

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV052720\
 Data File : VV016257.D
 Acq On : 28 May 2020 03:24
 Operator : SY/MD
 Sample : L2766-10
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BE4Z3

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.315	62	70	82	rVB	43894	48416	2.54%	0.648%
2	1.409	85	99	133	rBV2	27199	166772	8.76%	2.233%
3	1.582	146	153	170	rVB2	42634	65624	3.45%	0.879%
4	1.968	264	273	286	rVB	61127	86020	4.52%	1.152%
5	2.126	313	322	339	rVB	70002	102604	5.39%	1.374%
6	2.550	446	454	469	rVB	26490	49645	2.61%	0.665%
7	2.782	513	526	547	rVB2	53783	114879	6.03%	1.538%
8	3.061	600	613	636	rVB	51037	104834	5.50%	1.404%
9	3.315	676	692	725	rVB	223070	526639	27.65%	7.052%
10	3.788	822	839	860	rBV	65177	165370	8.68%	2.214%
11	3.936	873	885	924	rBV	37148	125767	6.60%	1.684%
12	4.380	1006	1023	1052	rBV3	61041	176124	9.25%	2.358%
13	4.701	1109	1123	1142	rBV4	8973	21663	1.14%	0.290%
14	5.074	1223	1239	1244	rBV2	132621	281438	14.78%	3.769%
15	5.126	1244	1255	1275	rVB	882436	1904630	100.00%	25.505%
16	5.643	1404	1416	1443	rVB	127164	253722	13.32%	3.398%
17	6.097	1544	1557	1569	rBV	74538	150764	7.92%	2.019%
18	7.010	1825	1841	1862	rBV2	33601	62099	3.26%	0.832%
19	7.338	1932	1943	1958	rBV	156845	267362	14.04%	3.580%
20	7.415	1958	1967	1985	rVB2	16412	28992	1.52%	0.388%
21	7.646	2030	2039	2062	rVB	20513	36552	1.92%	0.489%
22	8.116	2174	2185	2215	rBV	167271	308688	16.21%	4.134%
23	8.875	2408	2421	2425	rBV	200558	310192	16.29%	4.154%
24	8.900	2426	2429	2452	rVB	216963	305912	16.06%	4.096%
25	9.035	2461	2471	2481	rBV	20189	32270	1.69%	0.432%
26	9.158	2498	2509	2525	rBV	225413	364063	19.11%	4.875%
27	9.566	2626	2636	2653	rBV	23109	37623	1.98%	0.504%
28	9.952	2742	2756	2773	rBV	94500	144494	7.59%	1.935%
29	10.238	2836	2845	2858	rVB	51943	79492	4.17%	1.064%
30	10.376	2876	2888	2903	rBV3	29360	47801	2.51%	0.640%
31	10.469	2906	2917	2921	rVV2	15428	23810	1.25%	0.319%
32	10.502	2922	2927	2937	rVV6	21380	33045	1.73%	0.443%
33	10.560	2937	2945	2957	rVB	13783	20252	1.06%	0.271%
34	10.759	2995	3007	3019	rBV	59410	90865	4.77%	1.217%

LSC Area Percent Report

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Title : TRACE VOA SOM01.0

35	10.939	3054	3063	3080	rVB	37478	58275	3.06%	0.780%
36	11.270	3156	3166	3184	rBV	209097	351738	18.47%	4.710%
37	11.357	3184	3193	3209	rVB	51358	83840	4.40%	1.123%
38	11.550	3243	3253	3272	rVV2	37636	72141	3.79%	0.966%
39	11.646	3272	3283	3303	rVB2	167857	312990	16.43%	4.191%
40	12.039	3395	3405	3421	rVB2	15933	24707	1.30%	0.331%
41	12.347	3490	3501	3512	rBV4	16299	25605	1.34%	0.343%

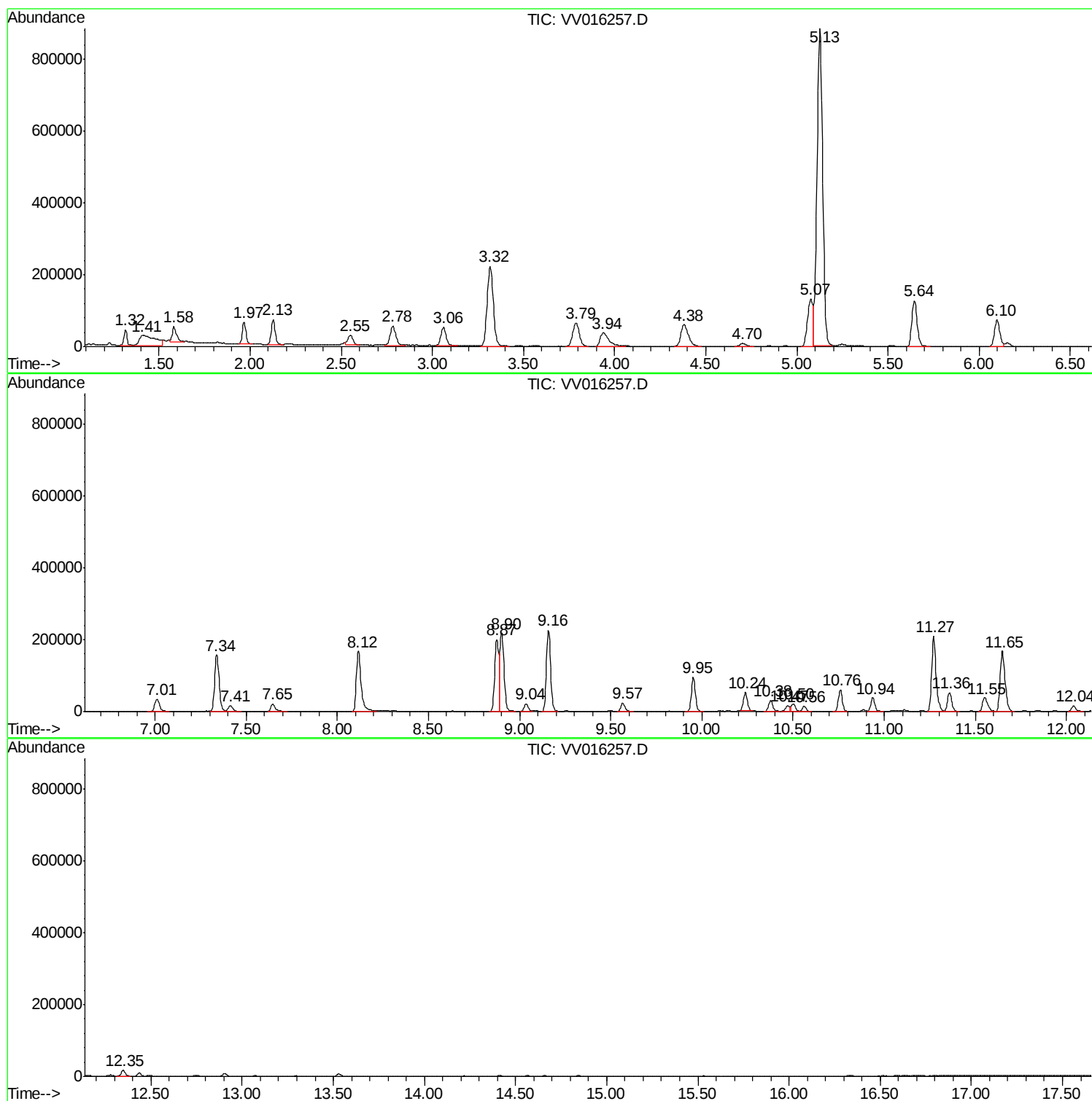
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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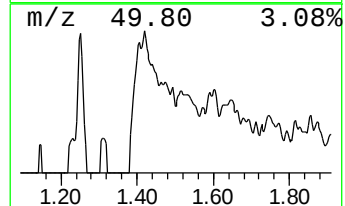
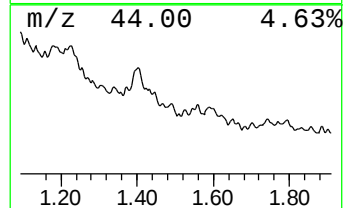
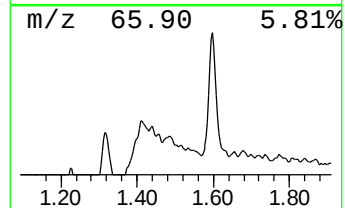
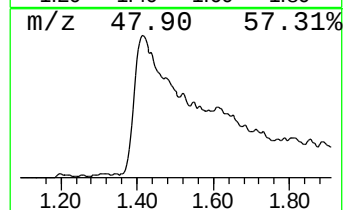
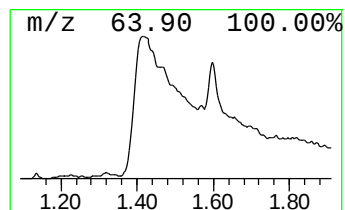
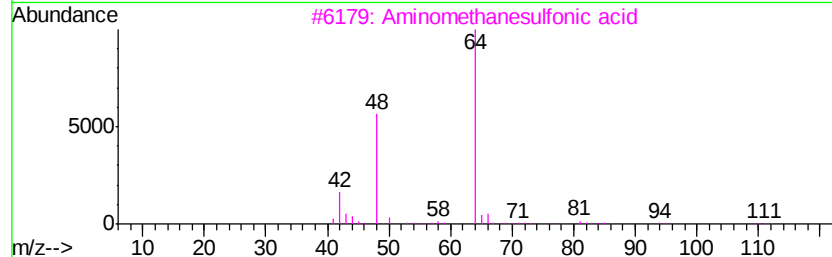
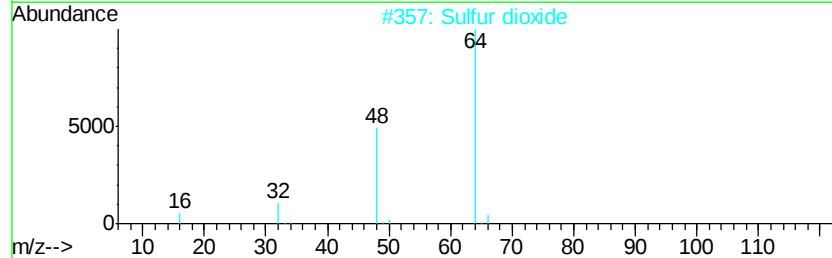
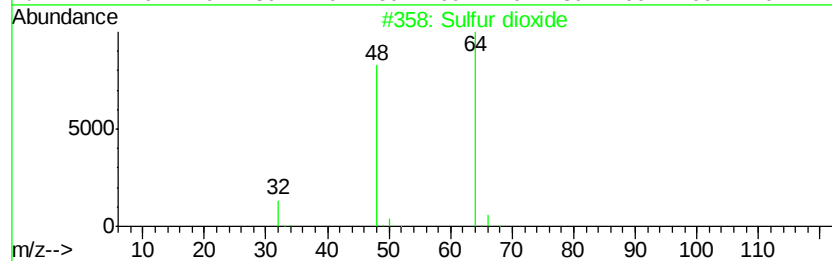
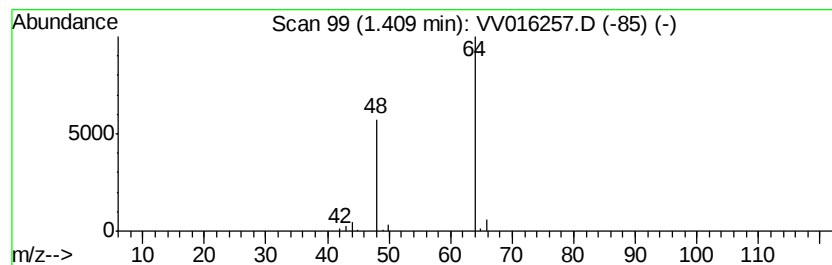
Instrument :
MSVOA_V
ClientSampled :
BE4Z3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.41	3.29 ug/L	166772	1,4-Difluorobenzene	5.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Sulfur dioxide	64 02S	007446-09-5	83
2	Sulfur dioxide	64 02S	007446-09-5	83
3	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	74
4	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	74
5	Aminomethanesulfonic acid	111 CH5N03S	013881-91-9	25



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ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BE4Z3

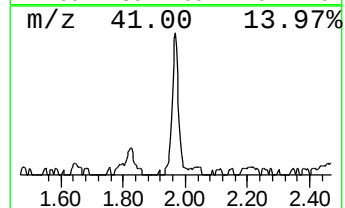
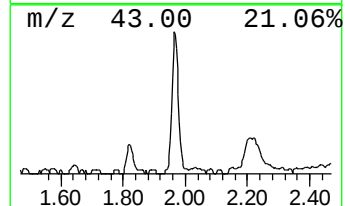
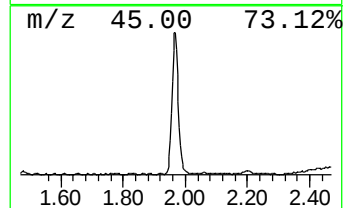
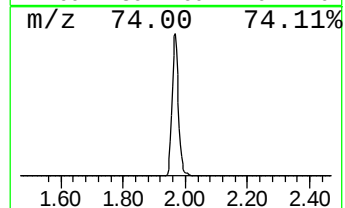
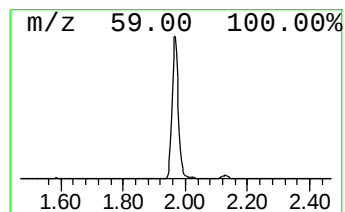
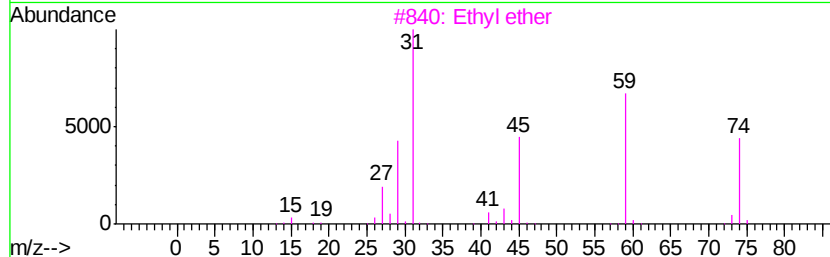
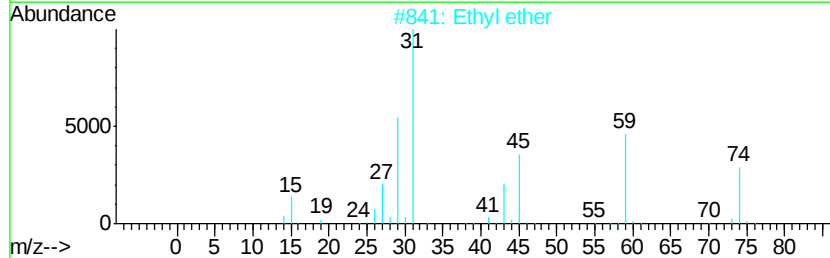
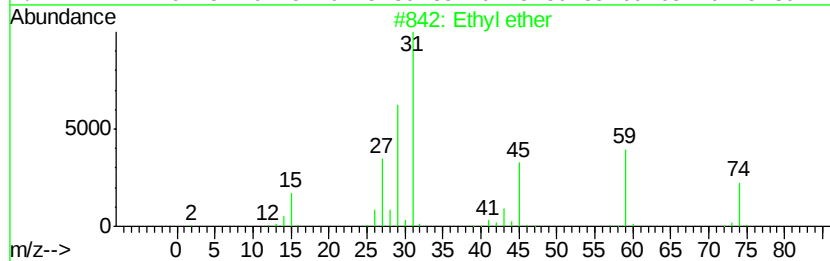
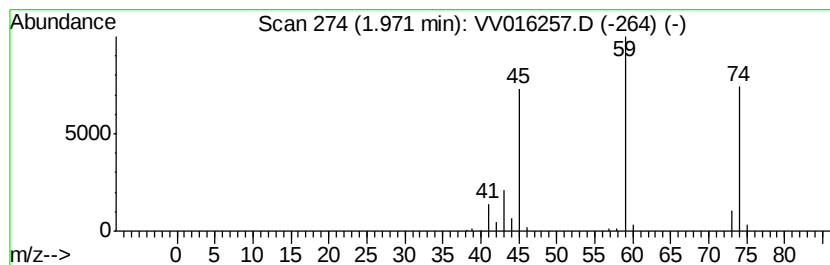
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.97	1.70 ug/L	86020	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	90
5		Ethyl ether	74	C4H10O	000060-29-7	83



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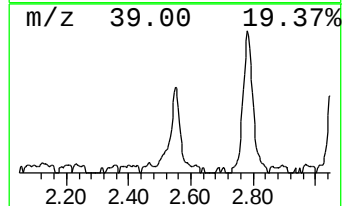
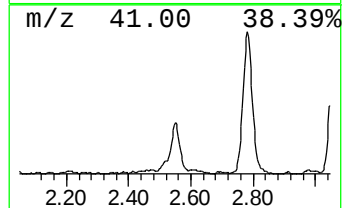
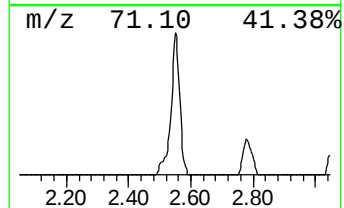
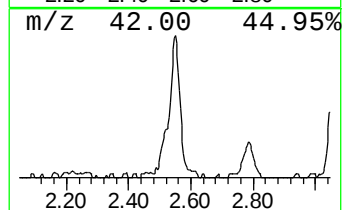
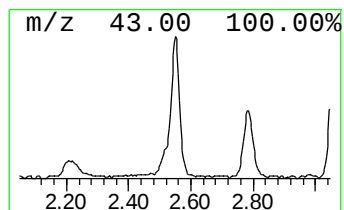
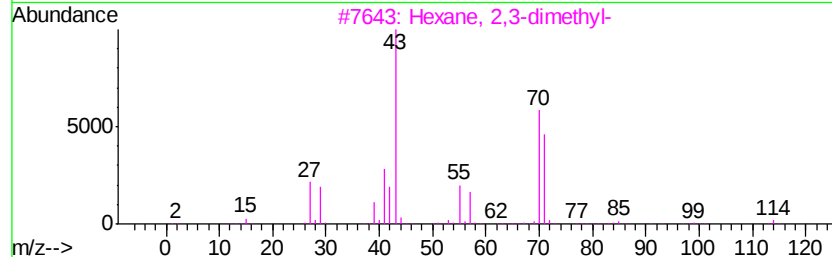
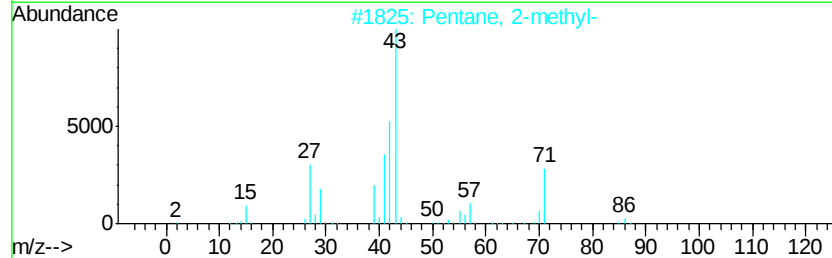
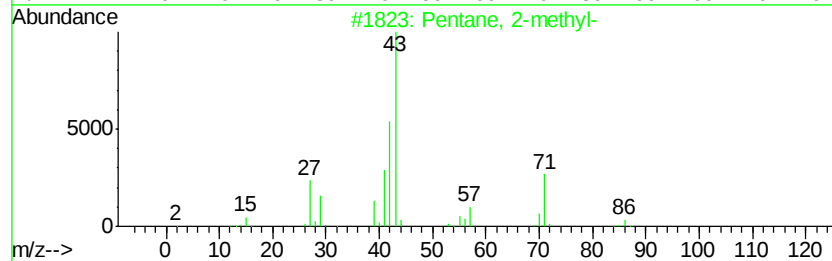
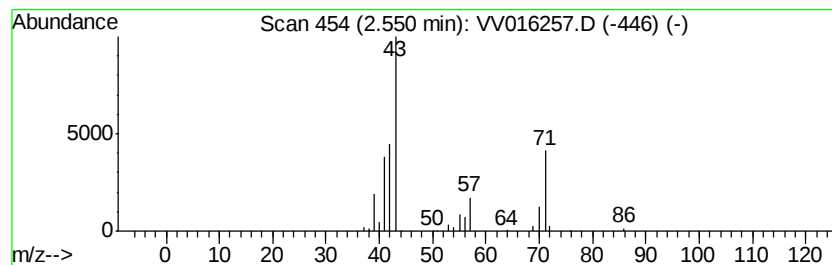
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.55	0.98 ug/L	49645	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	74
2		Pentane, 2-methyl-	86	C6H14	000107-83-5	74
3		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28
4		Pyrrolidine	71	C4H9N	000123-75-1	9
5		Pyrrolidine	71	C4H9N	000123-75-1	9



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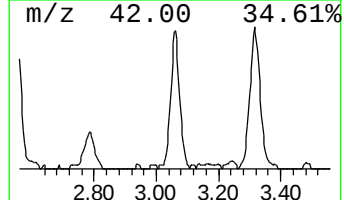
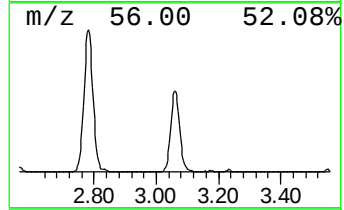
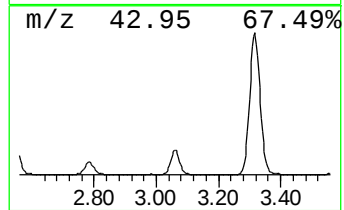
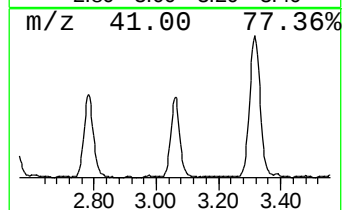
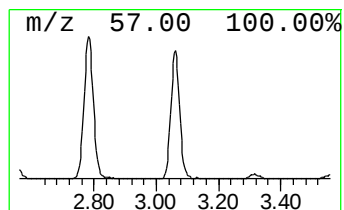
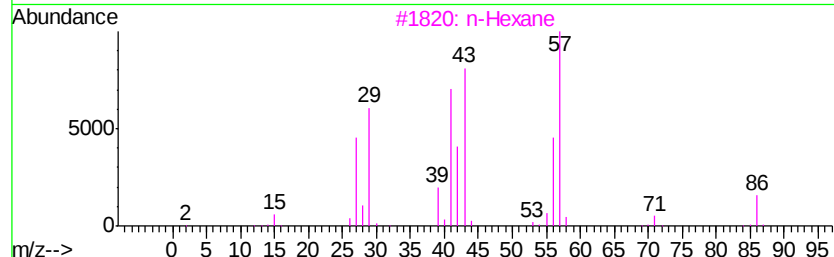
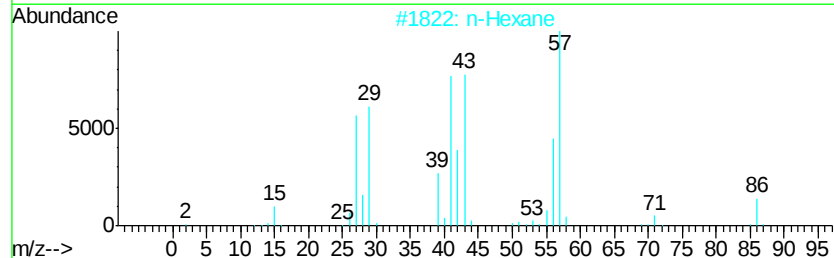
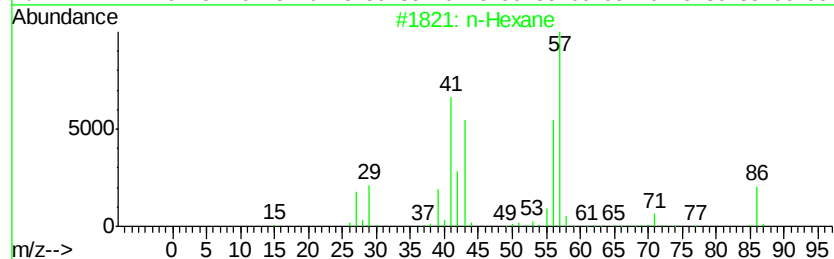
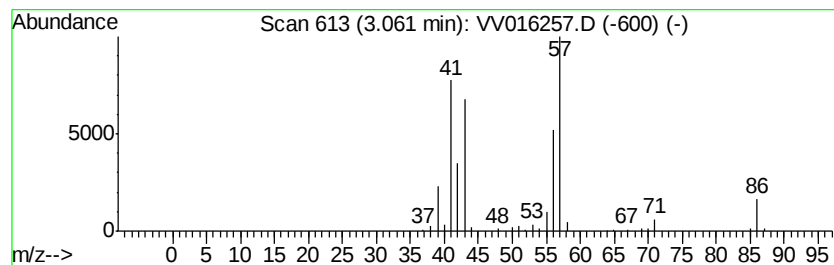
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	2.07 ug/L	104834	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	91
2		n-Hexane	86	C6H14	000110-54-3	91
3		n-Hexane	86	C6H14	000110-54-3	91
4		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
5		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	47



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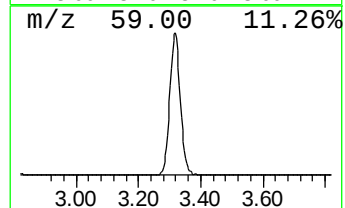
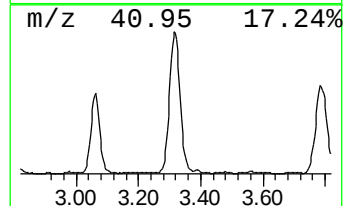
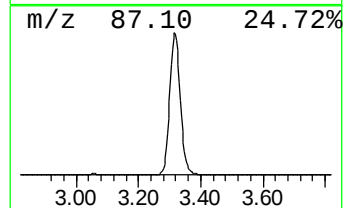
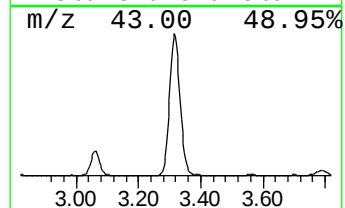
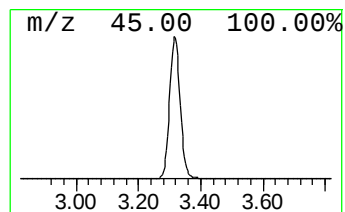
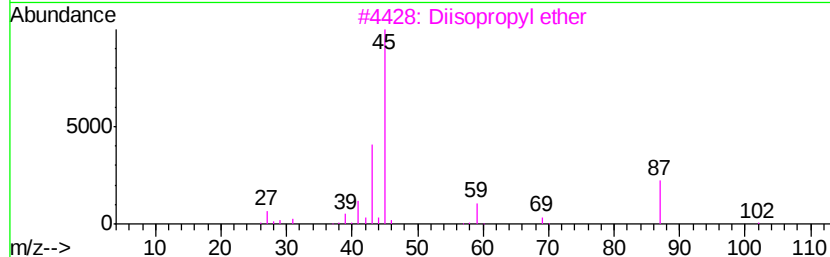
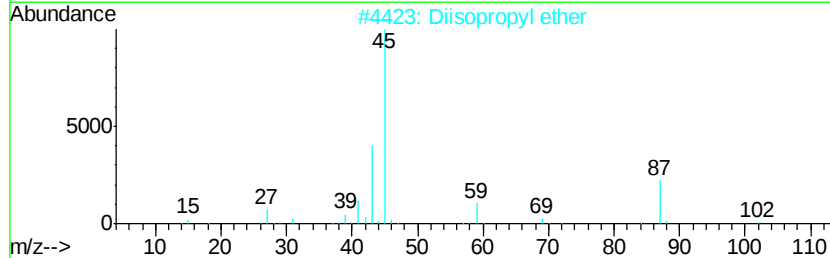
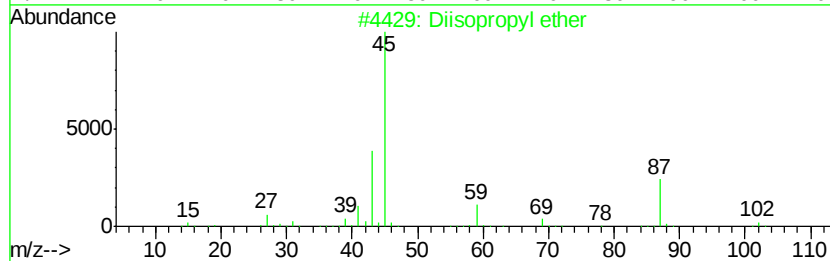
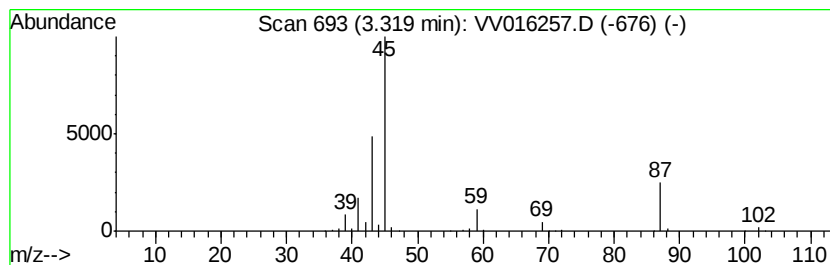
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Diisopropyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.32	10.38 ug/L	526639	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diisopropyl ether	102	C6H14O	000108-20-3	91
2		Diisopropyl ether	102	C6H14O	000108-20-3	83
3		Diisopropyl ether	102	C6H14O	000108-20-3	83
4		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	83
5		Ether, sec-butyl isopropyl	116	C7H16O	018641-81-1	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BE4Z3

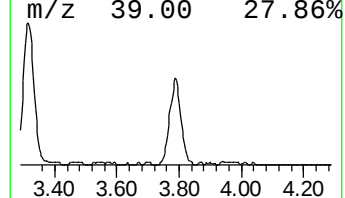
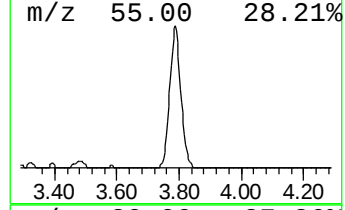
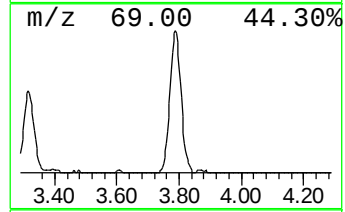
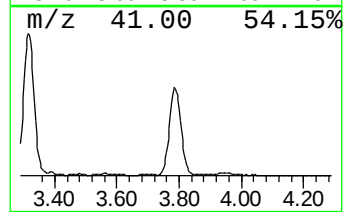
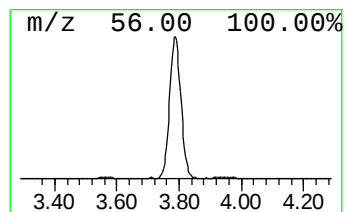
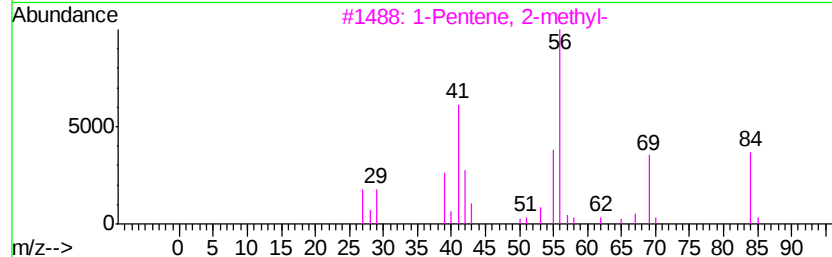
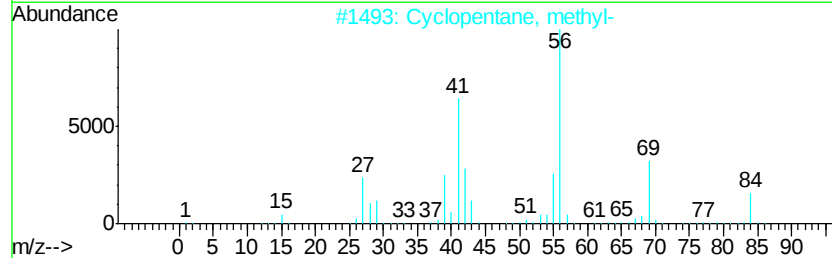
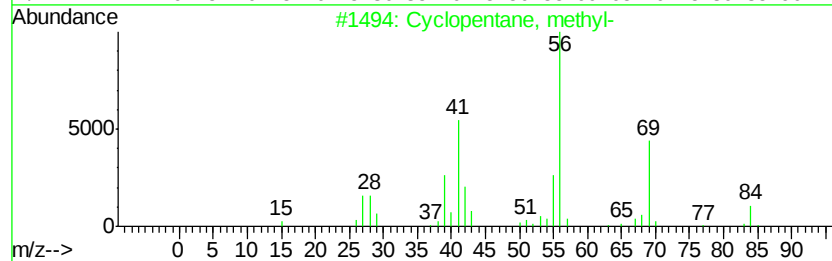
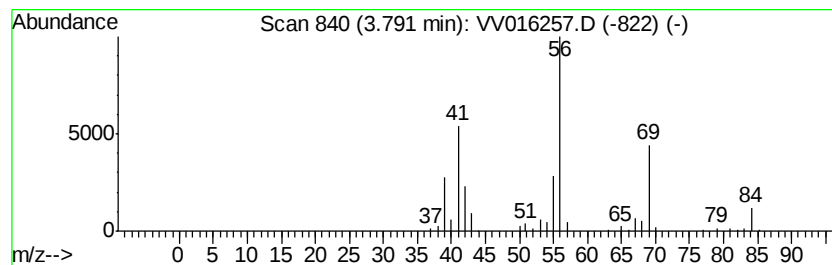
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic3.79 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.79	3.26 ug/L	165370	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	87
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampled :
BE4Z3

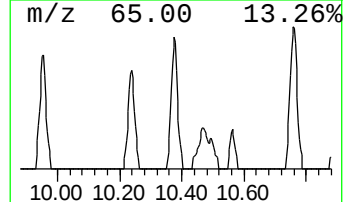
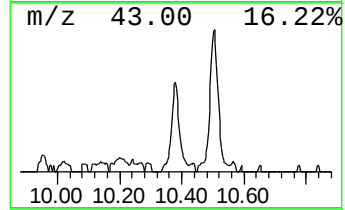
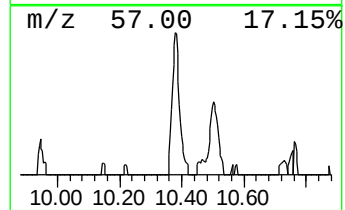
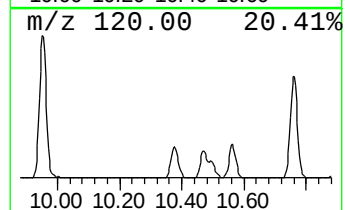
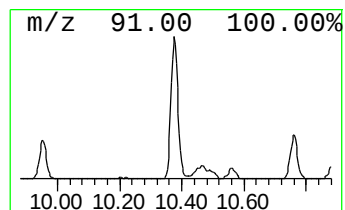
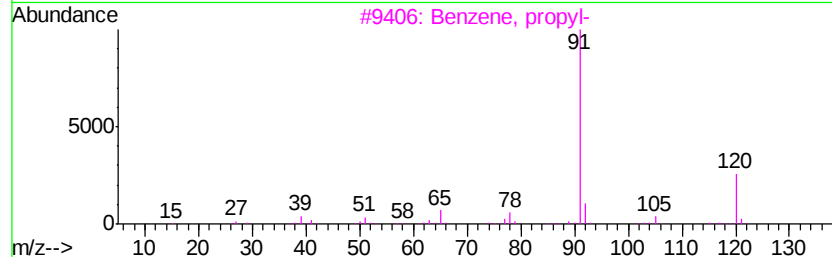
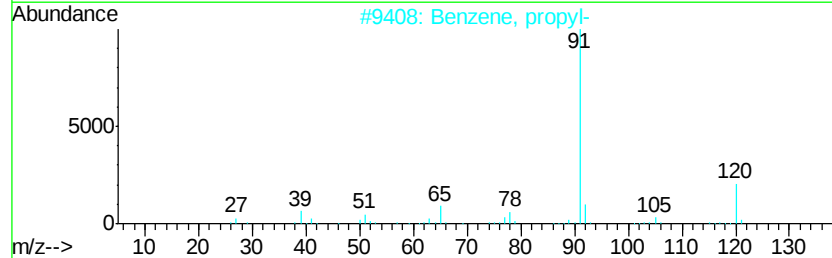
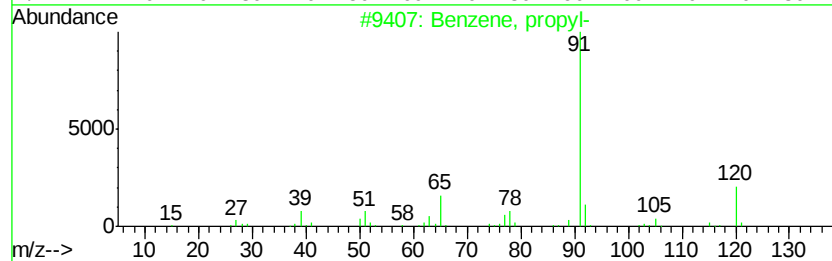
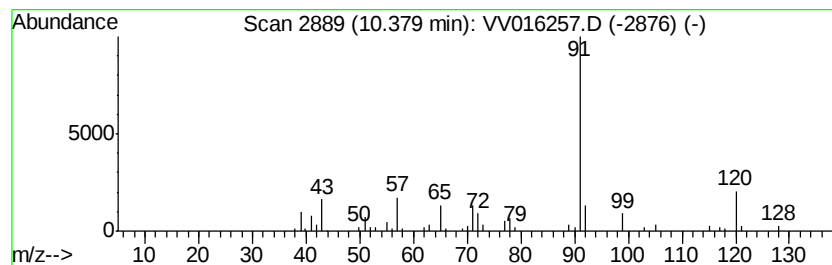
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	0.68 ug/L	47801	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	64
2		Benzene, propyl-	120	C9H12	000103-65-1	64
3		Benzene, propyl-	120	C9H12	000103-65-1	58
4		Benzeneacetaldehyde	120	C8H8O	000122-78-1	55
5		1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampled :
BE4Z3

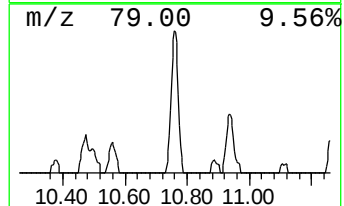
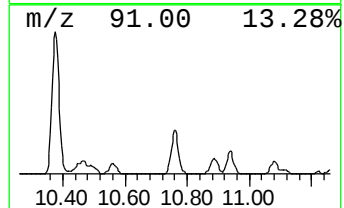
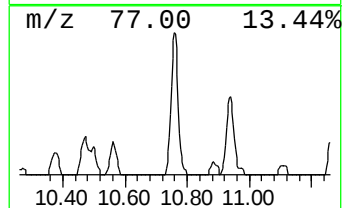
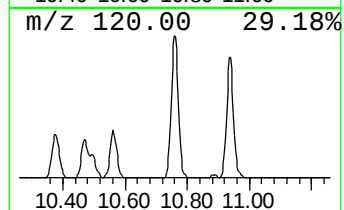
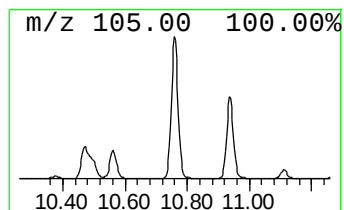
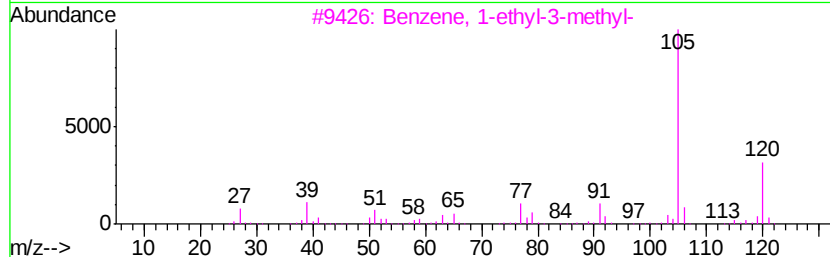
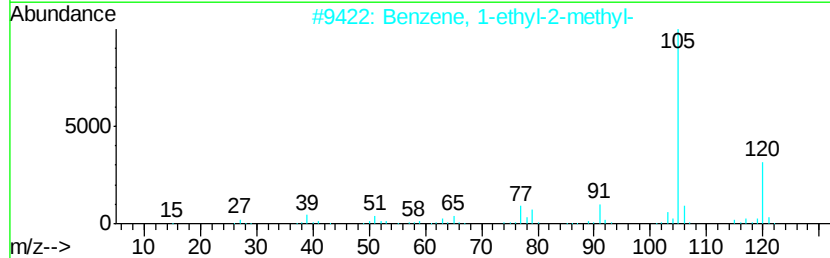
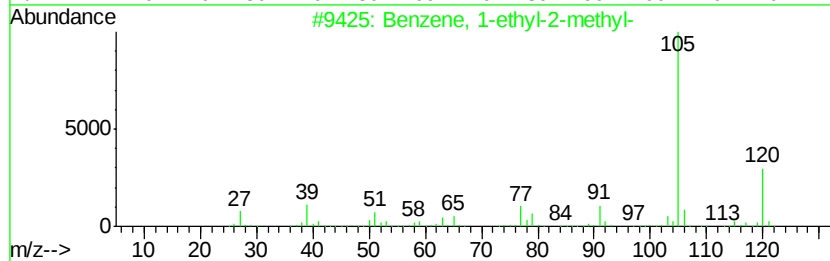
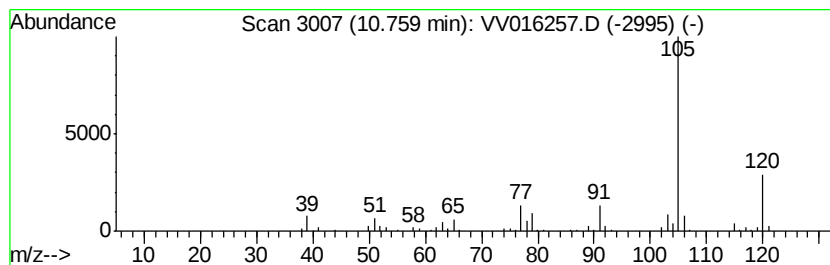
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	1.29 ug/L	90865	1,4-Dichlorobenzene-d4	11.27

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-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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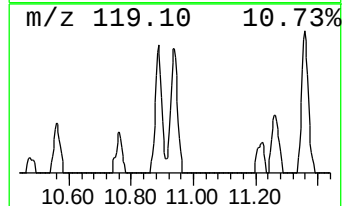
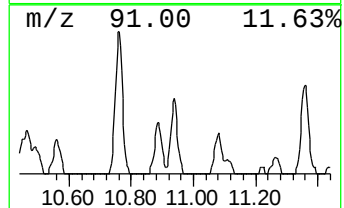
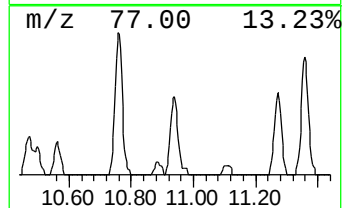
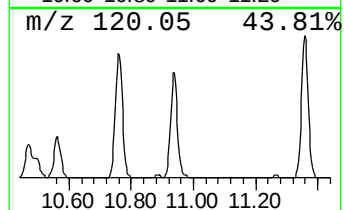
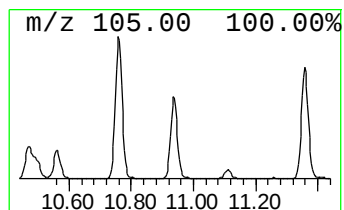
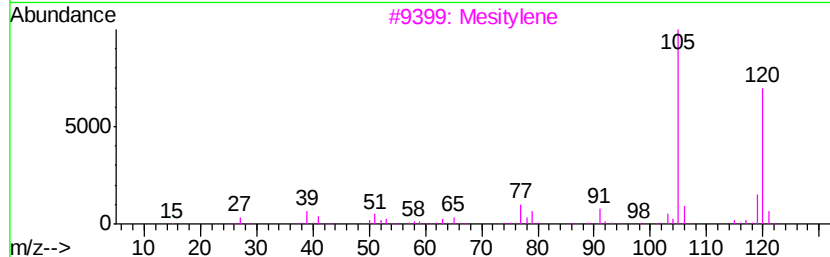
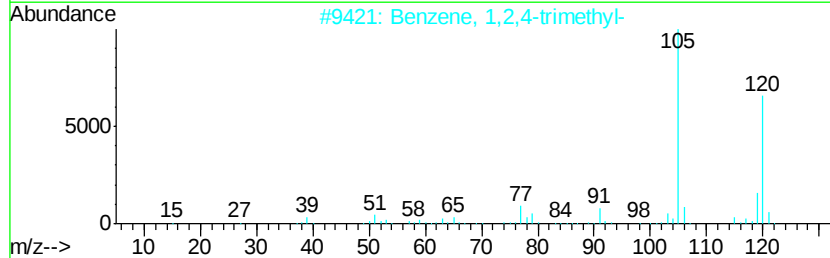
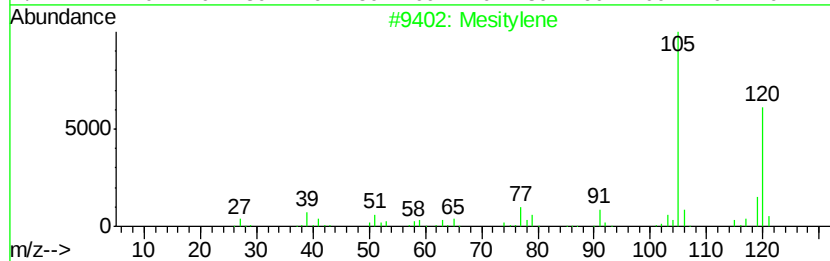
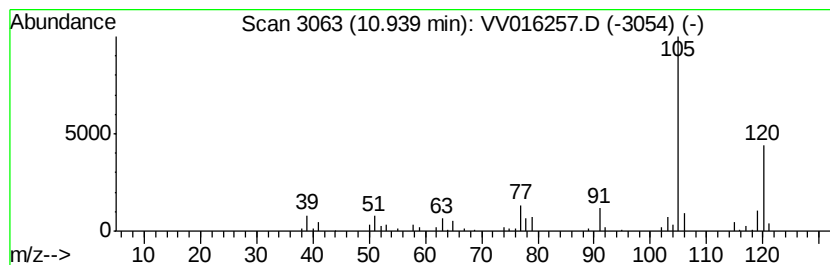
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Mesitylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	0.83 ug/L	58275	1,4-Dichlorobenzene-d4	11.27

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BE4Z3

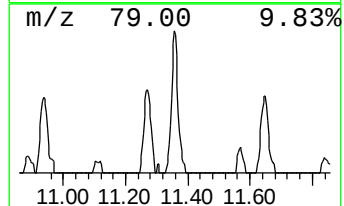
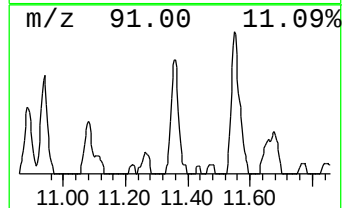
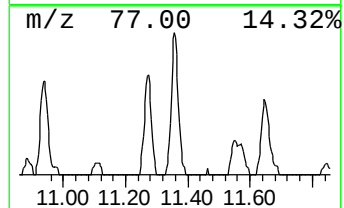
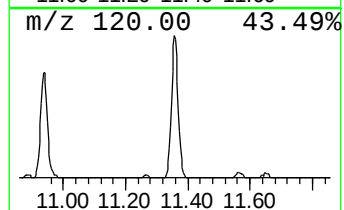
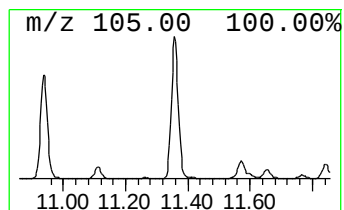
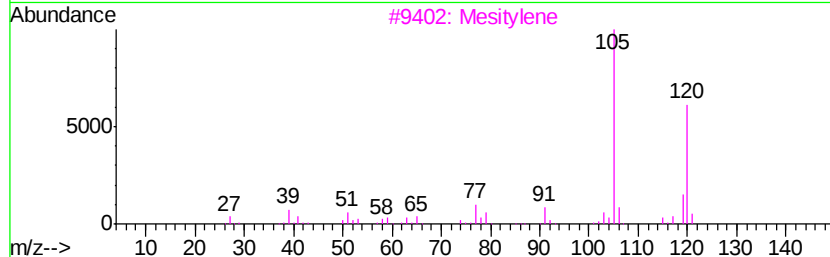
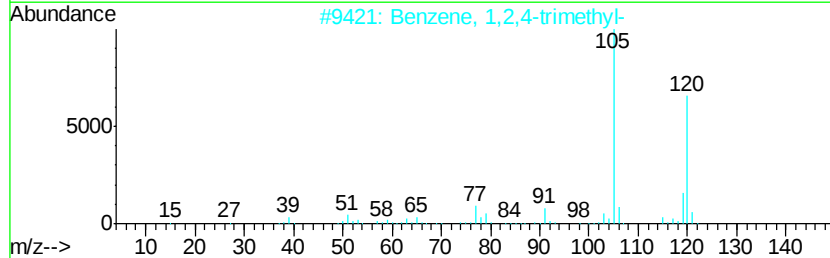
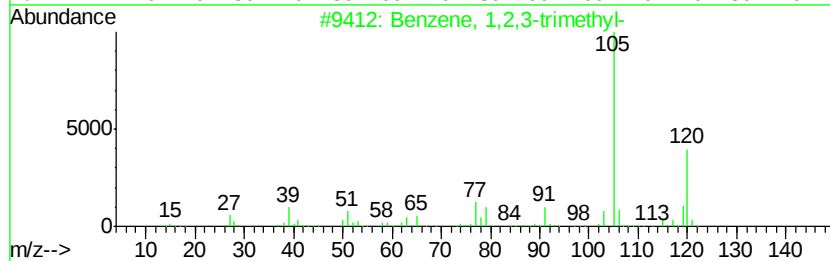
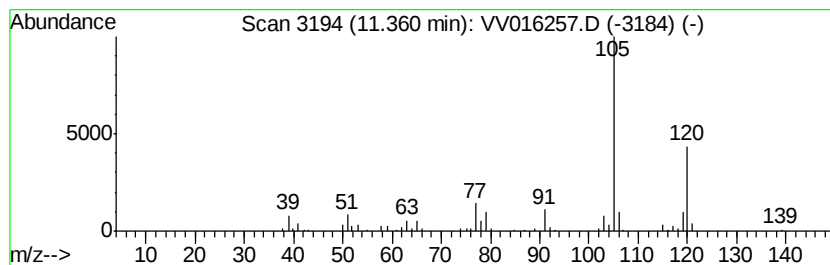
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	1.19 ug/L	83840	1,4-Dichlorobenzene-d4	11.27

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampled :
BE4Z3

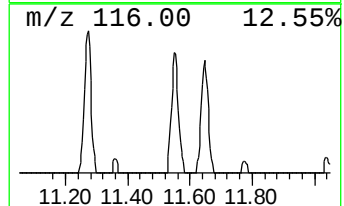
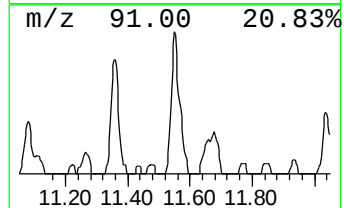
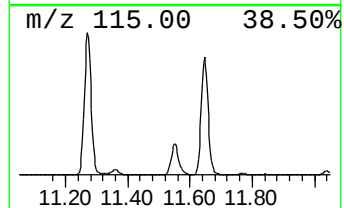
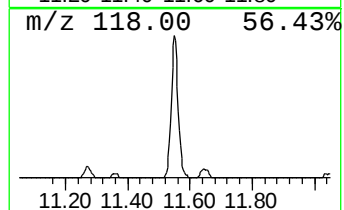
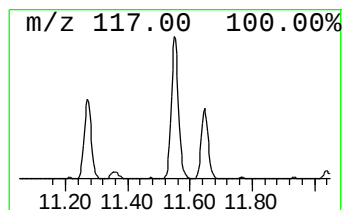
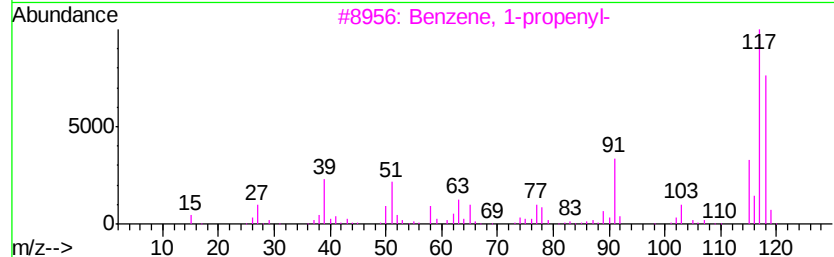
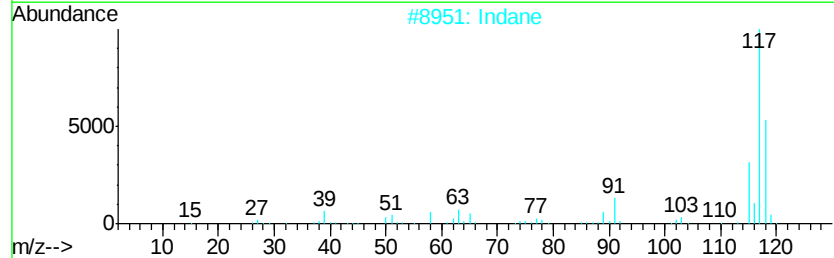
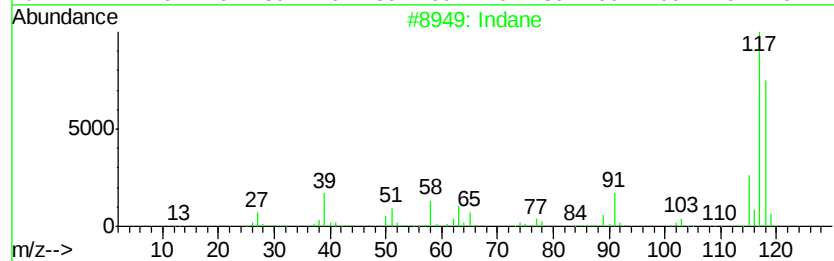
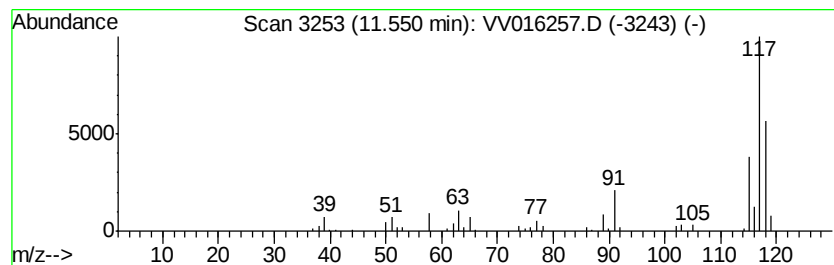
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	1.03 ug/L	72141	1,4-Dichlorobenzene-d4	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Indane		118	C9H10	000496-11-7	93
3	Benzene, 1-propenyl-		118	C9H10	000637-50-3	74
4	Indane		118	C9H10	000496-11-7	70
5	(Z)-1-Phenylpropene		118	C9H10	000766-90-5	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052720\
Data File : VV016257.D
Acq On : 28 May 2020 03:24
Operator : SY/MD
Sample : L2766-10
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BE4Z3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Sulfur dioxide	1.41	3.3	ug/L	166772	1	5.64	253722	5.0
Ethyl ether	1.97	1.7	ug/L	86020	1	5.64	253722	5.0
(DEL) Alkane: Str...	2.55	1.0	ug/L	49645	1	5.64	253722	5.0
(DEL) Alkane: Str...	3.06	2.1	ug/L	104834	1	5.64	253722	5.0
Diisopropyl ether	3.32	10.4	ug/L	526639	1	5.64	253722	5.0
(DEL) Alkane: Cyc...	3.79	3.3	ug/L	165370	1	5.64	253722	5.0
Benzene, propyl-	10.38	0.7	ug/L	47801	3	11.27	351738	5.0
Benzene, 1-ethyl-...	10.76	1.3	ug/L	90865	3	11.27	351738	5.0
Mesitylene	10.94	0.8	ug/L	58275	3	11.27	351738	5.0
Benzene, 1,2,3-tr...	11.36	1.2	ug/L	83840	3	11.27	351738	5.0
Indane	11.55	1.0	ug/L	72141	3	11.27	351738	5.0