

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100922\
 Data File : VU051304.D
 Acq On : 09 Oct 2022 16:51
 Operator : JC/MD
 Sample : N5013-03
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BHA40

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\SFAMUTR100722WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VU051304.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.375	310	322	339	rBV	1133306	1658500	2.87%	1.012%
2	3.989	805	824	852	rVB	1206530	3041075	5.25%	1.855%
3	4.665	1010	1034	1068	rVB2	1219513	3170125	5.48%	1.934%
4	5.057	1142	1156	1189	rVB2	275316	656318	1.13%	0.400%
5	5.774	1344	1379	1408	rBV	18296998	37364043	64.56%	22.790%
6	7.899	2022	2040	2048	rBV	335339	593480	1.03%	0.362%
7	7.970	2048	2062	2113	rVB2	33625487	57872945	100.00%	35.299%
8	8.632	2256	2268	2299	rBV2	505321	1017640	1.76%	0.621%
9	9.446	2499	2521	2546	rBV	4977863	8842195	15.28%	5.393%
10	9.568	2546	2559	2583	rVB	4722078	7495741	12.95%	4.572%
11	9.690	2583	2597	2614	rBV	12632837	20296697	35.07%	12.380%
12	10.099	2709	2724	2753	rVB	4114972	6421274	11.10%	3.917%
13	10.481	2826	2843	2857	rBV	1086378	1690256	2.92%	1.031%
14	10.902	2956	2974	2990	rBV2	441197	892591	1.54%	0.544%
15	10.996	2990	3003	3009	rBV2	577957	1011126	1.75%	0.617%
16	11.086	3022	3031	3046	rVB2	390521	608785	1.05%	0.371%
17	11.288	3084	3094	3117	rVB	517209	831922	1.44%	0.507%
18	11.465	3138	3149	3163	rVB	1716556	2545969	4.40%	1.553%
19	11.809	3243	3256	3272	rBV3	476333	1205381	2.08%	0.735%
20	11.893	3272	3282	3295	rVB	825899	1315951	2.27%	0.803%
21	12.092	3329	3344	3362	rBV2	1358462	2208851	3.82%	1.347%
22	12.192	3363	3375	3398	rVB4	584532	1479944	2.56%	0.903%
23	12.860	3575	3583	3597	rVB2	388919	642936	1.11%	0.392%
24	14.086	3951	3964	3987	rVB	632215	1089046	1.88%	0.664%

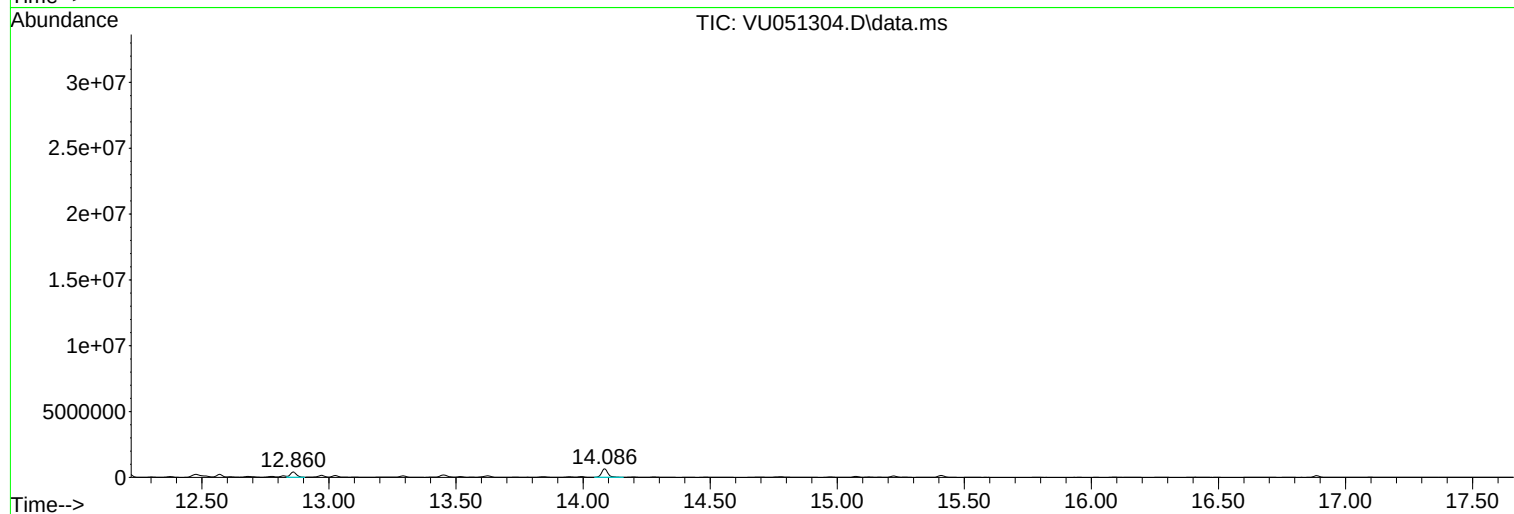
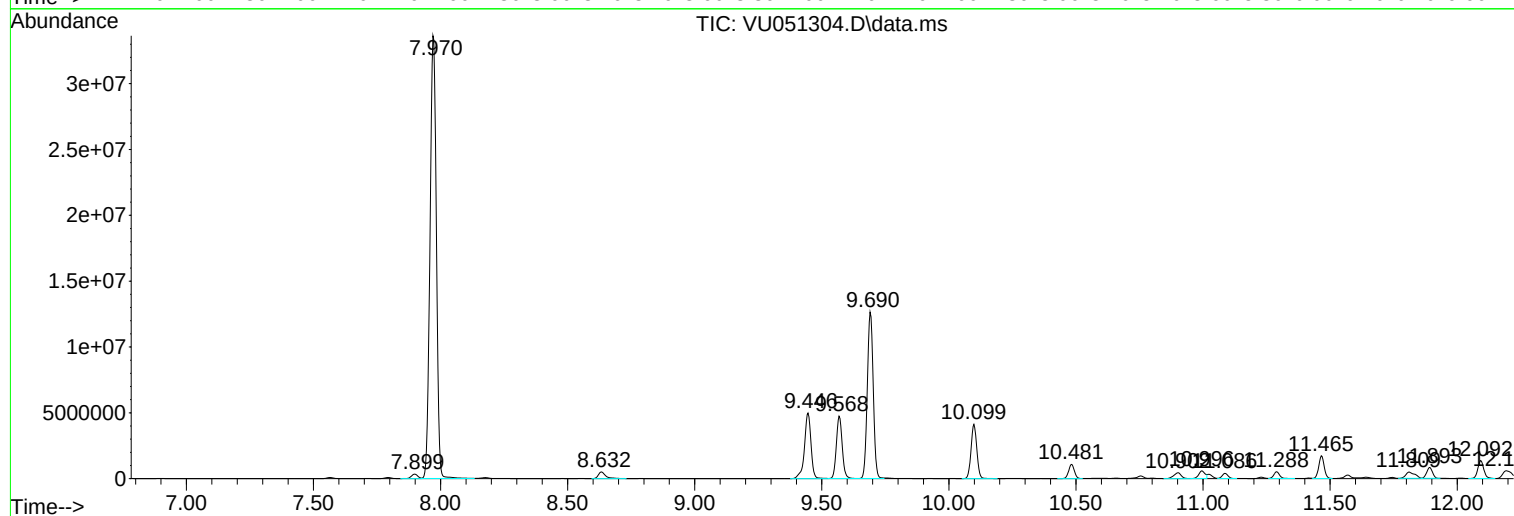
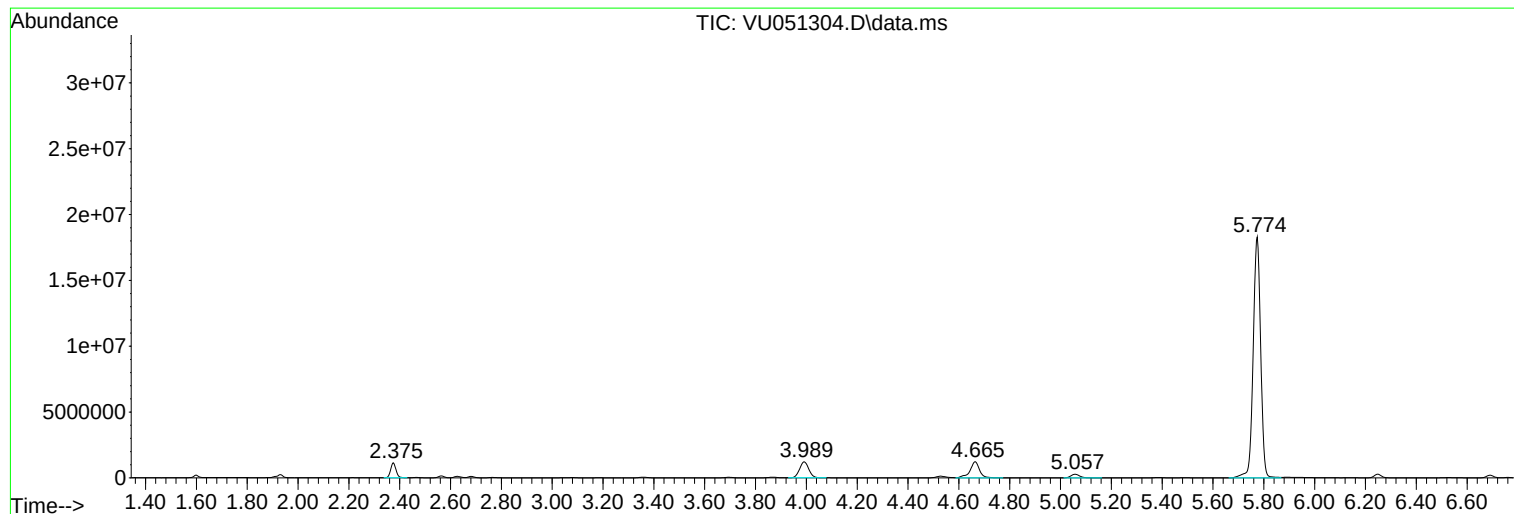
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ClientSampleId :
BHA40

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100722WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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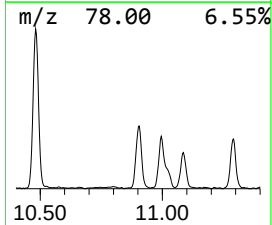
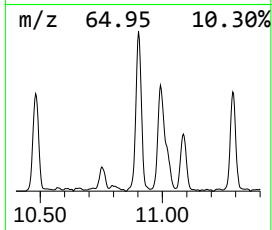
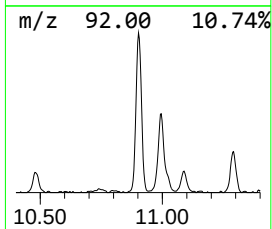
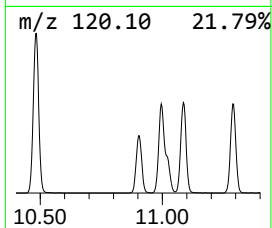
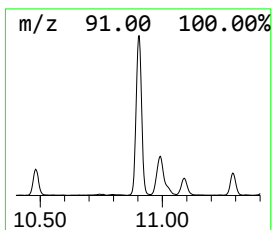
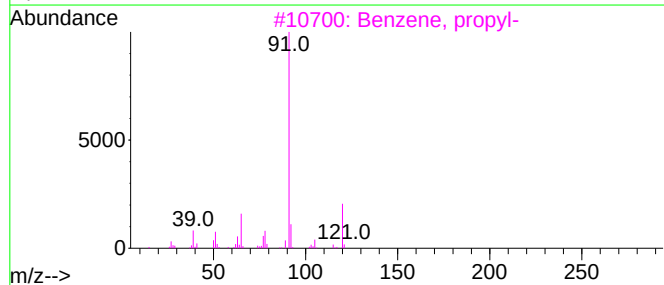
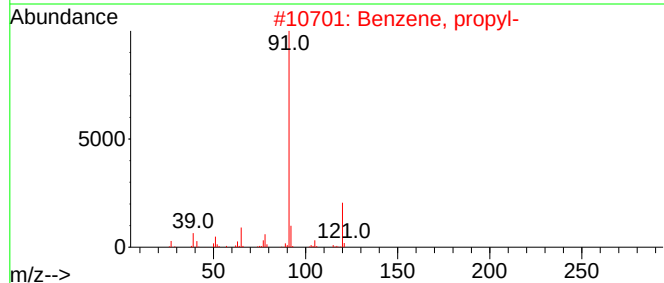
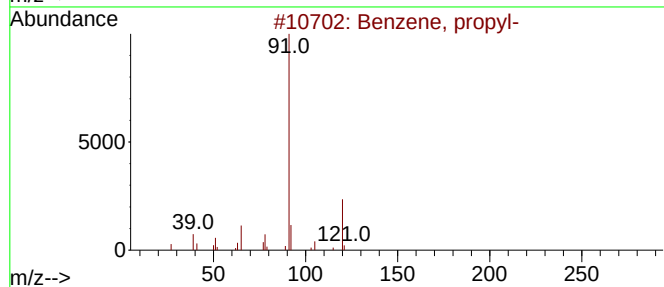
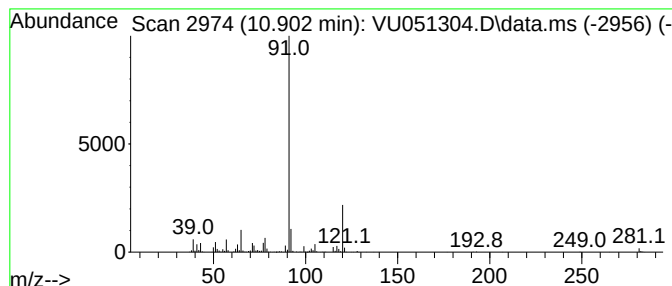
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.902	3.70 ug/L	892591	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	87
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	80
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64



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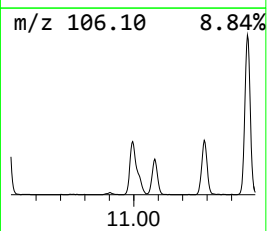
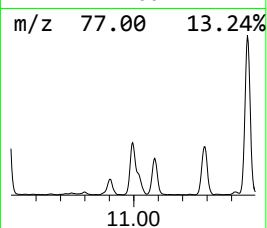
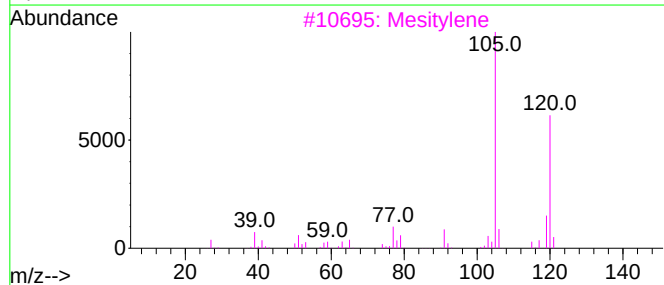
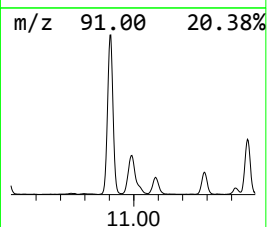
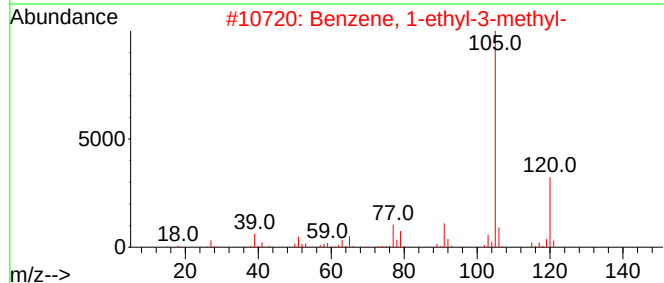
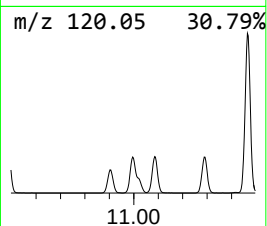
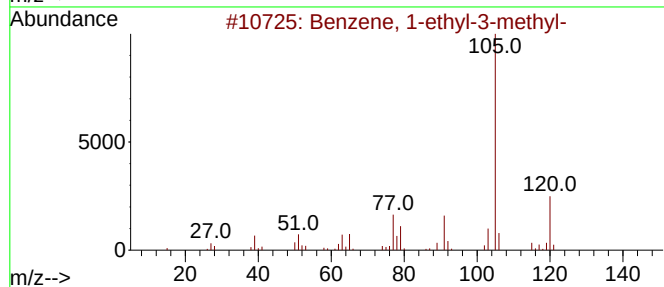
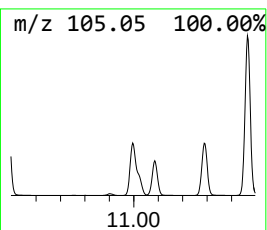
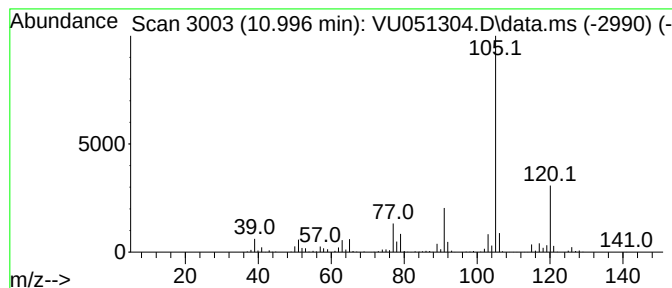
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100722WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.996	4.19 ug/L	1011130	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
3			Mesitylene	120	C9H12	000108-67-8	90
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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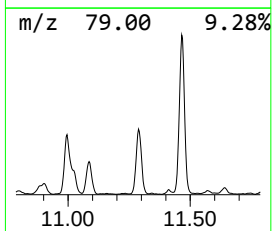
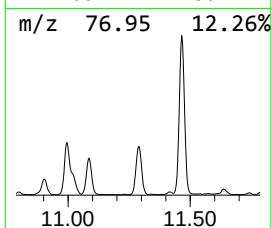
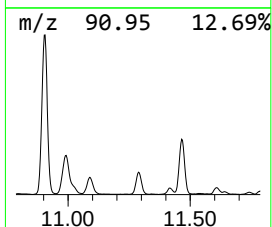
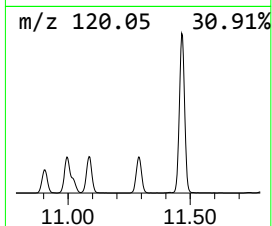
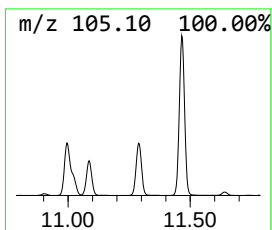
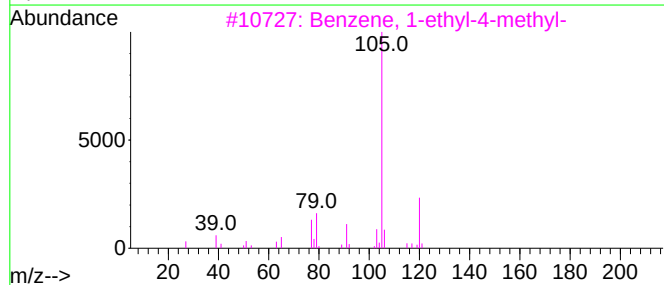
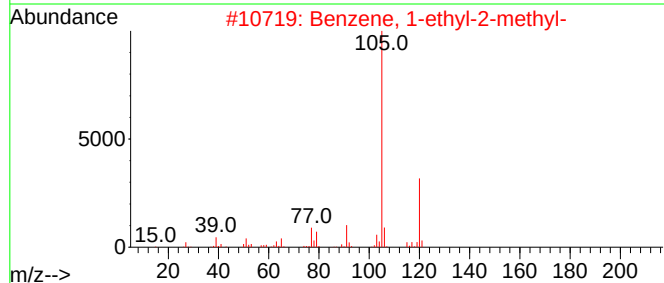
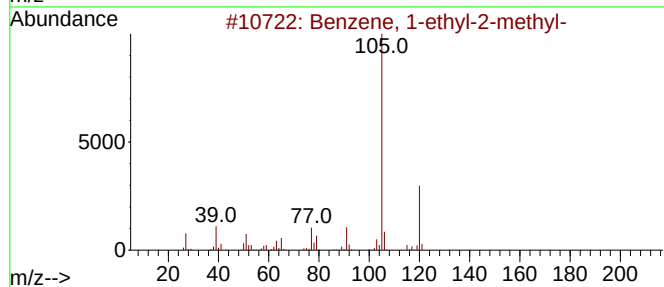
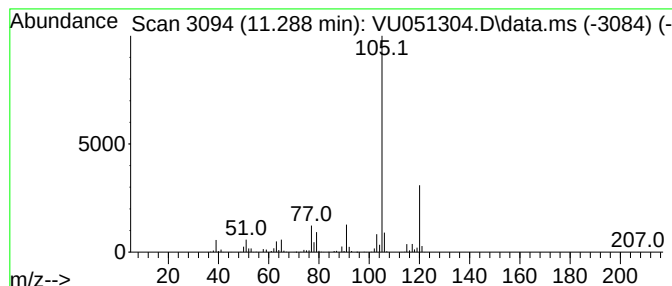
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100722WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.288	3.45 ug/L	831922	1,4-Dichlorobenzene-d4	11.809

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	93
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



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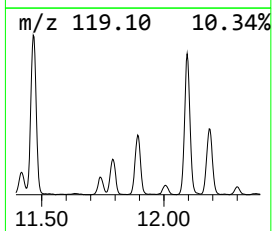
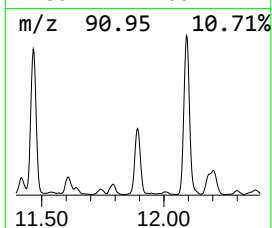
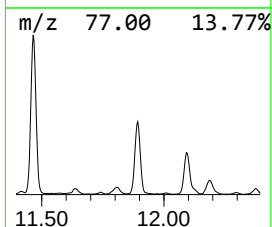
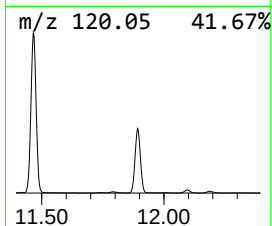
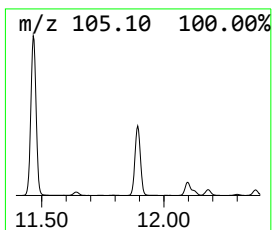
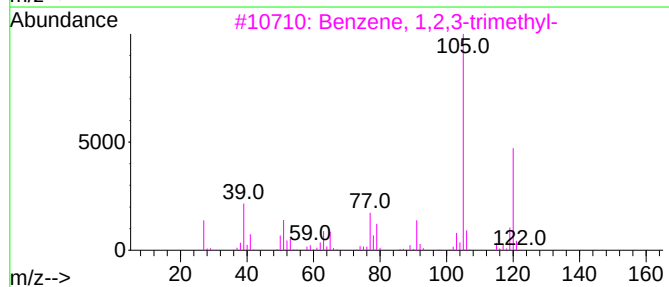
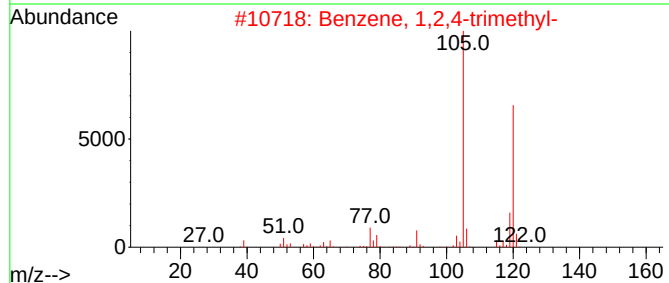
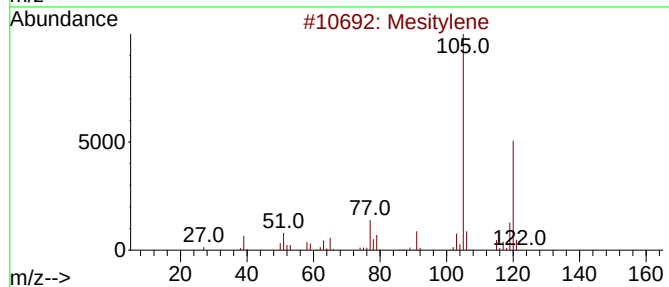
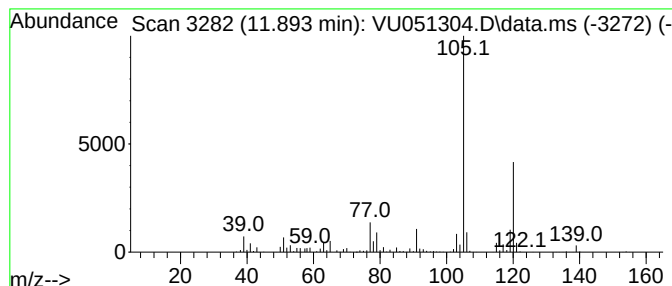
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.893	5.46 ug/L	1315950	1,4-Dichlorobenzene-d4	11.809

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



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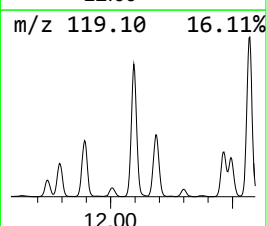
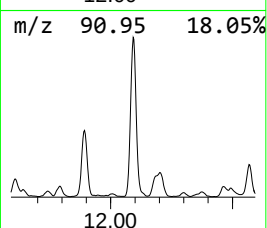
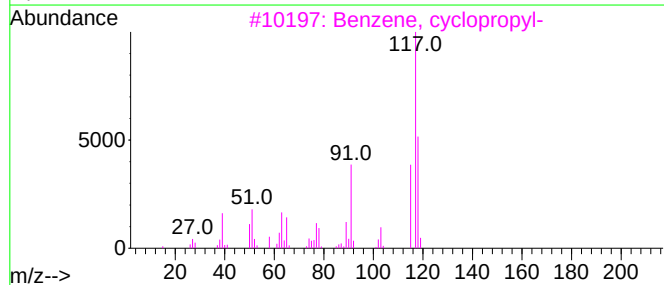
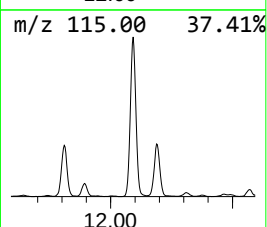
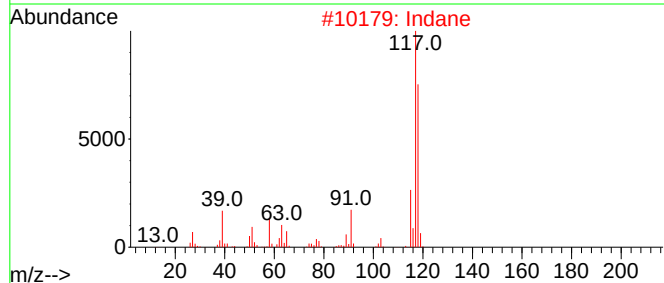
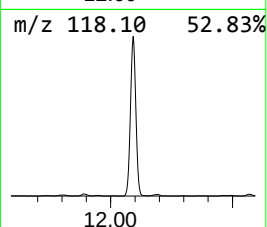
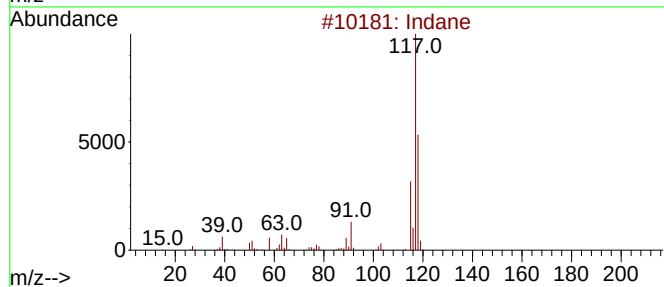
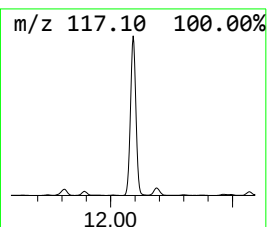
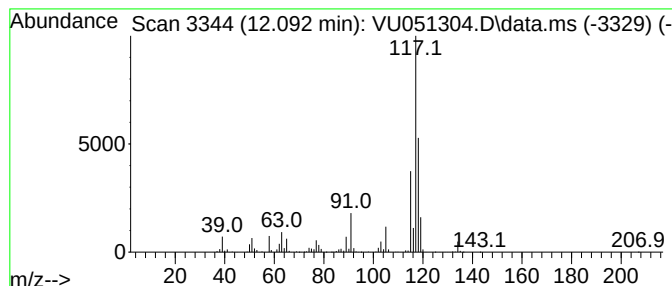
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.092	9.16 ug/L	2208850	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Indane	118	C9H10	000496-11-7	81
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
5			Deltacyclene	118	C9H10	007785-10-6	64



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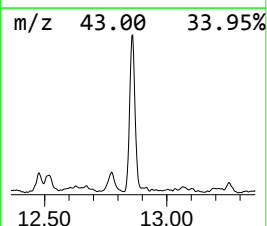
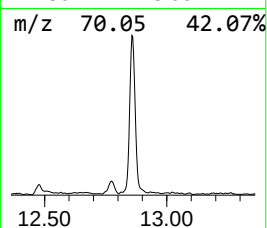
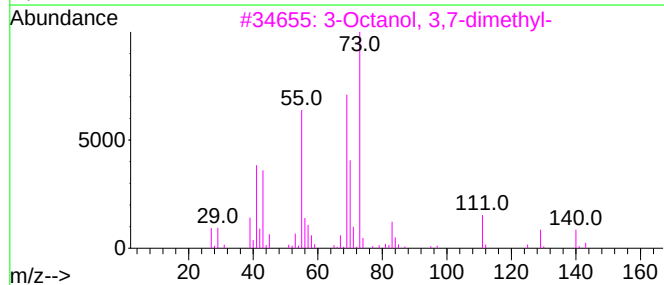
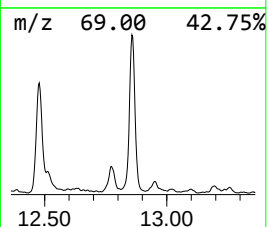
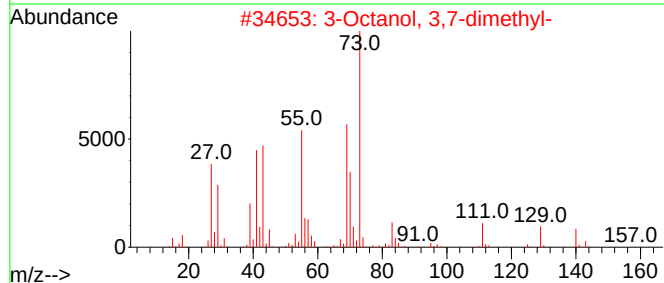
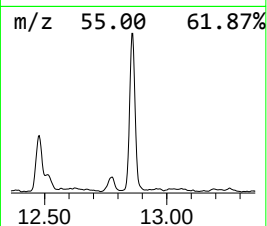
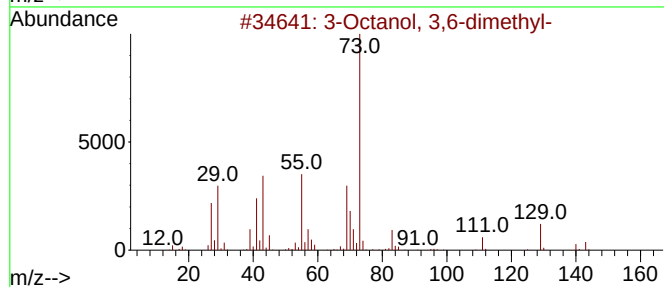
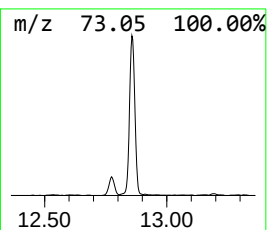
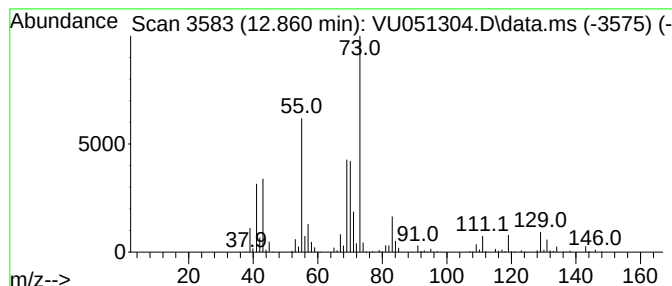
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.861	2.67 ug/L	642936	1,4-Dichlorobenzene-d4	11.809

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	43
2		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	42
3		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	38
4		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	35
5		Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100922\
Data File : VU051304.D
Acq On : 09 Oct 2022 16:51
Operator : JC/MD
Sample : N5013-03
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHA40

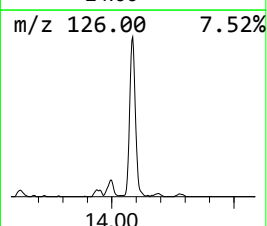
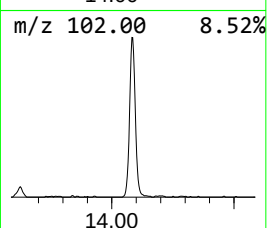
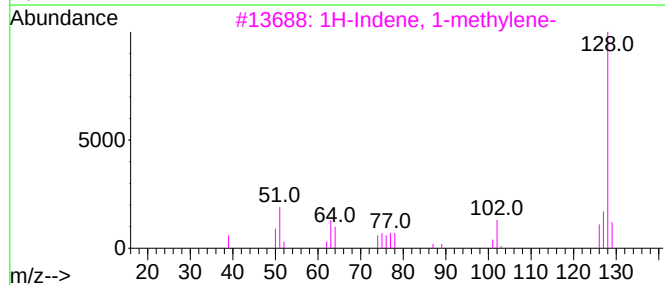
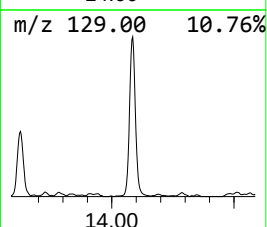
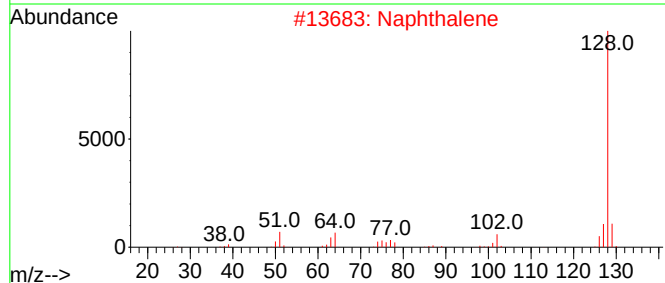
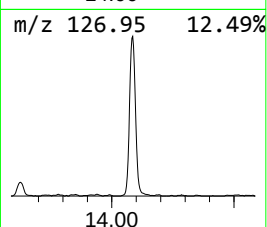
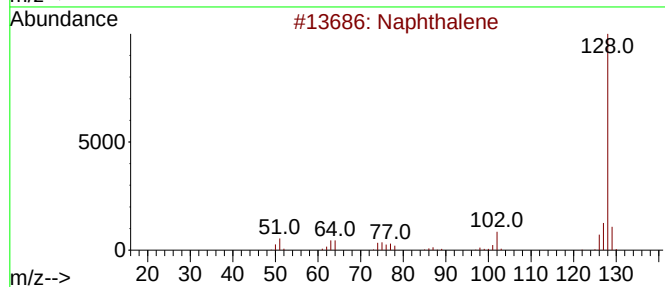
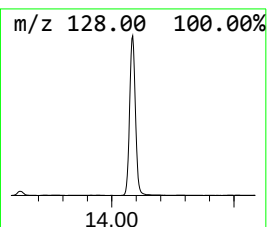
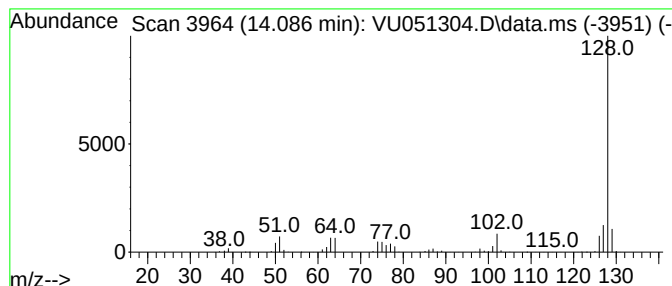
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 1H-Indene, 1-methylene- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	4.52 ug/L	1089050	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100922\
Data File : VU051304.D
Acq On : 09 Oct 2022 16:51
Operator : JC/MD
Sample : N5013-03
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHA40

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR100722WMA.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-	10.902	3.7	ug/L	892591	3	11.809	1205380	5.0
Benzene, 1-ethy...	10.996	4.2	ug/L	1011130	3	11.809	1205380	5.0
Benzene, 1-ethy...	11.288	3.5	ug/L	831922	3	11.809	1205380	5.0
Benzene, 1,2,4-...	11.893	5.5	ug/L	1315950	3	11.809	1205380	5.0
Indane	12.092	9.2	ug/L	2208850	3	11.809	1205380	5.0
unknown-01	12.861	2.7	ug/L	642936	3	11.809	1205380	5.0
1H-Indene, 1-me...	14.086	4.5	ug/L	1089050	3	11.809	1205380	5.0