

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
 Data File : VV026459.D
 Acq On : 20 Jun 2022 20:37
 Operator : SY/MD
 Sample : N3355-01
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EY0C3

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\SFAMVTR061522WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV026459.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.365	1017	1034	1069	rBV2	110671	311586	1.07%	0.265%
2	4.777	1146	1162	1190	rVB	342631	876246	3.02%	0.745%
3	5.060	1234	1250	1275	rBV2	264570	639924	2.20%	0.544%
4	5.243	1296	1307	1324	rVB	300023	766811	2.64%	0.652%
5	5.625	1414	1426	1461	rBV	241160	535866	1.85%	0.456%
6	6.079	1553	1567	1574	rBV2	142721	301133	1.04%	0.256%
7	6.658	1732	1747	1766	rVB	181855	381024	1.31%	0.324%
8	6.780	1766	1785	1816	rBV2	226455	665614	2.29%	0.566%
9	7.320	1944	1953	1967	rVB	323191	558876	1.92%	0.475%
10	8.095	2179	2194	2215	rBV	204576	409507	1.41%	0.348%
11	8.854	2419	2430	2445	rBV	344018	559749	1.93%	0.476%
12	9.011	2468	2479	2499	rBV	606750	934256	3.22%	0.795%
13	9.140	2510	2519	2535	rVB	268967	441028	1.52%	0.375%
14	9.931	2752	2765	2787	rBV	828883	1271162	4.38%	1.081%
15	10.352	2880	2896	2913	rBV	2696347	3961126	13.64%	3.370%
16	10.448	2914	2926	2927	rVV	1170256	1324944	4.56%	1.127%
17	10.471	2928	2933	2944	rVV	1990257	2866890	9.87%	2.439%
18	10.538	2944	2954	2968	rVB	748102	1109200	3.82%	0.944%
19	10.735	3002	3015	3041	rBV	5843620	8773702	30.22%	7.464%
20	10.915	3058	3071	3087	rBV	19640653	29032787	100.00%	24.698%
21	11.053	3104	3114	3119	rBV	209066	320915	1.11%	0.273%
22	11.085	3119	3124	3138	rVB	276166	418825	1.44%	0.356%
23	11.191	3147	3157	3165	rBV	234486	356656	1.23%	0.303%
24	11.246	3165	3174	3190	rVV2	508802	808156	2.78%	0.687%
25	11.333	3190	3201	3215	rVV	3879530	5601274	19.29%	4.765%
26	11.529	3249	3262	3273	rBV3	2149942	5099777	17.57%	4.338%
27	11.625	3283	3292	3312	rVB4	1837026	3818933	13.15%	3.249%
28	11.744	3318	3329	3340	rVB2	266864	419397	1.44%	0.357%
29	11.818	3340	3352	3366	rBV	1068338	1605621	5.53%	1.366%
30	11.908	3369	3380	3385	rVV	2254052	3345805	11.52%	2.846%
31	11.937	3385	3389	3401	rVV	2034714	2820813	9.72%	2.400%
32	12.014	3401	3413	3433	rVV	3970062	6304503	21.72%	5.363%
33	12.114	3433	3444	3476	rVB2	1737252	3319787	11.43%	2.824%
34	12.313	3495	3506	3519	rVB	728712	1144589	3.94%	0.974%
35	12.410	3519	3536	3544	rBV	2297207	3359236	11.57%	2.858%
36	12.464	3544	3553	3569	rVB	2921718	4343260	14.96%	3.695%

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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\SFAMVTR061522WMA.M
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37	12.548	3569	3579	3585	rBV	309610	465367	1.60%	0.396%
38	12.583	3585	3590	3602	rVB2	225205	324942	1.12%	0.276%
39	12.722	3624	3633	3648	rVB	1681625	2483863	8.56%	2.113%
40	12.879	3662	3682	3693	rBV2	3402675	5959615	20.53%	5.070%
41	12.998	3711	3719	3726	rBV	226208	322735	1.11%	0.275%
42	13.046	3726	3734	3754	rVB3	351543	651081	2.24%	0.554%
43	13.162	3758	3770	3777	rBV	330726	535770	1.85%	0.456%
44	13.214	3777	3786	3794	rVB	607709	916866	3.16%	0.780%
45	13.268	3794	3803	3809	rBV2	936828	1416899	4.88%	1.205%
46	13.365	3827	3833	3839	rVV	397786	522465	1.80%	0.444%
47	13.397	3839	3843	3856	rVB	319676	428777	1.48%	0.365%
48	13.500	3866	3875	3884	rBV	510704	736767	2.54%	0.627%
49	13.709	3929	3940	3950	rBV2	327928	627355	2.16%	0.534%
50	13.767	3950	3958	3968	rVB	306693	469757	1.62%	0.400%
51	13.908	3990	4002	4015	rBV	563092	848431	2.92%	0.722%
52	14.088	4047	4058	4072	rBV2	465290	1023095	3.52%	0.870%
53	14.291	4112	4121	4135	rVB3	251576	447705	1.54%	0.381%
54	14.812	4274	4283	4295	rVB	368351	562229	1.94%	0.478%

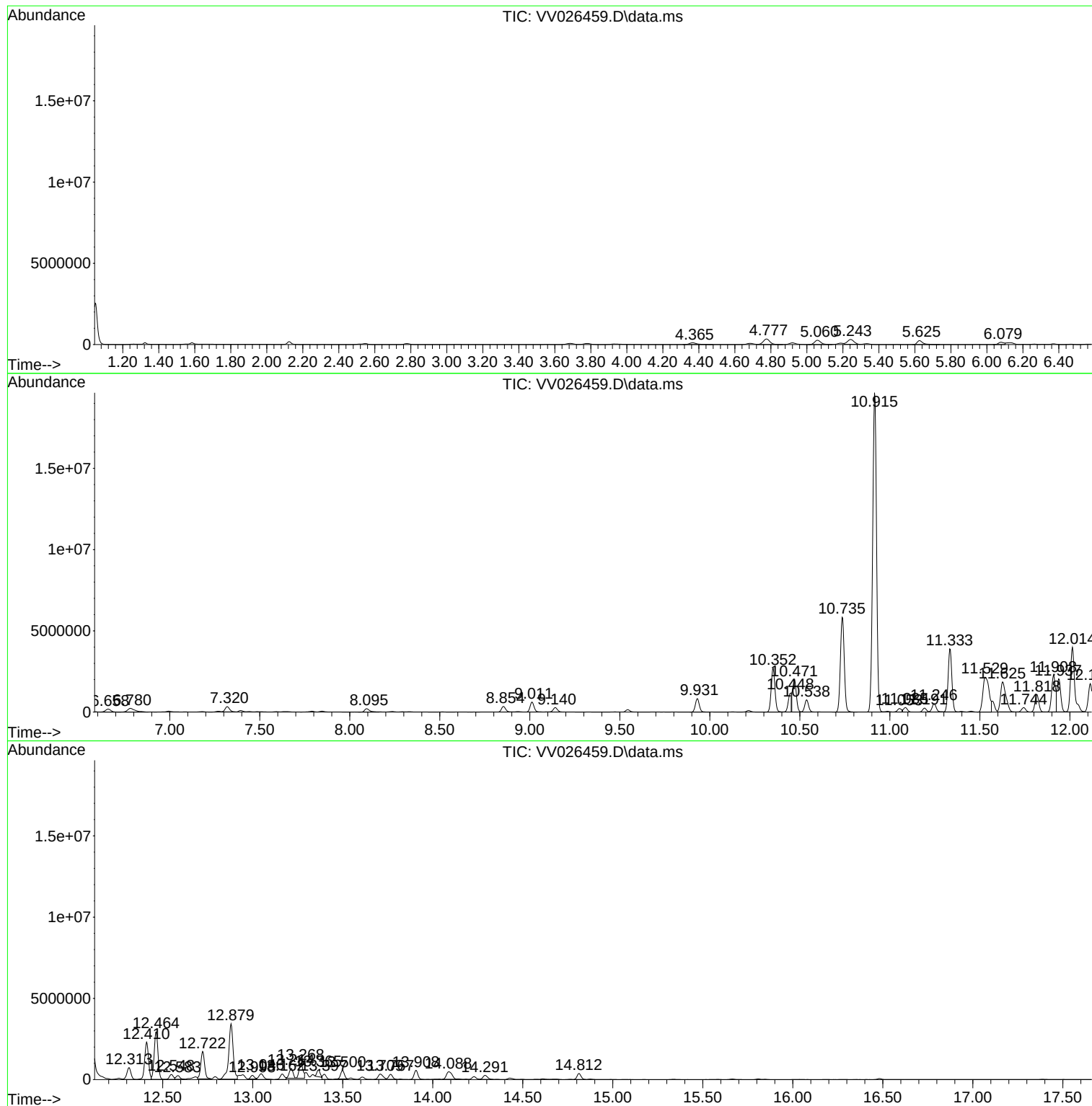
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EY0C3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR061522WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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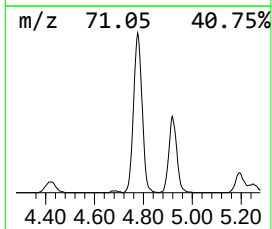
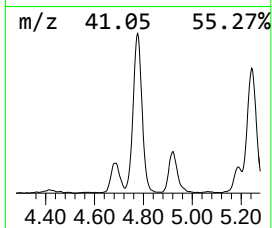
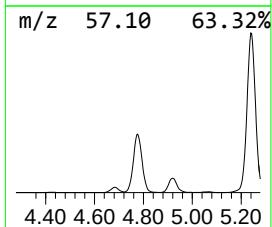
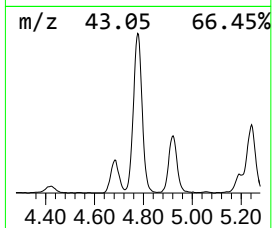
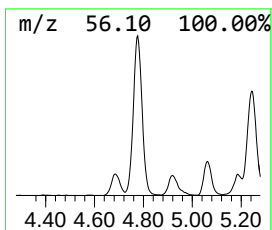
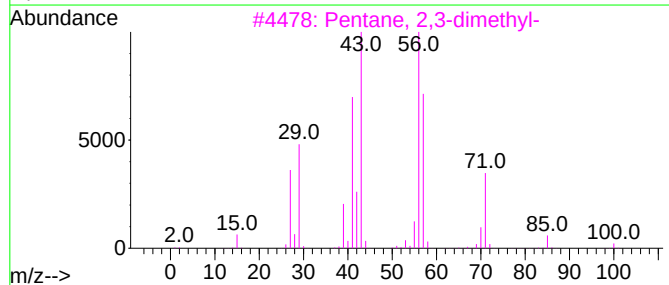
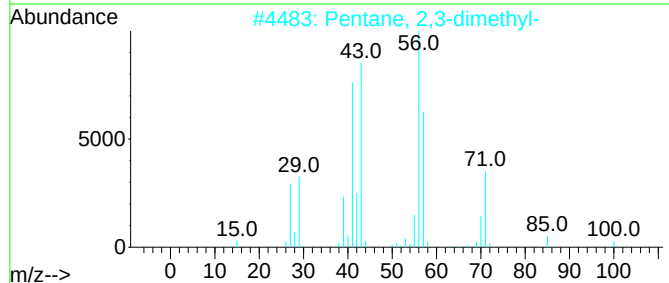
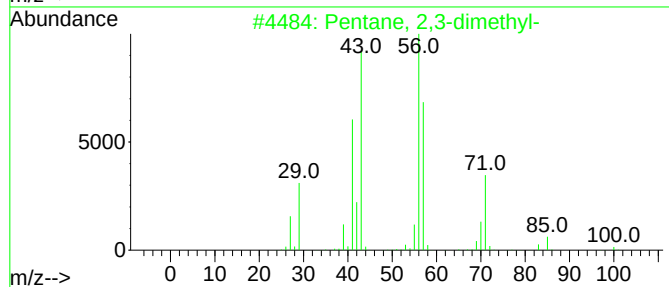
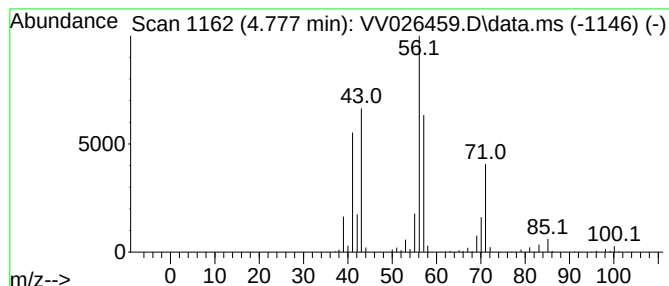
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.777	8.18 ug/L	876246	1,4-Difluorobenzene	5.625

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
3		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	78
4		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	72
5		Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	64



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ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

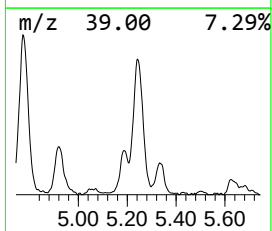
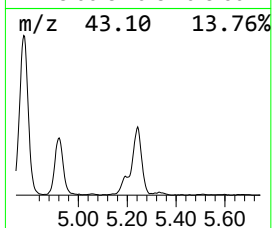
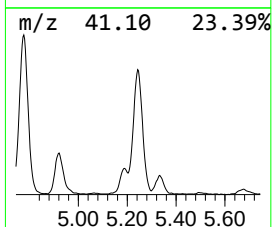
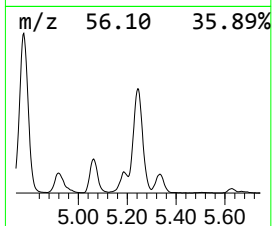
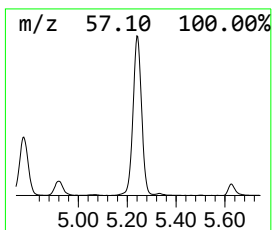
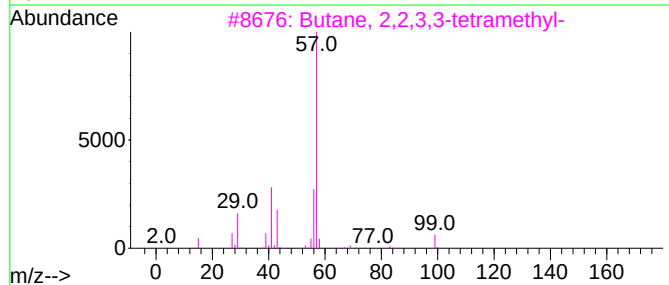
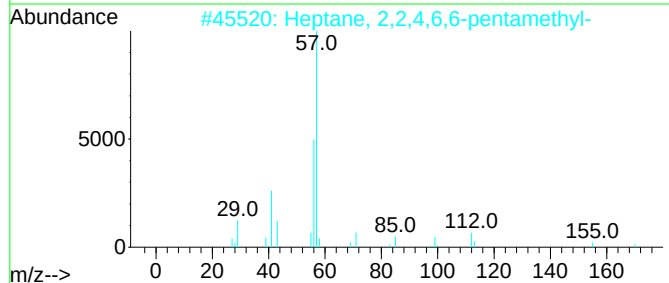
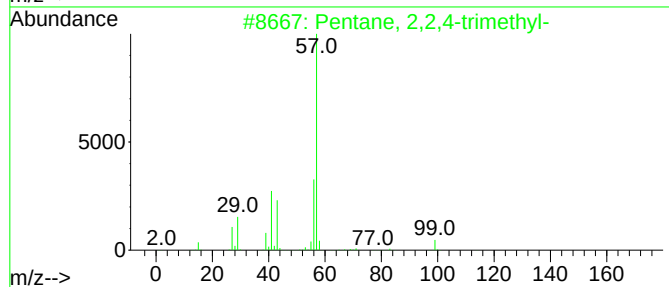
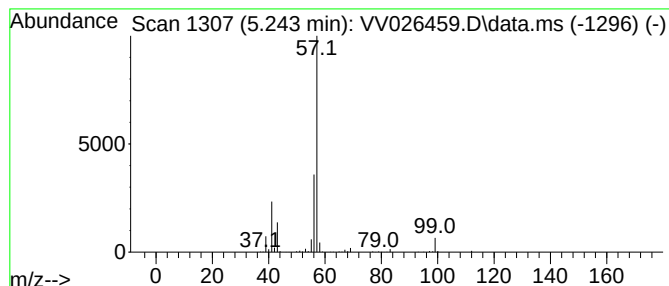
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR061522WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.243	7.15 ug/L	766811	1,4-Difluorobenzene	5.625

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	72
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
4			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
5			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72



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Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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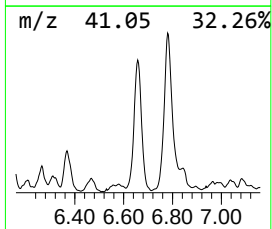
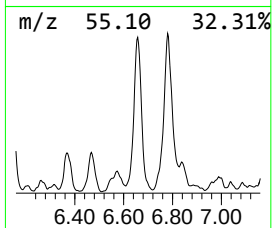
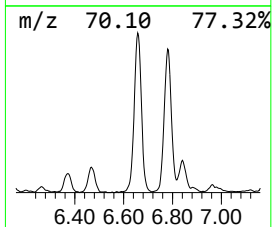
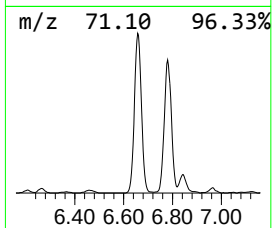
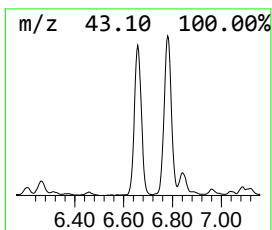
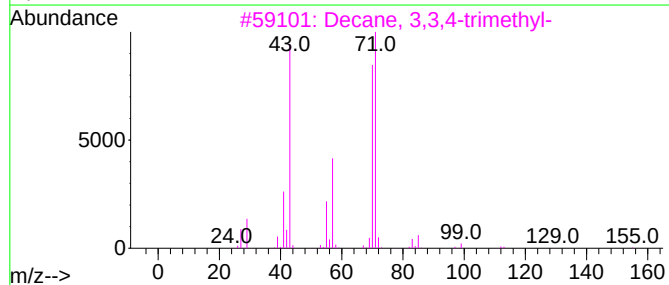
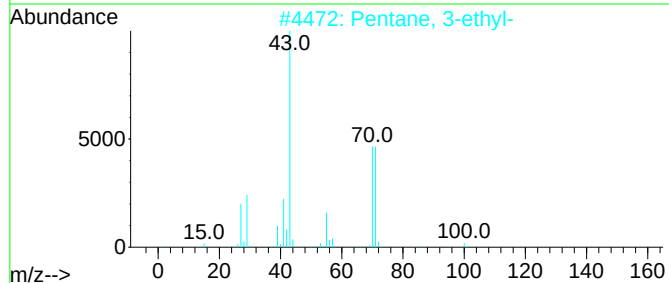
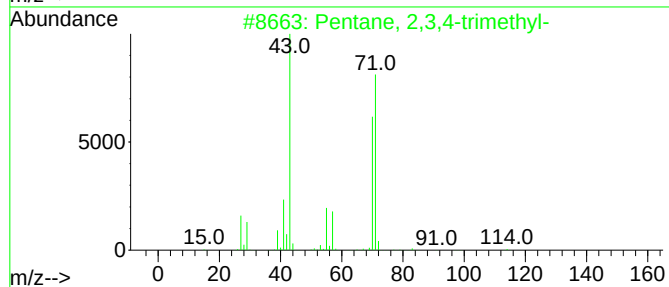
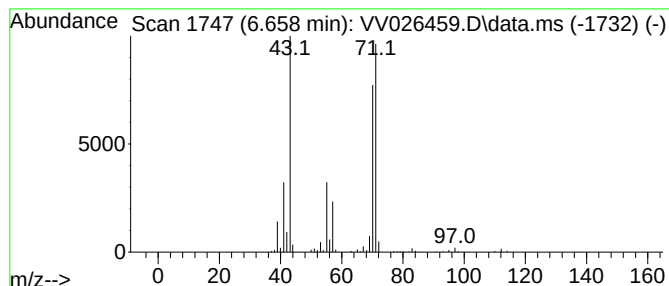
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR061522WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.658	3.56 ug/L	381024	1,4-Difluorobenzene	5.625

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	83
4		Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74



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Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

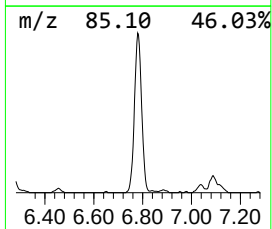
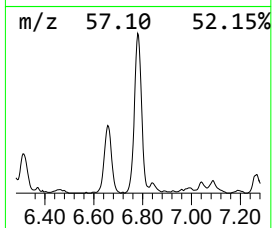
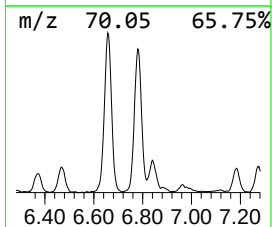
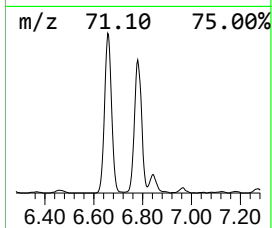
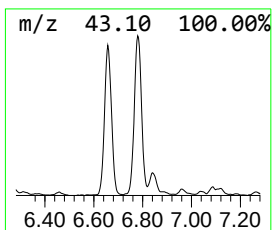
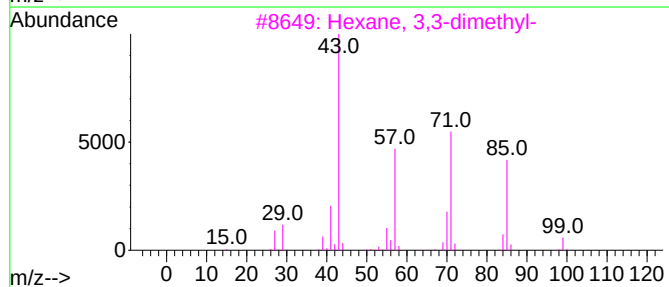
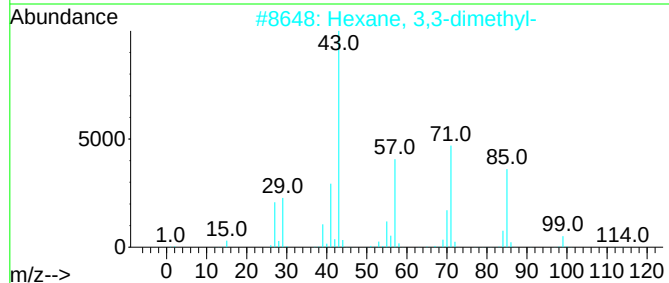
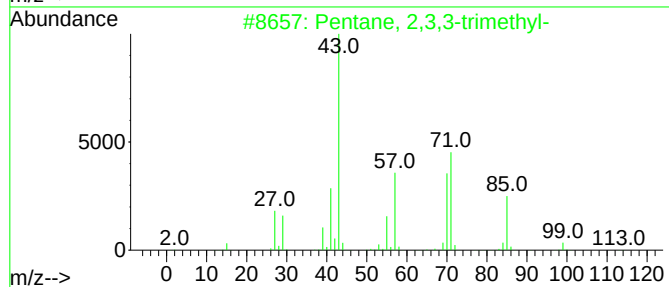
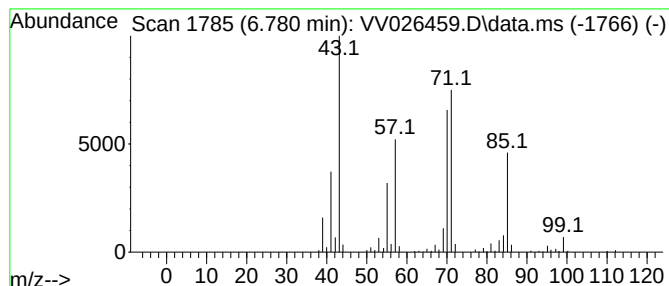
Instrument :
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.780	6.21 ug/L	665614	1,4-Difluorobenzene	5.625	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	72
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59



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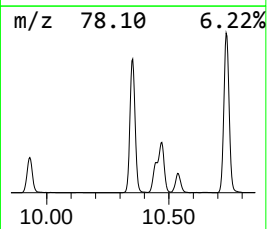
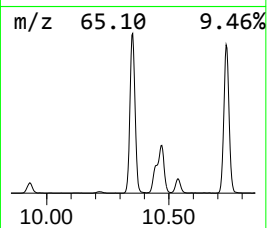
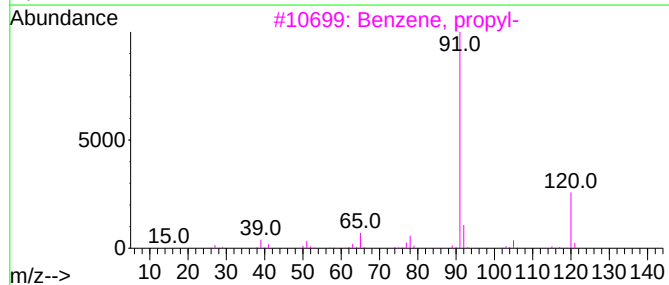
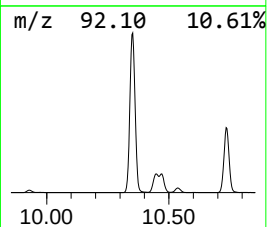
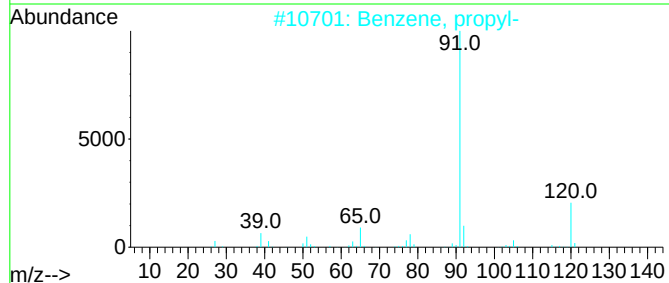
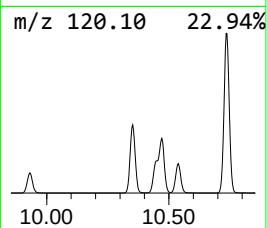
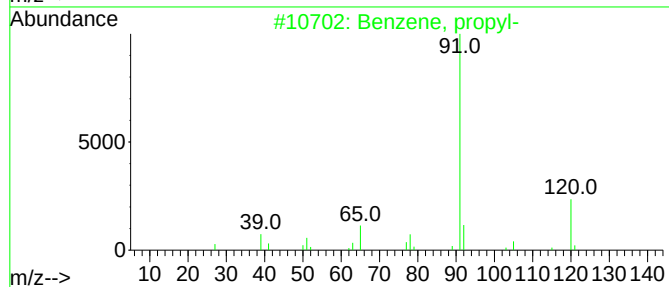
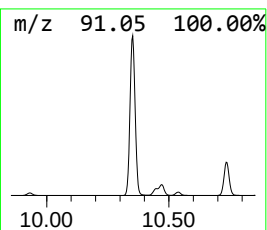
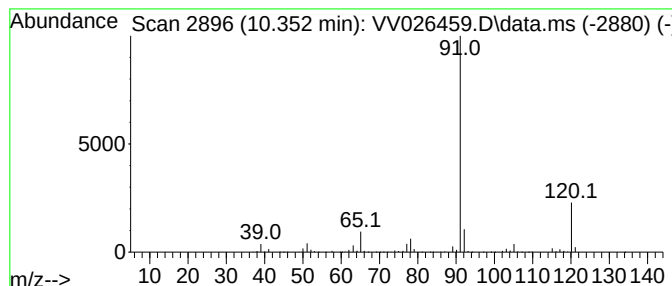
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.352	24.51 ug/L	3961130	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

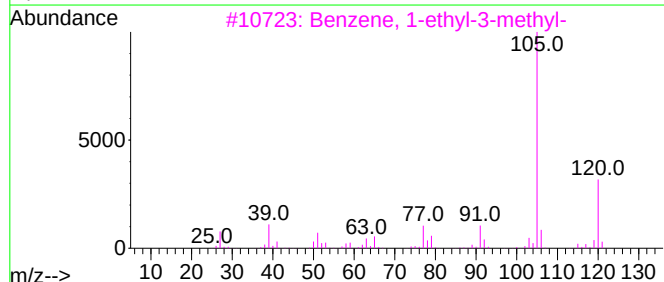
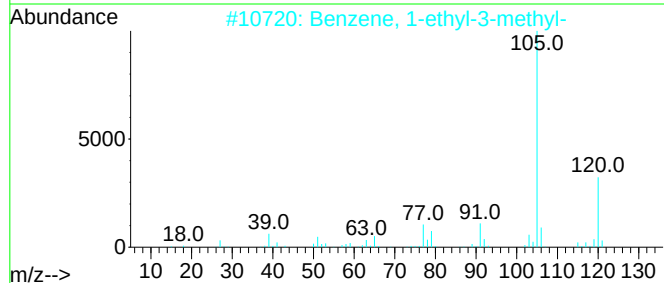
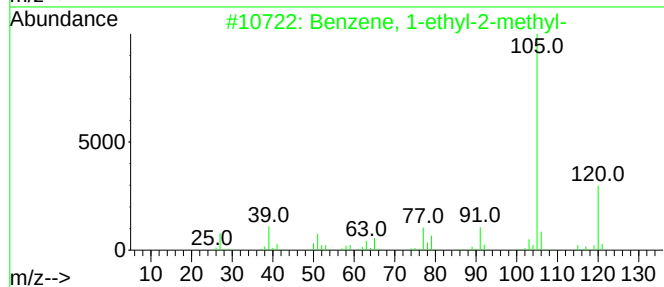
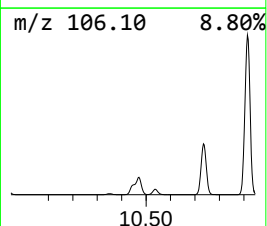
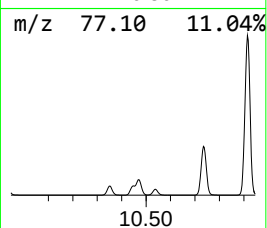
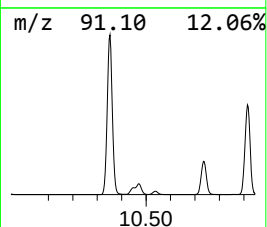
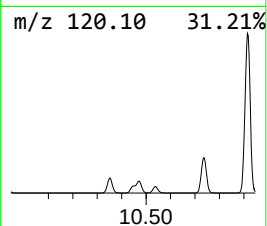
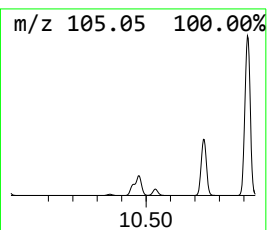
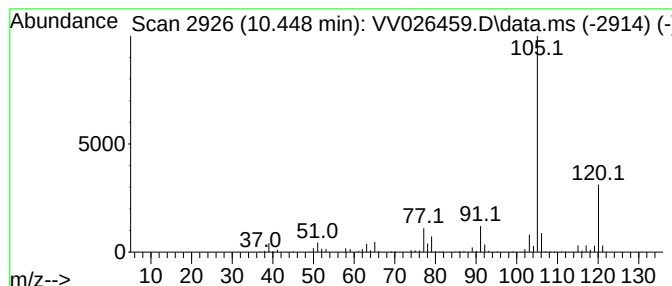
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	8.20 ug/L	1324940	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			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

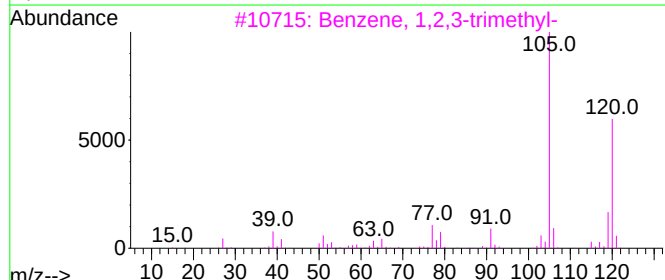
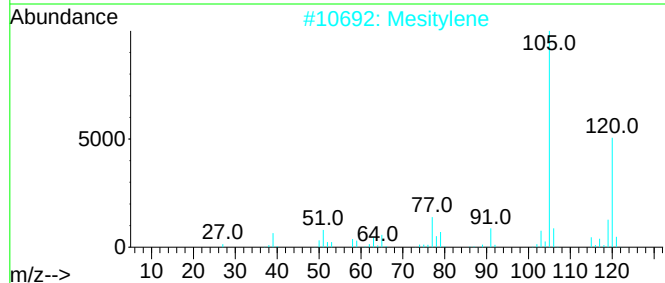
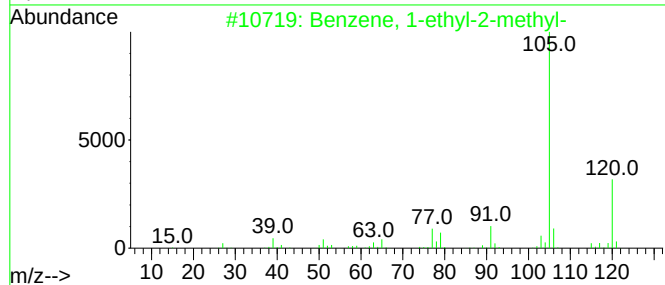
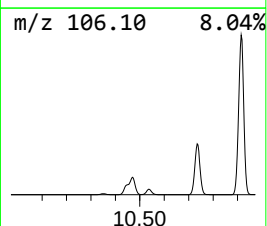
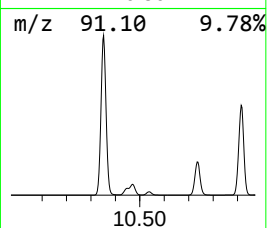
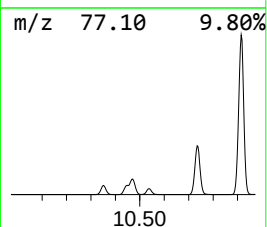
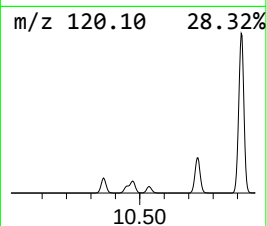
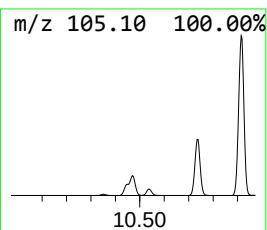
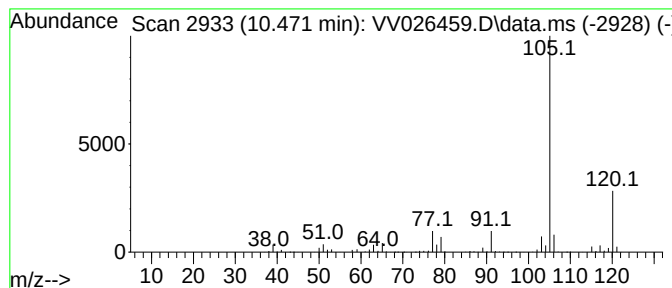
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.471	17.74 ug/L	2866890	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	94
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

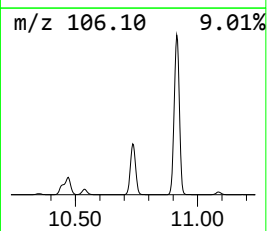
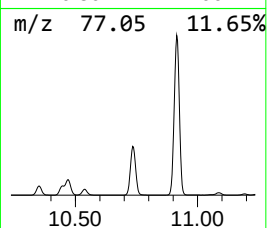
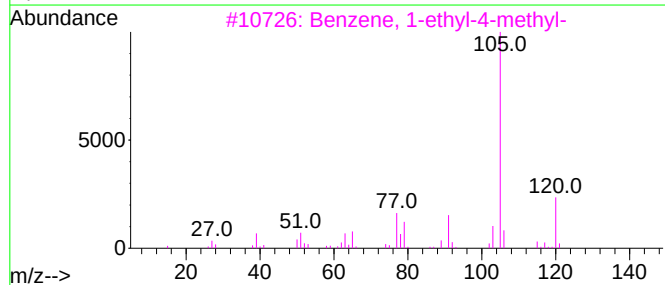
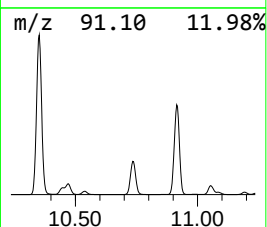
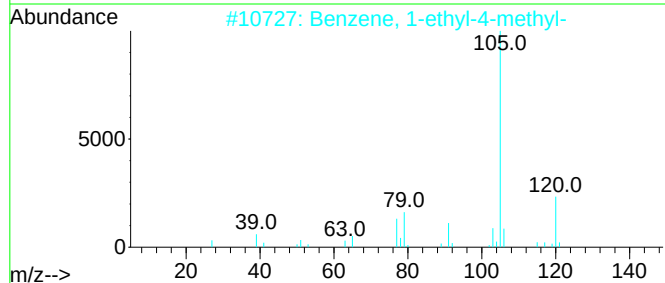
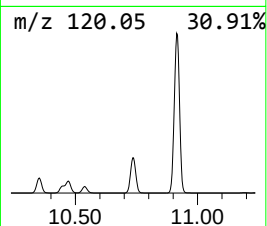
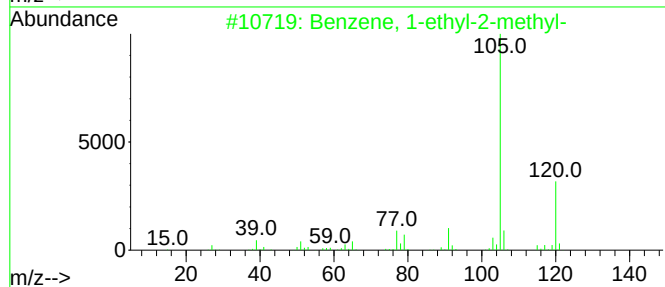
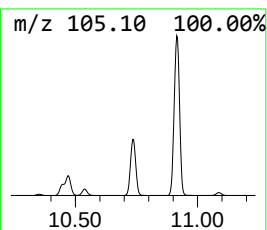
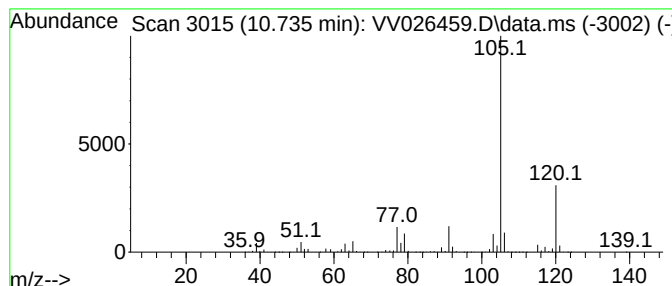
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.735	54.28 ug/L	8773700	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	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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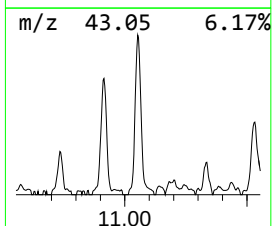
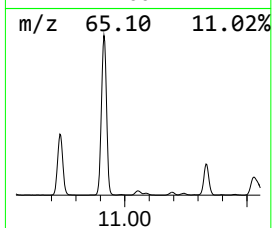
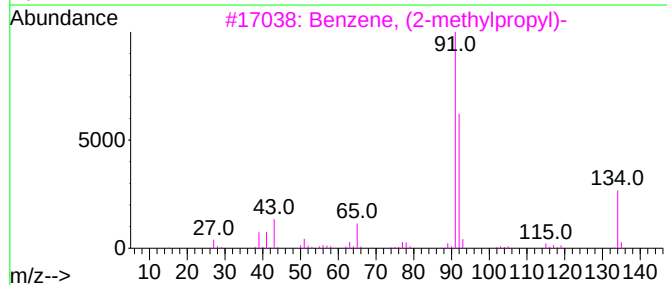
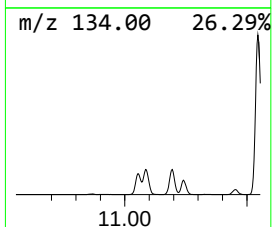
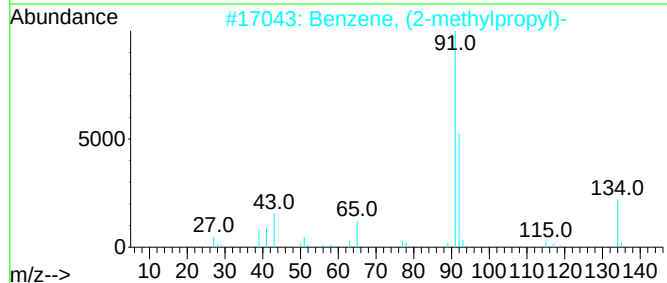
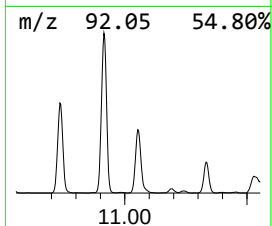
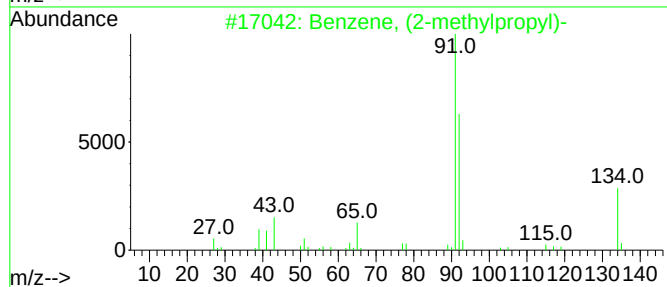
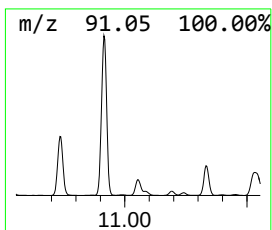
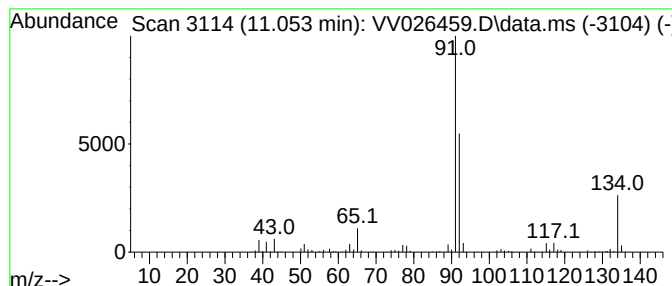
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, (2-methylpropyl)- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.053	1.99 ug/L	320915	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
4			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87
5			Benzene, n-butyl-	134	C10H14	000104-51-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

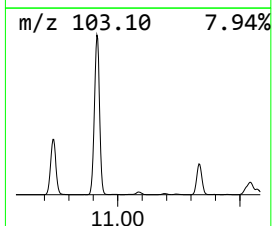
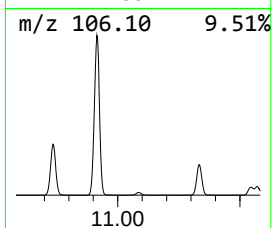
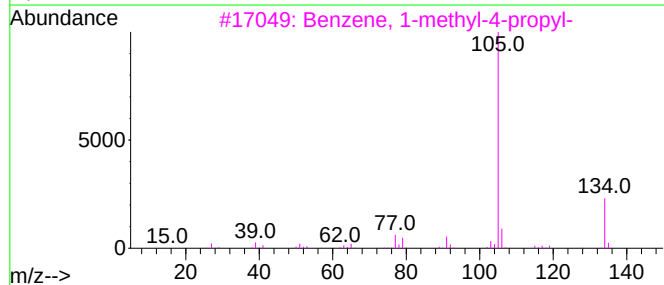
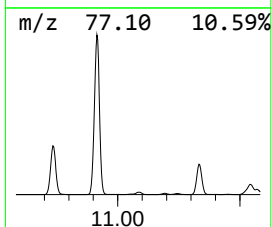
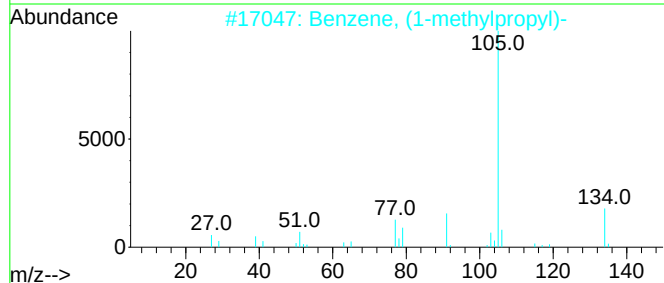
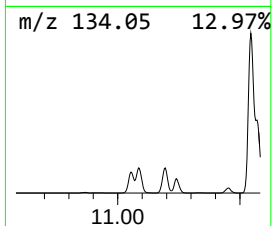
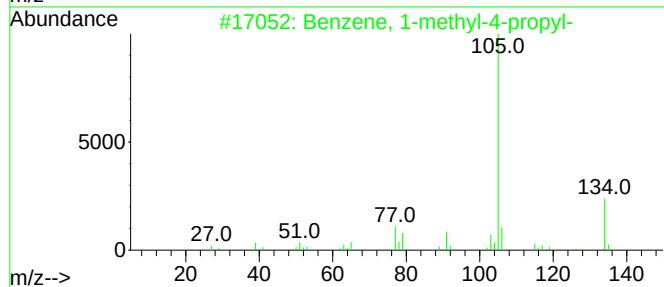
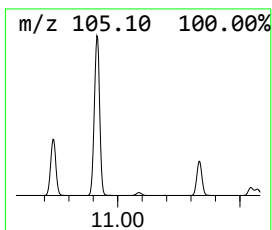
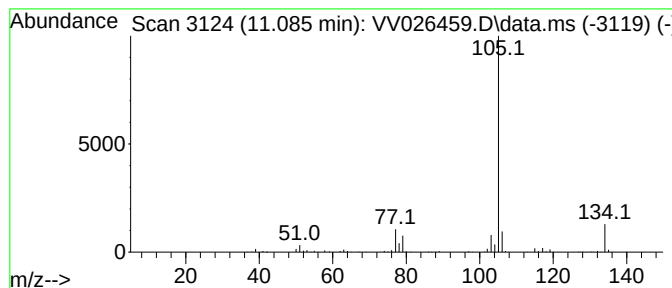
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 1-methyl-4-propyl- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.085	2.59 ug/L	418825	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	90
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	86
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	80
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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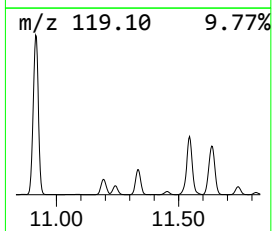
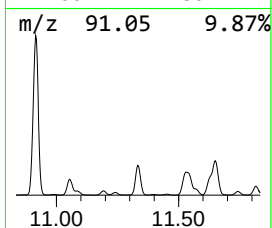
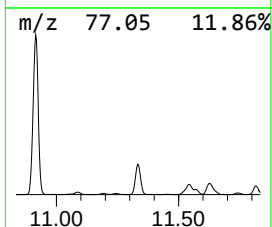
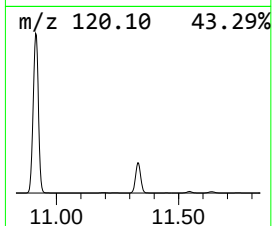
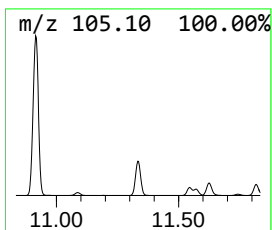
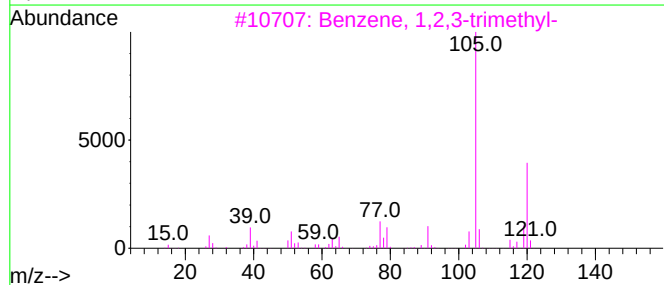
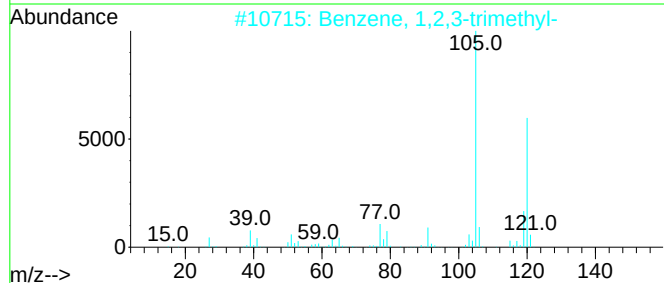
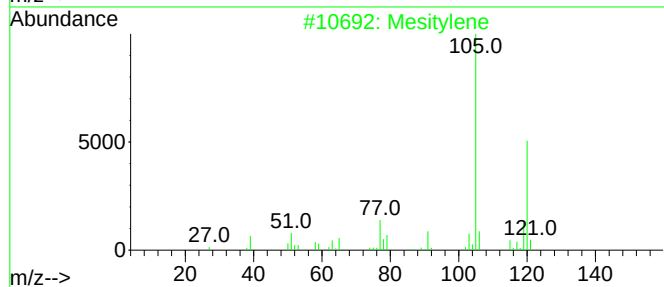
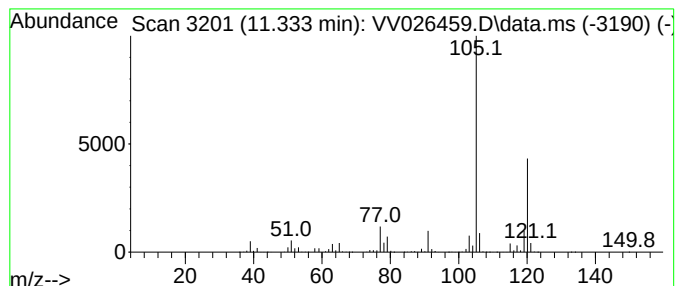
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.333	34.65 ug/L	5601270	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

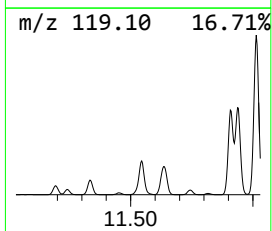
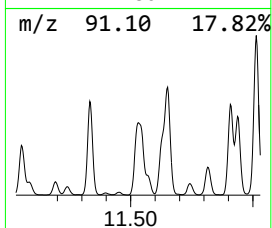
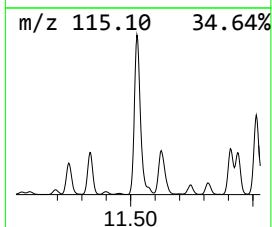
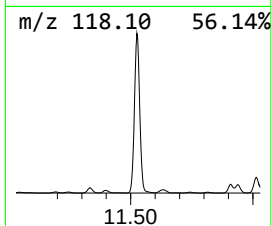
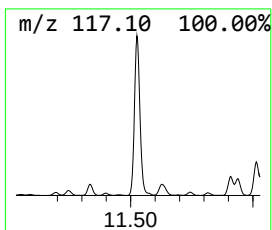
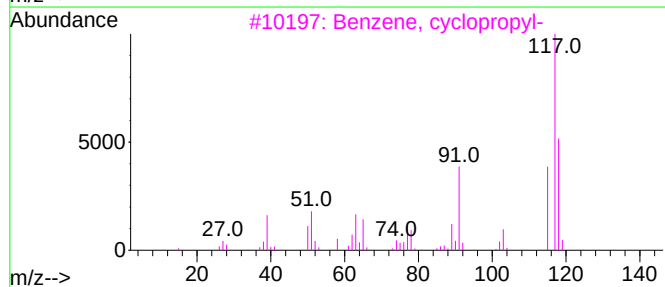
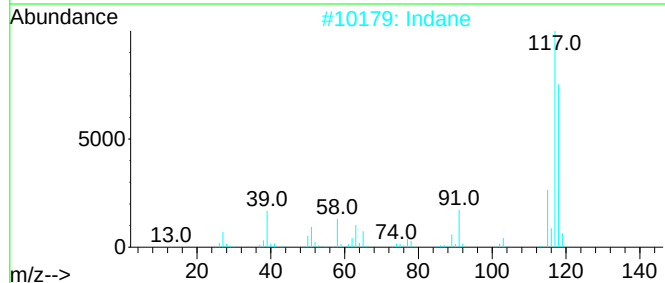
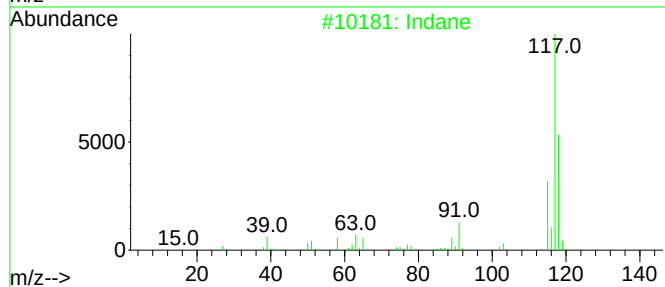
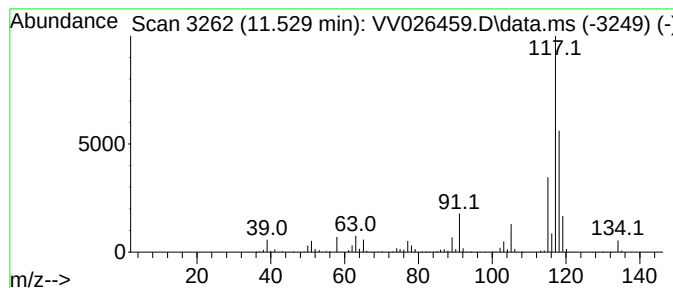
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	31.55 ug/L	5099780	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

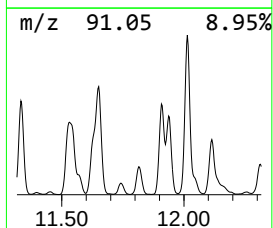
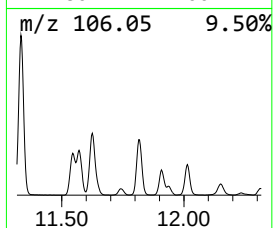
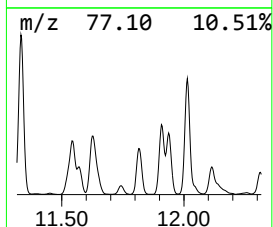
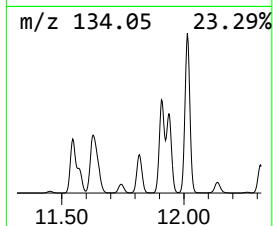
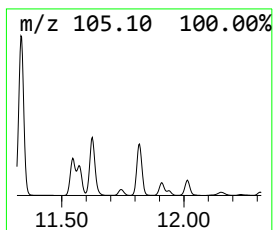
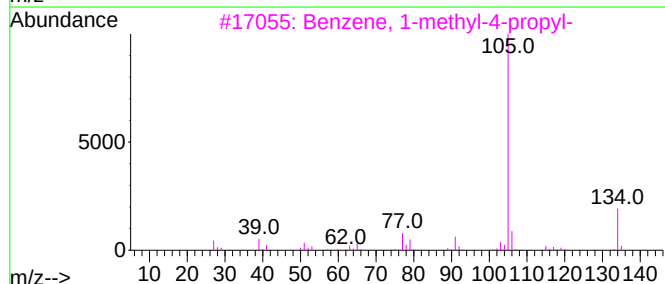
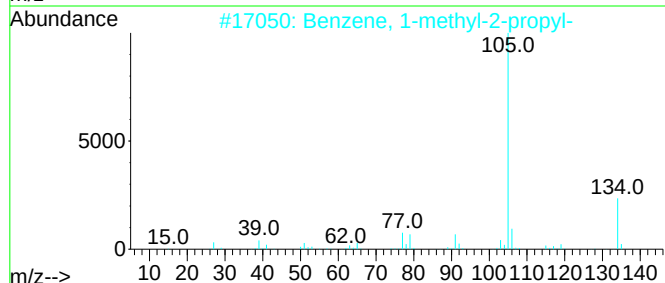
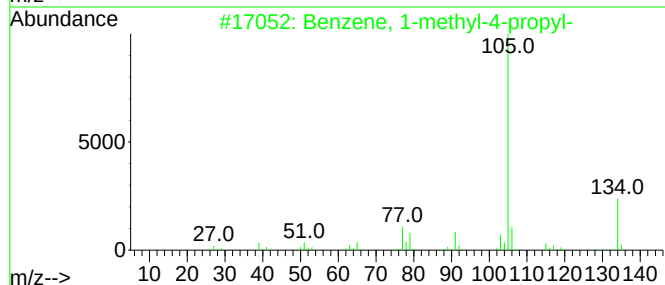
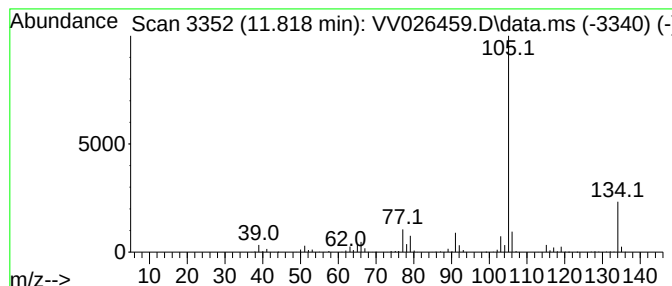
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 15 Benzene, 1-methyl-2-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.818	9.93 ug/L	1605620	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-2-propyl-	134	C10H14	001074-17-5	93
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
4			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

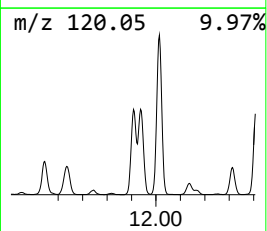
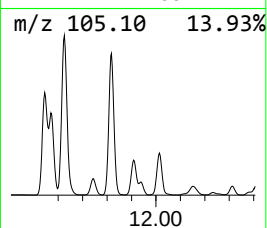
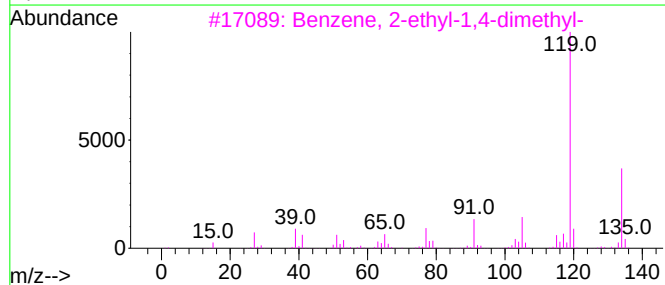
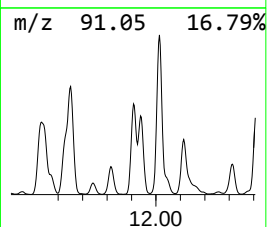
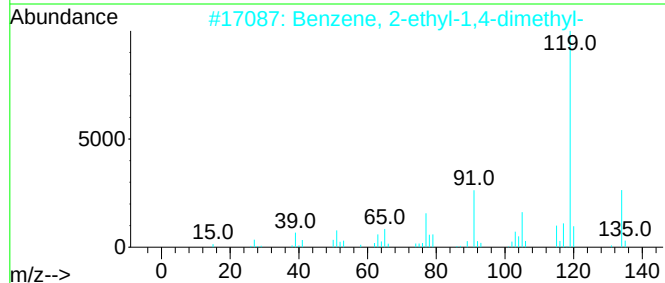
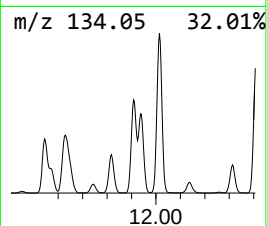
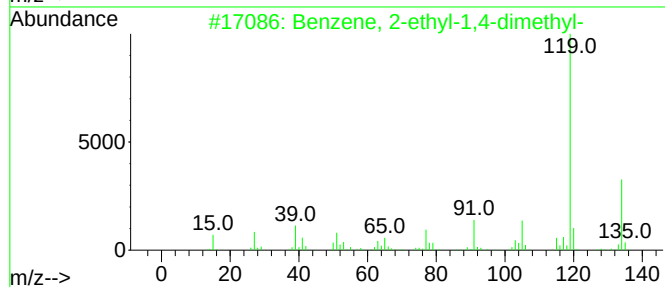
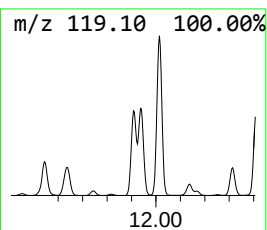
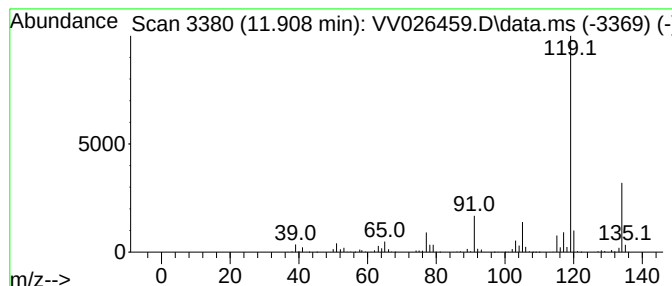
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.908	20.70 ug/L	3345810	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	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

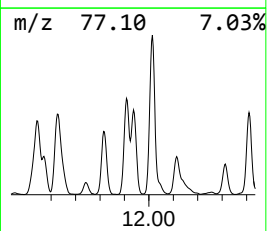
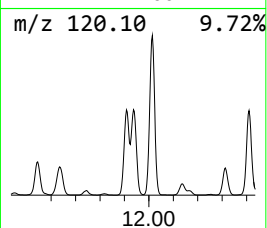
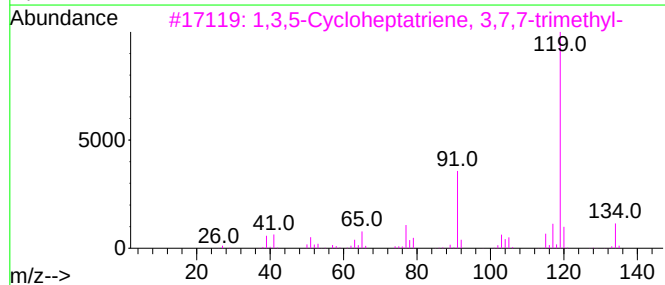
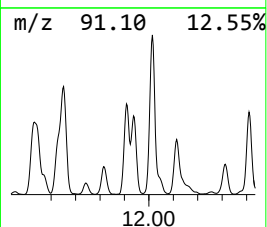
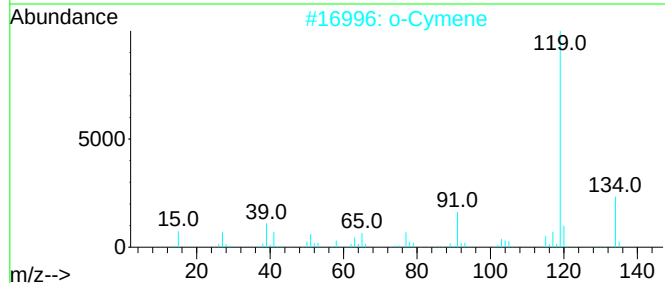
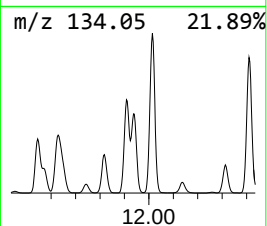
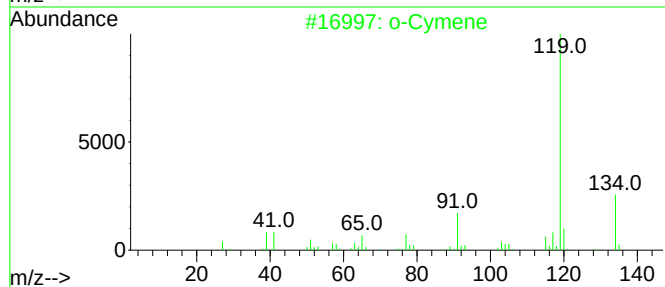
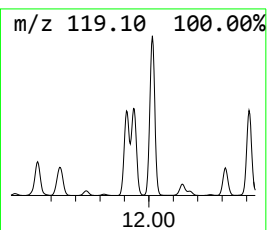
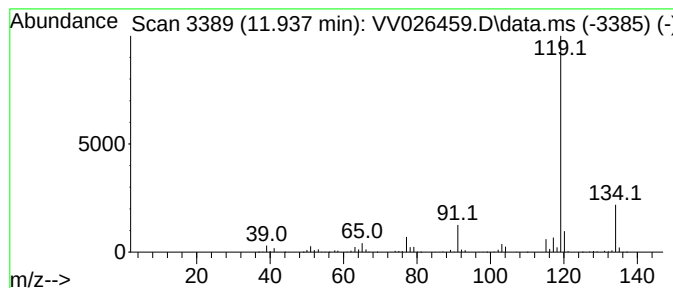
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 17 o-Cymene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.937	17.45 ug/L	2820810	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	94
2			o-Cymene	134	C10H14	000527-84-4	94
3			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	91
4			p-Cymene	134	C10H14	000099-87-6	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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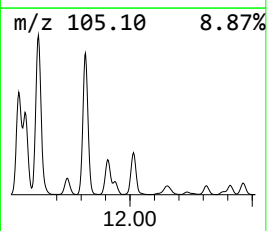
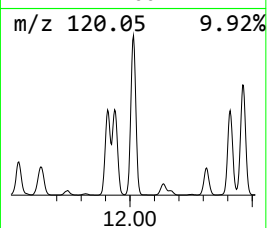
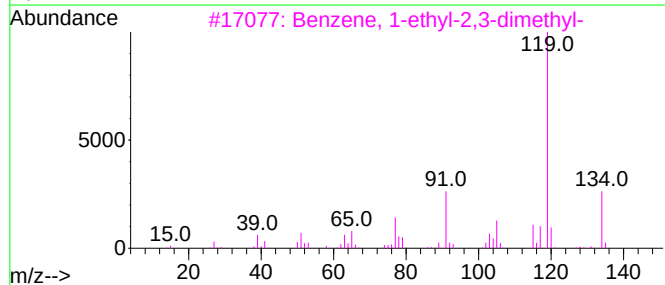
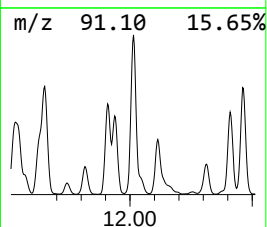
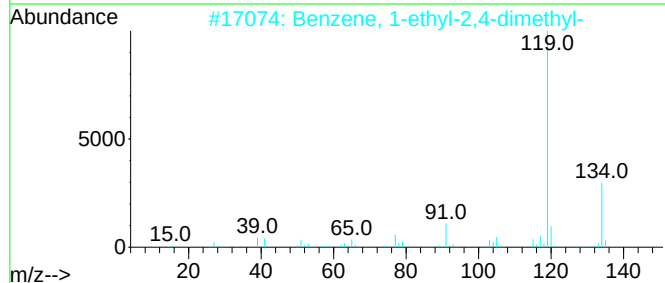
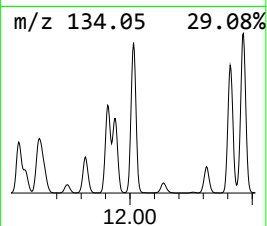
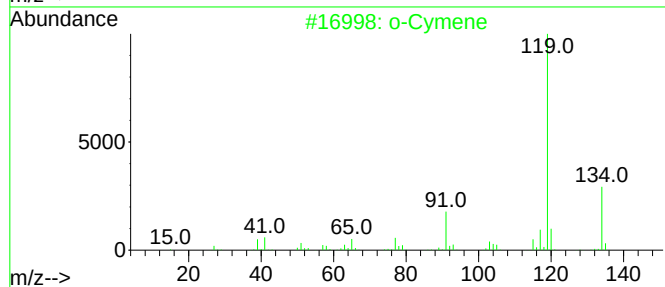
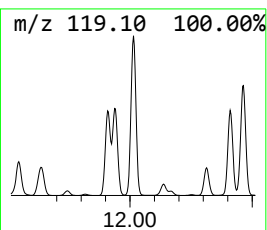
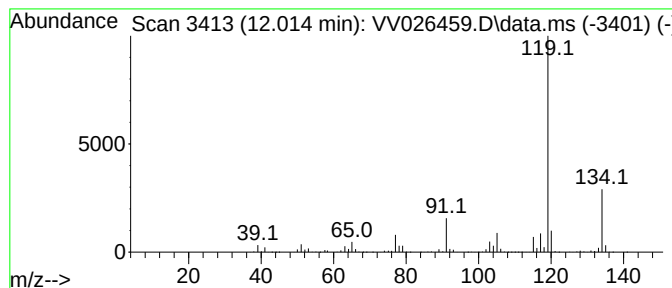
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 18 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.014	39.01 ug/L	6304500	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

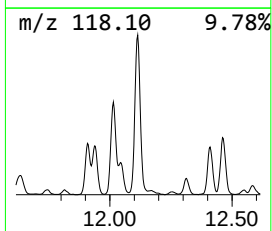
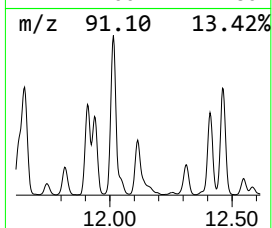
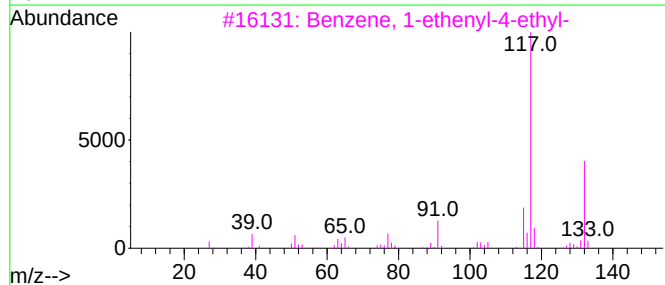
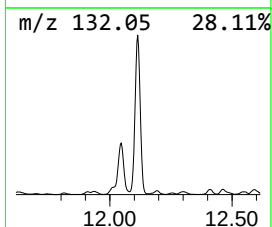
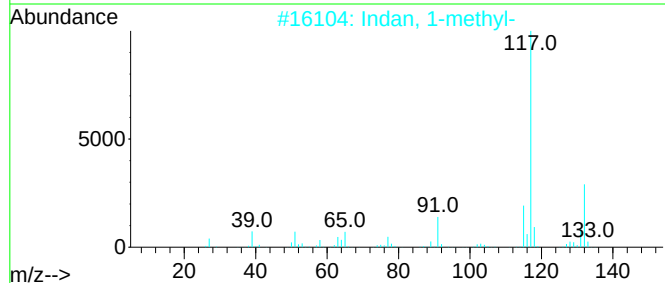
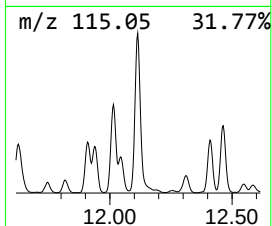
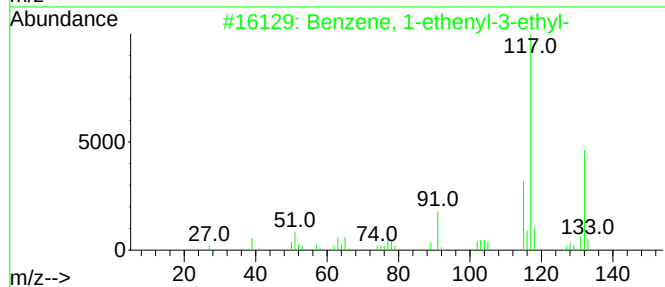
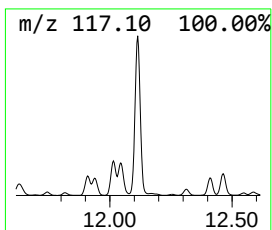
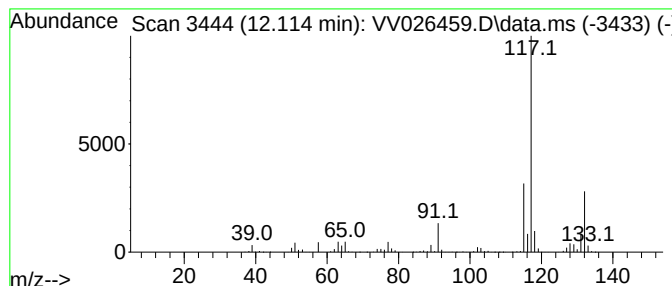
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 19 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.114	20.54 ug/L	3319790	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	91
2			Indan, 1-methyl-	132	C10H12	000767-58-8	90
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

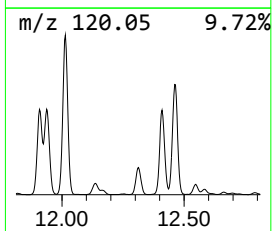
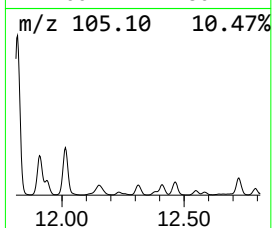
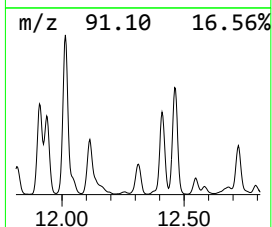
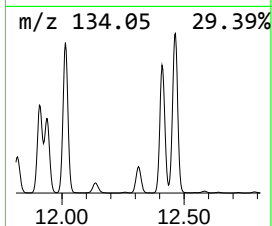
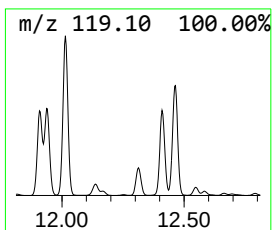
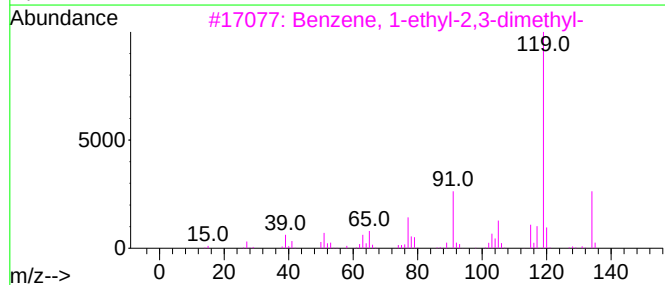
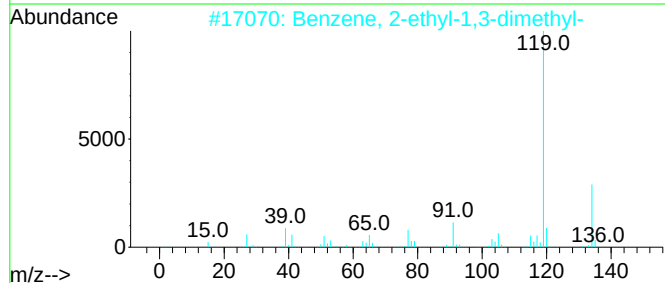
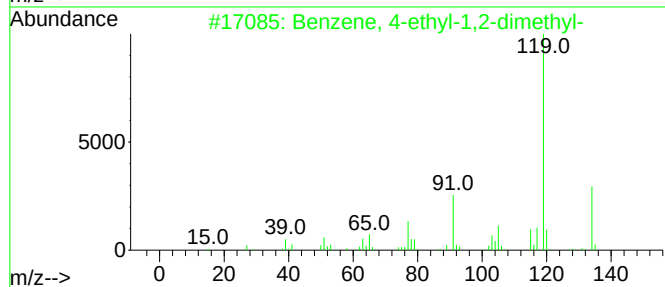
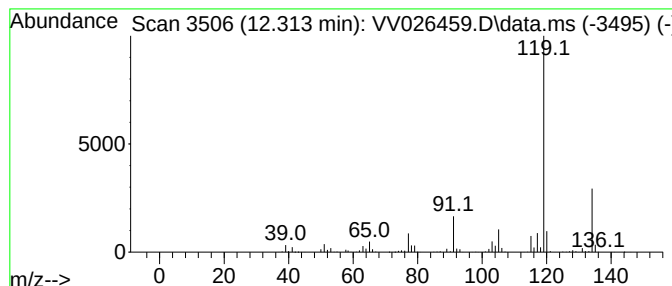
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 20 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.313	7.08 ug/L	1144590	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, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

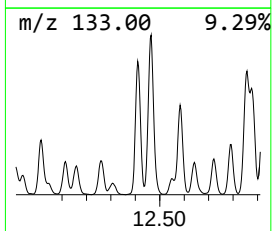
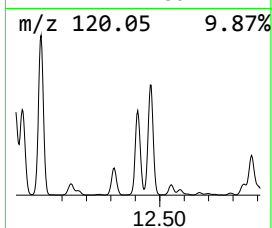
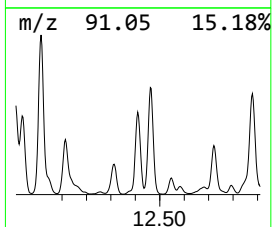
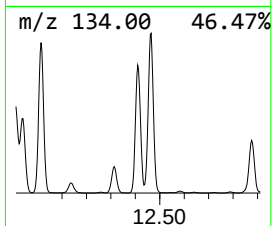
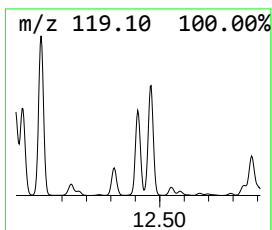
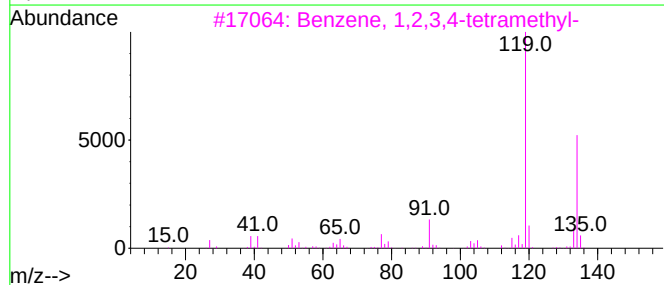
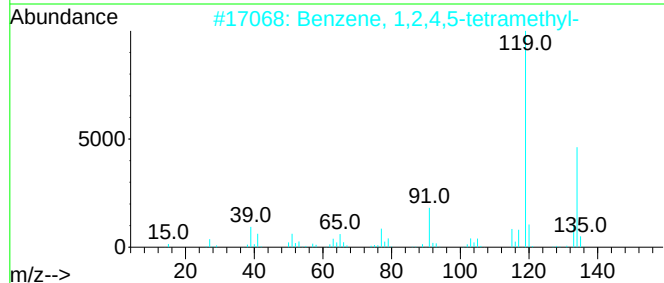
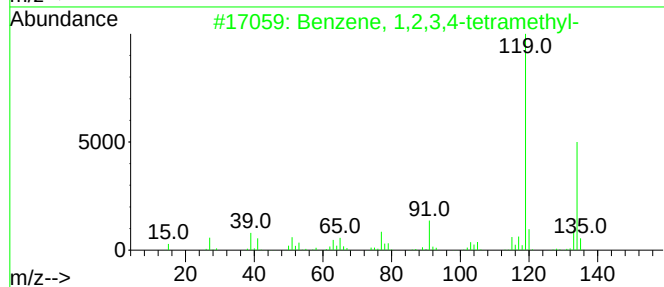
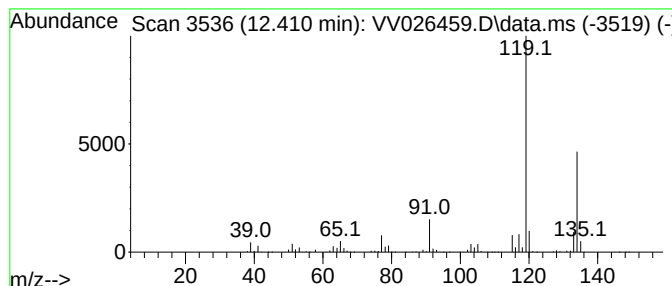
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 21 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	20.78 ug/L	3359240	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

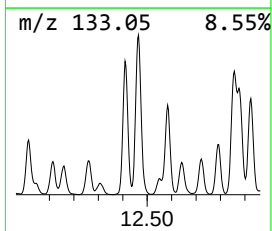
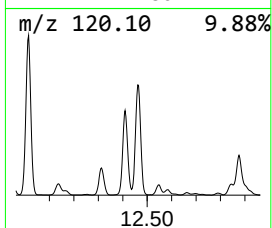
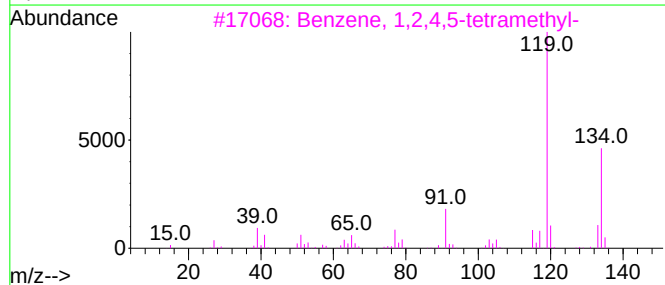
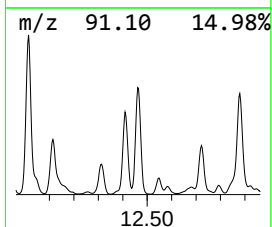
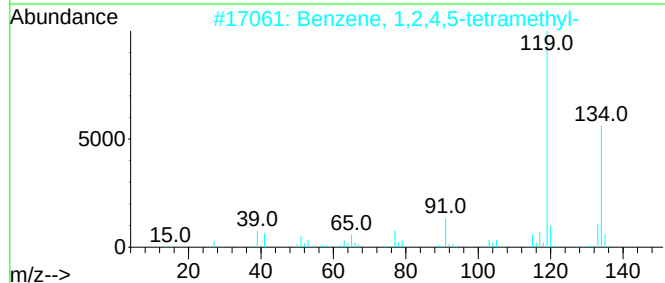
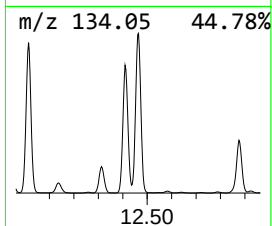
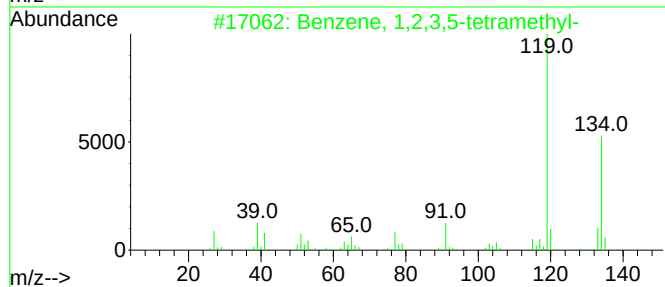
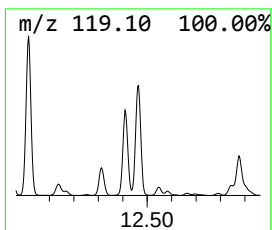
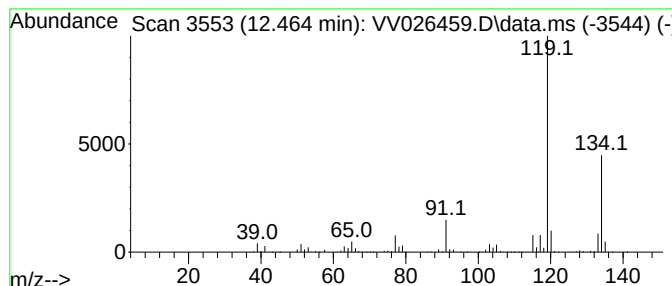
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 22 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	26.87 ug/L	4343260	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	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\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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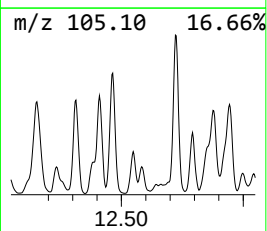
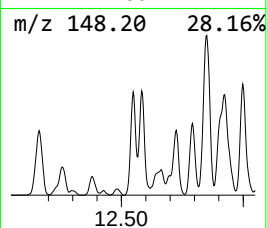
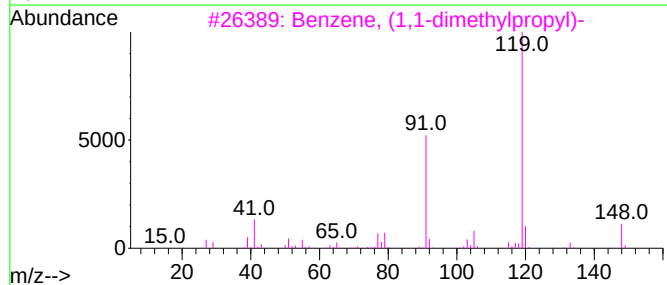
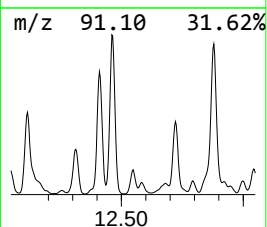
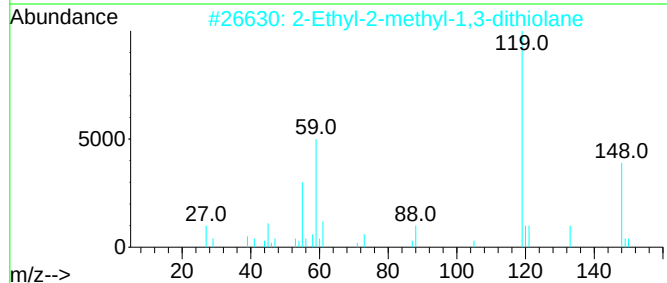
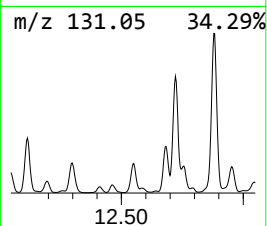
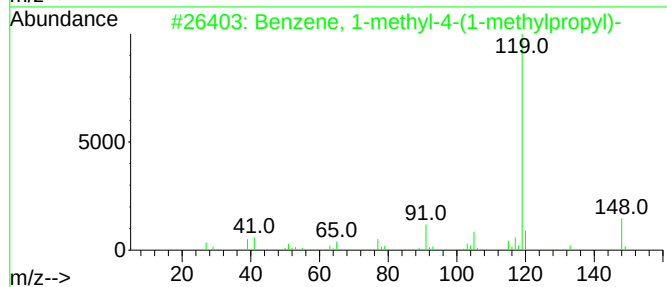
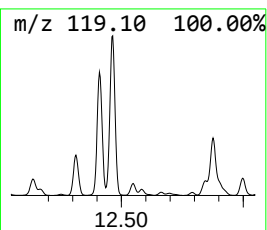
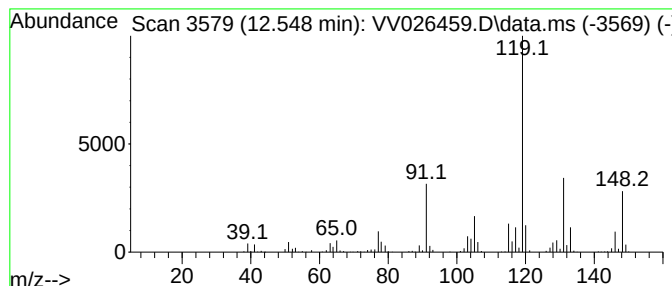
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 23 Benzene, 1-methyl-4-(1-meth... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.548	2.88 ug/L	465367	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	70
2			2-Ethyl-2-methyl-1,3-dithiolane	148	C6H12S2	006008-81-7	58
3			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	58
4			4'-Methylpropiophenone	148	C10H12O	005337-93-9	52
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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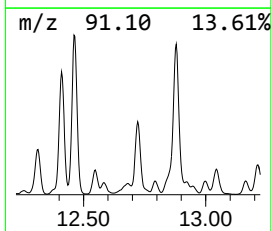
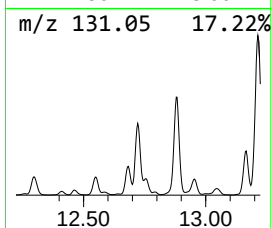
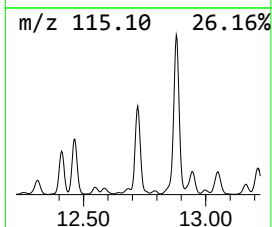
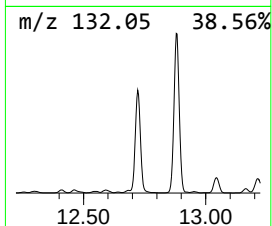
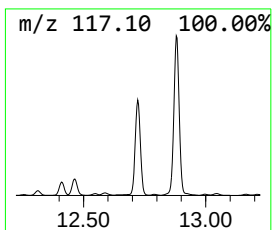
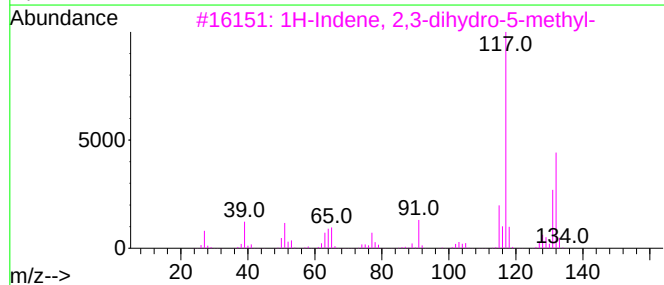
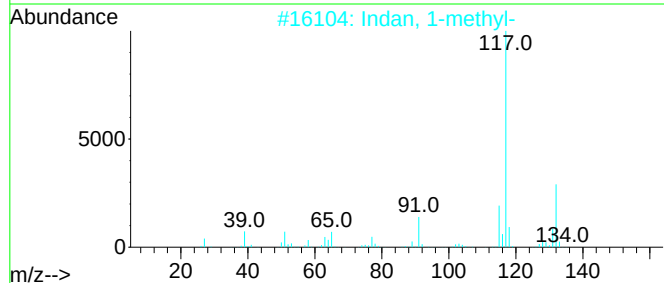
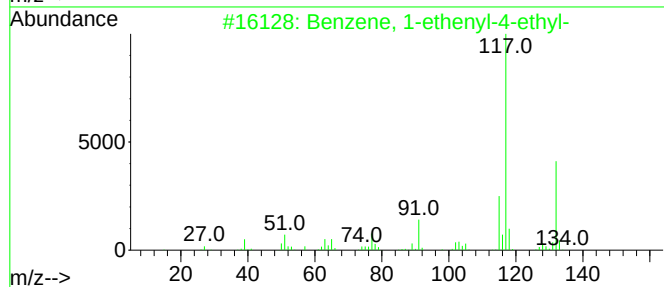
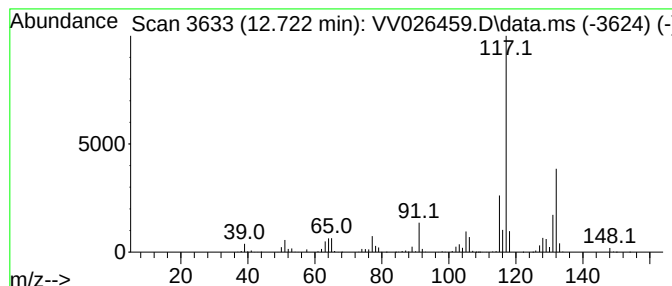
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 25 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.722	15.37 ug/L	2483860	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			Indan, 1-methyl-	132	C10H12	000767-58-8	93
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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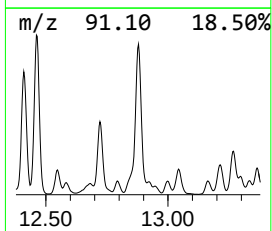
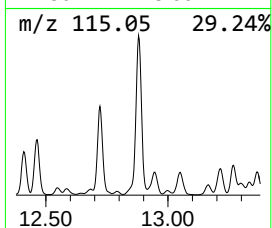
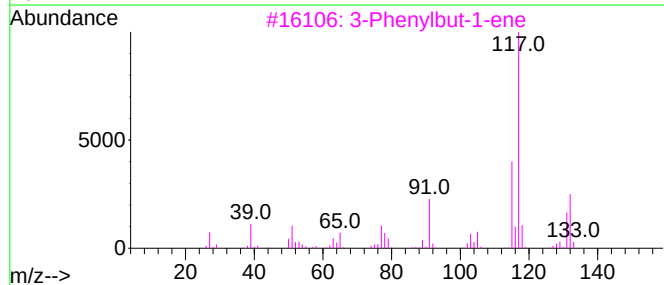
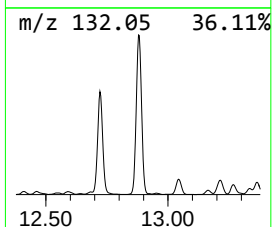
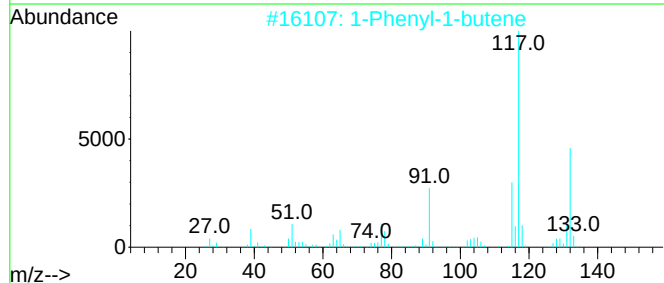
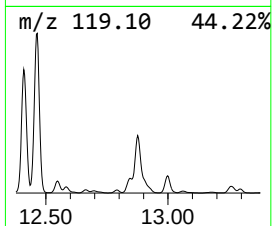
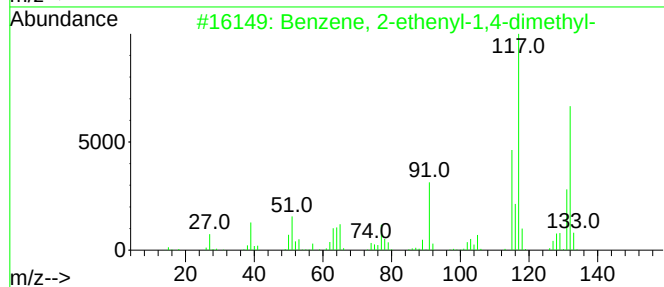
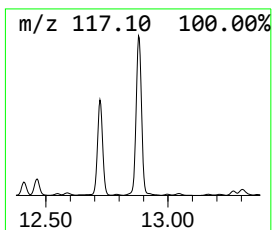
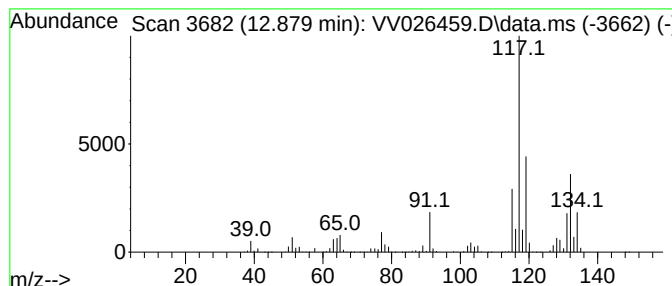
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 26 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.879	36.87 ug/L	5959620	1,4-Dichlorobenzene-d4	11.249

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	76
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	76
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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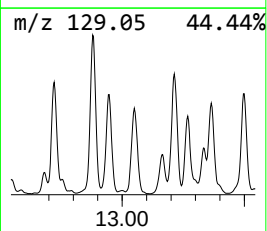
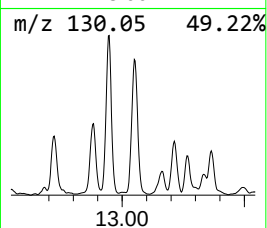
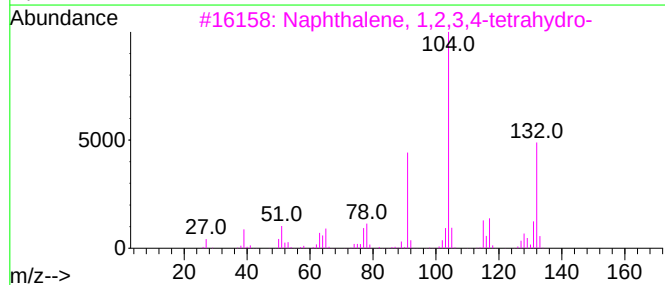
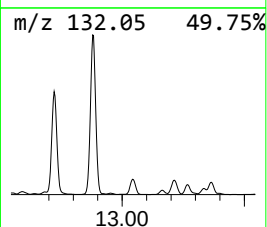
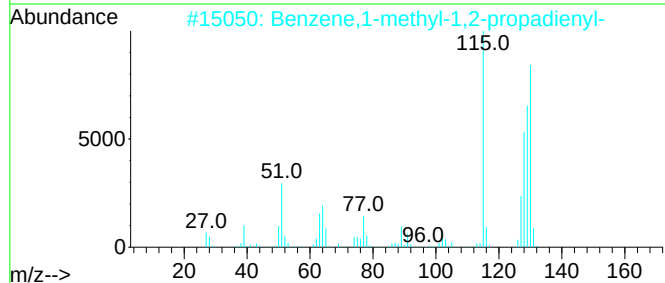
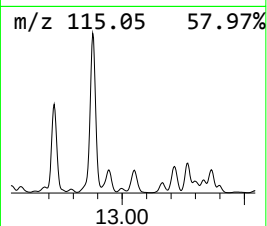
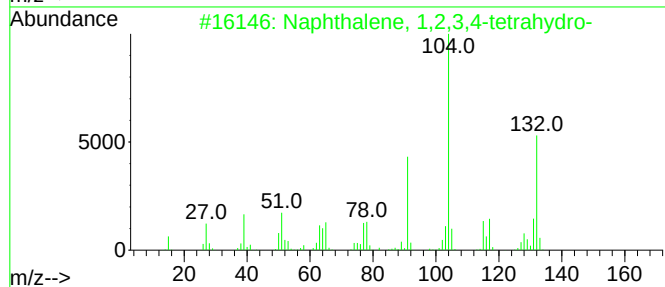
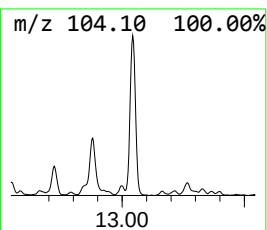
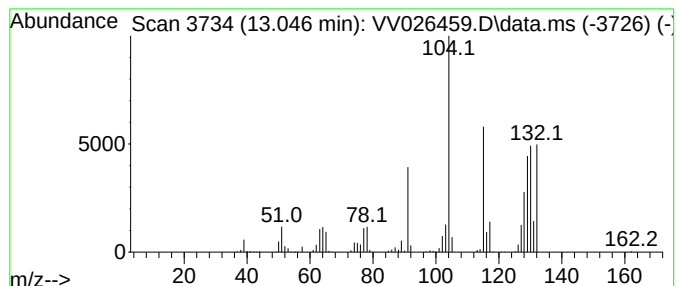
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 28 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.046	4.03 ug/L	651081	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	87
2			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	84
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	50
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	50
5			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

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ClientSampleId :
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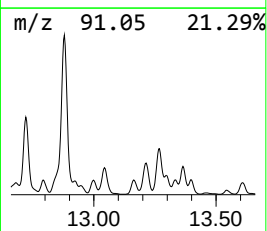
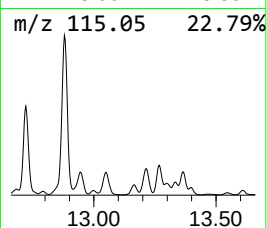
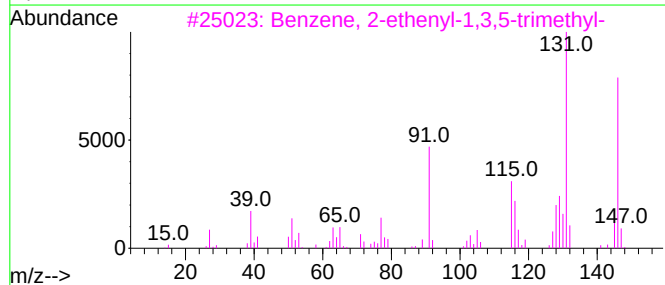
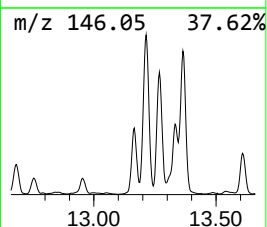
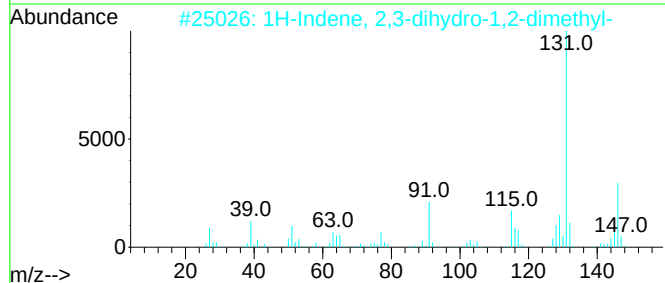
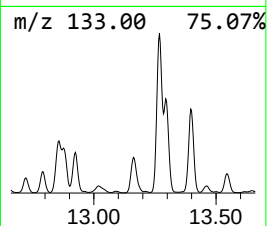
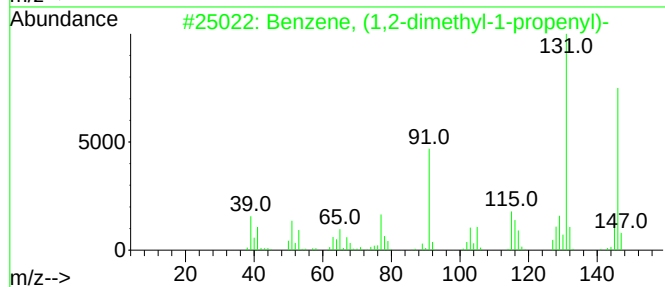
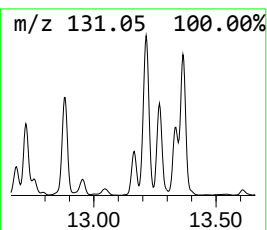
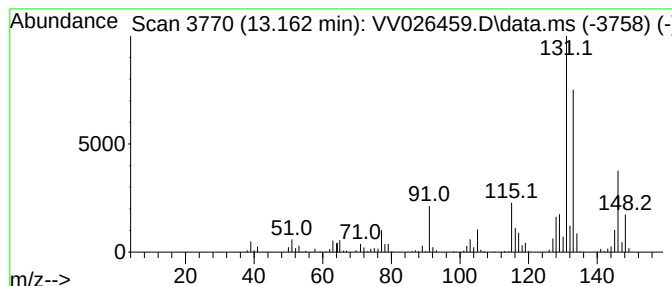
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 29 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.162	3.31 ug/L	535770	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	60
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

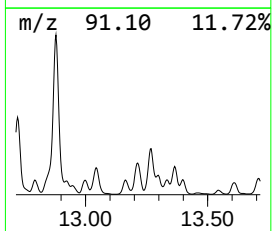
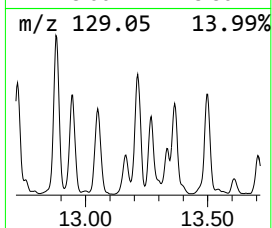
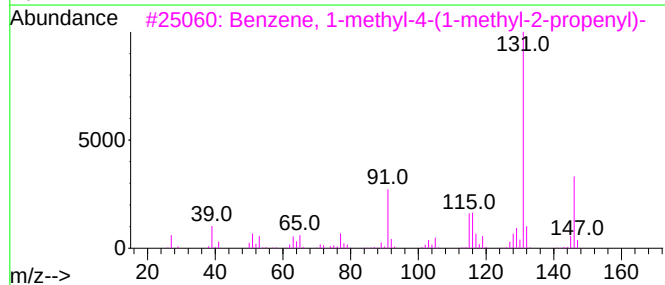
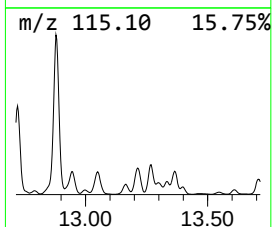
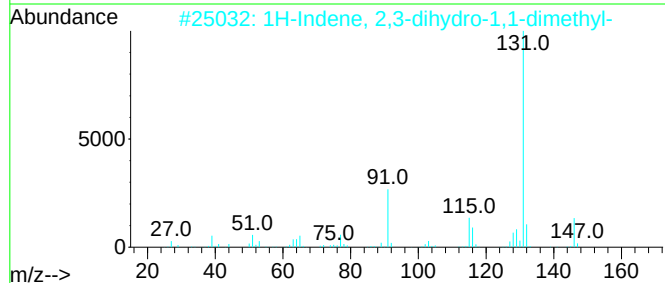
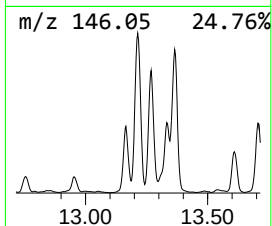
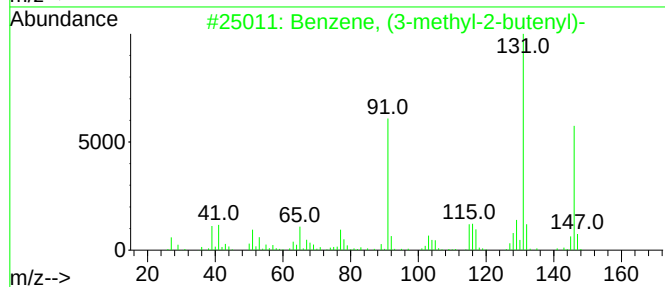
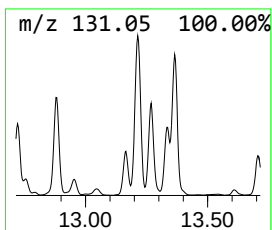
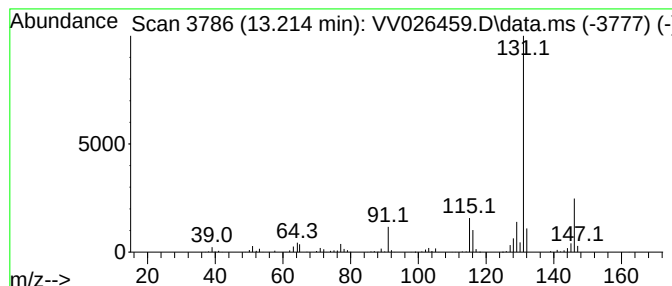
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 30 Benzene, (3-methyl-2-butenyl)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.214	5.67 ug/L	916866	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
2			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	90
3			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

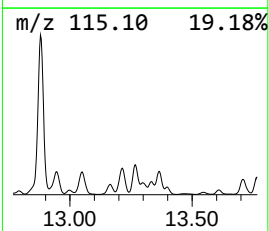
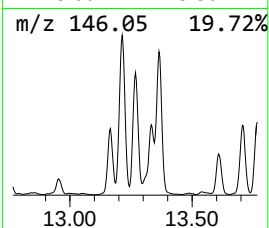
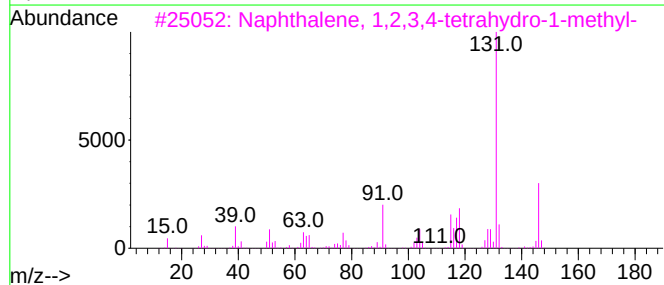
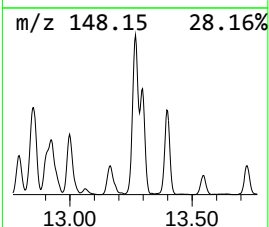
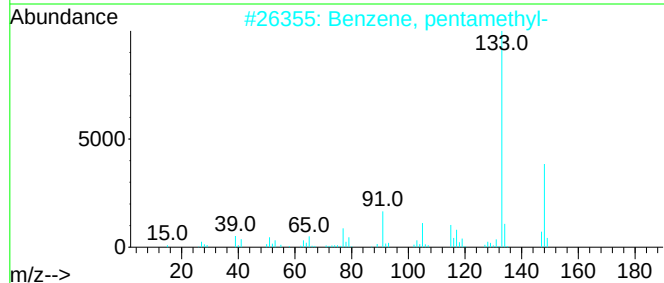
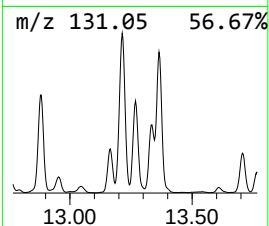
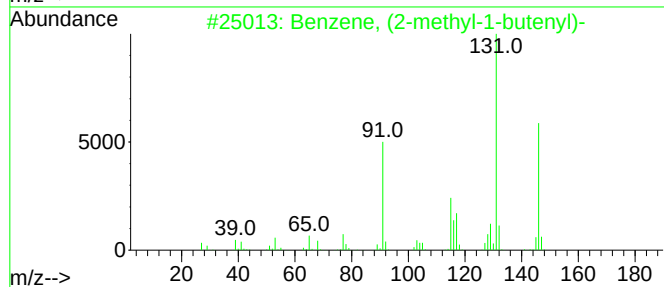
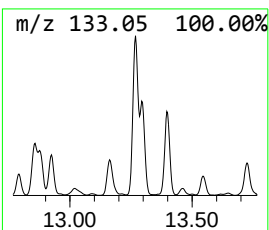
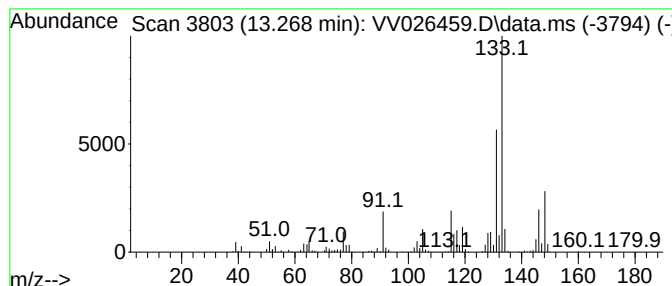
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 31 Benzene, (2-methyl-1-butenyl)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	8.77 ug/L	1416900	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	64
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	51
5			4-tert-Butyltoluene	148	C11H16	000098-51-1	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

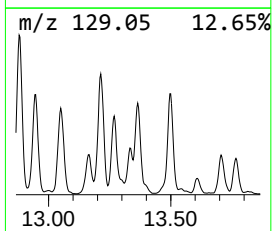
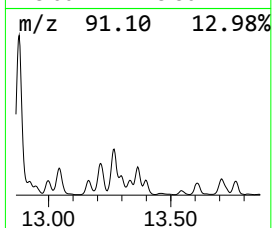
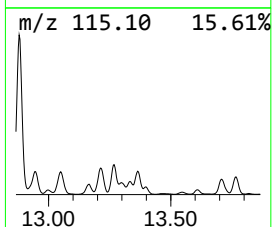
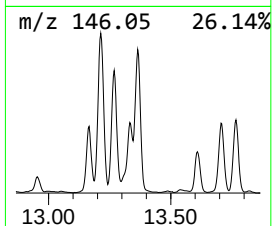
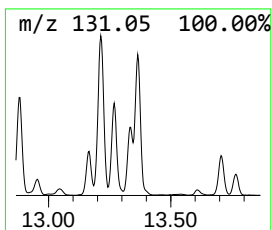
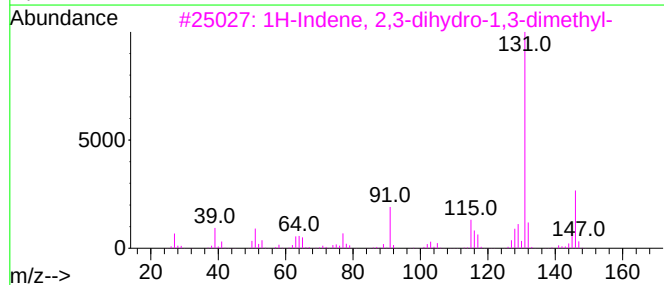
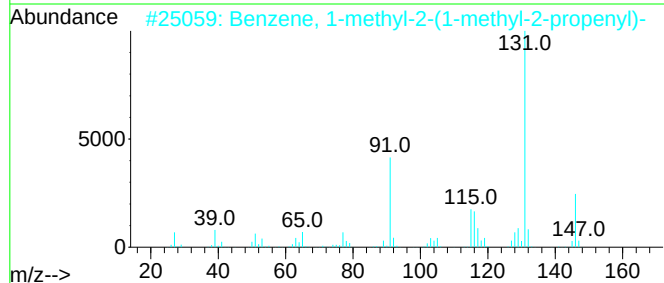
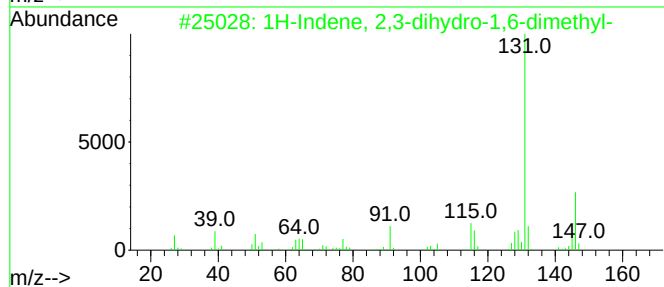
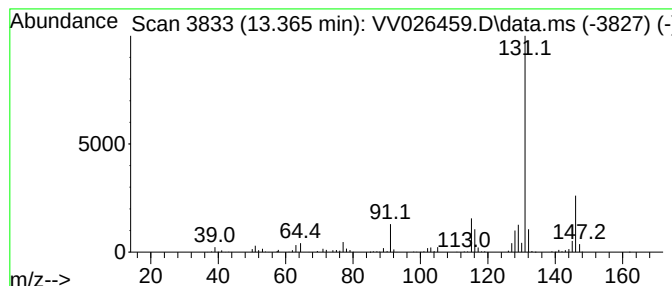
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 32 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.365	3.23 ug/L	522465	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	97
2			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

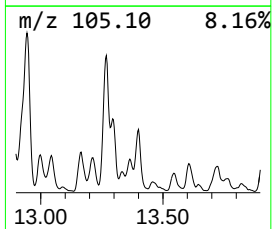
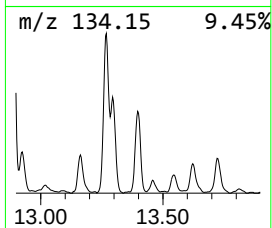
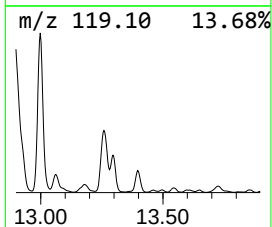
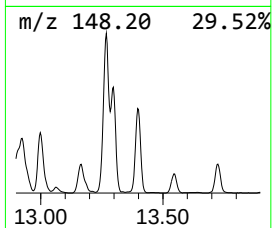
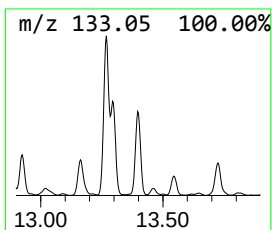
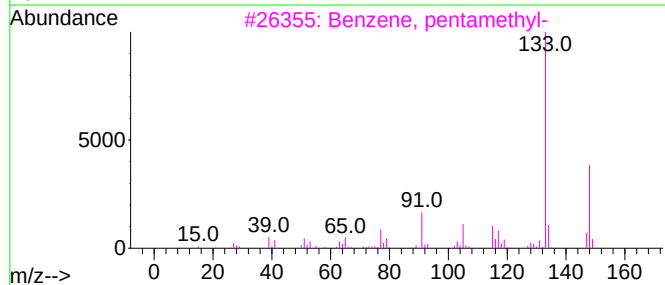
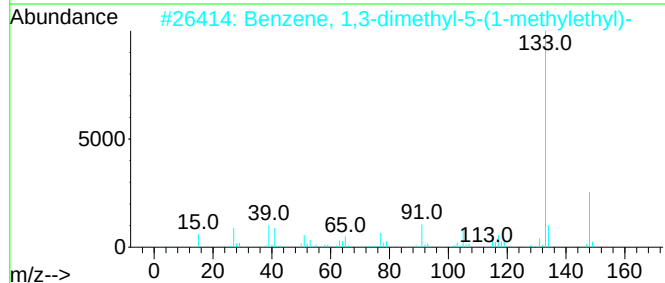
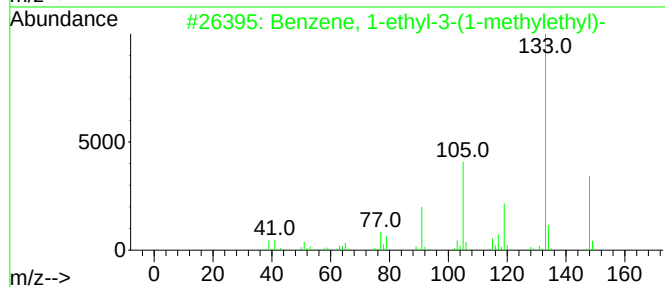
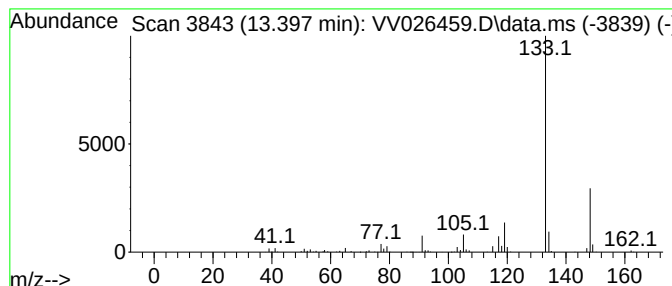
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 33 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.397	2.65 ug/L	428777	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	91
2			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	90
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	86
4			Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	78
5			Benzene, pentamethyl-	148	C11H16	000700-12-9	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

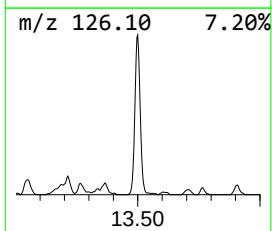
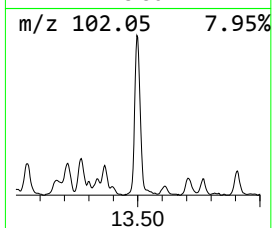
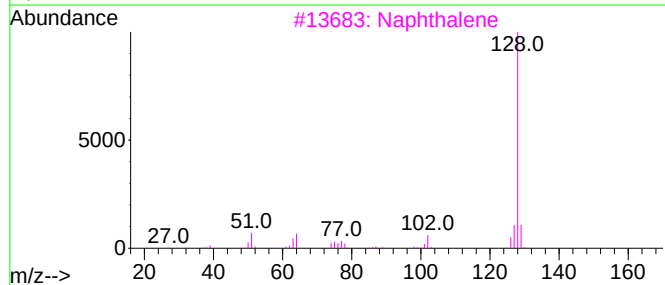
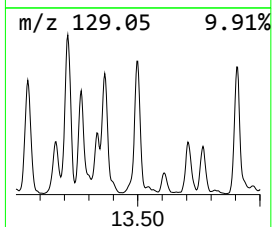
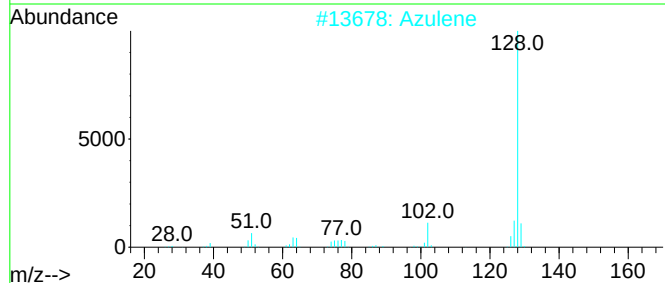
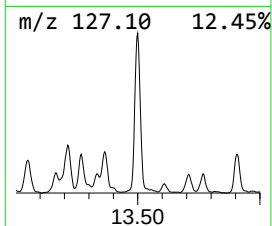
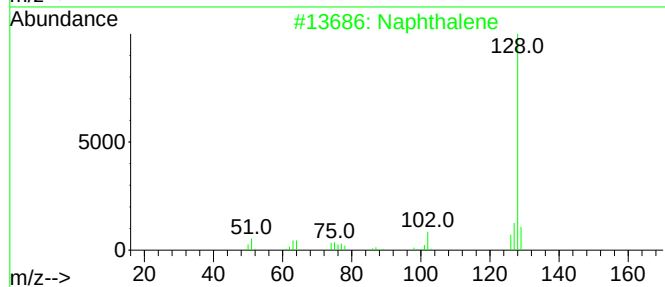
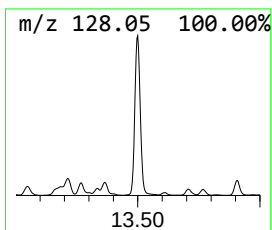
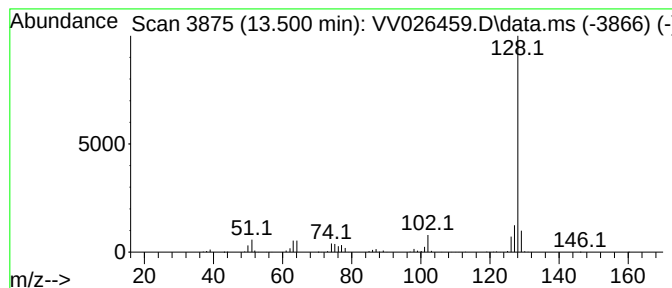
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 34 Azulene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	4.56 ug/L	736767	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	93
3			Naphthalene	128	C10H8	000091-20-3	93
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

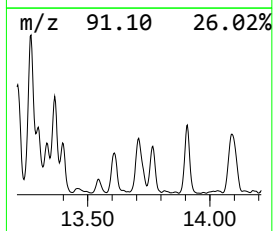
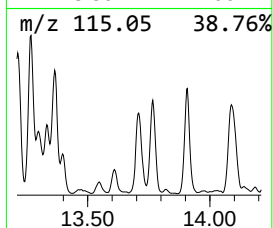
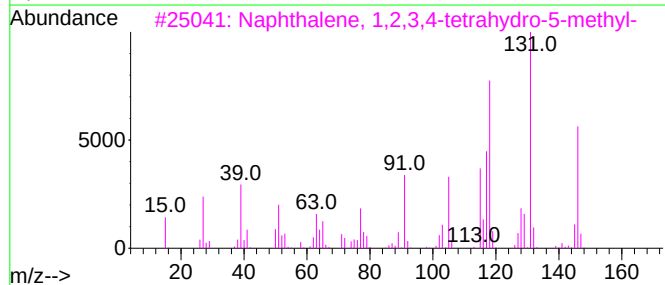
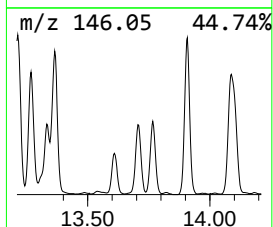
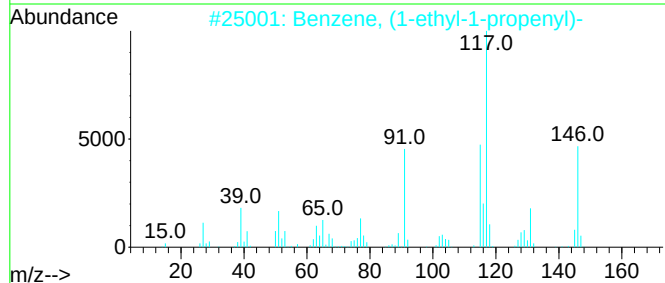
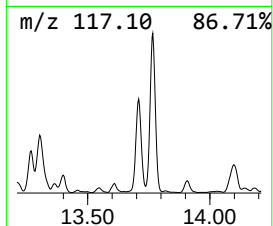
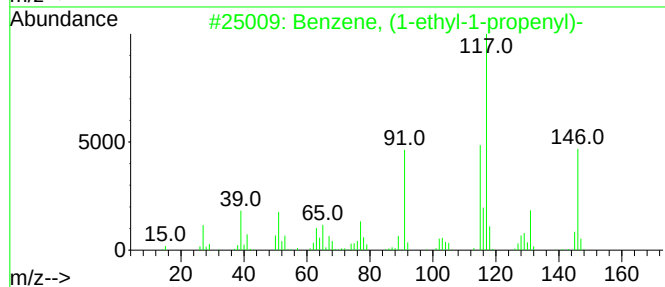
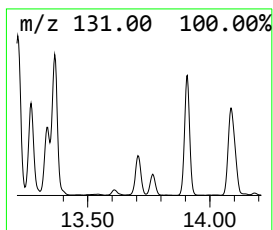
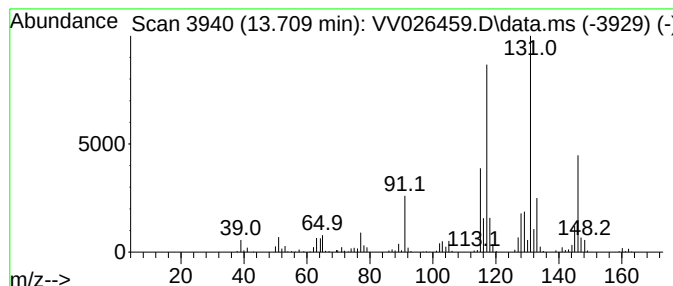
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 35 Benzene, (1-ethyl-1-propenyl)- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.709	3.88 ug/L	627355	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	78
2			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	78
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
4			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	58
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

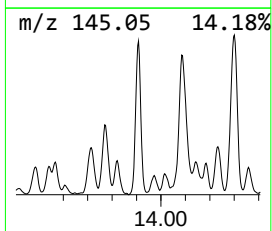
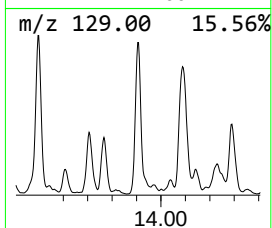
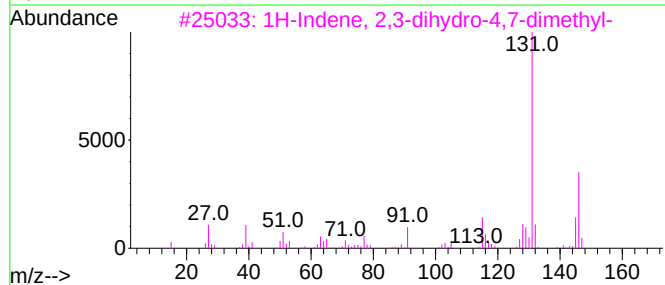
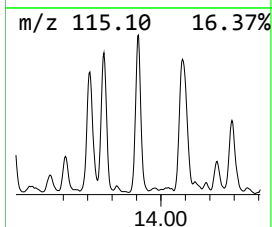
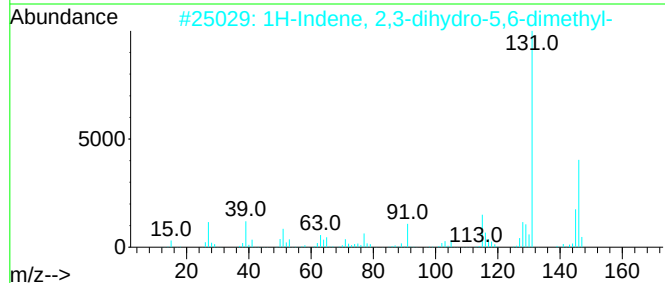
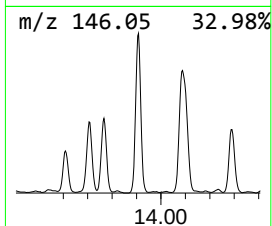
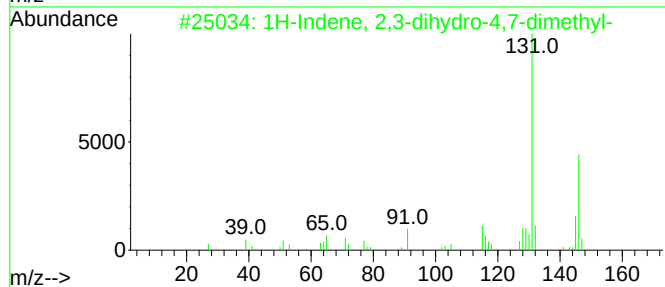
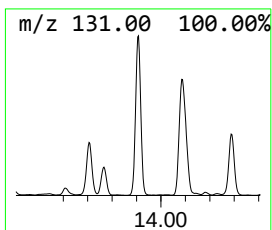
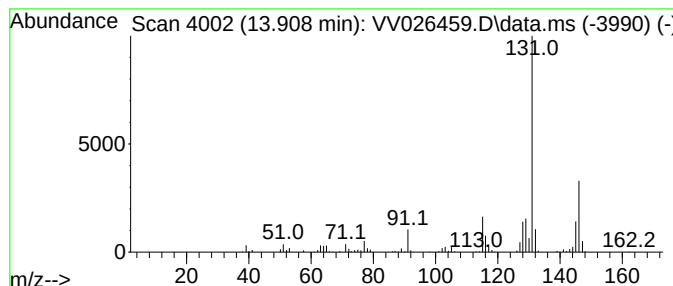
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 37 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.908	5.25 ug/L	848431	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	96
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
5			1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

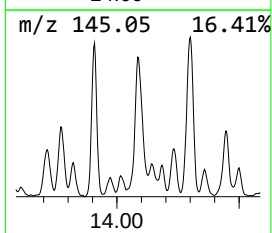
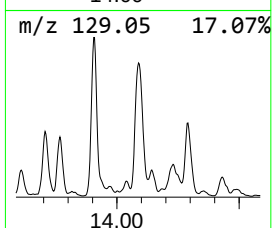
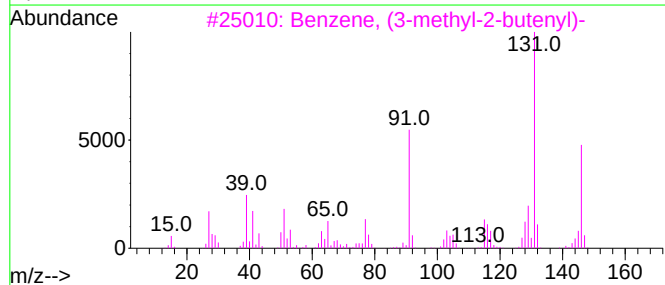
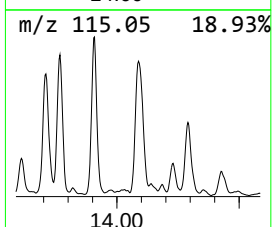
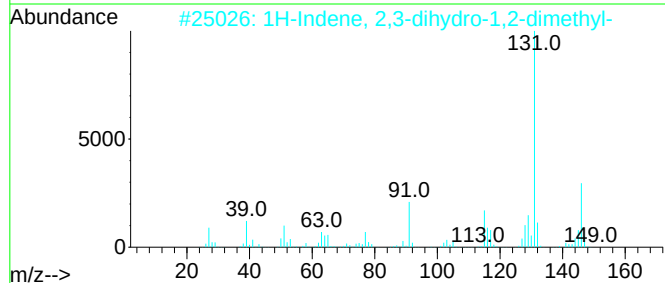
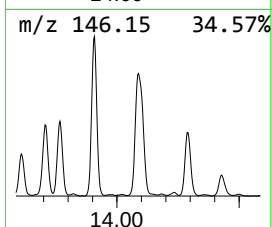
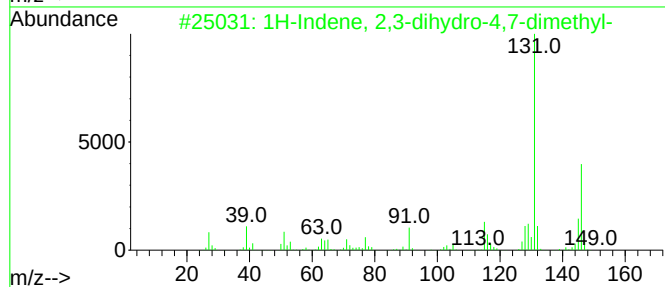
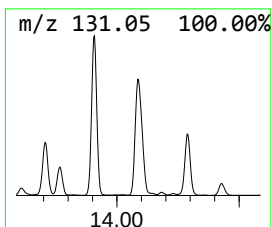
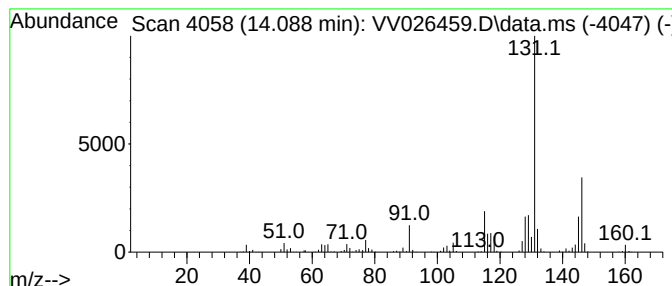
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 38 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.088	6.33 ug/L	1023100	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
Data File : VV026459.D
Acq On : 20 Jun 2022 20:37
Operator : SY/MD
Sample : N3355-01
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EY0C3

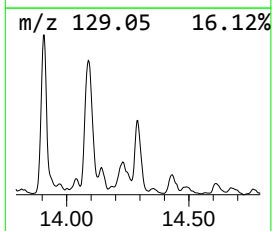
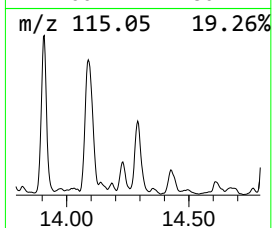
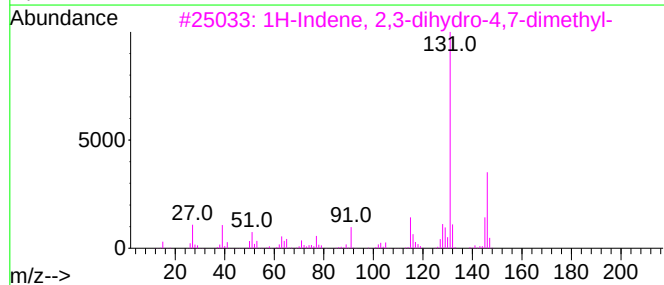
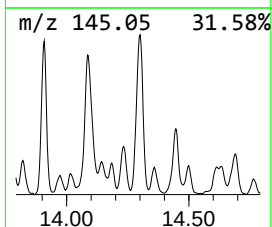
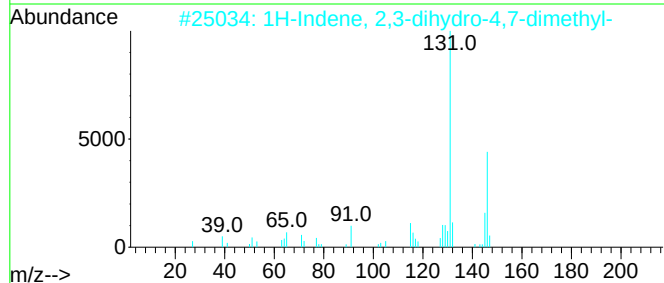
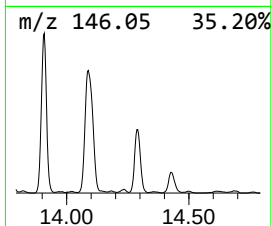
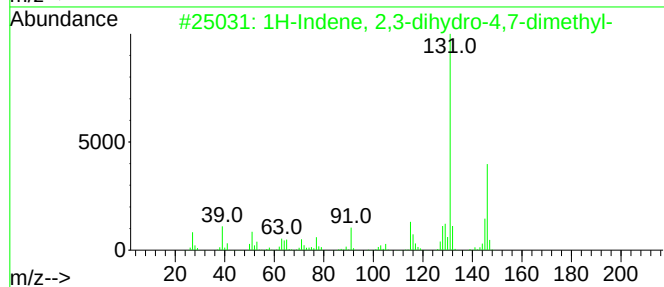
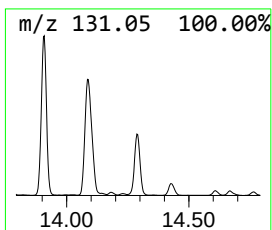
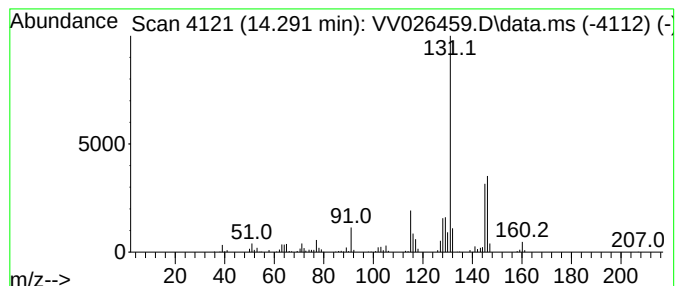
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR061522WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 39 1H-Indene, 2,3-dihydro-4,6-... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.291	2.77 ug/L	447705	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93		
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93		
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90		
4	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	87		
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062022\
 Data File : VV026459.D
 Acq On : 20 Jun 2022 20:37
 Operator : SY/MD
 Sample : N3355-01
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EY0C3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR061522WMA.M
 Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	4.777	8.2	ug/L	876246	1	5.625	535866	5.0
(DEL) Alkane: S...	5.243	7.2	ug/L	766811	1	5.625	535866	5.0
(DEL) Alkane: S...	6.658	3.6	ug/L	381024	1	5.625	535866	5.0
(DEL) Alkane: S...	6.780	6.2	ug/L	665614	1	5.625	535866	5.0
Benzene, propyl-	10.352	24.5	ug/L	3961130	3	11.249	808156	5.0
Benzene, 1-ethy...	10.448	8.2	ug/L	1324940	3	11.249	808156	5.0
Benzene, 1-ethy...	10.471	17.7	ug/L	2866890	3	11.249	808156	5.0
Benzene, 1-ethy...	10.735	54.3	ug/L	8773700	3	11.249	808156	5.0
Benzene, (2-met...	11.053	2.0	ug/L	320915	3	11.249	808156	5.0
Benzene, 1-meth...	11.085	2.6	ug/L	418825	3	11.249	808156	5.0
Benzene, 1,2,3-...	11.333	34.6	ug/L	5601270	3	11.249	808156	5.0
Indane	11.529	31.6	ug/L	5099780	3	11.249	808156	5.0
Benzene, 1-meth...	11.818	9.9	ug/L	1605620	3	11.249	808156	5.0
Benzene, 2-ethy...	11.908	20.7	ug/L	3345810	3	11.249	808156	5.0
o-Cymene	11.937	17.4	ug/L	2820810	3	11.249	808156	5.0
Benzene, 1-ethy...	12.014	39.0	ug/L	6304500	3	11.249	808156	5.0
Benzene, 1-ethe...	12.114	20.5	ug/L	3319790	3	11.249	808156	5.0
Benzene, 4-ethy...	12.313	7.1	ug/L	1144590	3	11.249	808156	5.0
Benzene, 1,2,3,...	12.410	20.8	ug/L	3359240	3	11.249	808156	5.0
Benzene, 1,2,3,...	12.464	26.9	ug/L	4343260	3	11.249	808156	5.0
Benzene, 1-meth...	12.548	2.9	ug/L	465367	3	11.249	808156	5.0
Benzene, 1-ethe...	12.722	15.4	ug/L	2483860	3	11.249	808156	5.0
Benzene, 2-ethe...	12.879	36.9	ug/L	5959620	3	11.249	808156	5.0
Naphthalene, 1,...	13.046	4.0	ug/L	651081	3	11.249	808156	5.0
Benzene, (1,2-d...	13.162	3.3	ug/L	535770	3	11.249	808156	5.0
Benzene, (3-met...	13.214	5.7	ug/L	916866	3	11.249	808156	5.0
Benzene, (2-met...	13.268	8.8	ug/L	1416900	3	11.249	808156	5.0
1H-Indene, 2,3-...	13.365	3.2	ug/L	522465	3	11.249	808156	5.0
Benzene, 1-ethy...	13.397	2.6	ug/L	428777	3	11.249	808156	5.0
Azulene	13.500	4.6	ug/L	736767	3	11.249	808156	5.0
Benzene, (1-eth...	13.709	3.9	ug/L	627355	3	11.249	808156	5.0
1H-Indene, 2,3-...	13.908	5.3	ug/L	848431	3	11.249	808156	5.0
1H-Indene, 2,3-...	14.088	6.3	ug/L	1023100	3	11.249	808156	5.0
1H-Indene, 2,3-...	14.291	2.8	ug/L	447705	3	11.249	808156	5.0