

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091718\
Data File : BG036777.D
Acq On : 18 Sep 2018 7:20
Operator : JU/SJ
Sample : J4929-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SP-PIT-12

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	310	319	326	rVB	118576	189244	4.72%	0.927%
2	5.620	389	397	407	rBV	947095	1529695	38.16%	7.495%
3	7.206	659	667	678	rBV	891111	1534397	38.28%	7.518%
4	7.582	722	731	741	rBV	988901	1674219	41.77%	8.204%
5	8.053	803	811	818	rBV	218671	365918	9.13%	1.793%
6	8.376	858	866	874	rVB	733828	1289709	32.18%	6.320%
7	9.210	999	1008	1020	rBV	635947	1129707	28.18%	5.536%
8	10.855	1279	1288	1295	rBV	281244	511899	12.77%	2.508%
9	13.299	1696	1704	1714	rBV	1607425	2552816	63.69%	12.509%
10	14.674	1931	1938	1946	rBV2	442193	716562	17.88%	3.511%
11	16.161	2184	2191	2205	rBV	1206740	1869195	46.63%	9.159%
12	17.418	2398	2405	2412	rBV	625578	930809	23.22%	4.561%
13	20.027	2842	2849	2858	rBV	2904638	4008370	100.00%	19.641%
14	21.684	3124	3131	3140	rBV	634912	1070729	26.71%	5.247%
15	24.892	3666	3677	3695	rVB2	339497	1035122	25.82%	5.072%

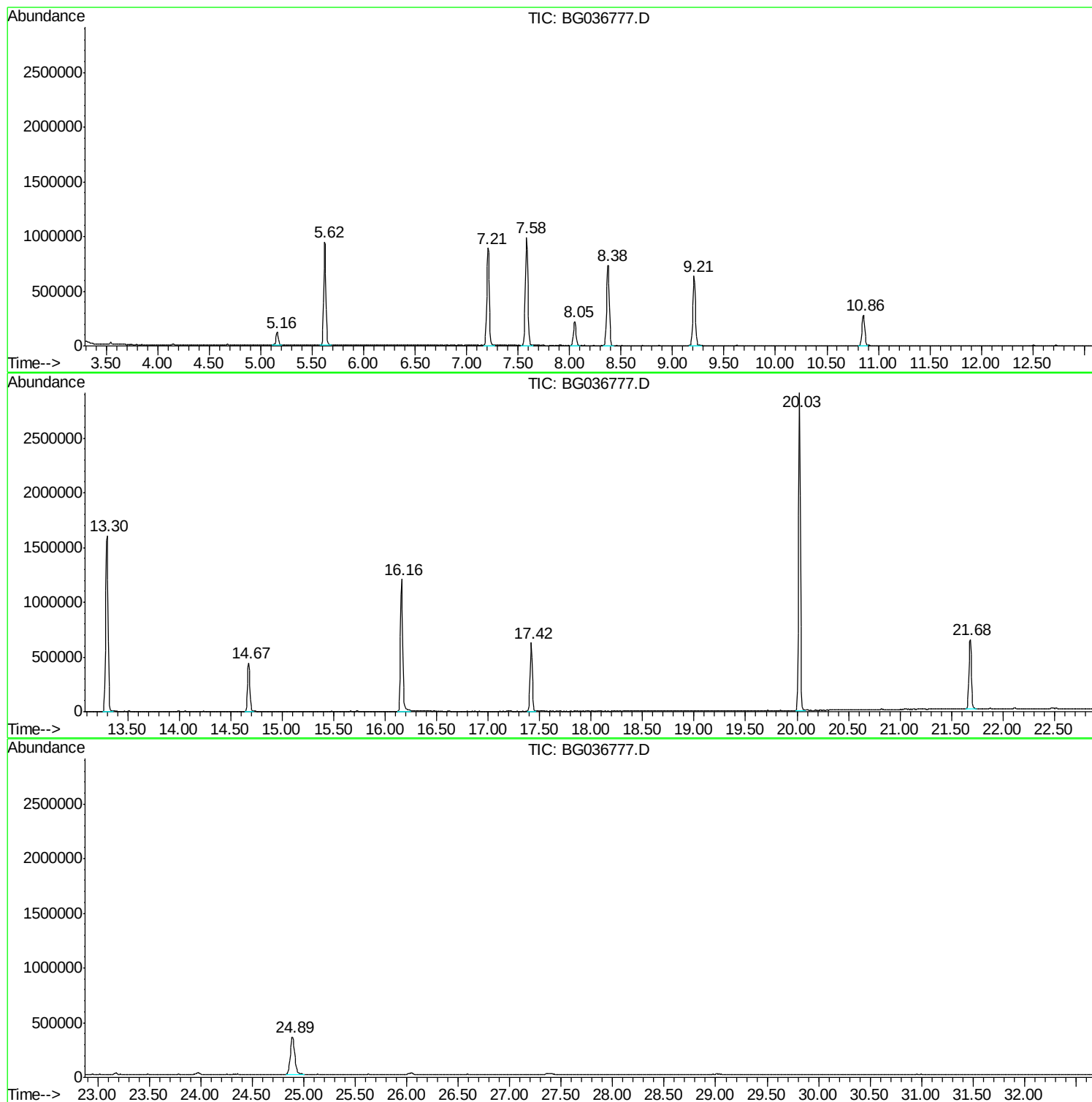
Sum of corrected areas: 20408391

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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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Instrument :
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ClientSampled :
SP-PIT-12

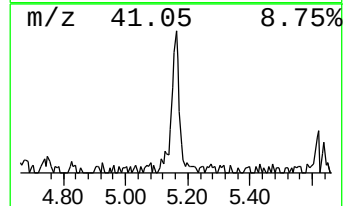
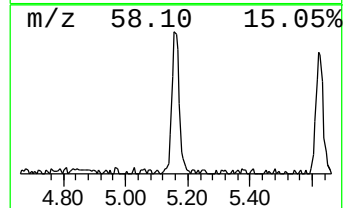
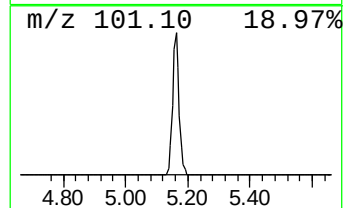
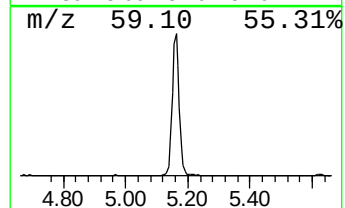
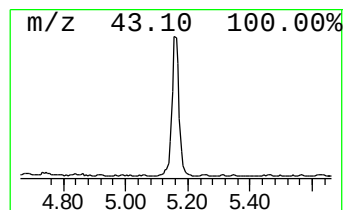
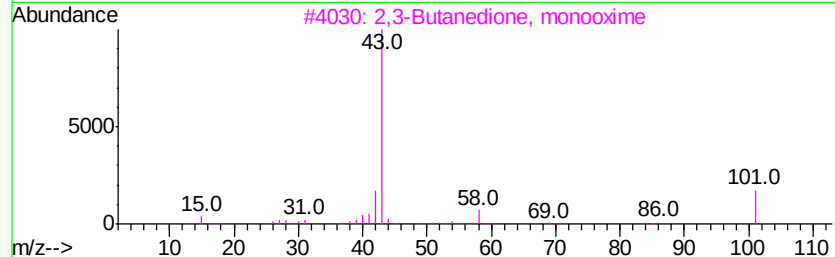
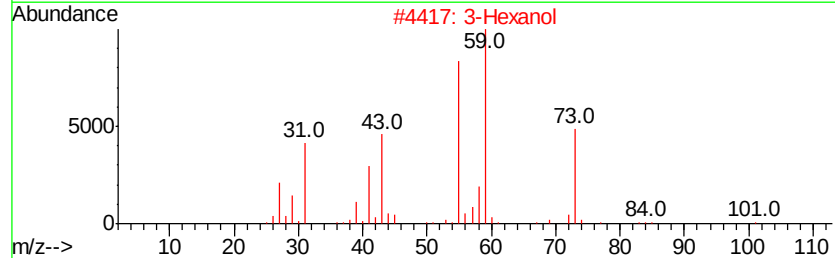
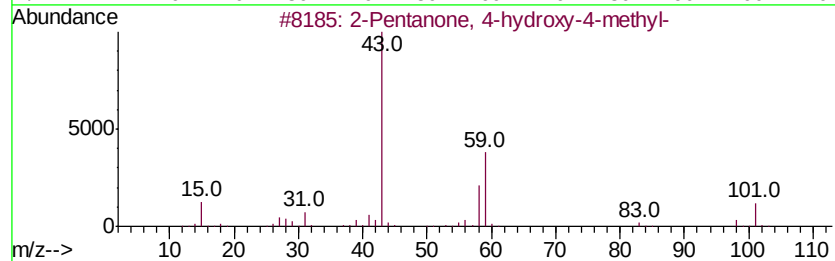
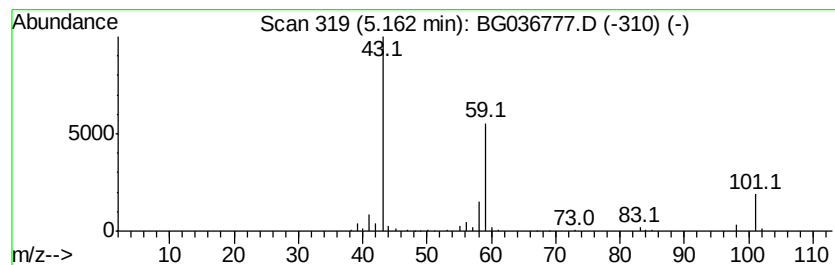
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	10.34 ng	189244	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		3-Hexanol	102	C6H14O	000623-37-0	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
5		Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	9



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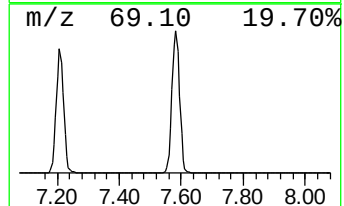
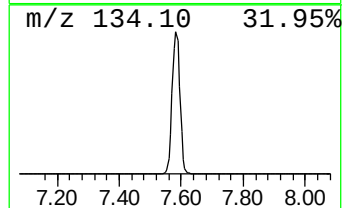
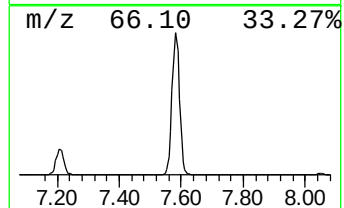
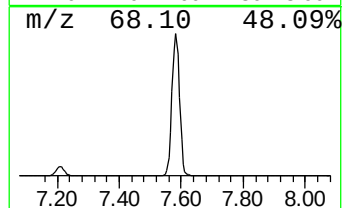
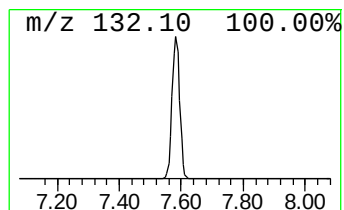
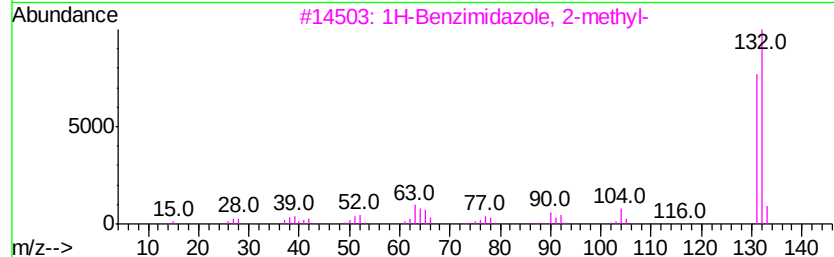
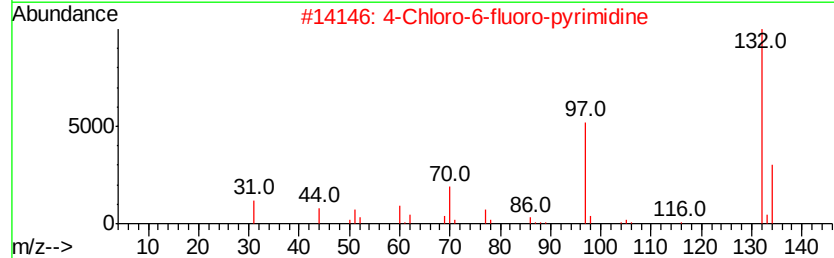
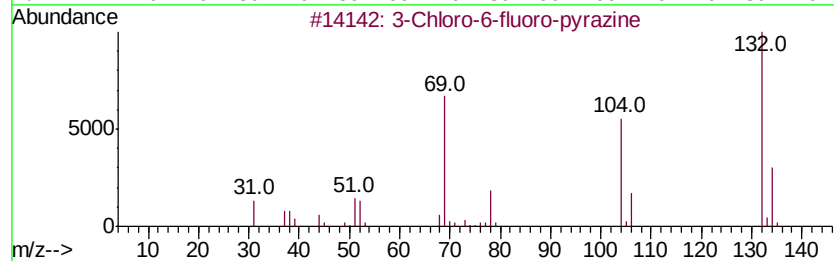
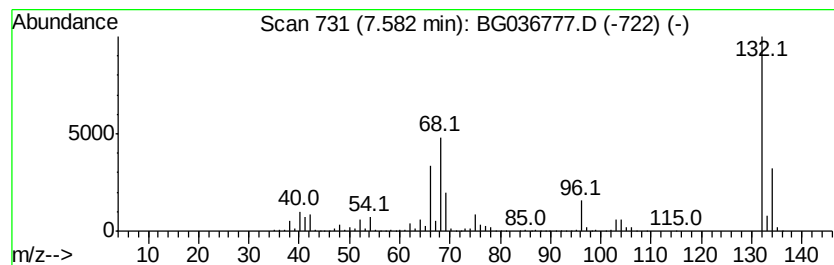
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	91.51 ng	1674220	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		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
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...	5.16	10.3	ng	189244	1	8.05	365918	20.0
unknown7.58	7.58	91.5	ng	1674220	1	8.05	365918	20.0