

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
 Data File : VU059971.D
 Acq On : 16 Jul 2024 13:52
 Operator : MD/SY
 Sample : P3217-14DL 5X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 YDZJ9DL

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: VU059971.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.358	14	21	35	rBV3	9389	14218	2.08%	0.218%
2	1.592	82	94	109	rBV2	49872	87436	12.78%	1.340%
3	1.904	183	191	206	rBV	32319	60188	8.80%	0.922%
4	2.547	381	391	411	rVB	84132	143179	20.93%	2.194%
5	3.030	527	541	559	rBV2	40375	85291	12.47%	1.307%
6	4.657	1033	1047	1086	rBV	41667	149912	21.91%	2.297%
7	5.055	1156	1171	1189	rBV	69818	154491	22.58%	2.367%
8	5.714	1358	1376	1388	rBV2	141362	349955	51.15%	5.362%
9	6.242	1528	1540	1563	rBV	156230	305180	44.60%	4.676%
10	6.563	1624	1640	1656	rBV2	29480	62467	9.13%	0.957%
11	6.685	1666	1678	1691	rBV	83348	165522	24.19%	2.536%
12	6.750	1692	1698	1713	rVB3	12468	23860	3.49%	0.366%
13	7.563	1939	1951	1967	rBV	37632	65647	9.59%	1.006%
14	7.894	2042	2054	2067	rBV	147329	254998	37.27%	3.907%
15	7.965	2067	2076	2096	rVB	37895	68245	9.97%	1.046%
16	8.181	2130	2143	2156	rBV	22679	39196	5.73%	0.601%
17	8.640	2275	2286	2328	rBV	173907	369179	53.96%	5.656%
18	9.415	2515	2527	2550	rBV	236942	420607	61.47%	6.444%
19	9.695	2600	2614	2631	rBV	17445	30580	4.47%	0.469%
20	10.097	2727	2739	2754	rVB2	28926	47120	6.89%	0.722%
21	10.483	2846	2859	2871	rBV2	10347	16308	2.38%	0.250%
22	10.756	2932	2944	2956	rVV2	76570	124944	18.26%	1.914%
23	10.904	2977	2990	3006	rBV3	30760	52358	7.65%	0.802%
24	10.997	3006	3019	3023	rBV	38737	59901	8.75%	0.918%
25	11.087	3038	3047	3059	rVB2	26536	42877	6.27%	0.657%
26	11.290	3099	3110	3123	rBV	75335	116602	17.04%	1.786%
27	11.467	3153	3165	3177	rBV2	86577	135362	19.78%	2.074%
28	11.637	3199	3218	3230	rVB10	8779	18876	2.76%	0.289%
29	11.746	3244	3252	3260	rBV5	6487	9104	1.33%	0.139%
30	11.811	3260	3272	3290	rVV2	264310	518418	75.77%	7.943%
31	11.894	3290	3298	3317	rVB	15254	28340	4.14%	0.434%
32	12.094	3348	3360	3378	rBV3	13604	25625	3.75%	0.393%
33	12.206	3378	3395	3413	rVV2	297828	684229	100.00%	10.483%
34	12.377	3440	3448	3462	rBV2	4595	7366	1.08%	0.113%
35	12.573	3498	3509	3515	rBV2	4858	7324	1.07%	0.112%
36	12.682	3531	3543	3548	rBV3	7142	10397	1.52%	0.159%

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Title : TRACE VOA SFAM1.0

37	12.946	3616	3625	3642	rBV3	12047	27213	3.98%	0.417%
38	13.026	3642	3650	3663	rVB2	6687	10212	1.49%	0.156%
39	13.405	3759	3768	3774	rBV6	5338	9151	1.34%	0.140%
40	13.450	3774	3782	3796	rVV3	20248	36131	5.28%	0.554%
41	13.624	3827	3836	3844	rBV4	8350	14751	2.16%	0.226%
42	13.733	3860	3870	3878	rBV8	10325	17394	2.54%	0.266%
43	13.782	3878	3885	3891	rBV6	6047	8226	1.20%	0.126%
44	13.833	3892	3901	3913	rBV6	15380	35119	5.13%	0.538%
45	13.942	3931	3935	3951	rVB7	15689	29885	4.37%	0.458%
46	14.087	3963	3980	4000	rVB	31891	70572	10.31%	1.081%
47	14.190	4002	4012	4024	rVB3	15965	26367	3.85%	0.404%
48	14.286	4026	4042	4053	rBV6	28919	59781	8.74%	0.916%
49	14.344	4053	4060	4069	rVV8	8715	12268	1.79%	0.188%
50	14.483	4091	4103	4111	rBV4	25271	44429	6.49%	0.681%
51	14.576	4123	4132	4136	rBV3	8169	12734	1.86%	0.195%
52	14.669	4148	4161	4178	rBV4	41819	100779	14.73%	1.544%
53	14.807	4195	4204	4213	rVV2	38072	60948	8.91%	0.934%
54	14.872	4215	4224	4235	rVB2	42685	70322	10.28%	1.077%
55	14.936	4236	4244	4257	rVB4	10000	16913	2.47%	0.259%
56	15.016	4257	4269	4281	rBV5	41244	78893	11.53%	1.209%
57	15.219	4317	4332	4342	rBV	203704	352446	51.51%	5.400%
58	15.341	4363	4370	4378	rBV6	7284	12035	1.76%	0.184%
59	15.409	4379	4391	4402	rBV	297738	483872	70.72%	7.413%
60	15.463	4403	4408	4414	rVB7	8925	12120	1.77%	0.186%
61	15.926	4541	4552	4566	rVB2	21151	38832	5.68%	0.595%
62	16.145	4606	4620	4626	rBV4	12591	20781	3.04%	0.318%
63	16.183	4626	4632	4641	rVB6	10541	17121	2.50%	0.262%
64	16.264	4647	4657	4666	rBV3	17854	29776	4.35%	0.456%
65	16.408	4692	4702	4709	rBV2	28522	46078	6.73%	0.706%
66	16.450	4710	4715	4724	rVB3	11016	16546	2.42%	0.254%

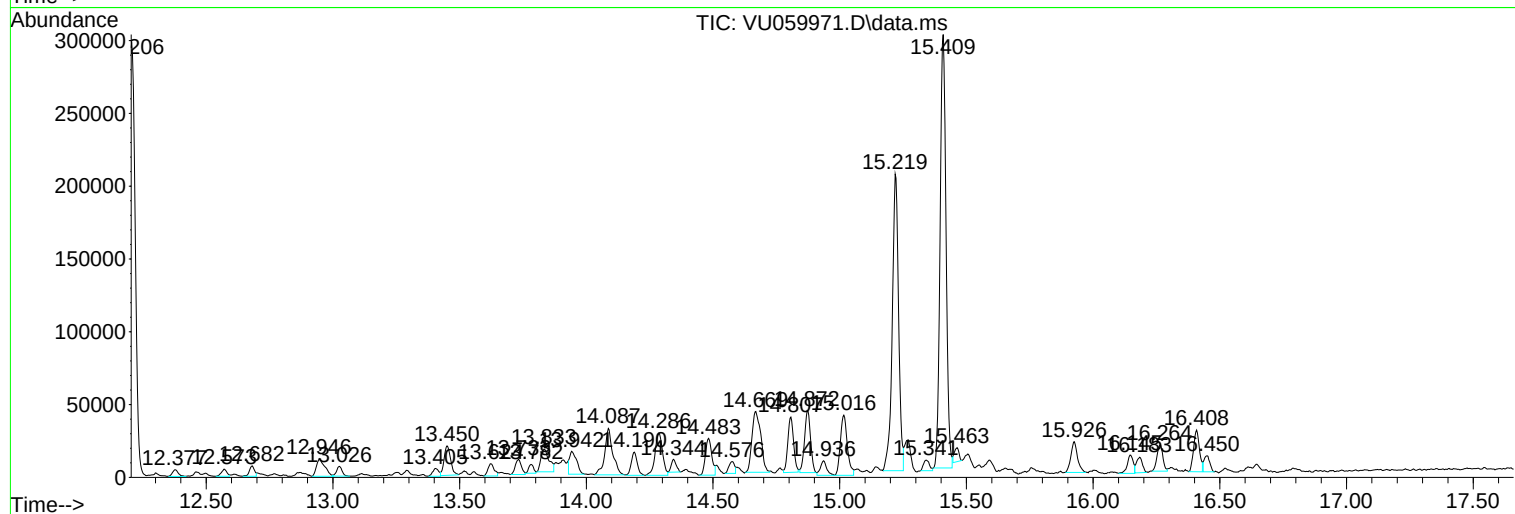
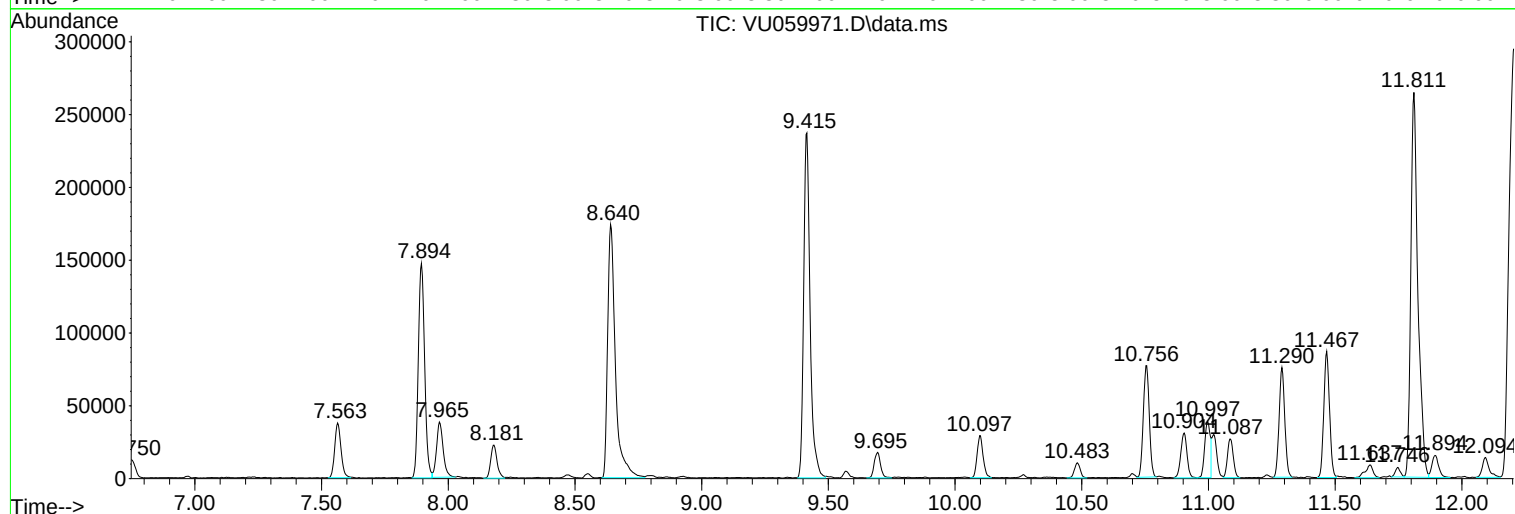
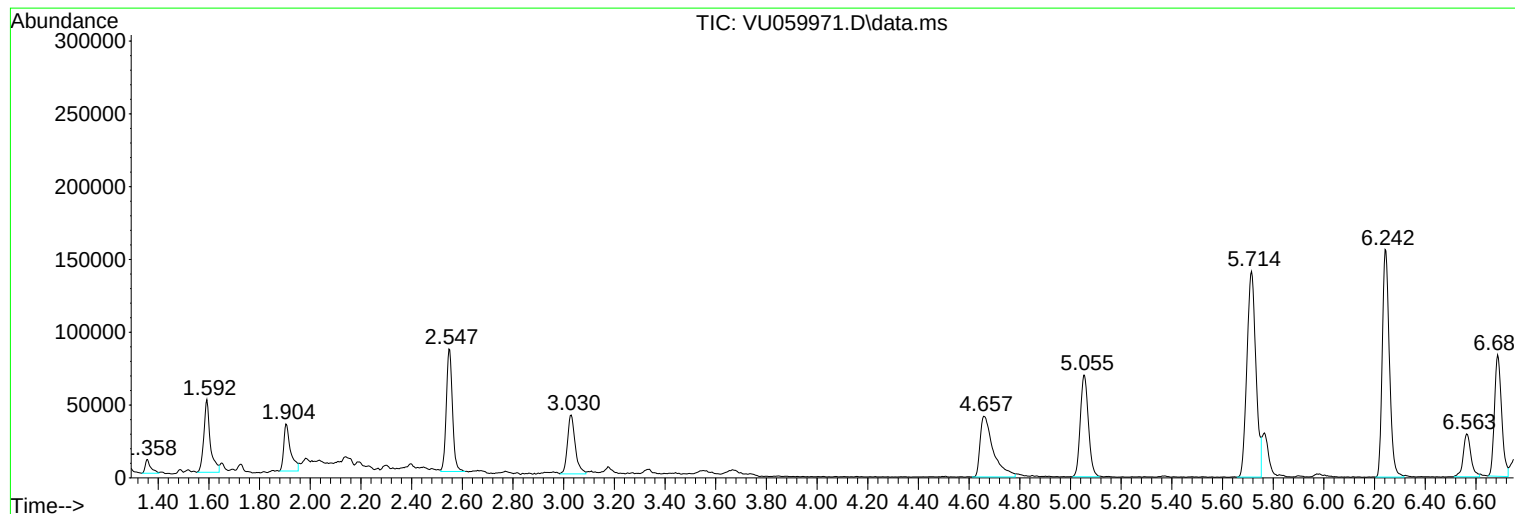
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

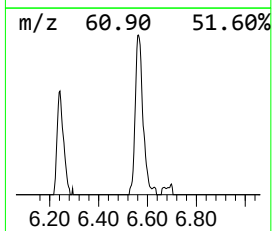
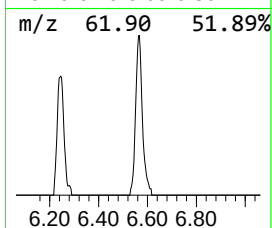
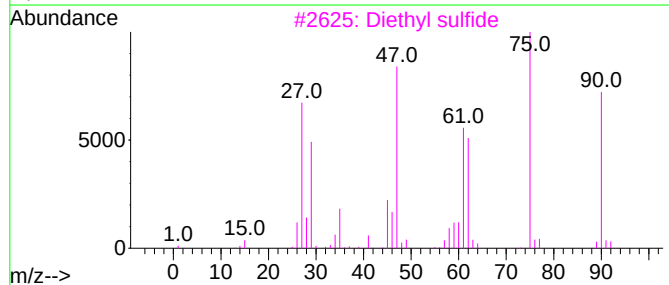
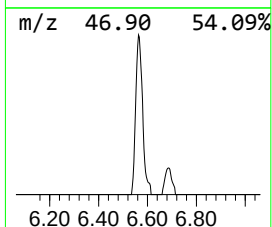
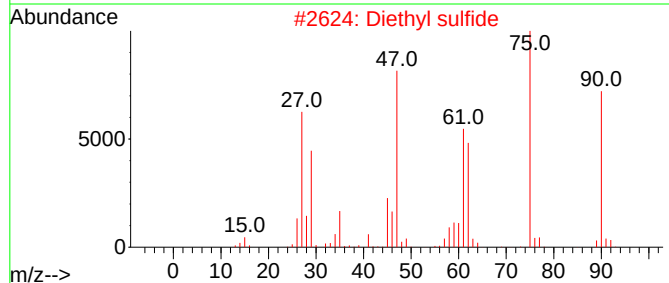
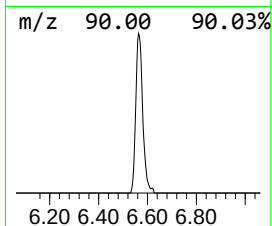
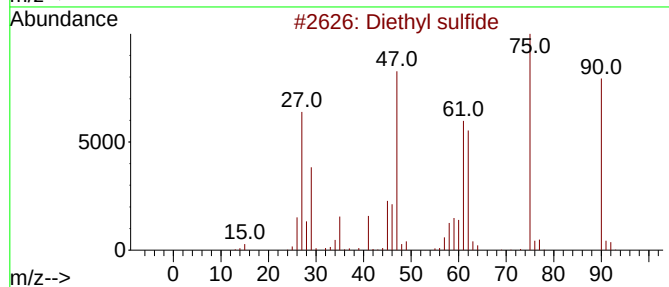
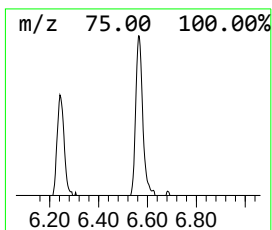
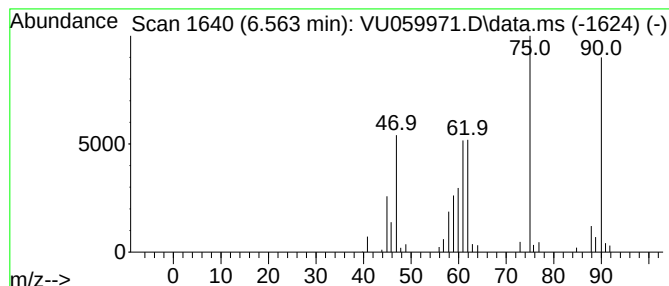
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Diethyl sulfide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.563	1.02 ug/L	62467	1,4-Difluorobenzene	6.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyl sulfide	90	C4H10S	000352-93-2	91
2			Diethyl sulfide	90	C4H10S	000352-93-2	91
3			Diethyl sulfide	90	C4H10S	000352-93-2	91
4			Diethyl sulfide	90	C4H10S	000352-93-2	91
5			Thiourea, methyl-	90	C2H6N2S	000598-52-7	38



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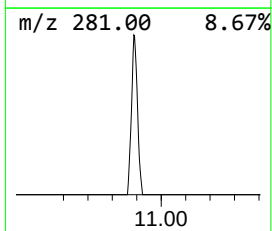
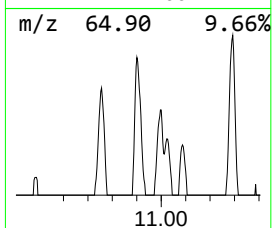
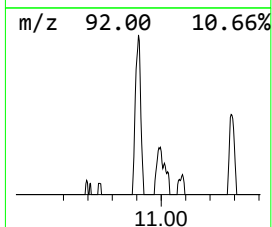
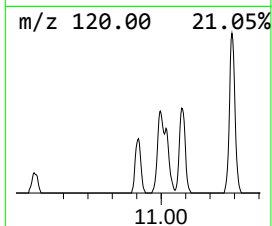
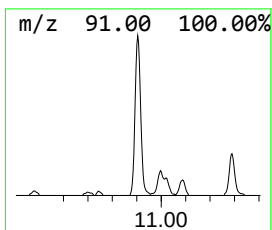
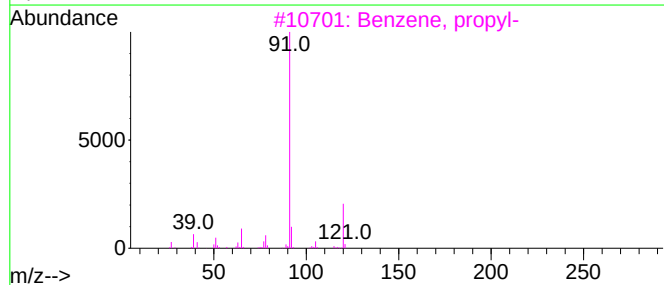
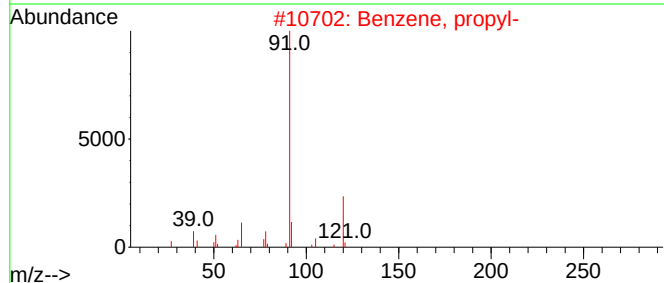
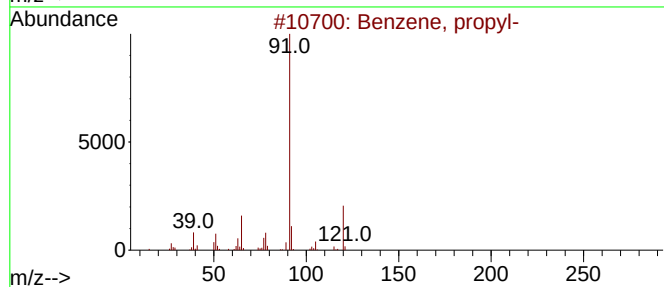
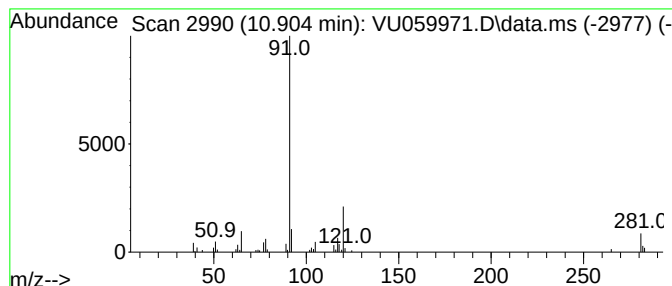
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.904	0.50 ug/L	52358	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	76
2			Benzene, propyl-	120	C9H12	000103-65-1	70
3			Benzene, propyl-	120	C9H12	000103-65-1	70
4			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64
5			1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	64



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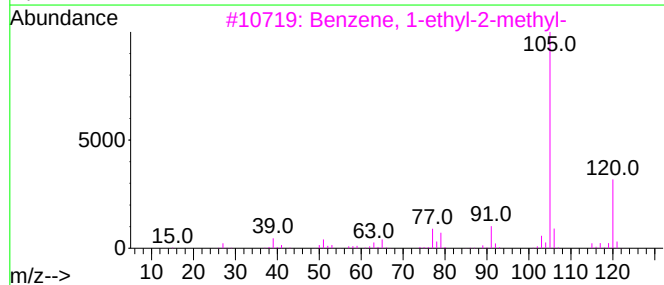
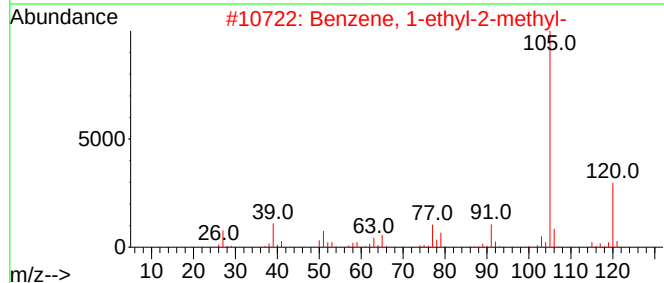
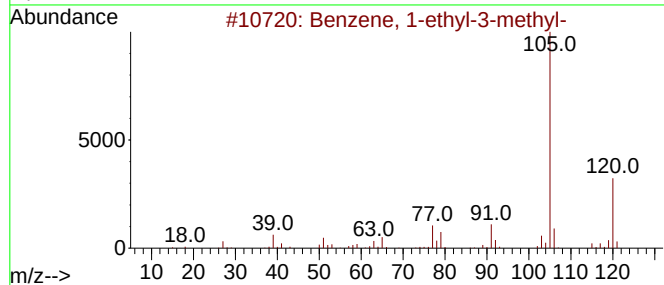
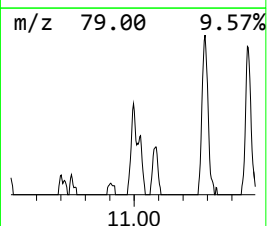
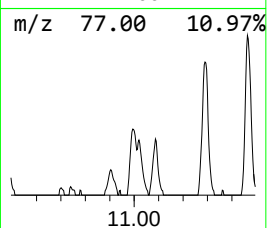
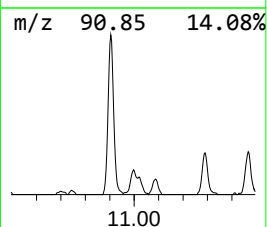
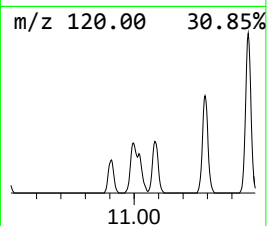
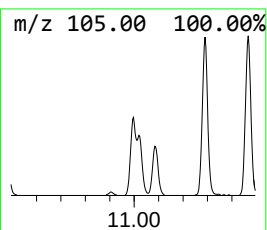
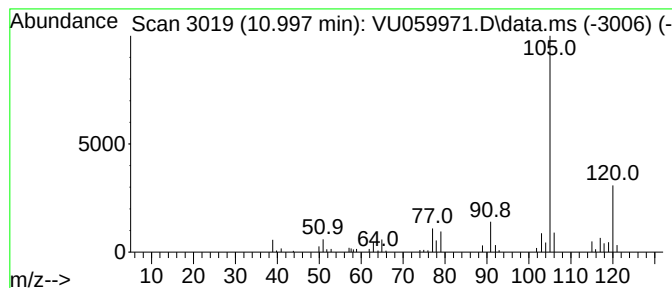
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.997	0.58 ug/L	59901	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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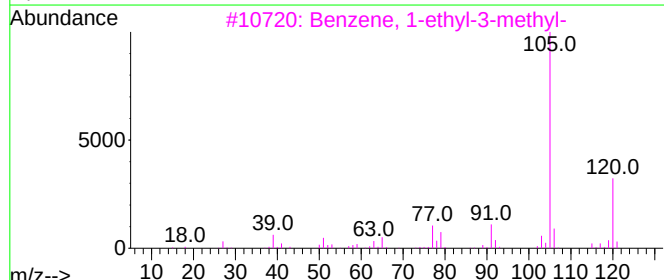
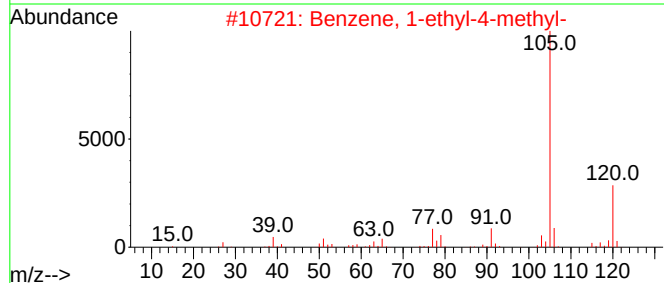
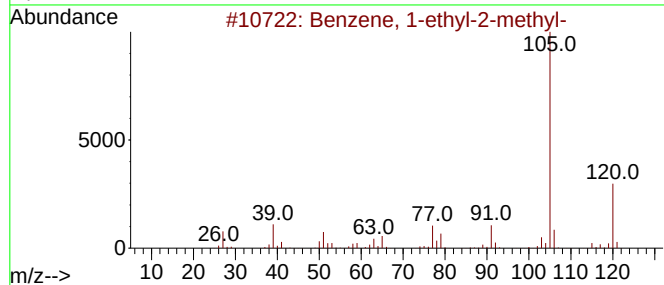
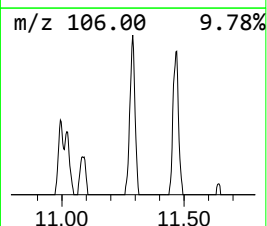
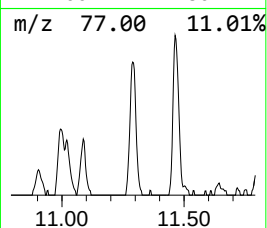
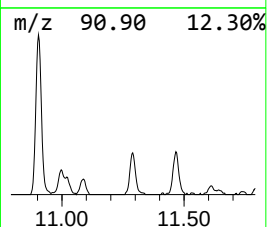
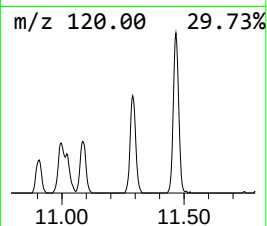
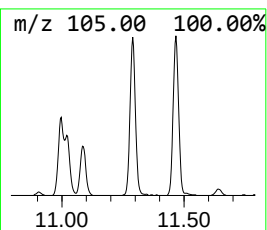
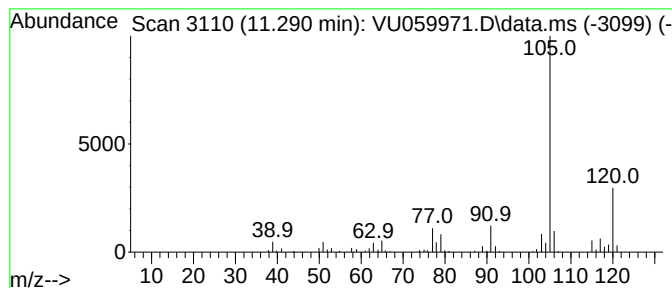
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.290	1.12 ug/L	116602	1,4-Dichlorobenzene-d4	11.811

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

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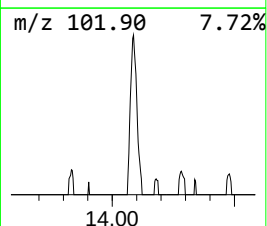
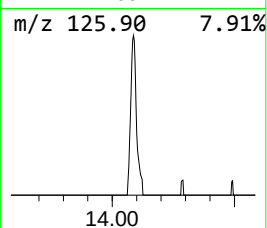
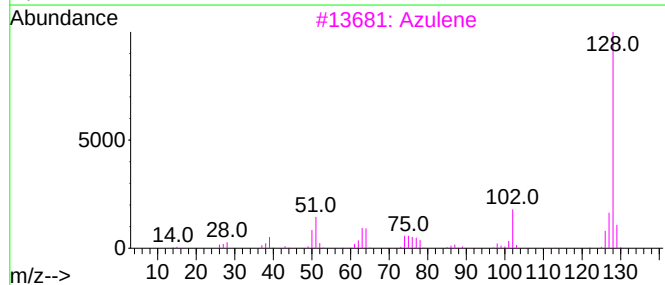
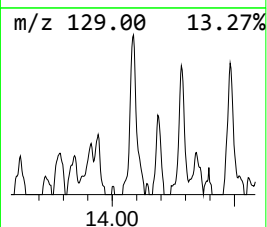
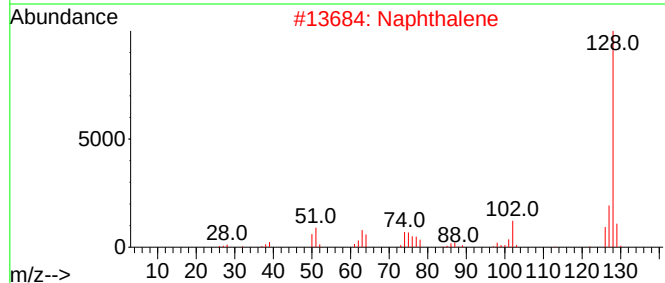
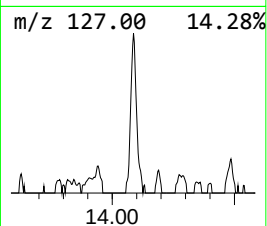
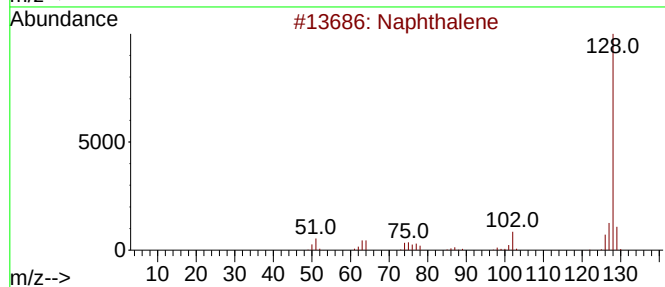
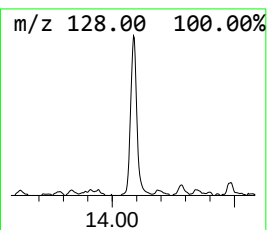
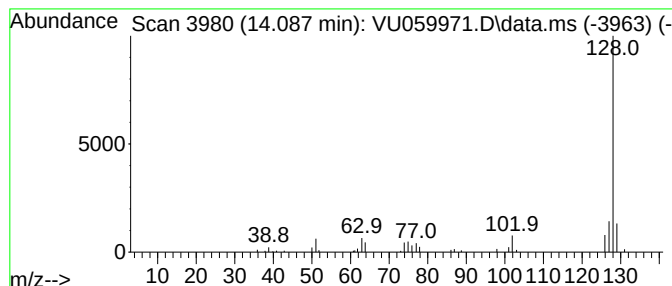
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.087	0.68 ug/L	70572	1,4-Dichlorobenzene-d4	11.811

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			Azulene	128	C10H8	000275-51-4	91
4			Naphthalene	128	C10H8	000091-20-3	90
5			Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

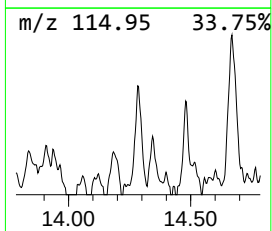
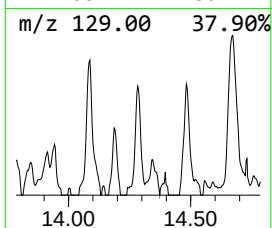
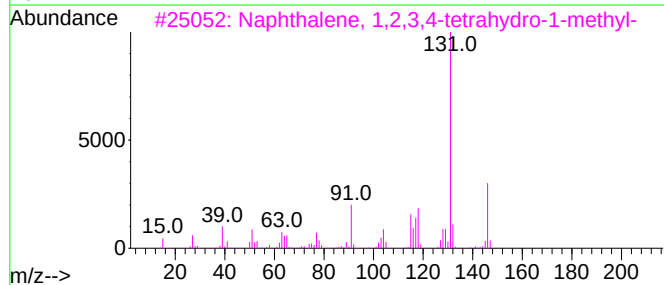
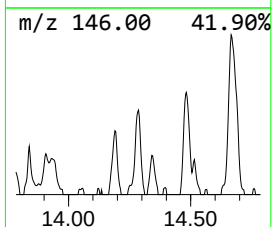
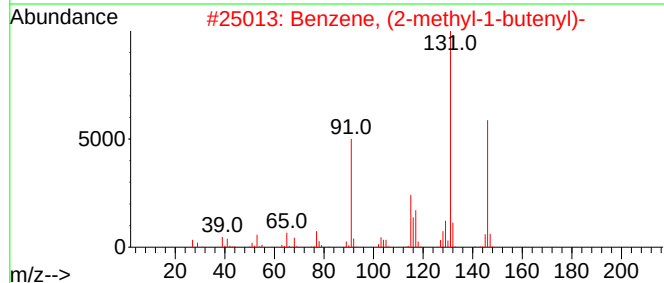
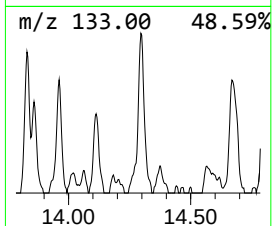
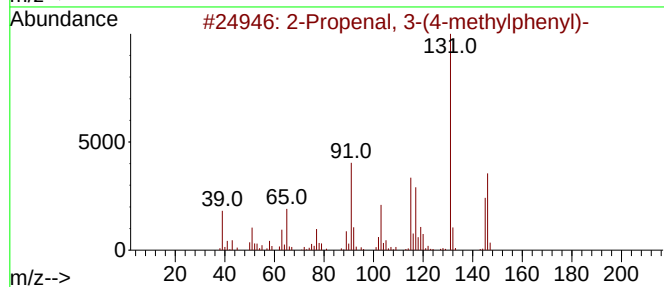
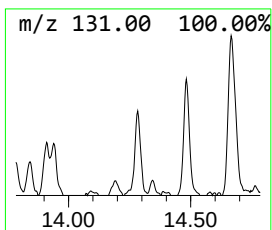
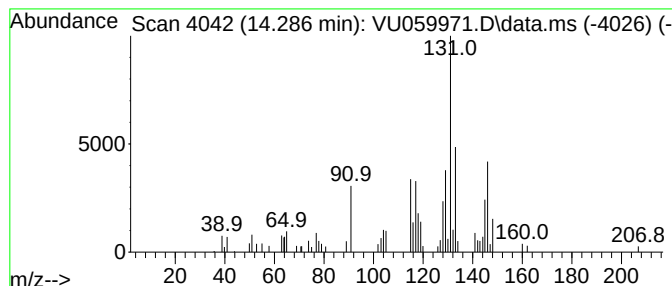
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 2-Propenal, 3-(4-methylphen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.286	0.58 ug/L	59781	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	64
2			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	60
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	55
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	55
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

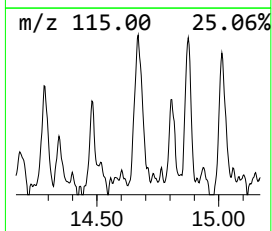
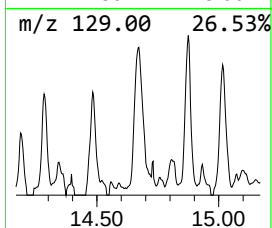
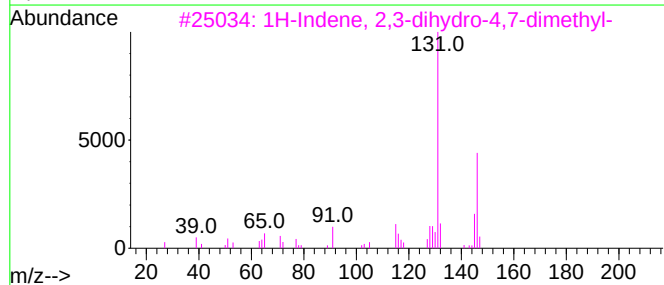
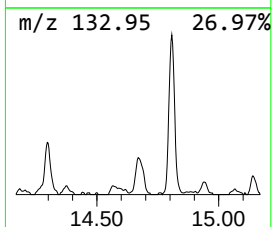
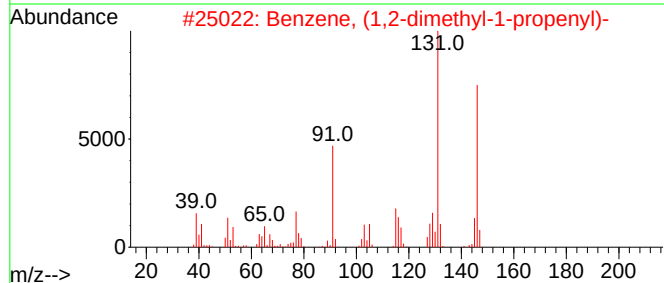
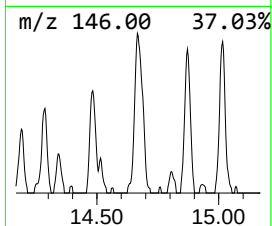
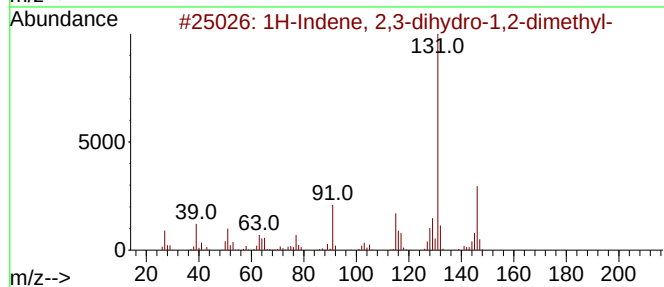
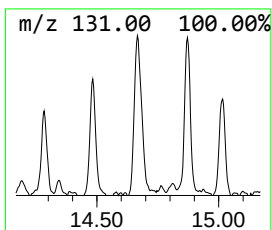
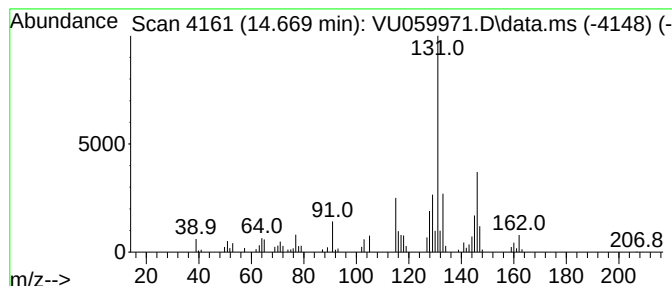
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.669	0.97 ug/L	100779	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	74
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

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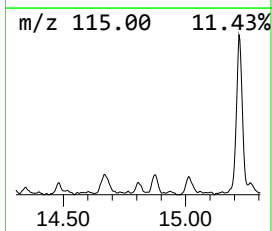
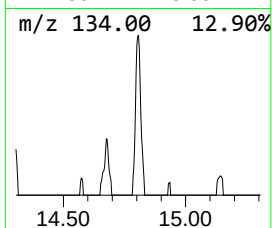
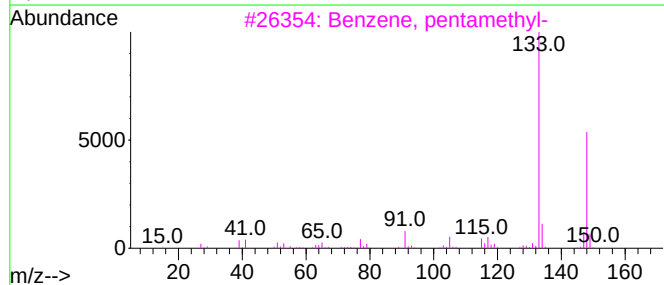
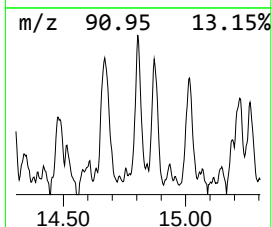
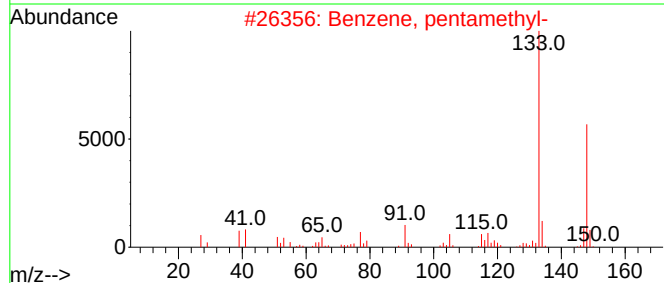
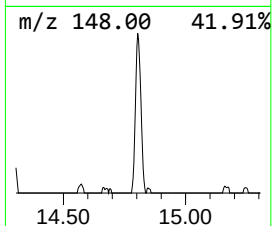
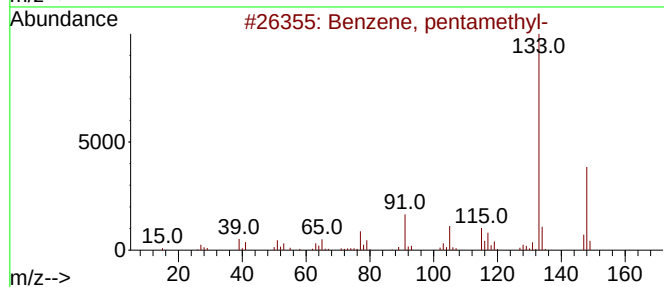
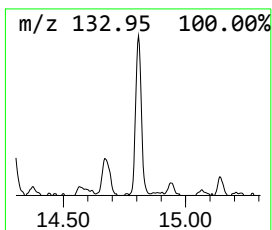
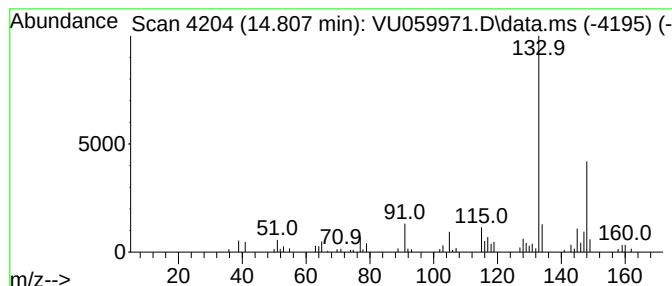
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, pentamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.807	0.59 ug/L	60948	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	95
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	87
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	87
4			Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	81
5			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

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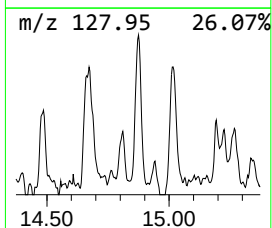
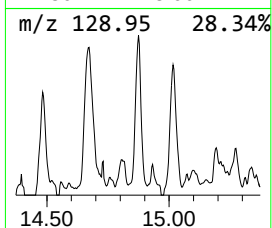
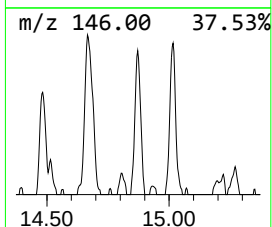
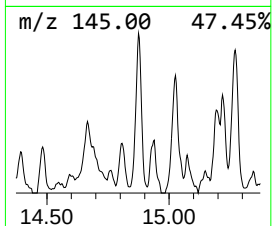
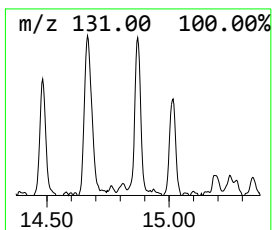
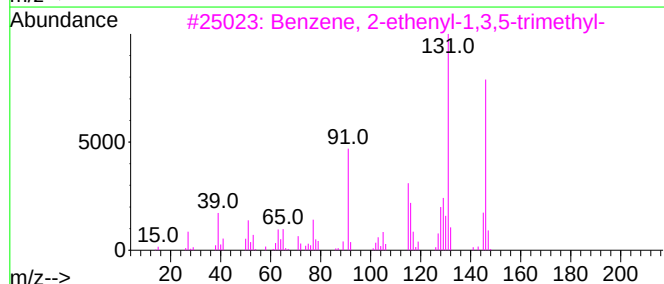
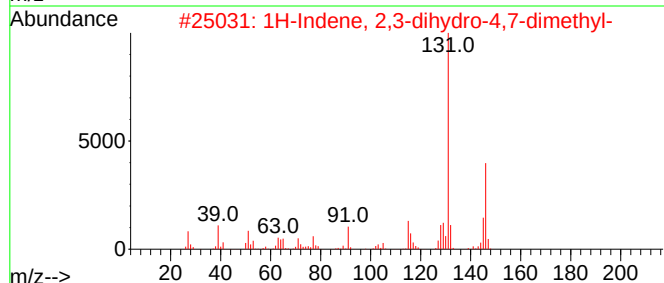
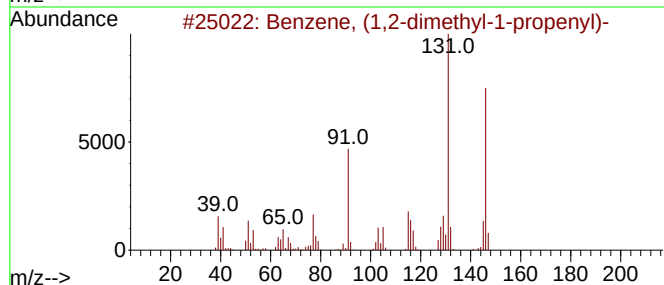
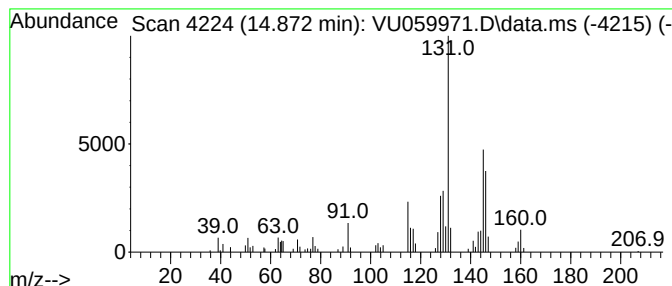
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.872	0.68 ug/L	70322	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	53
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	52
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	52
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	49
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

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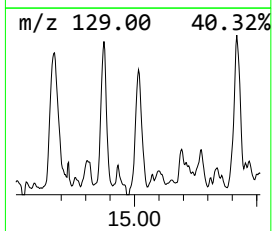
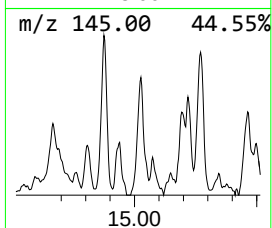
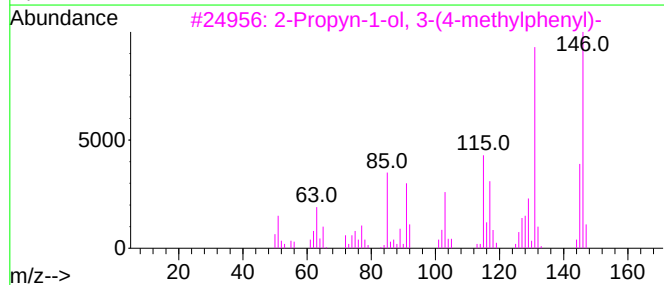
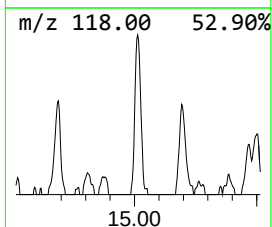
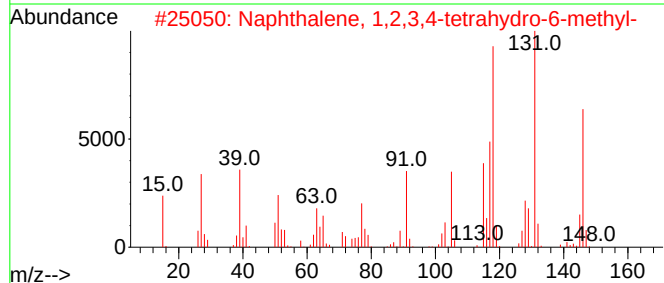
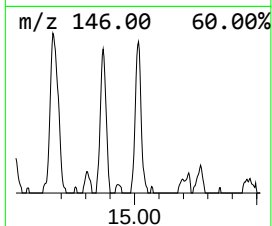
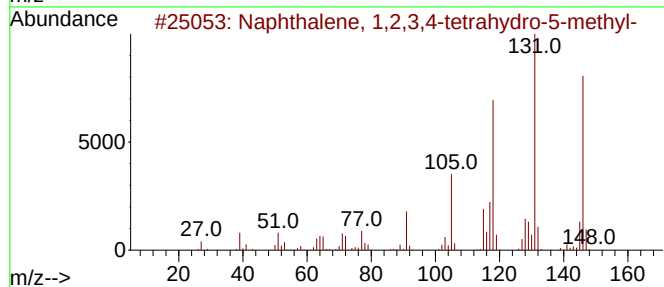
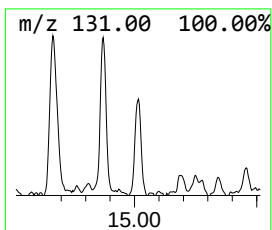
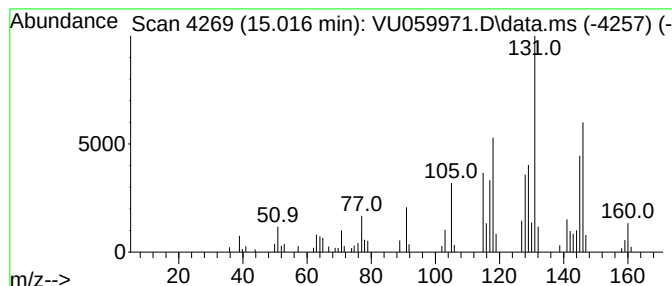
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.016	0.76 ug/L	78893	1,4-Dichlorobenzene-d4	11.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	60
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	55
5			1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071624\
Data File : VU059971.D
Acq On : 16 Jul 2024 13:52
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

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
Diethyl sulfide	6.563	1.0	ug/L	62467	1	6.245	305180	5.0
Benzene, propyl-	10.904	0.5	ug/L	52358	3	11.811	518418	5.0
Benzene, 1-ethy...	10.997	0.6	ug/L	59901	3	11.811	518418	5.0
Benzene, 1-ethy...	11.290	1.1	ug/L	116602	3	11.811	518418	5.0
Azulene	14.087	0.7	ug/L	70572	3	11.811	518418	5.0
2-Propenal, 3-(...	14.286	0.6	ug/L	59781	3	11.811	518418	5.0
1H-Indene, 2,3-...	14.669	1.0	ug/L	100779	3	11.811	518418	5.0
Benzene, pentam...	14.807	0.6	ug/L	60948	3	11.811	518418	5.0
Benzene, (1,2-d...	14.872	0.7	ug/L	70322	3	11.811	518418	5.0
Naphthalene, 1,...	15.016	0.8	ug/L	78893	3	11.811	518418	5.0