

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
 Data File : VU063030.D
 Acq On : 27 Jan 2025 21:04
 Operator : MD/SY
 Sample : Q1174-26
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 E2953

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU063030.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.355	14	20	32	rBV2	16683	20494	6.26%	0.414%
2	1.483	55	60	66	rVB	7478	7233	2.21%	0.146%
3	1.586	82	92	106	rVB2	36563	45813	13.99%	0.926%
4	1.721	124	134	145	rVB	82110	102735	31.37%	2.078%
5	1.901	181	190	200	rBV	25451	32614	9.96%	0.660%
6	1.982	206	215	222	rVB3	8724	11451	3.50%	0.232%
7	2.162	264	271	278	rBV2	4688	7923	2.42%	0.160%
8	2.355	321	331	336	rBV	3942	5327	1.63%	0.108%
9	2.551	379	392	409	rBB	44308	70760	21.61%	1.431%
10	2.667	414	428	437	rBV5	3237	6354	1.94%	0.128%
11	2.773	452	461	473	rVB2	4949	8664	2.65%	0.175%
12	3.117	559	568	577	rBV4	2150	3817	1.17%	0.077%
13	4.638	1028	1041	1074	rBV3	20400	70017	21.38%	1.416%
14	5.046	1154	1168	1186	rBV	55814	122115	37.29%	2.469%
15	5.711	1356	1375	1410	rBV3	94043	252518	77.10%	5.106%
16	6.236	1527	1538	1557	rBV2	111905	214196	65.40%	4.332%
17	6.679	1664	1676	1694	rBV2	58176	113550	34.67%	2.296%
18	7.554	1938	1948	1964	rBV3	32801	59509	18.17%	1.203%
19	7.888	2038	2052	2065	rBV	138983	242549	74.06%	4.905%
20	7.959	2065	2074	2091	rVB2	60311	108750	33.21%	2.199%
21	8.171	2130	2140	2154	rBV2	23120	39784	12.15%	0.805%
22	8.464	2219	2231	2242	rBB2	2578	4448	1.36%	0.090%
23	8.625	2269	2281	2320	rBV	129736	269739	82.36%	5.455%
24	9.406	2511	2524	2547	rBV	182282	310115	94.69%	6.271%
25	9.686	2602	2611	2619	rBV2	4965	7466	2.28%	0.151%
26	10.094	2728	2738	2747	rBV3	6821	10550	3.22%	0.213%
27	10.473	2850	2856	2866	rVB3	2454	3746	1.14%	0.076%
28	10.747	2928	2941	2954	rBV	63003	106059	32.38%	2.145%
29	10.882	2972	2983	2995	rVB4	5792	10580	3.23%	0.214%
30	10.991	3007	3017	3033	rBV2	5372	8416	2.57%	0.170%
31	11.081	3033	3045	3059	rVB	14215	22522	6.88%	0.455%
32	11.284	3100	3108	3121	rBV	9125	14086	4.30%	0.285%
33	11.460	3152	3163	3178	rBV	32005	53676	16.39%	1.085%
34	11.737	3239	3249	3257	rBV5	7193	10168	3.10%	0.206%
35	11.801	3257	3269	3286	rVV	197864	325825	99.49%	6.589%
36	11.885	3286	3295	3305	rVB	30013	47440	14.49%	0.959%

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 Integrator: RTE
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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
 Title : TRACE VOA SFAM1.0

37	12.084	3346	3357	3364	rBV3	13071	22299	6.81%	0.451%
38	12.184	3376	3388	3412	rVB	199898	327501	100.00%	6.623%
39	12.303	3416	3425	3427	rBV5	3061	4516	1.38%	0.091%
40	12.367	3436	3445	3457	rVB4	5322	8232	2.51%	0.166%
41	12.457	3464	3473	3477	rBV3	14030	20481	6.25%	0.414%
42	12.486	3478	3482	3493	rVB2	15663	21976	6.71%	0.444%
43	12.563	3495	3506	3515	rBV2	24755	38061	11.62%	0.770%
44	12.676	3533	3541	3544	rBV3	8152	11616	3.55%	0.235%
45	12.869	3587	3601	3613	rVB	15129	25414	7.76%	0.514%
46	12.965	3616	3631	3638	rVV	37938	58986	18.01%	1.193%
47	13.017	3638	3647	3663	rVB	75297	118396	36.15%	2.394%
48	13.103	3663	3674	3677	rBV6	4916	7854	2.40%	0.159%
49	13.126	3678	3681	3686	rVV4	5288	5485	1.67%	0.111%
50	13.158	3687	3691	3699	rVB3	4572	5215	1.59%	0.105%
51	13.284	3712	3730	3742	rBV3	27818	46848	14.30%	0.947%
52	13.338	3742	3747	3756	rVB9	3109	4770	1.46%	0.096%
53	13.396	3756	3765	3769	rBV4	7647	11389	3.48%	0.230%
54	13.441	3769	3779	3794	rVV4	107808	192496	58.78%	3.893%
55	13.512	3794	3801	3807	rVV5	8042	11375	3.47%	0.230%
56	13.550	3808	3813	3819	rVB2	4281	4846	1.48%	0.098%
57	13.615	3824	3833	3849	rVB3	20701	37217	11.36%	0.753%
58	13.721	3855	3866	3875	rBV6	6919	12664	3.87%	0.256%
59	13.775	3875	3883	3889	rVV6	14286	21569	6.59%	0.436%
60	13.824	3889	3898	3916	rVV5	22018	59744	18.24%	1.208%
61	13.901	3916	3922	3925	rVV5	4801	6374	1.95%	0.129%
62	13.952	3925	3938	3947	rVB4	14116	30088	9.19%	0.608%
63	14.078	3964	3977	3997	rVB	128076	227934	69.60%	4.609%
64	14.181	4002	4009	4022	rVB	4872	9234	2.82%	0.187%
65	14.283	4030	4041	4051	rBV8	12398	24056	7.35%	0.486%
66	14.335	4051	4057	4066	rVB5	7058	10058	3.07%	0.203%
67	14.476	4090	4101	4118	rVB3	26941	44316	13.53%	0.896%
68	14.566	4121	4129	4133	rBV7	2571	4084	1.25%	0.083%
69	14.608	4133	4142	4149	rVV6	8287	13817	4.22%	0.279%
70	14.660	4149	4158	4171	rVV6	26898	61286	18.71%	1.239%
71	14.721	4171	4177	4185	rVB2	11022	13872	4.24%	0.281%
72	14.801	4191	4202	4214	rBV2	31030	52599	16.06%	1.064%
73	14.865	4214	4222	4232	rVB3	22015	34605	10.57%	0.700%
74	15.007	4257	4266	4276	rBV4	16336	28239	8.62%	0.571%
75	15.148	4297	4310	4317	rBV10	2873	7358	2.25%	0.149%
76	15.216	4317	4331	4356	rVB	87410	157750	48.17%	3.190%

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Start Thrs: 0.1

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
 Title : TRACE VOA SFAM1.0

77	15.335	4358	4368	4376	rBV8	3774	5601	1.71%	0.113%
78	15.399	4376	4388	4400	rBV	117470	199976	61.06%	4.044%
79	15.560	4431	4438	4451	rVB10	5813	12328	3.76%	0.249%
80	15.679	4462	4475	4493	rVB10	7967	21647	6.61%	0.438%
81	15.817	4513	4518	4525	rBV8	2967	4374	1.34%	0.088%
82	15.869	4527	4534	4540	rVV8	2443	4360	1.33%	0.088%
83	15.914	4540	4548	4566	rVB8	11871	24988	7.63%	0.505%
84	16.004	4568	4576	4580	rBV7	2590	3772	1.15%	0.076%
85	16.068	4589	4596	4604	rBV10	3068	3884	1.19%	0.079%
86	16.139	4611	4618	4625	rBV6	7003	10495	3.20%	0.212%
87	16.171	4625	4628	4637	rVB7	4920	6528	1.99%	0.132%
88	16.258	4645	4655	4667	rBV5	11100	20681	6.31%	0.418%
89	16.402	4690	4700	4707	rBV2	25778	41734	12.74%	0.844%
90	16.441	4709	4712	4724	rVB4	11062	13846	4.23%	0.280%
91	16.528	4735	4739	4748	rVB7	2348	3471	1.06%	0.070%
92	16.608	4755	4764	4767	rBV7	3787	4951	1.51%	0.100%
93	16.637	4768	4773	4781	rVB7	4994	6806	2.08%	0.138%

Sum of corrected areas: 4945071

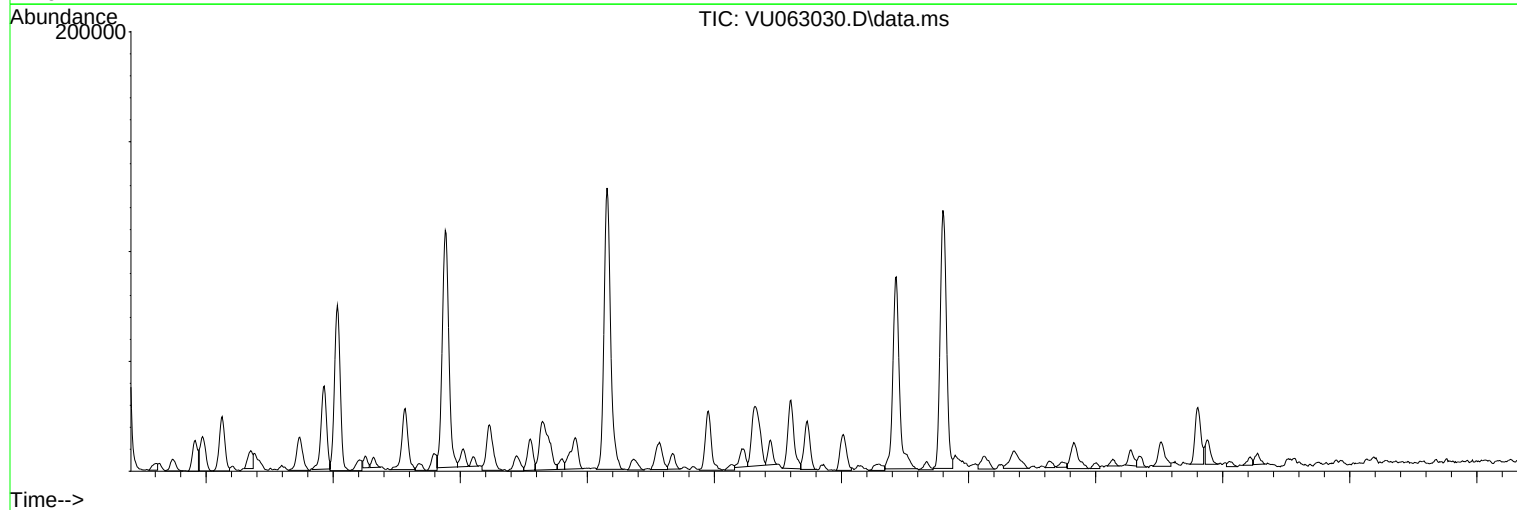
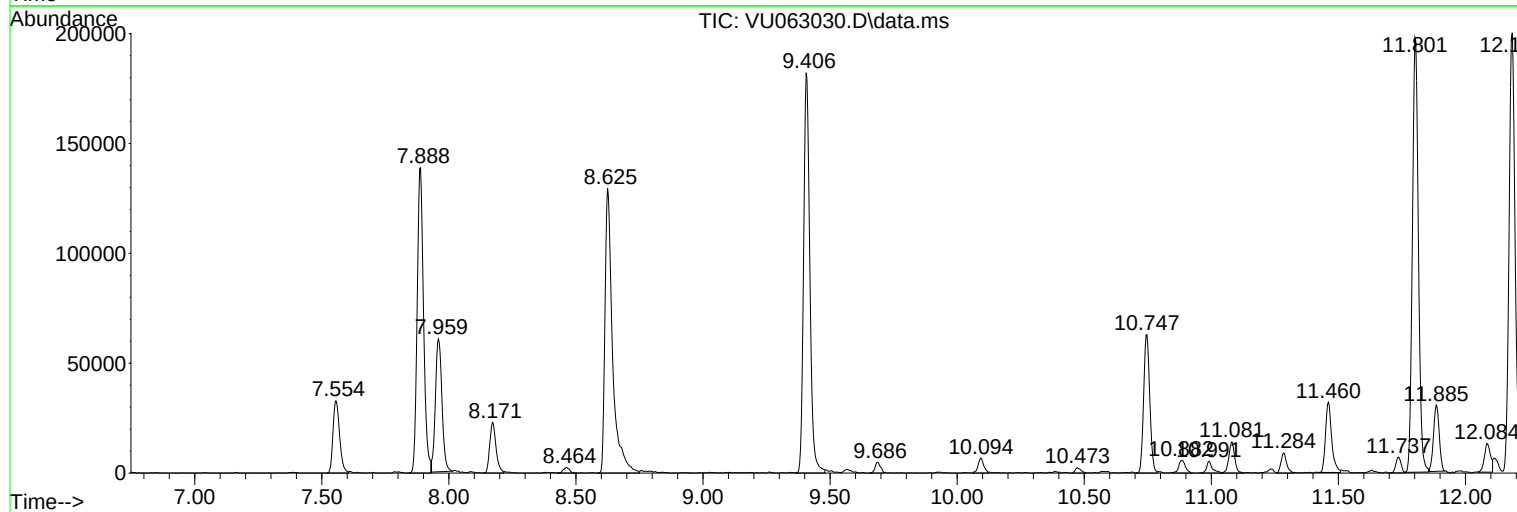
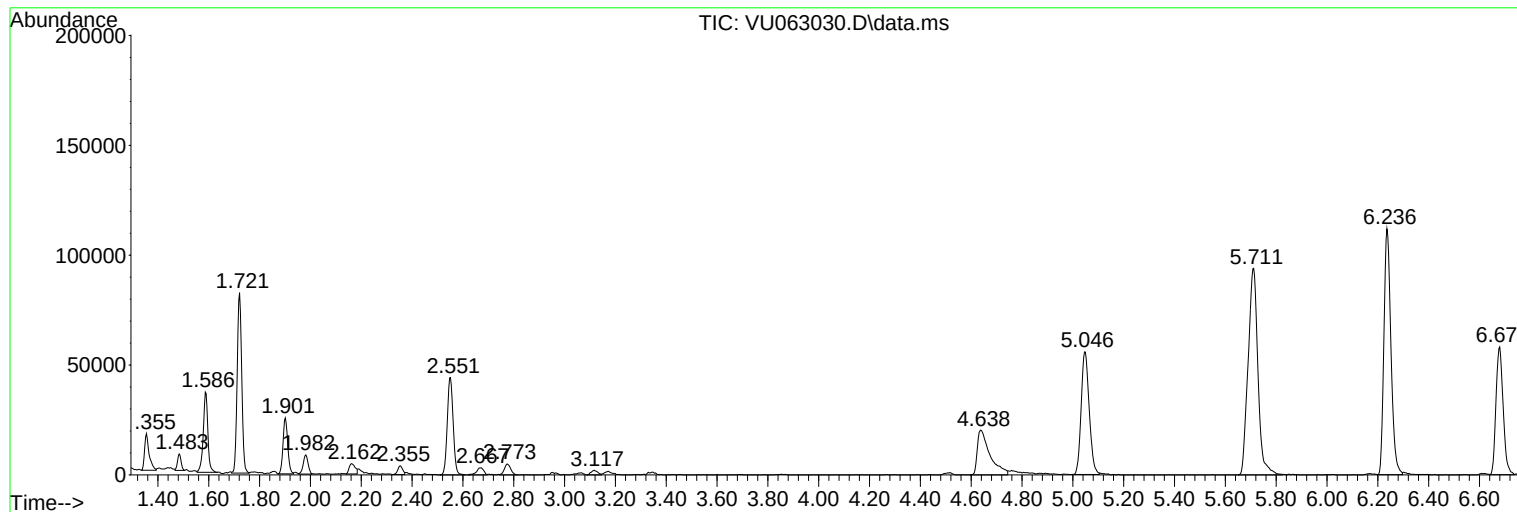
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

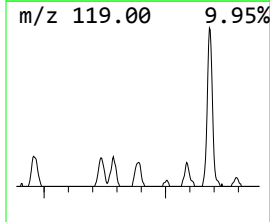
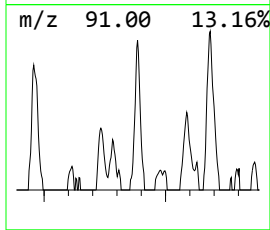
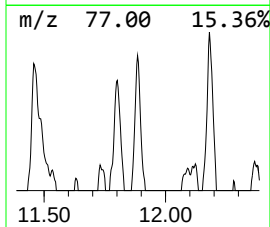
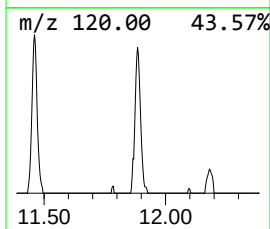
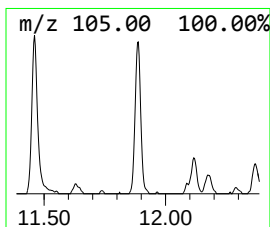
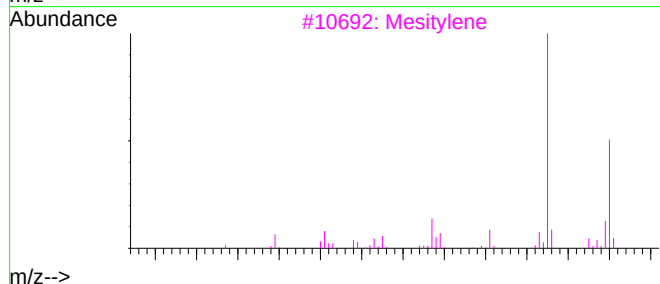
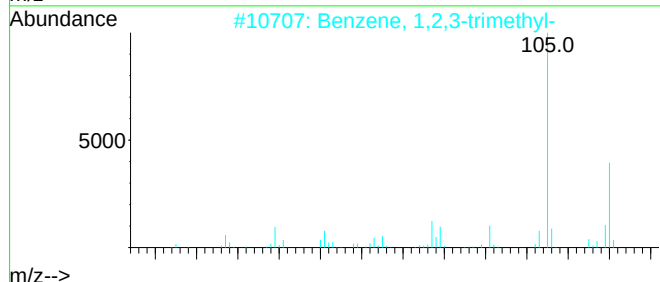
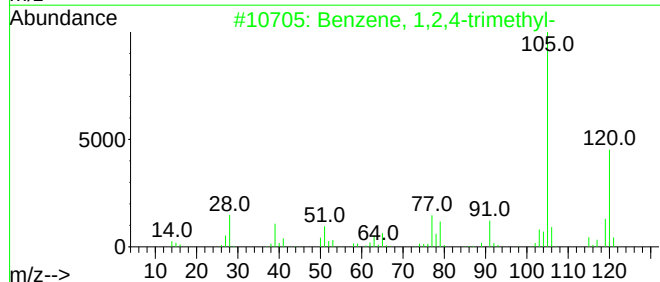
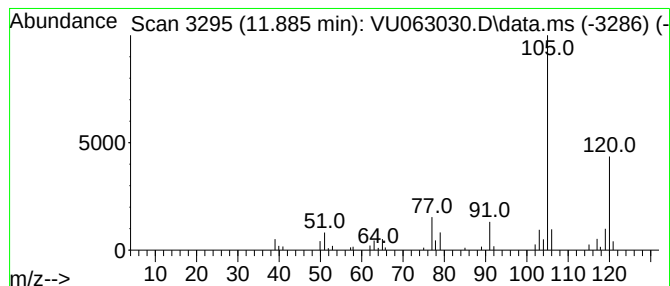
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.885	0.73 ug/L	47440	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Mesitylene	120	C9H12	000108-67-8	91



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ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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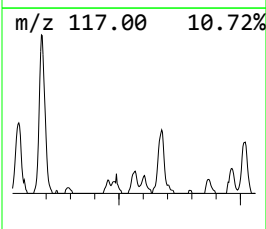
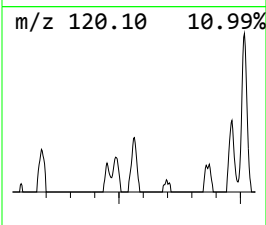
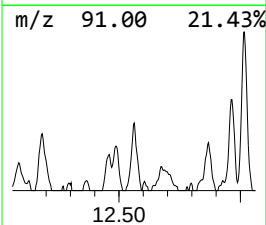
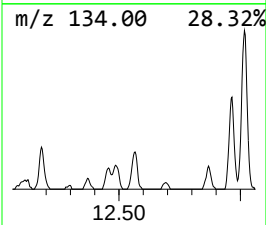
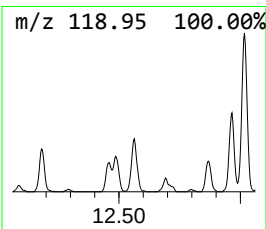
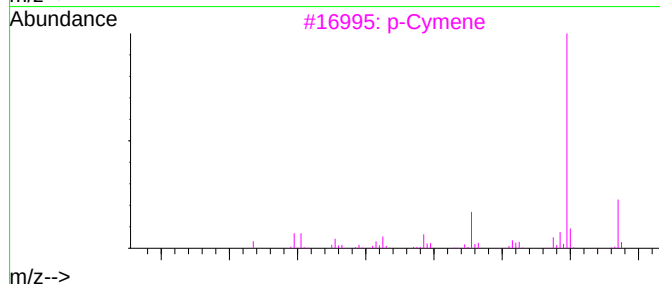
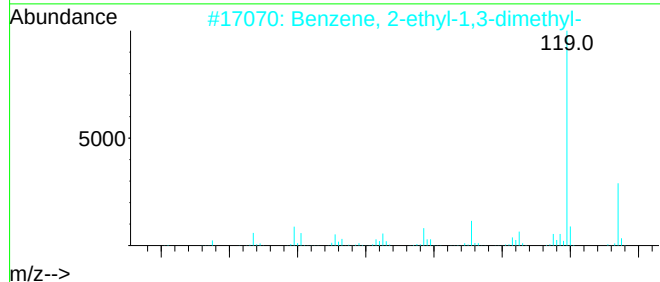
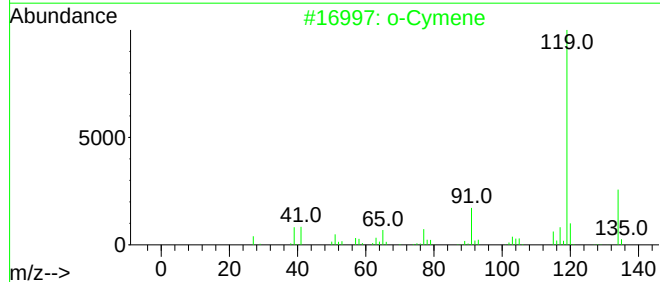
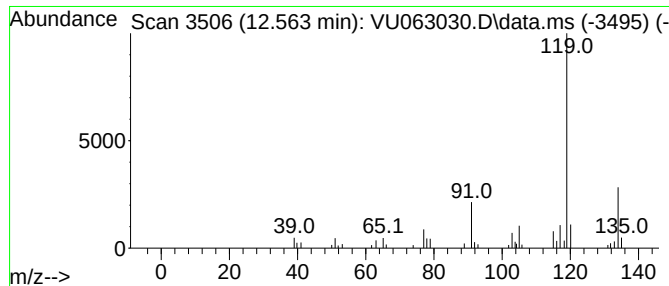
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 o-Cymene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.563	0.58 ug/L	38061	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
3	p-Cymene	134	C10H14	000099-87-6	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	o-Cymene	134	C10H14	000527-84-4	93



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Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

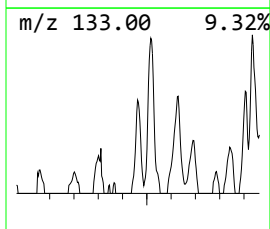
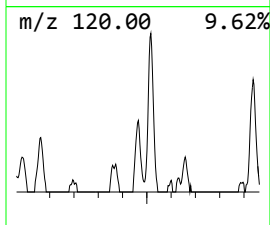
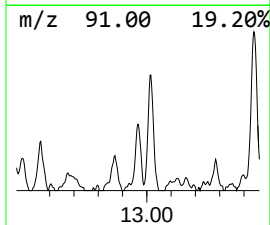
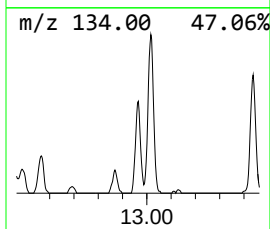
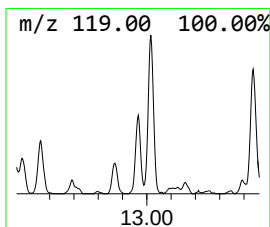
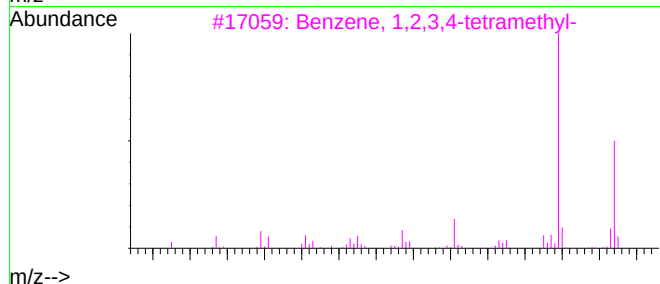
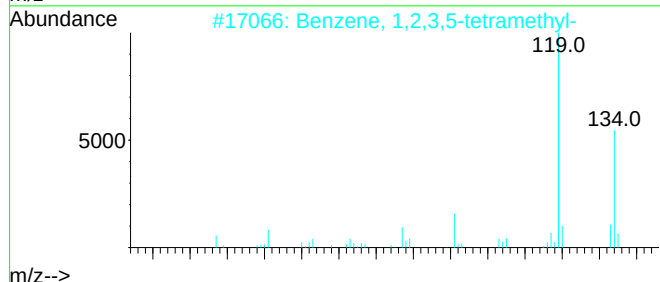
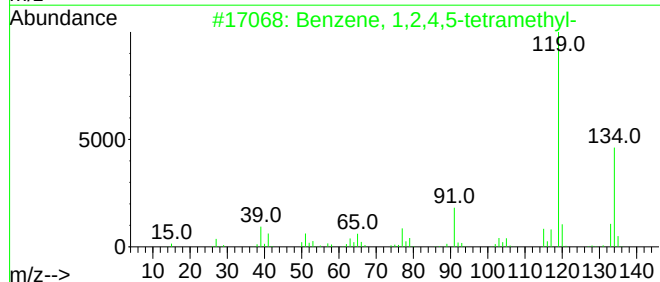
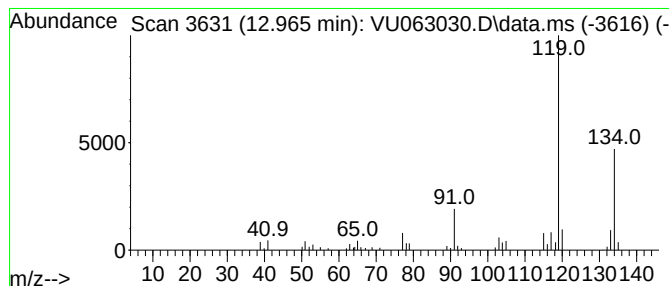
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.965	0.91 ug/L	58986	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	93
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

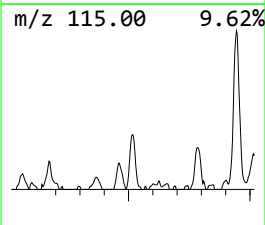
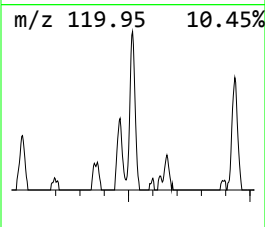
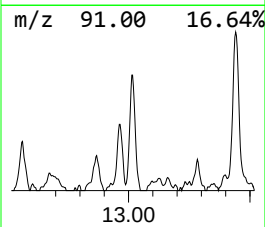
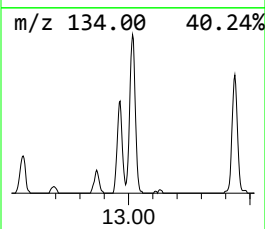
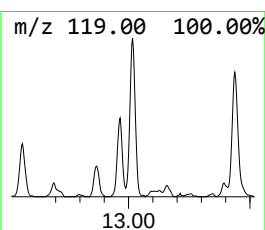
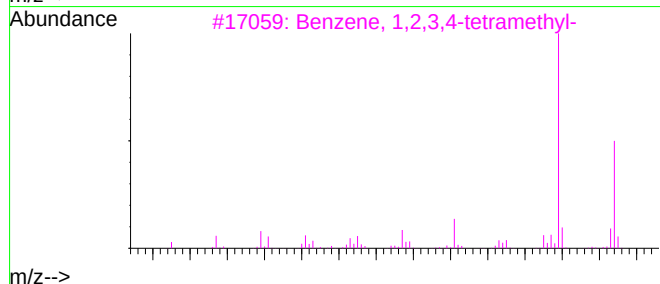
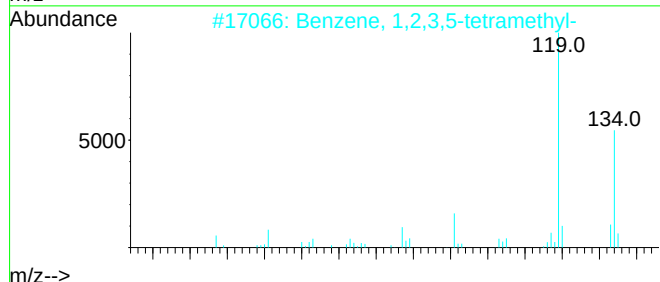
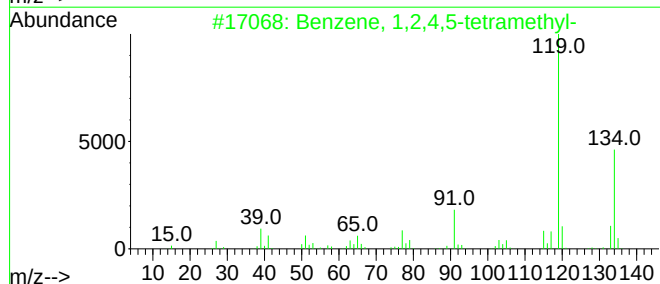
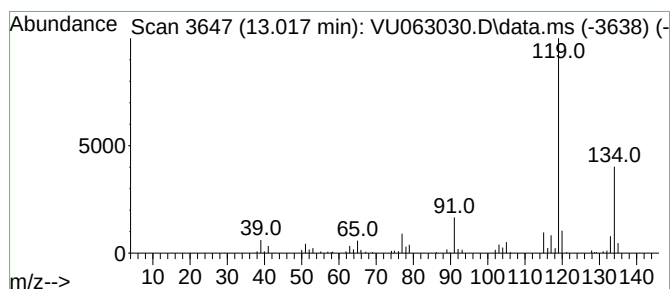
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.017	1.82 ug/L	118396	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

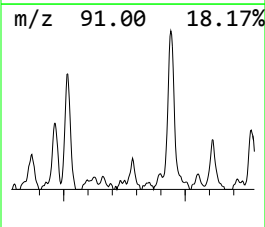
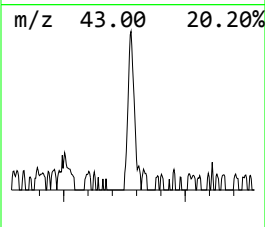
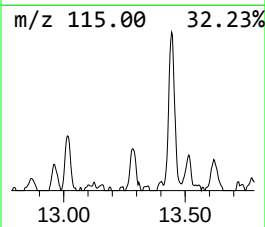
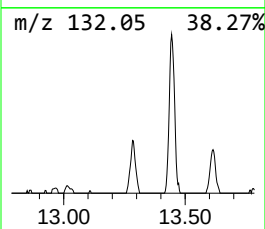
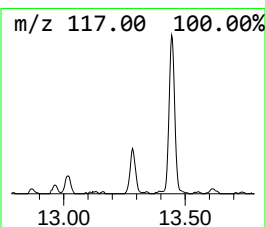
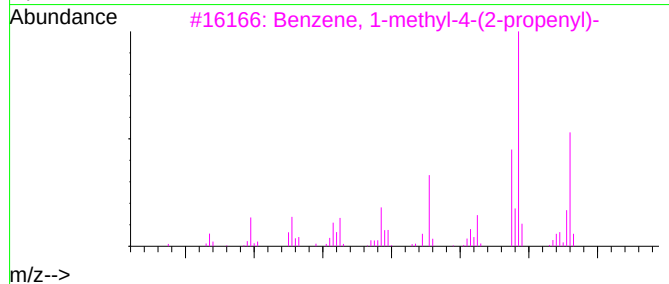
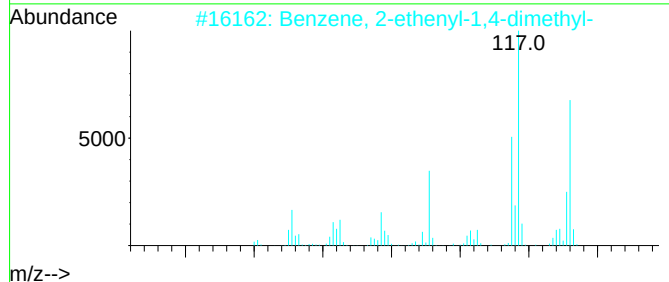
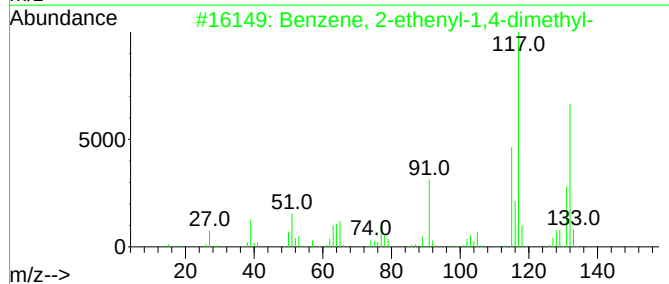
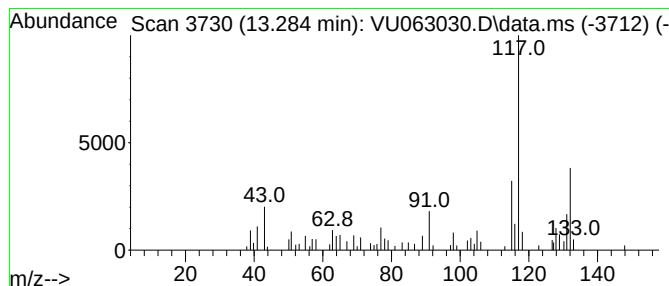
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.284	0.72 ug/L	46848	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	92
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

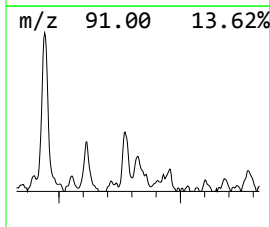
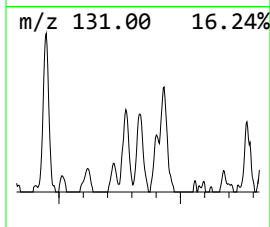
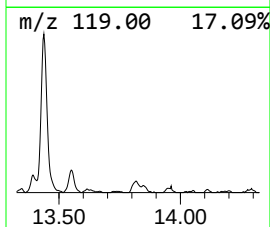
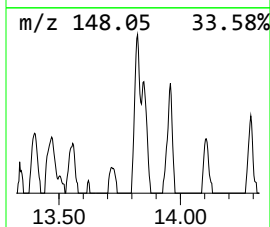
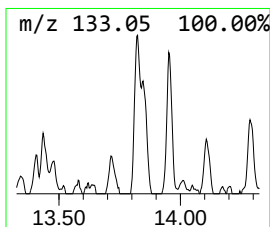
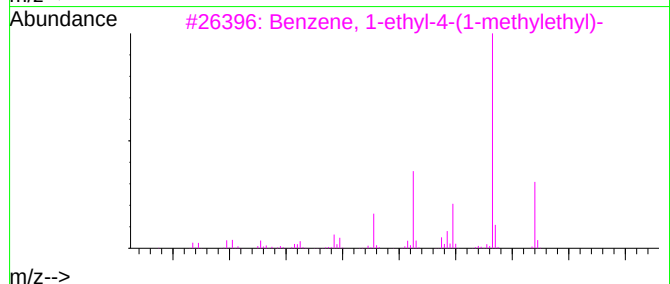
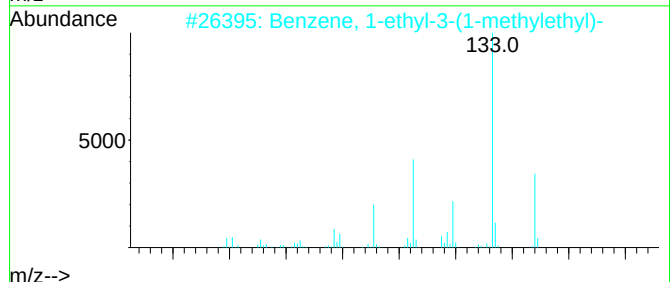
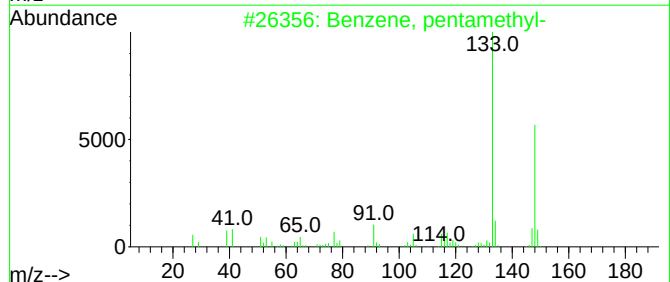
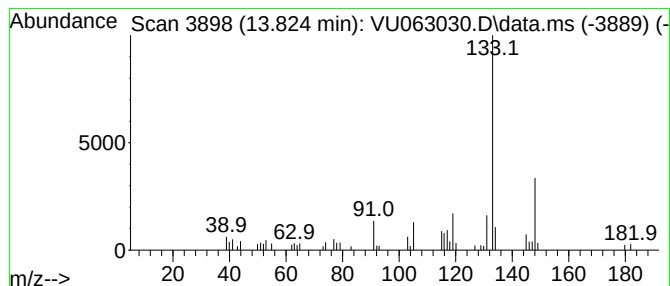
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, pentamethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.824	0.92 ug/L	59744	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	87
2	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	87
3	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	87
4	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	83
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

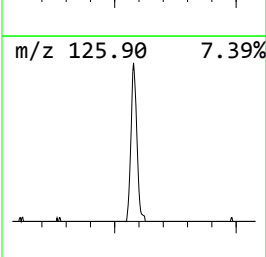
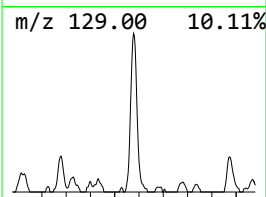
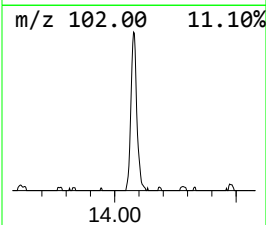
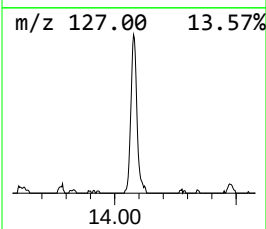
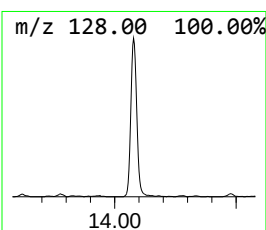
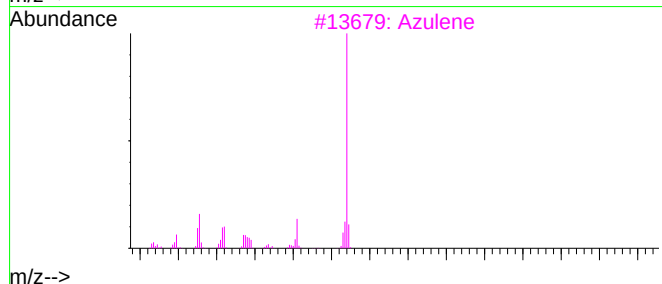
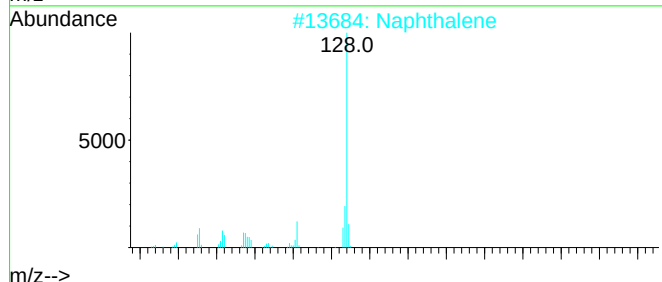
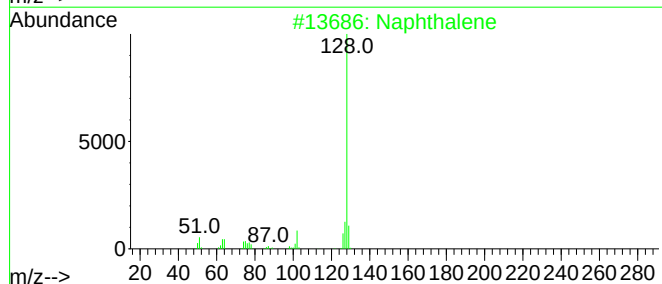
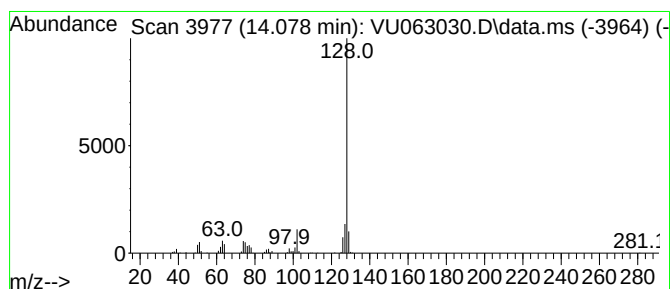
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 11 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.078	3.50 ug/L	227934	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	95
3	Azulene	128	C10H8	000275-51-4	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\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

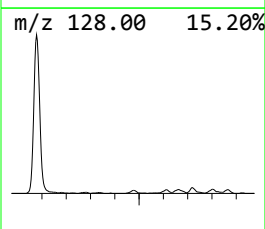
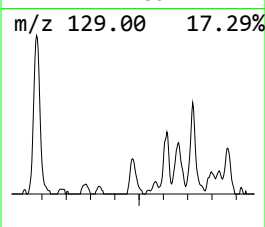
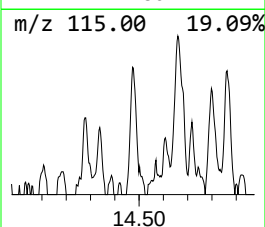
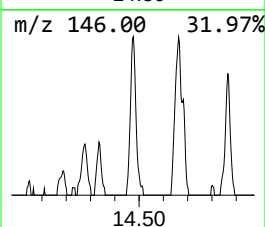
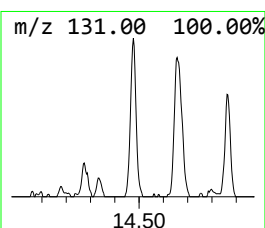
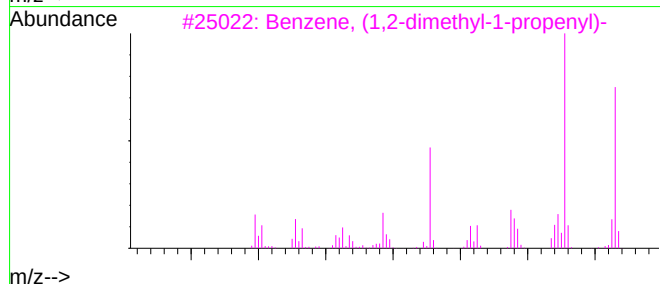
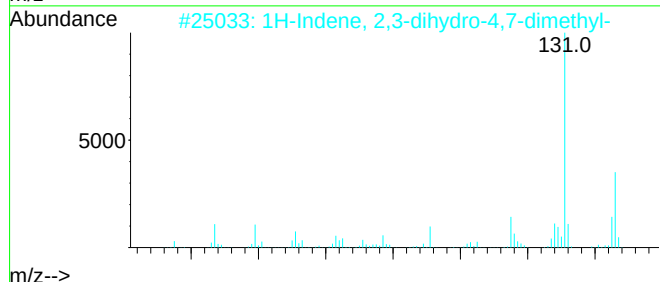
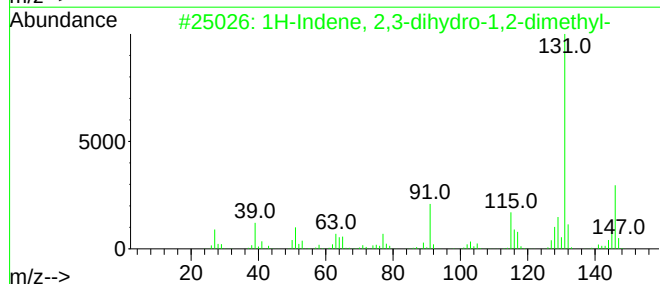
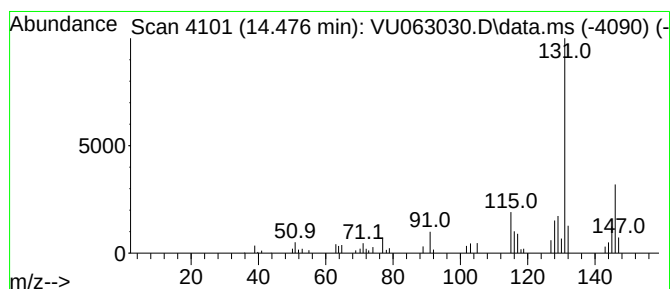
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 12 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.476	0.68 ug/L	44316	1,4-Dichlorobenzene-d4	11.805	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	95
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

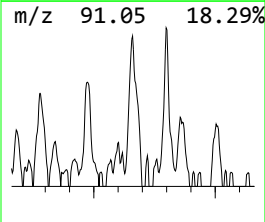
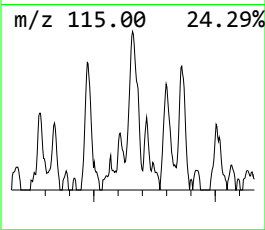
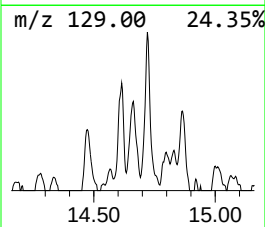
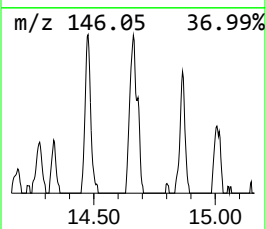
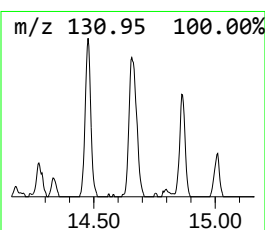
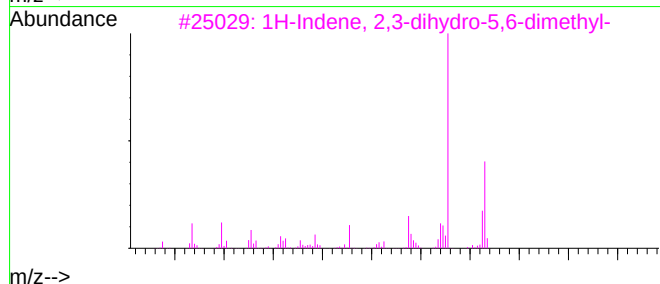
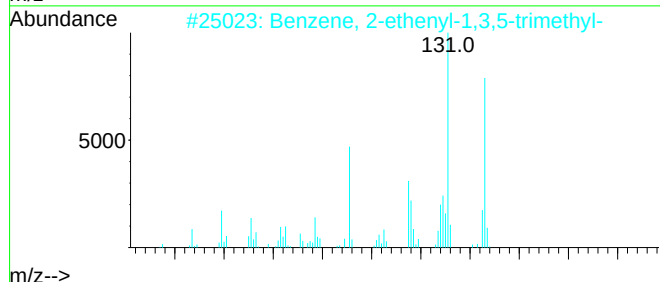
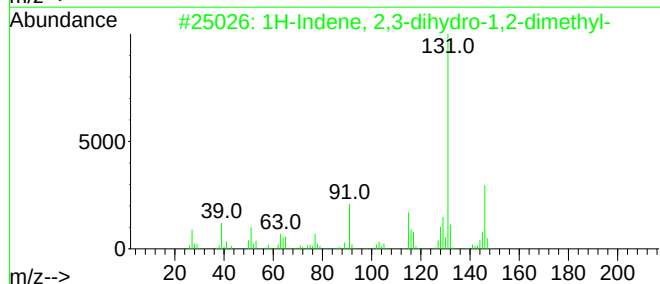
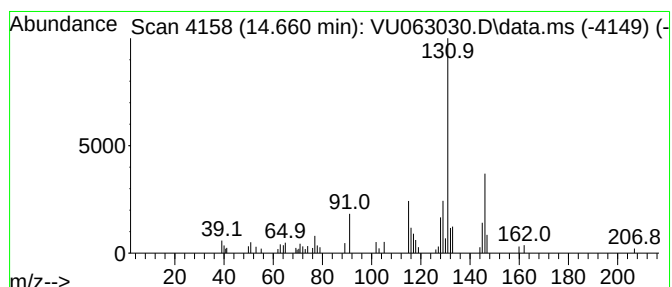
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.660	0.94 ug/L	61286	1,4-Dichlorobenzene-d4	11.805	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
2	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	87
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	87
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	83
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.M
Quant Title : TRACE VOA SFAM1.0

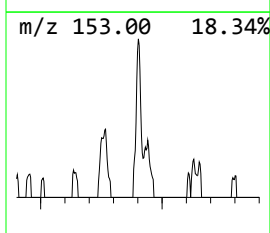
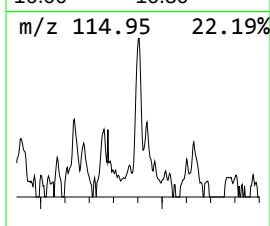
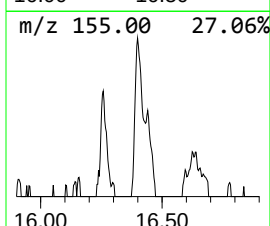
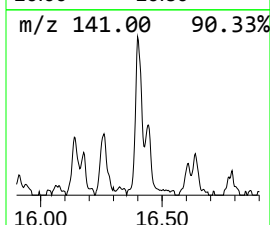
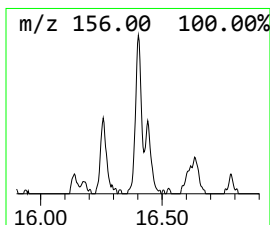
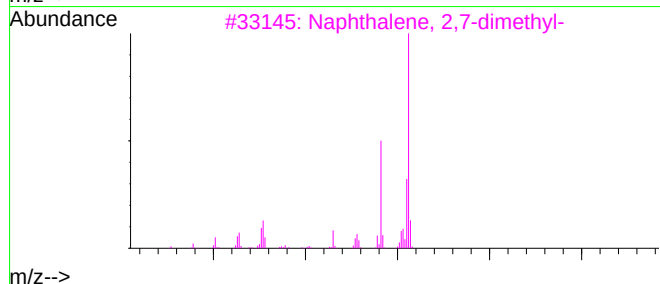
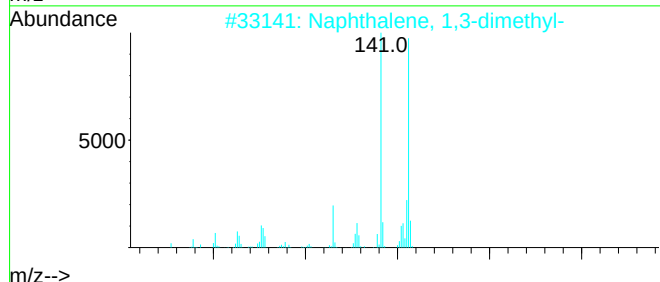
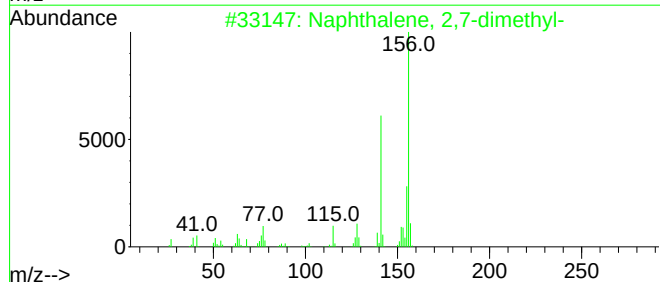
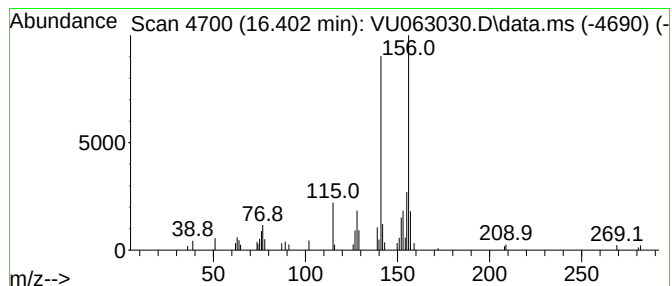
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 18 Naphthalene, 2,7-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.402	0.64 ug/L	41734	1,4-Dichlorobenzene-d4	11.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012725\
Data File : VU063030.D
Acq On : 27 Jan 2025 21:04
Operator : MD/SY
Sample : Q1174-26
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2953

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR012725WMA.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, 1,2,3-...	11.885	0.7	ug/L	47440	3	11.805	325825	5.0
o-Cymene	12.563	0.6	ug/L	38061	3	11.805	325825	5.0
Benzene, 1,2,4,...	12.965	0.9	ug/L	58986	3	11.805	325825	5.0
Benzene, 1,2,3,...	13.017	1.8	ug/L	118396	3	11.805	325825	5.0
Benzene, 2-ethe...	13.284	0.7	ug/L	46848	3	11.805	325825	5.0
Benzene, pentam...	13.824	0.9	ug/L	59744	3	11.805	325825	5.0
Azulene	14.078	3.5	ug/L	227934	3	11.805	325825	5.0
1H-Indene, 2,3-...	14.476	0.7	ug/L	44316	3	11.805	325825	5.0
Benzene, 2-ethe...	14.660	0.9	ug/L	61286	3	11.805	325825	5.0
Naphthalene, 2,...	16.402	0.6	ug/L	41734	3	11.805	325825	5.0