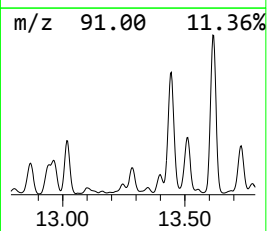
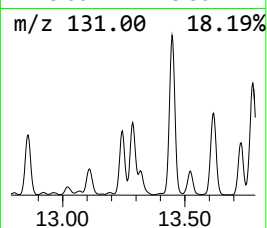
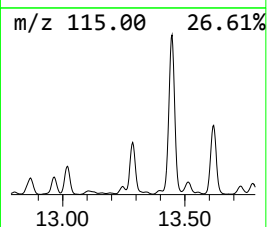
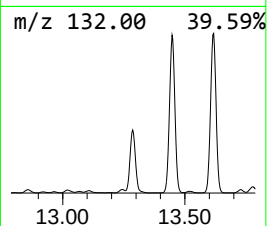
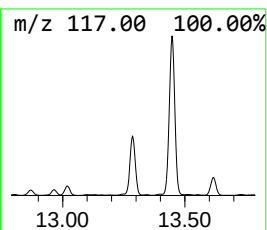
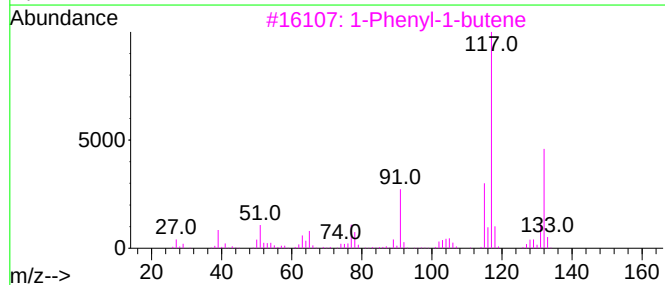
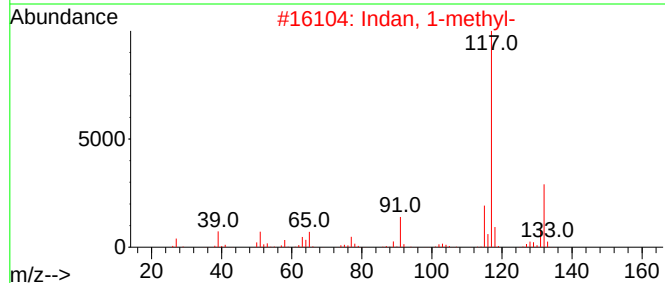
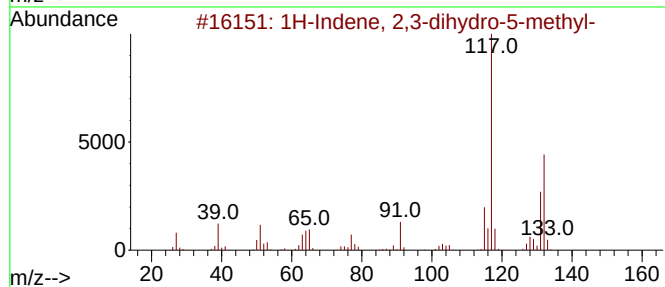
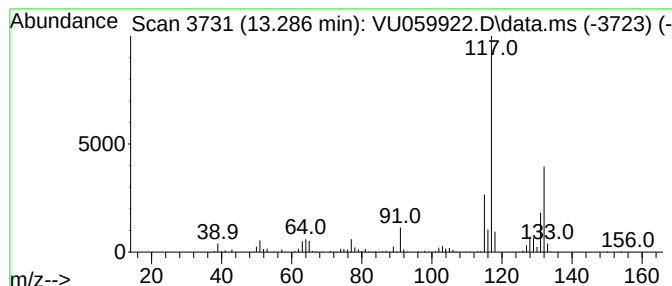
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.287	3.13 ug/L	1207540	1,4-Dichlorobenzene-d4	11.804

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Indan, 1-methyl-	132	C10H12	000767-58-8	91
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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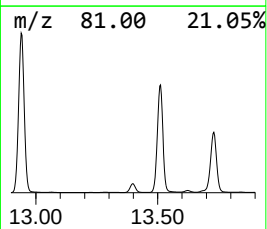
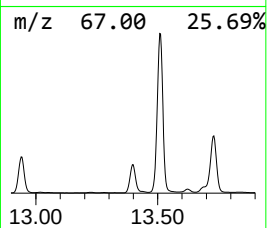
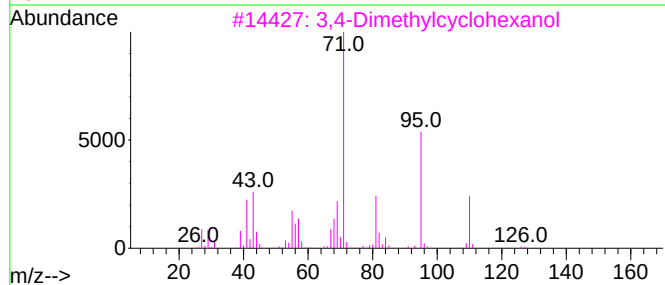
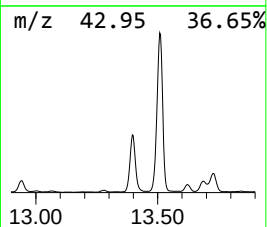
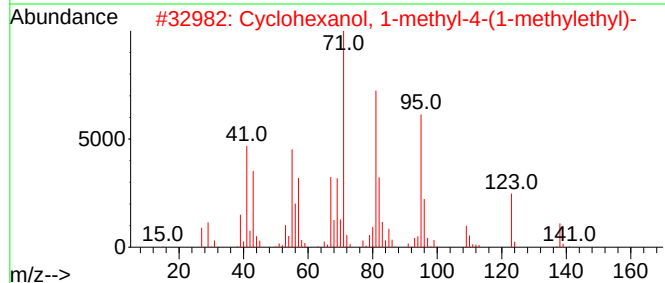
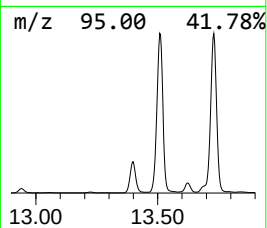
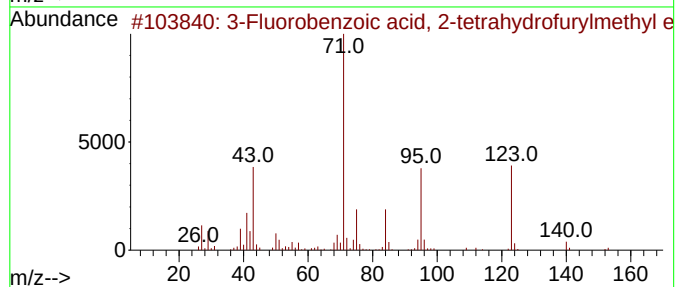
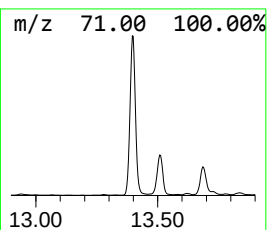
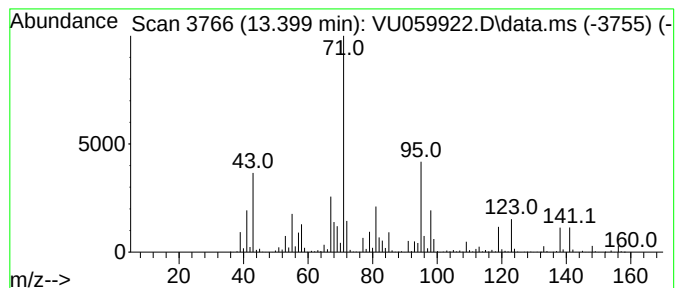
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 16 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.399	10.64 ug/L	4104330	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Fluorobenzoic acid, 2-tetrahyd...	224	C12H13F03	1000279-04-4	47
2			Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	021129-27-1	45
3			3,4-Dimethylcyclohexanol	128	C8H16O	005715-23-1	43
4			Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	003623-52-7	42
5			Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	023283-97-8	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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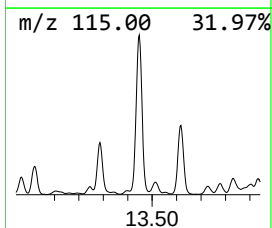
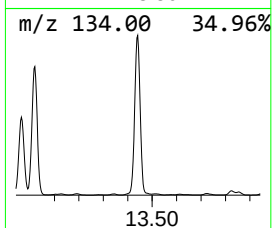
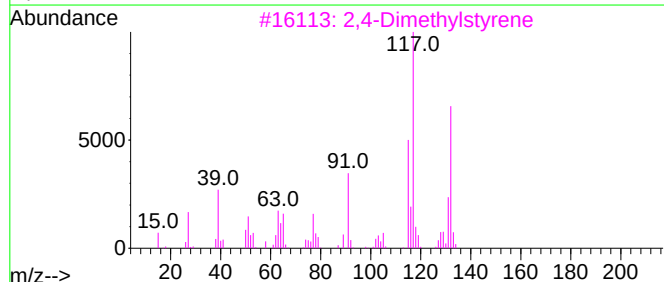
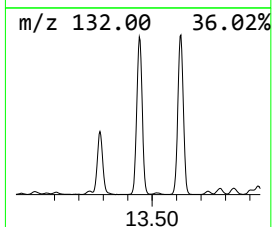
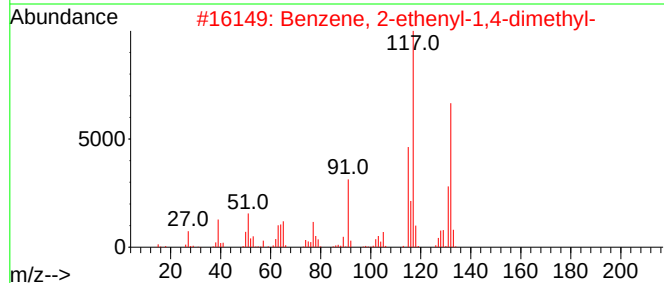
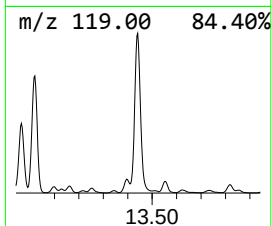
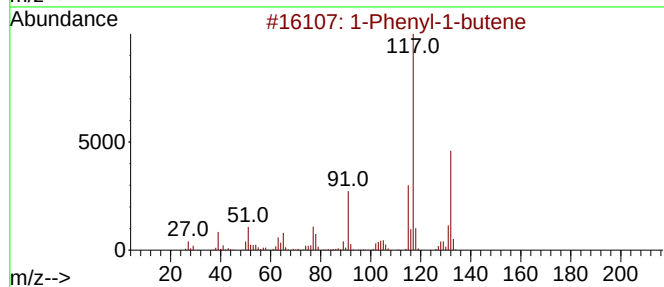
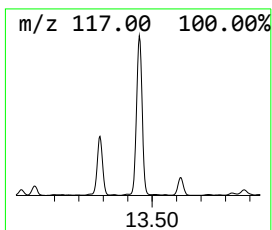
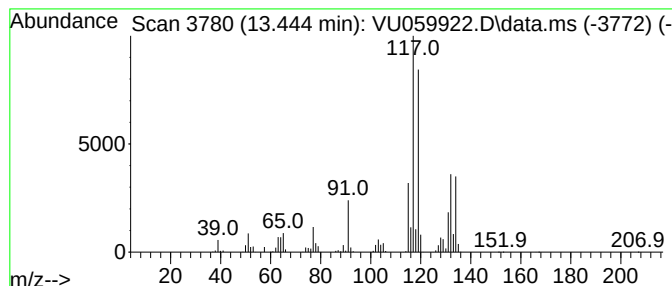
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 17 1-Phenyl-1-butene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.444	15.36 ug/L	5922050	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
3			2,4-Dimethylstyrene	132	C10H12	002234-20-0	83
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	78
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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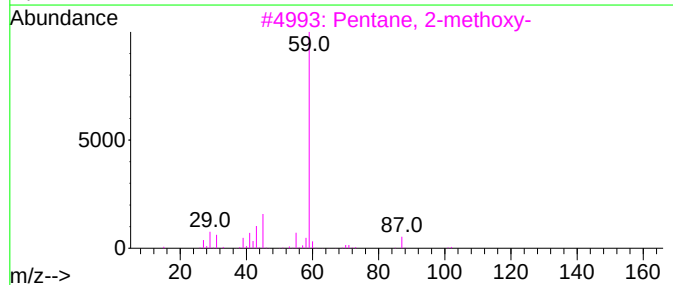
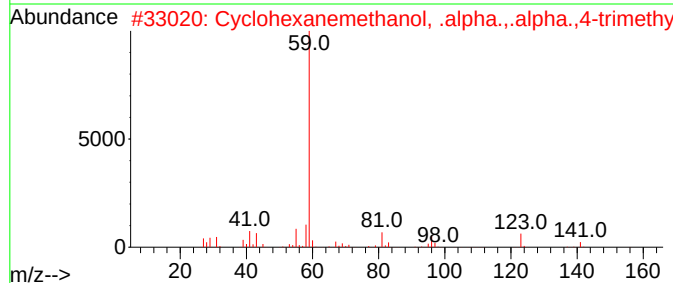
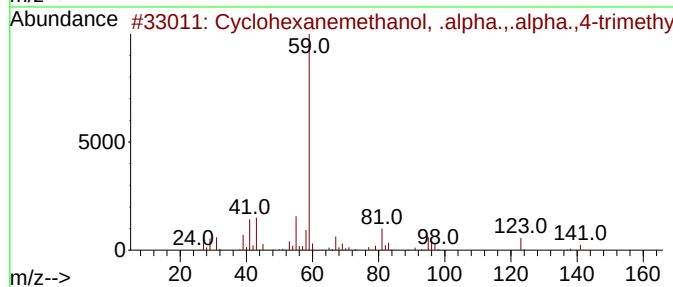
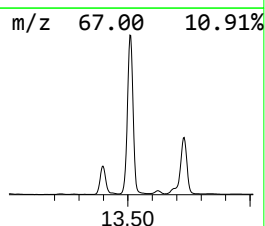
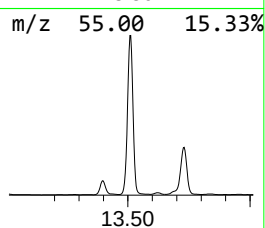
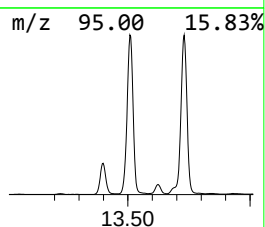
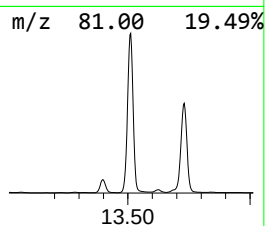
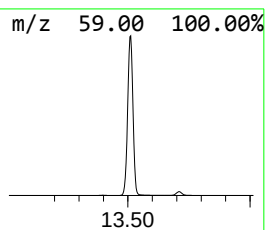
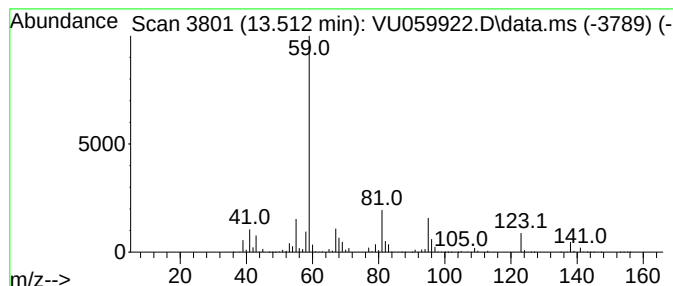
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 18 Cyclohexanemethanol, .alpha... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.512	73.14 ug/L	28204900	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	78
2			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	64
3			Pentane, 2-methoxy-	102	C6H14O	006795-88-6	53
4			Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	53
5			Erythro-2-methyl-3,4-dibromo-2-b...	244	C5H10Br2O	1000288-19-3	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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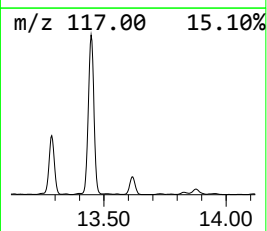
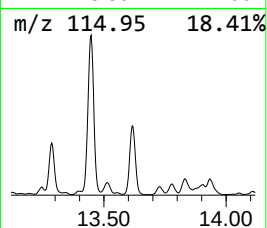
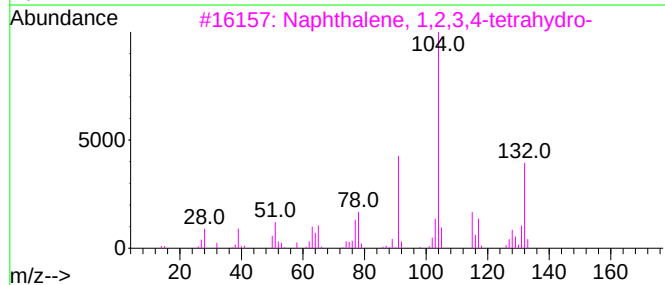
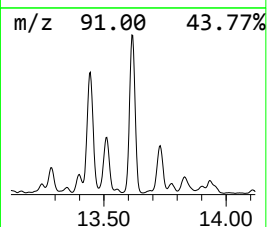
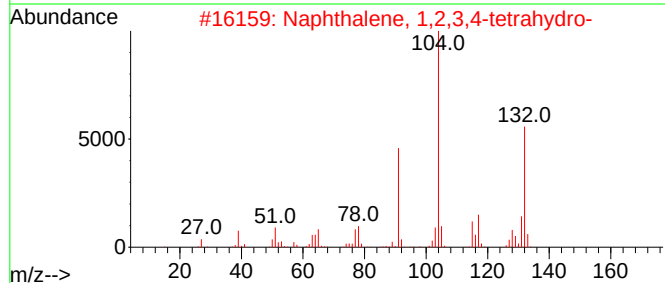
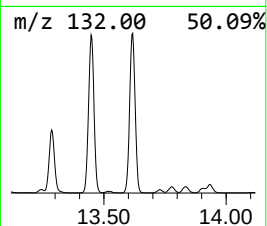
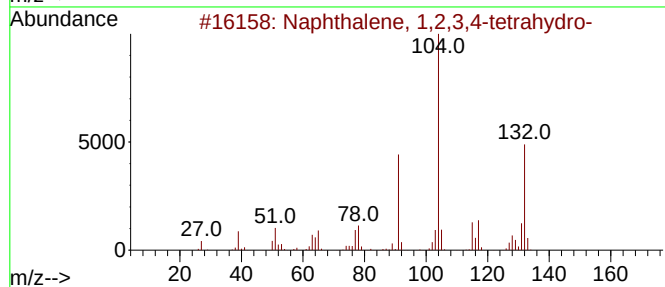
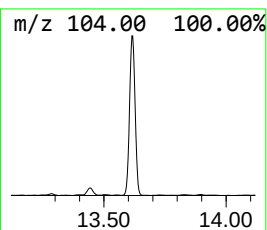
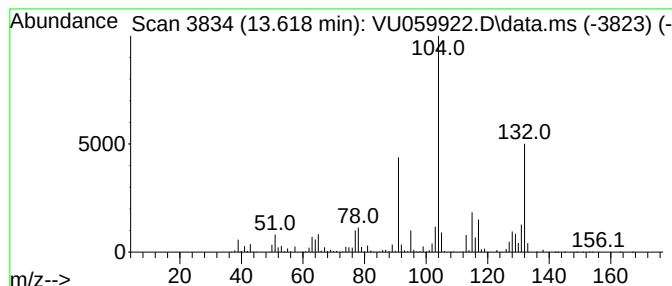
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 19 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.618	9.58 ug/L	3695050	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

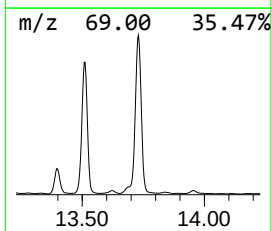
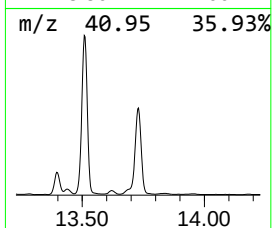
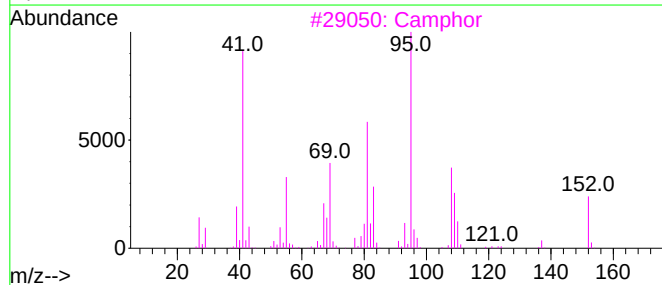
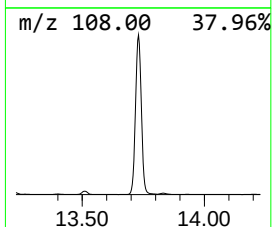
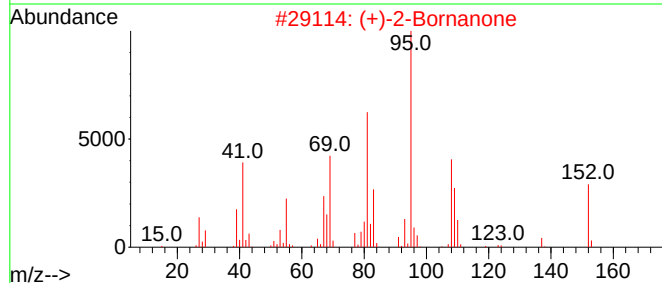
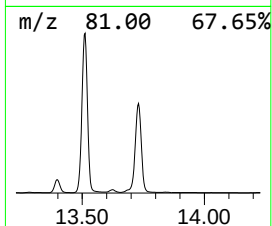
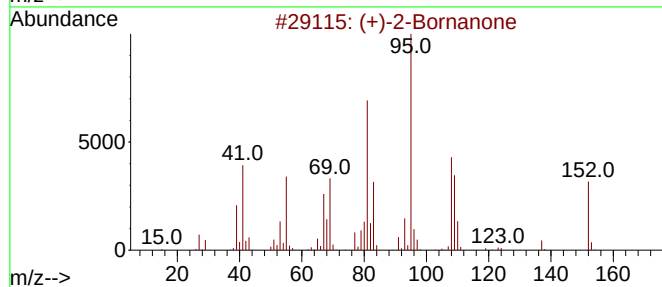
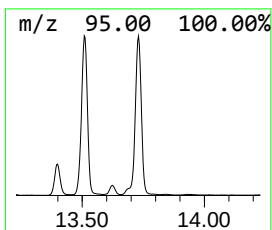
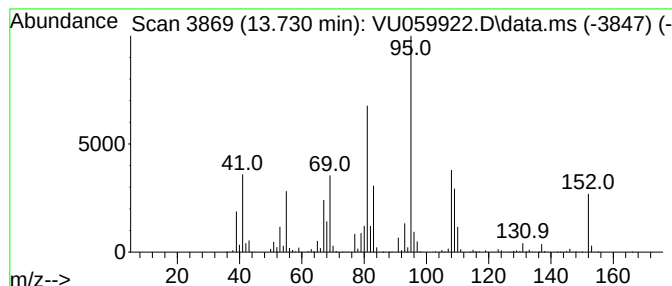
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 20 (+)-2-Bornanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.730	30.50 ug/L	11762700	1,4-Dichlorobenzene-d4			11.804
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone		152	C10H16O	000464-49-3	98
2	(+)-2-Bornanone		152	C10H16O	000464-49-3	98
3	Camphor		152	C10H16O	000076-22-2	98
4	(+)-2-Bornanone		152	C10H16O	000464-49-3	98
5	Camphor		152	C10H16O	000076-22-2	98



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

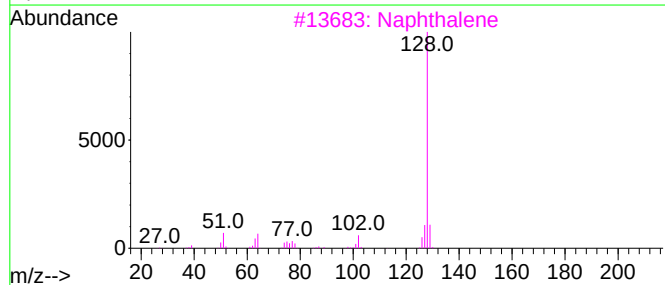
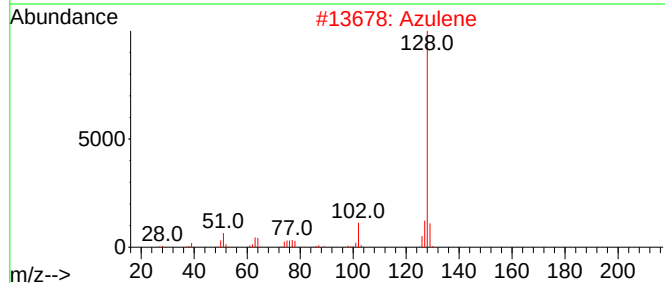
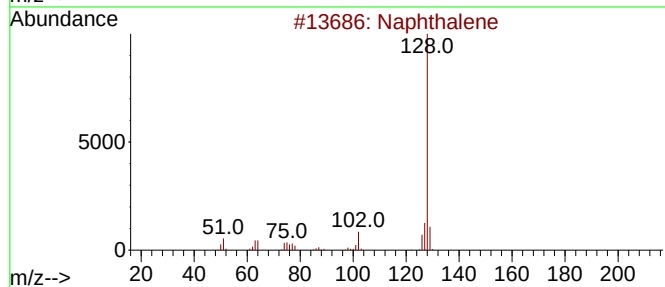
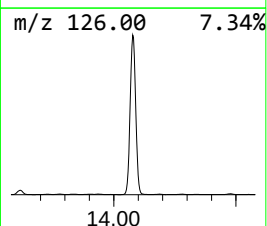
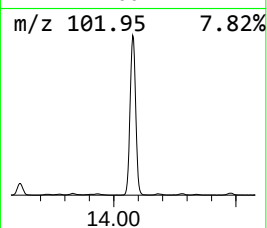
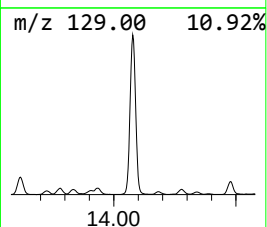
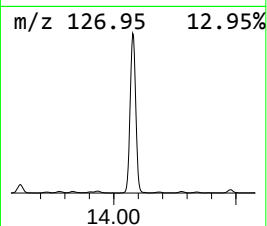
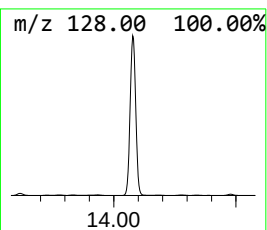
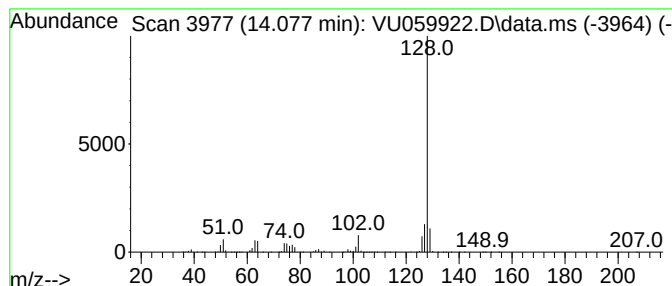
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 21 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.078	29.52 ug/L	11382100	1,4-Dichlorobenzene-d4	11.804

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

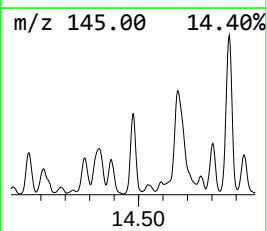
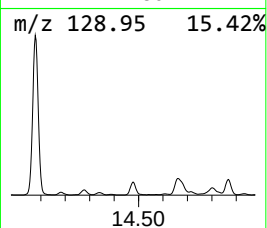
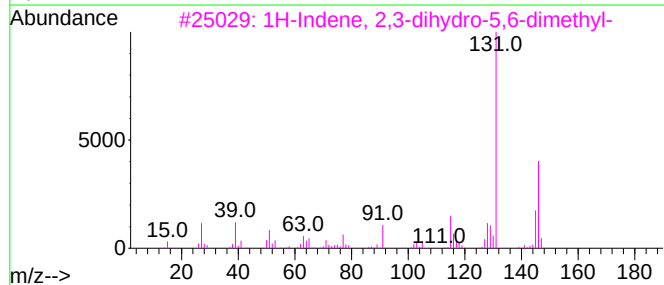
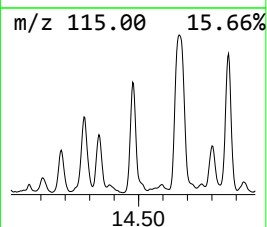
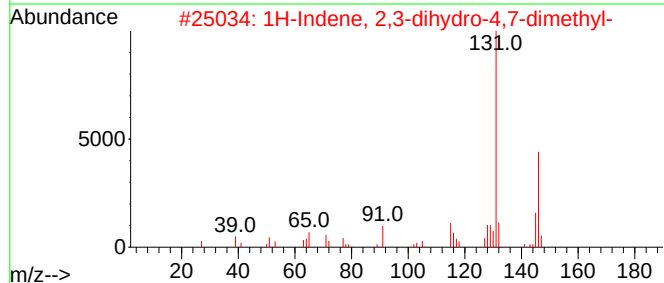
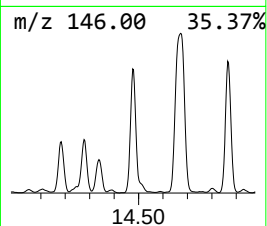
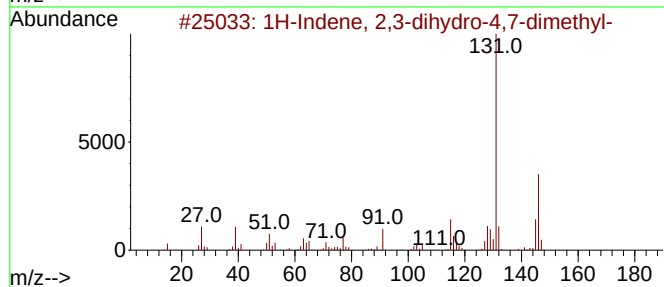
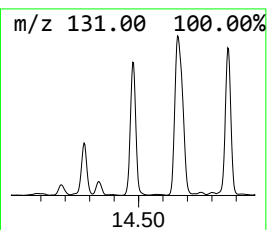
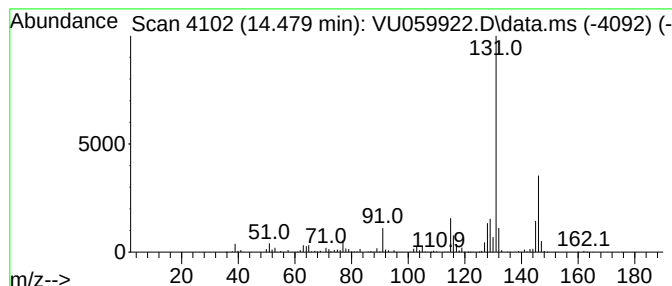
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 22 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.479	3.17 ug/L	1223730	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	97		
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96		
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	96		
4	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	95		
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

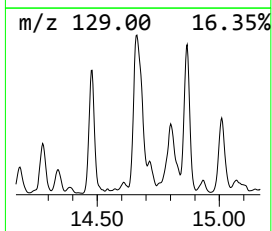
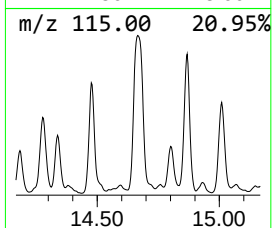
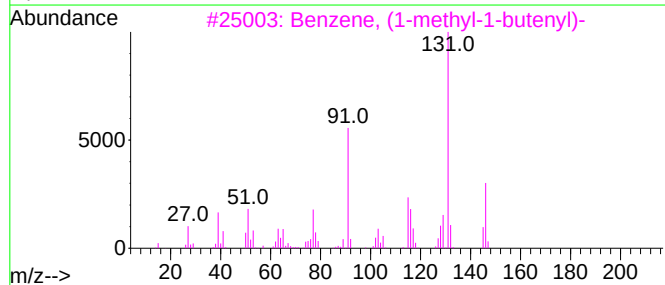
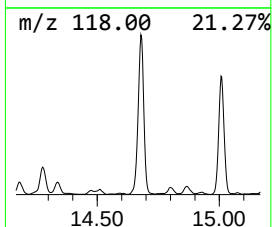
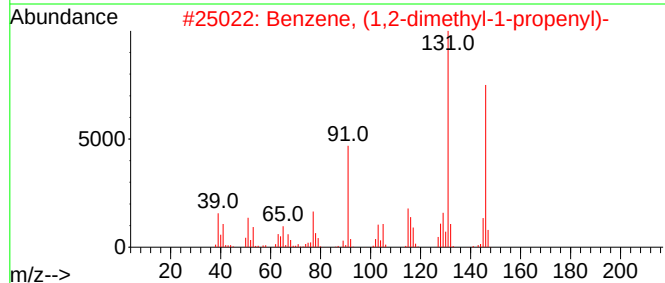
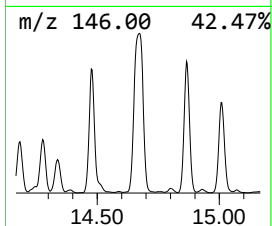
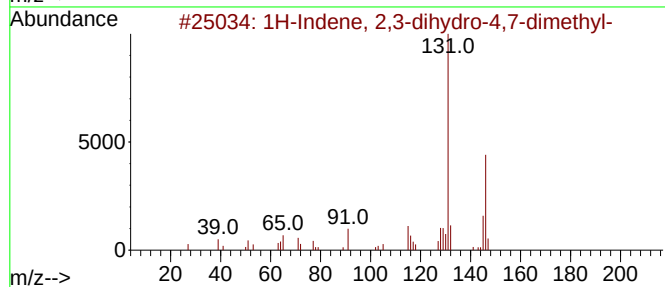
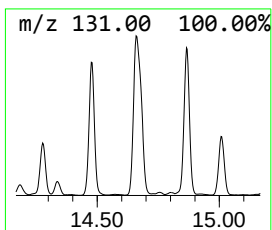
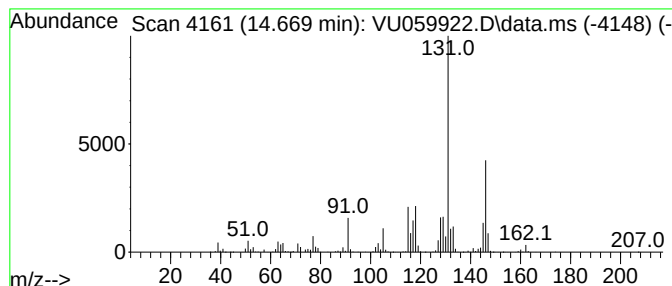
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 23 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.669	6.14 ug/L	2366140	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

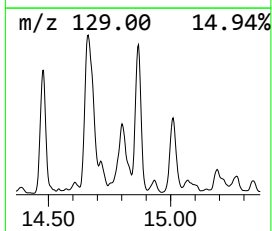
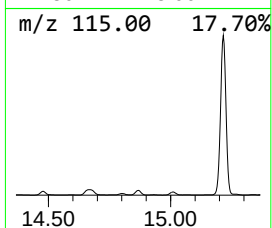
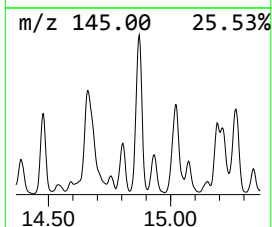
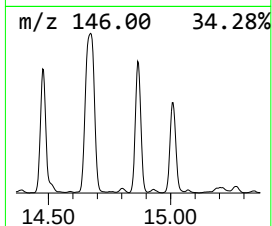
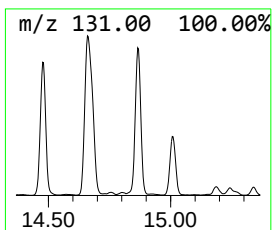
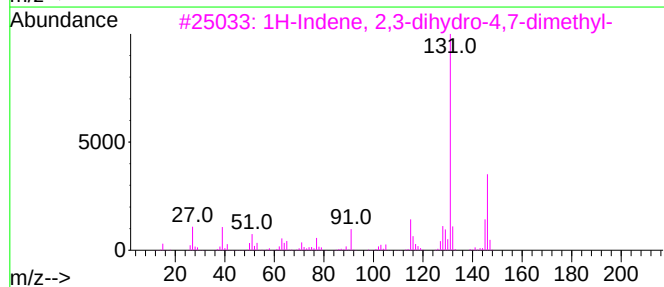
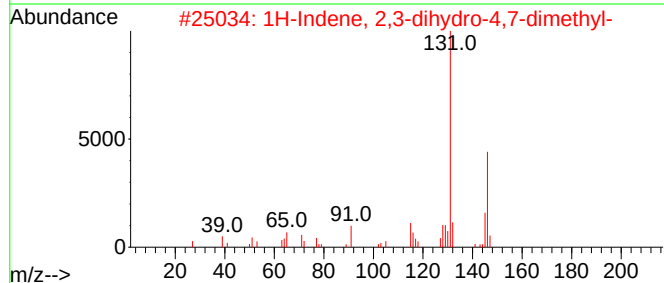
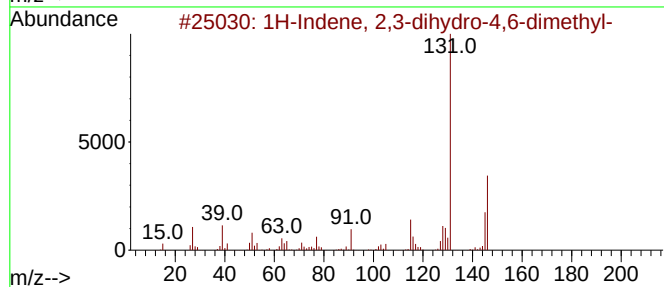
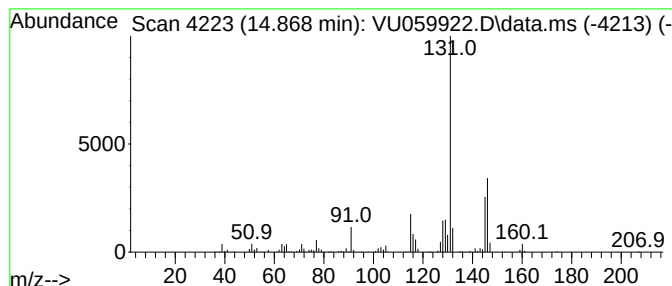
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 24 1H-Indene, 2,3-dihydro-4,6-... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	3.23 ug/L	1246300	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	94		
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93		
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93		
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90		
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

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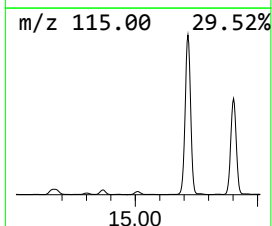
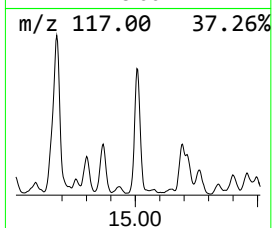
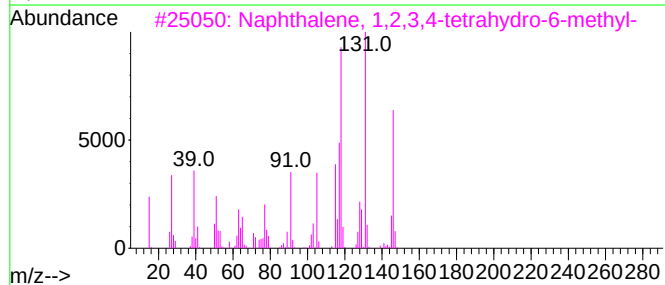
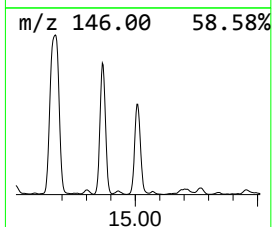
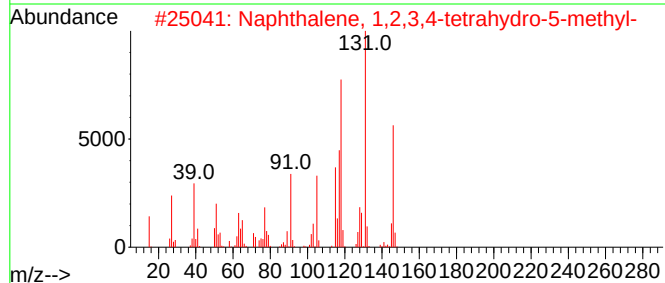
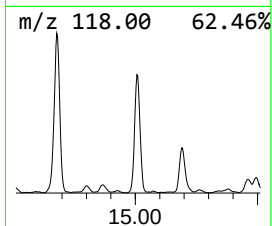
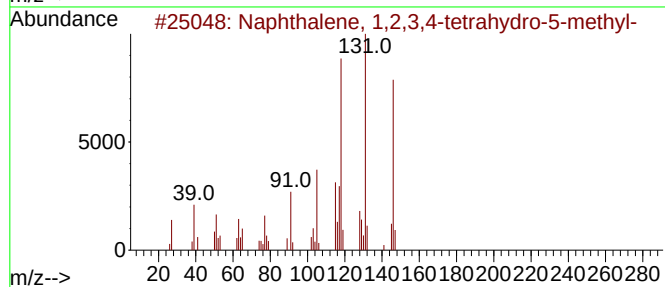
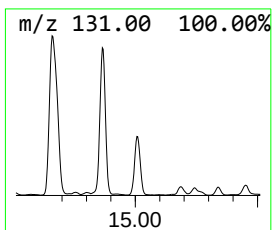
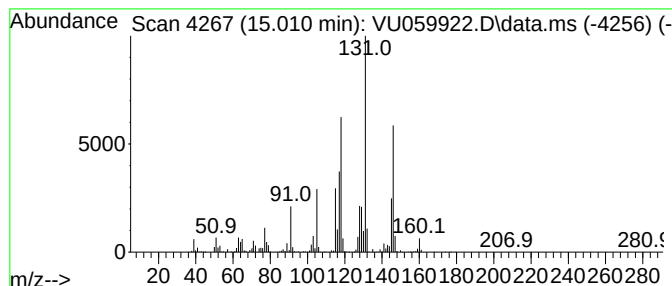
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 25 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.010	2.49 ug/L	960926	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	90
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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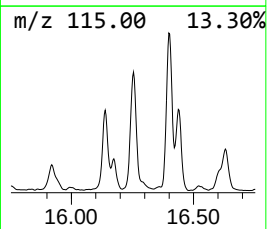
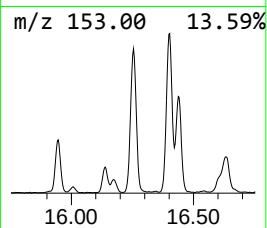
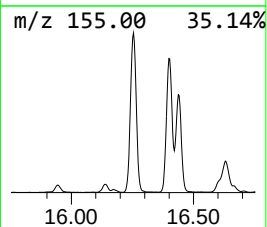
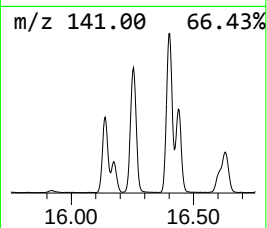
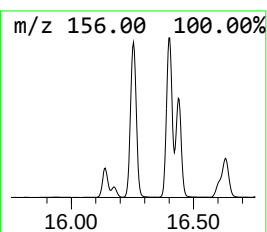
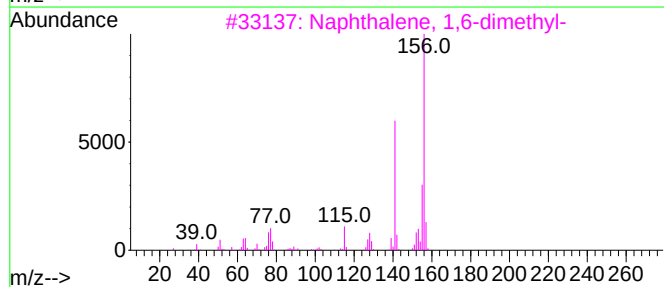
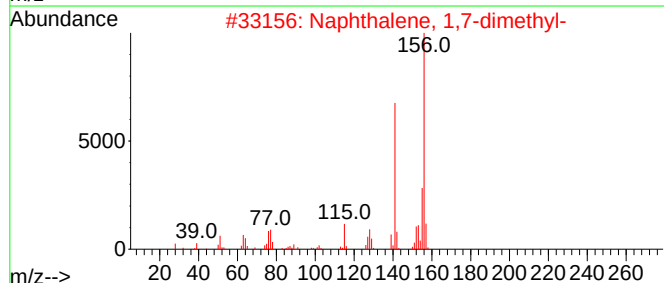
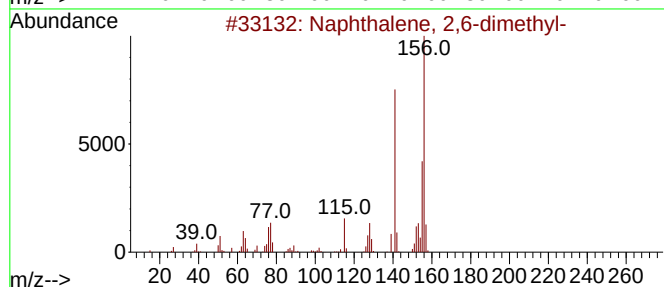
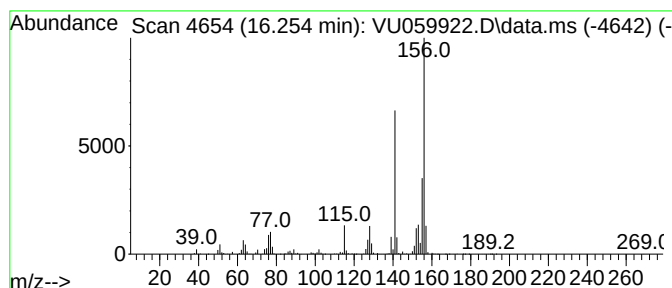
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 28 Naphthalene, 2,6-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.254	6.03 ug/L	2325710	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC7

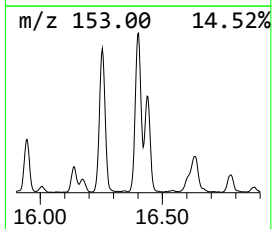
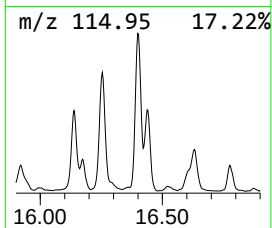
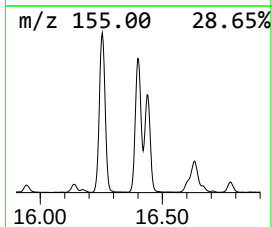
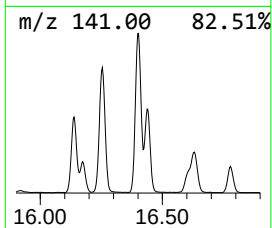
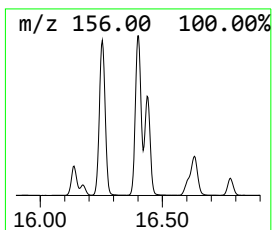
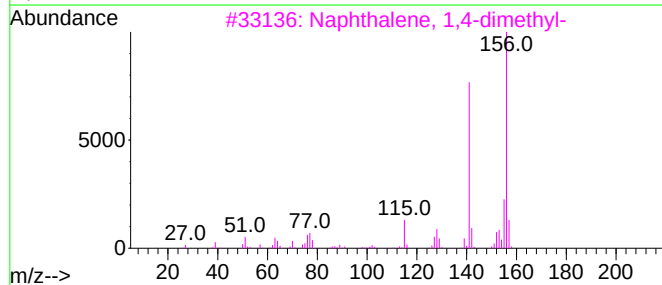
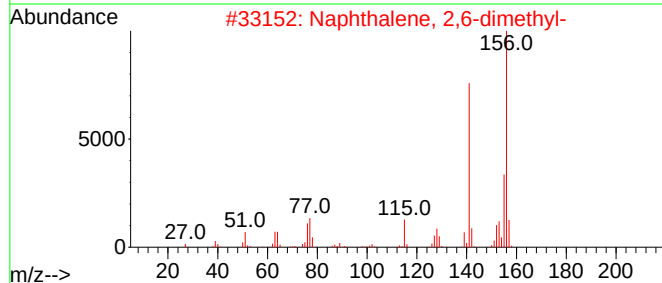
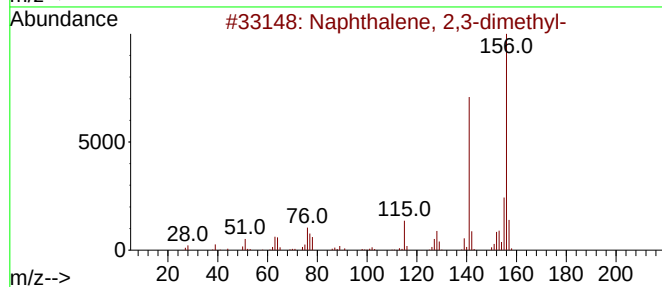
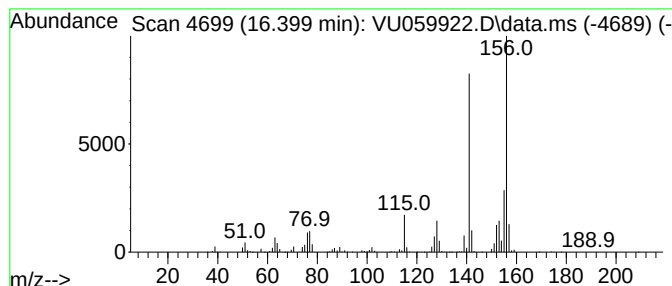
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 29 Naphthalene, 2,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.399	5.85 ug/L	2255350	1,4-Dichlorobenzene-d4	11.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
Data File : VU059922.D
Acq On : 12 Jul 2024 15:25
Operator : MD/SY
Sample : P3217-04
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

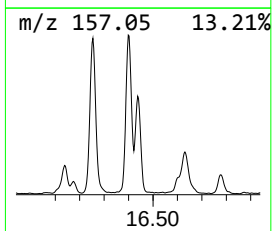
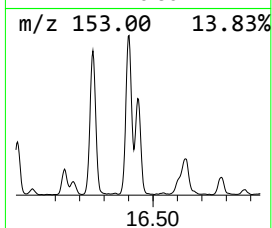
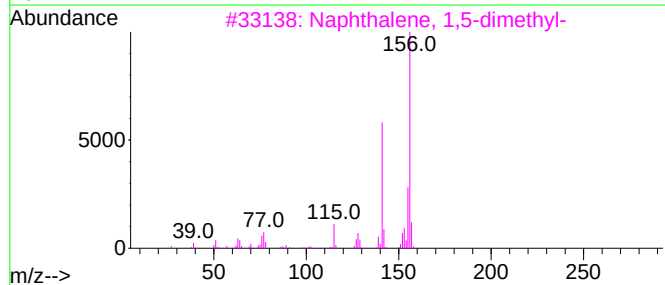
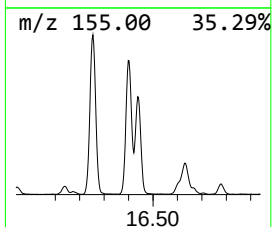
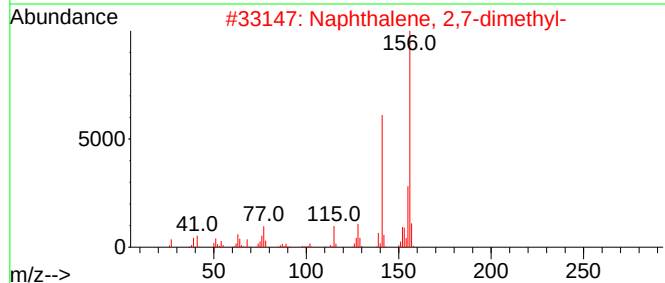
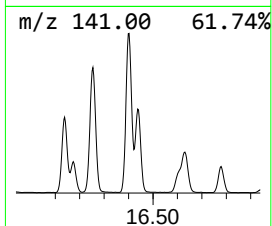
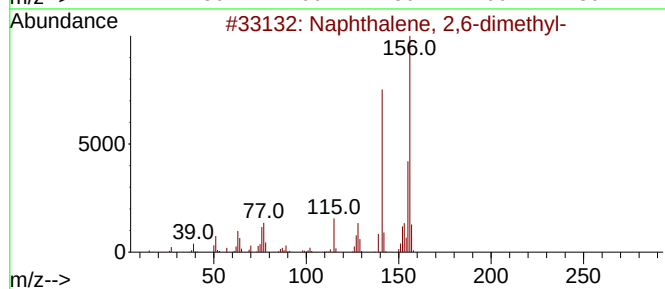
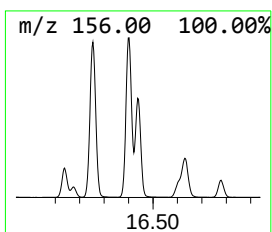
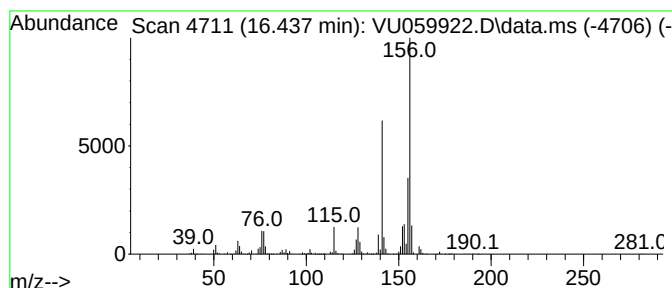
Instrument :
MSVOA_U
ClientSampleId :
YDZC7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 30 Naphthalene, 1,5-dimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.438	3.50 ug/L	1351270	1,4-Dichlorobenzene-d4	11.804	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071224\
 Data File : VU059922.D
 Acq On : 12 Jul 2024 15:25
 Operator : MD/SY
 Sample : P3217-04
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 YDZC7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
 Quant Title : TRACE VOA SFAM1.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, propyl-	10.898	5.3	ug/L	2057820	3	11.804	1928050	5.0
Benzene, 1-ethy...	10.991	29.6	ug/L	11411100	3	11.804	1928050	5.0
Cyclohexane, 1-...	11.155	12.0	ug/L	4640480	3	11.804	1928050	5.0
Benzene, 1-ethy...	11.283	14.0	ug/L	5385080	3	11.804	1928050	5.0
7-Oxabicyclo[2....	11.631	15.2	ug/L	5842140	3	11.804	1928050	5.0
Indane	12.087	13.9	ug/L	5379600	3	11.804	1928050	5.0
o-Cymene	12.460	4.0	ug/L	1533530	3	11.804	1928050	5.0
Benzene, 1-meth...	12.486	4.1	ug/L	1572110	3	11.804	1928050	5.0
Benzene, 1-ethy...	12.563	4.1	ug/L	1574810	3	11.804	1928050	5.0
Benzene, 1-ethe...	12.676	7.0	ug/L	2682340	3	11.804	1928050	5.0
Benzene, 4-ethy...	12.868	2.9	ug/L	1117720	3	11.804	1928050	5.0
Fenchone	12.939	25.0	ug/L	9637040	3	11.804	1928050	5.0
Benzene, 1,2,4,...	13.020	5.3	ug/L	2032510	3	11.804	1928050	5.0
1H-Indene, 2,3-...	13.287	3.1	ug/L	1207540	3	11.804	1928050	5.0
unknown-01	13.399	10.6	ug/L	4104330	3	11.804	1928050	5.0
1-Phenyl-1-butene	13.444	15.4	ug/L	5922050	3	11.804	1928050	5.0
Cyclohexanemeth...	13.512	73.1	ug/L	28204900	3	11.804	1928050	5.0
Naphthalene, 1,...	13.618	9.6	ug/L	3695050	3	11.804	1928050	5.0
(+)-2-Bornanone	13.730	30.5	ug/L	11762700	3	11.804	1928050	5.0
Azulene	14.078	29.5	ug/L	11382100	3	11.804	1928050	5.0
1H-Indene, 2,3-...	14.479	3.2	ug/L	1223730	3	11.804	1928050	5.0
Benzene, (1,2-d...	14.669	6.1	ug/L	2366140	3	11.804	1928050	5.0
1H-Indene, 2,3-...	14.868	3.2	ug/L	1246300	3	11.804	1928050	5.0
Naphthalene, 1,...	15.010	2.5	ug/L	960926	3	11.804	1928050	5.0
Naphthalene, 2,...	16.254	6.0	ug/L	2325710	3	11.804	1928050	5.0
Naphthalene, 2,...	16.399	5.8	ug/L	2255350	3	11.804	1928050	5.0
Naphthalene, 1,...	16.438	3.5	ug/L	1351270	3	11.804	1928050	5.0