

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036892.D
 Acq On : 22 Sep 2018 13:35
 Operator : JU/SJ
 Sample : J5019-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 091918-DBRI-E-D-M

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	5.149	312	317	324	rBV	27900	43389	1.09%	0.296%
2	5.620	390	397	411	rBV	165442	329343	8.25%	2.243%
3	7.206	660	667	682	rBV	101140	250443	6.27%	1.706%
4	7.576	723	730	744	rBV	349072	683785	17.12%	4.657%
5	8.046	802	810	828	rBV	120953	239251	5.99%	1.629%
6	8.369	857	865	884	rBV	528551	1001865	25.09%	6.823%
7	9.209	1000	1008	1026	rBV	442498	924271	23.15%	6.295%
8	10.849	1280	1287	1304	rBV	169643	356730	8.93%	2.430%
9	13.293	1696	1703	1718	rBV	1388555	2340936	58.62%	15.943%
10	14.121	1838	1844	1852	rBV	43520	74813	1.87%	0.510%
11	14.674	1931	1938	1949	rBV2	320891	561325	14.06%	3.823%
12	16.160	2183	2191	2204	rBV	721576	1234226	30.91%	8.406%
13	17.417	2398	2405	2416	rBV	471351	754545	18.90%	5.139%
14	20.026	2842	2849	2863	rBV	2829485	3993170	100.00%	27.196%
15	21.683	3125	3131	3143	rBV	528216	888646	22.25%	6.052%
16	23.352	3411	3415	3421	rVB	29426	55249	1.38%	0.376%
17	23.963	3514	3519	3527	rVB5	23808	48293	1.21%	0.329%
18	24.891	3666	3677	3690	rBV2	305481	902526	22.60%	6.147%

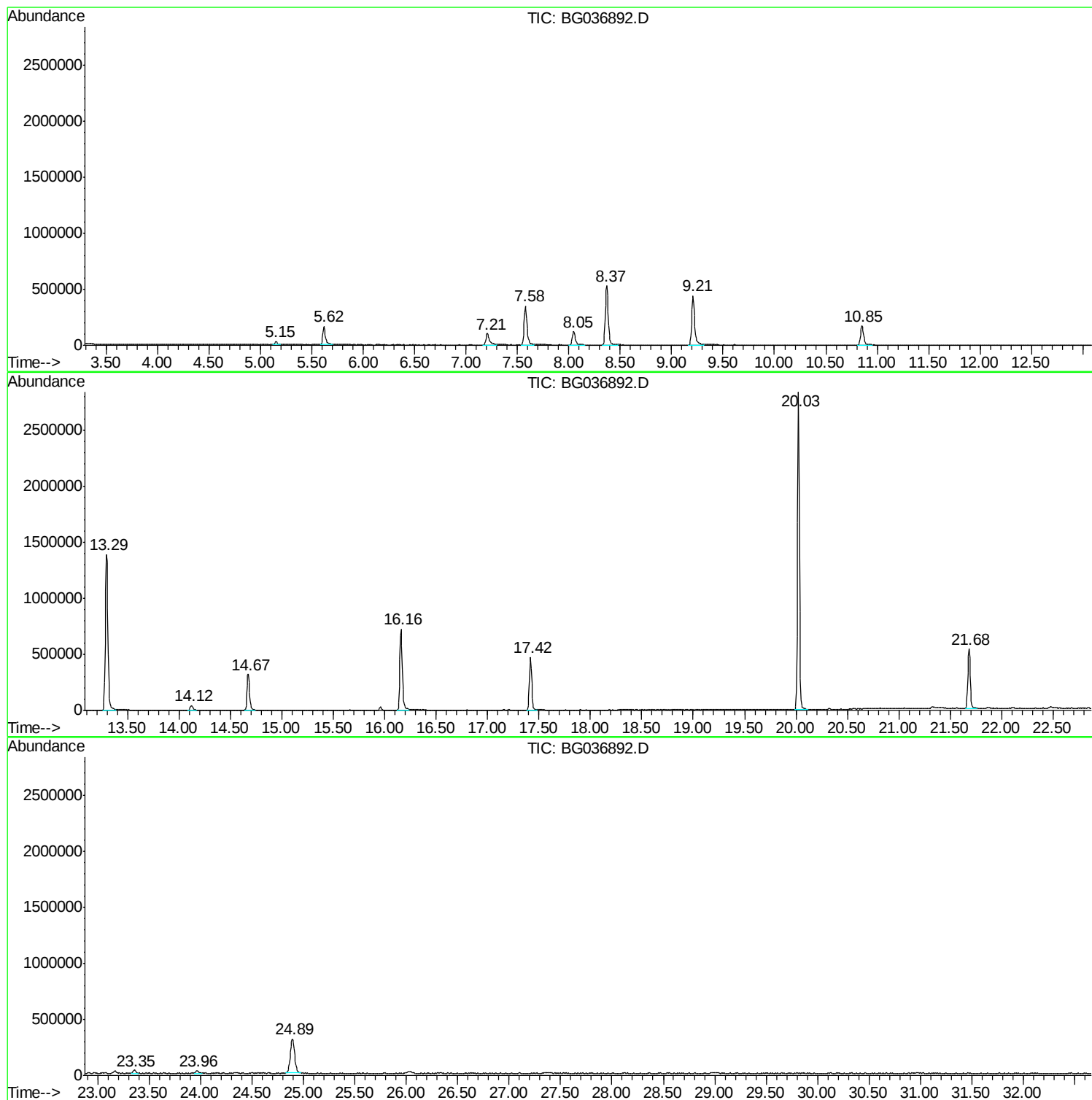
Sum of corrected areas: 14682806

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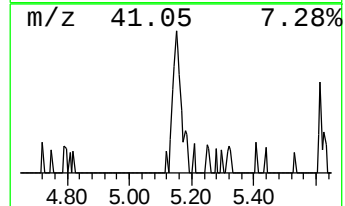
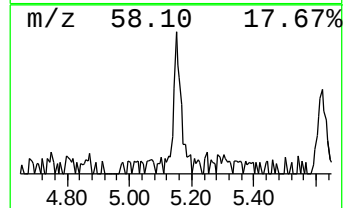
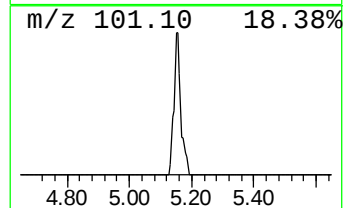
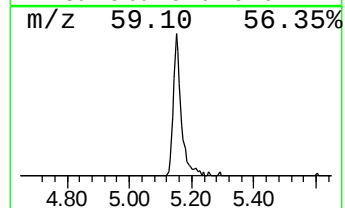
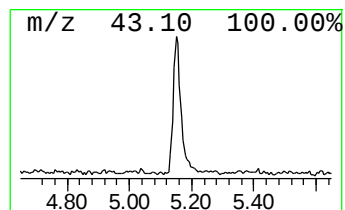
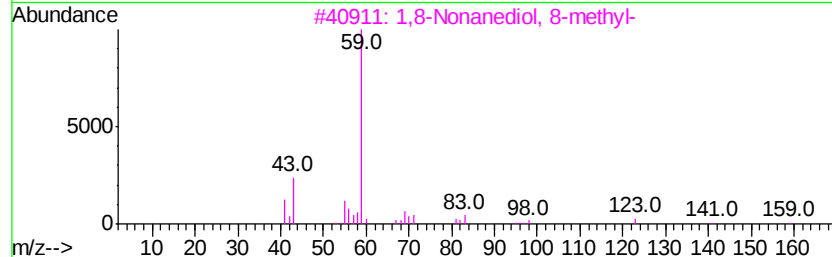
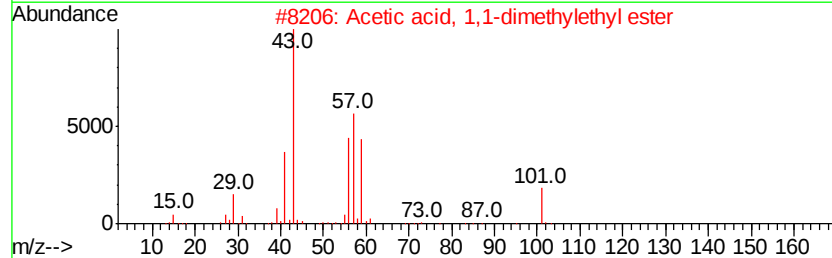
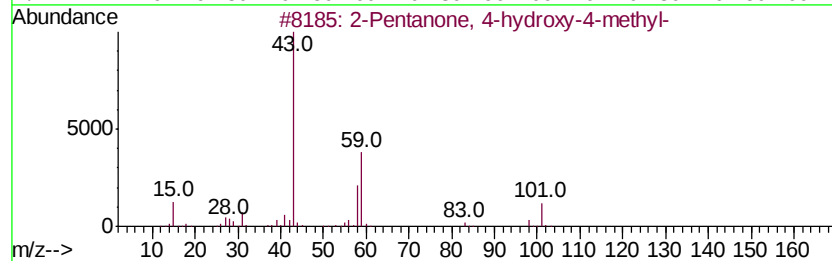
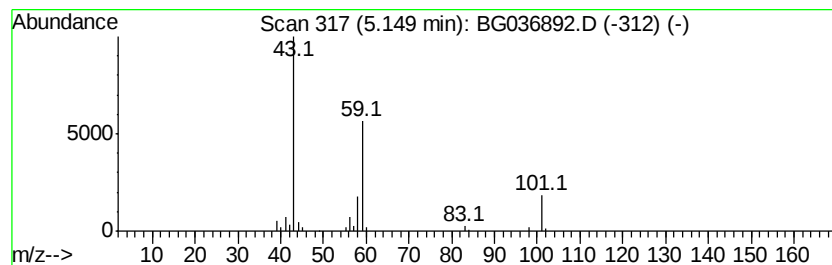
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.63 ng	43389	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	32
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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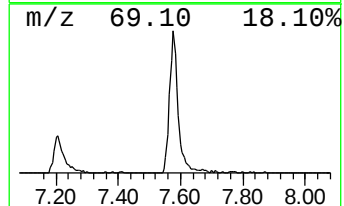
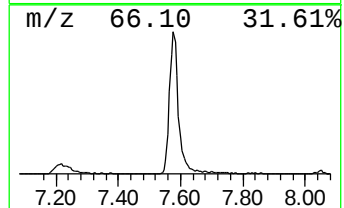
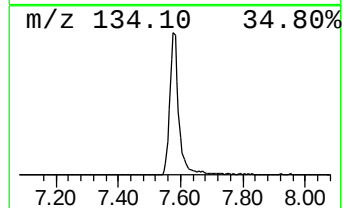
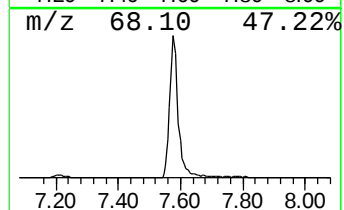
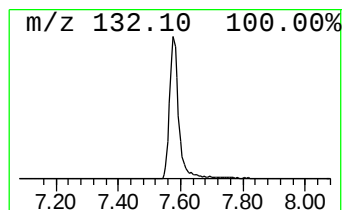
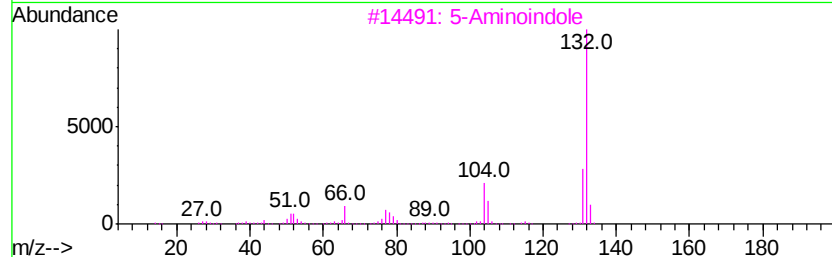
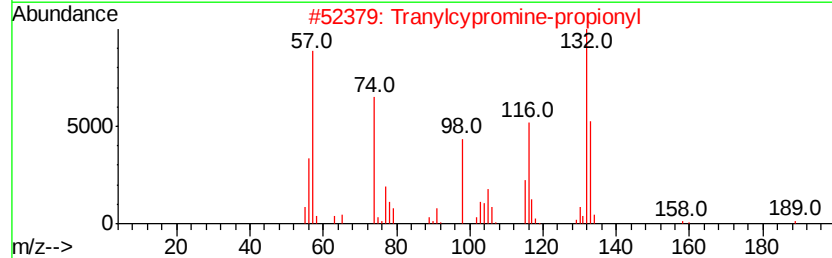
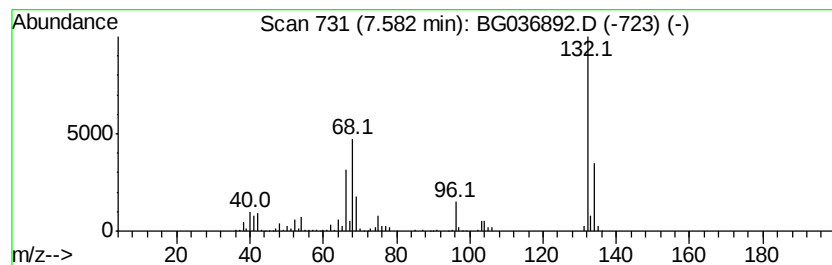
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	57.16 ng	683785	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3		5-Aminoindole	132	C8H8N2	005192-03-0	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	3.6	ng	43389	1	8.05	239251	20.0
unknown7.58	7.58	57.2	ng	683785	1	8.05	239251	20.0