

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036628.D
 Acq On : 10 Sep 2018 17:06
 Operator : JU/SJ
 Sample : J4830-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090618-TIPC-E-U-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-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	5.158	312	318	324	rBV	37149	57734	1.28%	0.307%
2	5.622	390	397	409	rBV	337857	539173	11.95%	2.864%
3	7.209	658	667	679	rBV	220317	385018	8.54%	2.045%
4	7.585	724	731	741	rBV	612044	1048099	23.24%	5.568%
5	8.055	804	811	818	rBV	206215	344085	7.63%	1.828%
6	8.378	857	866	874	rBV	824905	1429014	31.68%	7.592%
7	9.212	1000	1008	1020	rBV	701097	1271967	28.20%	6.758%
8	10.857	1280	1288	1298	rVB	291677	522149	11.58%	2.774%
9	13.302	1695	1704	1719	rBV	1978548	3183455	70.57%	16.913%
10	14.124	1838	1844	1851	rBV	86413	122134	2.71%	0.649%
11	14.676	1931	1938	1947	rBV2	492135	803306	17.81%	4.268%
12	15.969	2153	2158	2163	rVB	45655	59025	1.31%	0.314%
13	16.163	2184	2191	2211	rBV	947370	1484183	32.90%	7.885%
14	17.420	2398	2405	2414	rBV2	646193	995872	22.08%	5.291%
15	20.029	2843	2849	2861	rBV	3329846	4510791	100.00%	23.964%
16	21.686	3124	3131	3141	rBV	660993	1096552	24.31%	5.826%
17	24.900	3667	3678	3696	rBV	296773	970462	21.51%	5.156%

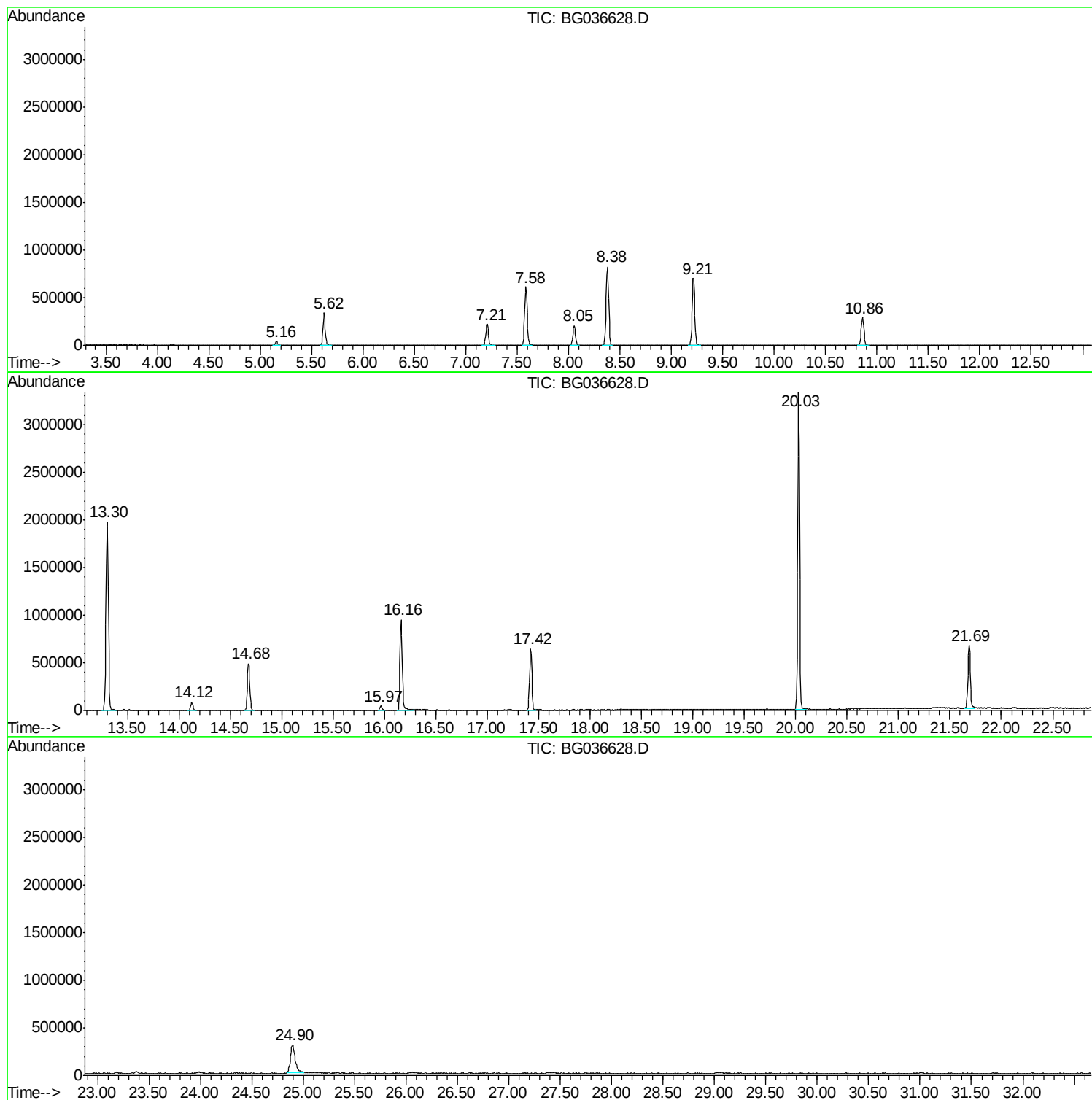
Sum of corrected areas: 18823019

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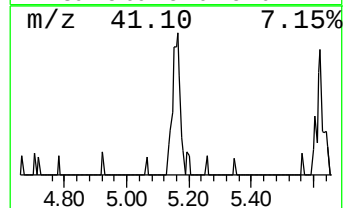
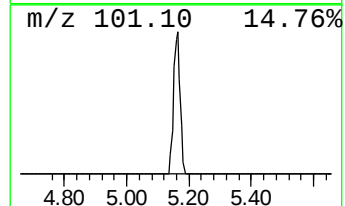
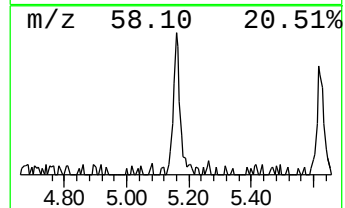
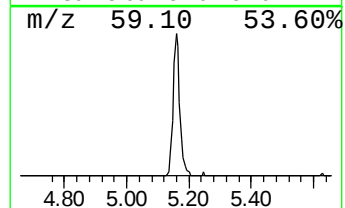
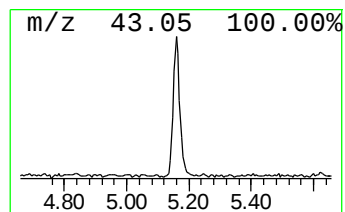
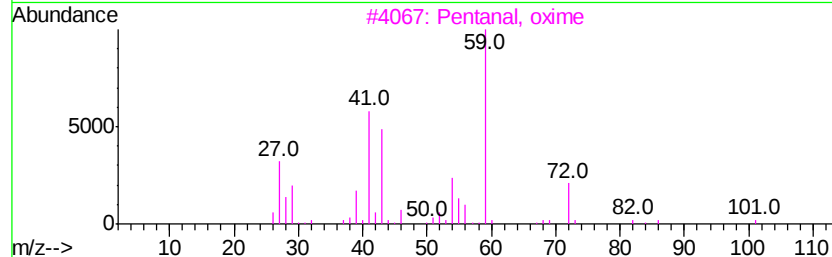
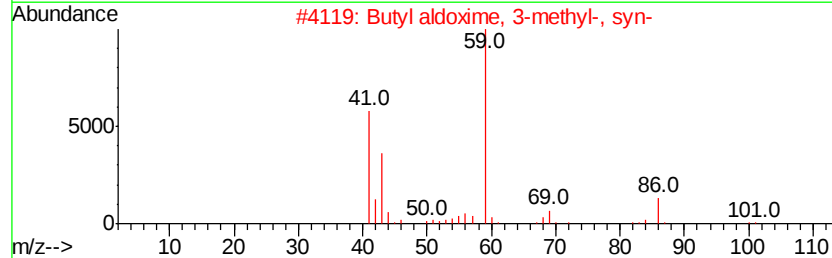
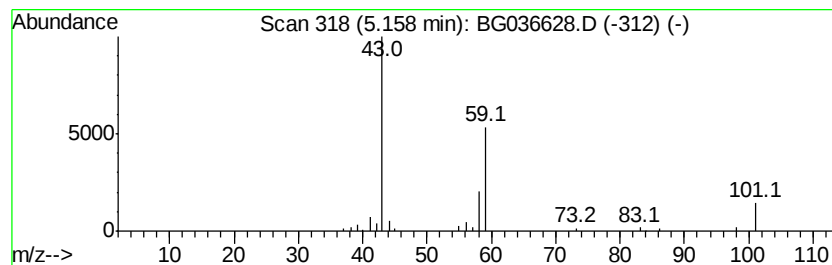
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	3.36 ng	57734	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	78
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
3			Pentanal, oxime	101	C5H11NO	000628-79-5	9
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	9



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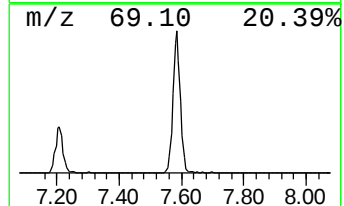
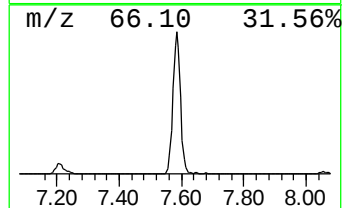
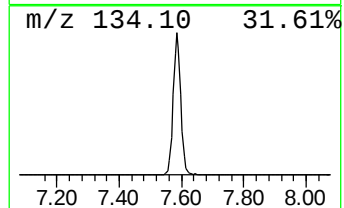
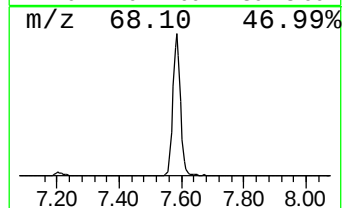
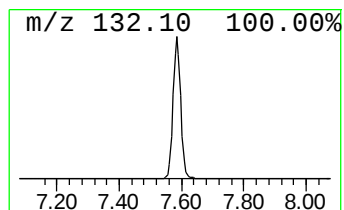
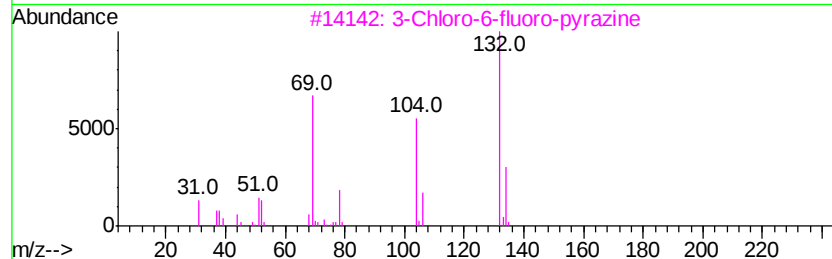
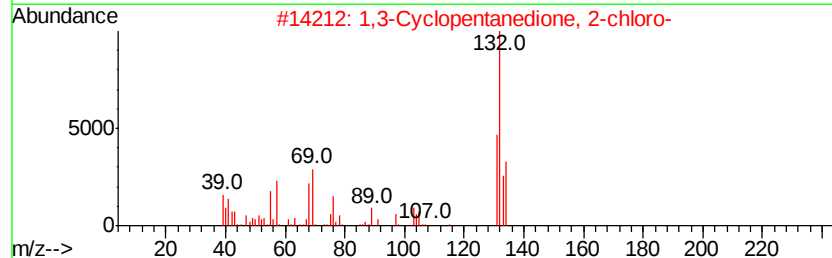
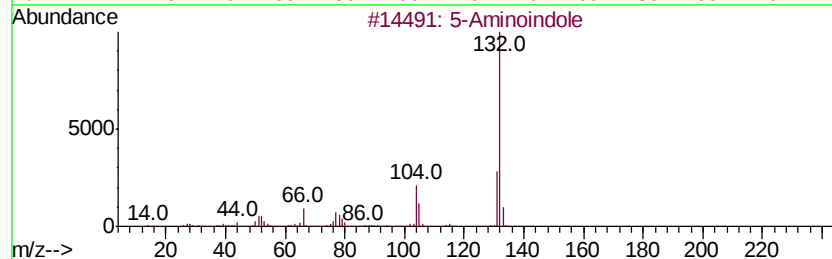
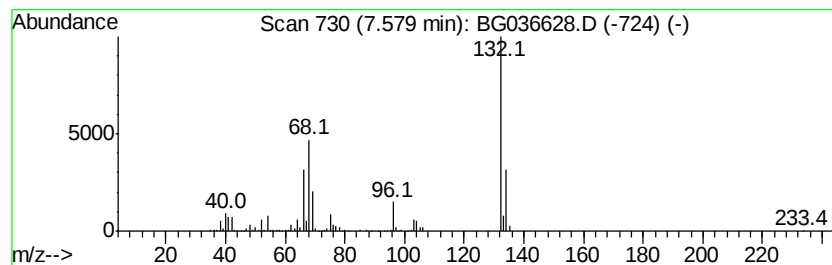
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	60.92 ng	1048100	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
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	3.4	ng	57734	1	8.05	344085	20.0
unknown7.58	7.58	60.9	ng	1048100	1	8.05	344085	20.0