

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010219\
 Data File : BG038899.D
 Acq On : 2 Jan 2019 19:35
 Operator : JU/SJ
 Sample : PB116088BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116088BL

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-BG121218.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	5.130	341	347	356	rVB2	24173	40417	1.69%	0.393%
2	5.600	419	427	437	rBV	445829	729312	30.43%	7.094%
3	7.210	691	701	714	rBV	449643	793959	33.12%	7.722%
4	7.574	754	763	777	rBV	435339	778129	32.46%	7.568%
5	8.044	835	843	853	rVB	82916	143947	6.01%	1.400%
6	8.368	891	898	907	rVB	336219	587563	24.51%	5.715%
7	9.225	1035	1044	1054	rBV	254858	473122	19.74%	4.602%
8	10.865	1314	1323	1332	rBV	101491	192592	8.04%	1.873%
9	13.303	1731	1738	1751	rBV	783997	1250769	52.18%	12.166%
10	14.684	1966	1973	1986	rBV2	188906	307551	12.83%	2.991%
11	16.176	2220	2227	2243	rBV2	634556	1090913	45.51%	10.611%
12	17.433	2434	2441	2455	rBV2	291104	451328	18.83%	4.390%
13	20.036	2878	2884	2897	rVB	1823824	2396901	100.00%	23.313%
14	21.711	3163	3169	3178	rBV2	290457	504263	21.04%	4.905%
15	23.150	3409	3414	3420	rBV4	15875	35766	1.49%	0.348%
16	23.967	3544	3553	3559	rBV4	14992	35227	1.47%	0.343%
17	25.001	3720	3729	3742	rVB2	155632	469413	19.58%	4.566%

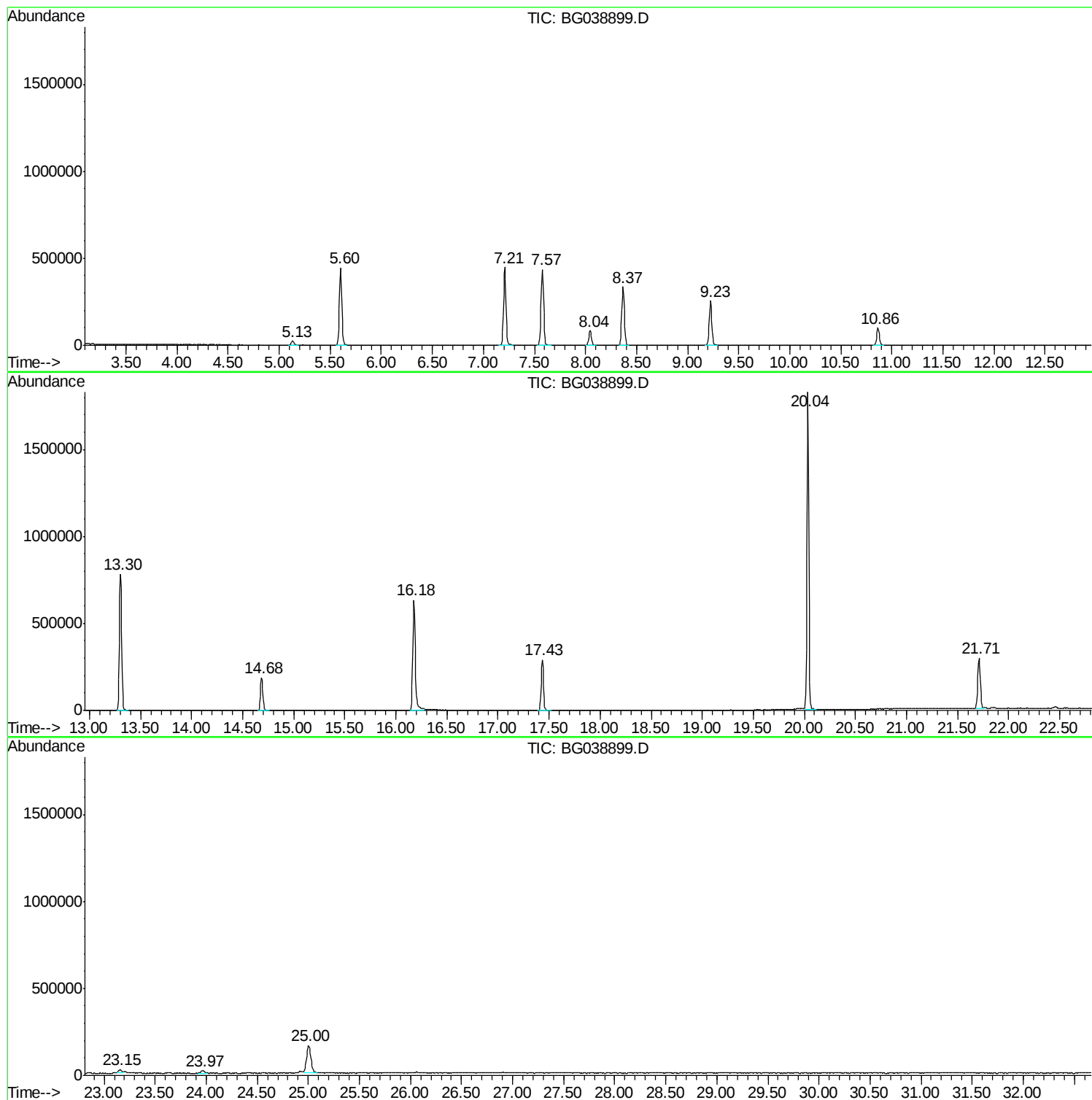
Sum of corrected areas: 10281172

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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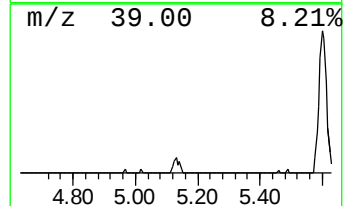
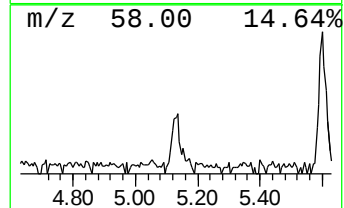
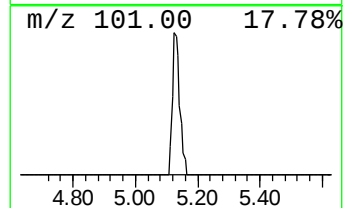
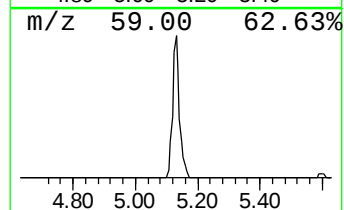
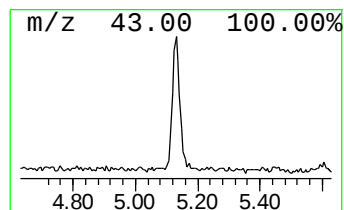
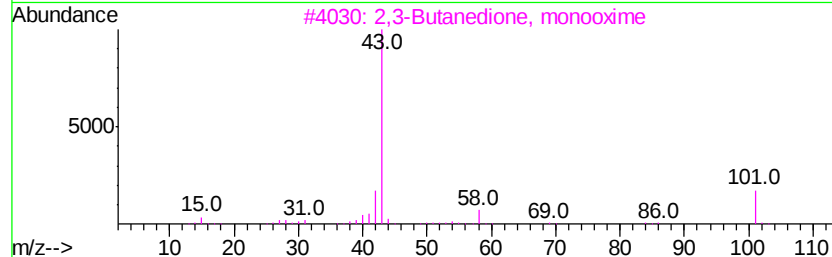
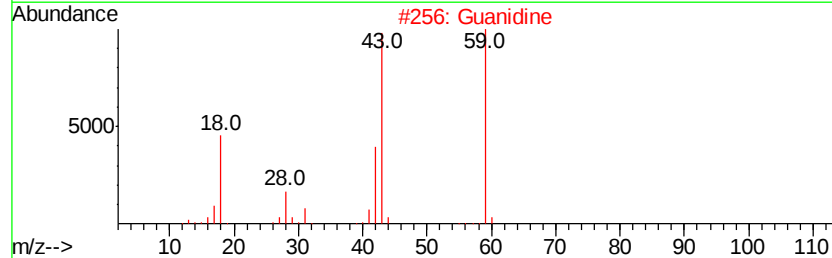
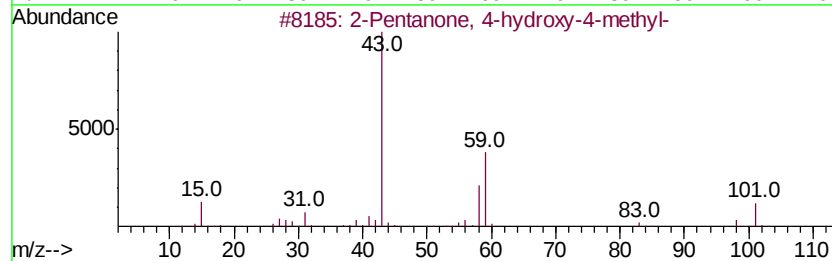
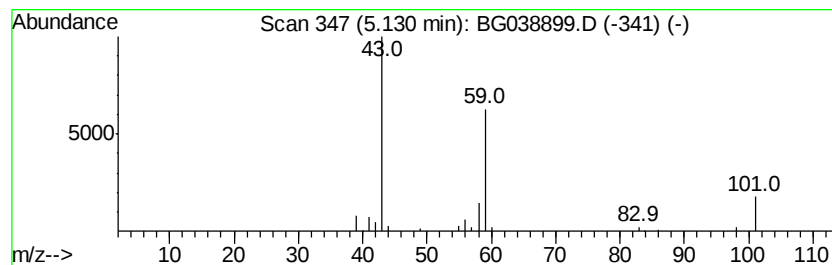
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	5.62 ng	40417	1,4-Dichlorobenzene-d4	8.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	53
2	Guanidine	59	CH5N3		000113-00-8	9
3	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	9
4	4-Penten-2-one, 4-methyl-	98	C6H10O		003744-02-3	9
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2		042781-12-4	9



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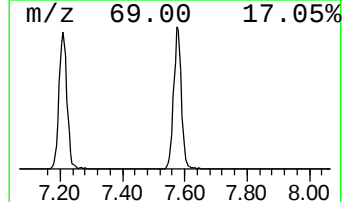
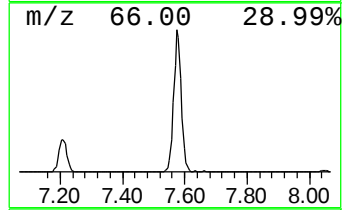
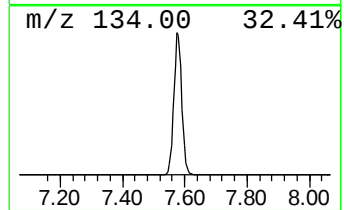
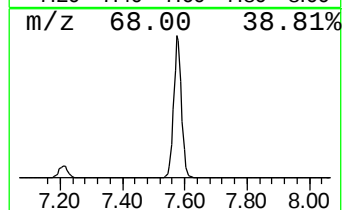
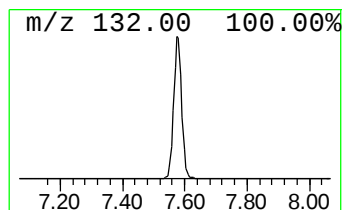
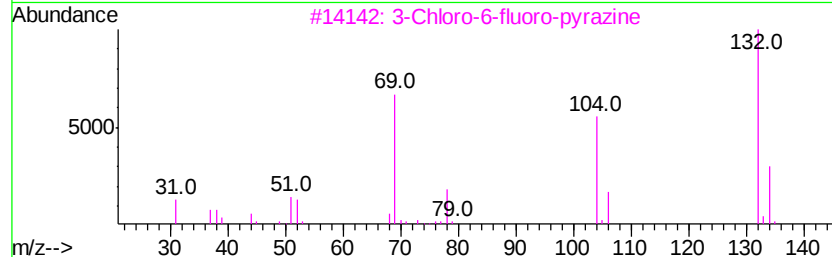
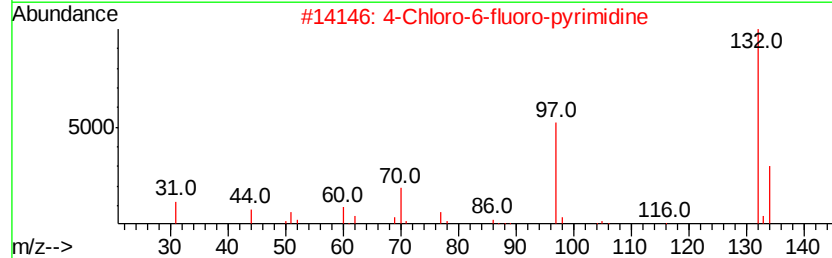
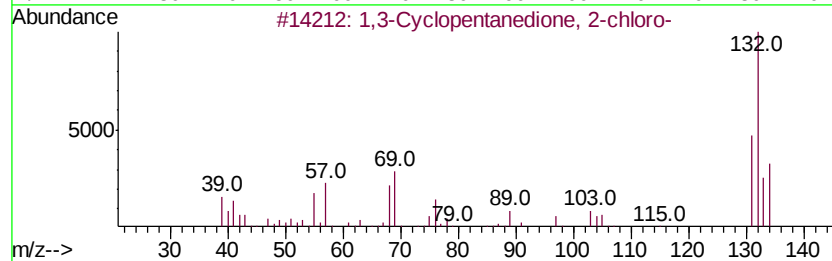
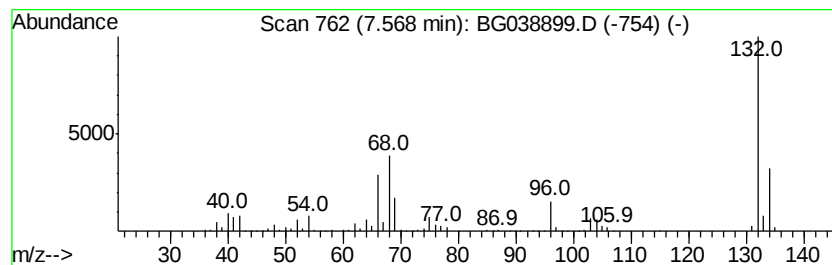
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	108.11 ng	778129	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	37
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	5.6	ng	40417	1	8.04	143947	20.0
unknown7.57	7.57	108.1	ng	778129	1	8.04	143947	20.0