

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
 Data File : VU060057.D
 Acq On : 19 Jul 2024 13:58
 Operator : MD/SY
 Sample : P3217-14DL 5X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 6 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\SFAMUTR071824WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VU060057.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	15	21	37	rBV3	13330	16812	2.17%	0.240%
2	1.592	83	94	107	rBV3	59900	80802	10.43%	1.152%
3	1.907	182	192	207	rBV	43468	62335	8.04%	0.889%
4	2.560	383	395	414	rBV	75587	122361	15.79%	1.745%
5	3.039	534	544	559	rBV3	5298	10427	1.35%	0.149%
6	4.647	1027	1044	1075	rBV	39194	162295	20.94%	2.315%
7	5.055	1156	1171	1198	rBV	79899	180573	23.30%	2.575%
8	5.717	1359	1377	1389	rBV2	160211	414067	53.43%	5.905%
9	6.242	1525	1540	1560	rBV	146023	287839	37.14%	4.105%
10	6.563	1624	1640	1659	rBV2	35705	70094	9.04%	1.000%
11	6.682	1664	1677	1692	rVV	97022	195760	25.26%	2.792%
12	6.753	1692	1699	1711	rVB5	13668	27216	3.51%	0.388%
13	7.560	1938	1950	1972	rBV2	43132	80606	10.40%	1.150%
14	7.894	2042	2054	2067	rBV	177721	314498	40.58%	4.485%
15	7.965	2067	2076	2093	rVB	41584	74891	9.66%	1.068%
16	8.177	2132	2142	2157	rBV	24607	44041	5.68%	0.628%
17	8.631	2272	2283	2315	rBV	194604	417231	53.84%	5.950%
18	9.412	2513	2526	2550	rBV	213551	391134	50.47%	5.578%
19	9.692	2602	2613	2626	rBV2	19288	34301	4.43%	0.489%
20	10.097	2729	2739	2756	rVB2	32395	54036	6.97%	0.771%
21	10.479	2849	2858	2872	rBV2	13426	20826	2.69%	0.297%
22	10.749	2931	2942	2959	rVV2	87870	141443	18.25%	2.017%
23	10.901	2974	2989	3005	rBV2	37571	66920	8.64%	0.954%
24	10.994	3005	3018	3022	rBV2	45991	72143	9.31%	1.029%
25	11.084	3037	3046	3064	rVB	33015	49922	6.44%	0.712%
26	11.286	3098	3109	3126	rVB	88356	139311	17.98%	1.987%
27	11.463	3151	3164	3176	rBV	101951	161051	20.78%	2.297%
28	11.637	3201	3218	3229	rBV9	8745	20154	2.60%	0.287%
29	11.743	3243	3251	3258	rBV3	6850	9707	1.25%	0.138%
30	11.807	3258	3271	3289	rVV2	232736	484545	62.53%	6.910%
31	11.891	3289	3297	3311	rVB	18599	32384	4.18%	0.462%
32	12.087	3343	3358	3367	rBV3	16621	29860	3.85%	0.426%
33	12.200	3377	3393	3417	rVV2	327596	774961	100.00%	11.052%
34	12.373	3439	3447	3457	rVB3	6151	8293	1.07%	0.118%
35	12.566	3499	3507	3514	rBV3	6130	8908	1.15%	0.127%
36	12.675	3530	3541	3549	rBV3	8438	14225	1.84%	0.203%

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 Title : TRACE VOA SFAM1.0

37	12.942	3613	3624	3630	rBV3	11303	20643	2.66%	0.294%
38	13.023	3641	3649	3657	rVB4	7815	11126	1.44%	0.159%
39	13.399	3757	3766	3773	rBV8	6280	10997	1.42%	0.157%
40	13.447	3773	3781	3794	rVV4	23443	43594	5.63%	0.622%
41	13.508	3795	3800	3810	rVV7	6301	10386	1.34%	0.148%
42	13.621	3826	3835	3844	rBV4	8842	15789	2.04%	0.225%
43	13.730	3858	3869	3876	rBV10	10939	18497	2.39%	0.264%
44	13.778	3876	3884	3891	rBV4	7664	11877	1.53%	0.169%
45	13.830	3891	3900	3911	rBV6	18415	39669	5.12%	0.566%
46	13.907	3917	3924	3927	rBV4	6488	9586	1.24%	0.137%
47	13.939	3928	3934	3953	rVB6	19042	42955	5.54%	0.613%
48	14.084	3963	3979	4000	rVB2	35821	85302	11.01%	1.217%
49	14.183	4001	4010	4022	rVB4	19595	32816	4.23%	0.468%
50	14.280	4022	4040	4051	rBV5	34487	72869	9.40%	1.039%
51	14.338	4051	4058	4069	rVV7	10644	20513	2.65%	0.293%
52	14.389	4069	4074	4091	rVB8	5604	12263	1.58%	0.175%
53	14.479	4091	4102	4120	rBV2	30110	60019	7.74%	0.856%
54	14.569	4120	4130	4147	rBV2	11506	24853	3.21%	0.354%
55	14.666	4148	4160	4176	rVV5	47982	116895	15.08%	1.667%
56	14.801	4194	4202	4213	rBV	42194	62478	8.06%	0.891%
57	14.868	4214	4223	4234	rVB2	47995	81074	10.46%	1.156%
58	14.933	4235	4243	4253	rVB5	10977	16836	2.17%	0.240%
59	15.010	4257	4267	4282	rBV4	45993	88958	11.48%	1.269%
60	15.142	4300	4308	4314	rBV6	6430	9176	1.18%	0.131%
61	15.215	4315	4331	4341	rBV	196251	352525	45.49%	5.028%
62	15.338	4362	4369	4376	rBV3	9289	12763	1.65%	0.182%
63	15.405	4378	4390	4401	rBV	292573	478055	61.69%	6.818%
64	15.495	4415	4418	4428	rVB7	10351	14506	1.87%	0.207%
65	15.585	4440	4446	4456	rVB8	8179	13575	1.75%	0.194%
66	15.920	4539	4550	4563	rVB3	21742	39349	5.08%	0.561%
67	16.145	4611	4620	4626	rBV5	11099	16685	2.15%	0.238%
68	16.177	4626	4630	4639	rVB3	8723	10687	1.38%	0.152%
69	16.260	4646	4656	4665	rBV4	14126	23627	3.05%	0.337%
70	16.402	4691	4700	4708	rBV2	22863	39359	5.08%	0.561%
71	16.444	4710	4713	4726	rVB3	10970	15580	2.01%	0.222%

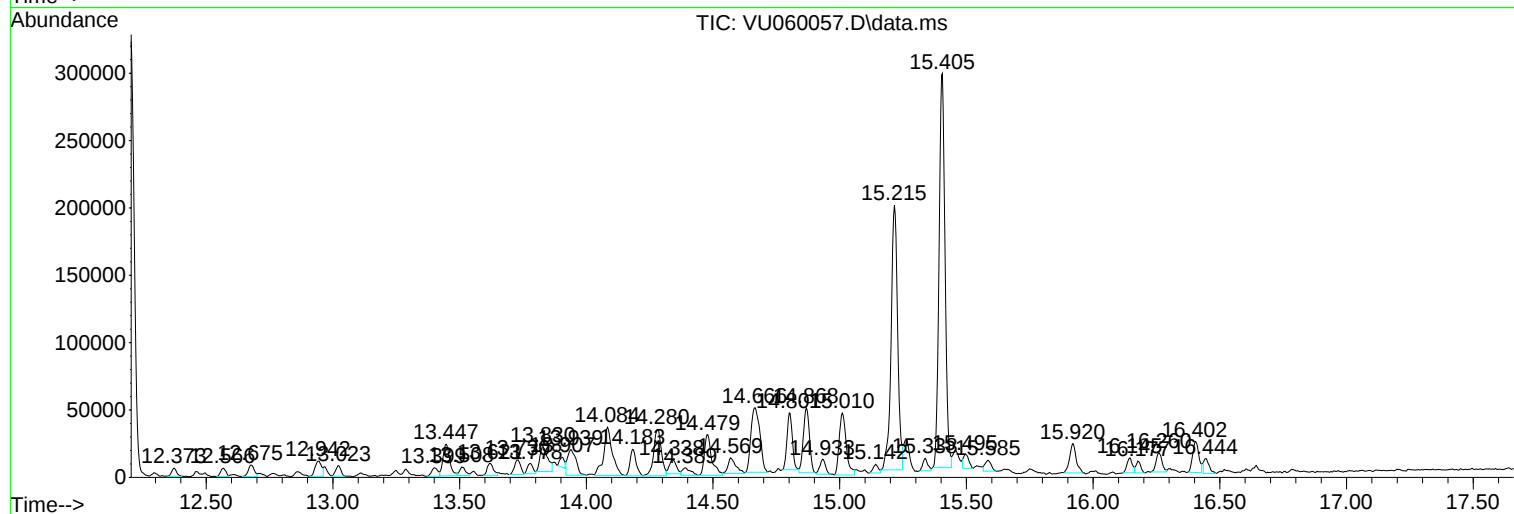
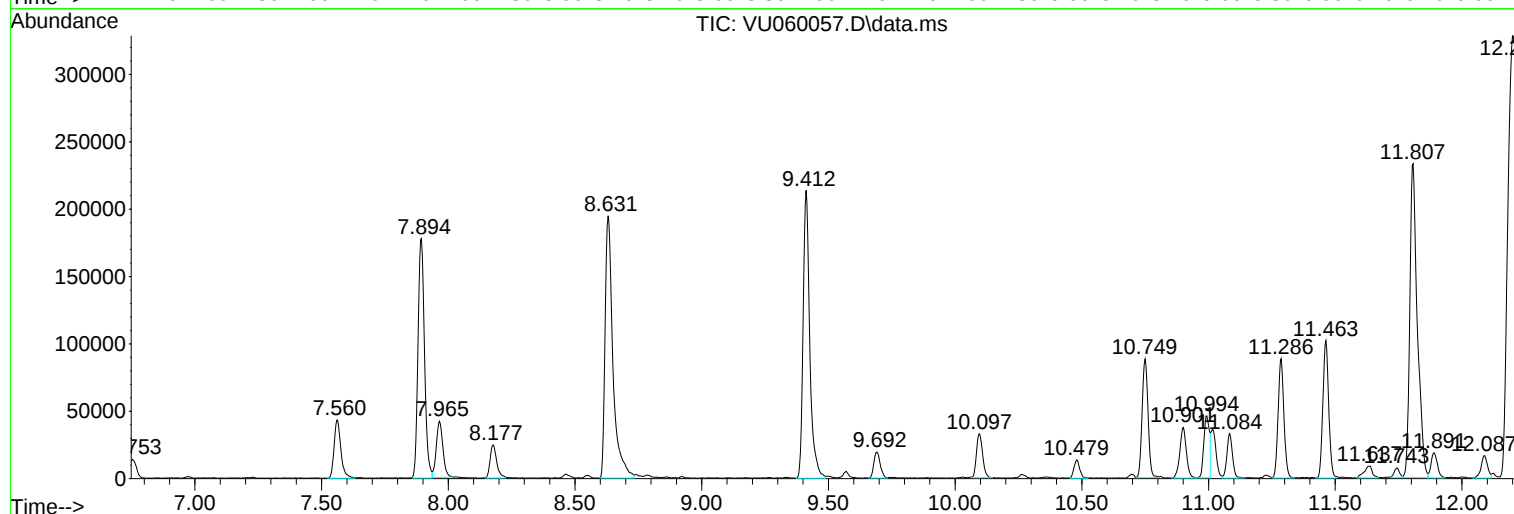
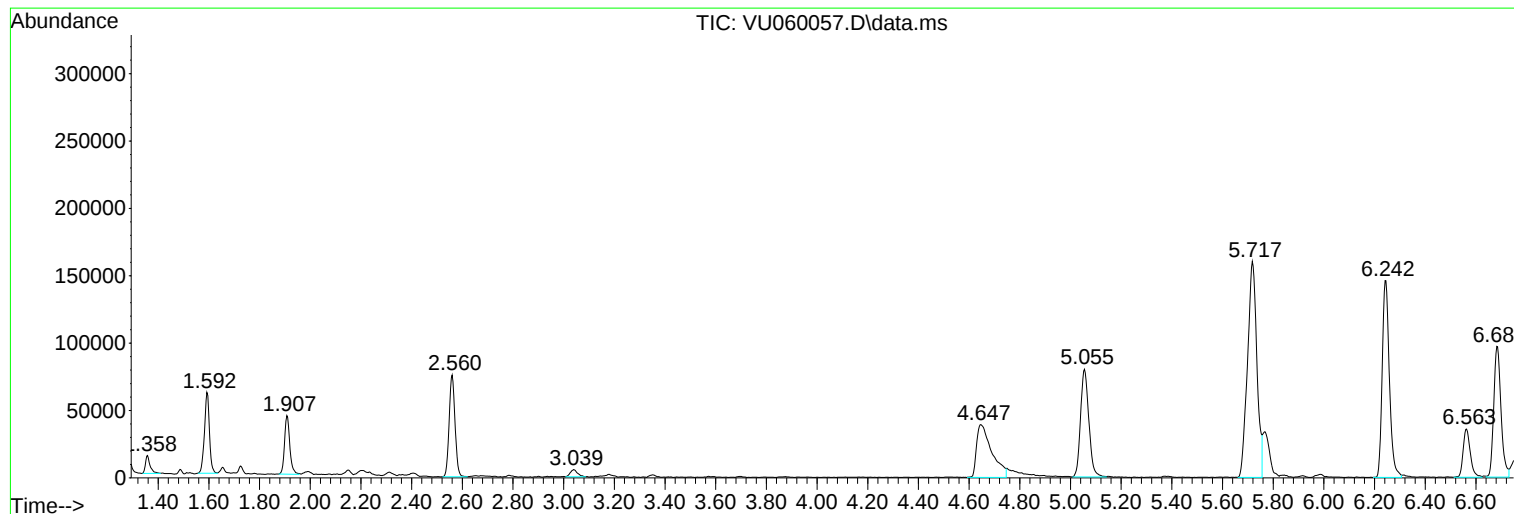
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

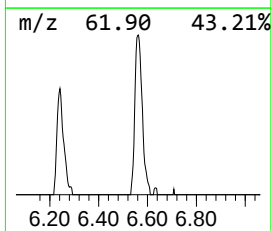
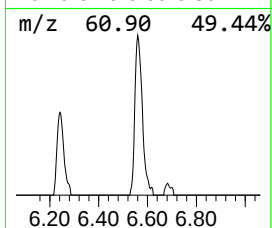
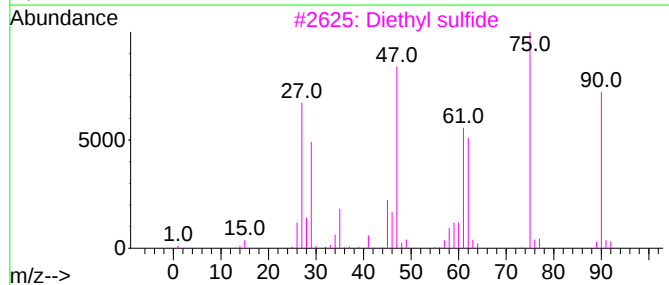
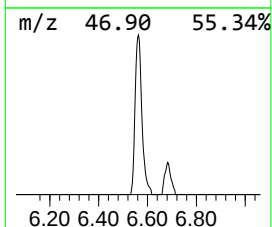
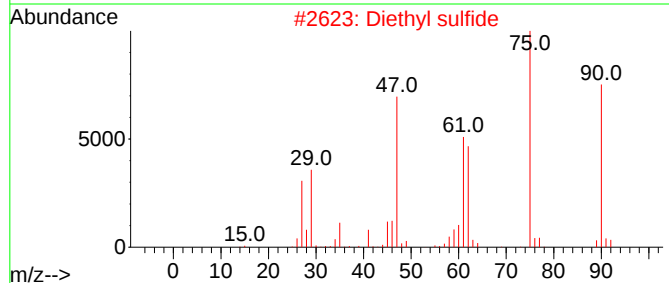
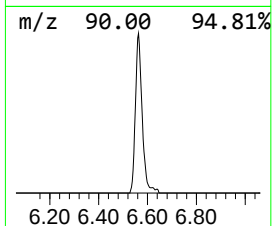
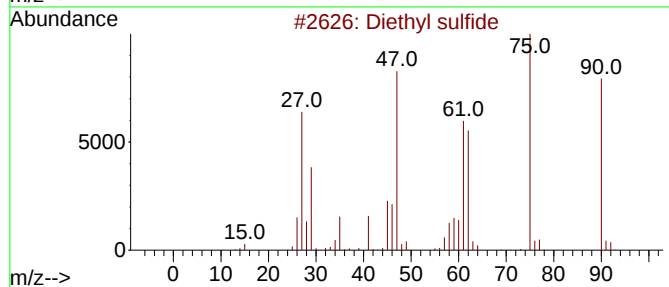
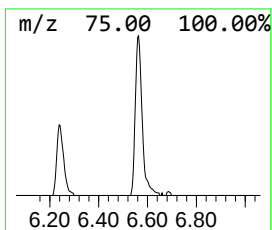
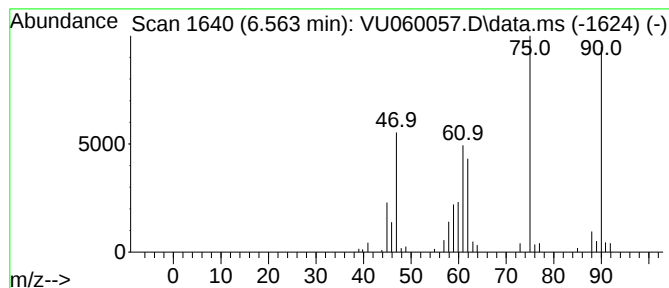
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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	1.22 ug/L	70094	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			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	47



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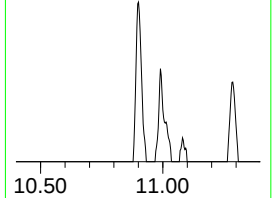
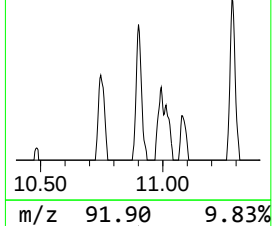
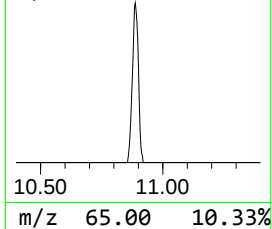
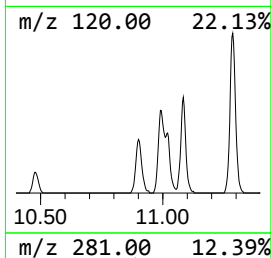
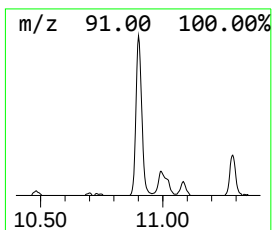
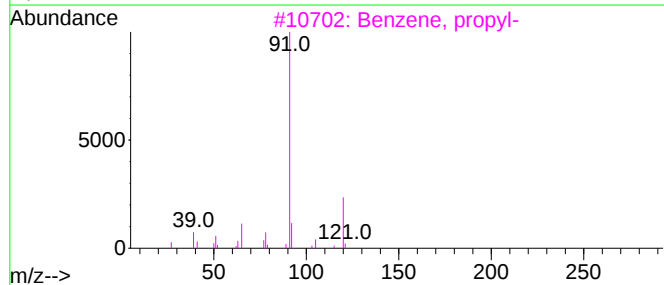
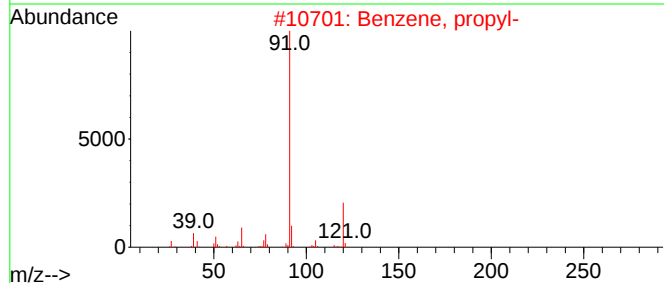
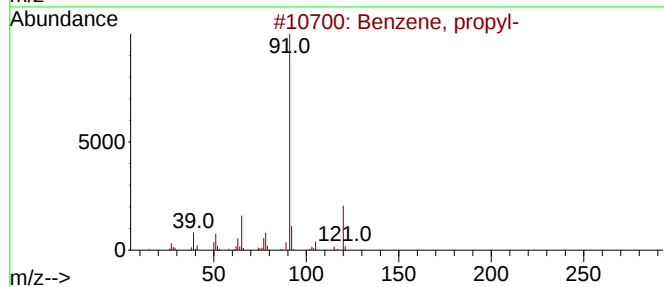
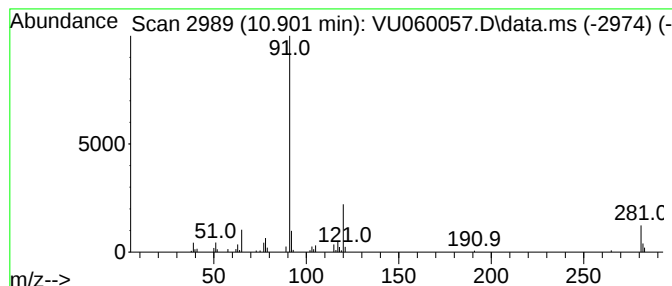
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.900	0.69 ug/L	66920	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	70
2			Benzene, propyl-	120	C9H12	000103-65-1	70
3			Benzene, propyl-	120	C9H12	000103-65-1	70
4			Benzene, propyl-	120	C9H12	000103-65-1	62
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	53



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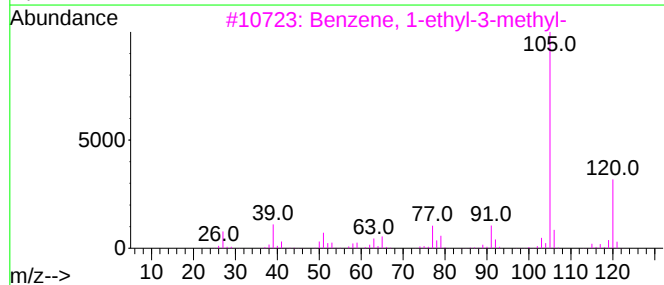
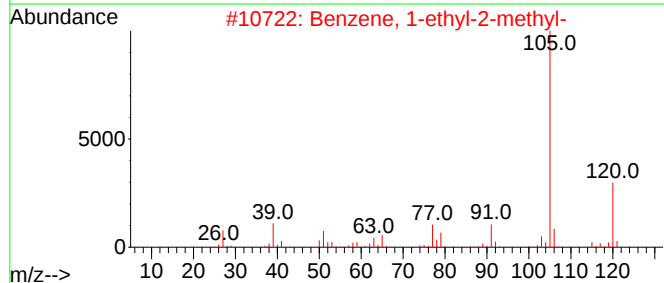
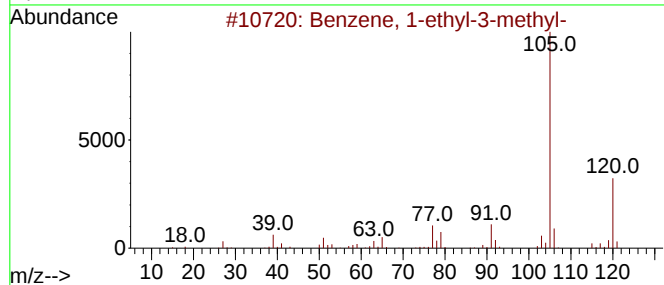
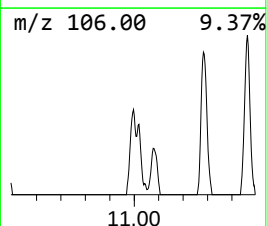
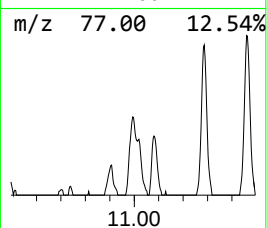
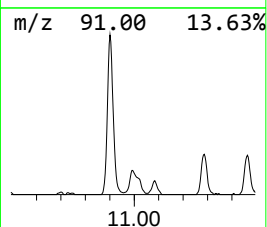
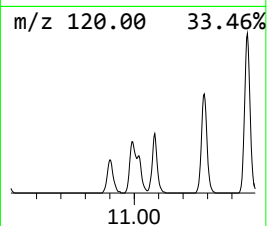
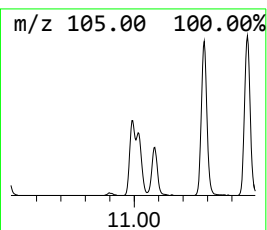
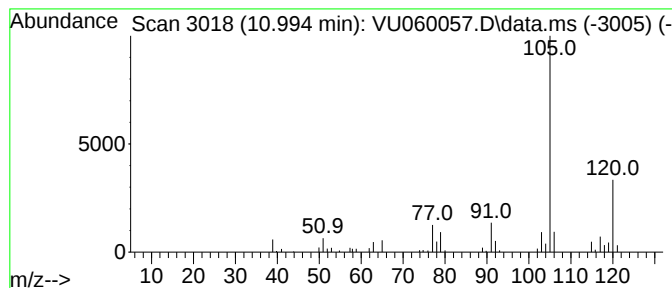
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Quant Title : TRACE VOA SFAM1.0

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.
10.994	0.74 ug/L	72143	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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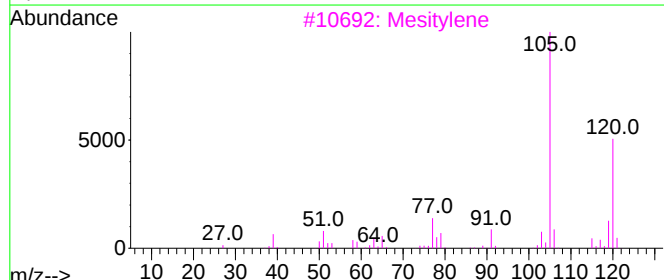
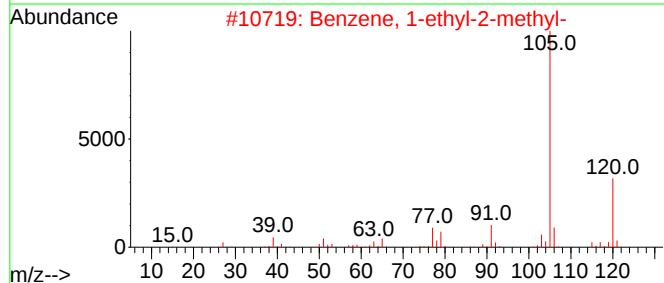
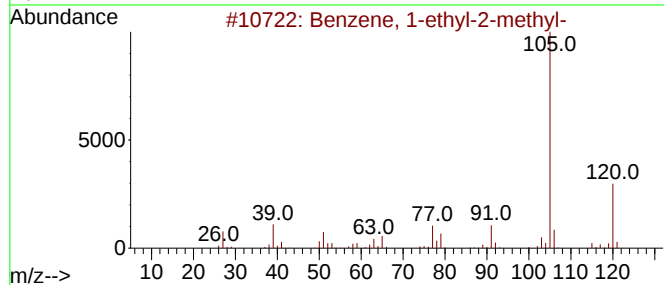
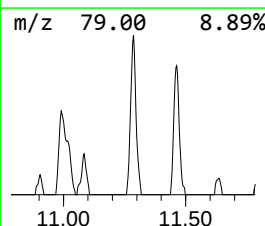
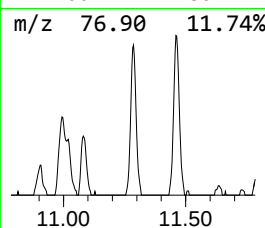
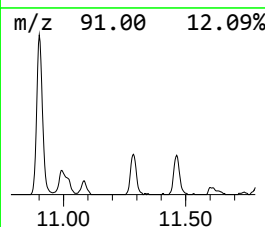
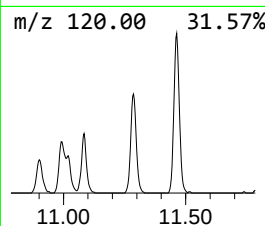
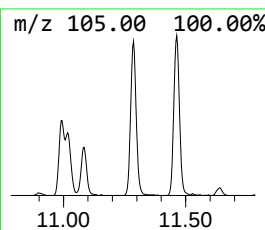
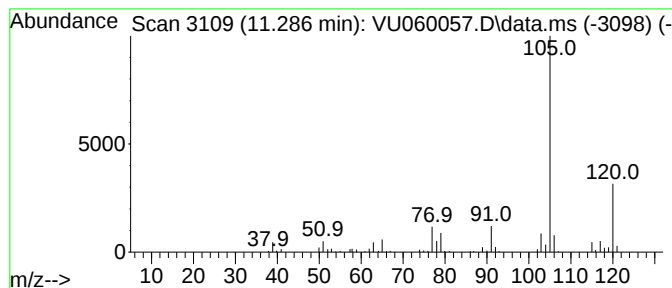
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.286	1.44 ug/L	139311	1,4-Dichlorobenzene-d4	11.807

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			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

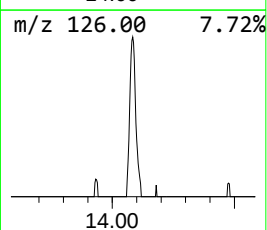
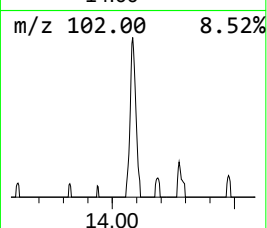
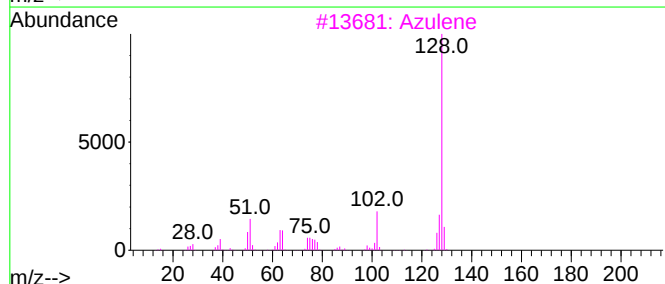
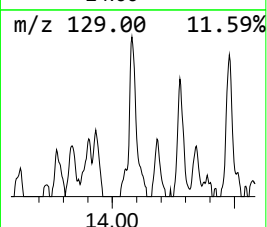
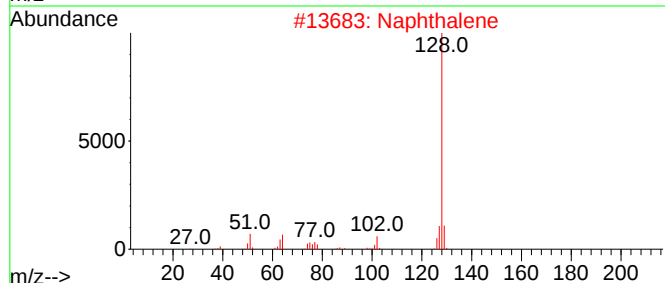
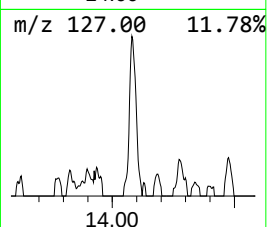
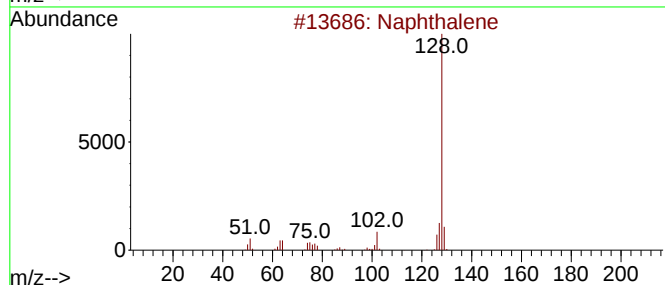
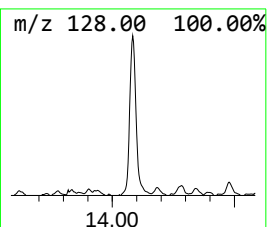
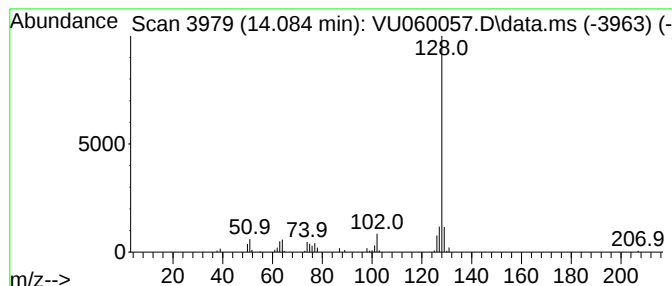
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.084	0.88 ug/L	85302	1,4-Dichlorobenzene-d4	11.807

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			Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

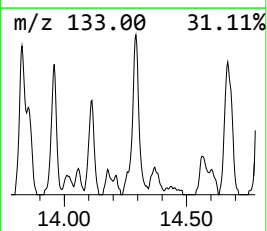
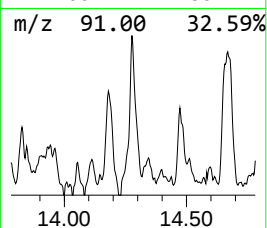
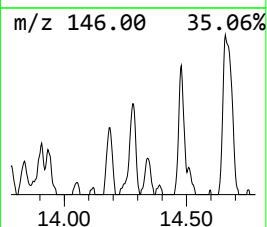
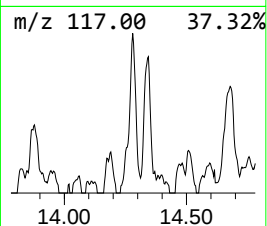
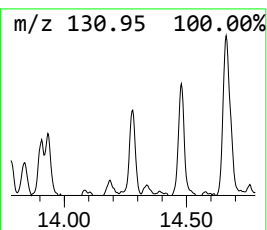
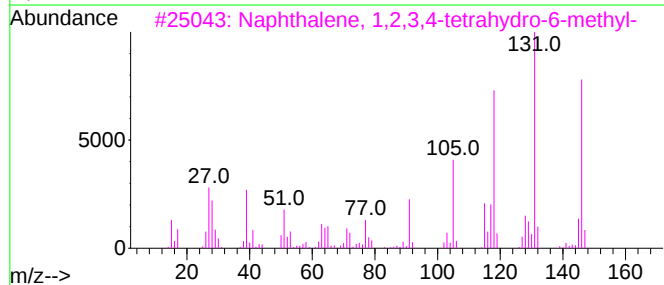
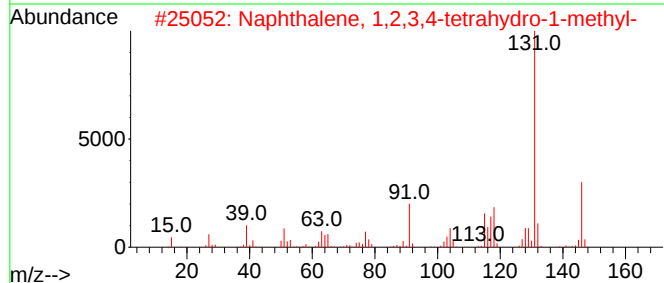
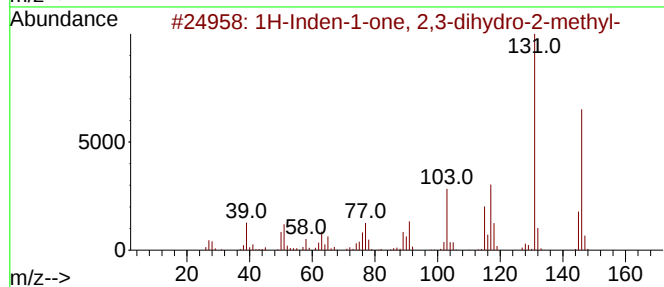
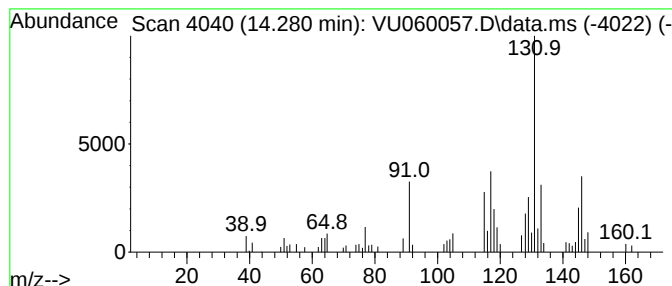
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.280	0.75 ug/L	72869	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	70
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	53
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

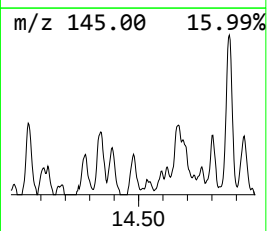
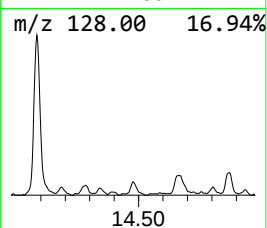
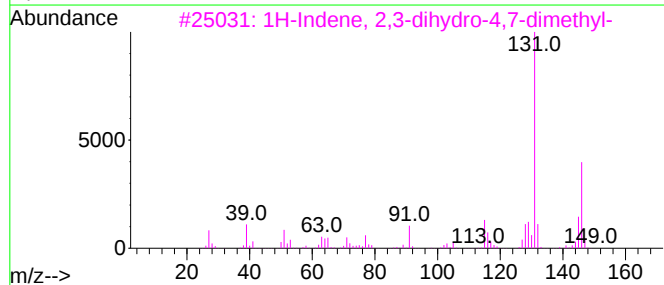
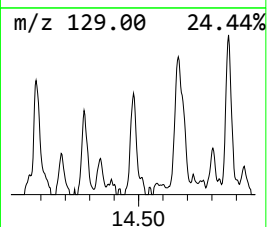
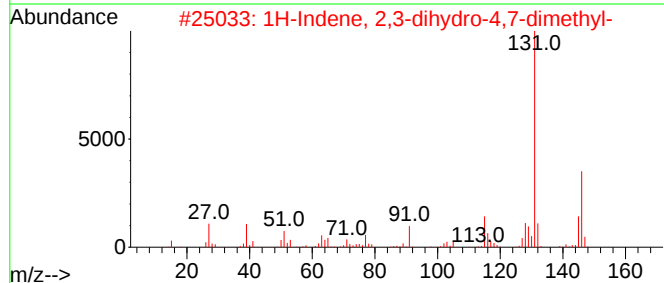
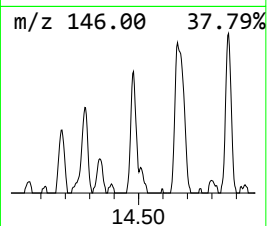
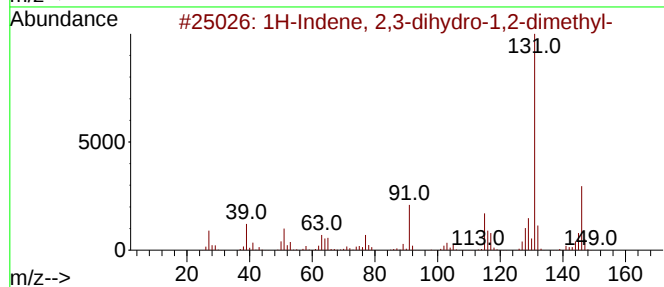
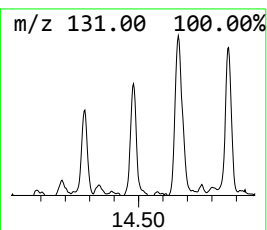
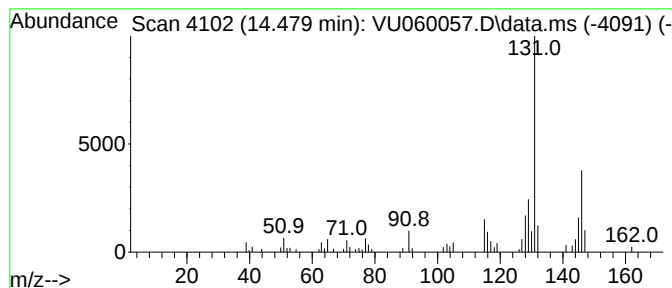
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.479	0.62 ug/L	60019	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

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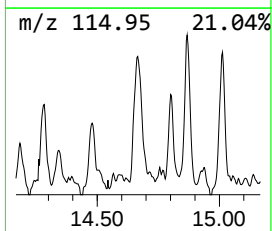
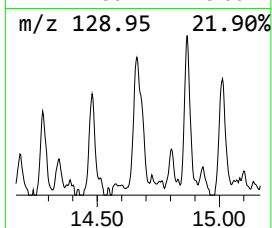
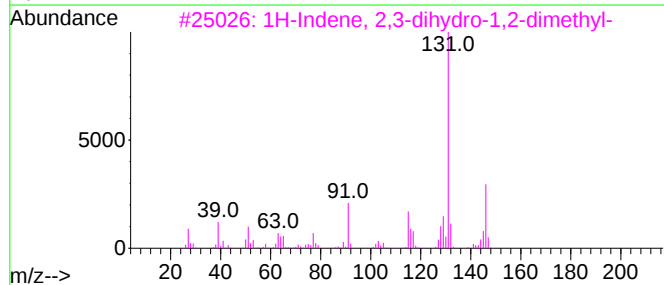
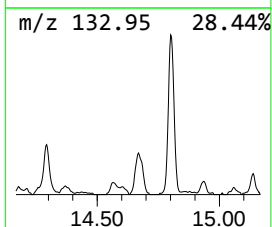
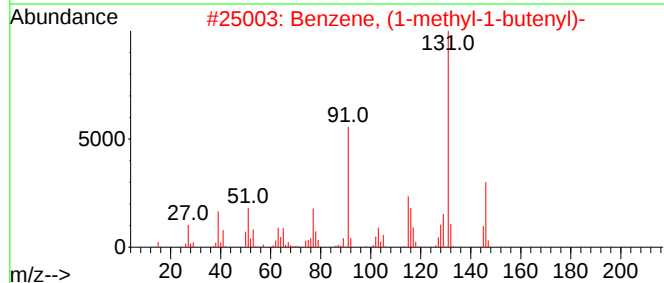
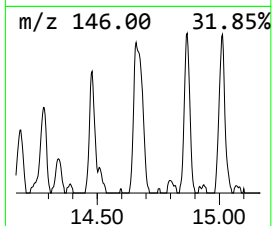
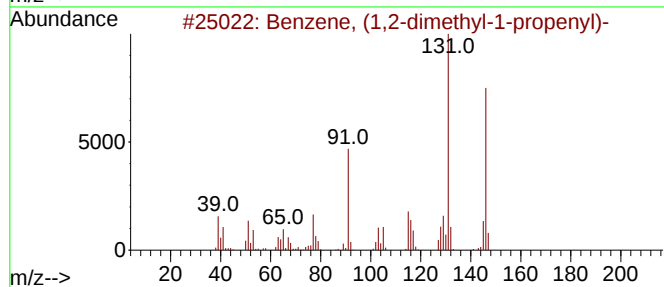
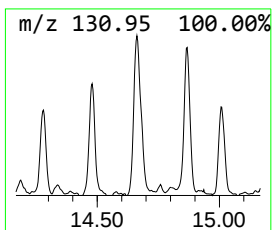
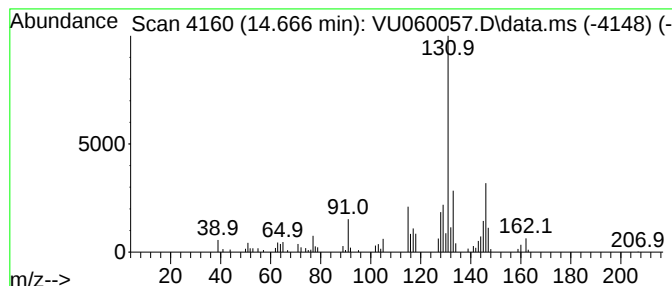
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.666	1.21 ug/L	116895	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
2			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	80
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

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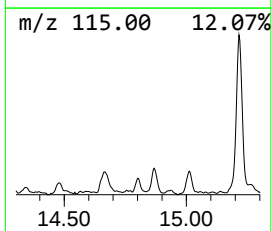
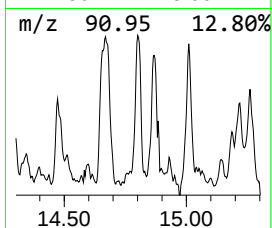
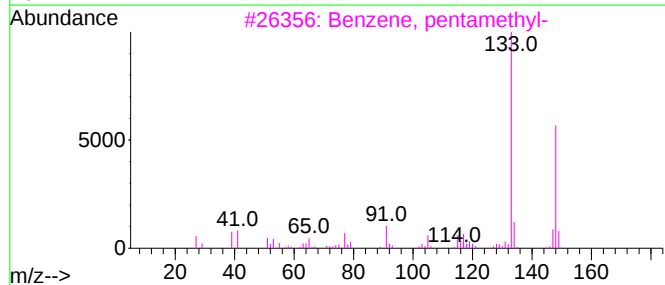
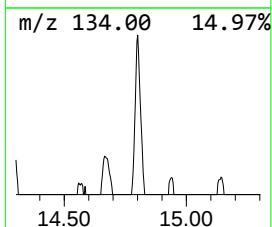
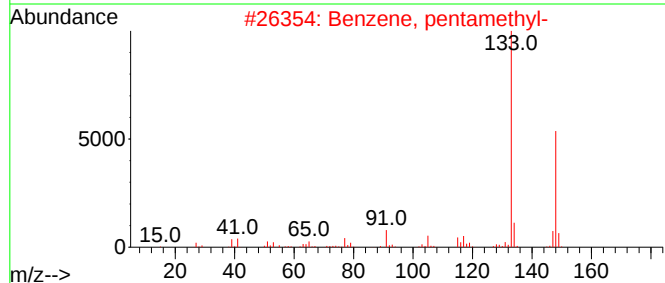
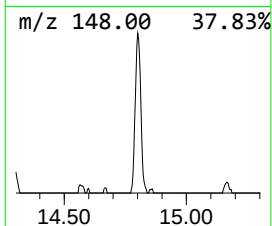
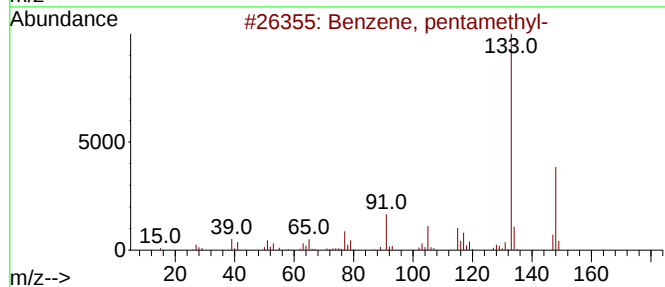
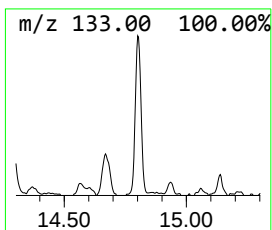
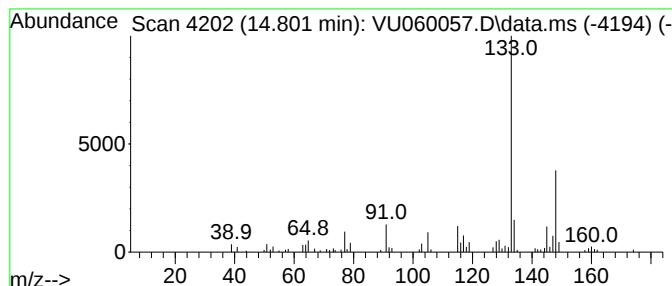
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, pentamethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.801	0.64 ug/L	62478	1,4-Dichlorobenzene-d4	11.807

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, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87
5			3,4-Dimethylcumene	148	C11H16	1000370-34-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

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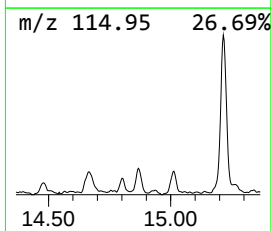
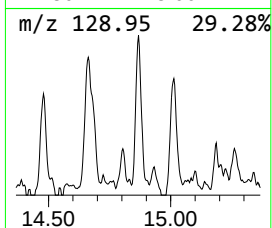
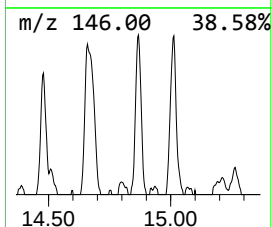
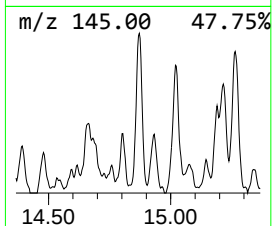
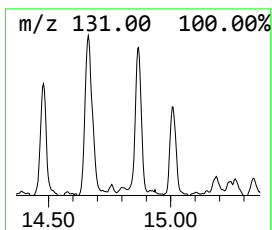
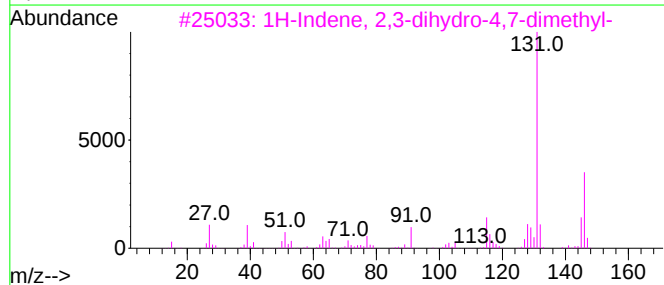
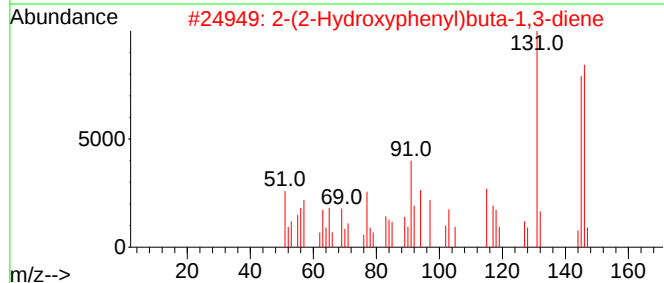
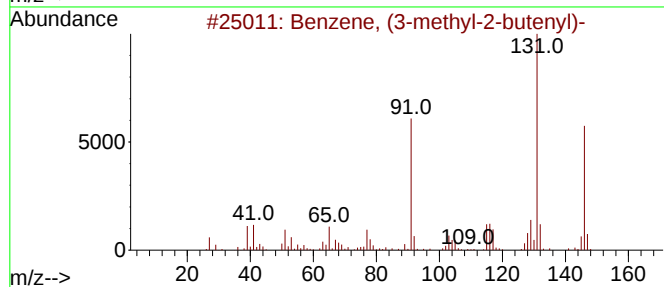
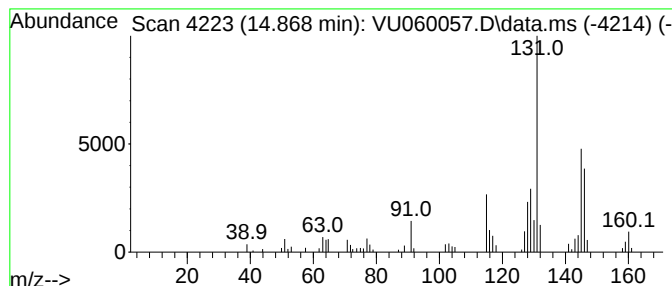
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Benzene, (3-methyl-2-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	0.84 ug/L	81074	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
2			2-(2-Hydroxyphenyl)buta-1,3-diene	146	C10H10O	038865-47-3	58
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	58
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	58
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

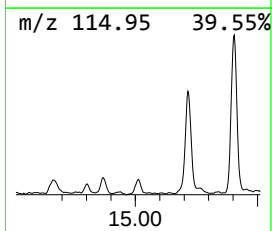
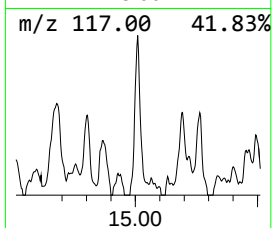
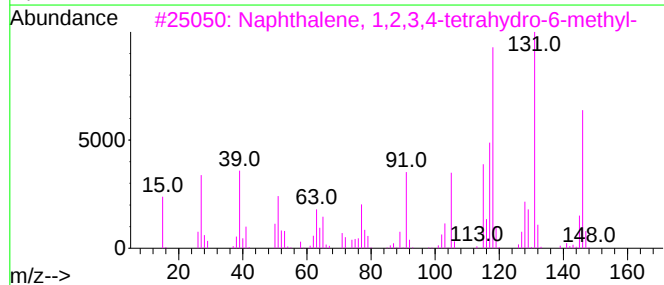
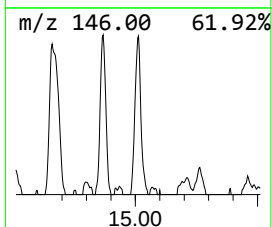
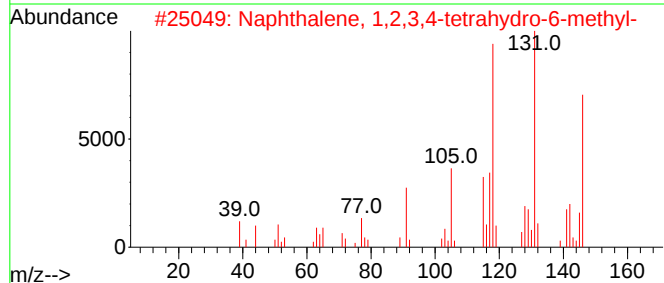
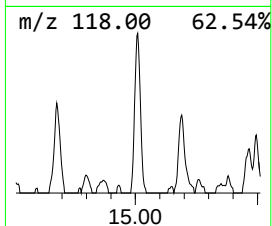
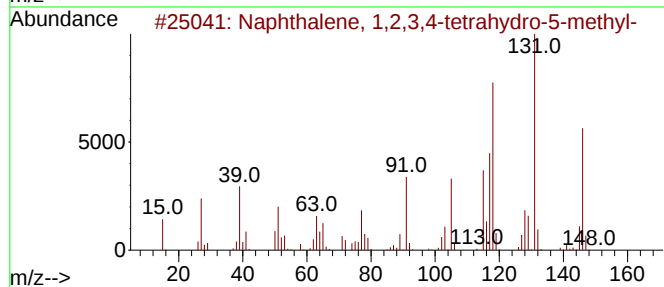
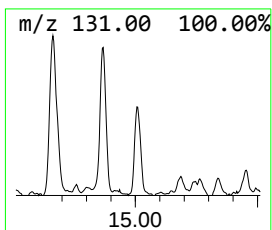
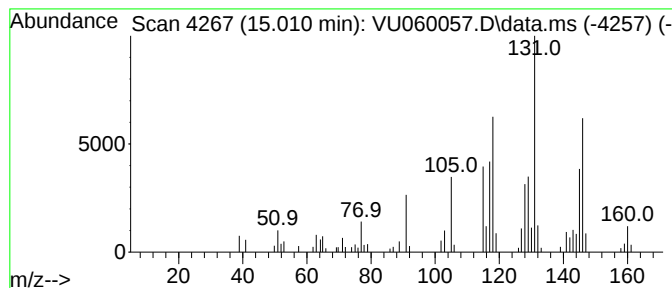
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.010	0.92 ug/L	88958	1,4-Dichlorobenzene-d4	11.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	89
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060057.D
Acq On : 19 Jul 2024 13:58
Operator : MD/SY
Sample : P3217-14DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZJ9DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.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	1.2	ug/L	70094	1	6.245	287839	5.0
Benzene, propyl-	10.900	0.7	ug/L	66920	3	11.807	484545	5.0
Benzene, 1-ethy...	10.994	0.7	ug/L	72143	3	11.807	484545	5.0
Benzene, 1-ethy...	11.286	1.4	ug/L	139311	3	11.807	484545	5.0
Azulene	14.084	0.9	ug/L	85302	3	11.807	484545	5.0
1H-Inden-1-one,...	14.280	0.8	ug/L	72869	3	11.807	484545	5.0
1H-Indene, 2,3-...	14.479	0.6	ug/L	60019	3	11.807	484545	5.0
Benzene, (1,2-d...	14.666	1.2	ug/L	116895	3	11.807	484545	5.0
Benzene, pentam...	14.801	0.6	ug/L	62478	3	11.807	484545	5.0
Benzene, (3-met...	14.868	0.8	ug/L	81074	3	11.807	484545	5.0
Naphthalene, 1,...	15.010	0.9	ug/L	88958	3	11.807	484545	5.0