

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100418\
 Data File : BG037134.D
 Acq On : 3 Oct 2018 15:27
 Operator : JU/SJ
 Sample : PB113454BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113454BL

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	5.151	312	317	326	rVB	31813	57165	1.99%	0.358%
2	5.615	390	396	417	rBV	696620	1292645	44.99%	8.106%
3	7.202	659	666	688	rBV	683645	1393907	48.52%	8.741%
4	7.572	722	729	751	rBV	692239	1355159	47.17%	8.498%
5	8.042	802	809	822	rBV	156528	331442	11.54%	2.078%
6	8.359	856	863	877	rBV	547347	1056610	36.78%	6.626%
7	9.199	999	1006	1026	rBV	397480	878579	30.58%	5.509%
8	10.844	1279	1286	1298	rBV	208230	443936	15.45%	2.784%
9	13.283	1694	1701	1721	rBV	1137523	1995810	69.47%	12.515%
10	14.663	1929	1936	1949	rBV2	356923	624016	21.72%	3.913%
11	16.150	2183	2189	2214	rBV	652027	1242994	43.27%	7.794%
12	17.407	2397	2403	2424	rBV	407725	753106	26.21%	4.722%
13	20.016	2841	2847	2866	rBV	1995742	2872919	100.00%	18.015%
14	21.673	3122	3129	3147	rBV2	429136	858031	29.87%	5.380%
15	24.881	3664	3675	3693	rBV2	240810	791247	27.54%	4.962%

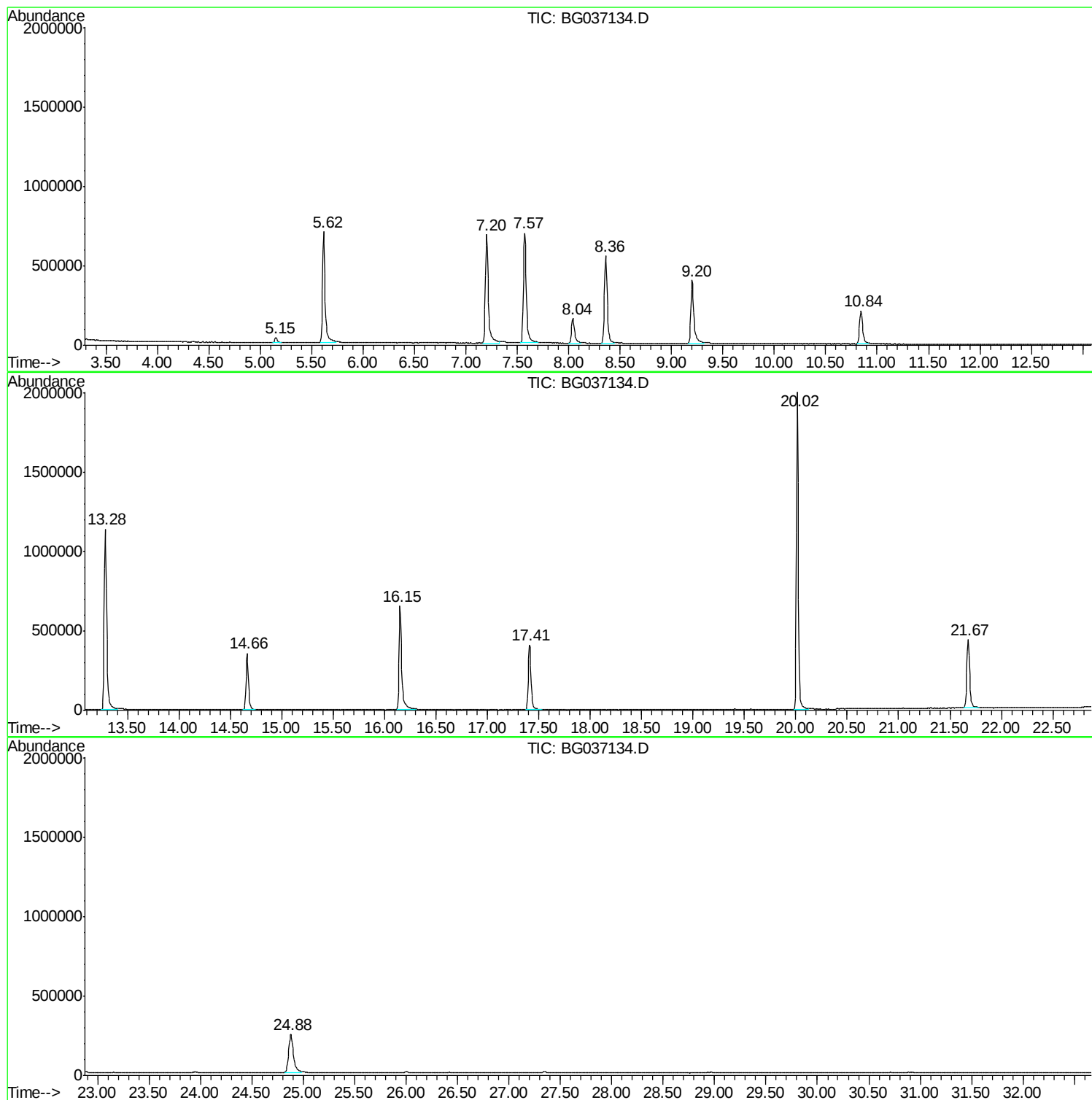
Sum of corrected areas: 15947566

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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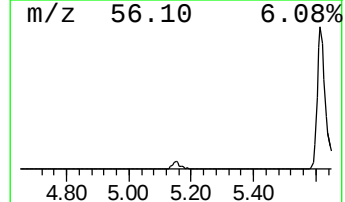
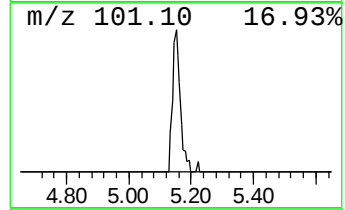
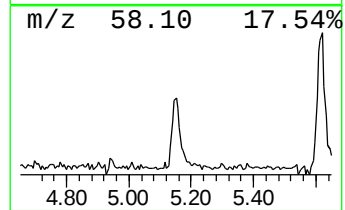
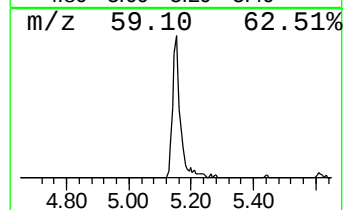
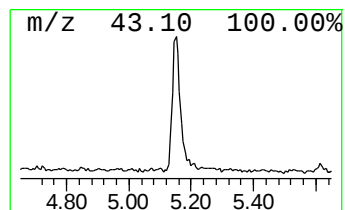
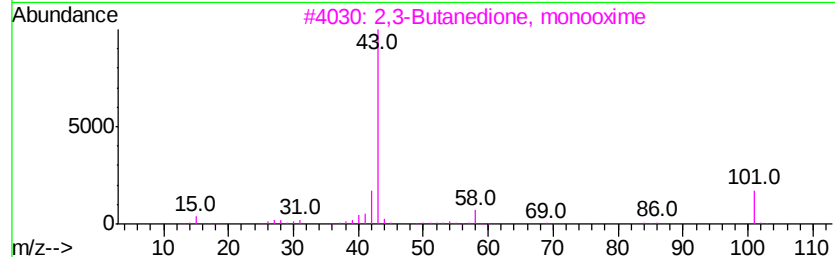
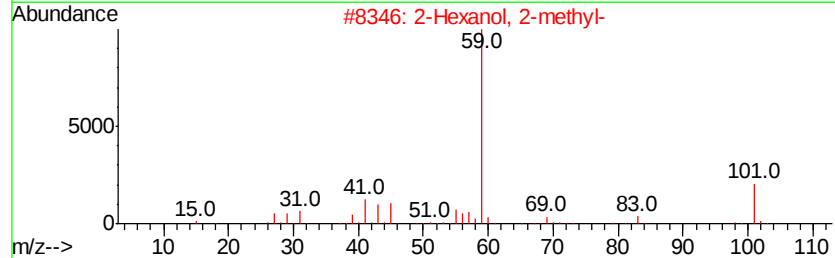
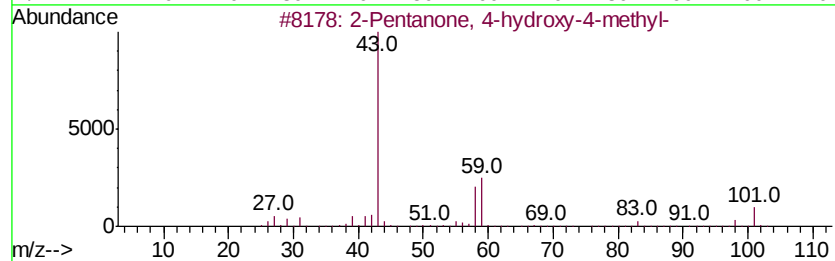
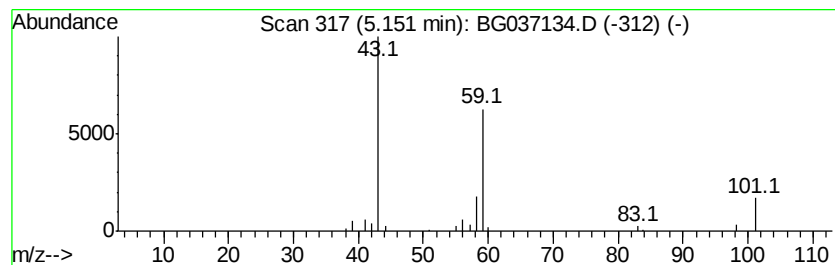
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.45 ng	57165	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	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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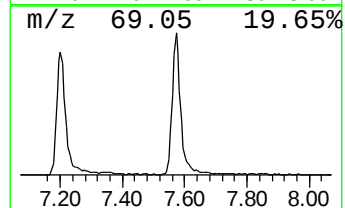
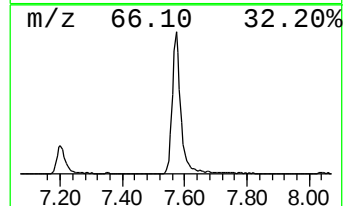
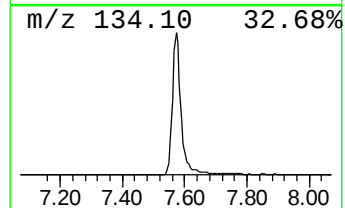
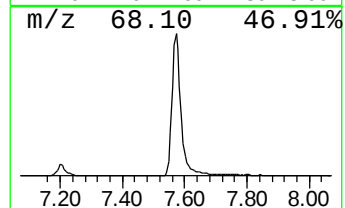
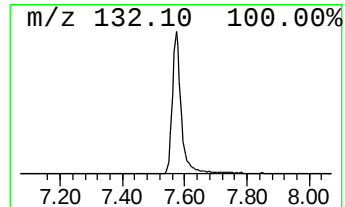
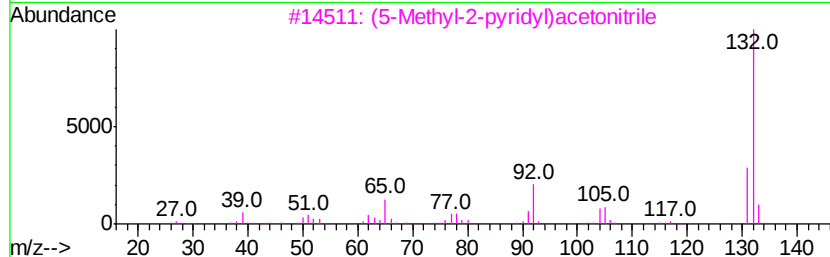
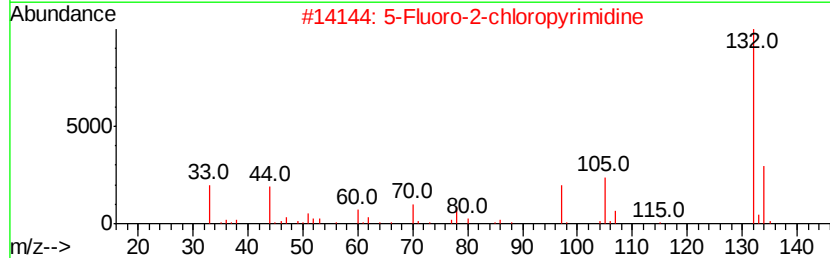
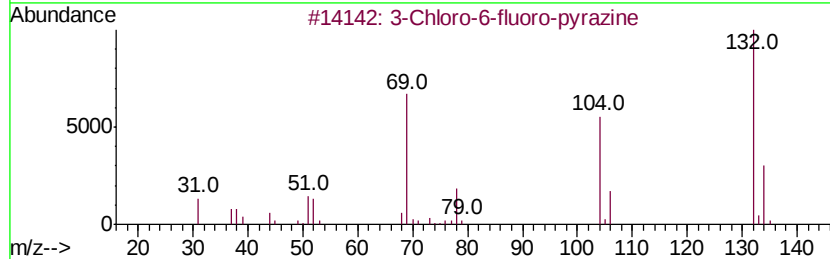
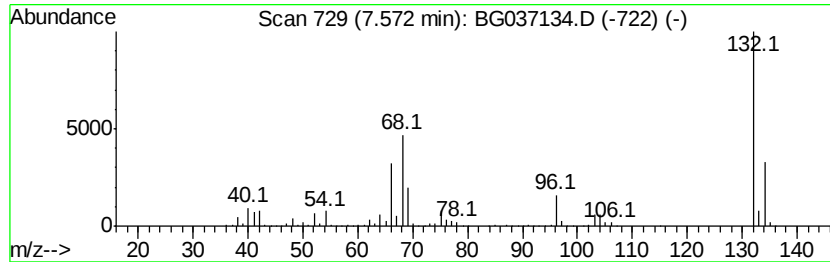
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	81.77 ng	1355160	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	3.5	ng	57165	1	8.04	331442	20.0
unknown7.57	7.57	81.8	ng	1355160	1	8.04	331442	20.0