

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092518\
 Data File : BG036959.D
 Acq On : 25 Sep 2018 21:33
 Operator : JU/SJ
 Sample : J4775-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OR-01-092418-A

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-BG092018.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	5.150	312	317	326	rVB	70478	113697	3.76%	0.776%
2	5.614	390	396	415	rBV	591242	1053926	34.84%	7.192%
3	7.201	660	666	683	rBV	516049	1033283	34.15%	7.051%
4	7.577	723	730	745	rBV	587237	1135520	37.53%	7.749%
5	8.041	803	809	818	rBV	132363	259188	8.57%	1.769%
6	8.364	856	864	877	rBV	487249	909930	30.08%	6.209%
7	9.204	999	1007	1029	rBV	369241	784714	25.94%	5.355%
8	10.844	1278	1286	1294	rBV	170752	356765	11.79%	2.435%
9	13.288	1695	1702	1717	rBV	1018773	1733097	57.29%	11.827%
10	14.669	1930	1937	1947	rBV2	292837	504572	16.68%	3.443%
11	16.155	2183	2190	2209	rBV	863603	1438176	47.54%	9.814%
12	17.412	2397	2404	2418	rBV	402730	691134	22.84%	4.716%
13	20.021	2841	2848	2864	rBV	2162299	3025390	100.00%	20.646%
14	21.678	3123	3130	3143	rVB2	459274	812030	26.84%	5.541%
15	24.886	3666	3676	3691	rVB	263832	802507	26.53%	5.476%

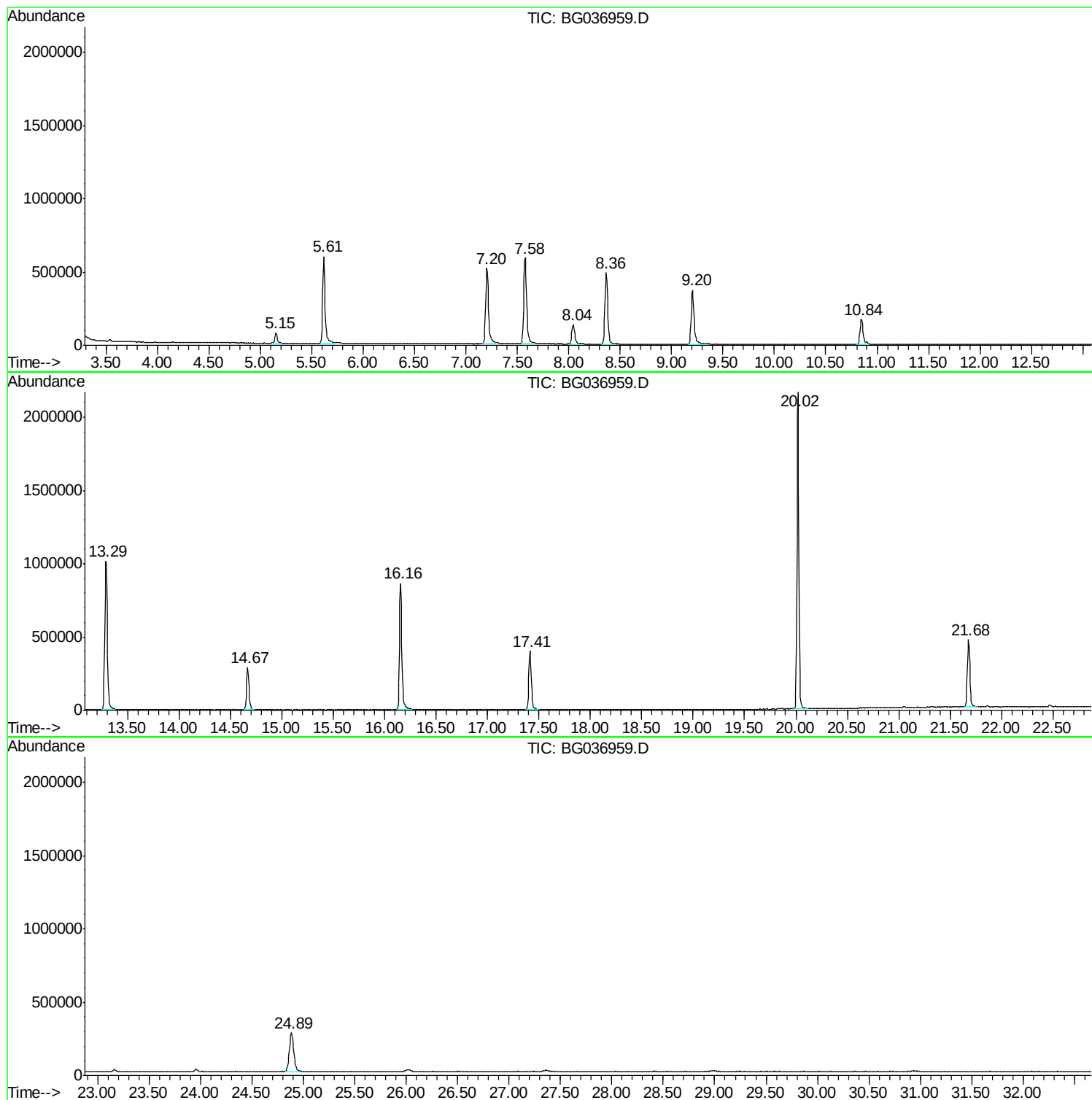
Sum of corrected areas: 14653929

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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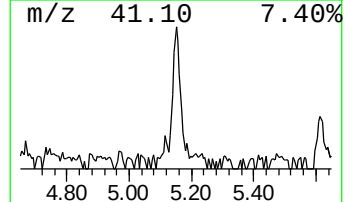
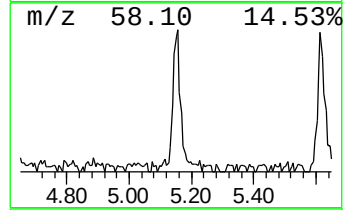
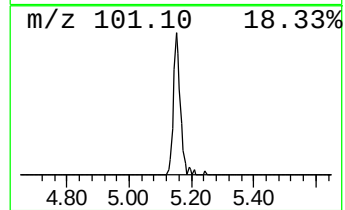
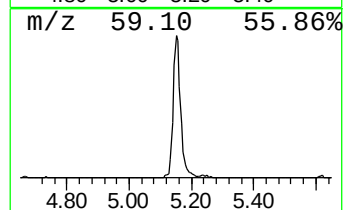
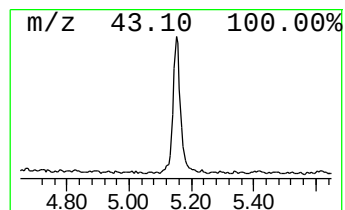
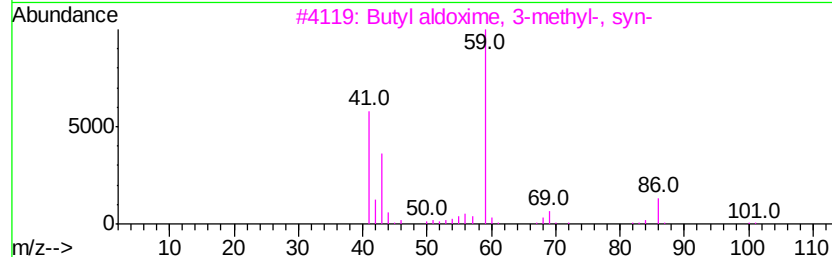
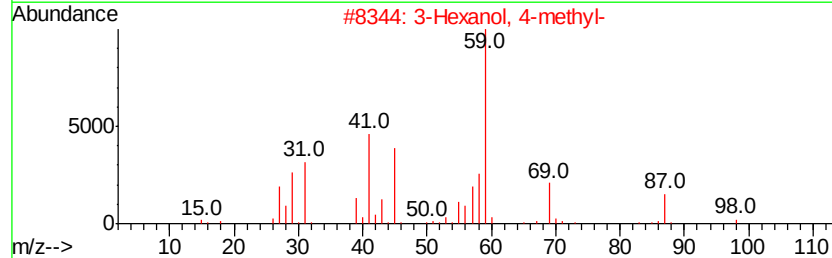
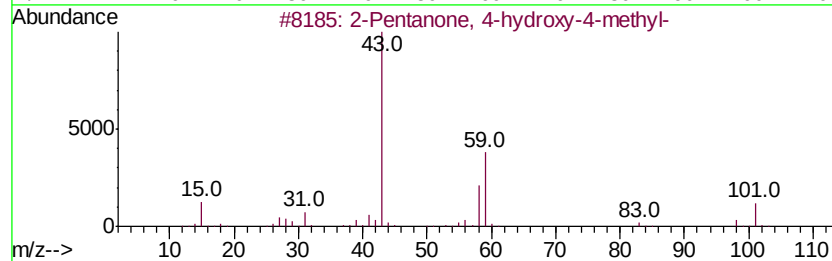
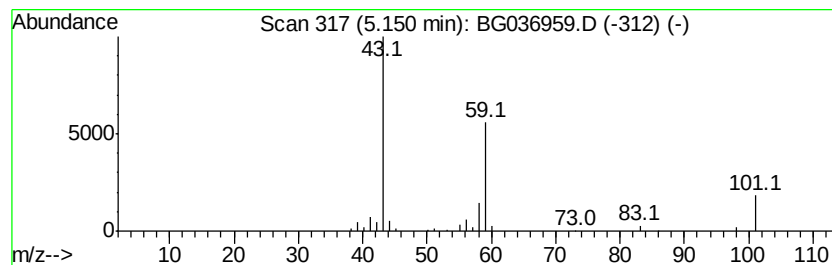
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	8.77 ng	113697	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Diacetamide	101	C4H7NO2	000625-77-4	9



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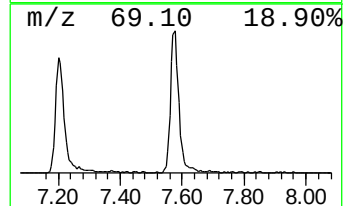
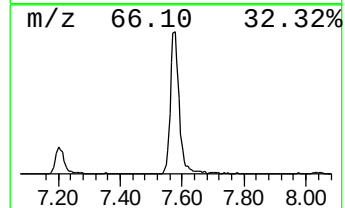
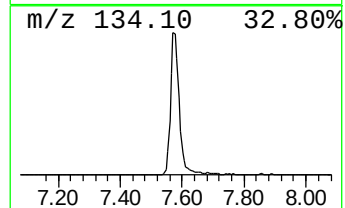
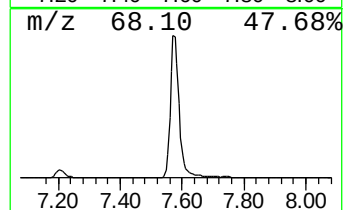
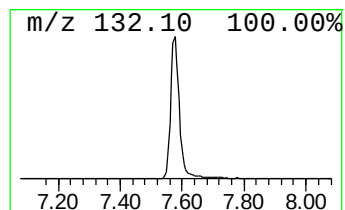
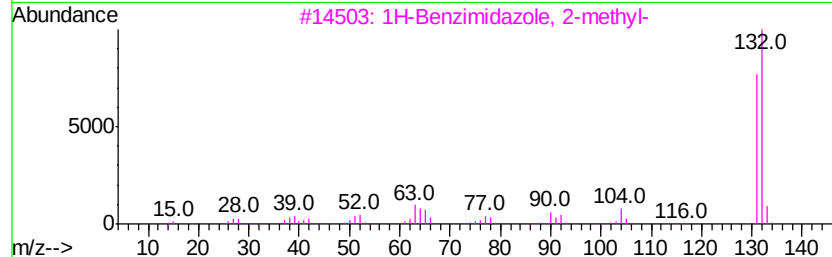
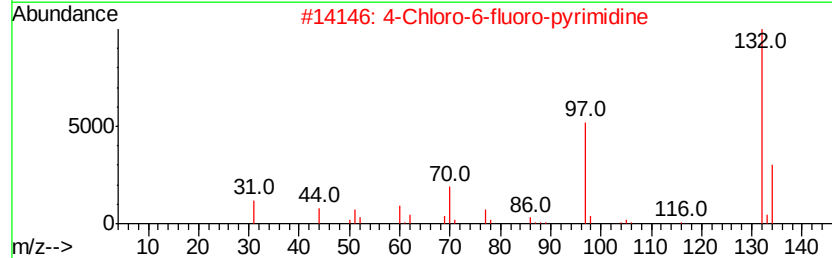
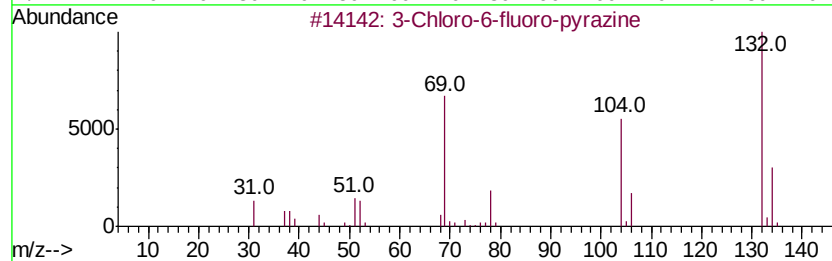
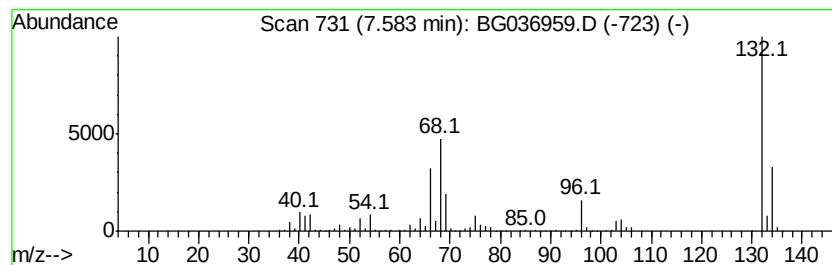
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	87.62 ng	1135520	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	8.8	ng	113697	1	8.04	259188	20.0
unknown7.58	7.58	87.6	ng	1135520	1	8.04	259188	20.0