

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061818\
 Data File : BG035215.D
 Acq On : 18 Jun 2018 12:40
 Operator : SJ/JU
 Sample : J3505-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-55-COMP

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-BG060718.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	5.173	315	321	334	rVB	118551	195199	2.81%	0.802%
2	5.637	393	400	413	rBV	916350	1503683	21.66%	6.176%
3	7.217	659	669	679	rBV	831614	1441830	20.77%	5.922%
4	7.593	726	733	741	rBV	1019509	1760926	25.37%	7.232%
5	8.063	806	813	820	rBV	190187	322056	4.64%	1.323%
6	8.386	859	868	880	rBV	762175	1345152	19.38%	5.525%
7	9.220	1000	1010	1020	rBV	655841	1179821	17.00%	4.845%
8	10.865	1283	1290	1296	rBV	228747	408451	5.88%	1.677%
9	13.309	1698	1706	1719	rBV	1743379	2752968	39.66%	11.306%
10	14.684	1933	1940	1947	rBV2	369096	603447	8.69%	2.478%
11	16.170	2186	2193	2207	rBV	1657617	2594921	37.38%	10.657%
12	17.427	2400	2407	2412	rBV	579877	910386	13.12%	3.739%
13	20.035	2844	2851	2857	rBV	5146245	6941491	100.00%	28.509%
14	21.692	3126	3133	3141	rVB2	800218	1283562	18.49%	5.272%
15	24.917	3668	3682	3696	rBV2	380003	1104906	15.92%	4.538%

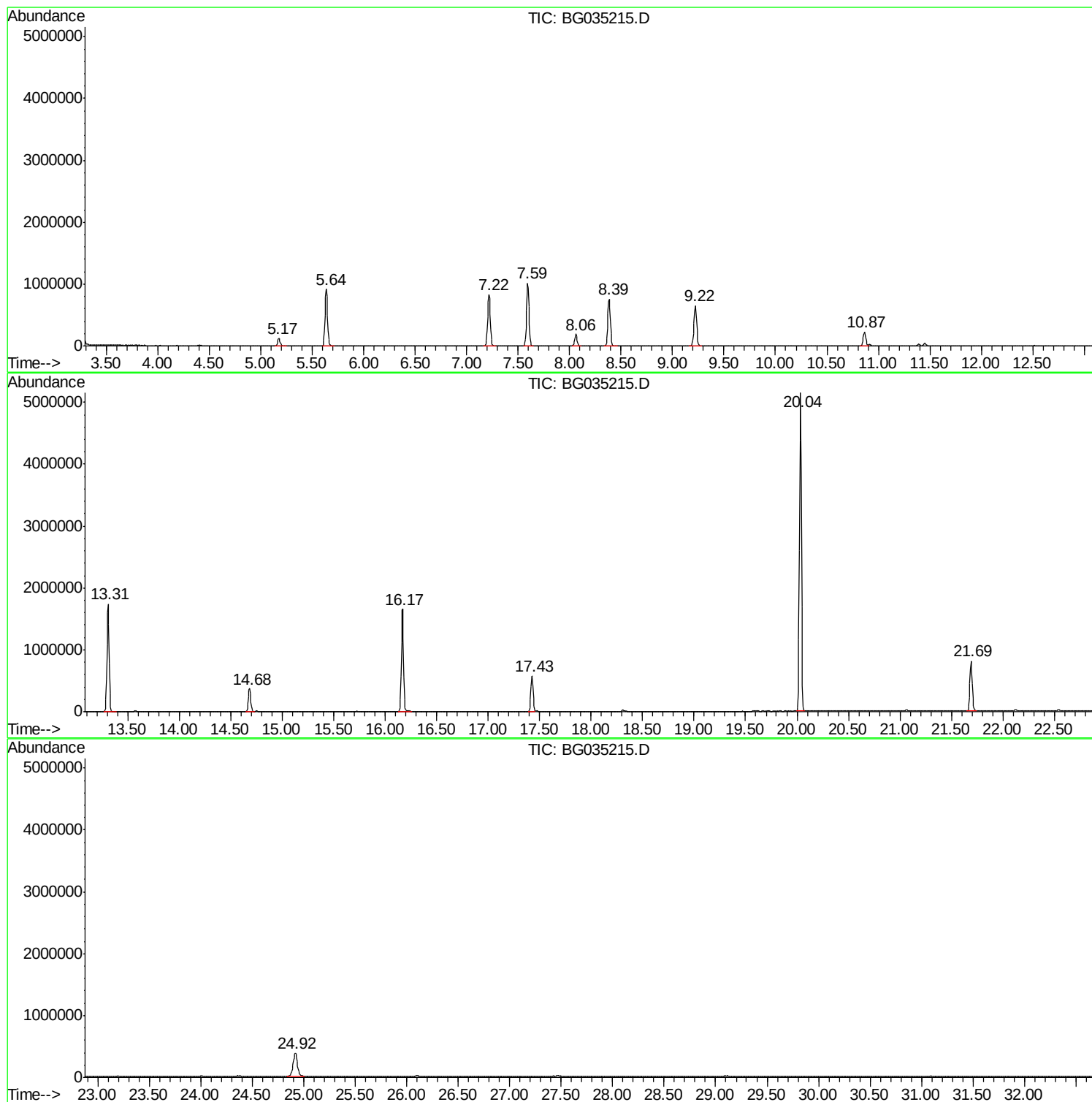
Sum of corrected areas: 24348799

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.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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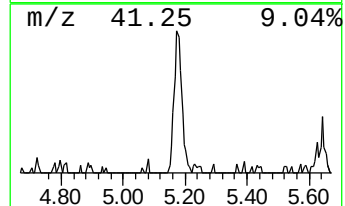
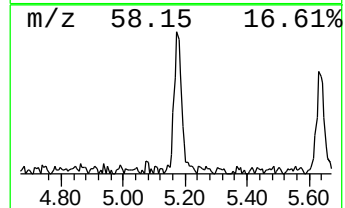
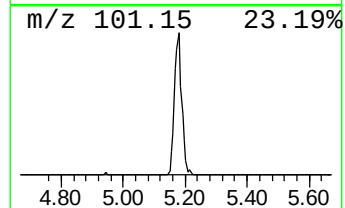
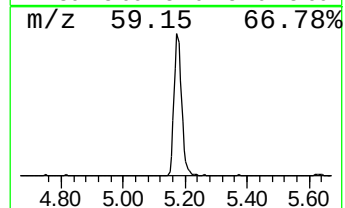
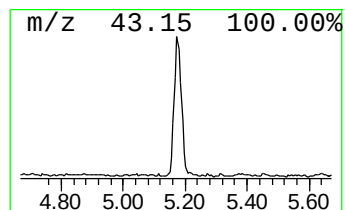
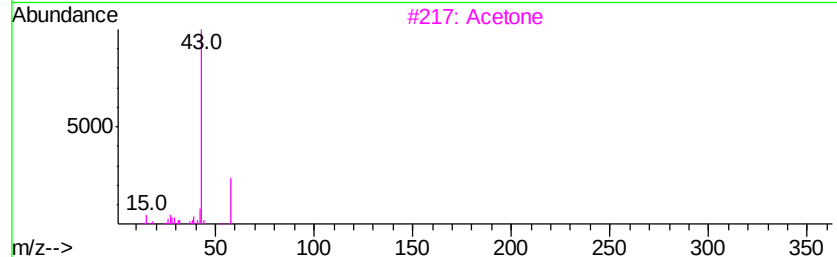
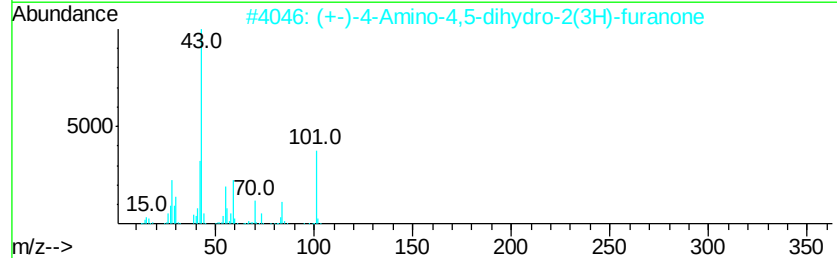
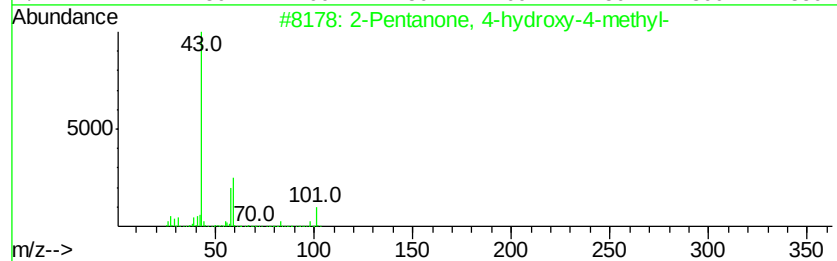
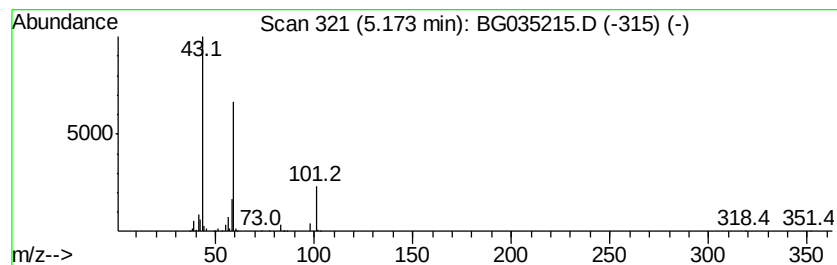
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	12.12 ng	195199	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
3		Acetone	58	C3H6O	000067-64-1	9
4		Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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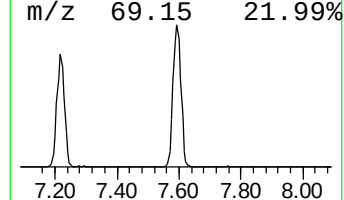
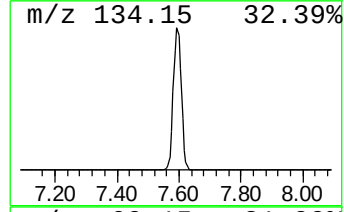
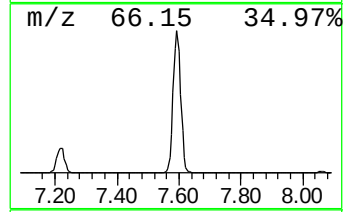
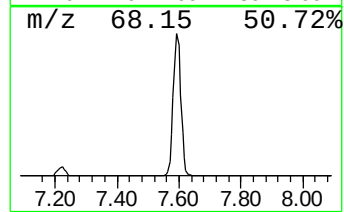
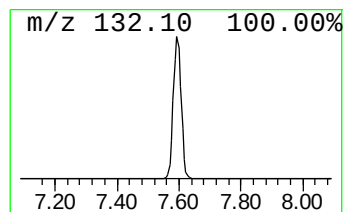
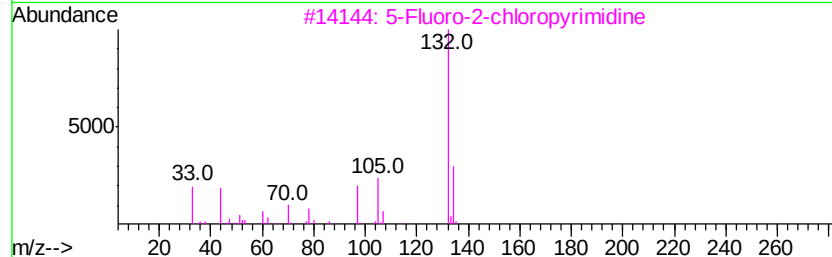
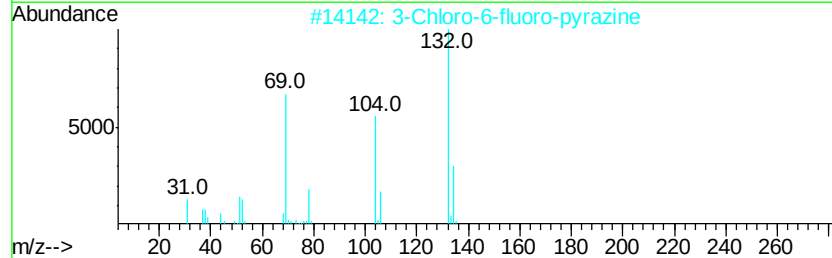
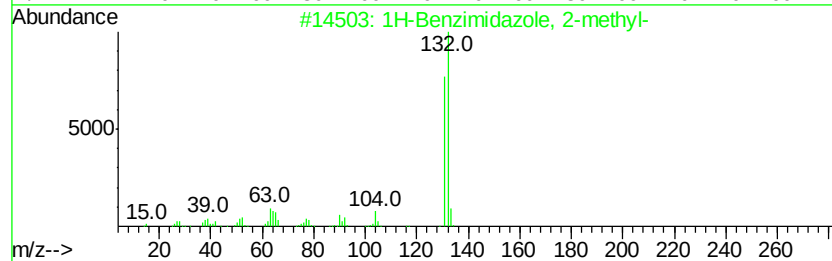
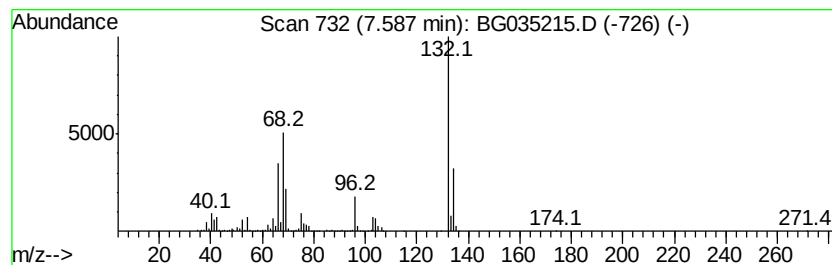
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	109.36 ng	1760930	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	12.1	ng	195199	1	8.06	322056	20.0
unknown7.59	7.59	109.4	ng	1760930	1	8.06	322056	20.0