

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100318\
 Data File : VX004963.D
 Acq On : 03 Oct 2018 12:27
 Operator : JC/MD
 Sample : J5170-07
 Misc : 6.63g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 HC-SOIL-2

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 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_X\METHOD\82X092618W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.666	62	84	88	rBV5	36566	167125	11.23%	2.130%
2	3.019	299	306	317	rVB	14345	34915	2.35%	0.445%
3	5.507	699	714	728	rBV	162035	481526	32.35%	6.137%
4	5.659	728	739	752	rBV	216058	631705	42.44%	8.051%
5	6.068	795	806	824	rVB2	176628	480753	32.30%	6.127%
6	6.854	925	935	970	rBV	366813	921800	61.94%	11.748%
7	8.714	1229	1240	1257	rBV	843035	1488310	100.00%	18.967%
8	10.116	1464	1470	1488	rBV	766078	1143416	76.83%	14.572%
9	10.360	1505	1510	1516	rVB	12961	18893	1.27%	0.241%
10	11.140	1632	1638	1651	rBV	812914	1083532	72.80%	13.809%
11	11.439	1682	1687	1695	rBV3	9156	22454	1.51%	0.286%
12	11.811	1744	1748	1755	rVB	14373	18452	1.24%	0.235%
13	12.079	1787	1792	1801	rBV	1089487	1353780	90.96%	17.253%

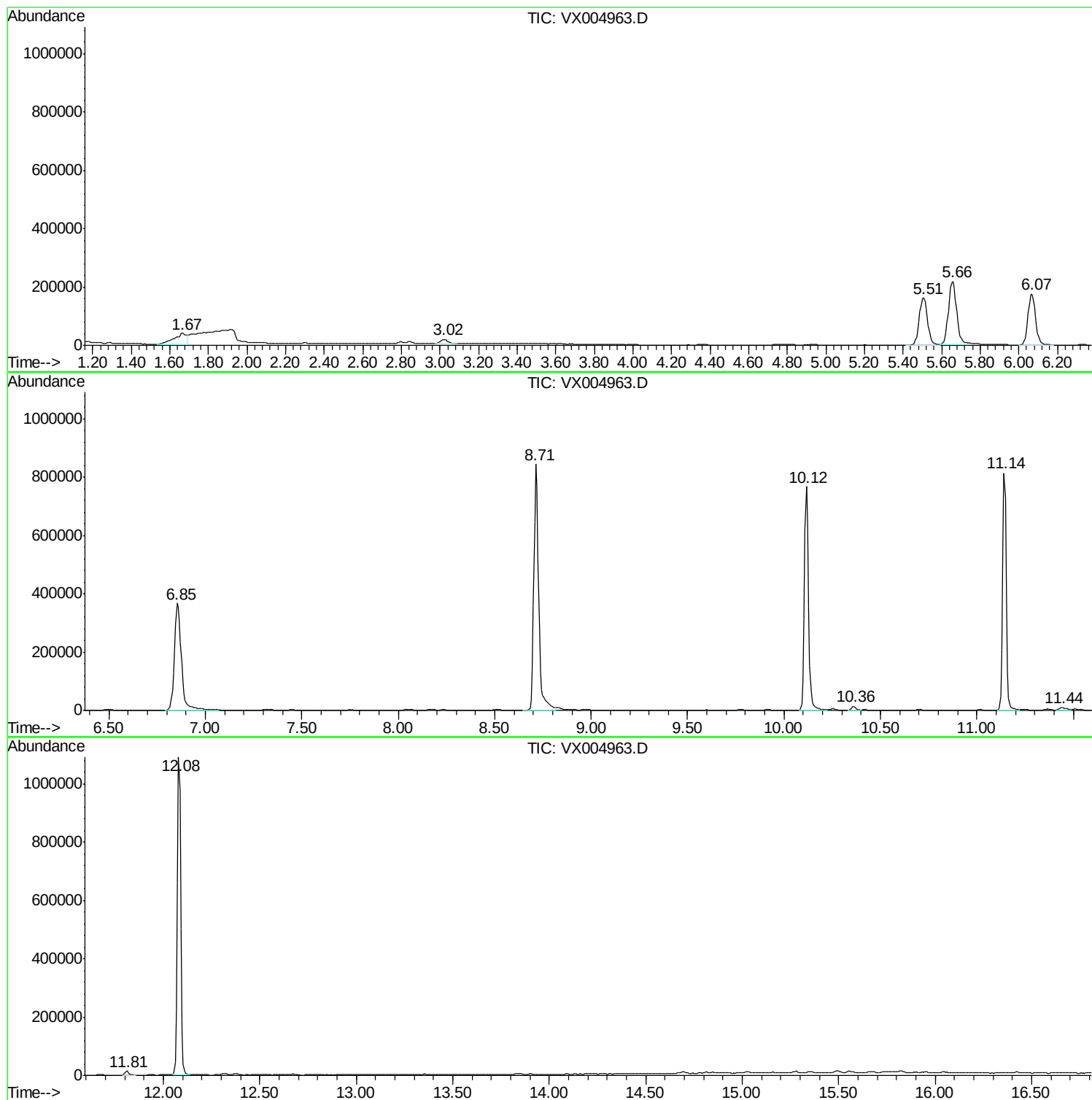
Sum of corrected areas: 7846661

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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



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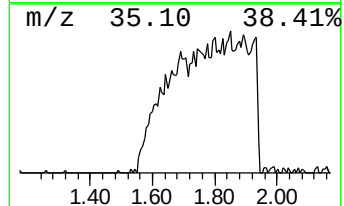
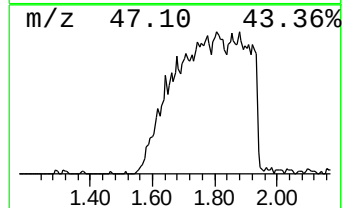
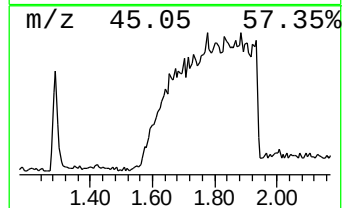
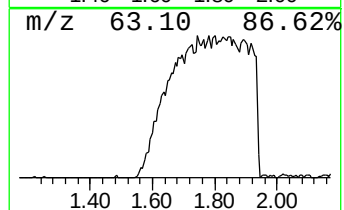
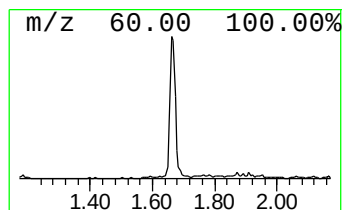
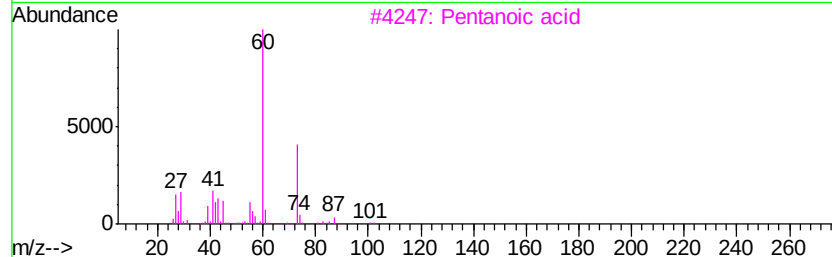
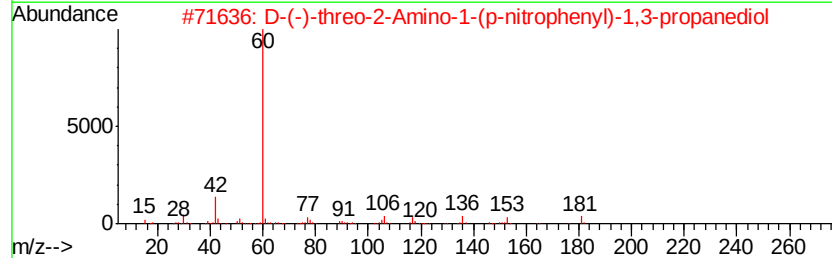
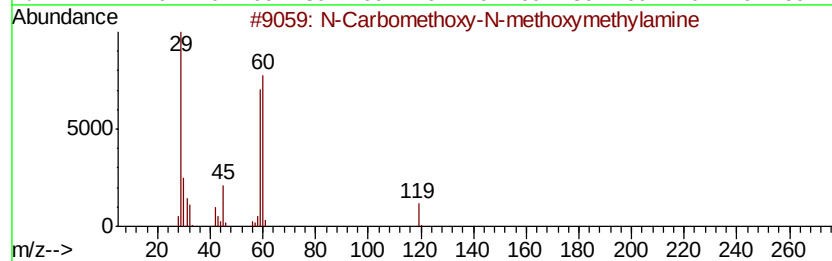
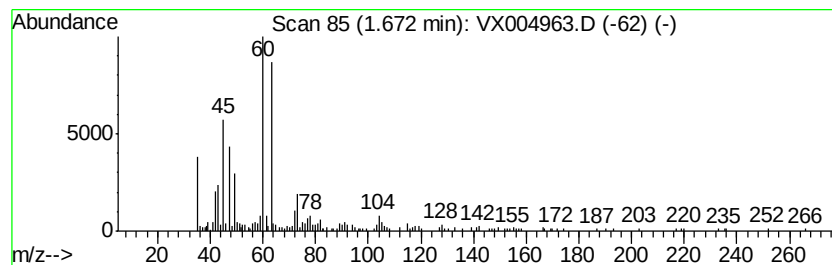
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Peak Number 1 unknown1.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.67	13.23 ug/l	167125	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-Carbomethoxy-N-methoxymethylamine	119	C4H9NO3	153654-07-0	11
2		D-(-)-threo-2-Amino-1-(p-nitroph...	212	C9H12N2O4	000716-61-0	10
3		Pentanoic acid	102	C5H10O2	000109-52-4	9
4		Hexanoic acid, 6-bromo-	194	C6H11BrO2	004224-70-8	9
5		Imidodicarbonic diamide	103	C2H5N3O2	000108-19-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.67	1.67	13.2	ug/l	167125	1	5.66	631705	50.0