

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
 Data File : VU060599.D
 Acq On : 04 Sep 2024 15:03
 Operator : MD/SY
 Sample : P3809-01 50X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 YE3E1

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\SFAMUTR082324WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU060599.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	87	96	111	rVB2	80538	99849	3.18%	0.838%
2	1.911	184	193	206	rBV	48098	63960	2.04%	0.537%
3	2.563	383	396	412	rBV	84630	139711	4.45%	1.173%
4	4.657	1029	1047	1087	rVV	1326237	3138860	100.00%	26.343%
5	5.055	1157	1171	1196	rBV2	82785	186209	5.93%	1.563%
6	5.721	1356	1378	1402	rBV2	164272	439414	14.00%	3.688%
7	6.242	1528	1540	1567	rBV	159706	316034	10.07%	2.652%
8	6.685	1660	1678	1695	rBV2	98728	198621	6.33%	1.667%
9	7.560	1940	1950	1971	rBV	37851	68684	2.19%	0.576%
10	7.891	2040	2053	2064	rBV	173062	308742	9.84%	2.591%
11	7.962	2064	2075	2109	rVB	1191326	2077517	66.19%	17.436%
12	8.177	2131	2142	2153	rBV	24581	44218	1.41%	0.371%
13	8.631	2273	2283	2322	rBV	186855	387814	12.36%	3.255%
14	9.412	2514	2526	2552	rBV	239613	421397	13.43%	3.537%
15	9.566	2562	2574	2587	rBV	44300	75947	2.42%	0.637%
16	9.685	2595	2611	2641	rBV	227038	413085	13.16%	3.467%
17	10.094	2725	2738	2758	rBV	140844	242968	7.74%	2.039%
18	10.750	2931	2942	2954	rBV	77500	127538	4.06%	1.070%
19	10.891	2973	2986	2999	rBV3	23173	49123	1.56%	0.412%
20	10.991	3006	3017	3036	rBV	83333	192359	6.13%	1.614%
21	11.084	3036	3046	3060	rVB	65717	111436	3.55%	0.935%
22	11.287	3098	3109	3126	rVB2	58520	97758	3.11%	0.820%
23	11.463	3153	3164	3184	rVB	188518	302139	9.63%	2.536%
24	11.631	3198	3216	3228	rBV	29396	80336	2.56%	0.674%
25	11.804	3255	3270	3286	rVV2	272376	533498	17.00%	4.477%
26	11.888	3286	3296	3312	rVB	165935	282997	9.02%	2.375%
27	12.090	3344	3359	3366	rBV3	49278	88876	2.83%	0.746%
28	12.187	3377	3389	3410	rVB	224578	422188	13.45%	3.543%
29	12.566	3497	3507	3514	rBV2	25420	37354	1.19%	0.313%
30	12.676	3529	3541	3560	rBV4	30966	71270	2.27%	0.598%
31	12.939	3614	3623	3639	rVV4	31038	82575	2.63%	0.693%
32	13.020	3640	3648	3664	rVB	34495	57113	1.82%	0.479%
33	13.447	3771	3781	3791	rVV3	76203	138432	4.41%	1.162%
34	13.508	3791	3800	3811	rVB3	43311	73886	2.35%	0.620%
35	13.618	3824	3834	3848	rVB2	42203	71500	2.28%	0.600%
36	13.727	3854	3868	3878	rBV4	48134	83425	2.66%	0.700%

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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
YE3E1

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\SFAMUTR082324WMA.M
Title : TRACE VOA SFAM1.0

37	14.084	3963	3979	3995	rBV	56988	107725	3.43%	0.904%
38	14.672	4148	4162	4174	rVB6	15795	39334	1.25%	0.330%
39	15.216	4316	4331	4354	rBV2	69628	129888	4.14%	1.090%
40	15.405	4378	4390	4403	rBV	65709	111543	3.55%	0.936%

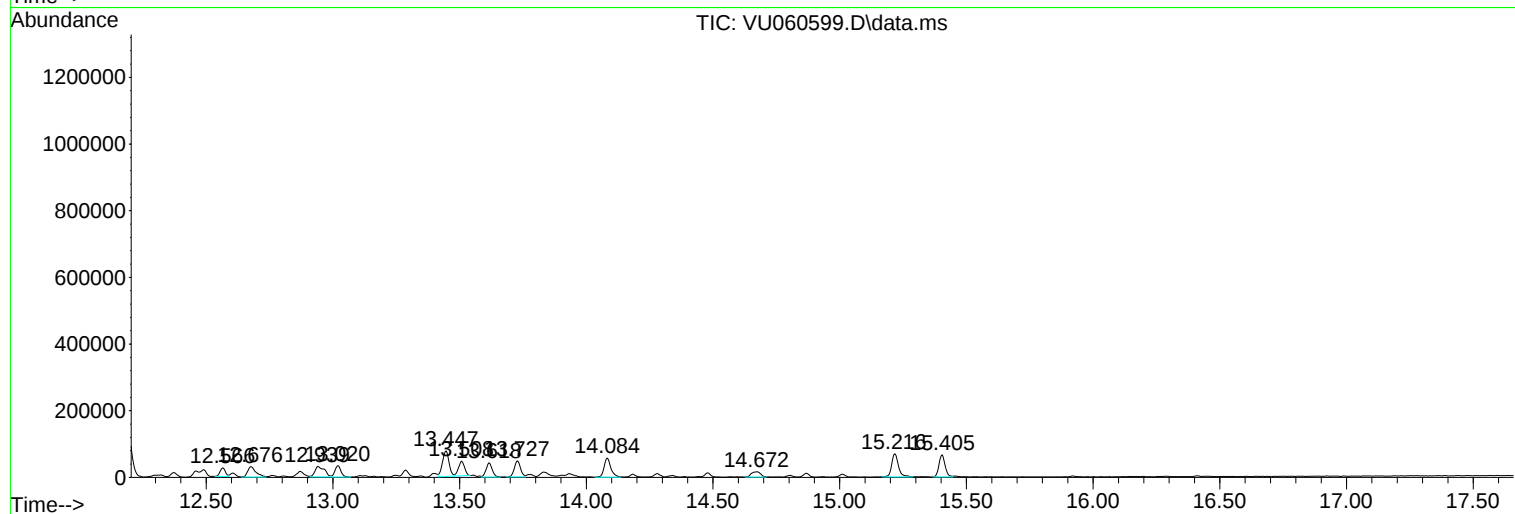
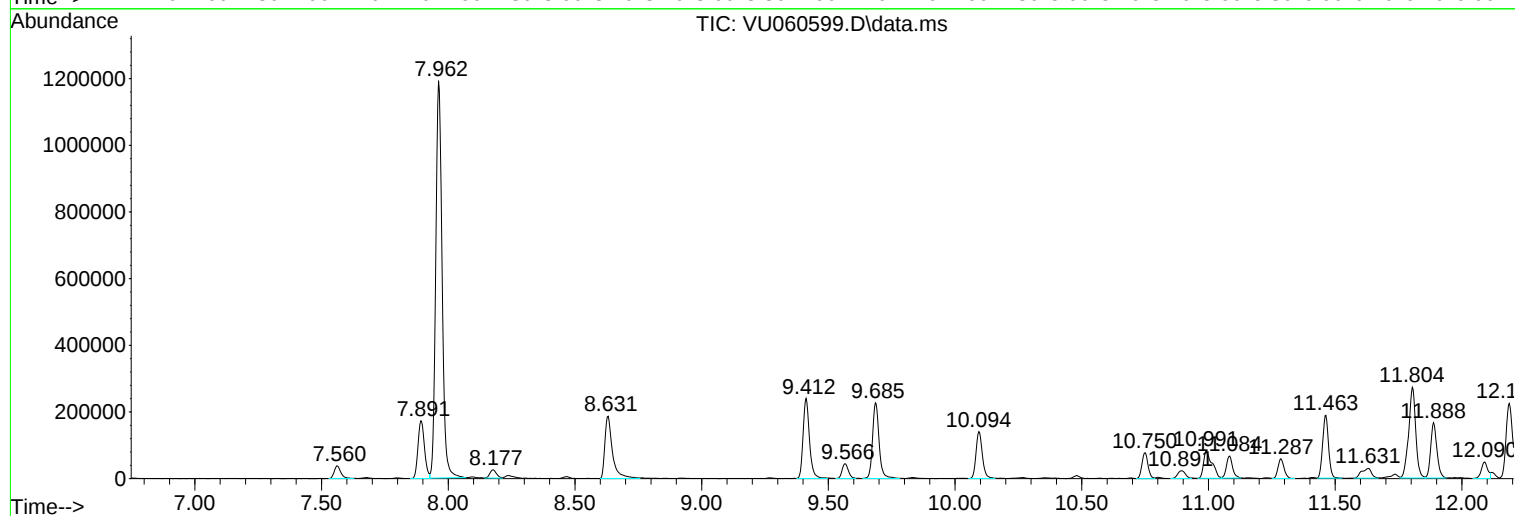
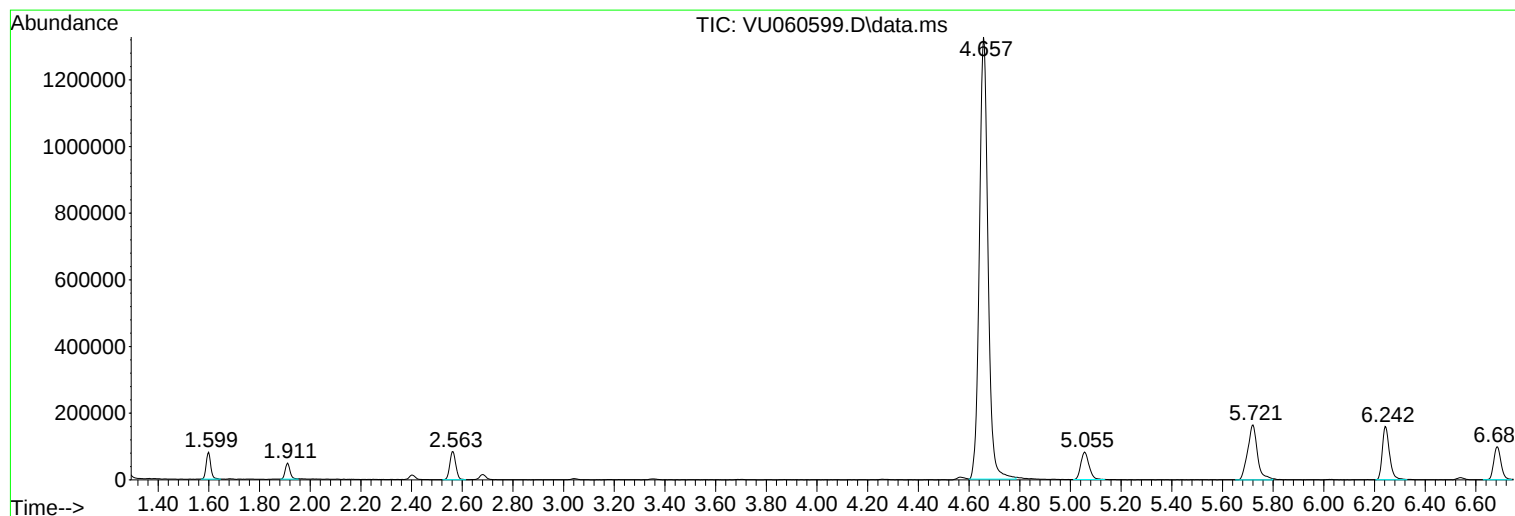
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR082324WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Misc : 25.0mL/MSVOA_U/WATER
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Instrument :
MSVOA_U
ClientSampleId :
YE3E1

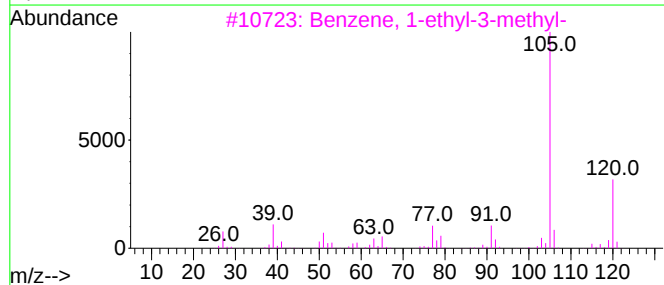
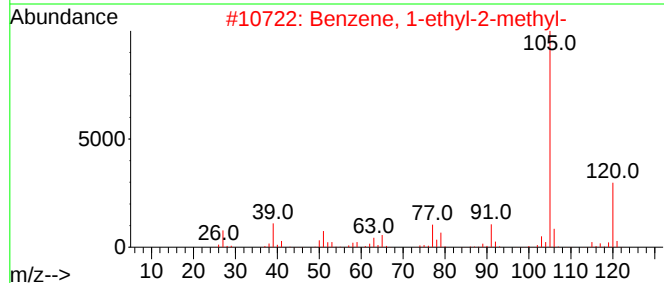
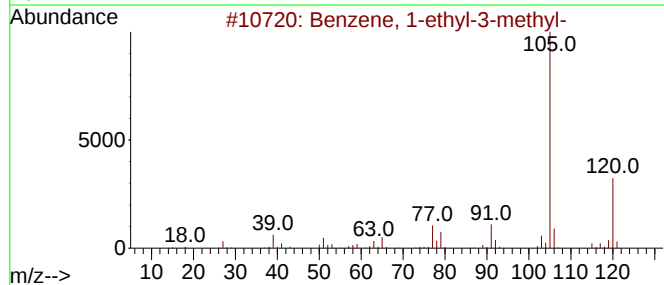
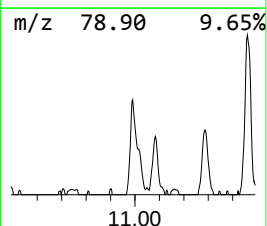
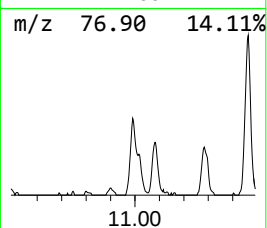
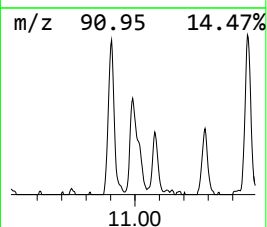
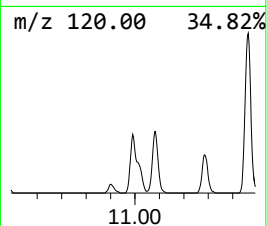
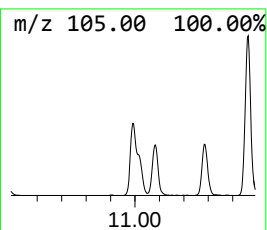
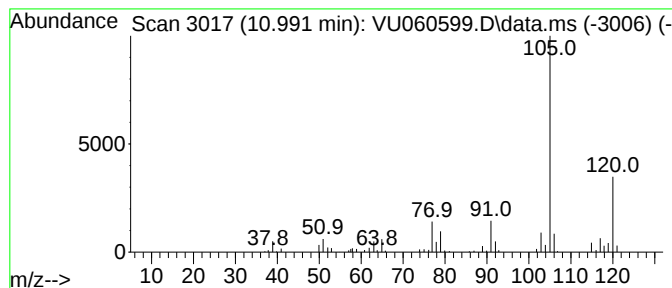
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.991	1.80 ug/L	192359	1,4-Dichlorobenzene-d4	11.807

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	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



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Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

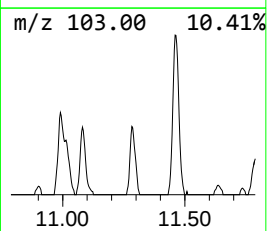
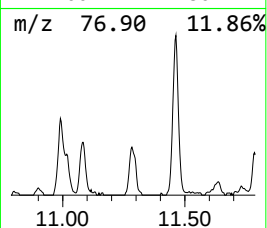
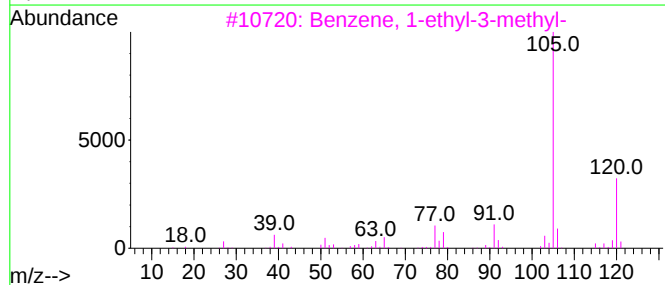
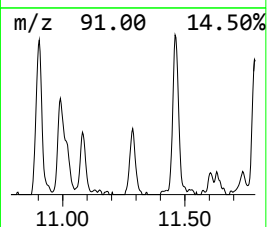
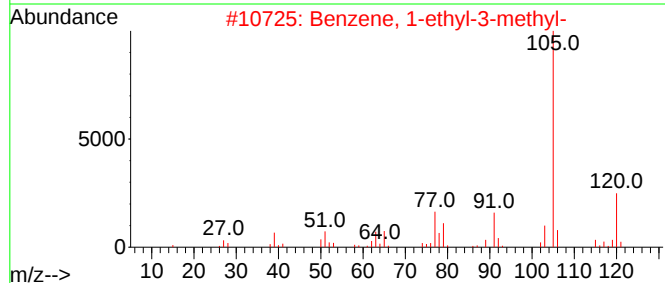
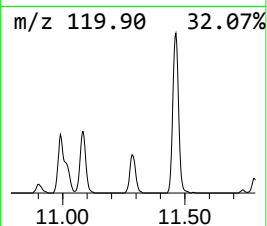
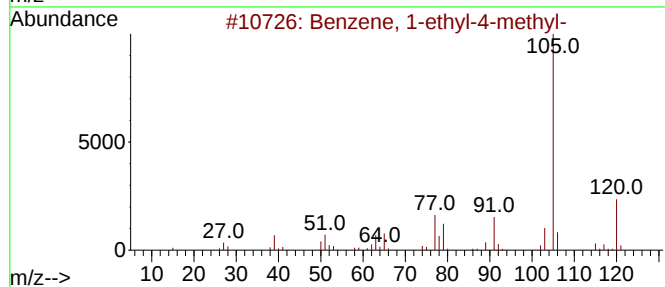
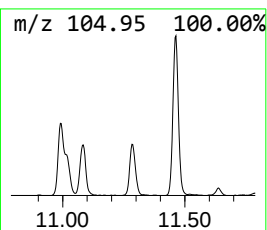
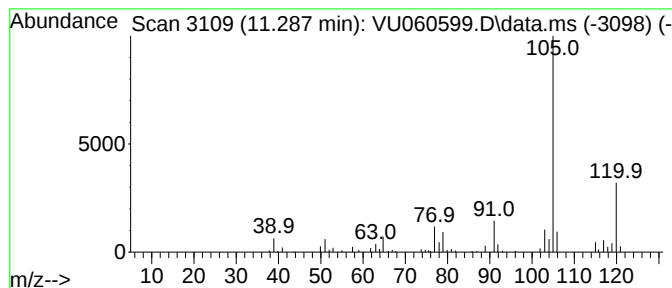
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR082324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.287	0.92 ug/L	97758	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

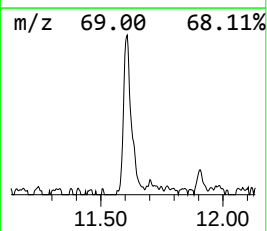
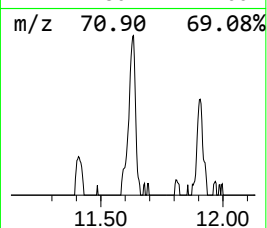
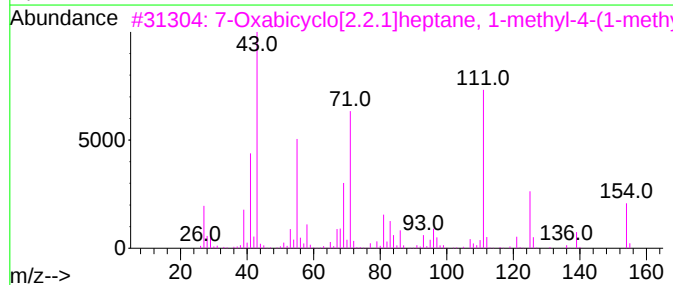
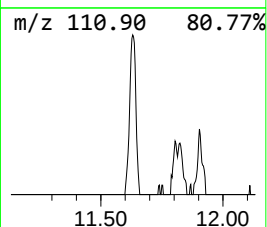
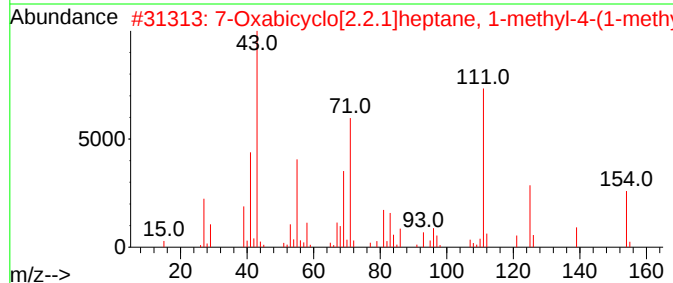
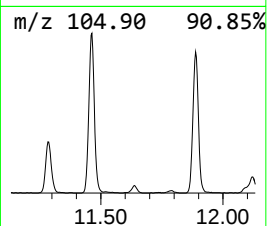
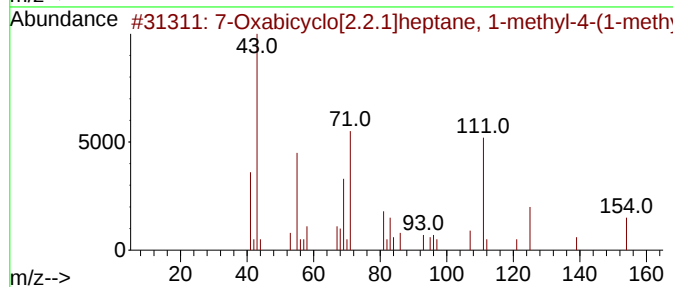
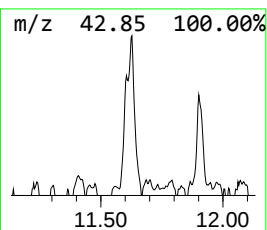
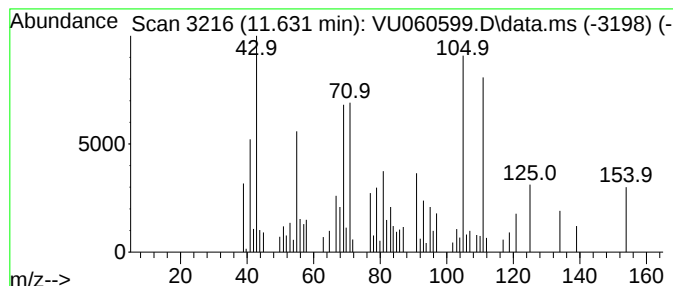
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR082324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 7-Oxabicyclo[2.2.1]heptane,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.631	0.75 ug/L	80336	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	90
2			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	76
3			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	76
4			Butanoic acid, tridec-2-ynyl ester	266	C17H30O2	1010299-12-6	58
5			1-Propanone, 2-methyl-1-[2-(1-me...	154	C10H18O	056259-15-5	38



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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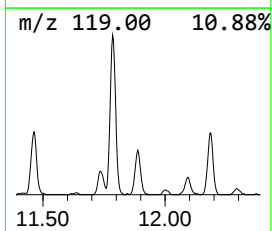
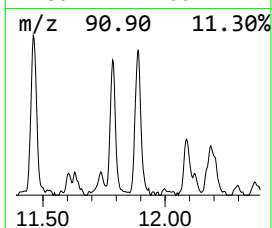
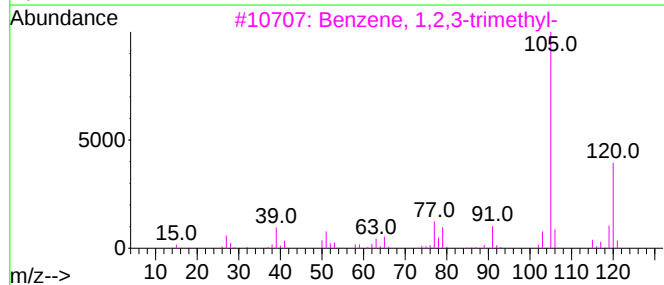
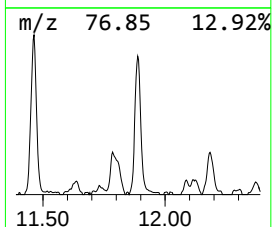
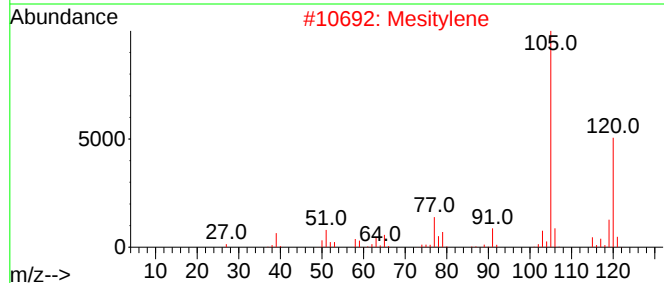
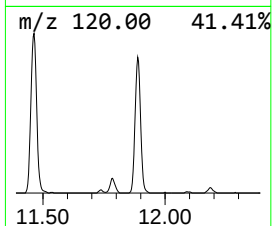
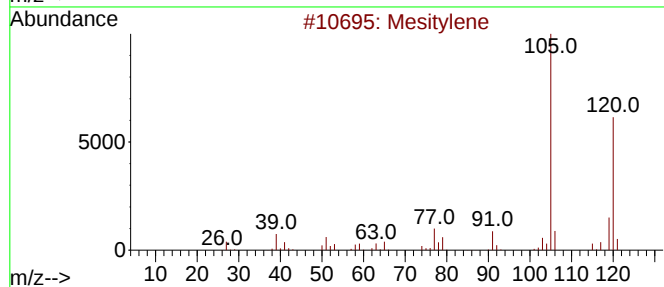
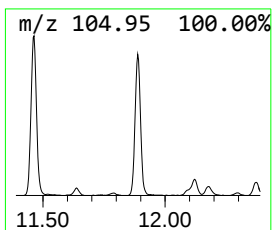
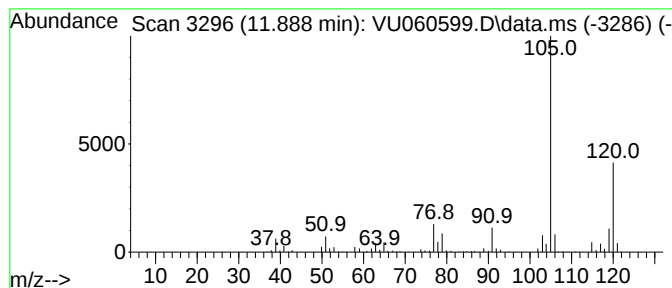
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR082324WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.888	2.65 ug/L	282997	1,4-Dichlorobenzene-d4	11.807

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



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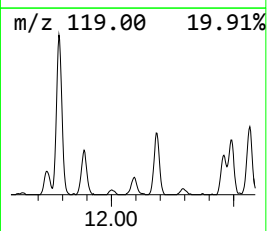
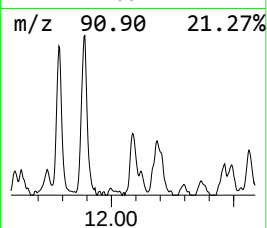
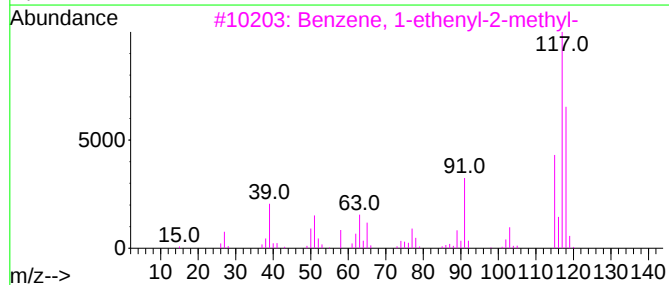
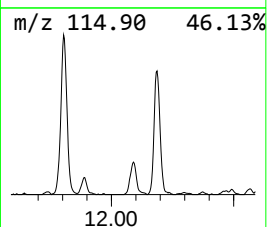
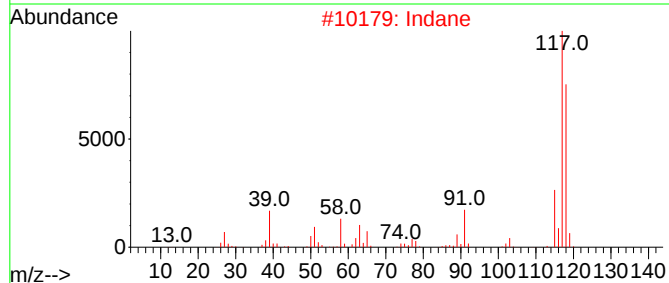
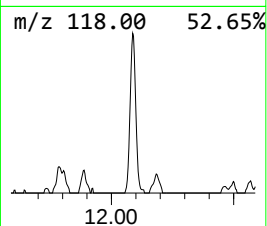
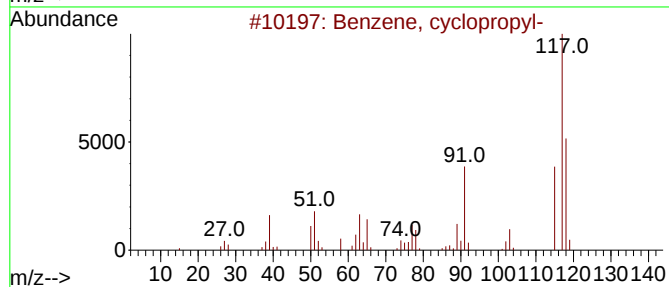
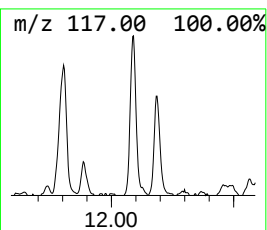
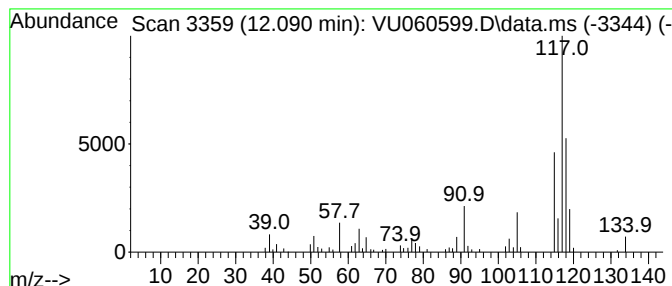
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, cyclopropyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.090	0.83 ug/L	88876	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
2			Indane	118	C9H10	000496-11-7	64
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
4			Indane	118	C9H10	000496-11-7	64
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

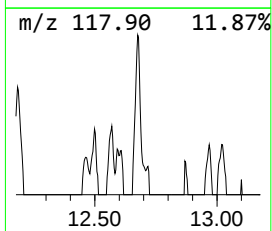
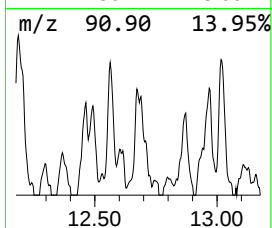
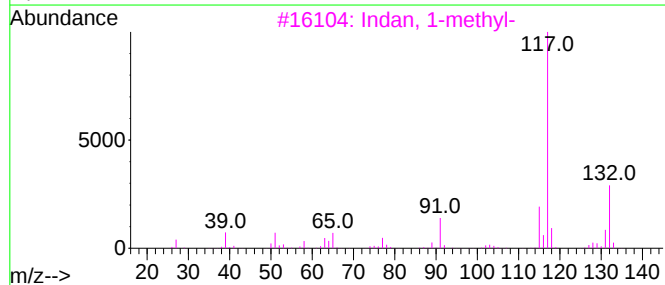
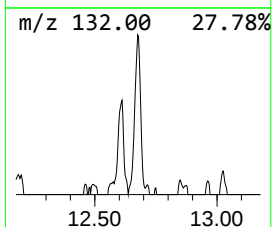
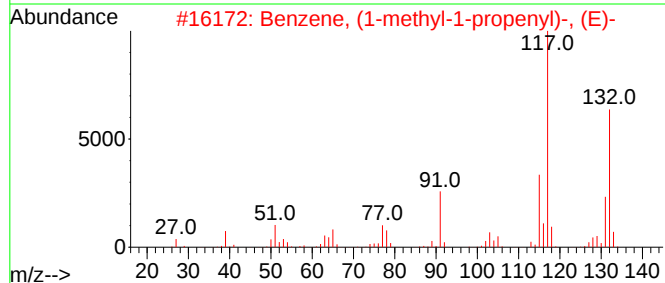
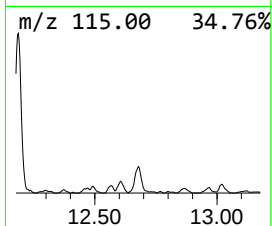
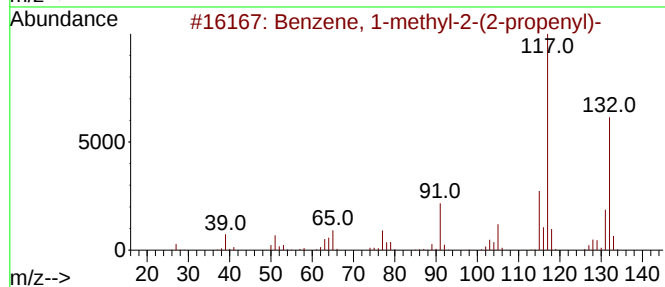
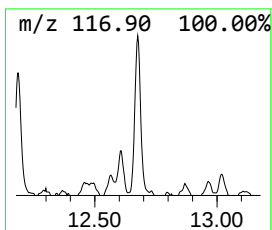
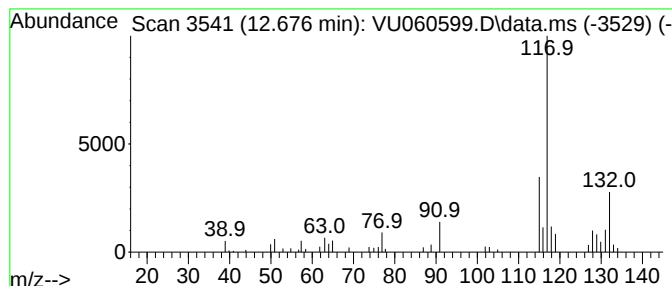
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.676	0.67 ug/L	71270	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

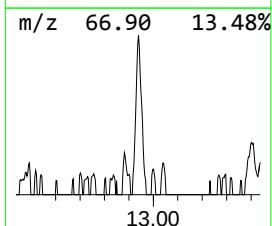
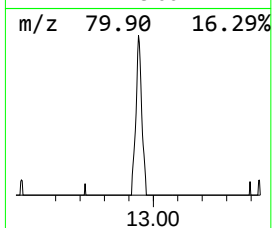
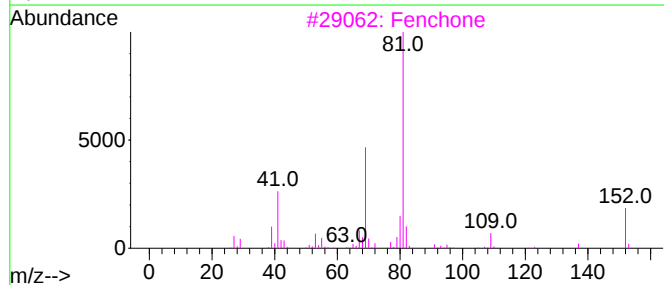
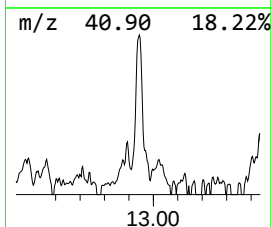
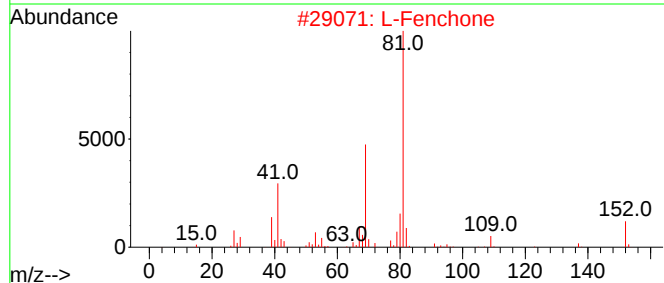
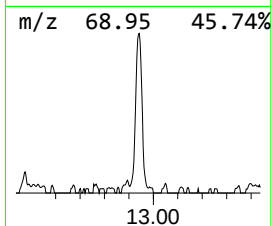
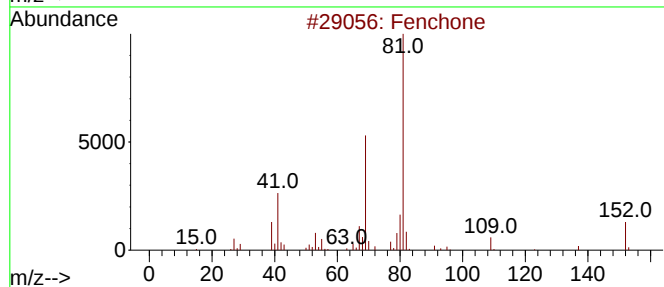
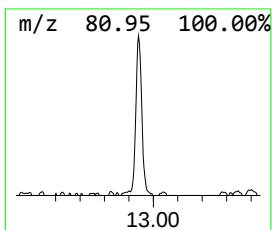
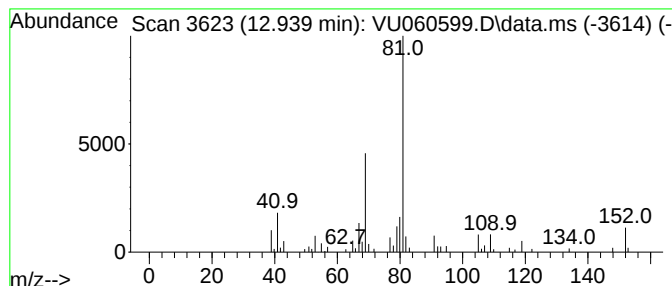
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Fenchone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.939	0.77 ug/L	82575	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchone	152	C10H16O	001195-79-5	94
2			L-Fenchone	152	C10H16O	007787-20-4	91
3			Fenchone	152	C10H16O	001195-79-5	87
4			Fenchone	152	C10H16O	001195-79-5	80
5			D-Fenchone	152	C10H16O	004695-62-9	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

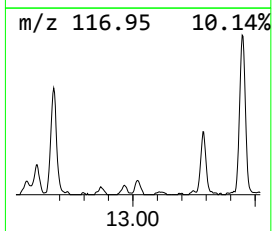
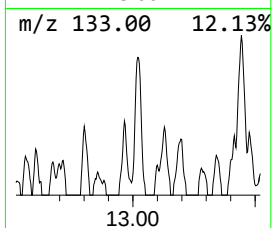
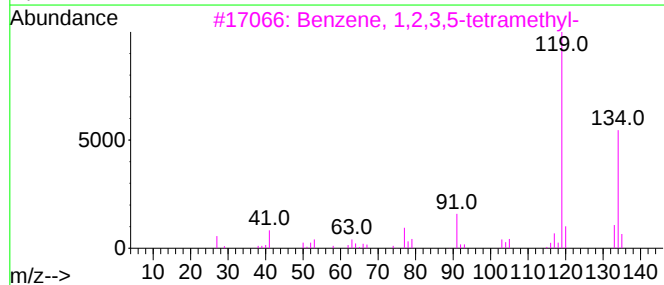
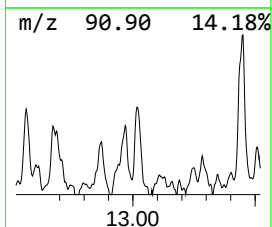
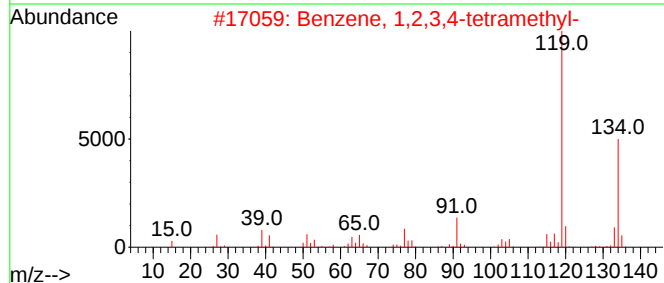
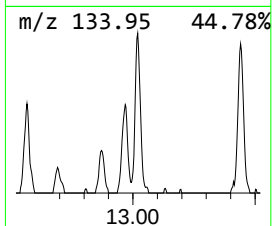
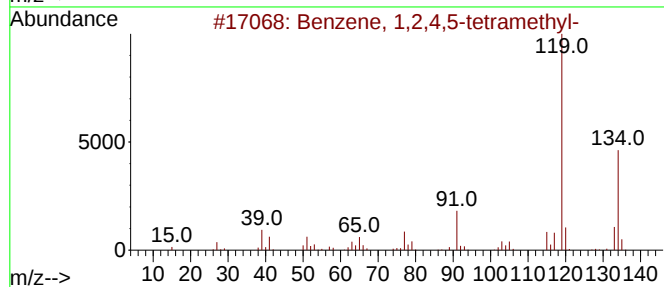
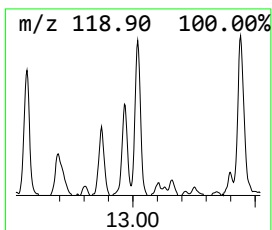
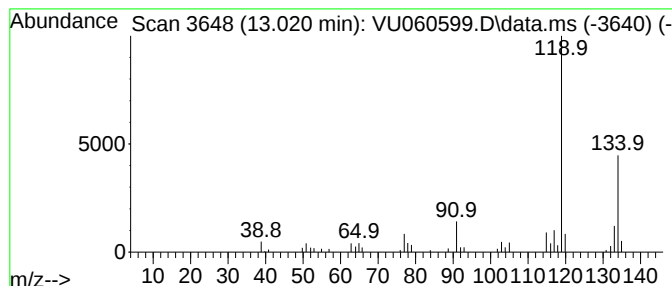
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.020	0.54 ug/L	57113	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

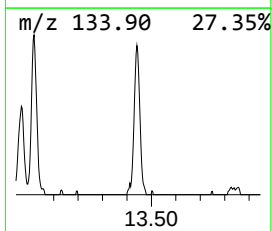
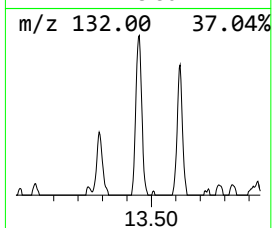
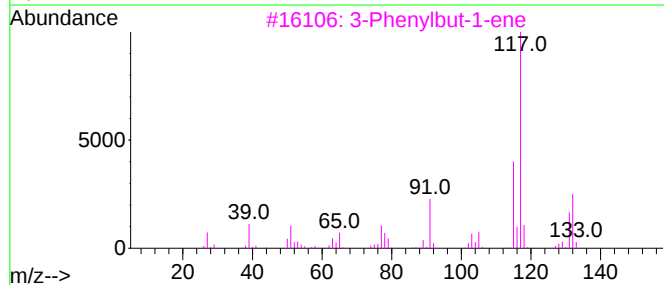
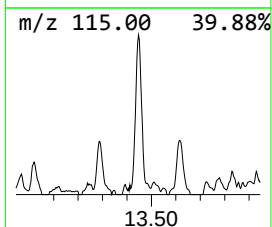
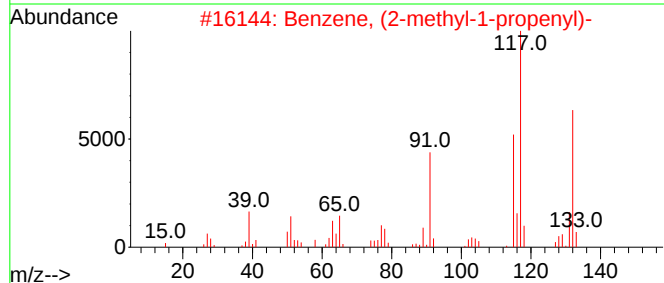
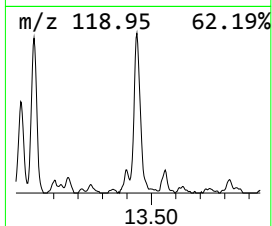
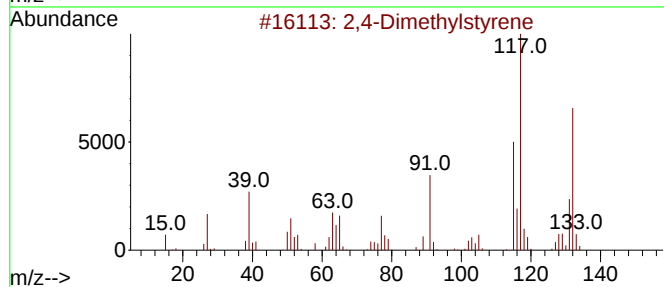
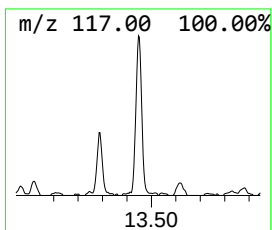
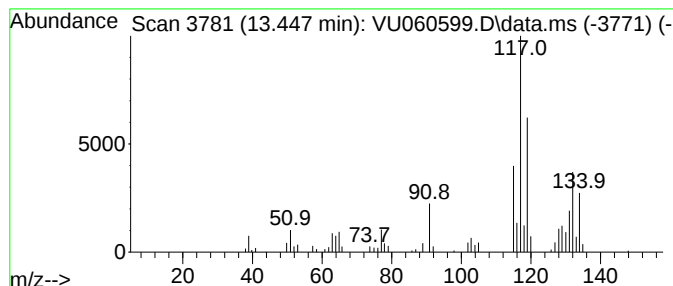
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 2,4-Dimethylstyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	1.30 ug/L	138432	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylstyrene	132	C10H12	002234-20-0	89
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	60
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

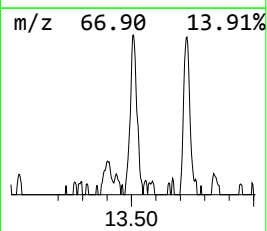
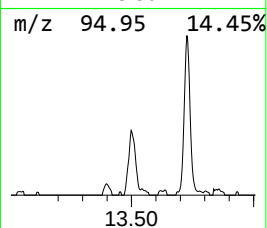
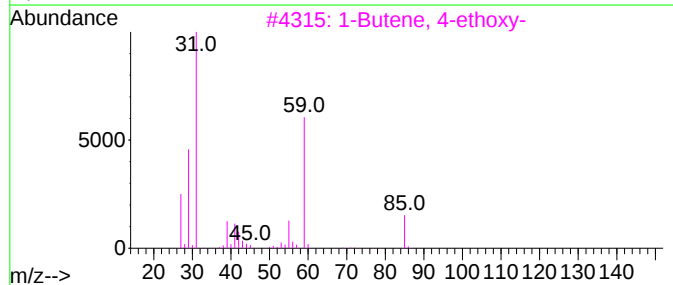
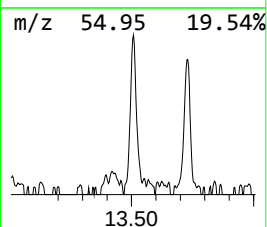
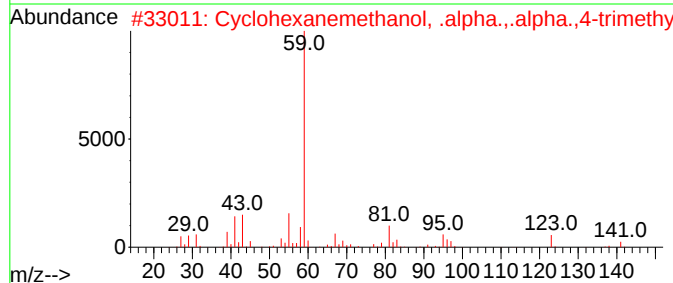
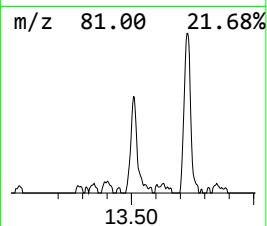
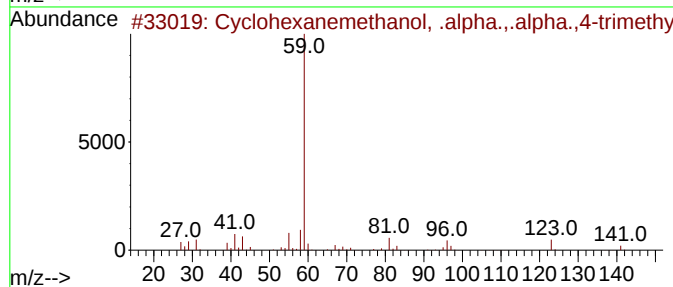
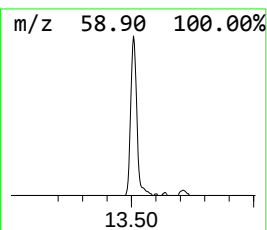
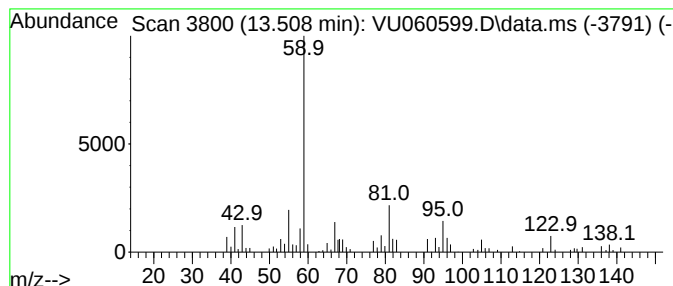
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Cyclohexanemethanol, .alpha... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.508	0.69 ug/L	73886	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	59
2			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	43
3			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	43
4			Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	42
5			2-Hydroxy-2-methylundec-5-en-3-one	198	C12H22O2	1000191-46-2	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

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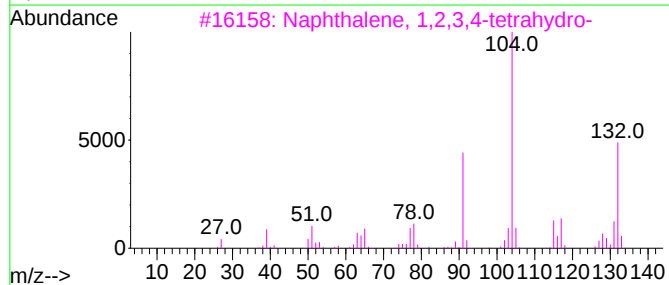
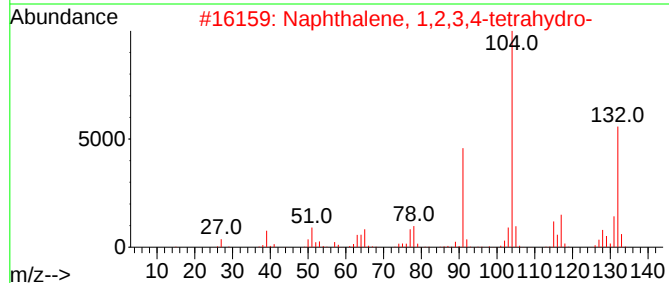
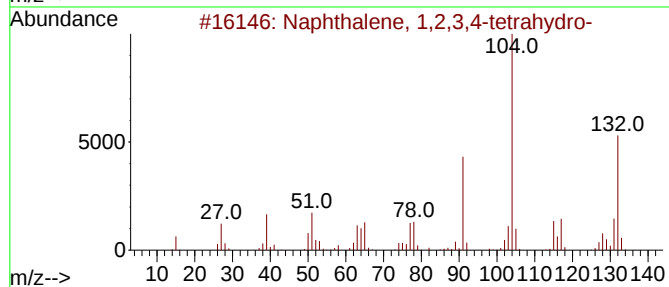
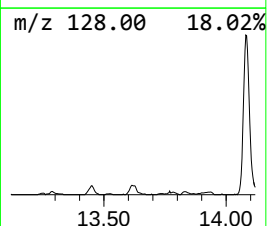
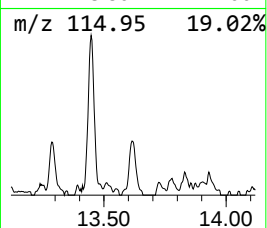
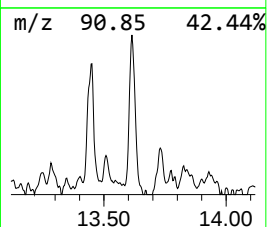
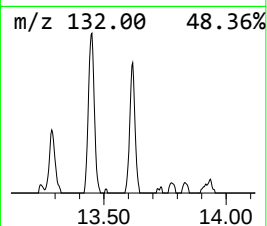
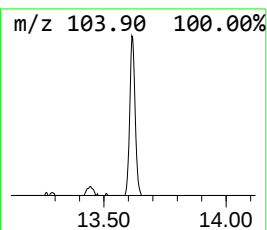
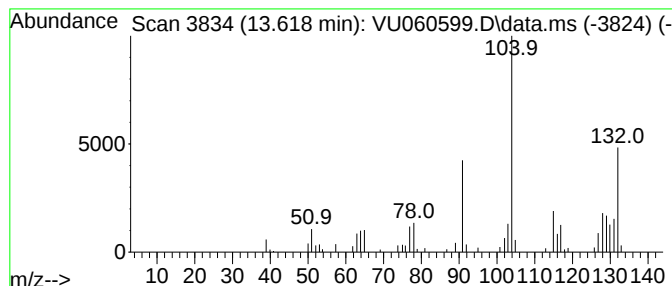
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.618	0.67 ug/L	71500	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	81
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

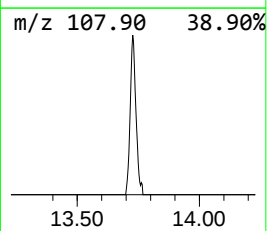
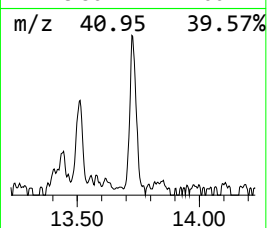
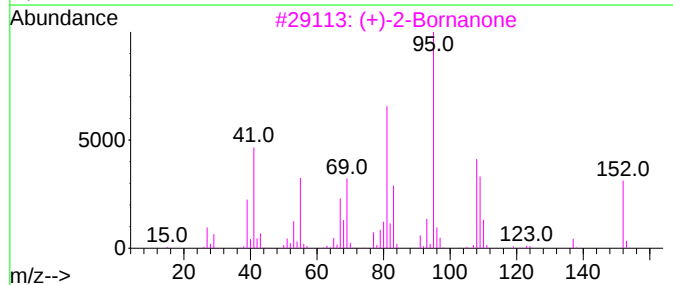
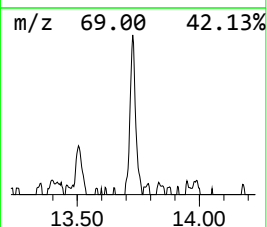
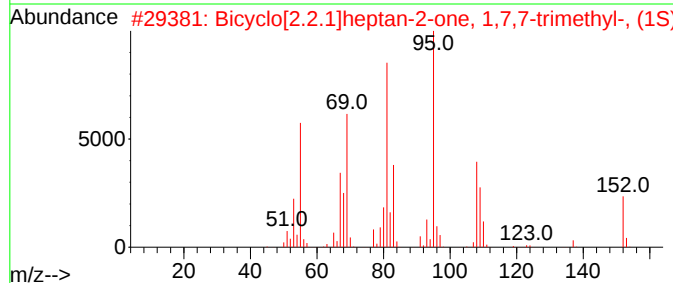
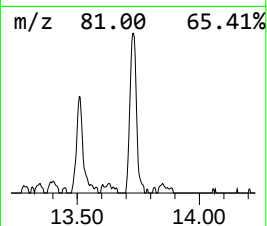
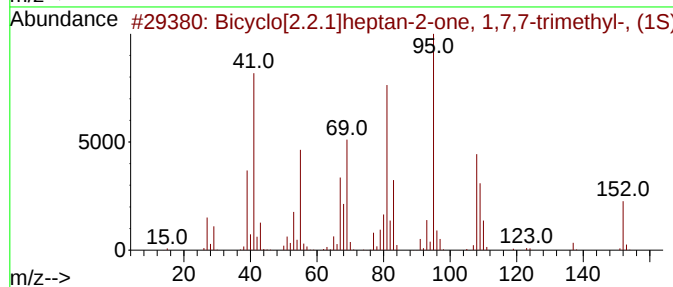
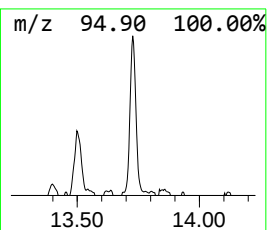
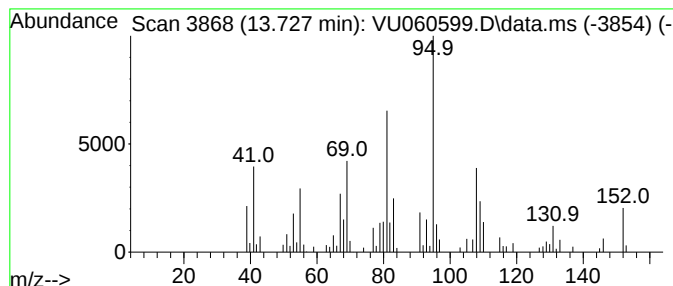
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.727	0.78 ug/L	83425	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
2			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	97
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	97
4			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	97
5			Camphor	152	C10H16O	000076-22-2	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090424\
Data File : VU060599.D
Acq On : 04 Sep 2024 15:03
Operator : MD/SY
Sample : P3809-01 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YE3E1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR082324WMA.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, 1-ethy...	10.991	1.8	ug/L	192359	3	11.807	533498	5.0
Benzene, 1-ethy...	11.287	0.9	ug/L	97758	3	11.807	533498	5.0
7-Oxabicyclo[2...	11.631	0.8	ug/L	80336	3	11.807	533498	5.0
Benzene, 1,2,3-...	11.888	2.6	ug/L	282997	3	11.807	533498	5.0
Benzene, cyclop...	12.090	0.8	ug/L	88876	3	11.807	533498	5.0
Benzene, 1-meth...	12.676	0.7	ug/L	71270	3	11.807	533498	5.0
Fenchone	12.939	0.8	ug/L	82575	3	11.807	533498	5.0
Benzene, 1,2,4,...	13.020	0.5	ug/L	57113	3	11.807	533498	5.0
2,4-Dimethylsty...	13.447	1.3	ug/L	138432	3	11.807	533498	5.0
Cyclohexanemeth...	13.508	0.7	ug/L	73886	3	11.807	533498	5.0
Naphthalene, 1,...	13.618	0.7	ug/L	71500	3	11.807	533498	5.0
Bicyclo[2.2.1]h...	13.727	0.8	ug/L	83425	3	11.807	533498	5.0