

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036637.D
 Acq On : 10 Sep 2018 22:55
 Operator : JU/SJ
 Sample : J4777-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PL-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	5.161	309	319	327	rBV	100039	158492	3.83%	0.831%
2	5.620	391	397	413	rBV	839561	1368138	33.03%	7.171%
3	7.212	660	668	681	rBV	796246	1387038	33.49%	7.270%
4	7.582	724	731	744	rBV	864651	1516159	36.60%	7.947%
5	8.052	803	811	819	rVB	193950	323451	7.81%	1.695%
6	8.375	859	866	874	rVB	660164	1136420	27.44%	5.956%
7	9.216	1001	1009	1017	rBV	585979	1030482	24.88%	5.401%
8	10.861	1281	1289	1295	rBV	266502	480314	11.60%	2.517%
9	13.305	1697	1705	1716	rBV	1549143	2478500	59.84%	12.991%
10	14.680	1932	1939	1948	rBV2	435363	704840	17.02%	3.694%
11	16.166	2185	2192	2203	rBV	935746	1587522	38.33%	8.321%
12	17.424	2399	2406	2420	rBV2	599808	939757	22.69%	4.926%
13	20.026	2843	2849	2859	rBV	3019673	4142106	100.00%	21.710%
14	21.689	3124	3132	3145	rBV	554612	1052992	25.42%	5.519%
15	24.897	3667	3678	3695	rBV2	198644	772960	18.66%	4.051%

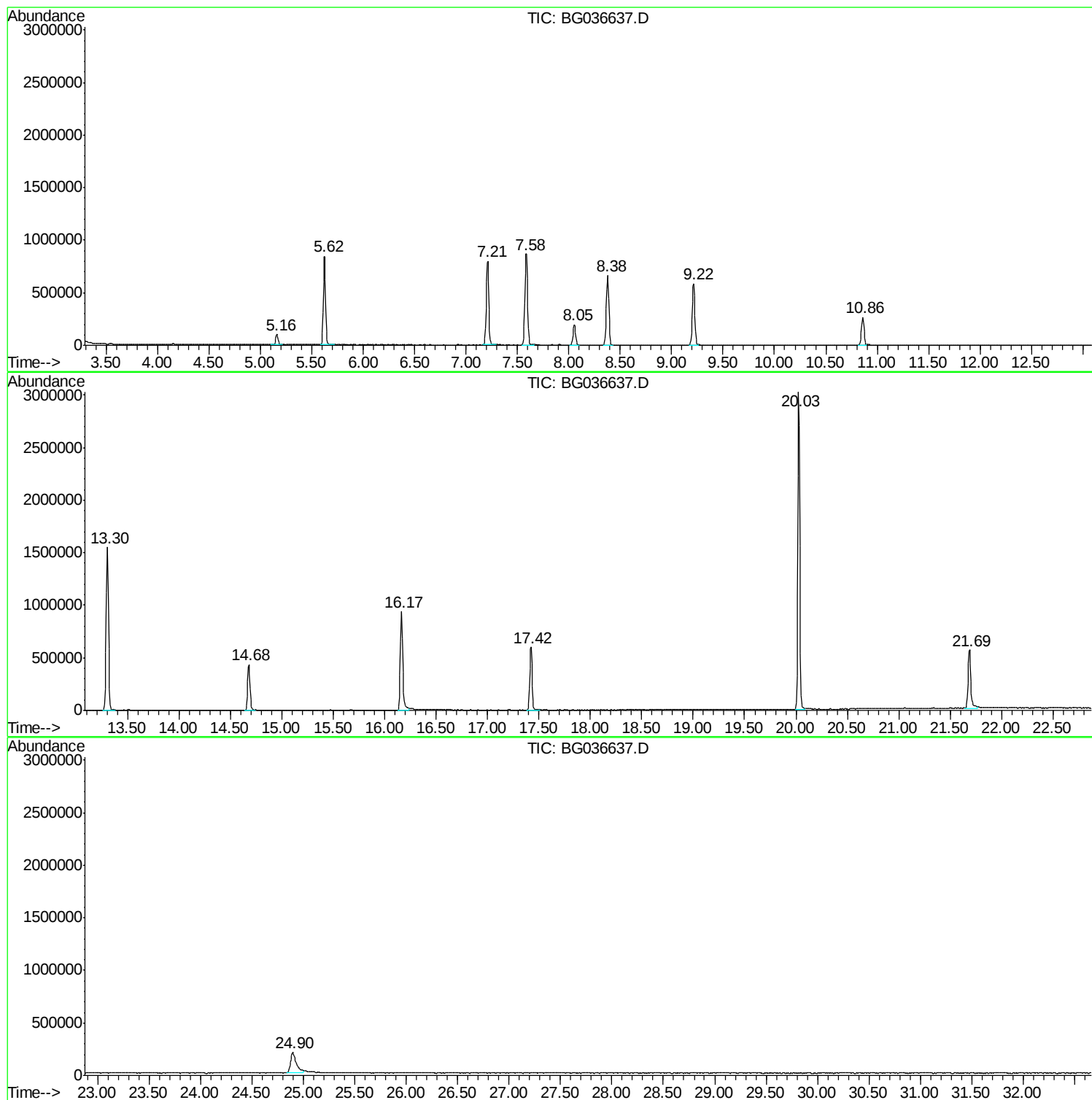
Sum of corrected areas: 19079171

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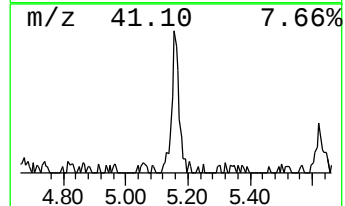
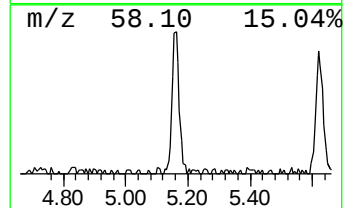
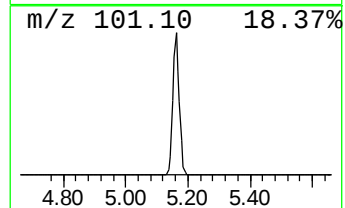
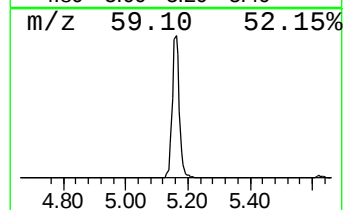
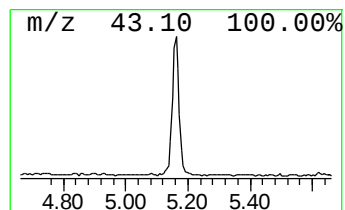
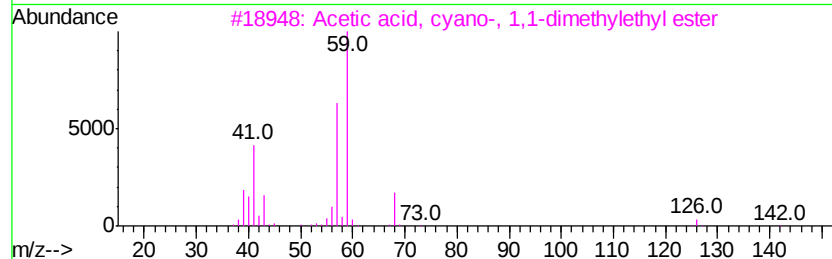
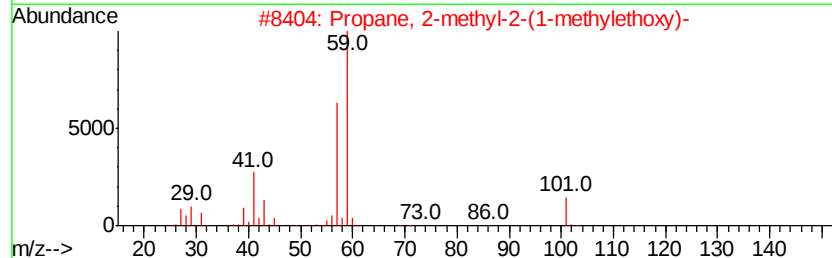
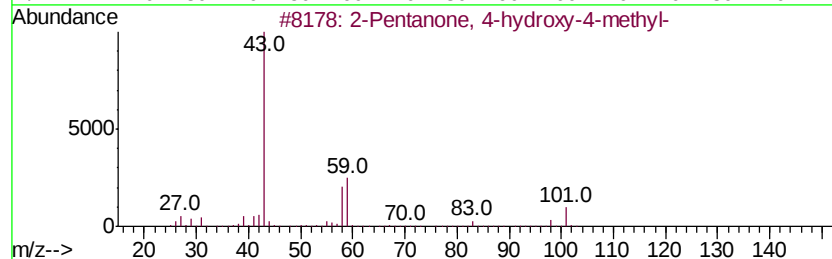
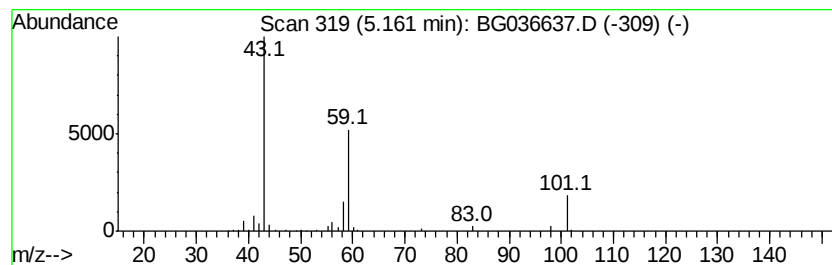
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	9.80 ng	158492	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	64
2			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	28
3			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5			Silane, dimethyl-	60	C2H8Si	001111-74-6	9



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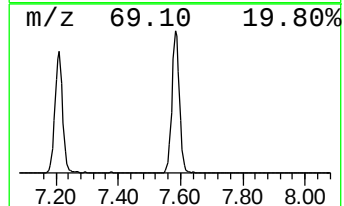
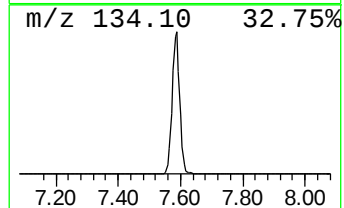
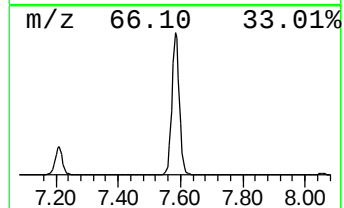
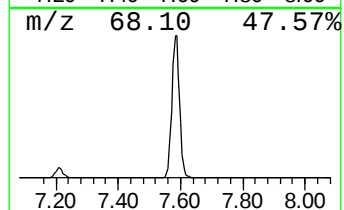
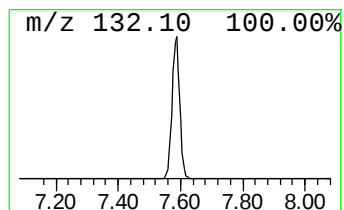
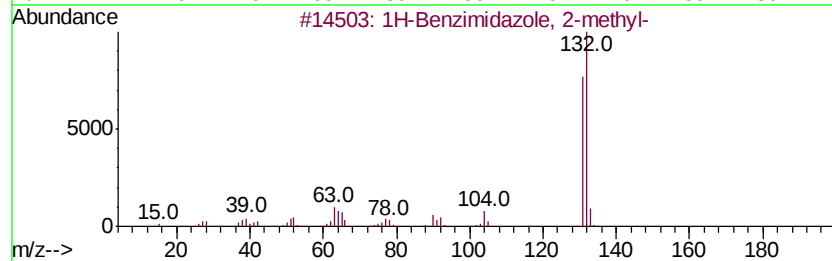
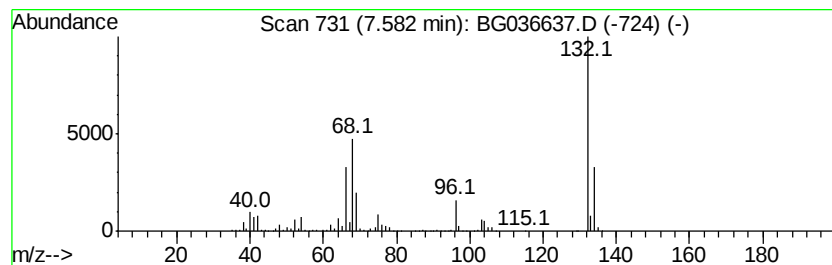
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	93.75 ng	1516160	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	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...	5.16	9.8	ng	158492	1	8.05	323451	20.0
unknown7.58	7.58	93.8	ng	1516160	1	8.05	323451	20.0