

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU060120\
 Data File : VU038466.D
 Acq On : 01 Jun 2020 13:24
 Operator : JC/MD
 Sample : L2789-04
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFX77

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.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	1.369	3	9	17	rVB	67864	65636	4.17%	0.573%
2	1.610	74	84	94	rBV	114556	133604	8.49%	1.167%
3	1.932	175	184	196	rBV	100607	132446	8.41%	1.157%
4	2.594	377	390	406	rBV2	234686	432414	27.47%	3.776%
5	3.395	624	639	653	rVB4	10424	24351	1.55%	0.213%
6	3.735	731	745	761	rBV4	9137	19940	1.27%	0.174%
7	4.571	989	1005	1019	rVB4	10844	26368	1.67%	0.230%
8	4.684	1019	1040	1077	rBV	187704	703341	44.68%	6.142%
9	4.848	1077	1091	1124	rVB	155143	392305	24.92%	3.426%
10	5.102	1152	1170	1198	rBV4	250076	649428	41.25%	5.671%
11	5.761	1353	1375	1397	rBV2	409681	1088959	69.17%	9.509%
12	6.279	1521	1536	1556	rBV	365920	704454	44.75%	6.152%
13	6.571	1612	1627	1645	rBV	107864	203808	12.95%	1.780%
14	6.722	1656	1674	1698	rBV	256444	498269	31.65%	4.351%
15	7.070	1768	1782	1802	rBV	173447	317221	20.15%	2.770%
16	7.590	1931	1944	1961	rBV2	132936	235834	14.98%	2.059%
17	7.922	2033	2047	2083	rBV	471833	854260	54.26%	7.460%
18	8.201	2117	2134	2151	rBV	87514	149149	9.47%	1.302%
19	8.574	2236	2250	2261	rBV2	56847	97775	6.21%	0.854%
20	8.658	2261	2276	2317	rVV	837578	1574331	100.00%	13.748%
21	8.960	2357	2370	2389	rBV4	8480	22325	1.42%	0.195%
22	9.436	2504	2518	2555	rVB	522593	889606	56.51%	7.769%
23	10.770	2919	2933	2948	rBV	222430	358738	22.79%	3.133%
24	11.825	3247	3261	3276	rBV	554427	867013	55.07%	7.571%
25	12.082	3329	3341	3351	rBV5	9458	19479	1.24%	0.170%
26	12.204	3365	3379	3395	rVB	516956	835178	53.05%	7.293%
27	17.388	4980	4991	5004	rBV2	79743	155132	9.85%	1.355%

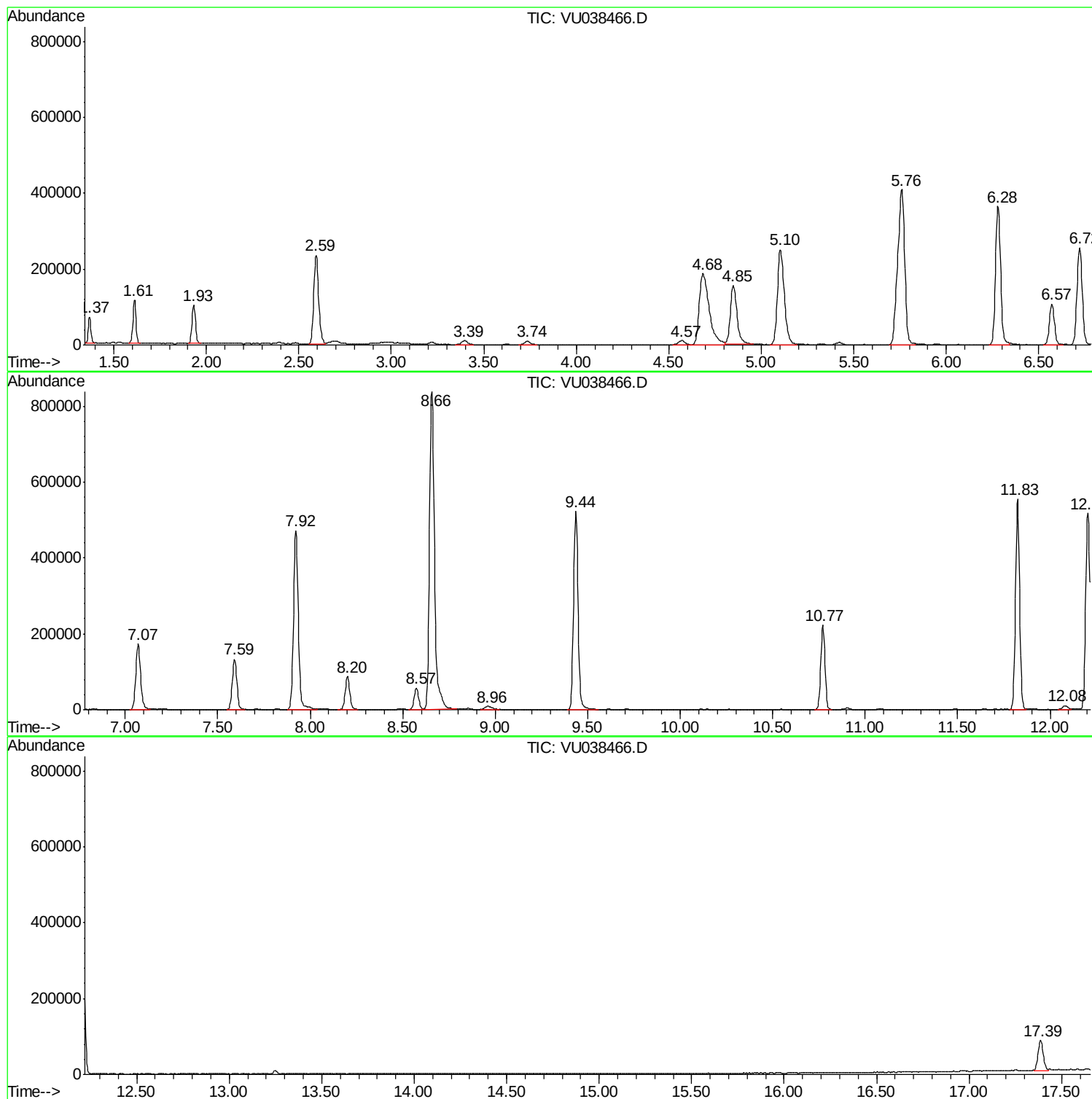
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Instrument :
MSVOA_U
ClientSampleId :
BFX77

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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



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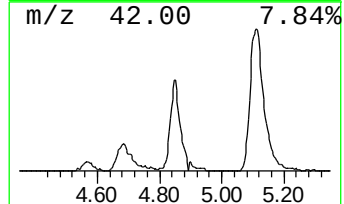
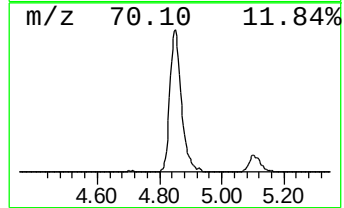
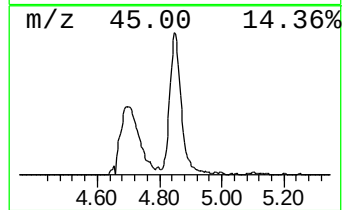
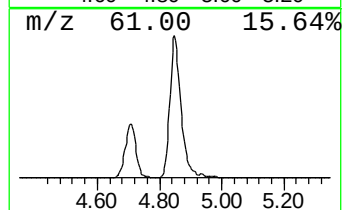
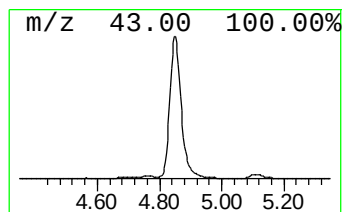
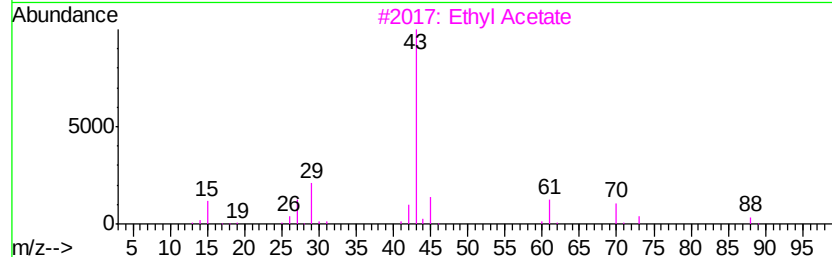
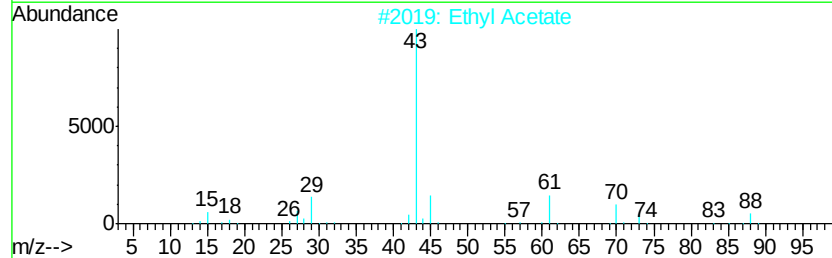
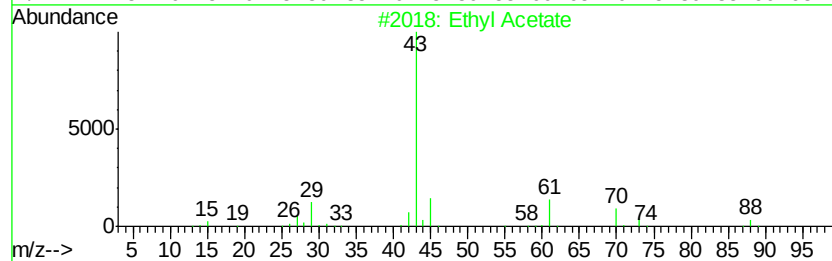
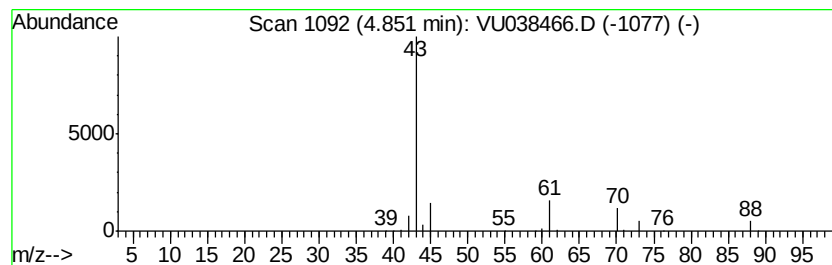
Instrument :
MSVOA_U
ClientSampleId :
BFX77

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Ethyl Acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
4.85	2.78 ug/L	392305	1,4-Difluorobenzene	6.28		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl Acetate		88	C4H8O2	000141-78-6	86
2	Ethyl Acetate		88	C4H8O2	000141-78-6	86
3	Ethyl Acetate		88	C4H8O2	000141-78-6	86
4	Ethyl Acetate		88	C4H8O2	000141-78-6	78
5	Propanoic acid, 2-methyl-		88	C4H8O2	000079-31-2	37



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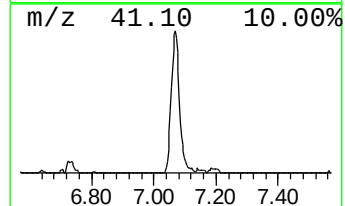
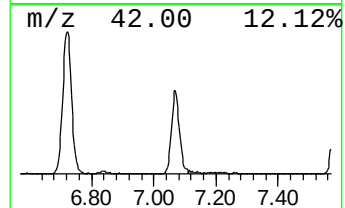
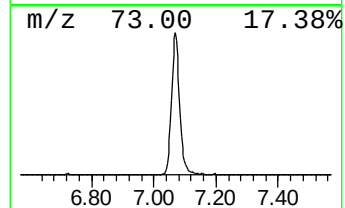
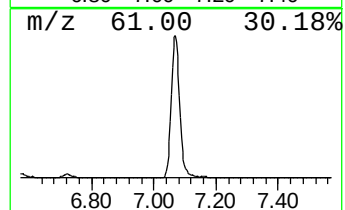
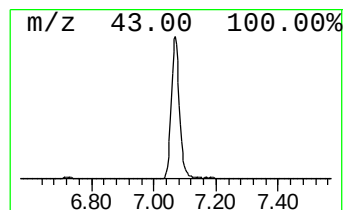
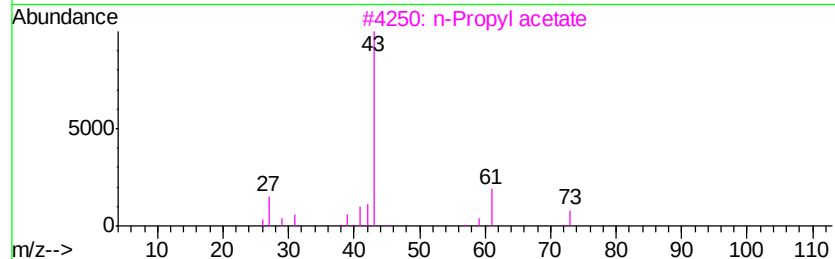
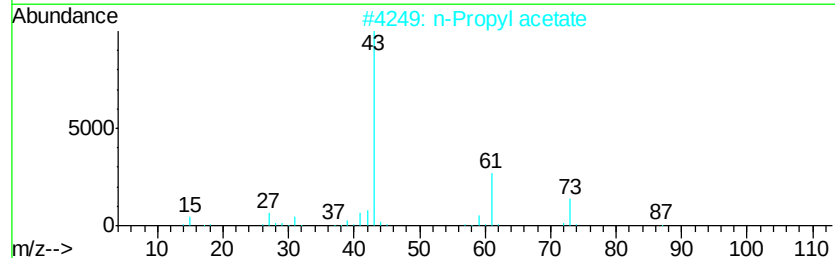
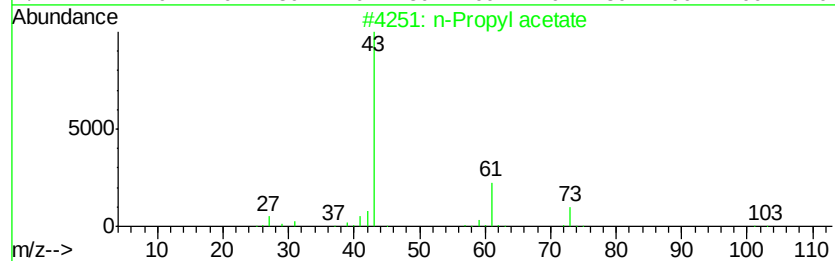
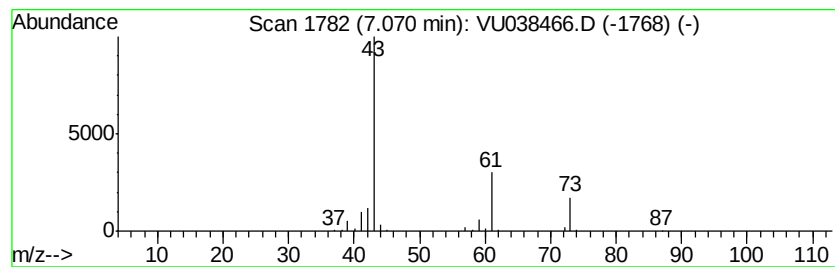
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 n-Propyl acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	2.25 ug/L	317221	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		n-Propyl acetate	102	C5H10O2	000109-60-4	83
3		n-Propyl acetate	102	C5H10O2	000109-60-4	78
4		Isopropyl acetate	102	C5H10O2	000108-21-4	36
5		Isopropyl acetate	102	C5H10O2	000108-21-4	33



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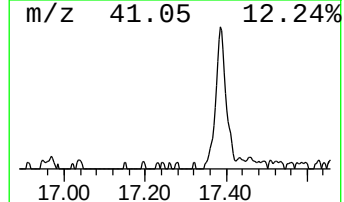
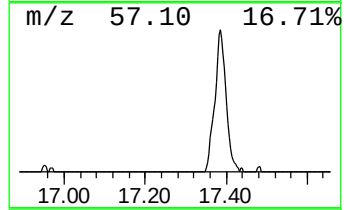
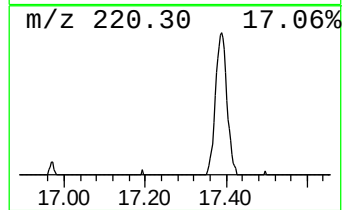
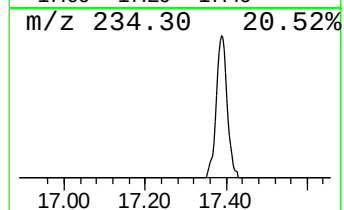
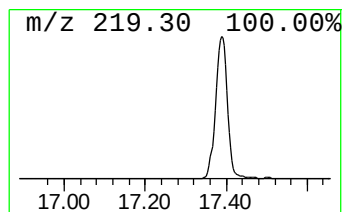
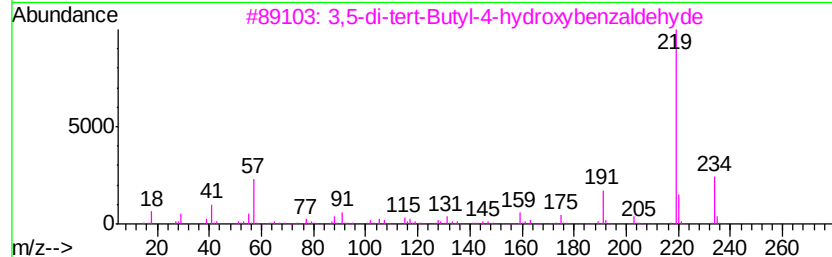
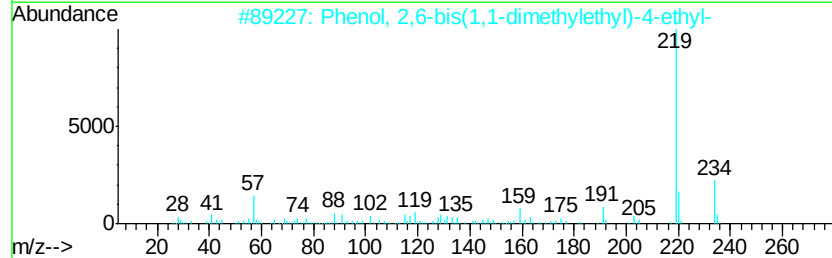
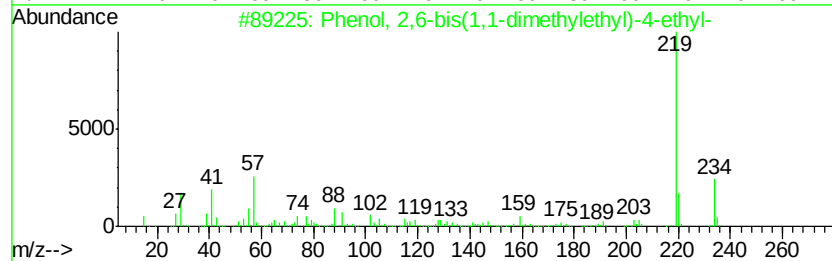
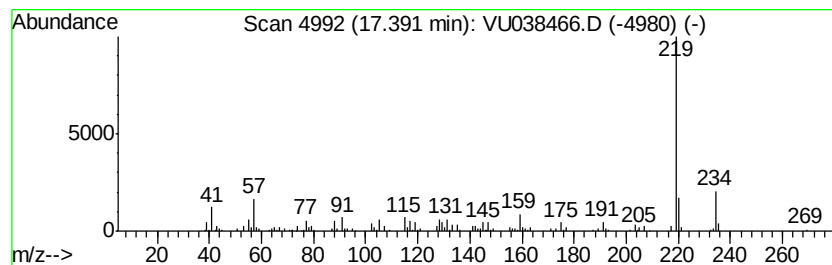
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.39	0.89 ug/L	155132	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,6-bis(1,1-dimethylethy...	234	C16H26O	004130-42-1	94
2		Phenol, 2,6-bis(1,1-dimethylethy...	234	C16H26O	004130-42-1	93
3		3,5-di-tert-Butyl-4-hydroxybenza...	234	C15H22O2	001620-98-0	90
4		3,5-di-tert-Butyl-4-hydroxybenza...	234	C15H22O2	001620-98-0	90
5		4-tert-Butyl-2,6-diisopropylphenol	234	C16H26O	057354-65-1	74



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethyl Acetate	4.85	2.8	ug/L	392305	1	6.28	704454	5.0
n-Propyl acetate	7.07	2.3	ug/L	317221	1	6.28	704454	5.0
Phenol, 2,6-bis(1...	17.39	0.9	ug/L	155132	3	11.82	867013	5.0