

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102722\
 Data File : VV028715.D
 Acq On : 27 Oct 2022 13:12
 Operator : SY/MD
 Sample : N5232-16
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BHAB9

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\SFAMVTR102522WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV028715.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.581	155	168	178	rVB2	132196	224073	1.23%	0.708%
2	1.947	271	282	301	rBV	2012519	2737328	15.00%	8.643%
3	3.282	677	697	725	rBV	979966	2275664	12.47%	7.185%
4	3.915	883	894	929	rBV3	42476	186091	1.02%	0.588%
5	4.343	1010	1027	1047	rBV	82669	213514	1.17%	0.674%
6	5.044	1231	1245	1252	rBV2	182149	424135	2.32%	1.339%
7	5.095	1253	1261	1290	rVB	245991	567100	3.11%	1.791%
8	5.616	1412	1423	1445	rBV	138397	291538	1.60%	0.921%
9	6.066	1551	1563	1590	rBV	118282	258835	1.42%	0.817%
10	7.314	1933	1951	1969	rBV	128707	236968	1.30%	0.748%
11	8.092	2178	2193	2230	rBV2	242401	653623	3.58%	2.064%
12	8.876	2418	2437	2473	rBV	11285367	18246156	100.00%	57.613%
13	9.564	2636	2651	2675	rBV3	86489	215784	1.18%	0.681%
14	9.924	2751	2763	2779	rBV	1230265	1893027	10.37%	5.977%
15	10.349	2878	2895	2909	rBV2	247566	496635	2.72%	1.568%
16	11.243	3164	3173	3192	rVB4	180385	452632	2.48%	1.429%
17	11.342	3193	3204	3223	rBV3	211998	365849	2.01%	1.155%
18	11.519	3248	3259	3278	rBV	280570	485821	2.66%	1.534%
19	11.619	3279	3290	3316	rVB3	162804	378608	2.08%	1.195%
20	12.008	3400	3411	3425	rBV	141019	231530	1.27%	0.731%
21	12.403	3522	3534	3543	rBV	141558	219381	1.20%	0.693%
22	12.873	3663	3680	3690	rBV2	176192	289239	1.59%	0.913%
23	16.297	4734	4745	4758	rBV	214942	326896	1.79%	1.032%

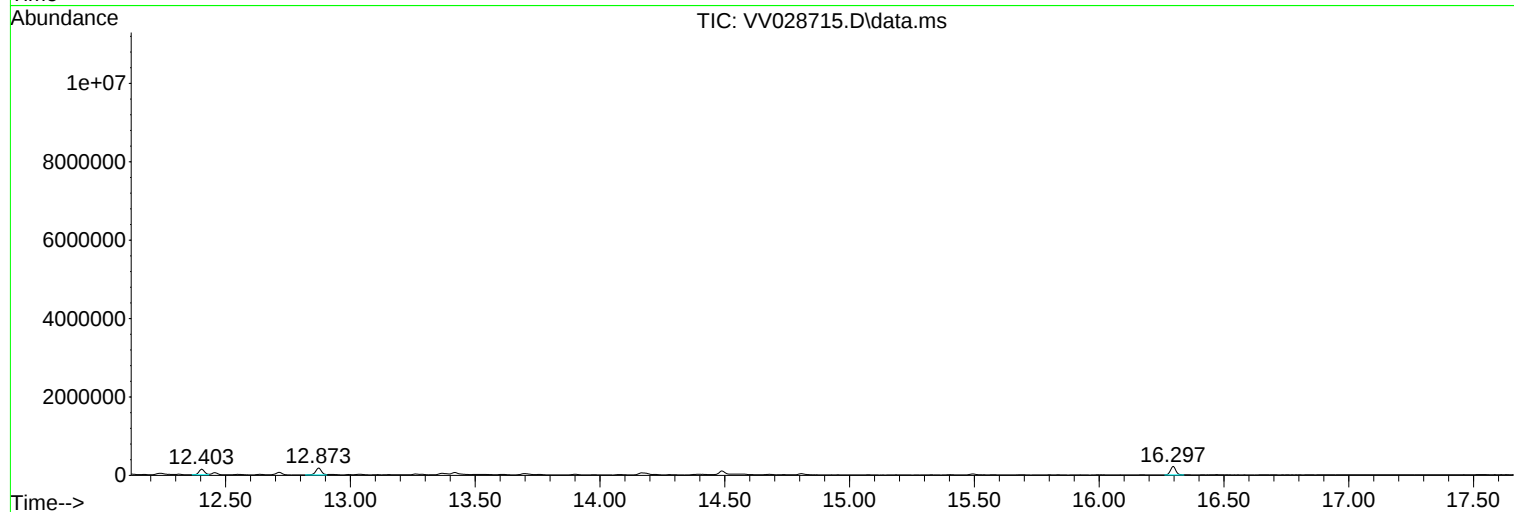
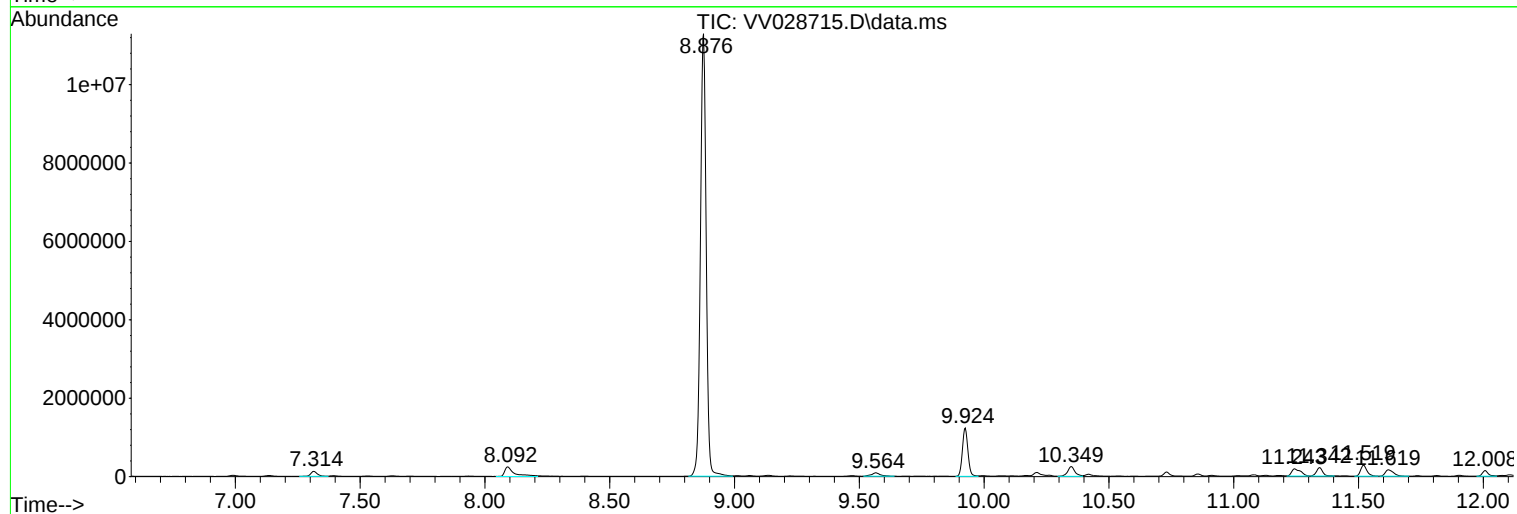
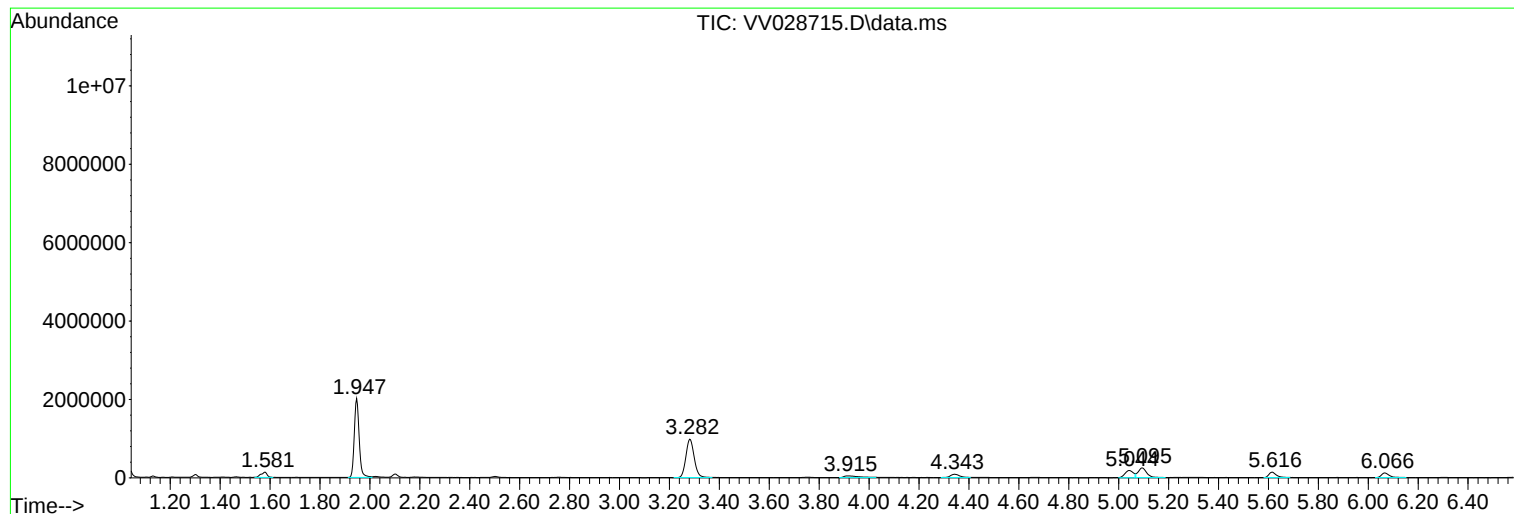
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102522WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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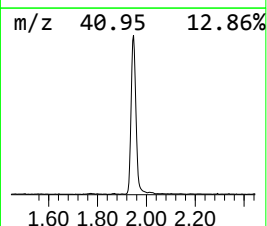
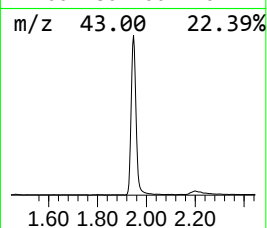
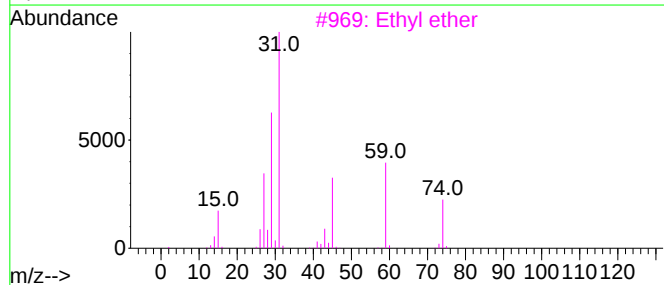
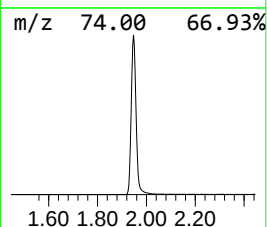
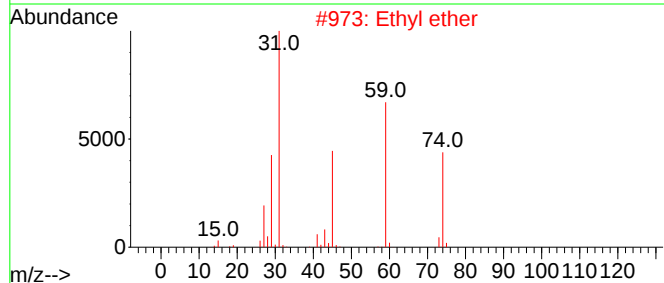
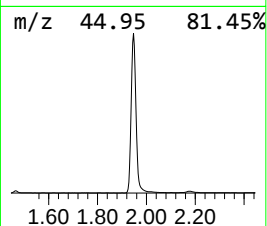
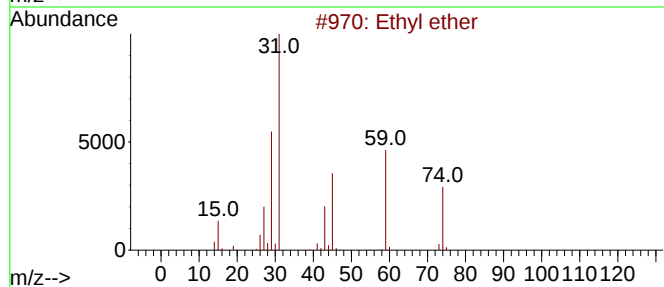
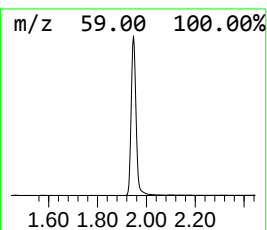
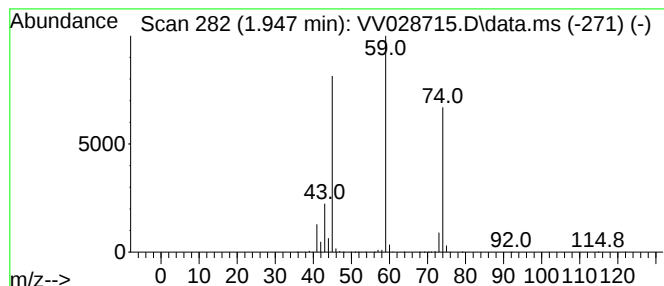
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 Ethyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.947	46.95 ug/L	2737330	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether			74	C4H10O	000060-29-7	90
2	Ethyl ether			74	C4H10O	000060-29-7	90
3	Ethyl ether			74	C4H10O	000060-29-7	90
4	Ethyl ether			74	C4H10O	000060-29-7	72
5	Ethyl ether			74	C4H10O	000060-29-7	64



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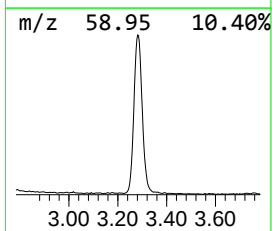
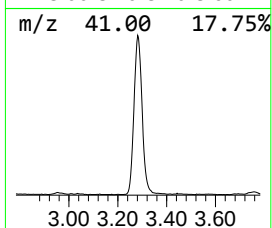
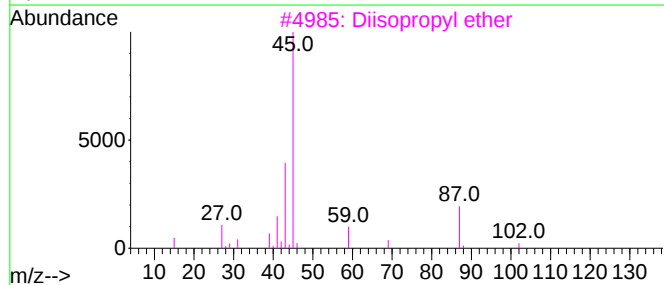
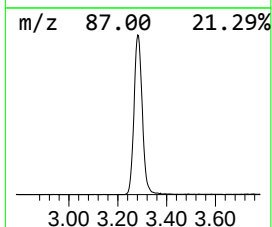
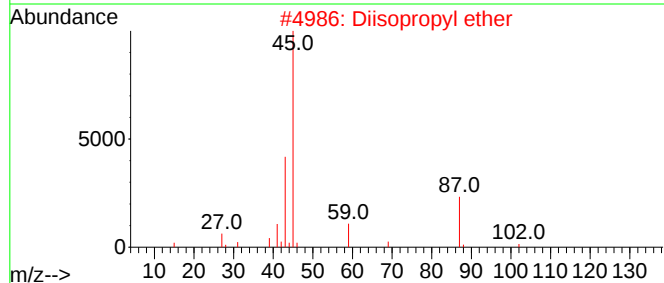
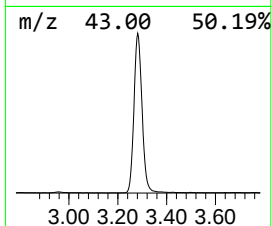
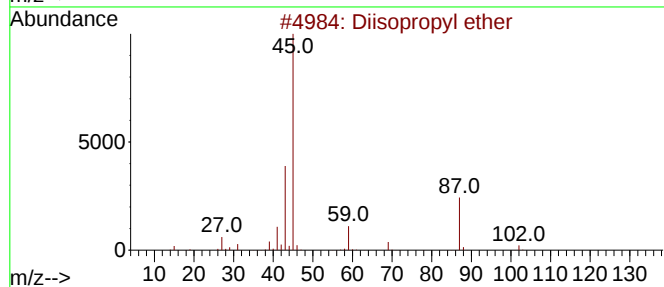
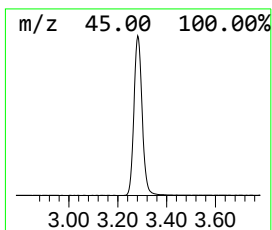
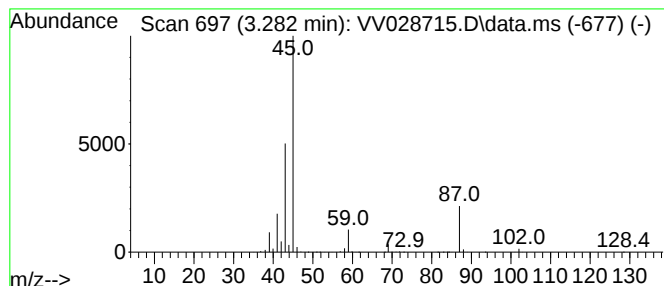
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102522WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Diisopropyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.282	39.03 ug/L	2275660	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisopropyl ether	102	C6H14O	000108-20-3	91
2			Diisopropyl ether	102	C6H14O	000108-20-3	91
3			Diisopropyl ether	102	C6H14O	000108-20-3	83
4			Diisopropyl ether	102	C6H14O	000108-20-3	83
5			Diisopropyl ether	102	C6H14O	000108-20-3	83



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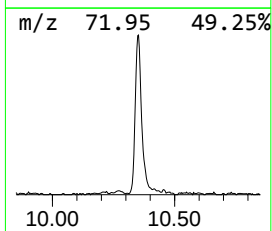
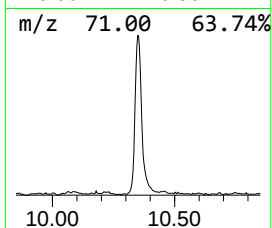
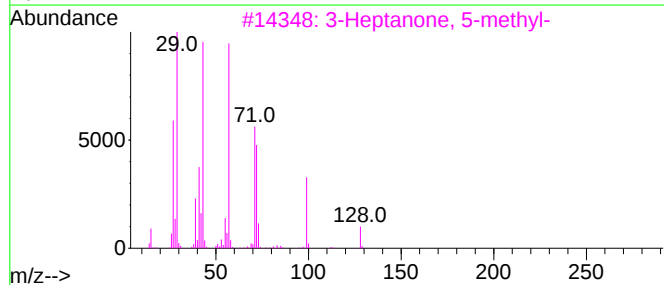
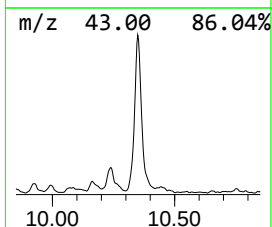
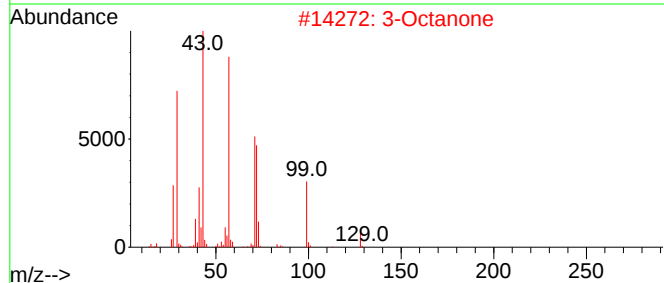
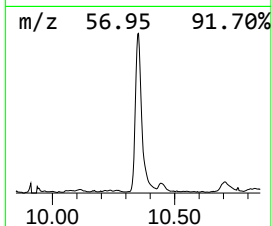
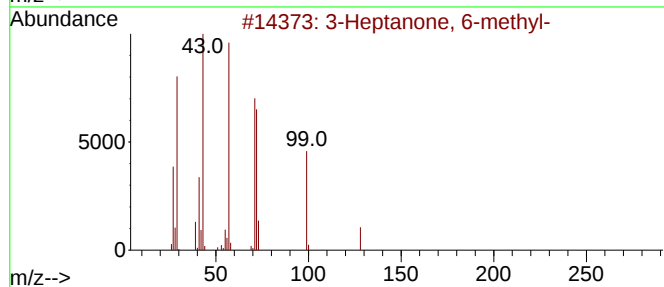
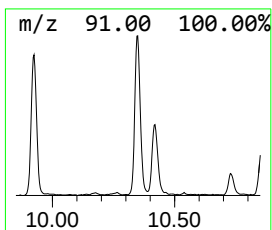
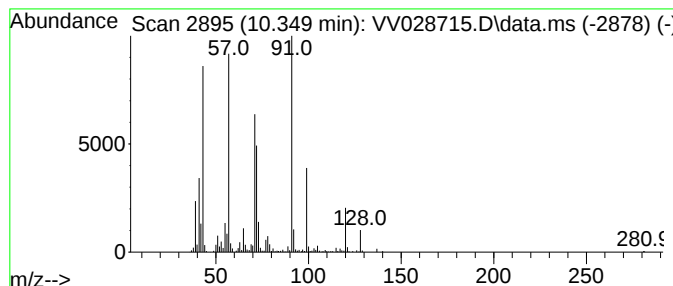
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102522WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
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Peak Number 3 3-Heptanone, 6-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.349	5.49 ug/L	496635	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Heptanone, 6-methyl-	128	C8H16O	000624-42-0	93
2			3-Octanone	128	C8H16O	000106-68-3	76
3			3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	76
4			3-Octanone	128	C8H16O	000106-68-3	70
5			3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	58



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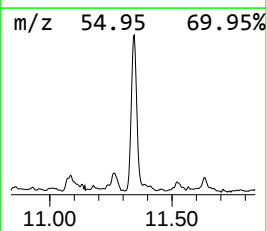
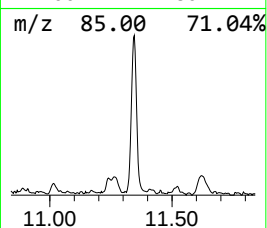
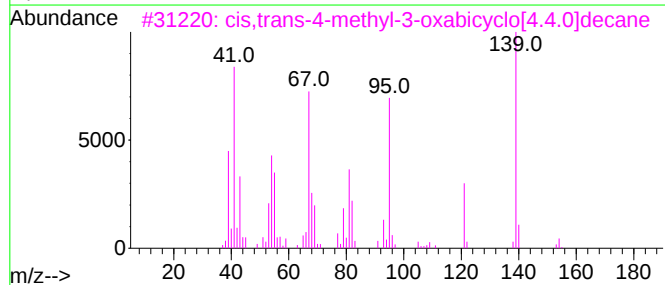
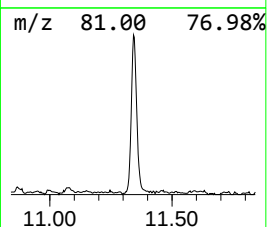
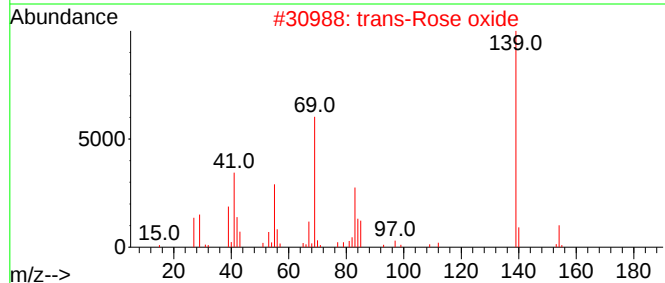
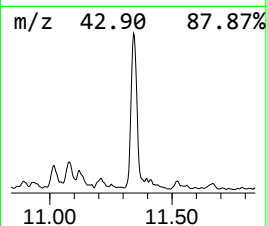
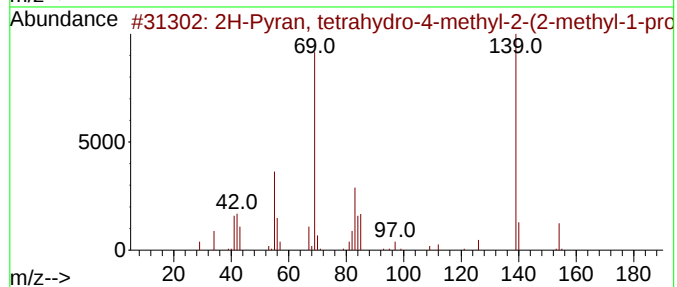
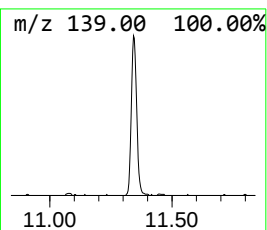
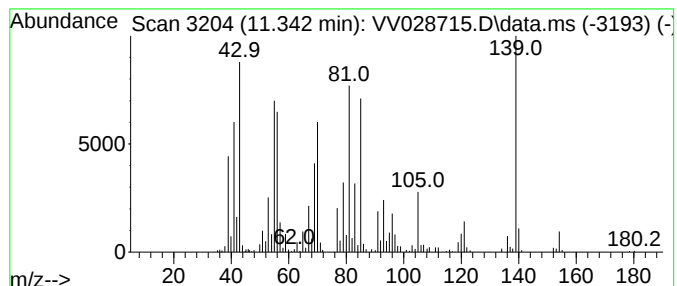
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.342	4.04 ug/L	365849	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran, tetrahydro-4-methyl-2-...	154	C10H18O	016409-43-1	25
2			trans-Rose oxide	154	C10H18O	000876-18-6	18
3			cis,trans-4-methyl-3-oxabicyclo[...]	154	C10H18O	1000216-89-8	18
4			(2S,4R)-4-Methyl-2-(2-methylprop...	154	C10H18O	003033-23-6	18
5			4-Decene, 5-methyl-, (E)-	154	C11H22	062338-51-6	15



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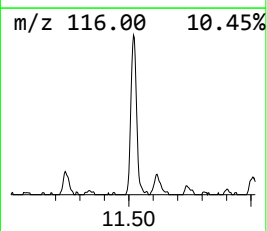
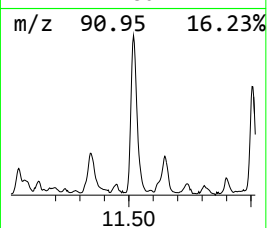
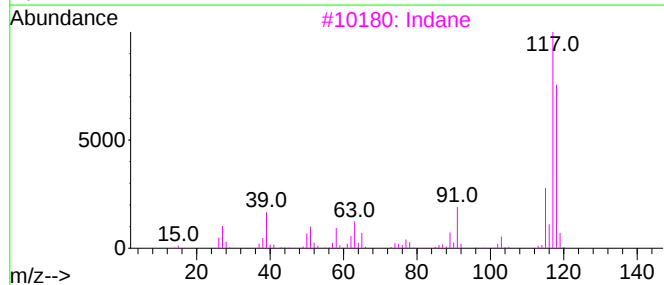
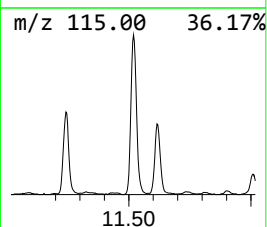
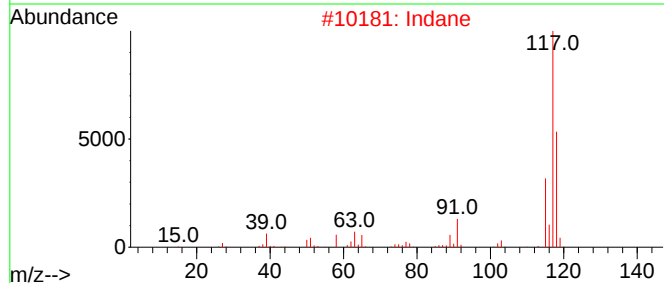
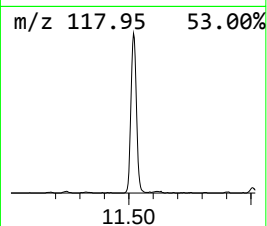
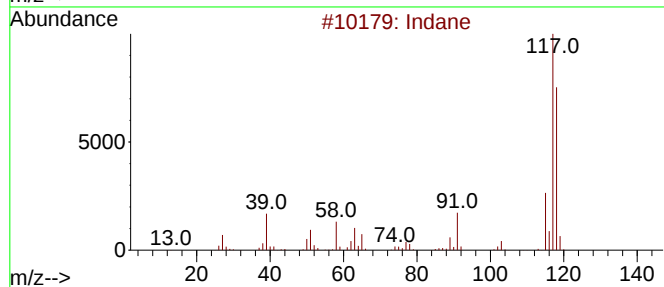
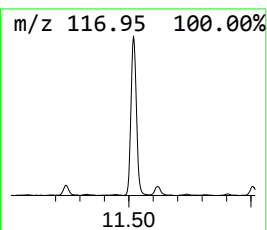
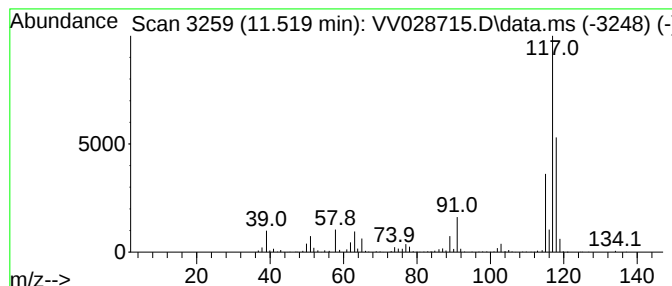
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.519	5.37 ug/L	485821	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	96
2	Indane			118	C9H10	000496-11-7	95
3	Indane			118	C9H10	000496-11-7	90
4	Benzene, 2-propenyl-			118	C9H10	000300-57-2	81
5	Indane			118	C9H10	000496-11-7	81



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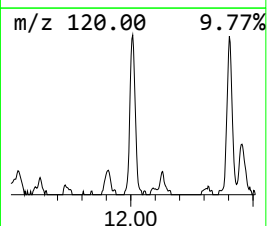
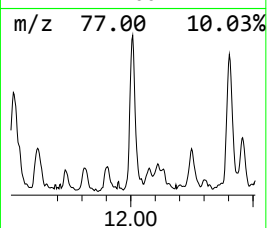
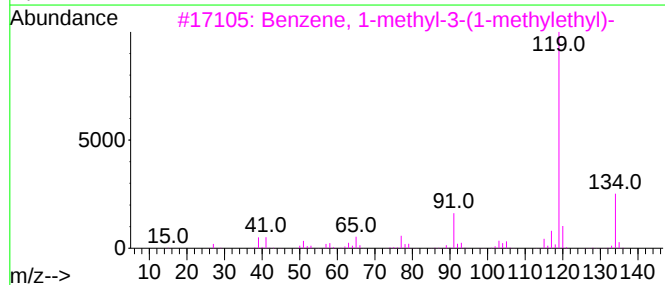
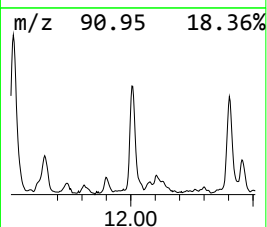
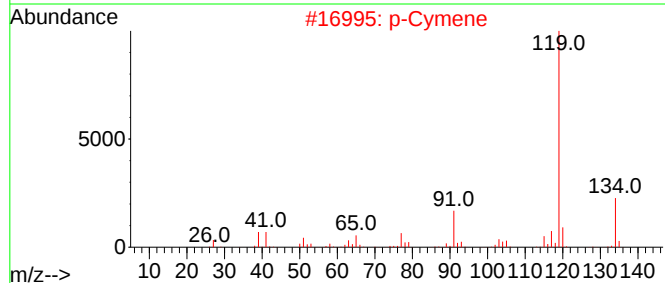
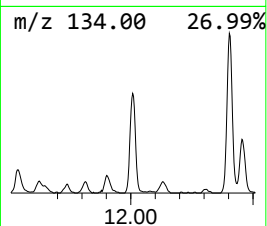
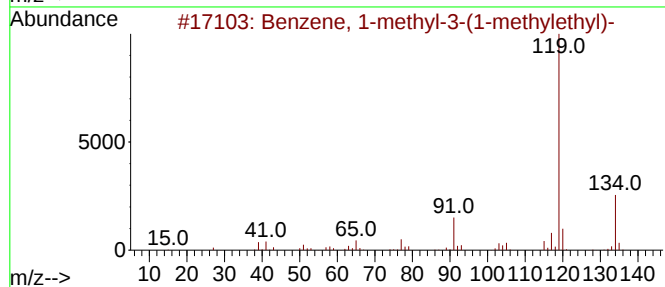
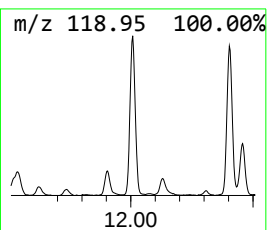
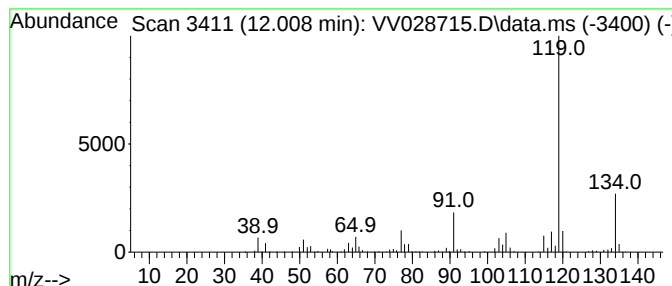
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.008	2.56 ug/L	231530	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2			p-Cymene	134	C10H14	000099-87-6	95
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	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\VV102722\
Data File : VV028715.D
Acq On : 27 Oct 2022 13:12
Operator : SY/MD
Sample : N5232-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHAB9

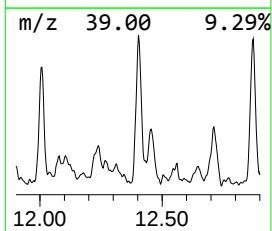
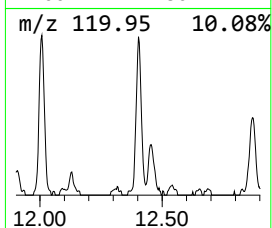
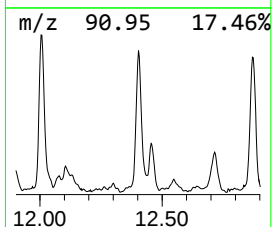
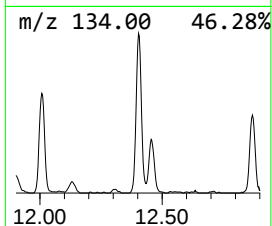
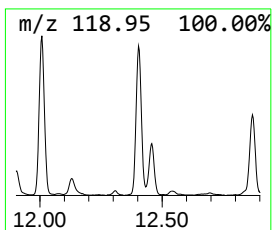
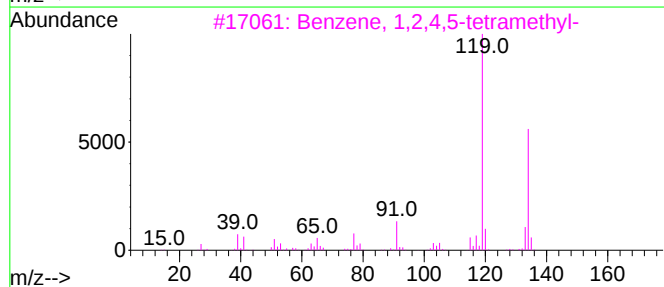
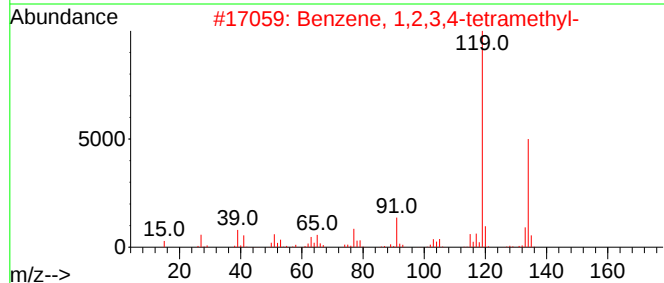
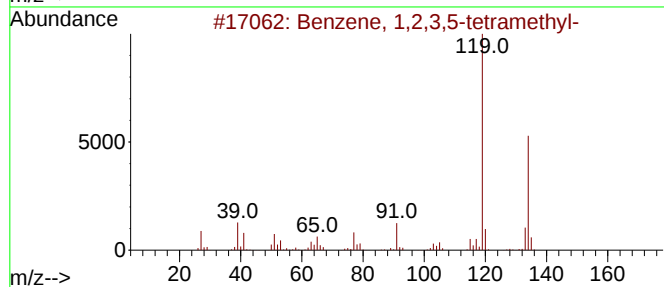
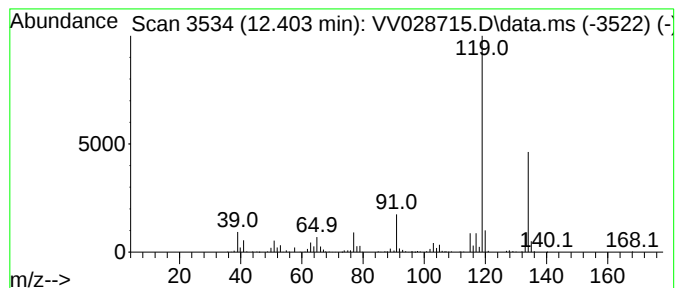
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.403	2.42 ug/L	219381	1,4-Dichlorobenzene-d4	11.243

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102722\
Data File : VV028715.D
Acq On : 27 Oct 2022 13:12
Operator : SY/MD
Sample : N5232-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHAB9

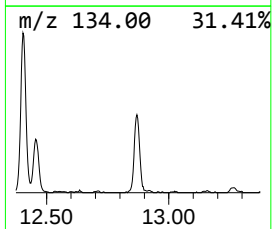
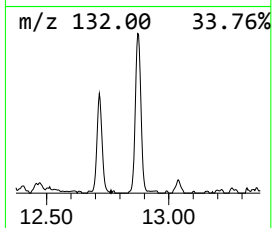
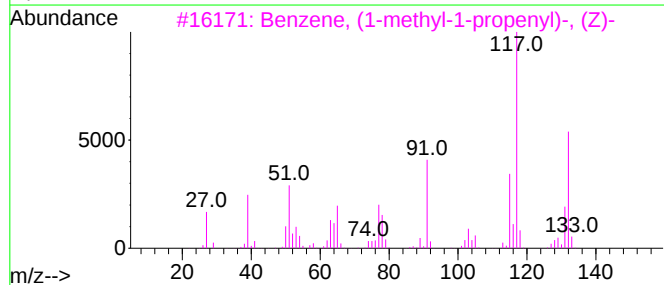
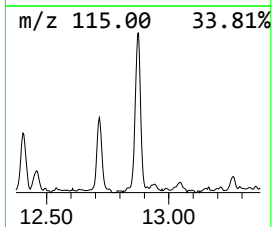
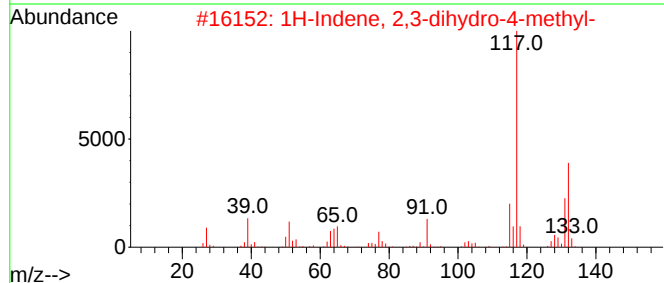
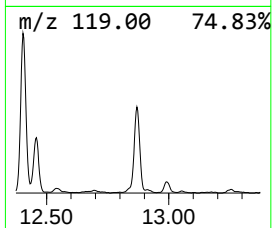
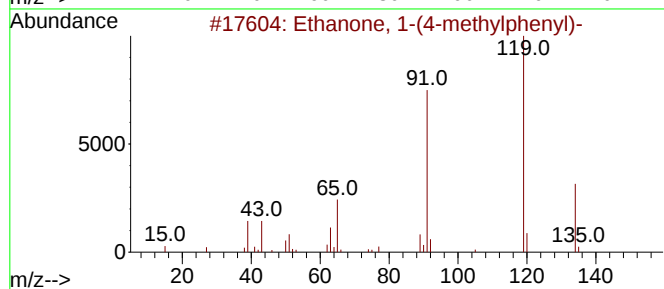
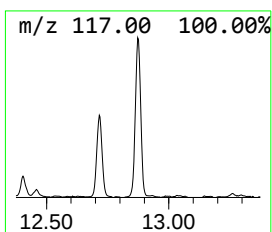
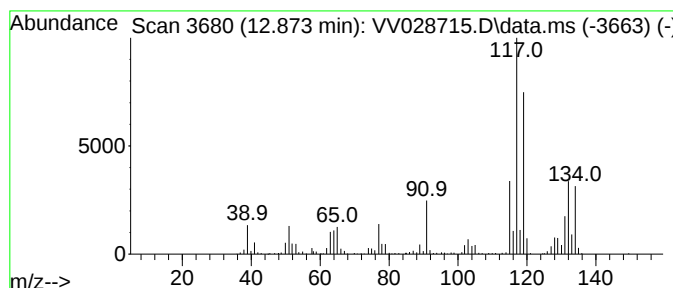
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 Ethanone, 1-(4-methylphenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	3.20 ug/L	289239	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-(4-methylphenyl)-	134	C9H10O	000122-00-9	70
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	60
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	60
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	55
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102722\
Data File : VV028715.D
Acq On : 27 Oct 2022 13:12
Operator : SY/MD
Sample : N5232-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHAB9

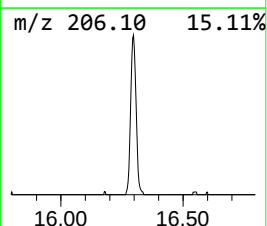
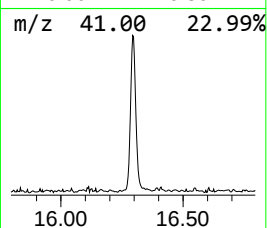
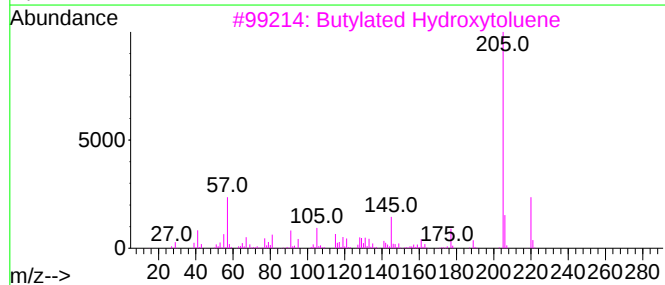
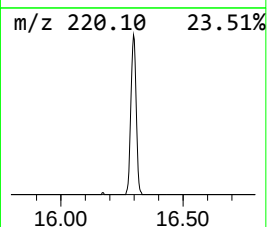
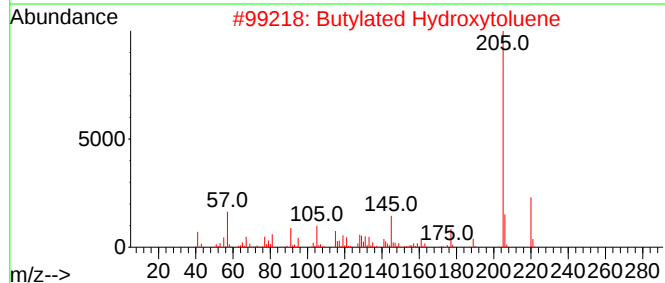
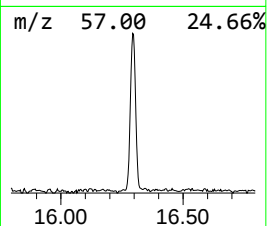
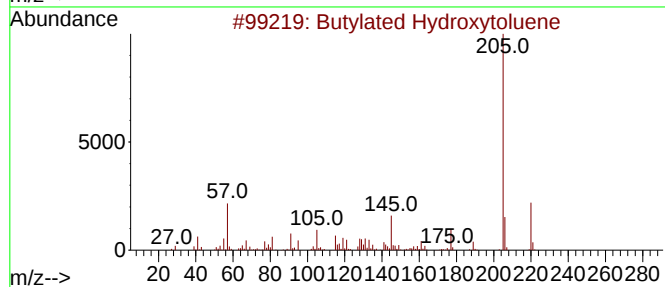
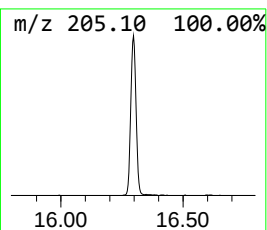
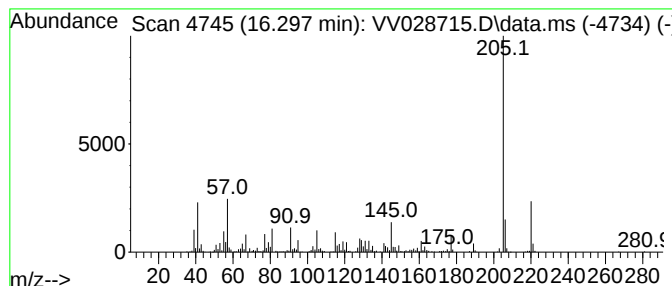
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Butylated Hydroxytoluene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.297	3.61 ug/L	326896	1,4-Dichlorobenzene-d4	11.243

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butylated Hydroxytoluene	220	C15H24O	000128-37-0	99
2			Butylated Hydroxytoluene	220	C15H24O	000128-37-0	98
3			Butylated Hydroxytoluene	220	C15H24O	000128-37-0	98
4			Butylated Hydroxytoluene	220	C15H24O	000128-37-0	97
5			Ethyl 4-oxo-2-phenylpentanoate	220	C13H16O3	1000427-11-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102722\
Data File : VV028715.D
Acq On : 27 Oct 2022 13:12
Operator : SY/MD
Sample : N5232-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHAB9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102522WMA.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
Ethyl ether	1.947	47.0	ug/L	2737330	1	5.616	291538	5.0
Diisopropyl ether	3.282	39.0	ug/L	2275660	1	5.616	291538	5.0
3-Heptanone, 6-...	10.349	5.5	ug/L	496635	3	11.243	452632	5.0
unknown-01	11.342	4.0	ug/L	365849	3	11.243	452632	5.0
Indane	11.519	5.4	ug/L	485821	3	11.243	452632	5.0
Benzene, 1-meth...	12.008	2.6	ug/L	231530	3	11.243	452632	5.0
Benzene, 1,2,3,...	12.403	2.4	ug/L	219381	3	11.243	452632	5.0
Ethanone, 1-(4-...	12.873	3.2	ug/L	289239	3	11.243	452632	5.0
Butylated Hydro...	16.297	3.6	ug/L	326896	3	11.243	452632	5.0