

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091118\
 Data File : BG036662.D
 Acq On : 11 Sep 2018 18:50
 Operator : JU/SJ
 Sample : J4841-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090718-DBRI-F-U-M

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	5.156	313	318	329	rBV	47546	73705	1.57%	0.392%
2	5.620	391	397	408	rVB	373478	593095	12.65%	3.158%
3	7.206	660	667	678	rBV	248935	429153	9.15%	2.285%
4	7.582	723	731	739	rBV	644379	1089298	23.23%	5.800%
5	8.052	805	811	818	rBV	187308	315621	6.73%	1.681%
6	8.376	858	866	874	rBV	853470	1457479	31.08%	7.760%
7	9.216	1001	1009	1022	rBV	726255	1295660	27.63%	6.899%
8	10.861	1281	1289	1297	rBV	257006	465838	9.93%	2.480%
9	13.299	1697	1704	1717	rBV	1934494	3132585	66.79%	16.680%
10	14.122	1839	1844	1850	rVB	81744	117217	2.50%	0.624%
11	14.680	1932	1939	1948	rVB2	437052	714741	15.24%	3.806%
12	16.161	2185	2191	2206	rBV	994106	1563522	33.34%	8.325%
13	17.418	2399	2405	2415	rBV2	598462	917339	19.56%	4.884%
14	20.027	2843	2849	2861	rBV	3524350	4689924	100.00%	24.972%
15	21.684	3124	3131	3149	rBV	596959	1008661	21.51%	5.371%
16	24.892	3667	3677	3692	rBV2	257732	869642	18.54%	4.630%
17	26.049	3867	3874	3884	rBV7	16329	47354	1.01%	0.252%

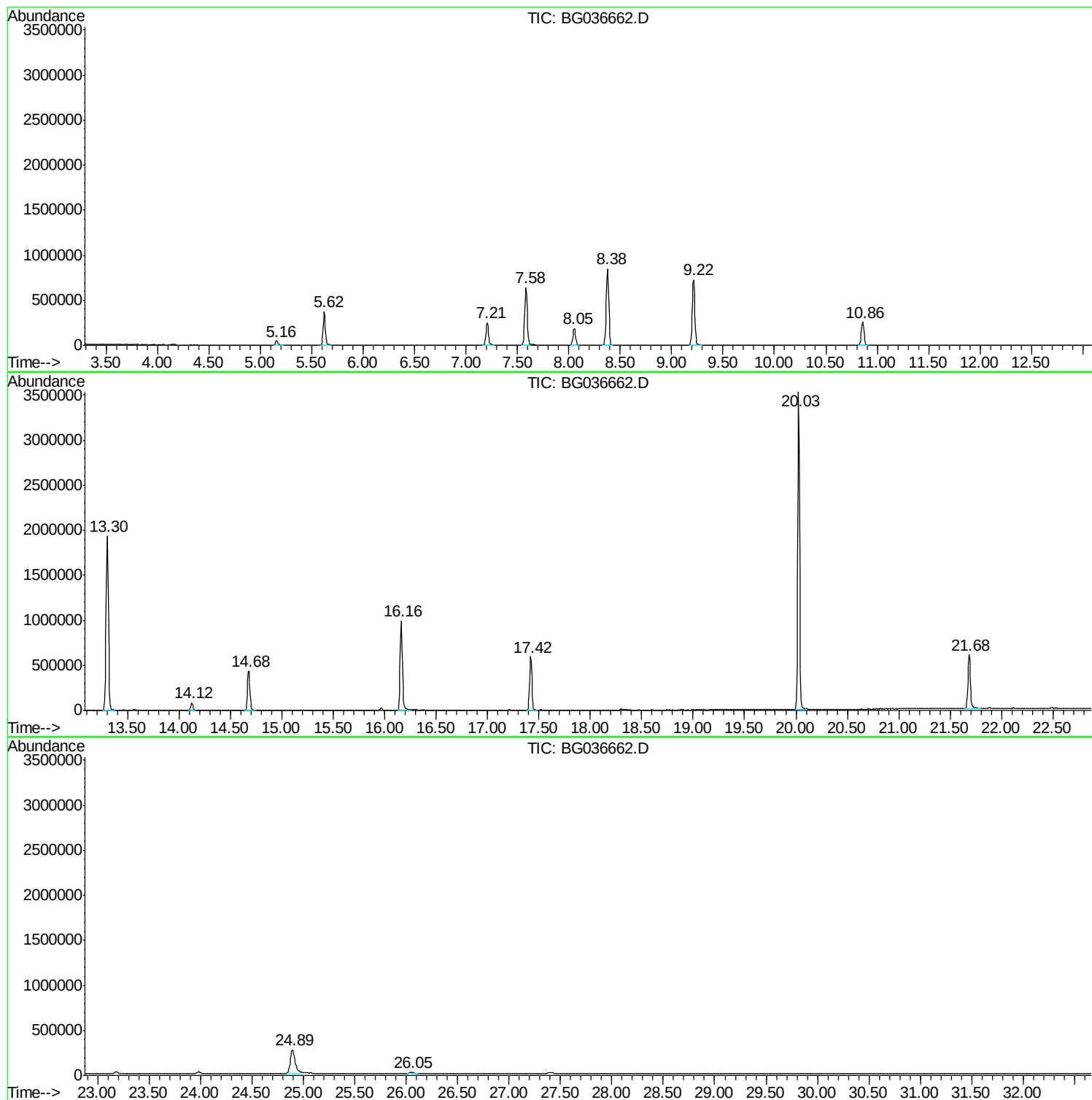
Sum of corrected areas: 18780834

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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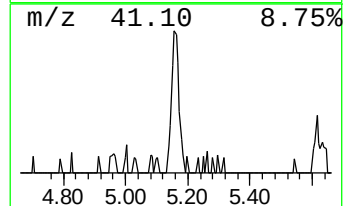
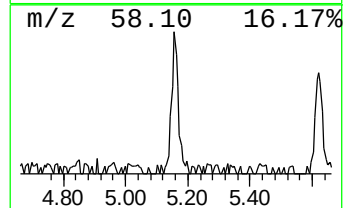
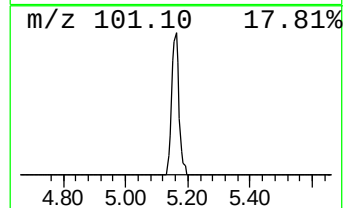
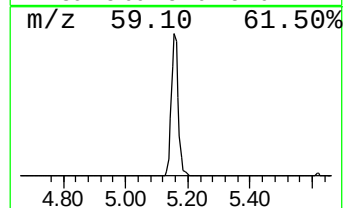
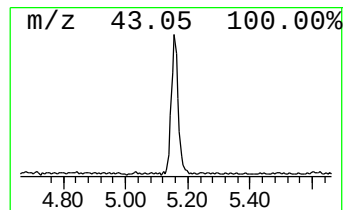
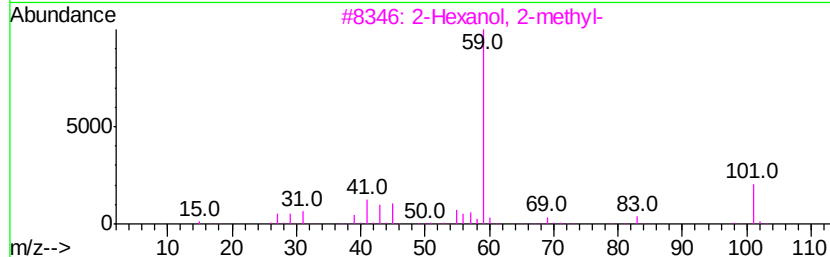
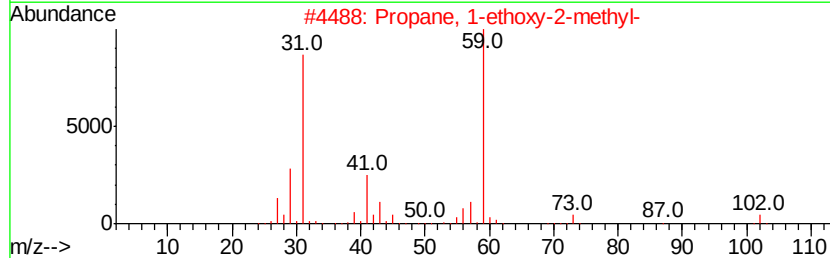
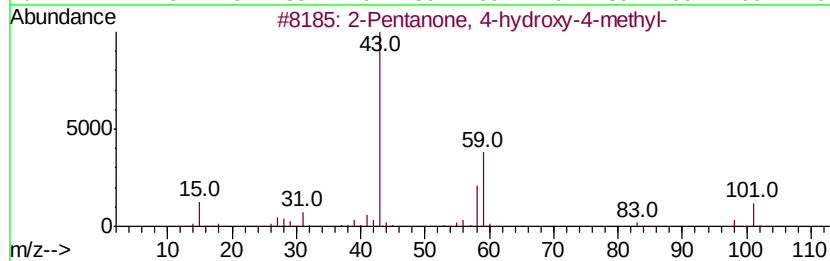
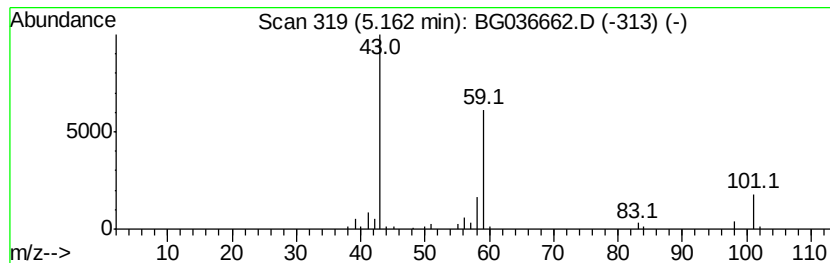
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.67 ng	73705	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	53
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17



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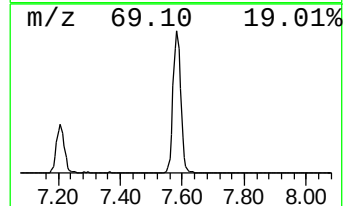
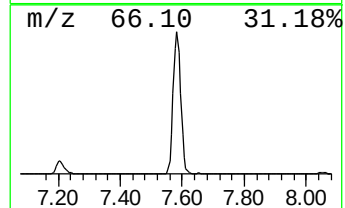
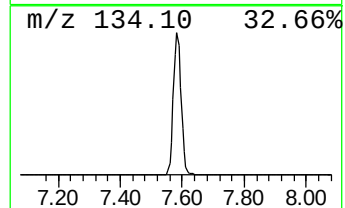
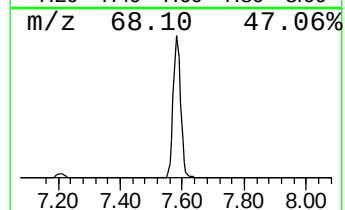
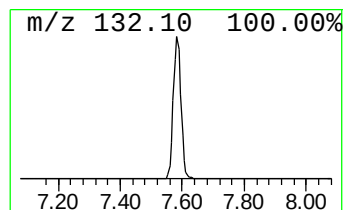
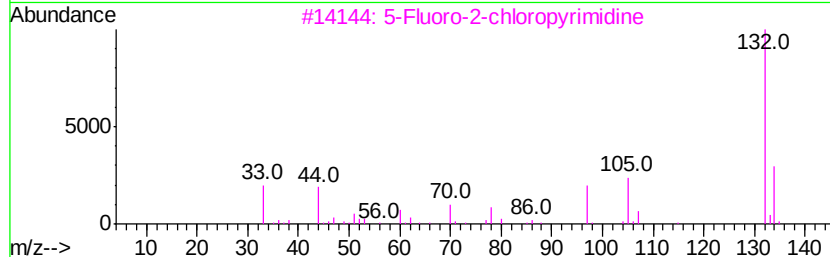
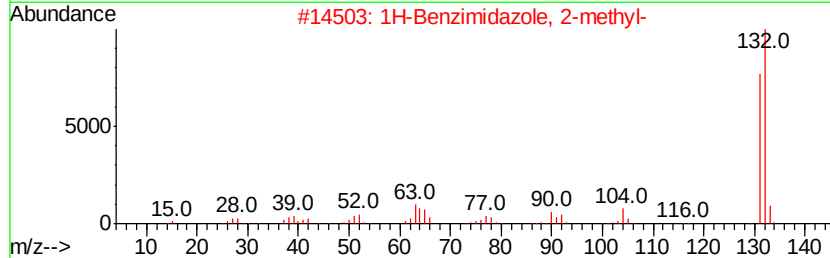
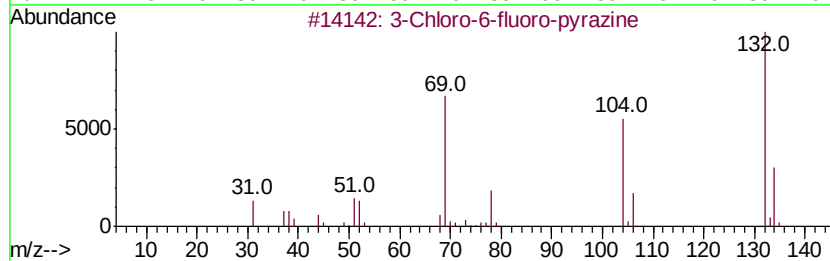
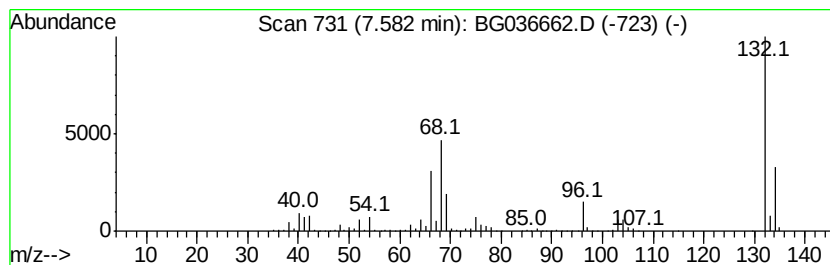
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	69.03 ng	1089300	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	4.7	ng	73705	1	8.05	315621	20.0
unknown7.58	7.58	69.0	ng	1089300	1	8.05	315621	20.0