

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090418\
 Data File : BG036560.D
 Acq On : 5 Sep 2018 10:44
 Operator : JU/SJ
 Sample : J4759-16
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 083018-JETT-F-D-M

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082318.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.158	313	318	330	rBV	73345	113135	2.59%	0.728%
2	8.055	805	811	819	rVB	219041	370250	8.48%	2.384%
3	8.378	858	866	879	rBV	821126	1449910	33.20%	9.334%
4	9.218	1000	1009	1021	rBV	699913	1260297	28.85%	8.113%
5	10.863	1281	1289	1299	rBV	292997	525875	12.04%	3.385%
6	13.308	1696	1705	1715	rBV	1920967	3010678	68.93%	19.382%
7	14.124	1839	1844	1852	rBV	153087	225438	5.16%	1.451%
8	14.683	1932	1939	1947	rBV2	476118	772950	17.70%	4.976%
9	17.421	2398	2405	2414	rBV2	664807	1022564	23.41%	6.583%
10	20.029	2843	2849	2859	rBV	3349216	4367717	100.00%	28.118%
11	21.686	3124	3131	3140	rBV2	720302	1184198	27.11%	7.624%
12	23.989	3518	3523	3529	rBV5	21144	47419	1.09%	0.305%
13	24.900	3667	3678	3695	rBV2	358999	1133479	25.95%	7.297%
14	26.081	3873	3879	3888	rVB7	18403	49449	1.13%	0.318%

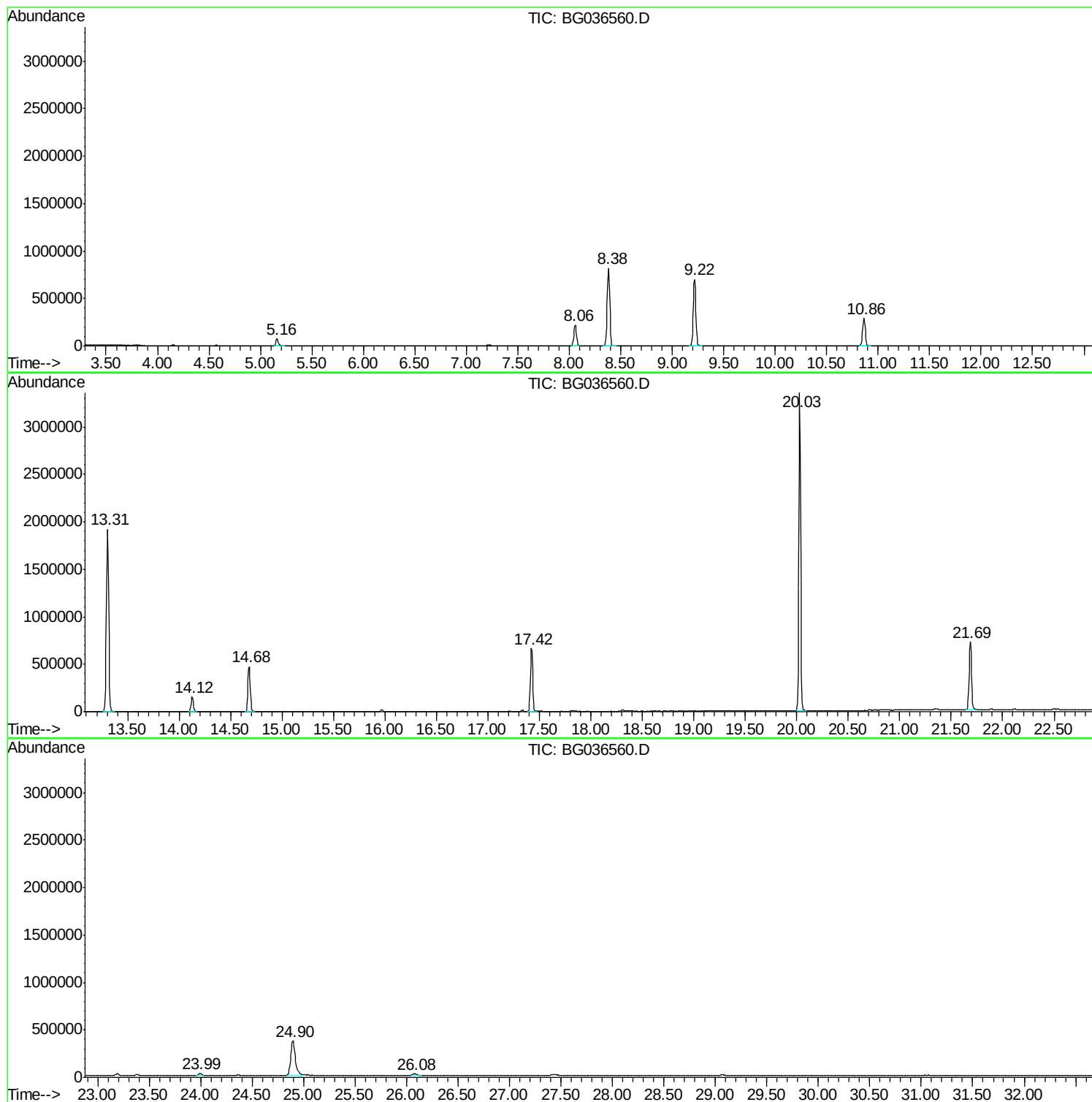
Sum of corrected areas: 15533359

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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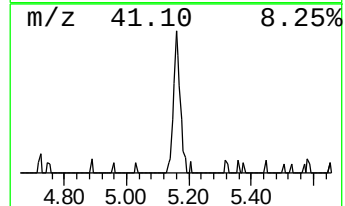
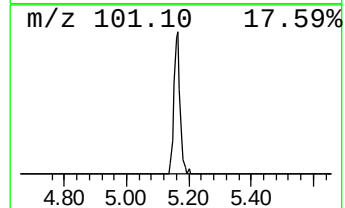
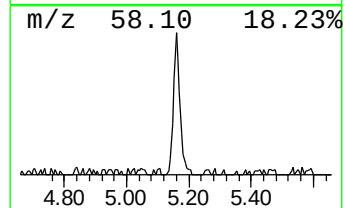
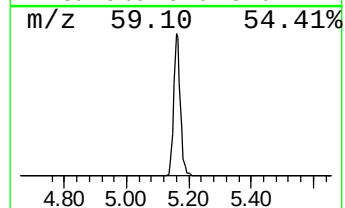
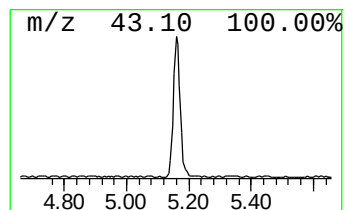
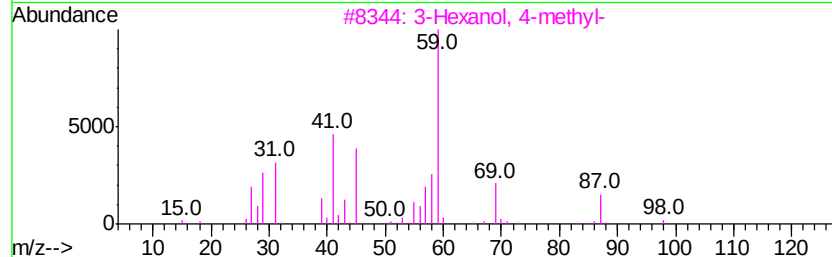
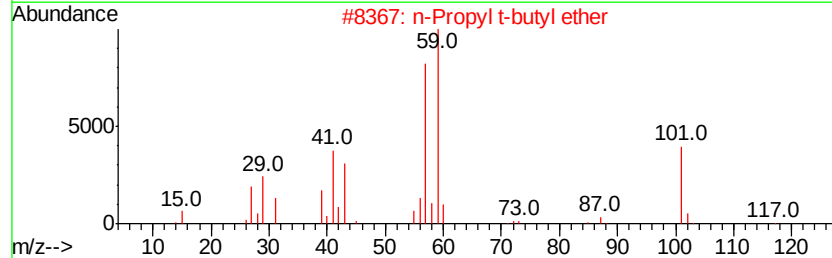
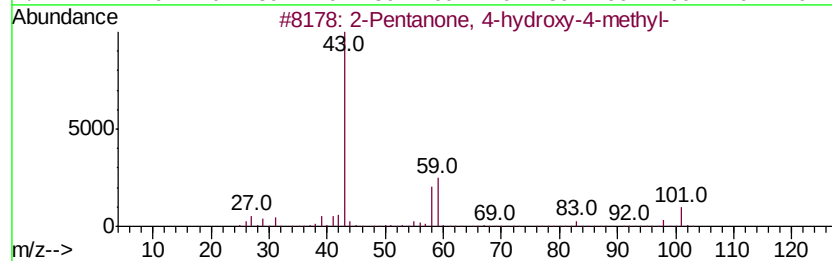
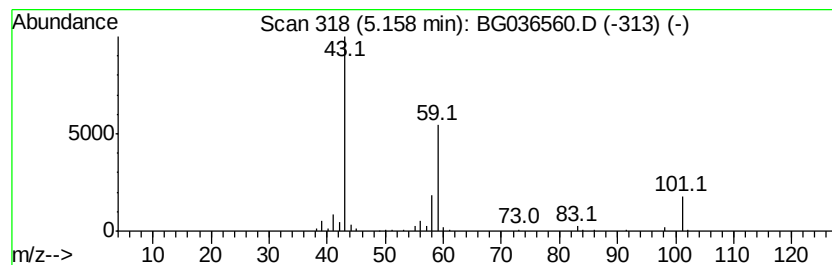
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.16	6.11 ng	113135	1,4-Dichlorobenzene-d4	8.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	6.1 ng		113135	1	8.06	370250	20.0