

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036631.D
 Acq On : 10 Sep 2018 19:02
 Operator : JU/SJ
 Sample : J4769-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-05-090718-B

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	5.163	309	319	326	rVB	130891	198484	4.48%	0.938%
2	5.627	390	398	410	rBV	939539	1511397	34.14%	7.139%
3	7.213	659	668	678	rBV	900660	1558839	35.21%	7.363%
4	7.583	722	731	740	rBV	974287	1675157	37.84%	7.912%
5	8.053	805	811	819	rBV	206555	340049	7.68%	1.606%
6	8.377	858	866	875	rVB	758340	1294086	29.23%	6.113%
7	9.217	1001	1009	1024	rBV	636214	1144128	25.85%	5.404%
8	10.862	1279	1289	1295	rBV	282589	511439	11.55%	2.416%
9	13.300	1697	1704	1716	rBV	1745592	2764506	62.45%	13.058%
10	14.681	1932	1939	1946	rBV2	451774	744548	16.82%	3.517%
11	16.162	2185	2191	2208	rBV	1251454	1993446	45.03%	9.416%
12	17.419	2396	2405	2417	rBV	637501	986383	22.28%	4.659%
13	20.028	2843	2849	2862	rBV	3192737	4426876	100.00%	20.910%
14	21.685	3125	3131	3145	rBV2	646956	1098928	24.82%	5.191%
15	24.893	3667	3677	3701	rBV	258315	922848	20.85%	4.359%

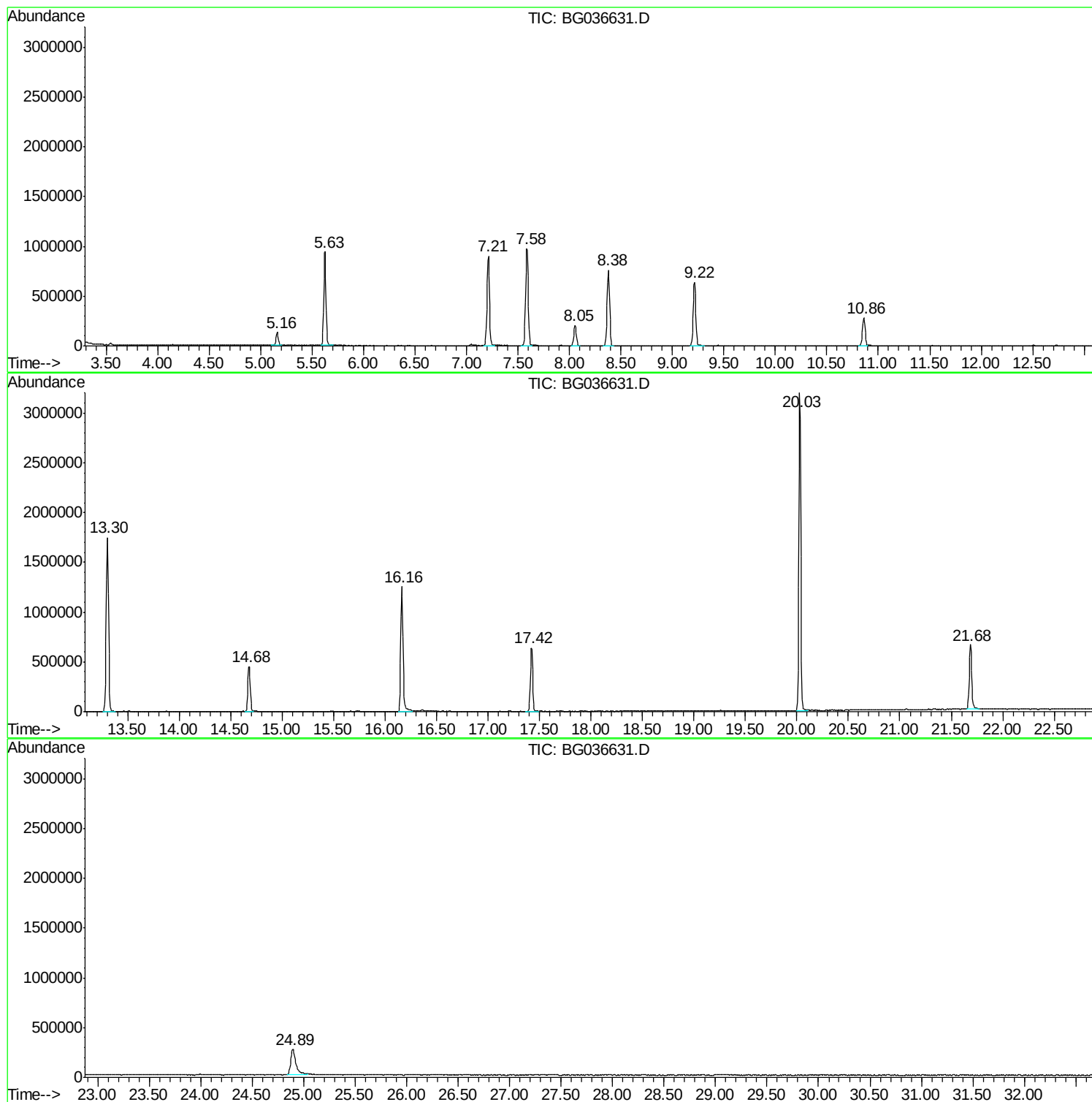
Sum of corrected areas: 21171114

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



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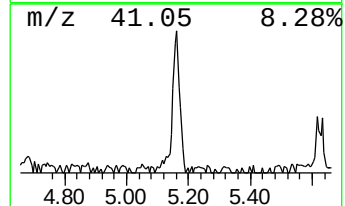
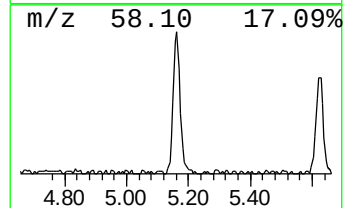
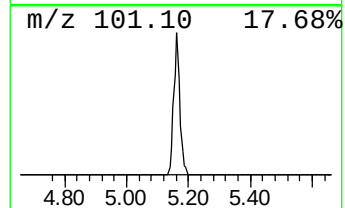
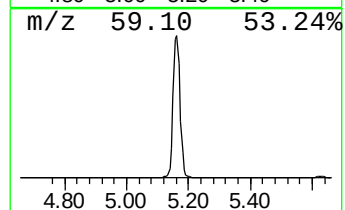
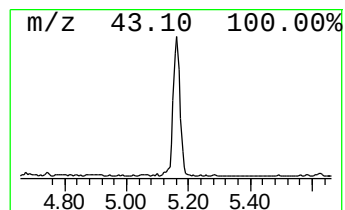
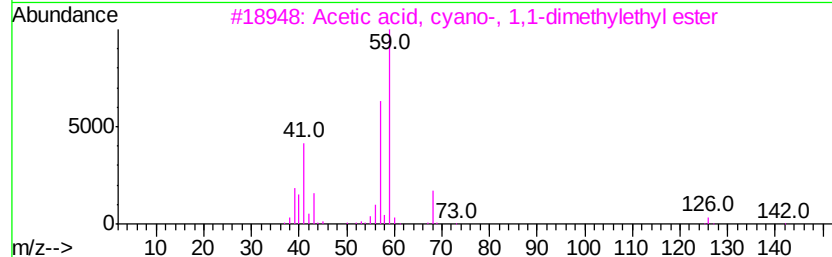
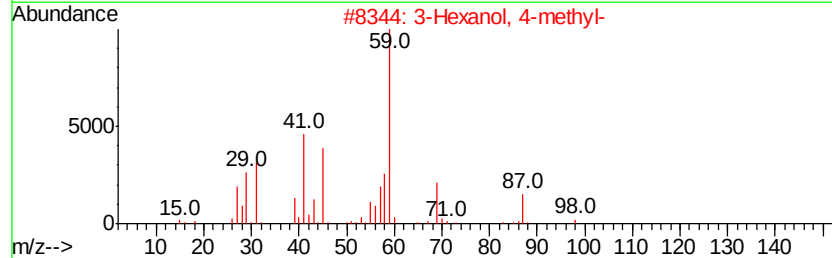
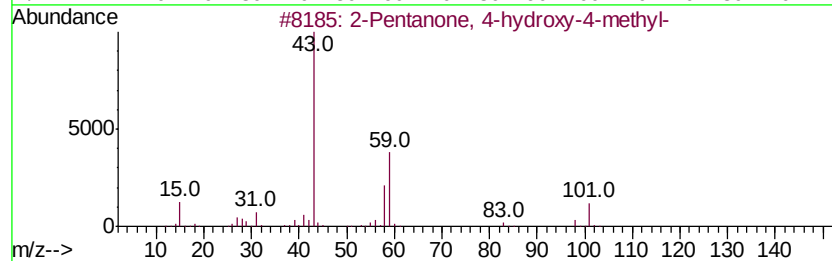
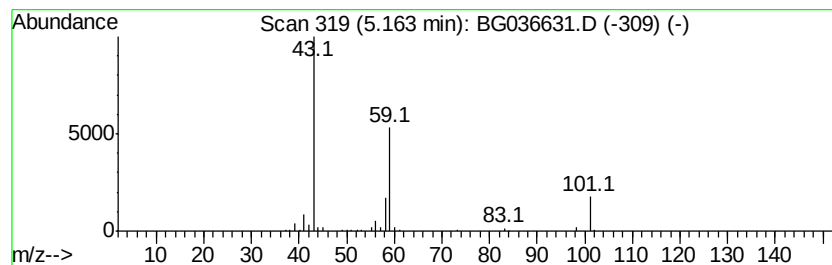
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	11.67 ng	198484	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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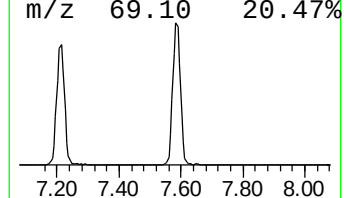
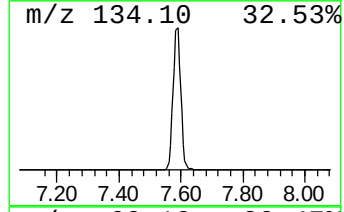
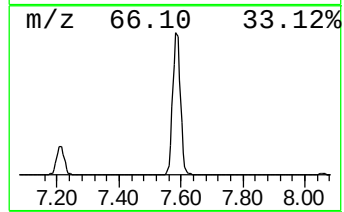
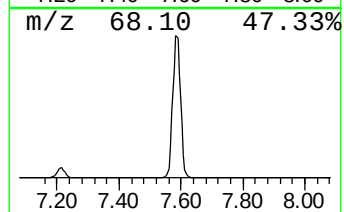
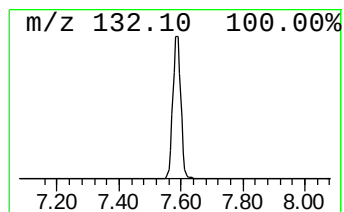
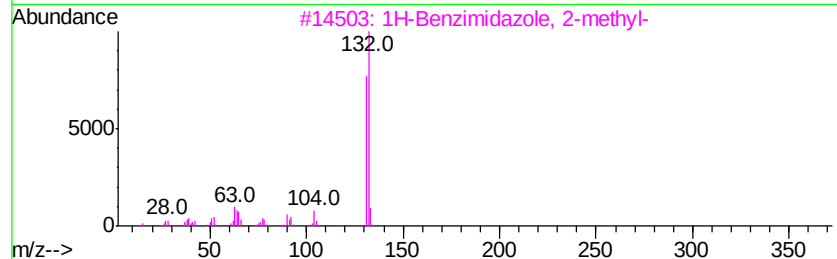
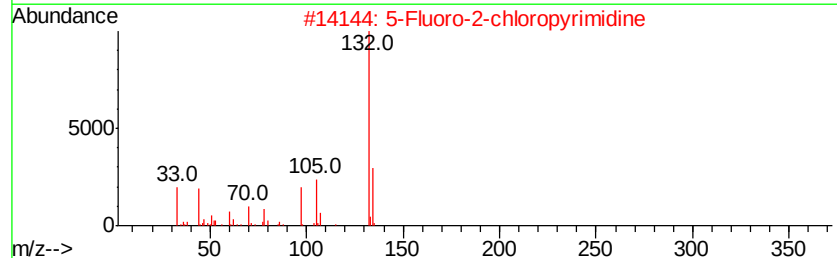
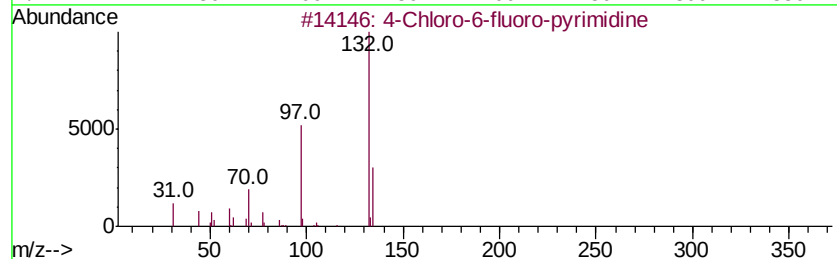
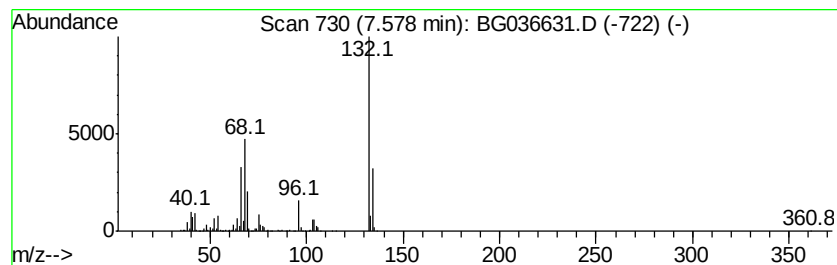
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	98.52 ng	1675160	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	11.7	ng	198484	1	8.05	340049	20.0
unknown7.58	7.58	98.5	ng	1675160	1	8.05	340049	20.0