

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
Data File : BG036630.D
Acq On : 10 Sep 2018 18:23
Operator : JU/SJ
Sample : J4769-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
HR-05-090718-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-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	5.162	309	319	329	rVB	116871	180831	4.50%	0.961%
2	5.626	391	398	408	rBV	813079	1299980	32.38%	6.907%
3	7.207	659	667	679	rBV	778840	1338290	33.34%	7.111%
4	7.583	724	731	742	rBV	833152	1460283	36.38%	7.759%
5	8.053	803	811	818	rVB	192907	330890	8.24%	1.758%
6	8.376	858	866	874	rBV	641595	1114788	27.77%	5.923%
7	9.216	1000	1009	1018	rBV	558951	989861	24.66%	5.260%
8	10.861	1280	1289	1295	rBV	270519	496749	12.37%	2.639%
9	13.305	1697	1705	1713	rBV	1507474	2420984	60.31%	12.864%
10	14.680	1931	1939	1946	rBV2	443673	736466	18.35%	3.913%
11	16.161	2185	2191	2213	rBV	921130	1574262	39.22%	8.365%
12	17.418	2398	2405	2419	rBV2	628922	986857	24.58%	5.244%
13	20.027	2843	2849	2863	rBV	2983002	4014321	100.00%	21.330%
14	21.690	3125	3132	3141	rBV	582217	1058913	26.38%	5.626%
15	24.898	3669	3678	3700	rBV	224099	816691	20.34%	4.339%

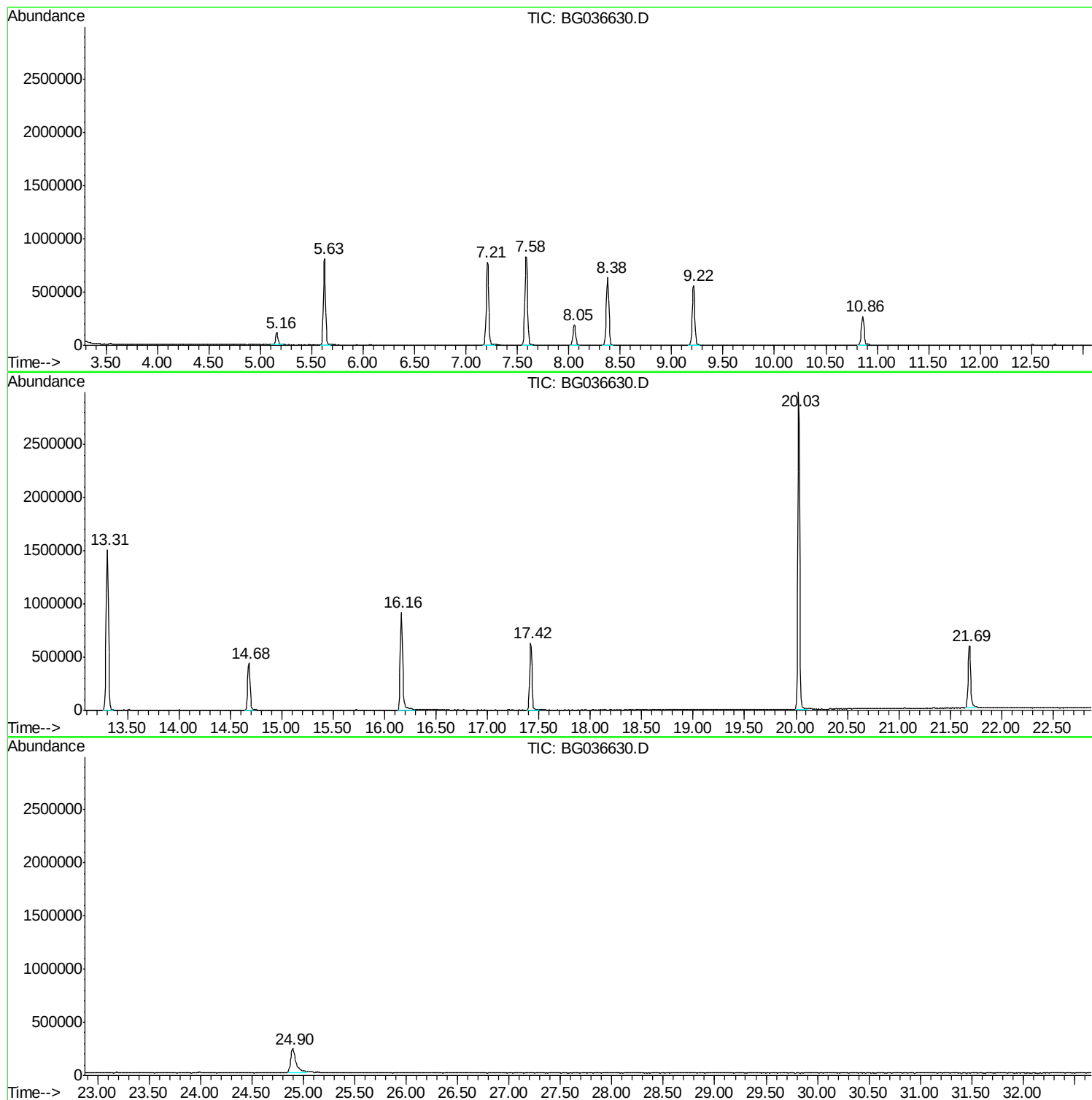
Sum of corrected areas: 18820166

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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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ClientSampled :
HR-05-090718-A

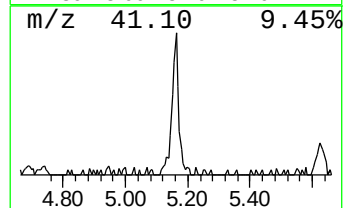
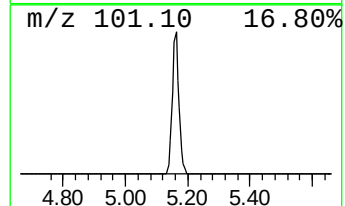
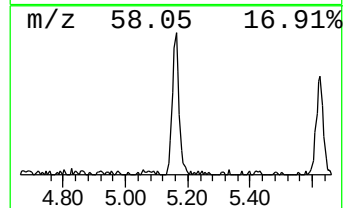
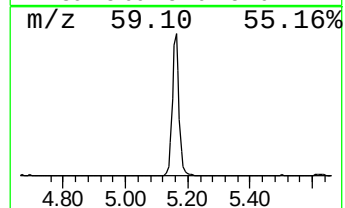
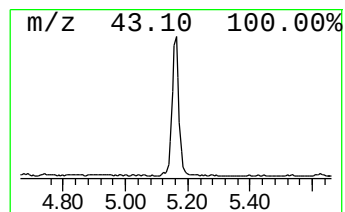
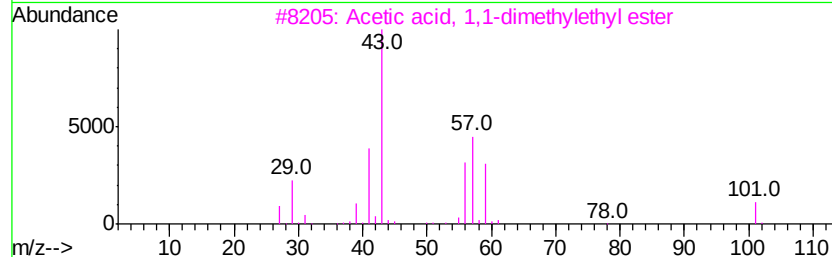
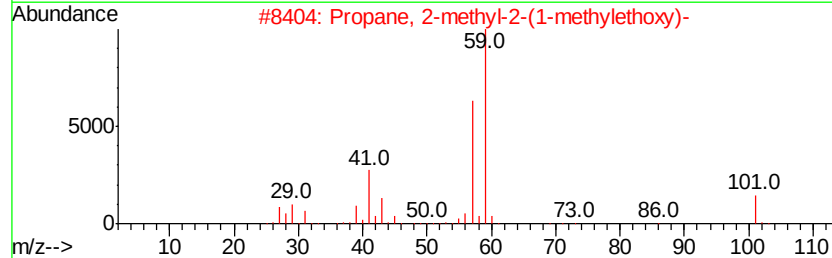
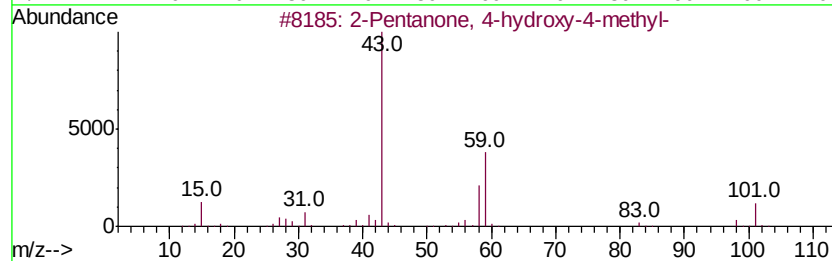
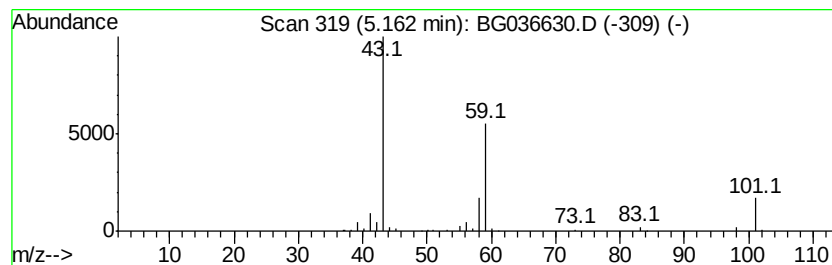
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	10.93 ng	180831	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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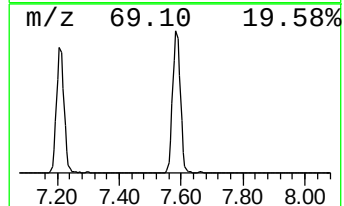
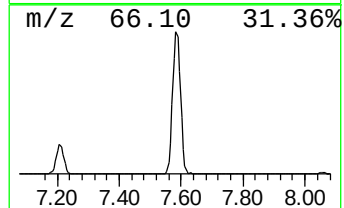
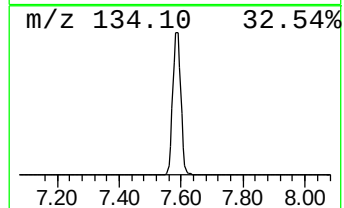
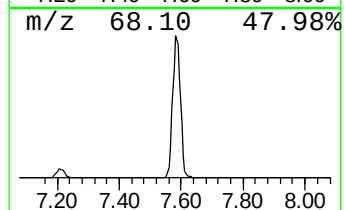
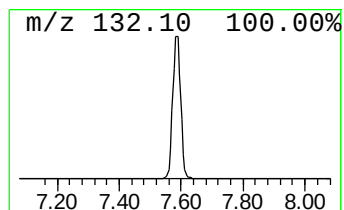
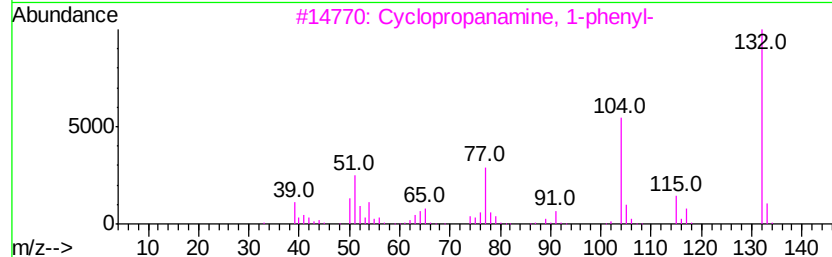
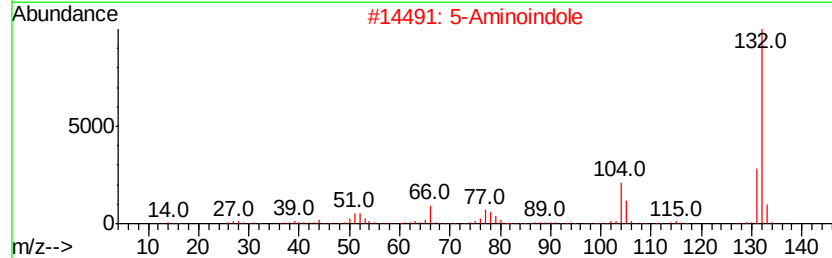
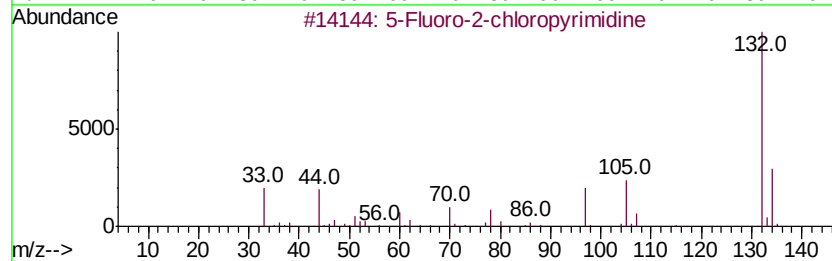
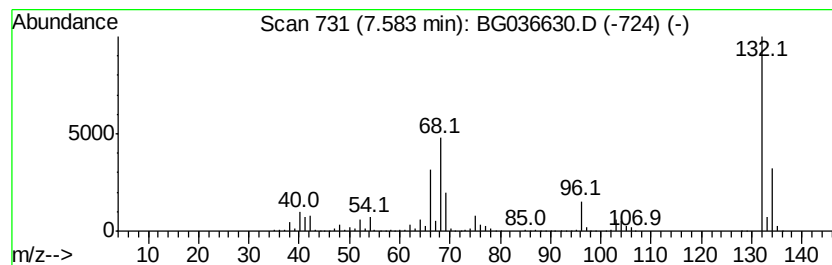
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	88.26 ng	1460280	1,4-Dichlorobenzene-d4	8.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
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	10.9	ng	180831	1	8.05	330890	20.0
unknown7.58	7.58	88.3	ng	1460280	1	8.05	330890	20.0