

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
 Data File : VV026683.D
 Acq On : 07 Jul 2022 17:42
 Operator : SY/MD
 Sample : N3601-10
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBUB5

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_V\Method\SFAMVTR062322WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VV026683.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.137	23	30	38	rVB2	13821	20791	1.05%	0.104%
2	1.320	78	87	91	rBV	88613	96944	4.91%	0.486%
3	1.368	91	102	129	rVV	203389	907536	45.99%	4.545%
4	1.478	131	136	161	rVV2	178064	483807	24.52%	2.423%
5	1.581	163	168	185	rVB	96474	155900	7.90%	0.781%
6	2.044	307	312	325	rVB2	29435	43084	2.18%	0.216%
7	2.118	327	335	354	rVB2	165915	266756	13.52%	1.336%
8	2.220	357	367	383	rVB	63051	96441	4.89%	0.483%
9	2.304	383	393	407	rBV	101075	148306	7.52%	0.743%
10	3.928	886	898	937	rBV2	46969	170718	8.65%	0.855%
11	4.362	1017	1033	1058	rBV	99705	250530	12.70%	1.255%
12	5.056	1234	1249	1256	rBV2	213445	472941	23.97%	2.369%
13	5.108	1257	1265	1299	rVB	339053	755704	38.30%	3.785%
14	5.375	1340	1348	1365	rVB	14855	32176	1.63%	0.161%
15	5.625	1414	1426	1459	rBV	203935	424109	21.49%	2.124%
16	5.921	1507	1518	1548	rVB	95985	201097	10.19%	1.007%
17	6.076	1554	1566	1577	rBV	119143	241624	12.25%	1.210%
18	6.133	1579	1584	1600	rVB2	37457	71814	3.64%	0.360%
19	6.995	1834	1852	1872	rBV	34663	68584	3.48%	0.343%
20	7.320	1942	1953	1969	rBV	227107	396477	20.09%	1.986%
21	7.397	1969	1977	2001	rVB	24583	51313	2.60%	0.257%
22	7.526	2006	2017	2035	rBV	179984	320647	16.25%	1.606%
23	7.632	2042	2050	2067	rVB	18519	35765	1.81%	0.179%
24	7.979	2134	2158	2177	rBV2	123180	216989	11.00%	1.087%
25	8.095	2184	2194	2223	rBV	169275	355201	18.00%	1.779%
26	8.879	2417	2438	2470	rBV2	901827	1973156	100.00%	9.882%
27	9.014	2470	2480	2499	rVV	78565	130581	6.62%	0.654%
28	9.143	2501	2520	2542	rVV3	144228	329722	16.71%	1.651%
29	9.545	2637	2645	2664	rVB	81305	132680	6.72%	0.664%
30	9.931	2752	2765	2777	rBV	37670	62916	3.19%	0.315%
31	10.217	2835	2854	2869	rVB	87062	160381	8.13%	0.803%
32	10.355	2887	2897	2915	rBV2	63291	122890	6.23%	0.615%
33	10.448	2916	2926	2944	rVV	132120	260177	13.19%	1.303%
34	10.542	2944	2955	2966	rVV2	227582	358165	18.15%	1.794%
35	10.606	2967	2975	2984	rVB3	43206	66228	3.36%	0.332%
36	10.738	3006	3016	3035	rBV	126122	212516	10.77%	1.064%

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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_V\Method\SFAMVTR062322WMA.M
 Title : TRACE VOA SFAM1.0

37	10.915	3060	3071	3088	rVB	326996	498140	25.25%	2.495%
38	11.085	3107	3124	3139	rBV4	107433	206167	10.45%	1.033%
39	11.162	3139	3148	3150	rBV4	19087	24441	1.24%	0.122%
40	11.191	3151	3157	3165	rVB	49609	68271	3.46%	0.342%
41	11.249	3165	3175	3179	rBV2	464917	748841	37.95%	3.750%
42	11.336	3193	3202	3225	rVB	300316	586942	29.75%	2.940%
43	11.452	3229	3238	3249	rVB2	22325	37105	1.88%	0.186%
44	11.532	3249	3263	3271	rBV5	109633	240840	12.21%	1.206%
45	11.571	3272	3275	3283	rVB	60437	74162	3.76%	0.371%
46	11.632	3283	3294	3316	rVB4	368456	883060	44.75%	4.423%
47	11.744	3316	3329	3343	rBV3	33409	58259	2.95%	0.292%
48	11.818	3343	3352	3368	rBV	59060	94001	4.76%	0.471%
49	11.911	3371	3381	3385	rBV	86482	137199	6.95%	0.687%
50	11.940	3385	3390	3400	rVB	109254	151186	7.66%	0.757%
51	12.014	3404	3413	3419	rBV	148430	220608	11.18%	1.105%
52	12.114	3431	3444	3466	rVB2	119257	295196	14.96%	1.478%
53	12.255	3476	3488	3493	rBV6	16000	33637	1.70%	0.168%
54	12.313	3495	3506	3517	rVB3	85987	166778	8.45%	0.835%
55	12.410	3531	3536	3544	rVB	152653	225170	11.41%	1.128%
56	12.464	3544	3553	3571	rVB	247321	378461	19.18%	1.895%
57	12.551	3572	3580	3588	rBV2	32085	48412	2.45%	0.242%
58	12.725	3626	3634	3651	rVB2	113858	208839	10.58%	1.046%
59	12.879	3666	3682	3693	rBV2	538892	841513	42.65%	4.214%
60	12.950	3698	3704	3714	rVB5	51627	83129	4.21%	0.416%
61	13.043	3724	3733	3753	rVB2	146085	252010	12.77%	1.262%
62	13.159	3758	3769	3779	rBV4	86059	152547	7.73%	0.764%
63	13.214	3779	3786	3794	rBV	36069	48877	2.48%	0.245%
64	13.268	3794	3803	3811	rBV4	81041	139835	7.09%	0.700%
65	13.336	3818	3824	3828	rBV2	34044	50917	2.58%	0.255%
66	13.365	3828	3833	3856	rVB2	98391	191603	9.71%	0.960%
67	13.500	3861	3875	3884	rBV	466322	736010	37.30%	3.686%
68	13.545	3885	3889	3899	rVB	46278	64525	3.27%	0.323%
69	13.616	3903	3911	3921	rVB3	22516	44259	2.24%	0.222%
70	13.712	3922	3941	3951	rBV6	33417	94717	4.80%	0.474%
71	13.767	3952	3958	3969	rVB2	29975	44973	2.28%	0.225%
72	13.915	3992	4004	4005	rBV3	33918	48099	2.44%	0.241%
73	14.088	4049	4058	4074	rBV5	53479	138911	7.04%	0.696%
74	14.230	4096	4102	4113	rVV3	17948	31152	1.58%	0.156%
75	14.291	4114	4121	4135	rVB3	35763	64454	3.27%	0.323%
76	14.429	4154	4164	4177	rBV3	42161	77641	3.93%	0.389%

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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_V\Method\SFAMVTR062322WMA.M
Title : TRACE VOA SFAM1.0

77	14.574	4201	4209	4215	rBV3	19474	29305	1.49%	0.147%
78	14.628	4216	4226	4238	rBV	457029	731432	37.07%	3.663%
79	14.812	4273	4283	4297	rVB	352307	570958	28.94%	2.859%
80	15.808	4582	4593	4600	rBV3	15049	26769	1.36%	0.134%
81	16.178	4701	4708	4720	rVB9	14481	23235	1.18%	0.116%

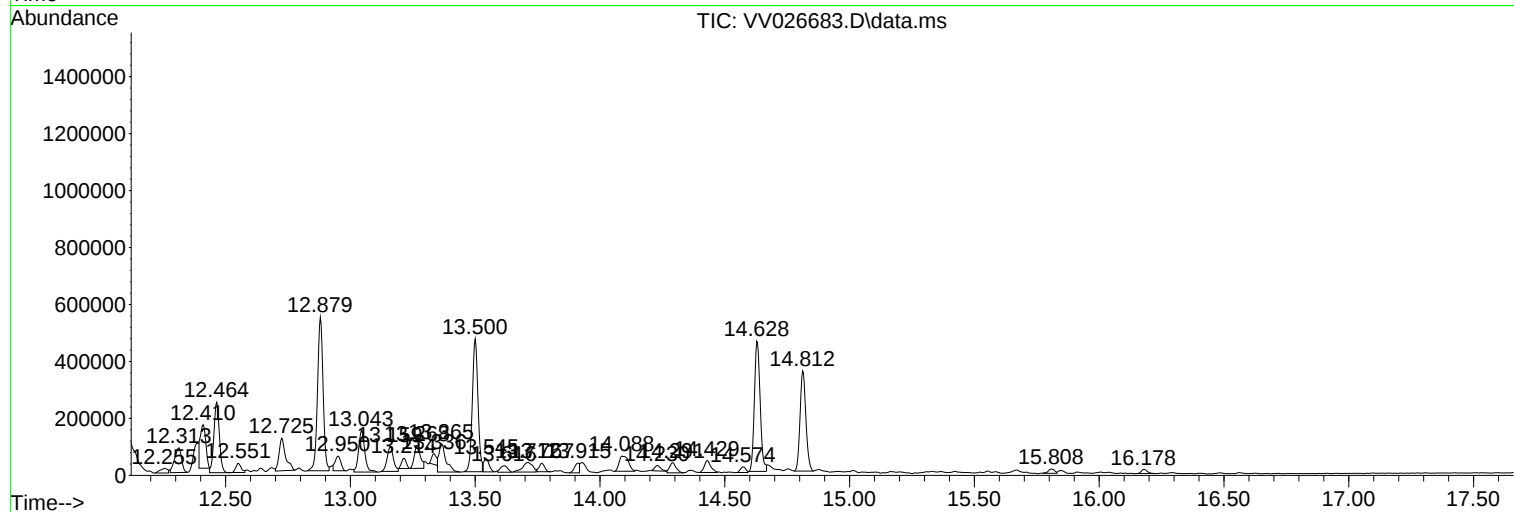
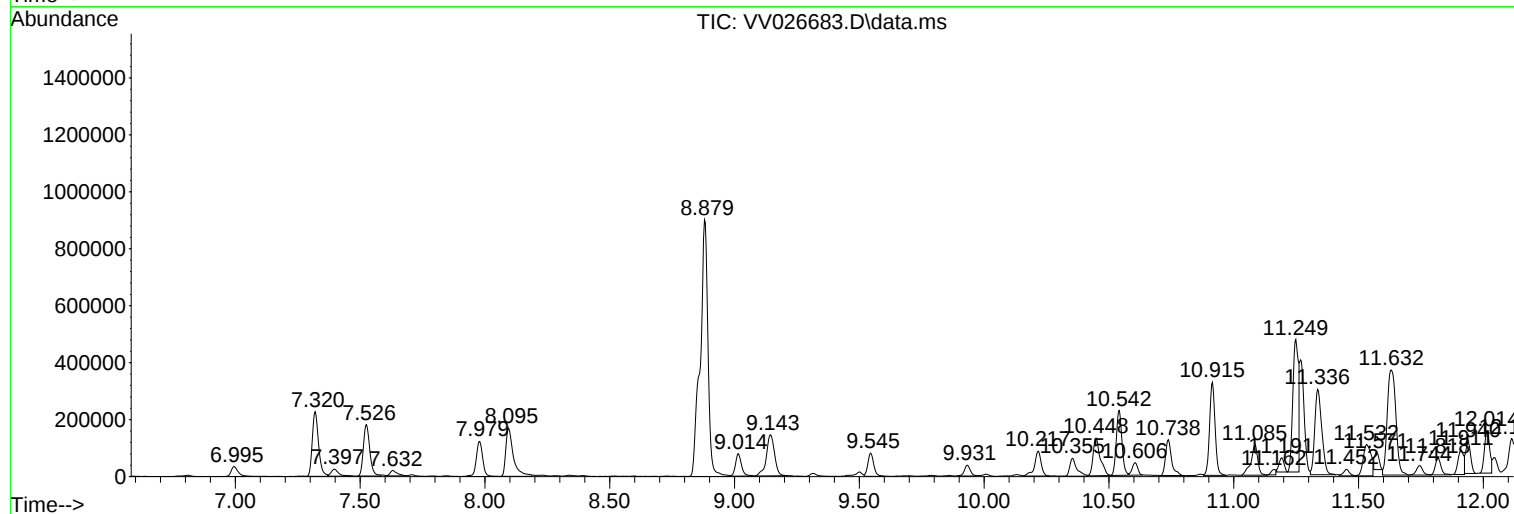
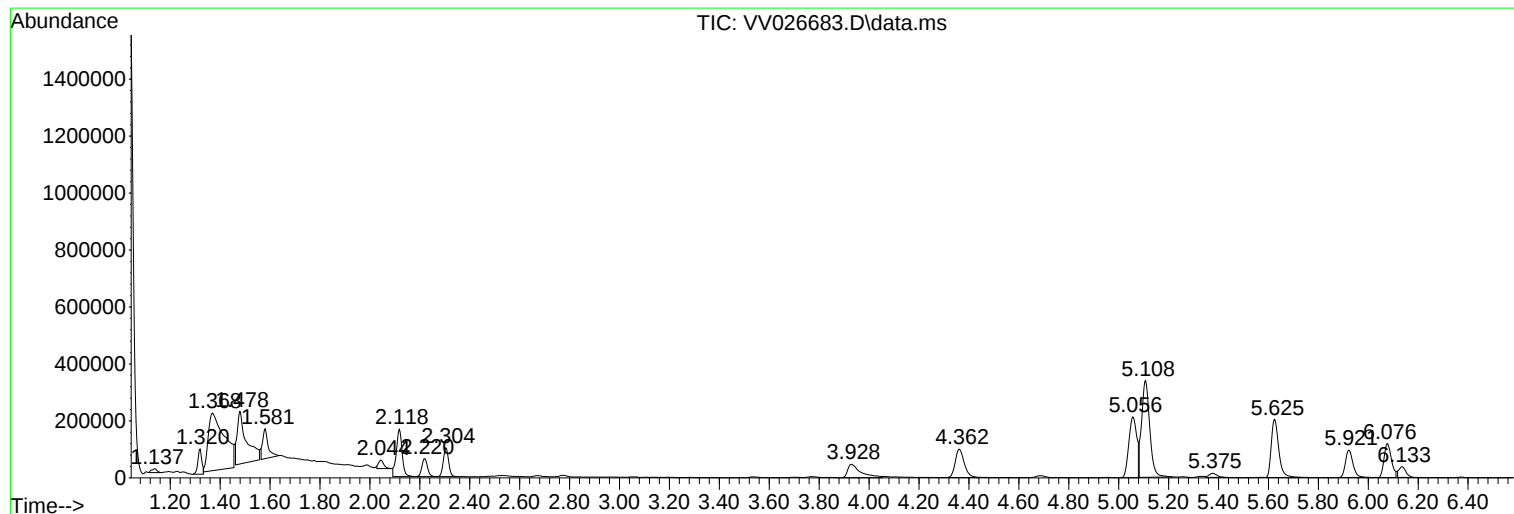
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MSVOA_V
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

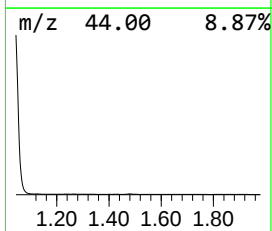
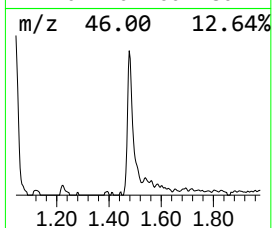
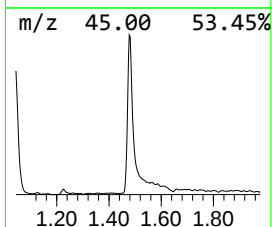
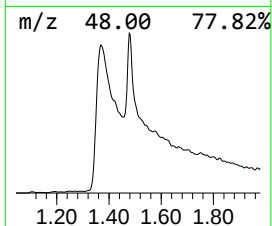
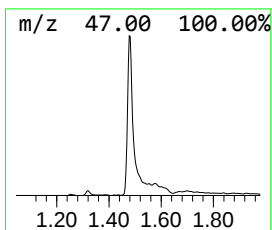
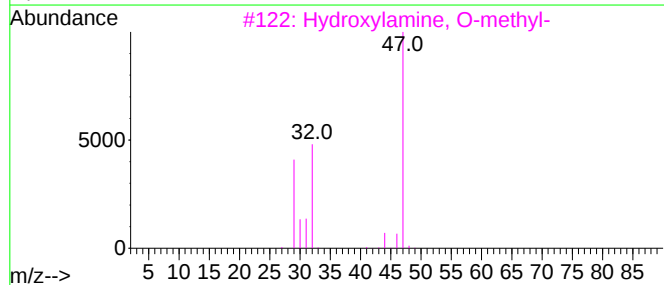
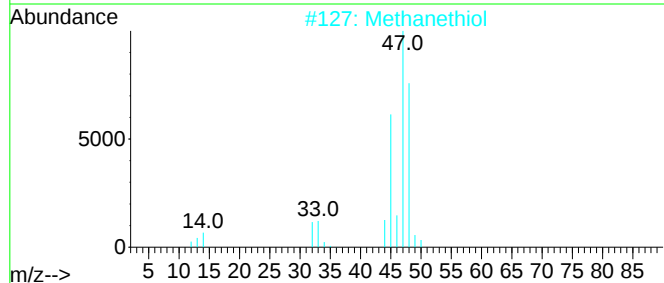
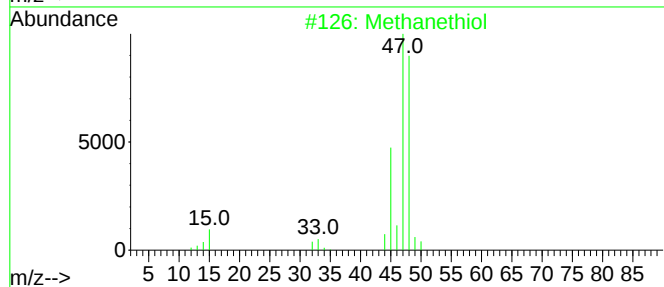
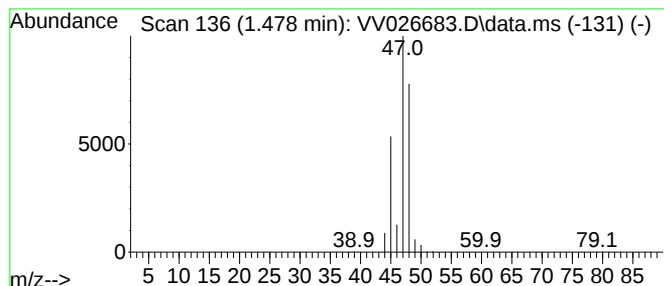
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Methanethiol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.478	5.70 ug/L	483807	1,4-Difluorobenzene	5.625

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanethiol	48	CH4S	000074-93-1	90
2			Methanethiol	48	CH4S	000074-93-1	90
3			Hydroxylamine, O-methyl-	47	CH5NO	000067-62-9	7
4			Hydroxylamine, O-methyl-	47	CH5NO	000067-62-9	7
5			Methane, nitroso-	45	CH3NO	000865-40-7	5



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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

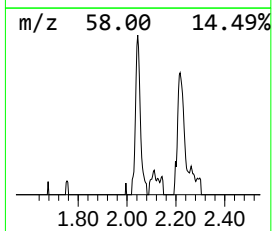
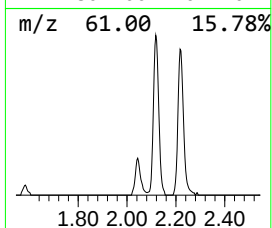
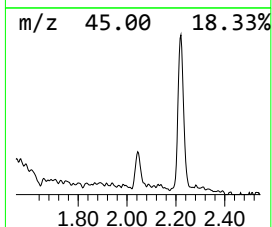
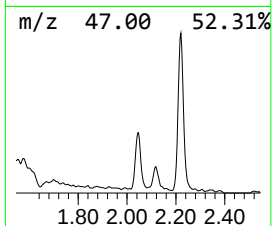
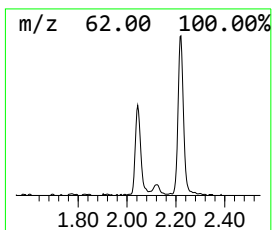
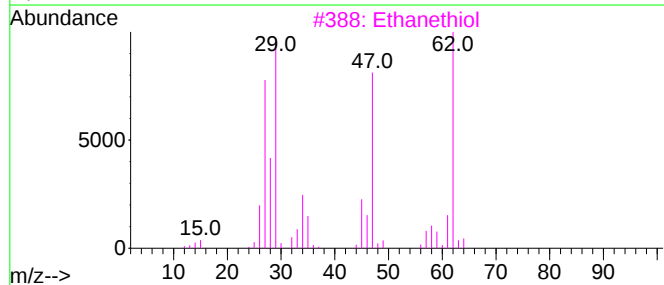
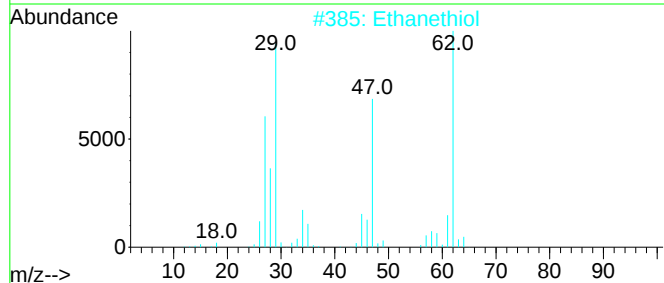
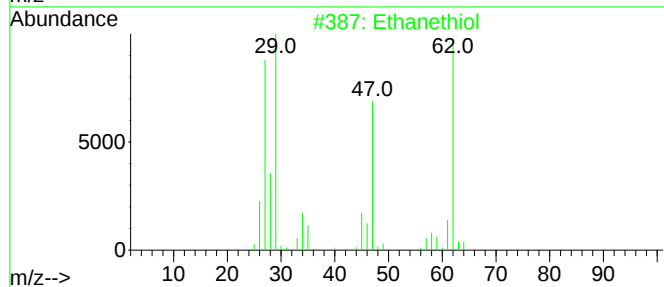
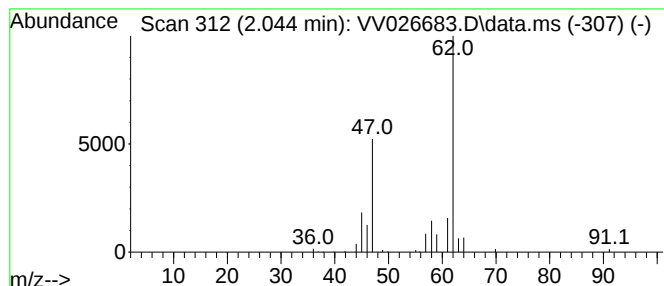
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Ethanethiol Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.044	0.51 ug/L	43084	1,4-Difluorobenzene	5.625

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanethiol			62	C2H6S	000075-08-1	90
2	Ethanethiol			62	C2H6S	000075-08-1	90
3	Ethanethiol			62	C2H6S	000075-08-1	83
4	Ethanethiol			62	C2H6S	000075-08-1	83
5	Ethanethiol			62	C2H6S	000075-08-1	83



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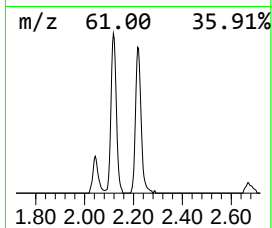
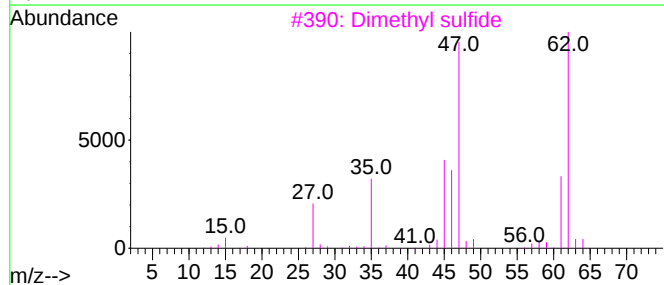
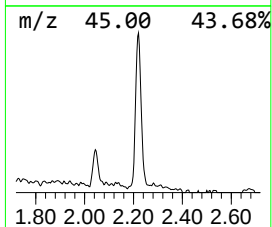
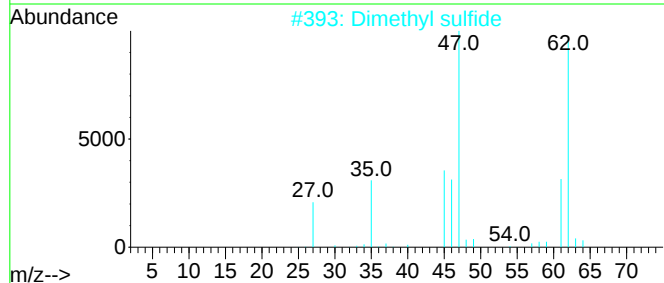
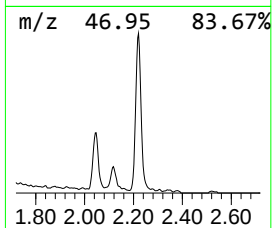
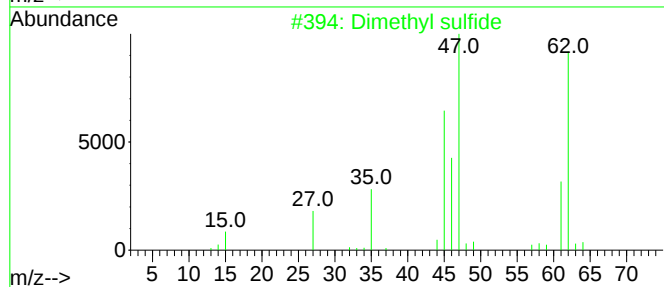
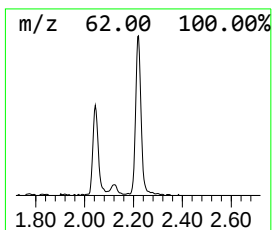
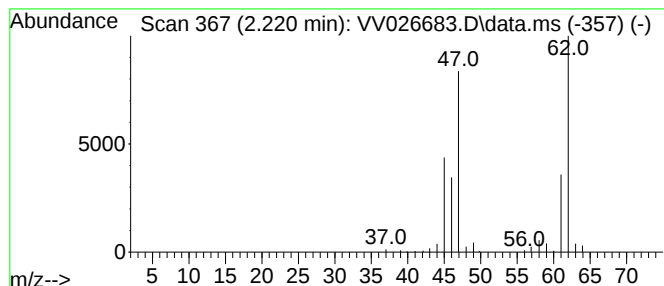
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Dimethyl sulfide Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.220	1.14 ug/L	96441	1,4-Difluorobenzene	5.625

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyl sulfide	62	C2H6S	000075-18-3	91
2			Dimethyl sulfide	62	C2H6S	000075-18-3	91
3			Dimethyl sulfide	62	C2H6S	000075-18-3	91
4			Borane-methyl sulfide complex	76	C2H9BS	013292-87-0	83
5			Ethanethiol	62	C2H6S	000075-08-1	78



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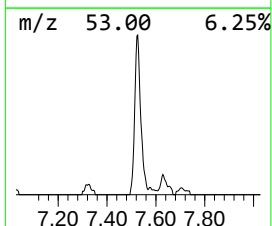
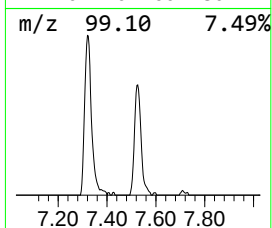
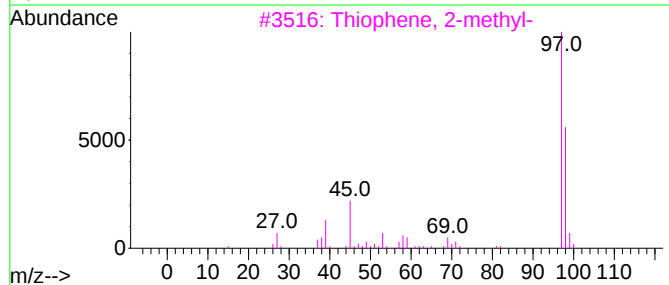
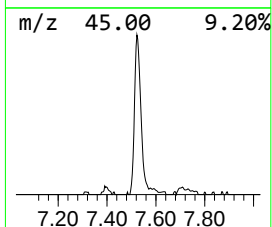
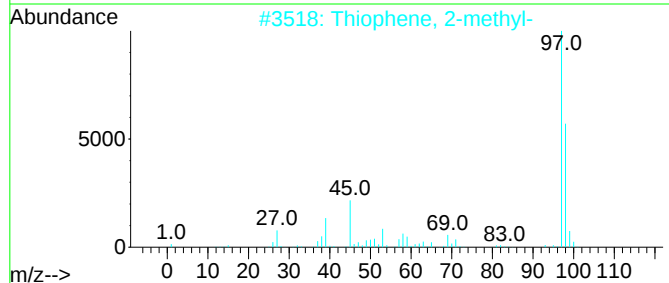
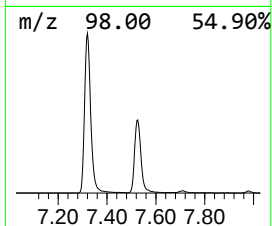
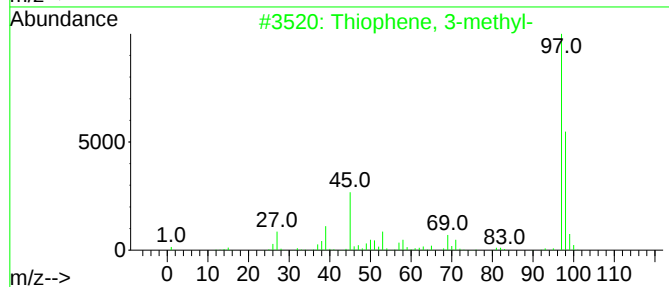
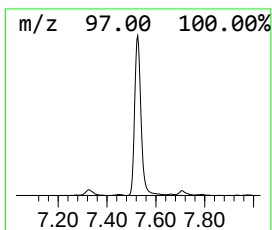
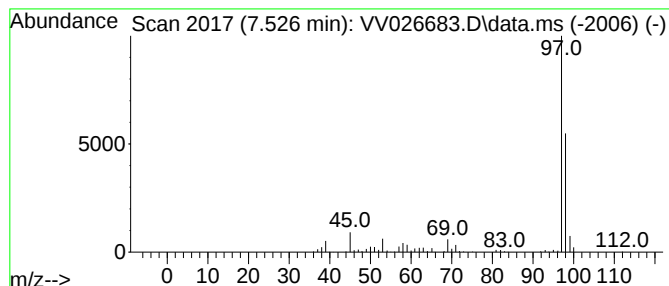
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Thiophene, 3-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.526	0.81 ug/L	320647	Chlorobenzene-d5	8.854

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiophene, 3-methyl-	98	C5H6S	000616-44-4	91
2			Thiophene, 2-methyl-	98	C5H6S	000554-14-3	91
3			Thiophene, 2-methyl-	98	C5H6S	000554-14-3	91
4			Thiophene, 3-methyl-	98	C5H6S	000616-44-4	91
5			Thiophene, 3-methyl-	98	C5H6S	000616-44-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

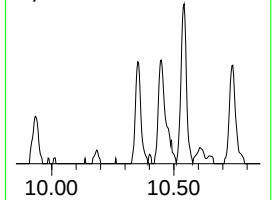
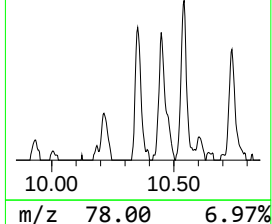
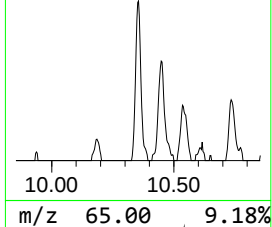
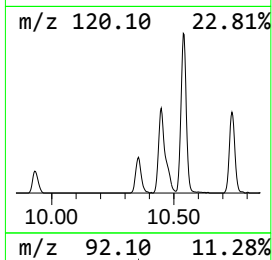
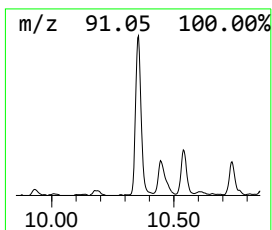
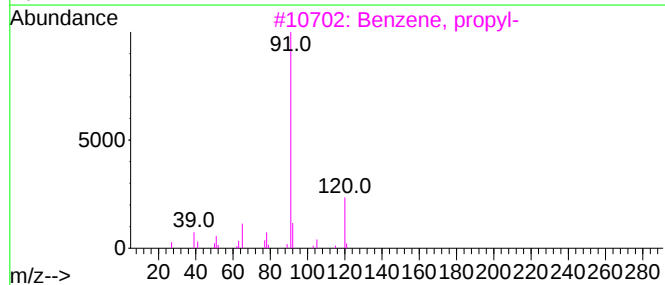
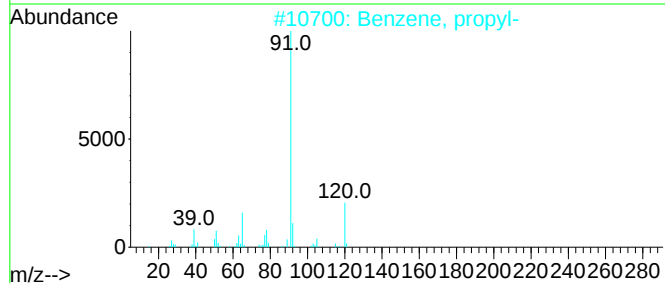
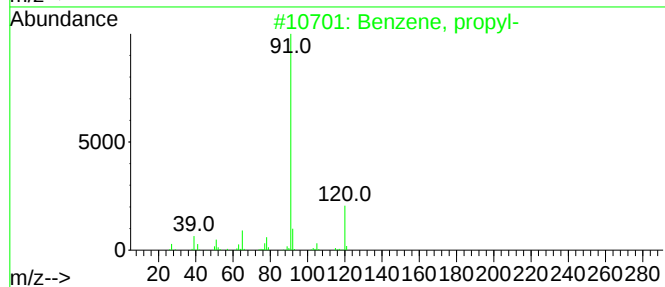
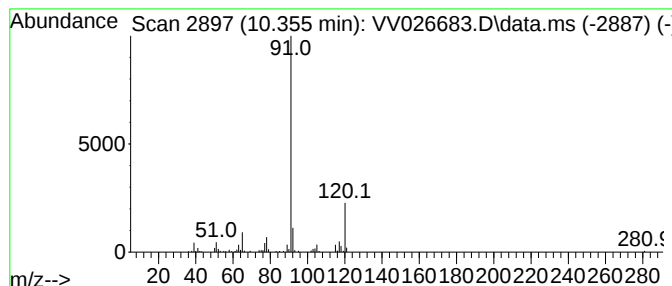
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, propyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	0.82 ug/L	122890	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	81
2			Benzene, propyl-	120	C9H12	000103-65-1	81
3			Benzene, propyl-	120	C9H12	000103-65-1	76
4			Benzene, propyl-	120	C9H12	000103-65-1	76
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

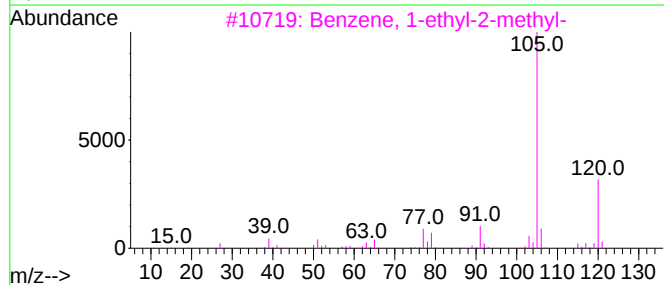
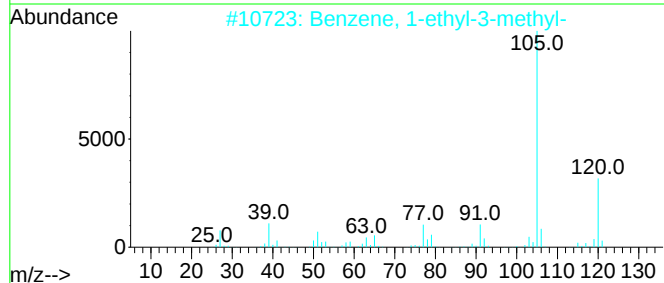
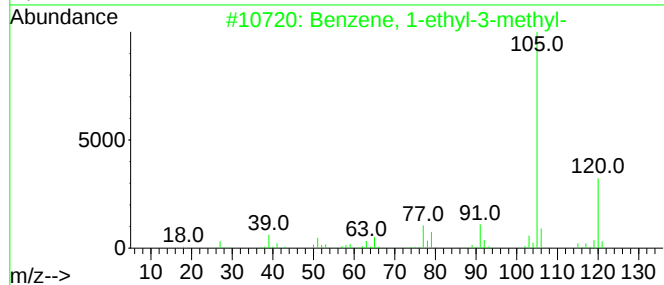
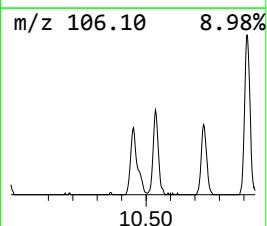
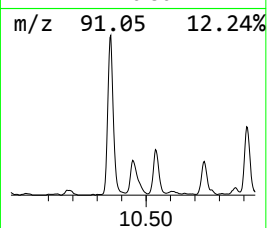
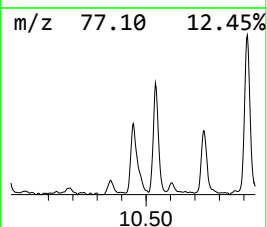
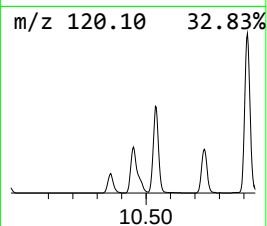
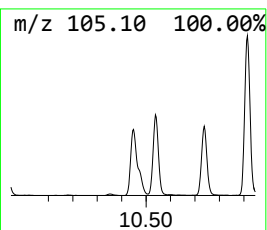
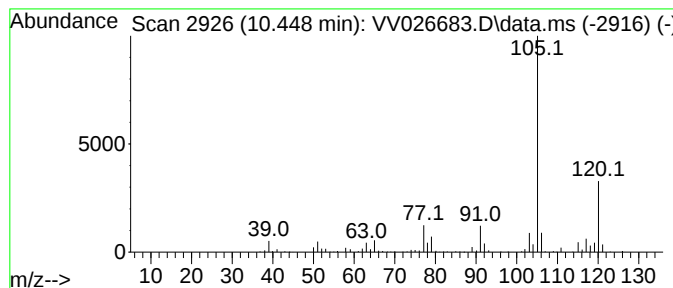
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	1.74 ug/L	260177	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

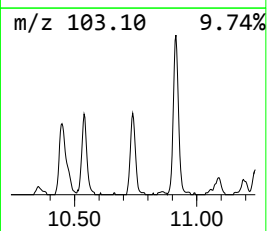
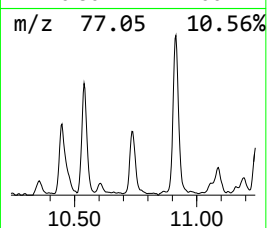
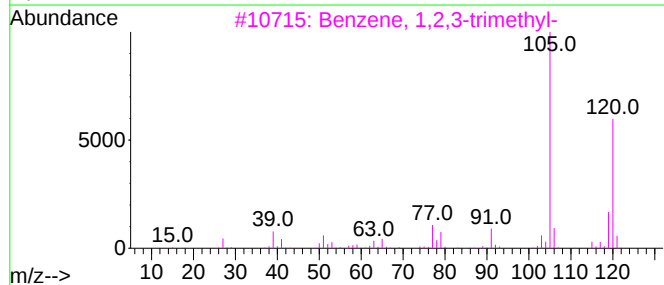
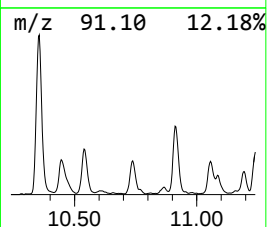
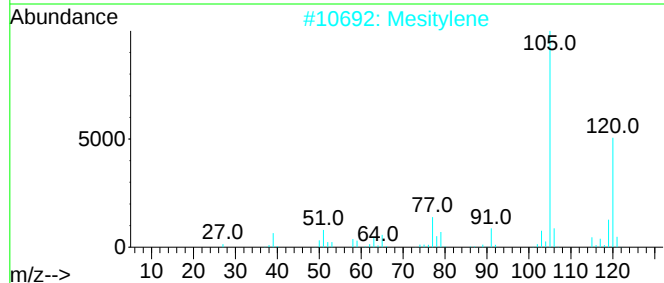
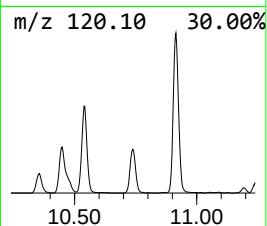
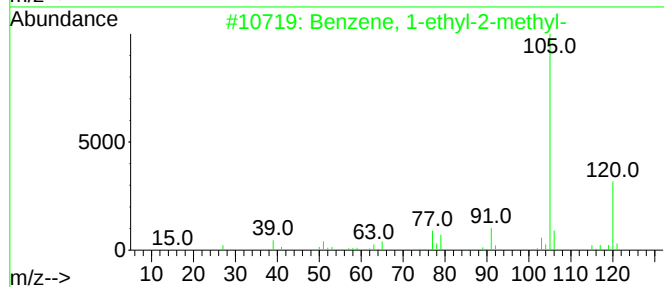
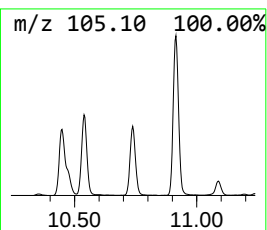
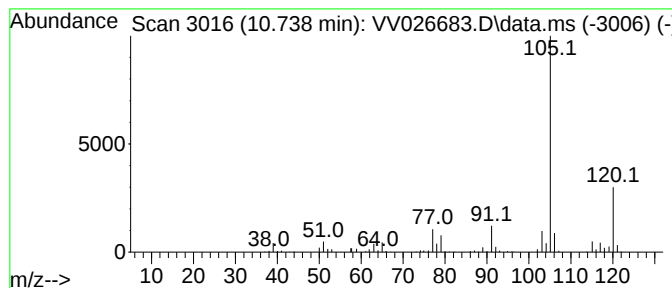
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.738	1.42 ug/L	212516	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	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_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

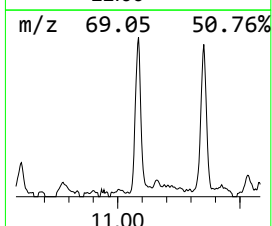
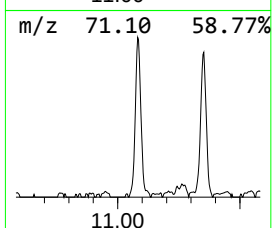
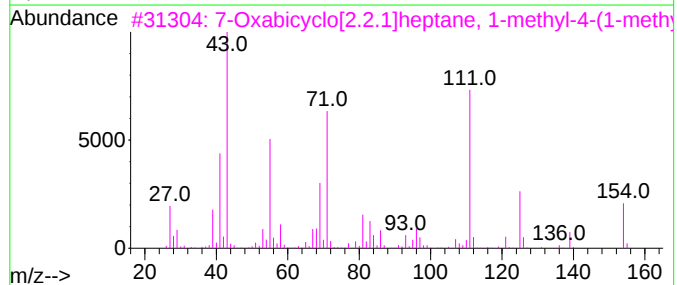
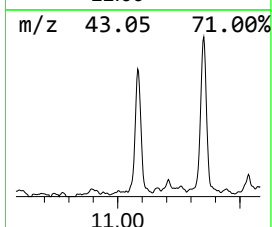
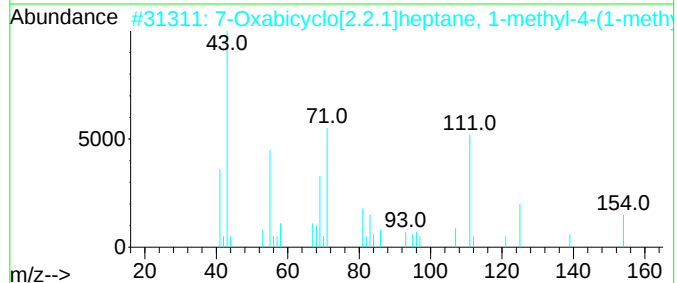
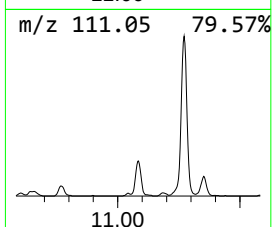
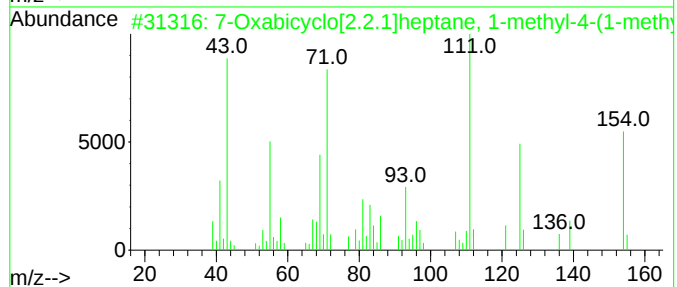
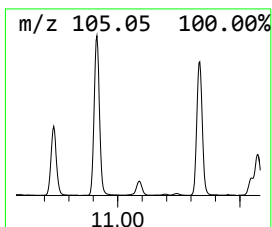
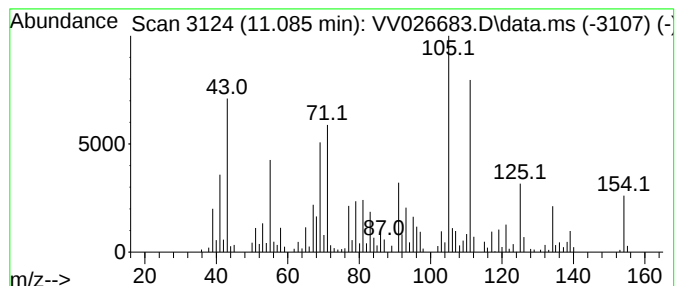
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 7-Oxabicyclo[2.2.1]heptane,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.085	1.38 ug/L	206167	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	94
2			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	93
3			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	90
4			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	89
5			7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

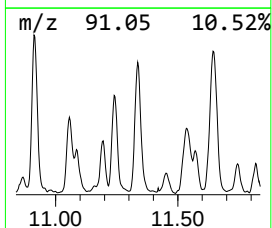
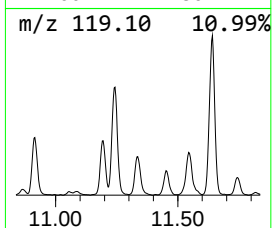
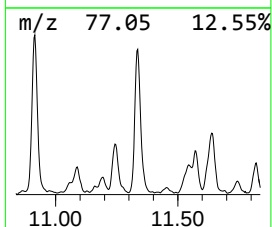
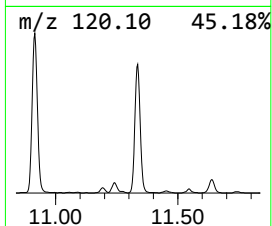
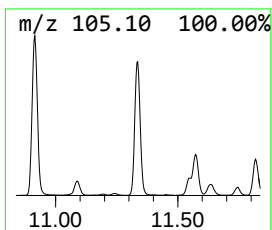
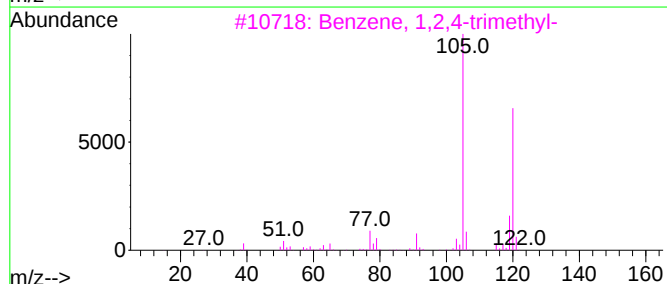
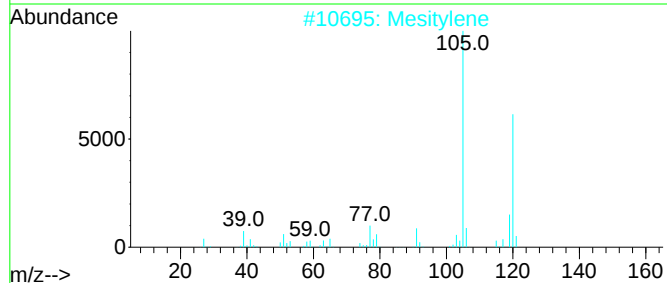
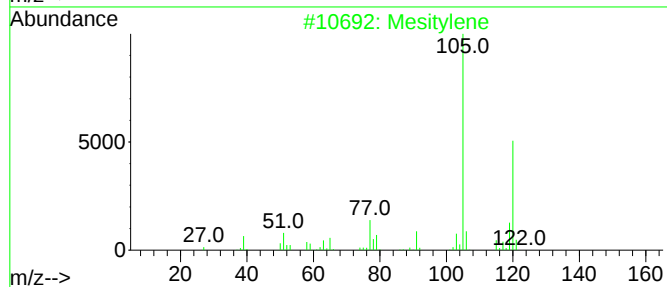
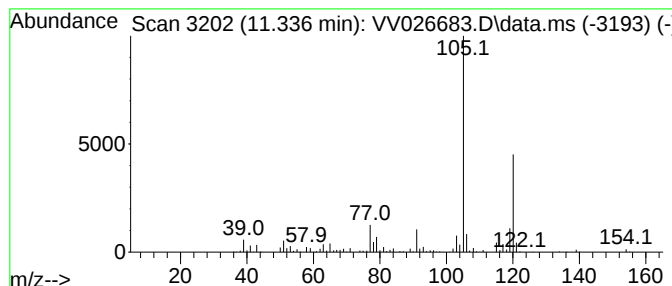
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	3.92 ug/L	586942	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

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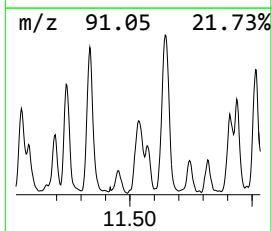
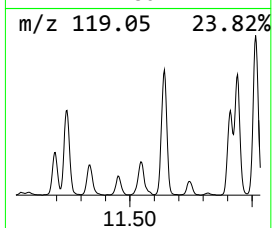
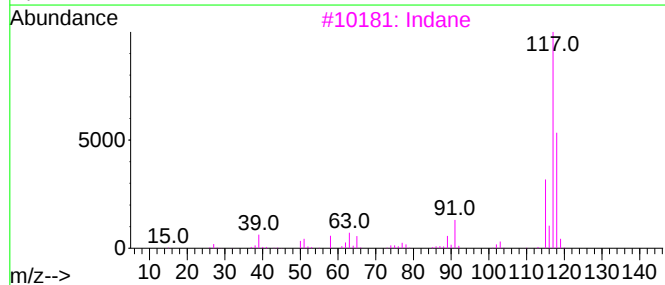
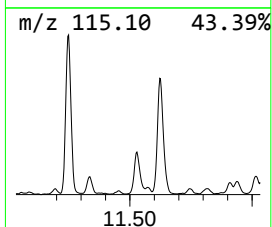
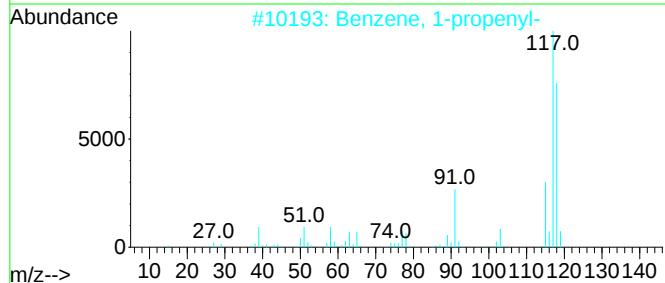
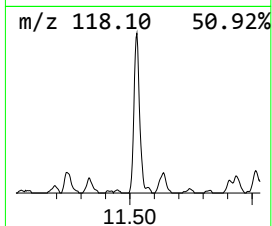
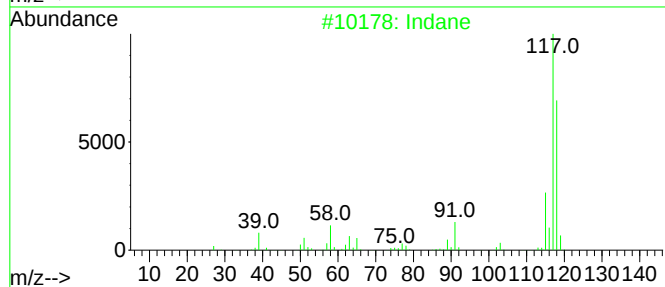
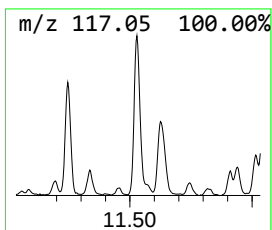
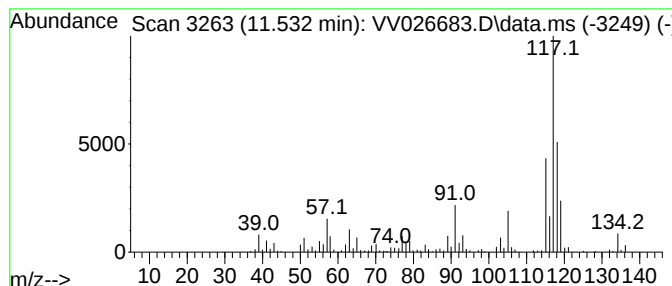
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Indane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.532	1.61 ug/L	240840	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	64
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
3			Indane	118	C9H10	000496-11-7	60
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	60
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

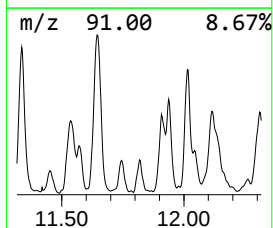
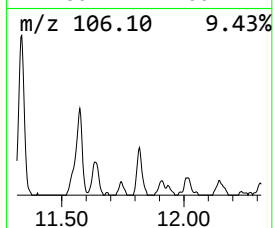
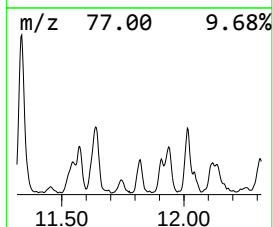
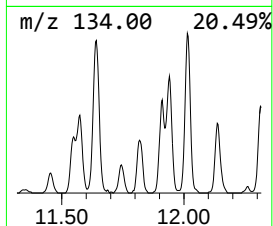
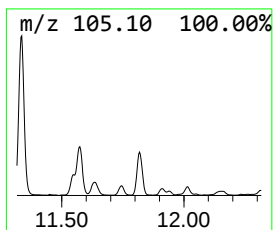
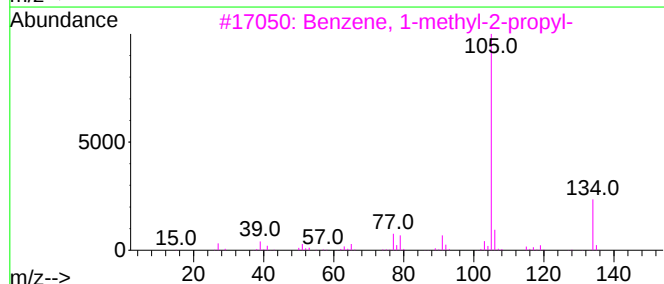
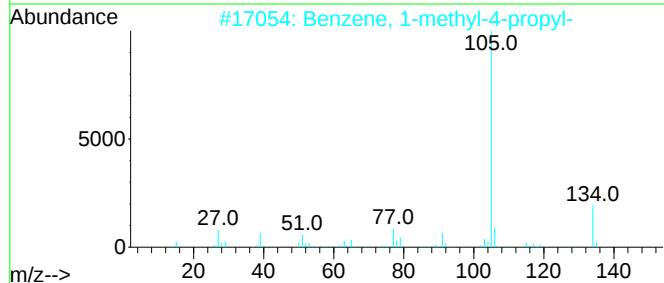
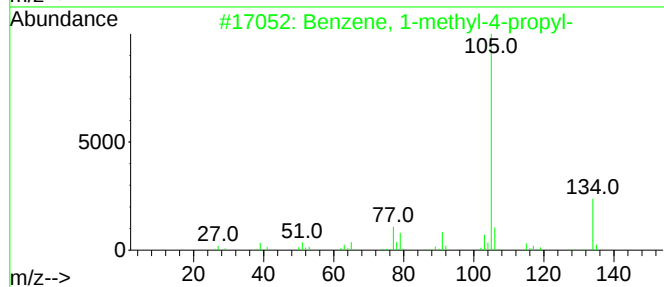
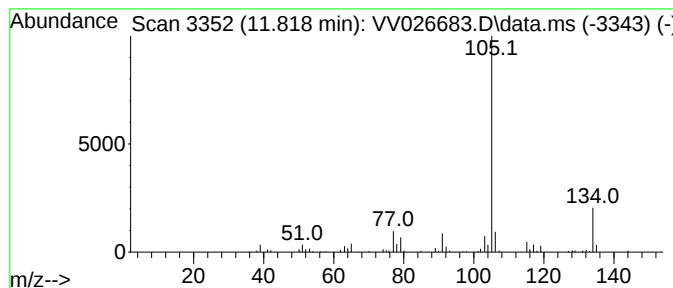
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzene, 1-methyl-4-propyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.818	0.63 ug/L	94001	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
4			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

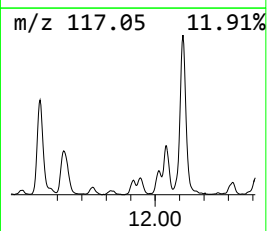
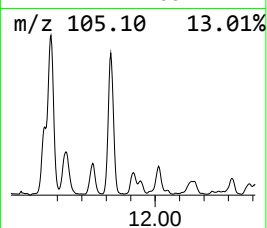
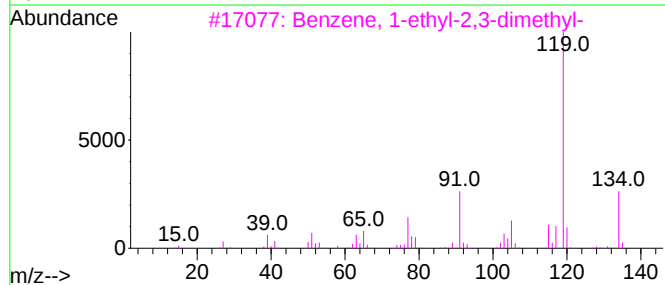
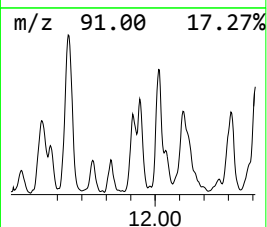
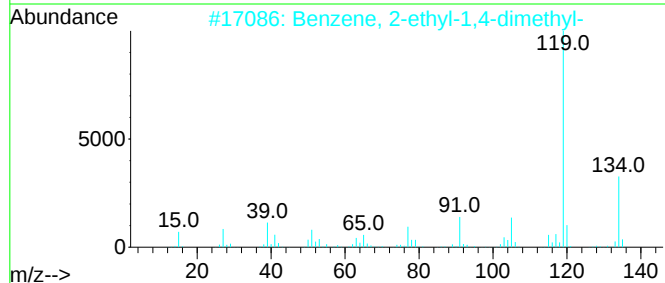
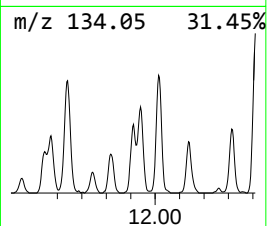
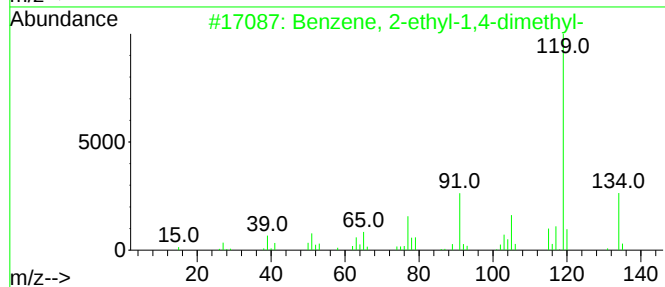
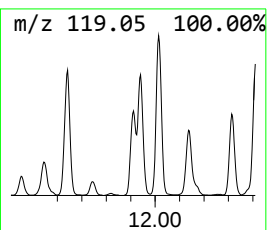
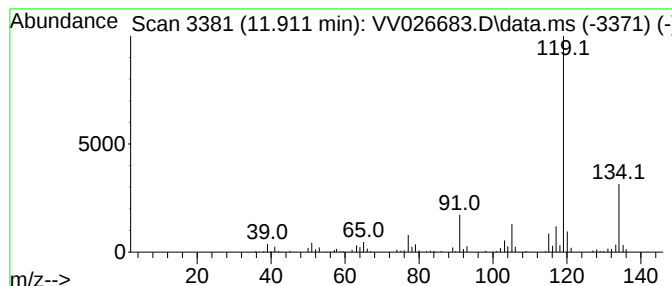
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	0.92 ug/L	137199	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

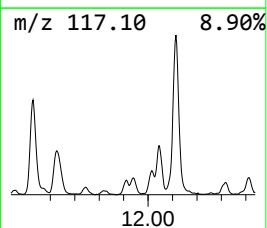
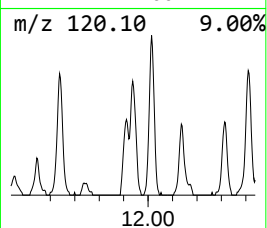
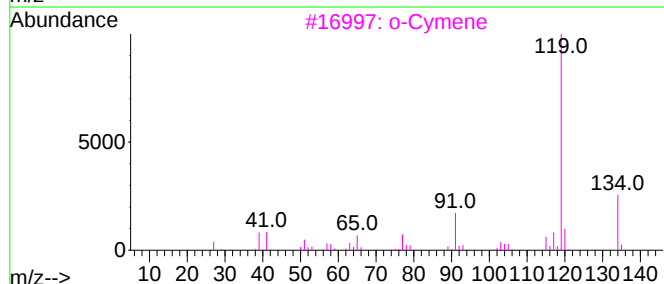
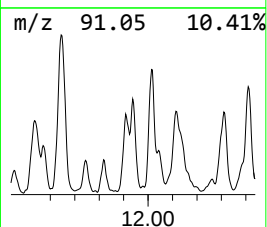
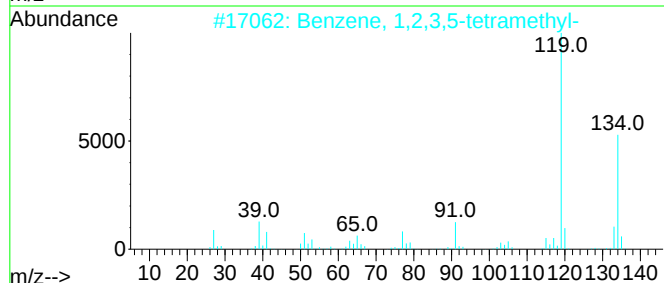
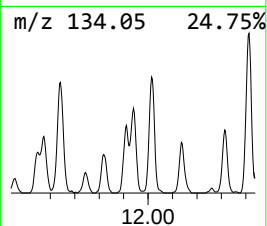
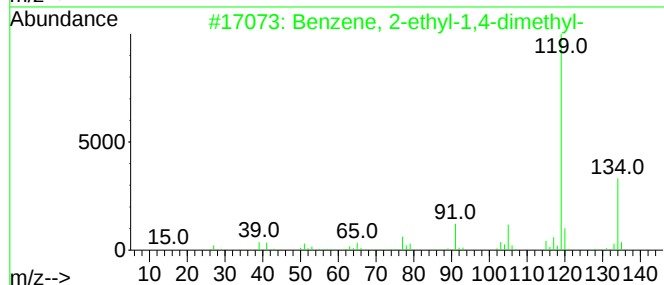
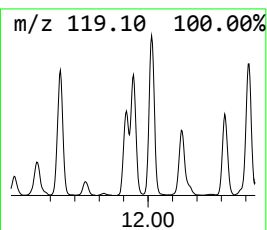
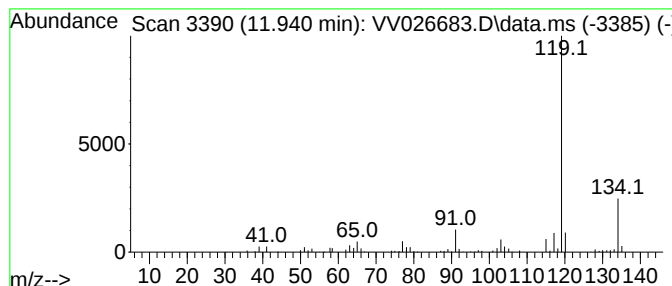
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 14 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.940	1.01 ug/L	151186	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3			o-Cymene	134	C10H14	000527-84-4	91
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

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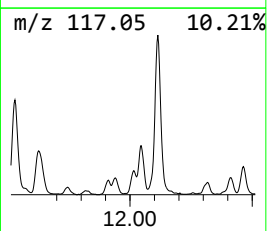
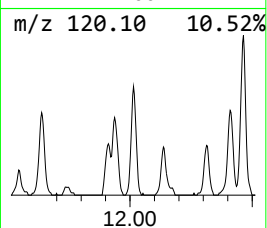
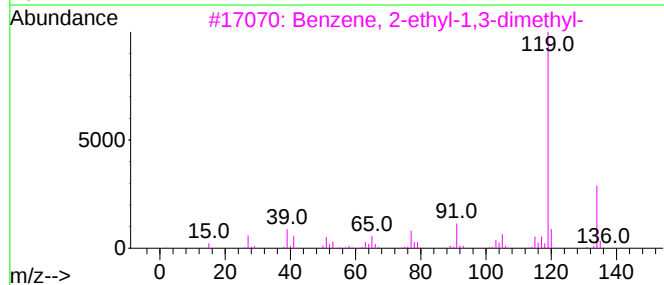
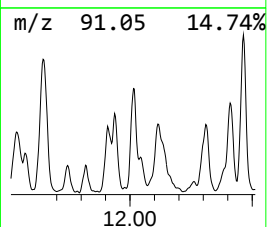
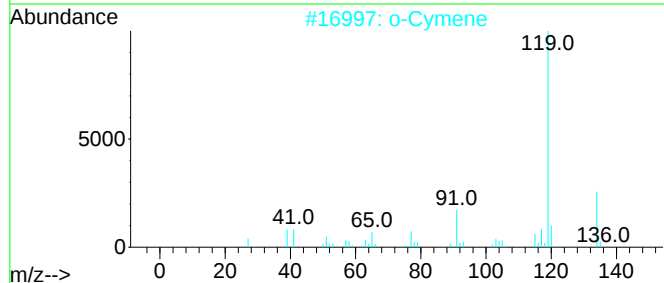
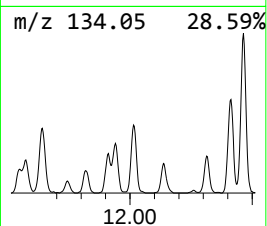
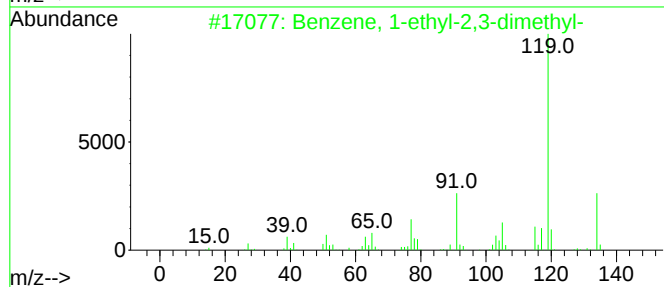
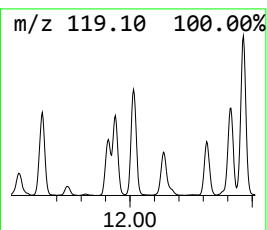
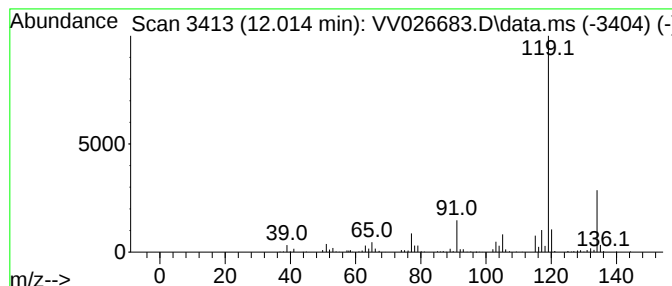
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.014	1.47 ug/L	220608	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

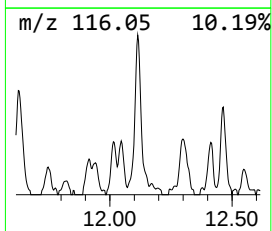
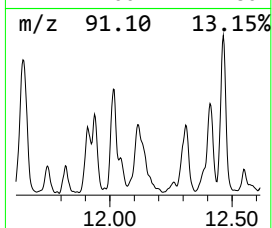
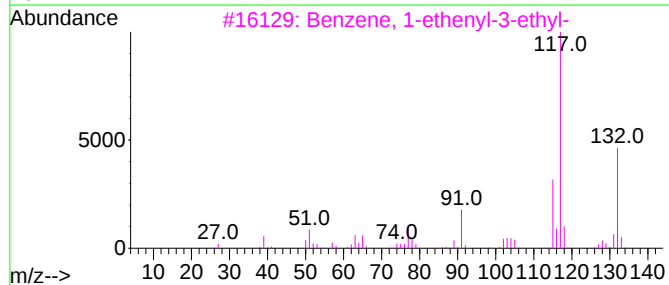
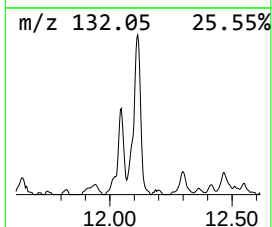
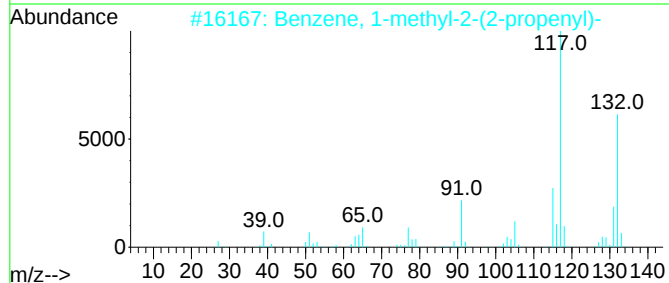
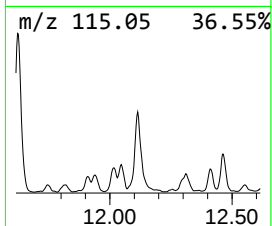
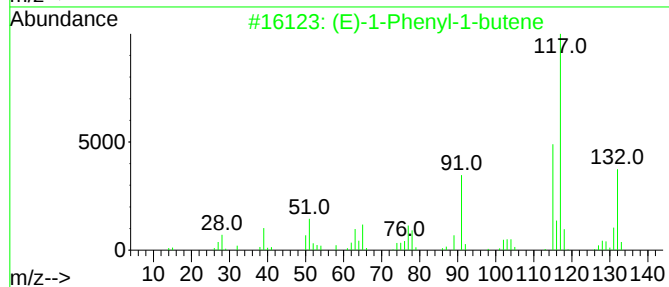
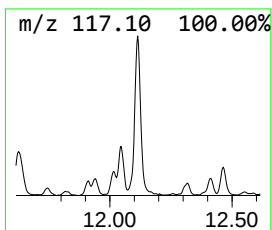
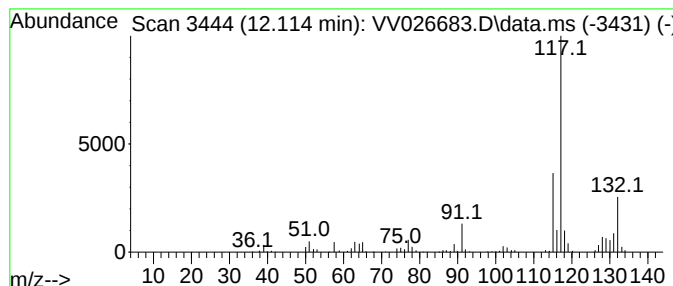
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 16 (E)-1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.114	1.97 ug/L	295196	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
4			Indan, 1-methyl-	132	C10H12	000767-58-8	81
5			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

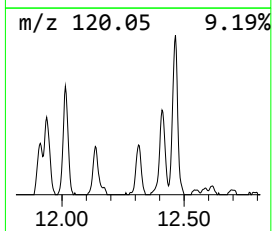
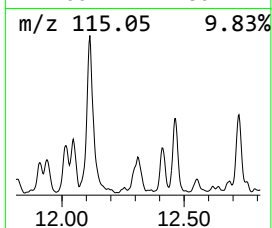
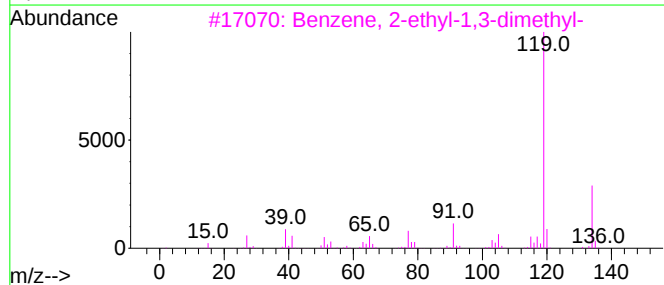
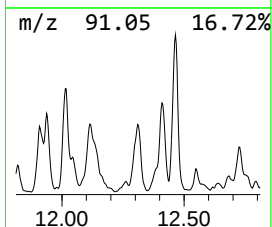
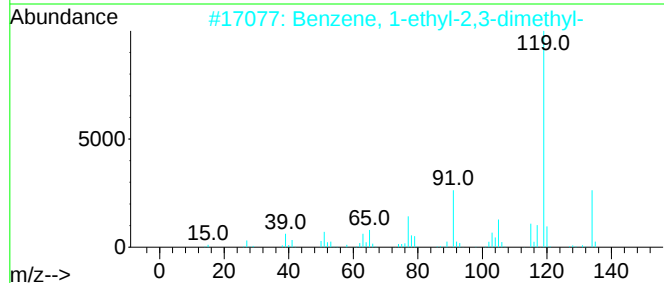
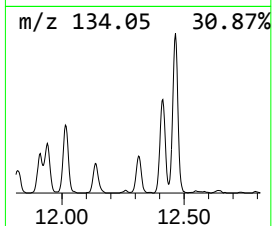
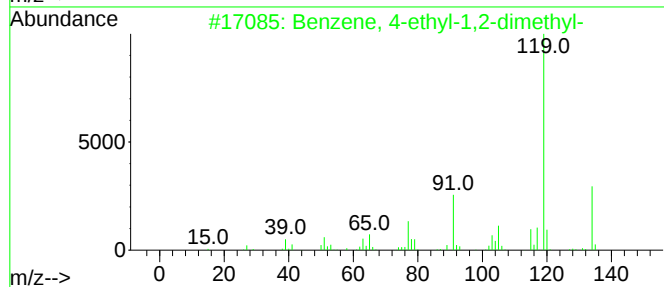
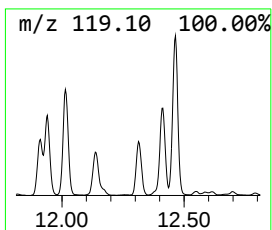
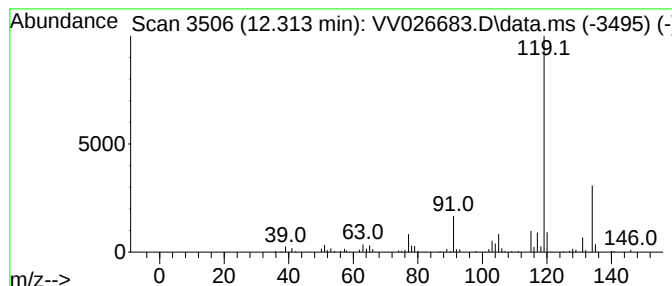
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.313	1.11 ug/L	166778	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

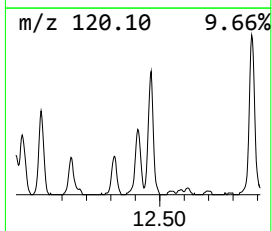
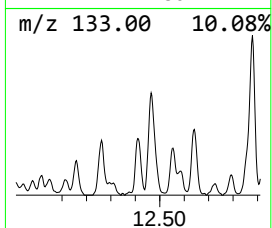
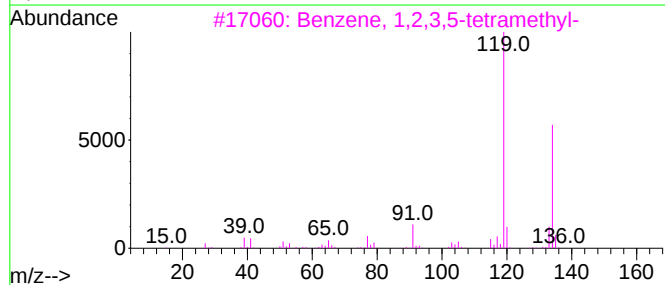
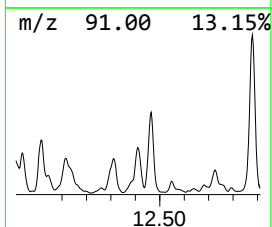
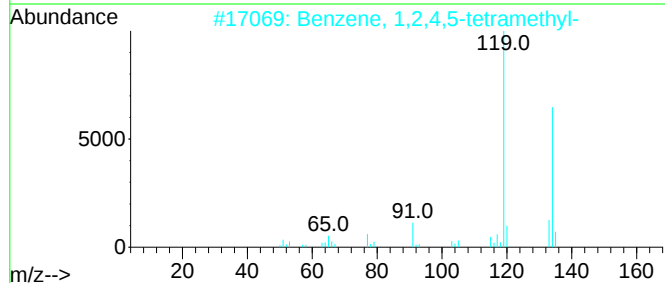
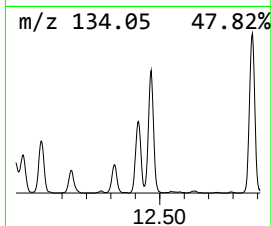
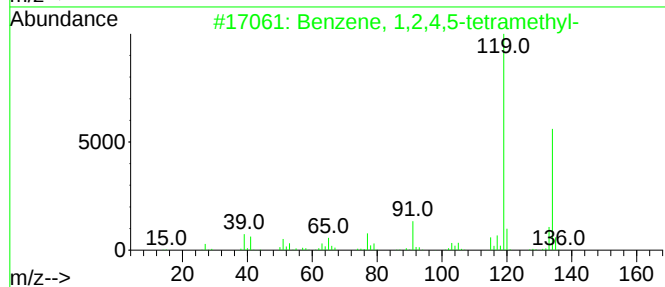
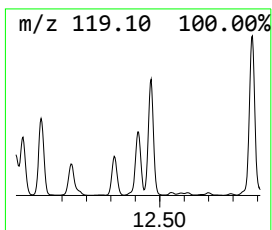
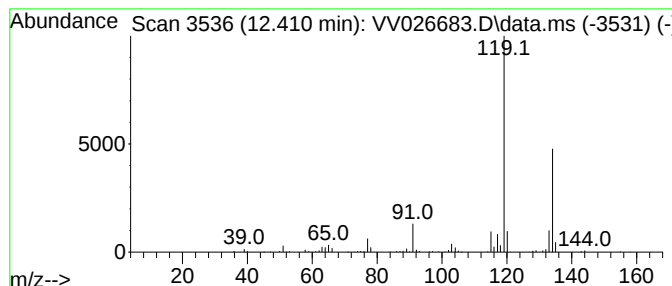
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 18 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	1.50 ug/L	225170	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

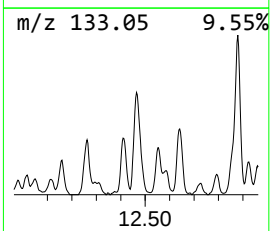
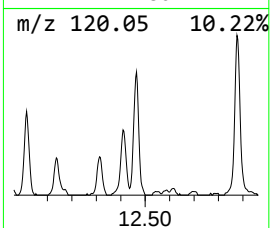
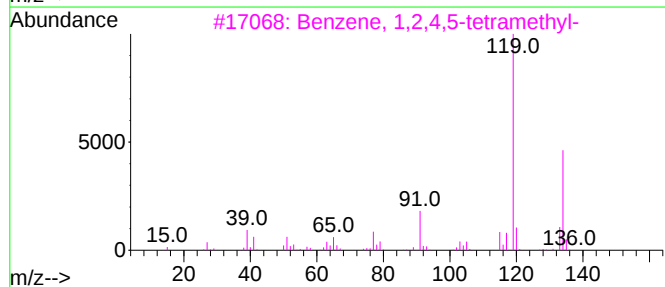
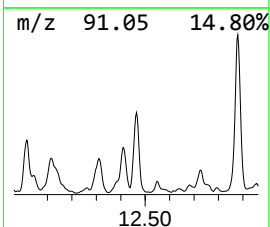
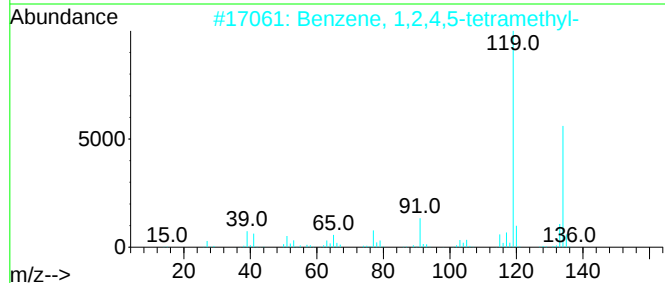
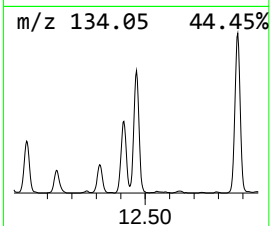
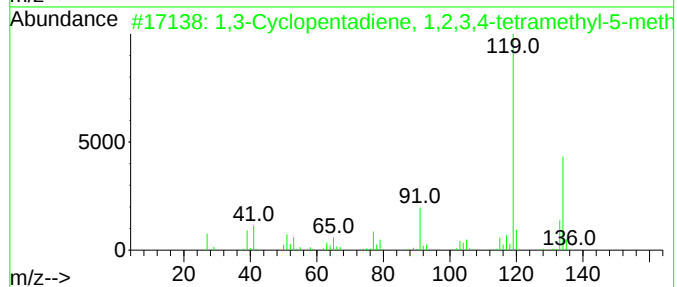
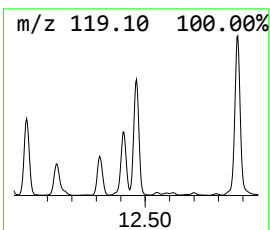
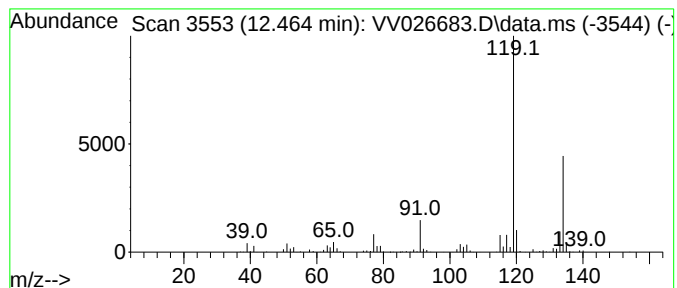
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 19 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	2.53 ug/L	378461	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

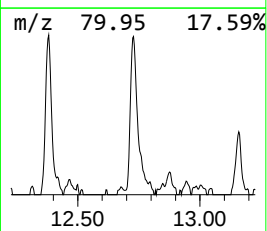
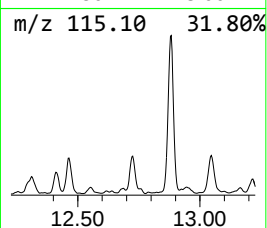
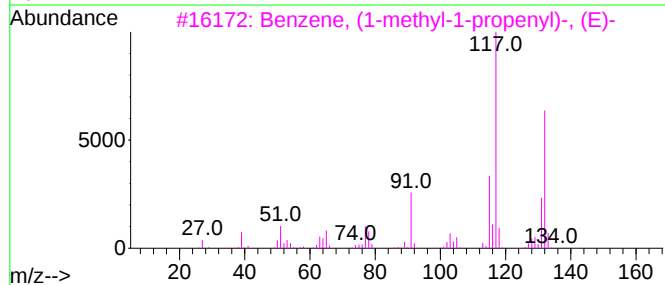
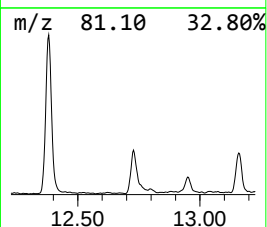
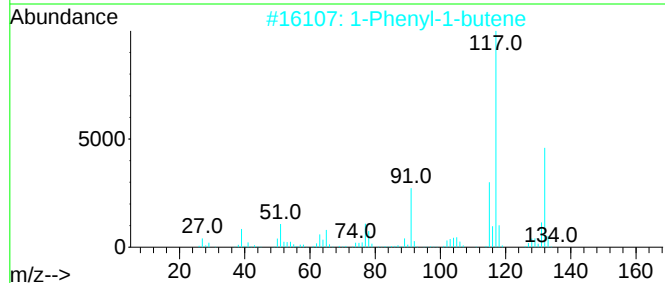
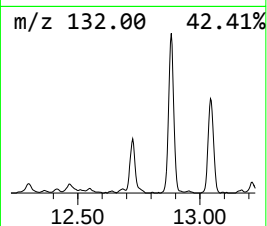
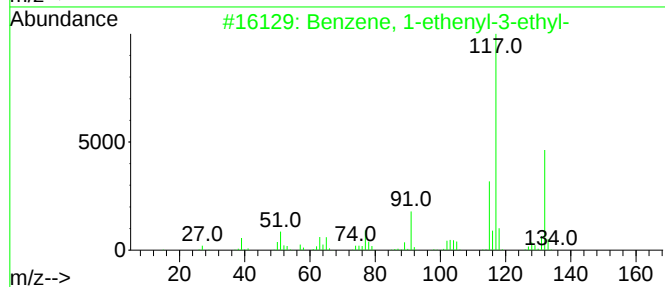
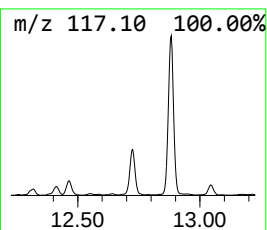
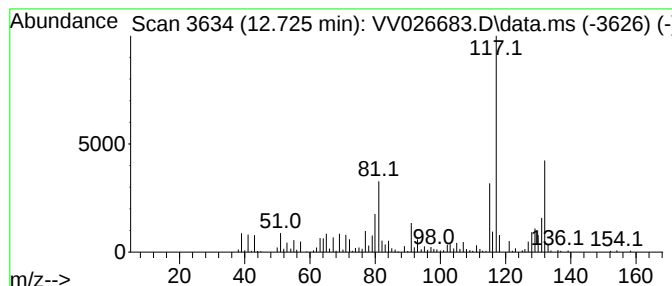
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 20 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.725	1.39 ug/L	208839	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

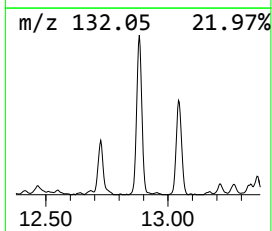
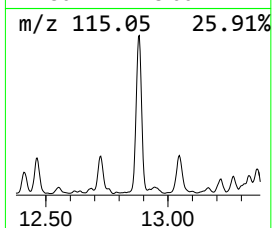
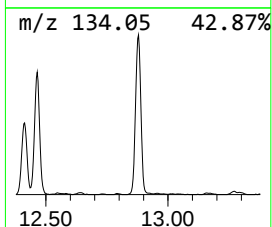
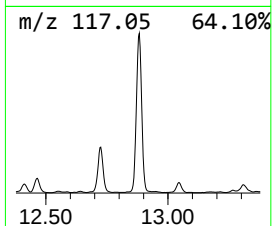
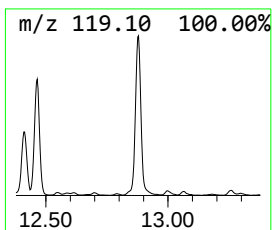
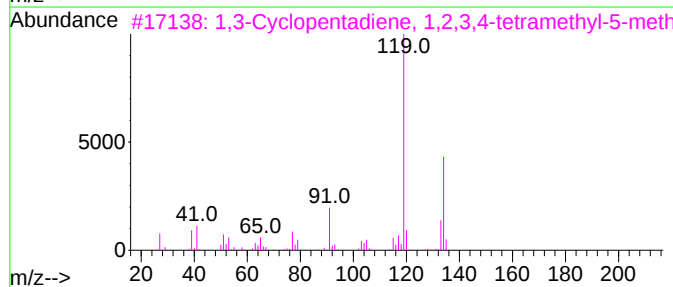
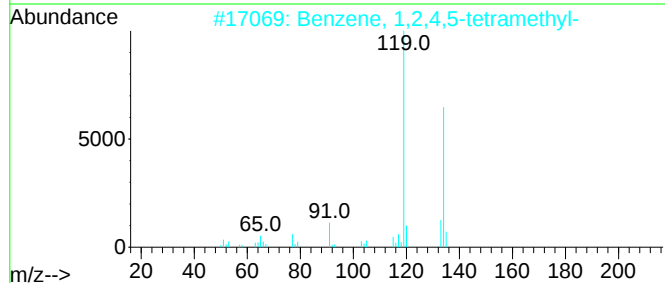
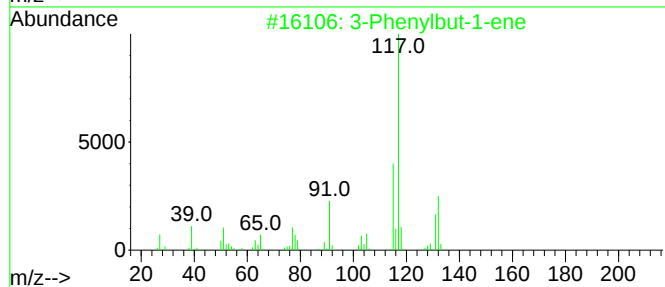
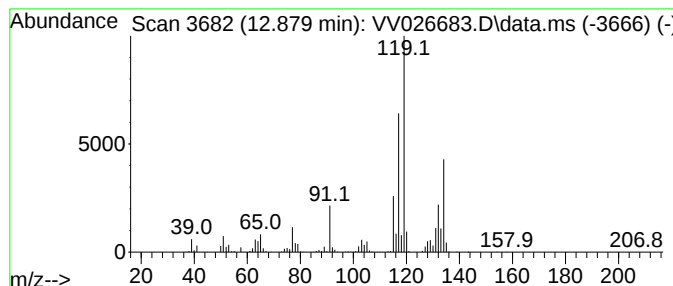
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 21 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.879	5.62 ug/L	841513	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	80
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
3			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	64
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

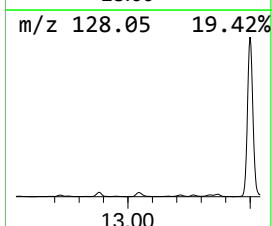
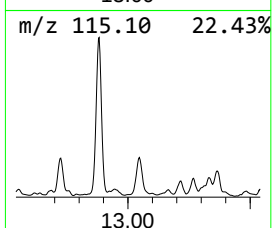
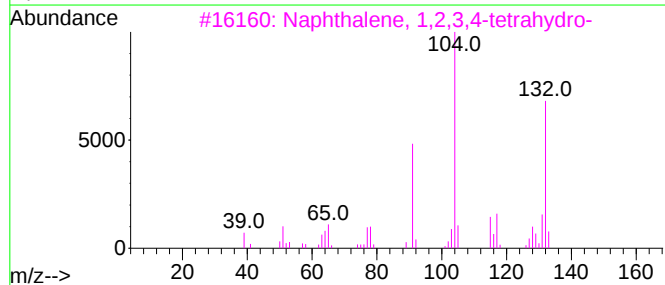
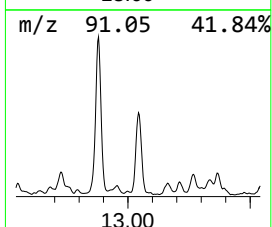
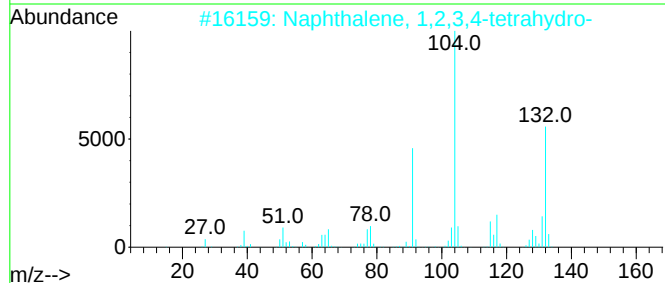
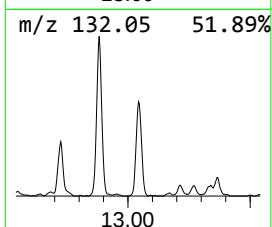
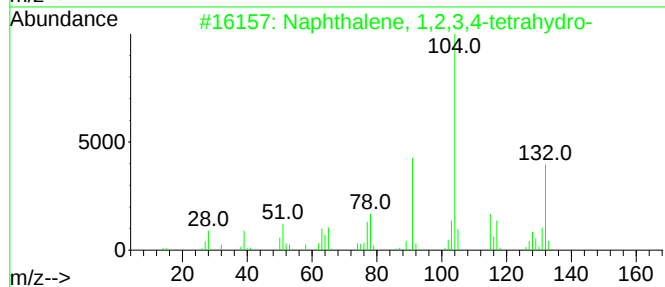
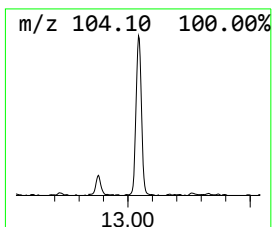
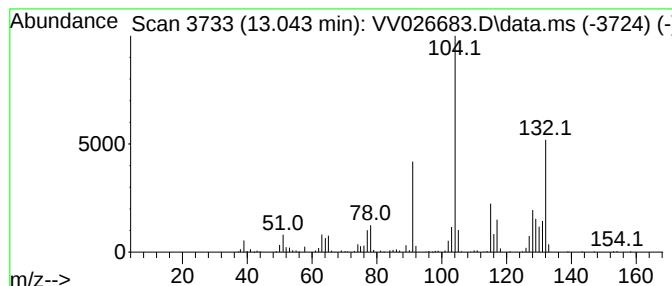
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 23 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.043	1.68 ug/L	252010	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	76
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

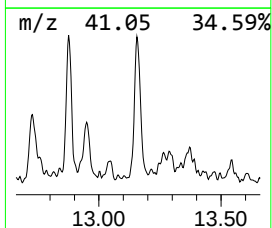
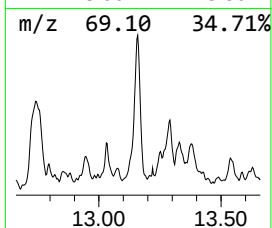
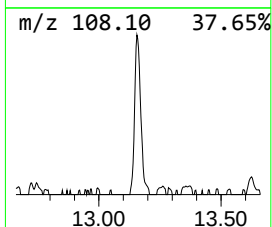
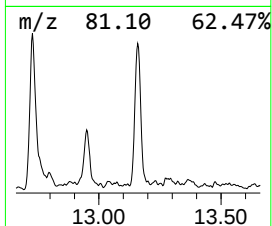
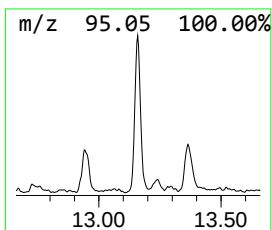
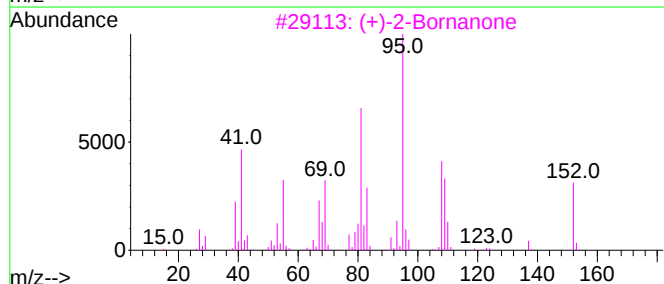
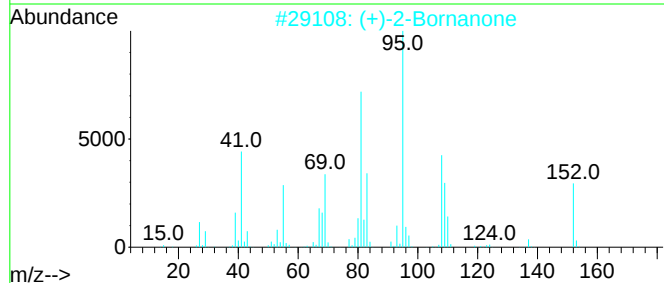
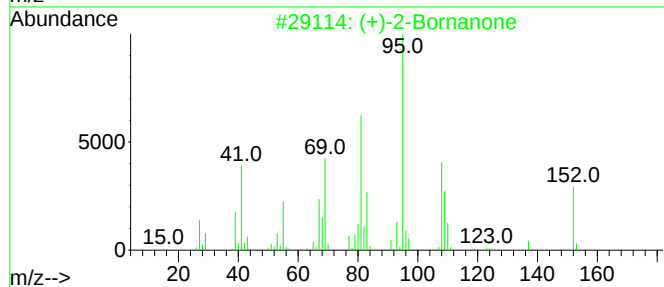
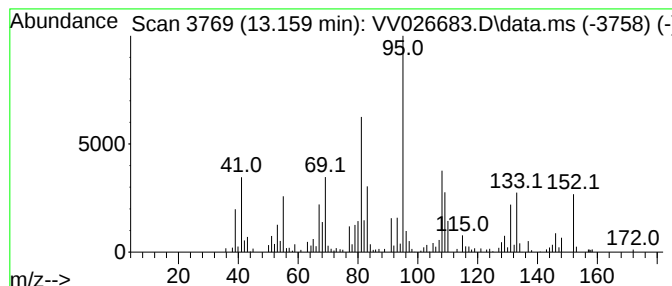
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 24 (+)-2-Bornanone Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.159	1.02 ug/L	152547	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone			152	C10H16O	000464-49-3	98
2	(+)-2-Bornanone			152	C10H16O	000464-49-3	98
3	(+)-2-Bornanone			152	C10H16O	000464-49-3	98
4	Camphor			152	C10H16O	000076-22-2	98
5	Camphor			152	C10H16O	000076-22-2	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

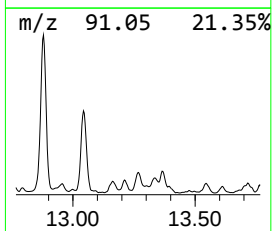
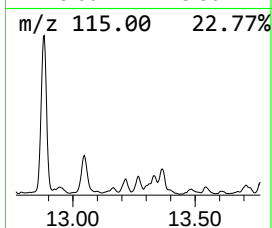
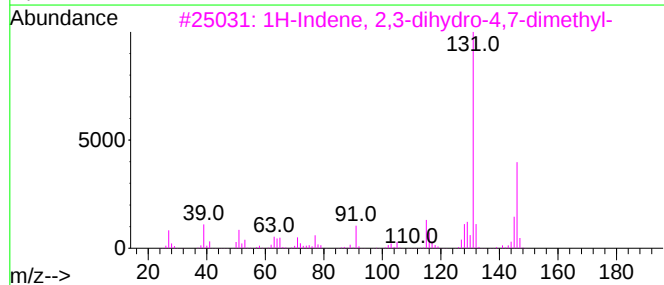
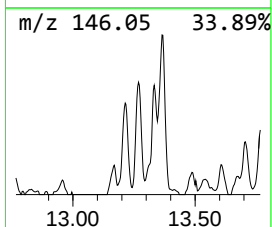
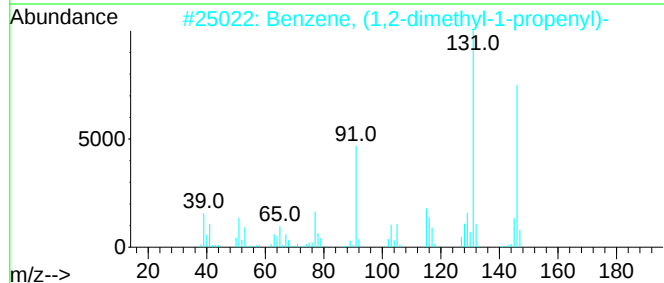
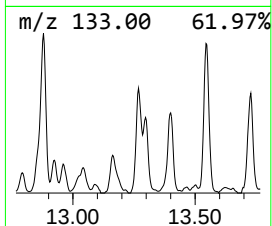
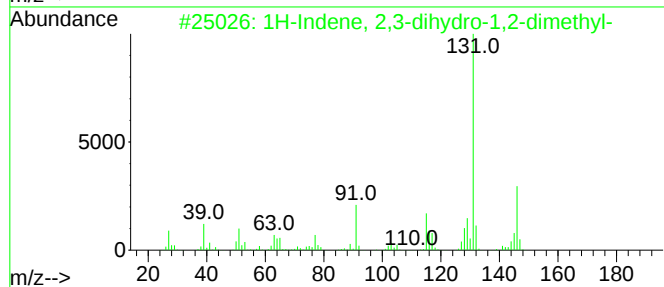
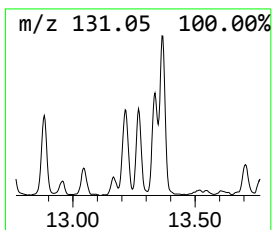
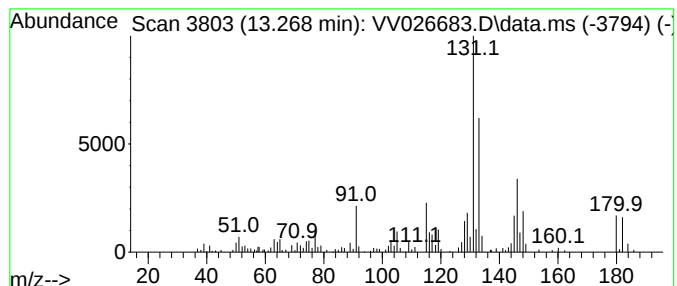
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 25 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	0.93 ug/L	139835	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	78		
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70		
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60		
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60		
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

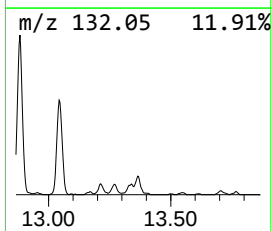
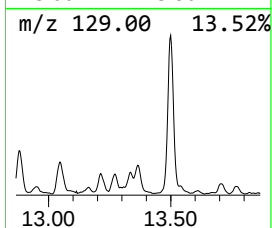
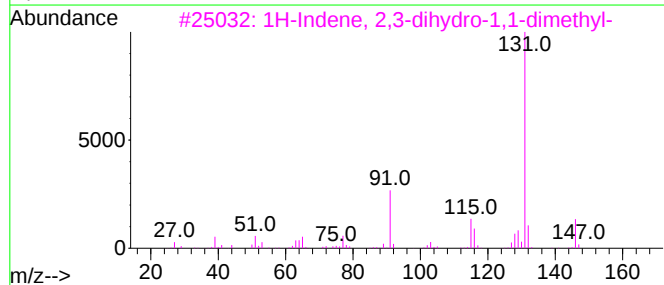
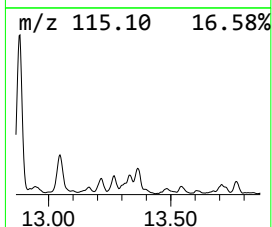
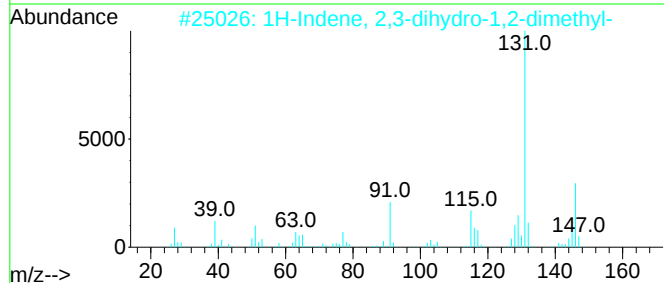
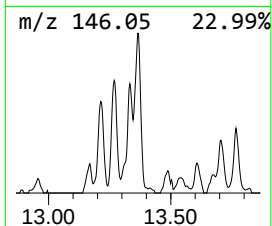
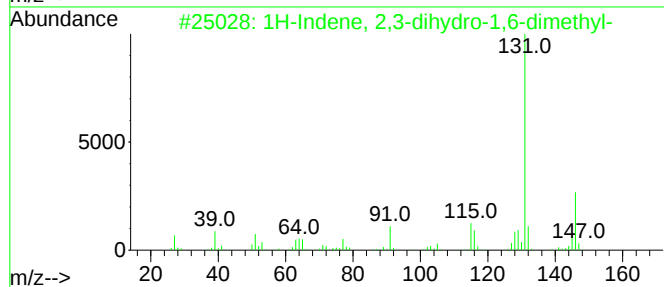
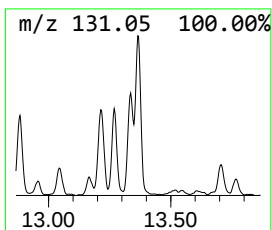
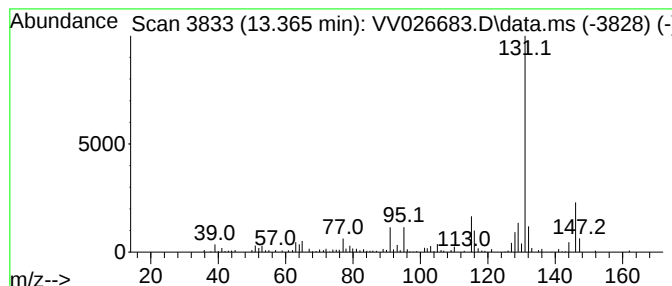
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 26 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.365	1.28 ug/L	191603	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
3			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	90
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

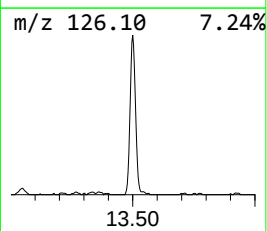
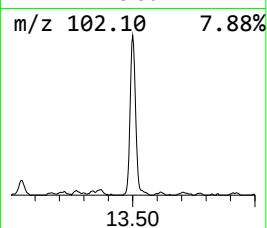
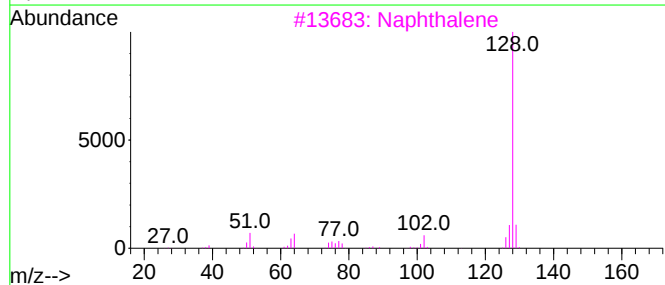
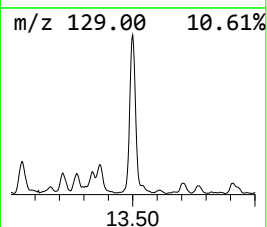
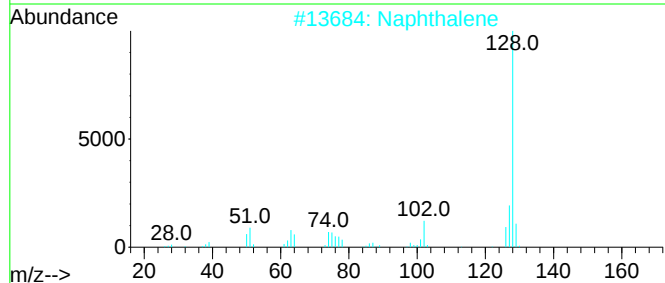
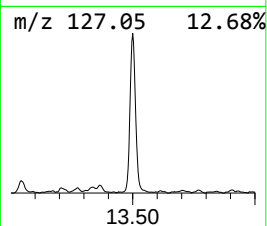
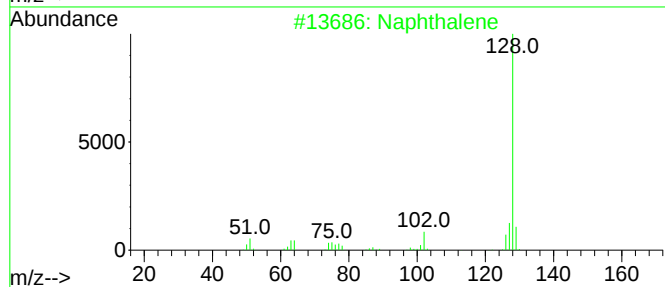
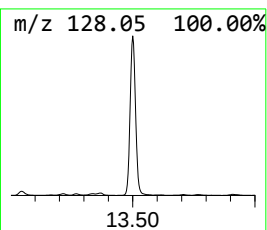
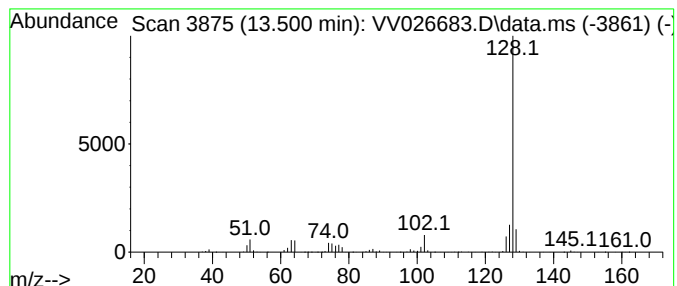
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 27 Azulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	4.91 ug/L	736010	1,4-Dichlorobenzene-d4	11.249

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			Naphthalene	128	C10H8	000091-20-3	93
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

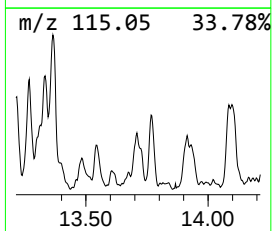
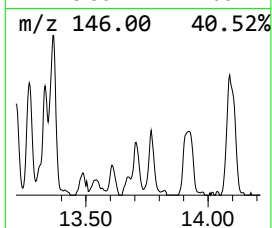
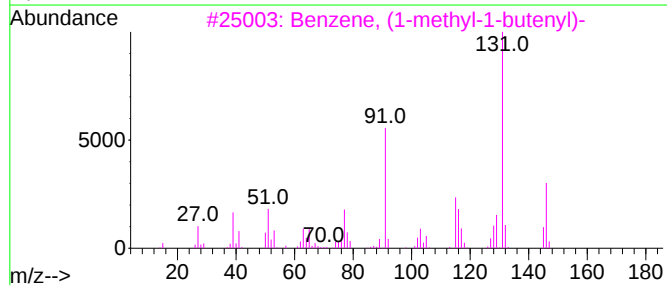
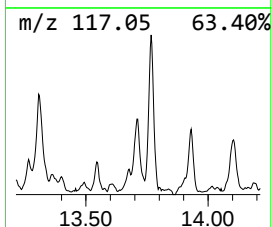
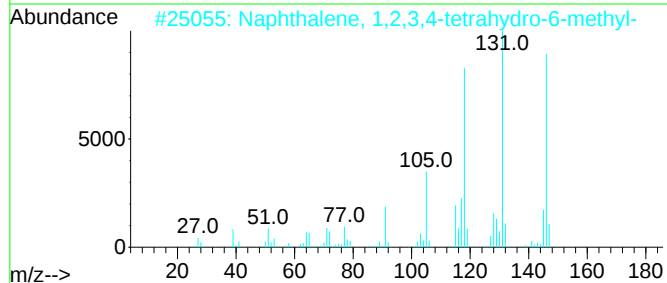
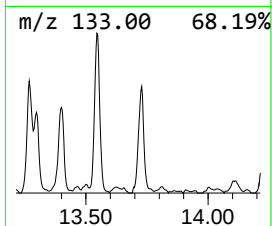
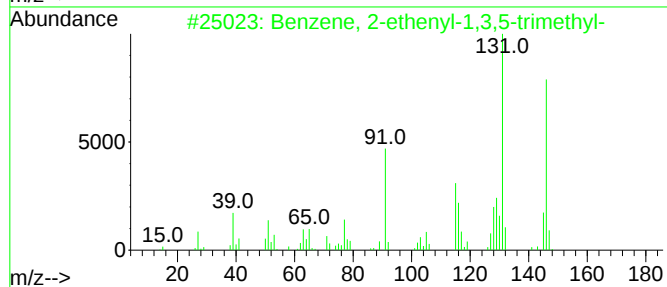
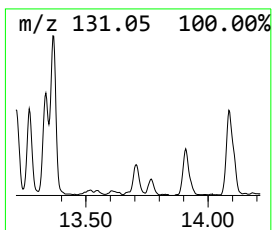
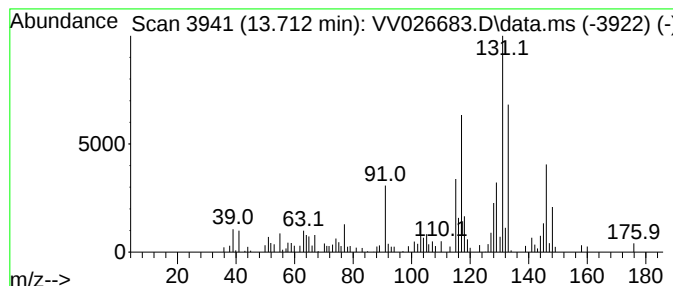
Instrument :
MSVOA_V
ClientSampleId :
DBUB5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 28 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.712	0.63 ug/L	94717	1,4-Dichlorobenzene-d4			11.249
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,3,5-trimethyl-		146	C11H14	000769-25-5	78
2	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001680-51-9	64
3	Benzene, (1-methyl-1-butenyl)-		146	C11H14	053172-84-2	55
4	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	55
5	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

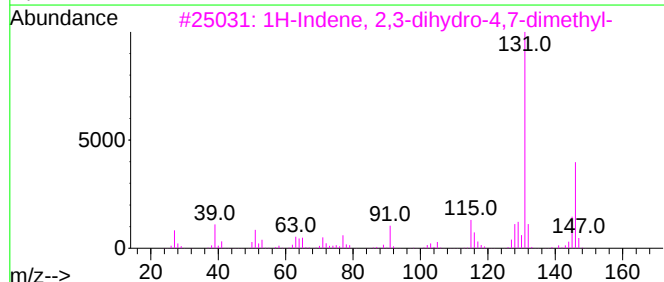
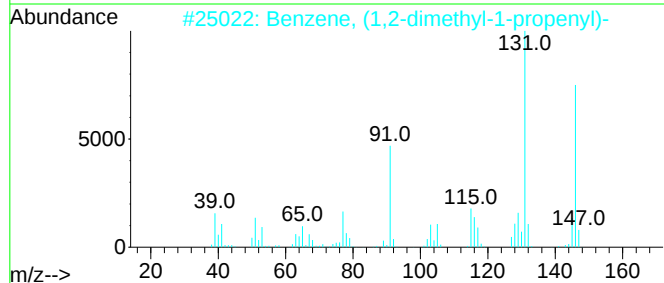
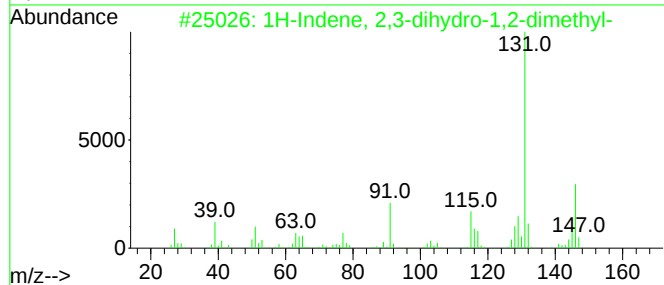
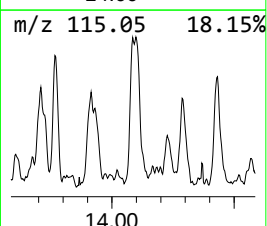
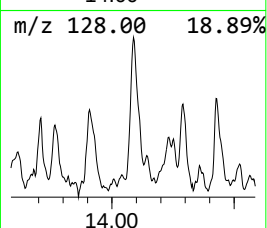
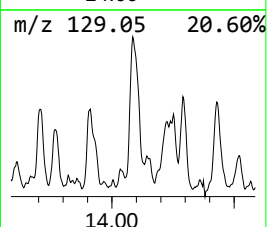
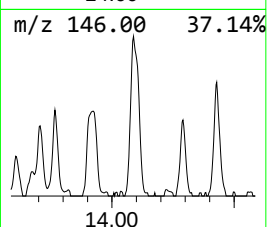
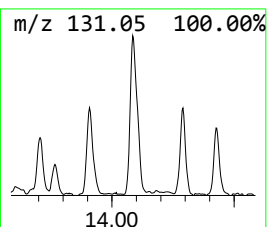
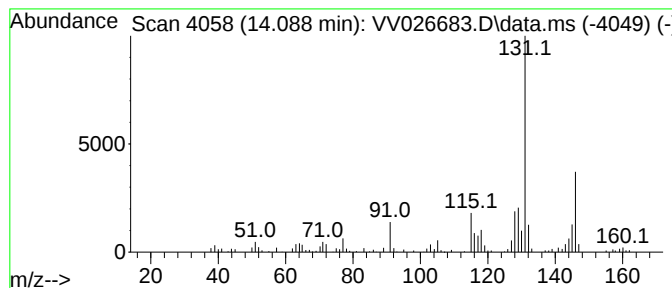
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 29 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.088	0.93 ug/L	138911	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93		
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93		
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87		
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83		
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

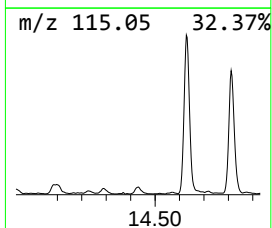
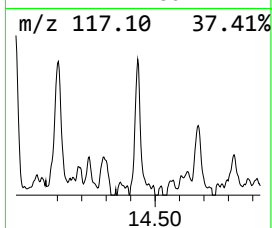
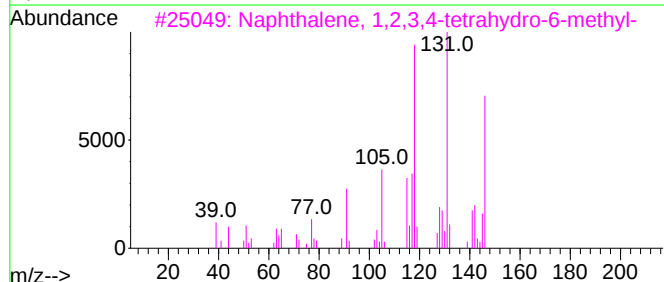
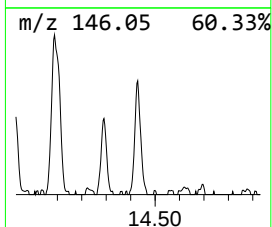
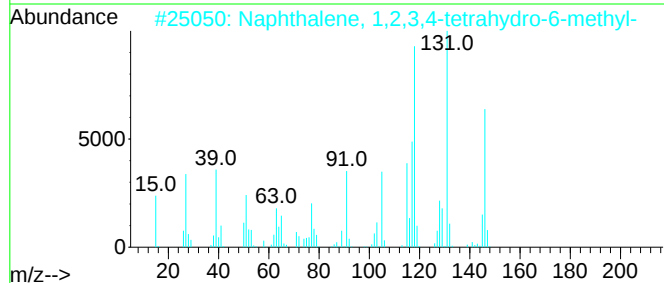
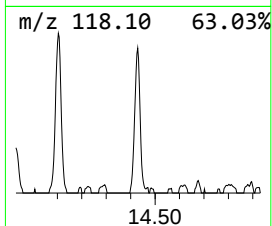
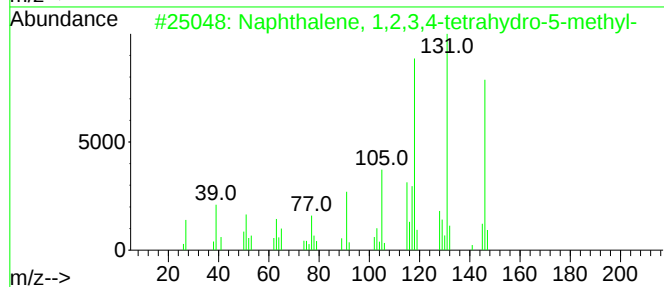
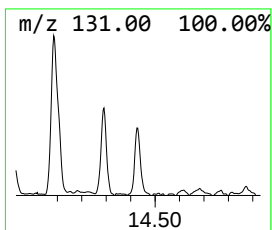
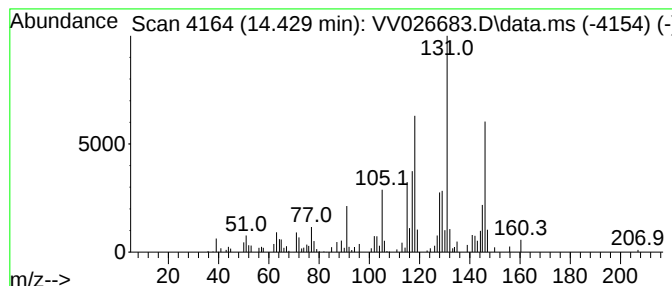
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 30 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.429	0.52 ug/L	77641	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

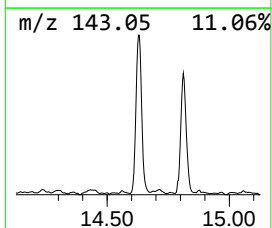
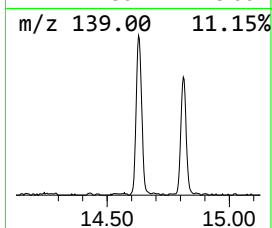
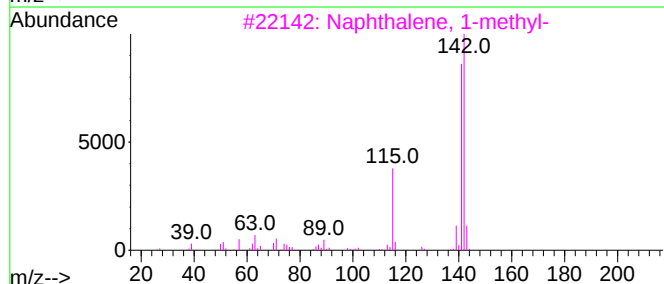
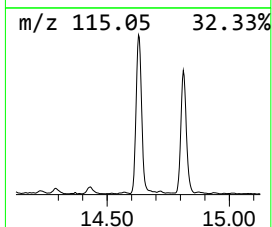
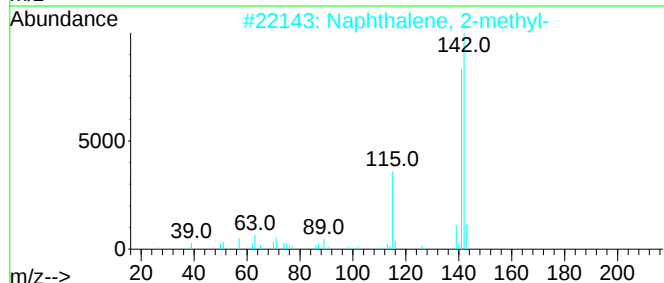
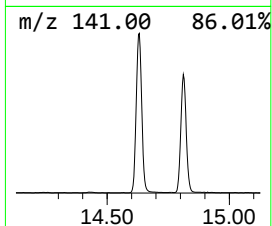
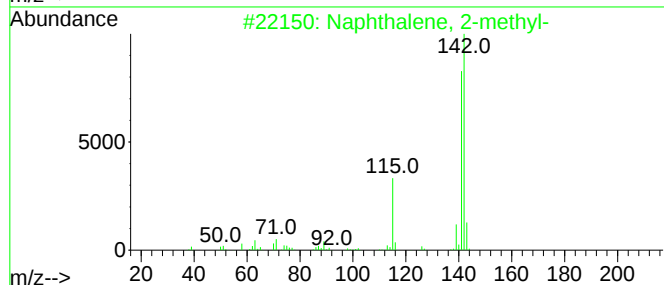
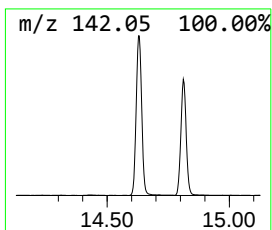
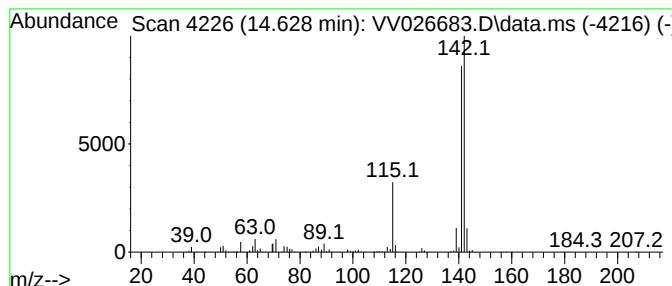
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 31 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.628	4.88 ug/L	731432	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
Data File : VV026683.D
Acq On : 07 Jul 2022 17:42
Operator : SY/MD
Sample : N3601-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBUB5

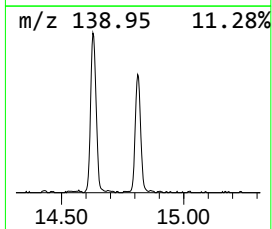
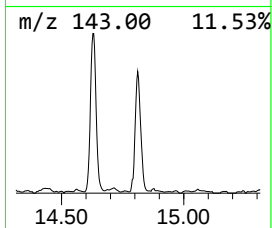
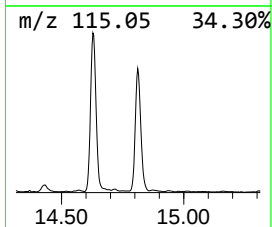
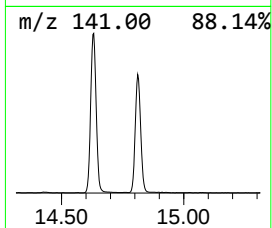
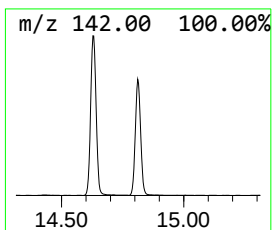
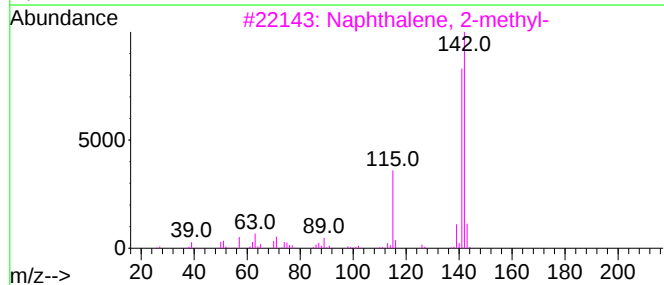
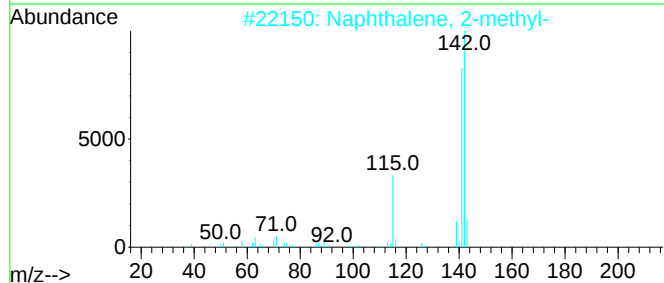
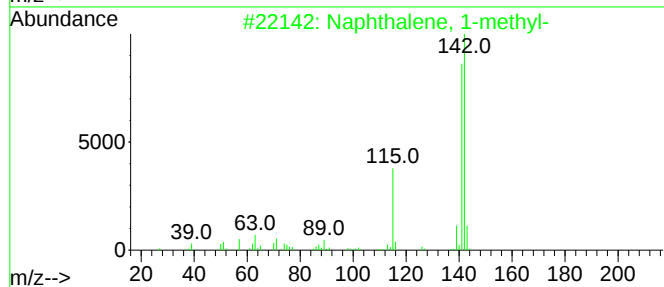
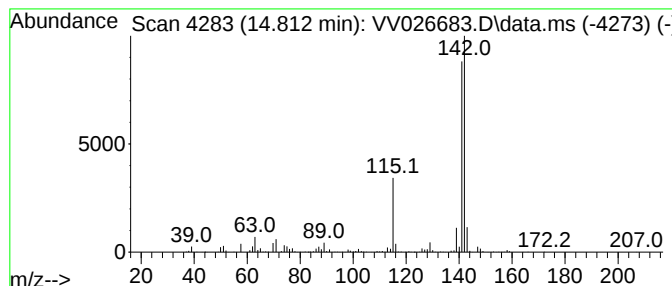
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 32 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.812	3.81 ug/L	570958	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV070722\
 Data File : VV026683.D
 Acq On : 07 Jul 2022 17:42
 Operator : SY/MD
 Sample : N3601-10
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBUB5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.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
Methanethiol	1.478	5.7	ug/L	483807	1	5.625	424109	5.0
Ethanethiol	2.044	0.5	ug/L	43084	1	5.625	424109	5.0
Dimethyl sulfide	2.220	1.1	ug/L	96441	1	5.625	424109	5.0
Thiophene, 3-me...	7.526	0.8	ug/L	320647	2	8.854	1973160	5.0
Benzene, propyl-	10.355	0.8	ug/L	122890	3	11.249	748841	5.0
Benzene, 1-ethy...	10.448	1.7	ug/L	260177	3	11.249	748841	5.0
Benzene, 1-ethy...	10.738	1.4	ug/L	212516	3	11.249	748841	5.0
7-Oxabicyclo[2....	11.085	1.4	ug/L	206167	3	11.249	748841	5.0
Benzene, 1,2,3-...	11.336	3.9	ug/L	586942	3	11.249	748841	5.0
Indane	11.532	1.6	ug/L	240840	3	11.249	748841	5.0
Benzene, 1-meth...	11.818	0.6	ug/L	94001	3	11.249	748841	5.0
Benzene, 2-ethy...	11.911	0.9	ug/L	137199	3	11.249	748841	5.0
Benzene, 1,2,3,...	11.940	1.0	ug/L	151186	3	11.249	748841	5.0
Benzene, 1-ethy...	12.014	1.5	ug/L	220608	3	11.249	748841	5.0
(E)-1-Phenyl-1-...	12.114	2.0	ug/L	295196	3	11.249	748841	5.0
Benzene, 4-ethy...	12.313	1.1	ug/L	166778	3	11.249	748841	5.0
Benzene, 1,2,4,...	12.410	1.5	ug/L	225170	3	11.249	748841	5.0
1,3-Cyclopentad...	12.464	2.5	ug/L	378461	3	11.249	748841	5.0
Benzene, 1-ethe...	12.725	1.4	ug/L	208839	3	11.249	748841	5.0
3-Phenylbut-1-ene	12.879	5.6	ug/L	841513	3	11.249	748841	5.0
Naphthalene, 1,...	13.043	1.7	ug/L	252010	3	11.249	748841	5.0
(+)-2-Bornanone	13.159	1.0	ug/L	152547	3	11.249	748841	5.0
1H-Indene, 2,3-...	13.268	0.9	ug/L	139835	3	11.249	748841	5.0
1H-Indene, 2,3-...	13.365	1.3	ug/L	191603	3	11.249	748841	5.0
Azulene	13.500	4.9	ug/L	736010	3	11.249	748841	5.0
Benzene, 2-ethe...	13.712	0.6	ug/L	94717	3	11.249	748841	5.0
Benzene, (1,2-d...	14.088	0.9	ug/L	138911	3	11.249	748841	5.0
Naphthalene, 1,...	14.429	0.5	ug/L	77641	3	11.249	748841	5.0
Naphthalene, 2-...	14.628	4.9	ug/L	731432	3	11.249	748841	5.0
Naphthalene, 1-...	14.812	3.8	ug/L	570958	3	11.249	748841	5.0