

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080718\
 Data File : BG036170.D
 Acq On : 7 Aug 2018 20:21
 Operator : JU/SJ
 Sample : PB111837BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111837BL

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-BG071218.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	5.126	308	313	324	rBV	80791	134167	4.09%	1.020%
2	5.596	386	393	409	rBV	541946	950074	28.97%	7.226%
3	7.182	655	663	680	rBV	557105	1016751	31.00%	7.733%
4	7.547	717	725	734	rBV	582976	1005863	30.67%	7.650%
5	8.011	798	804	811	rBV	103110	177591	5.41%	1.351%
6	8.334	851	859	868	rBV	433156	788668	24.05%	5.998%
7	9.174	993	1002	1015	rBV	287258	611998	18.66%	4.655%
8	10.819	1274	1282	1295	rBV2	86465	210230	6.41%	1.599%
9	13.256	1690	1697	1710	rBV	925610	1550242	47.27%	11.791%
10	14.637	1925	1932	1947	rBV	173436	337164	10.28%	2.564%
11	16.123	2178	2185	2207	rBV2	789353	1393585	42.49%	10.599%
12	17.380	2392	2399	2411	rBV	277803	497537	15.17%	3.784%
13	19.989	2835	2843	2853	rBV	2283617	3279780	100.00%	24.945%
14	21.639	3117	3124	3137	rBV2	311515	623865	19.02%	4.745%
15	24.829	3655	3667	3681	rBV	174194	570541	17.40%	4.339%

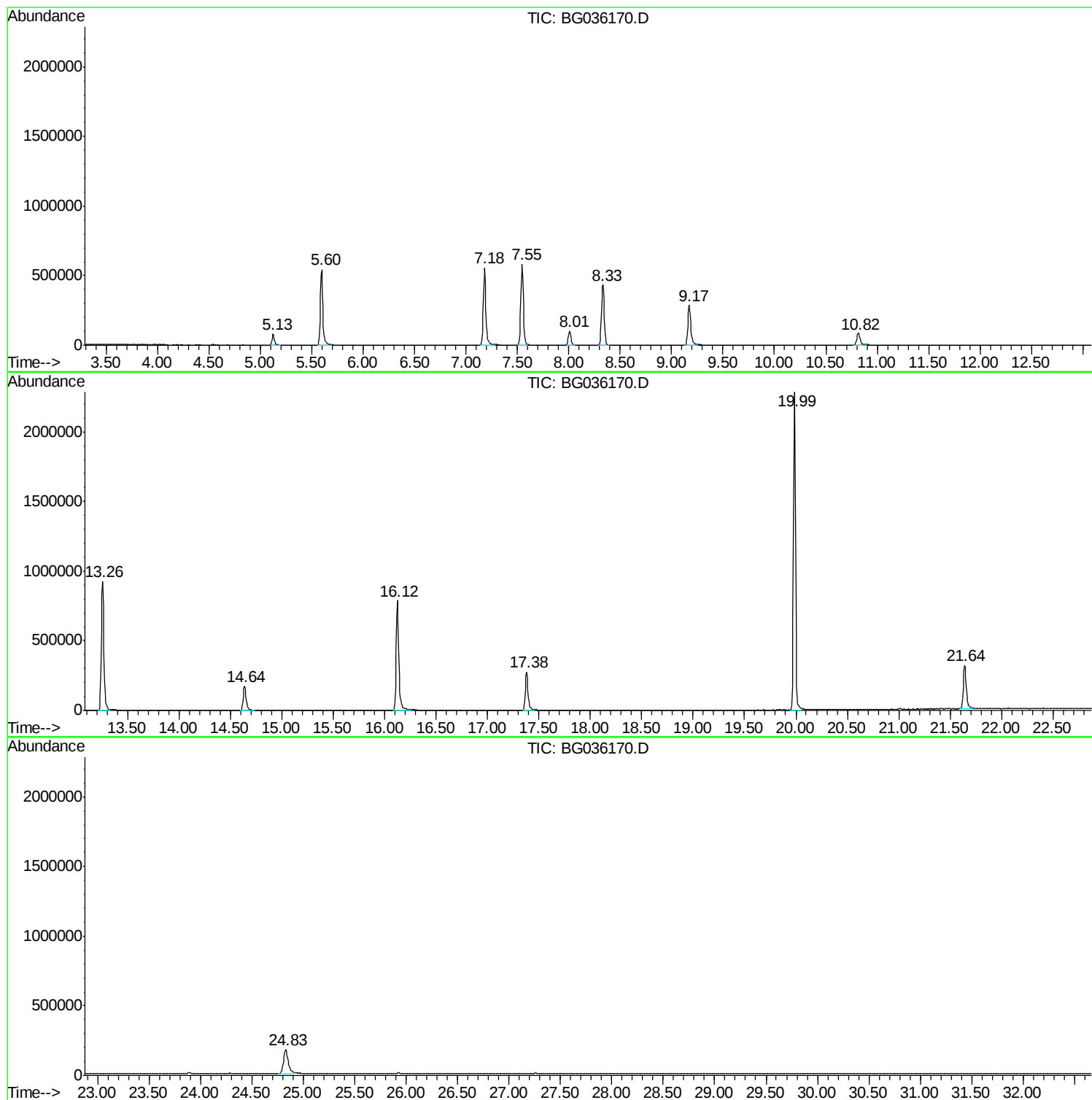
Sum of corrected areas: 13148056

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



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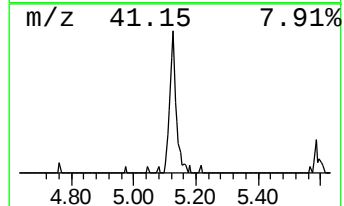
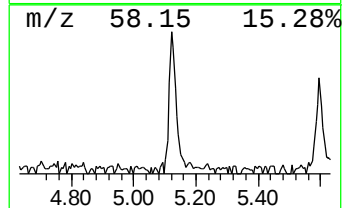
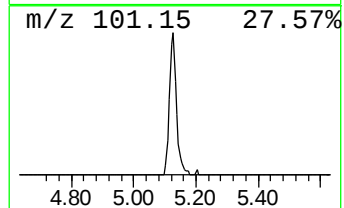
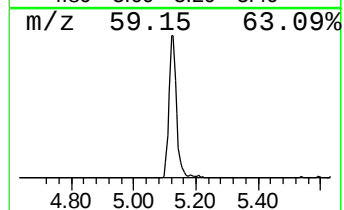
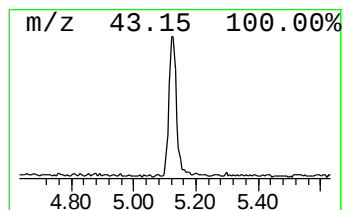
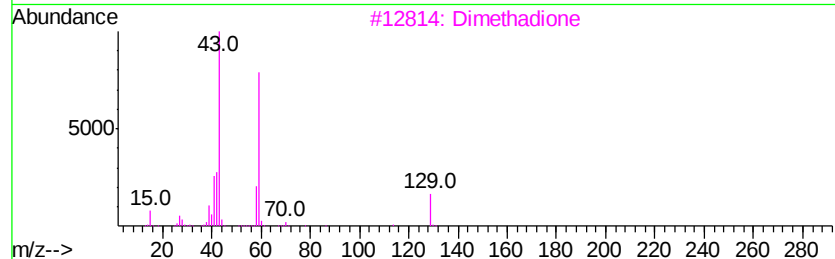
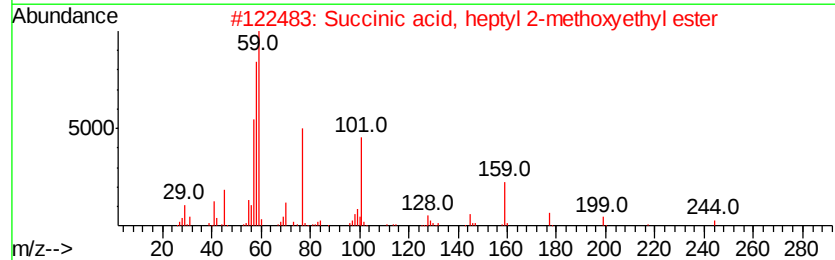
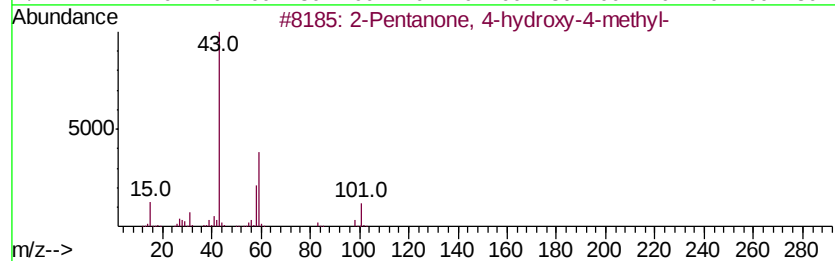
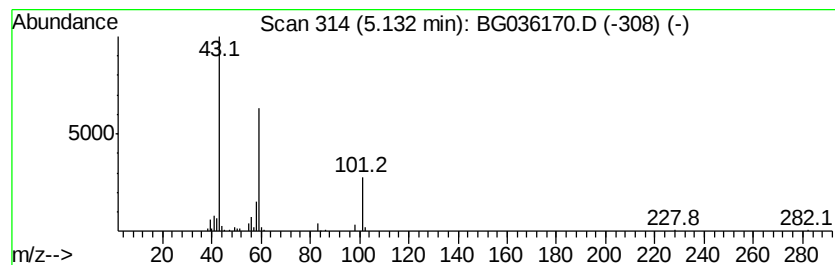
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.13	15.11 ng	134167	1,4-Dichlorobenzene-d4	8.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	28
3	Dimethadione	129	C5H7NO3	000695-53-4	28
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	2-Octanol, 2-methyl-	144	C9H20O	000628-44-4	25



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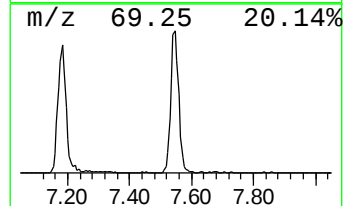
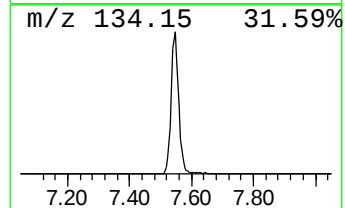
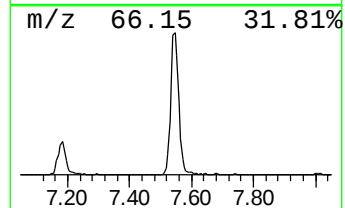
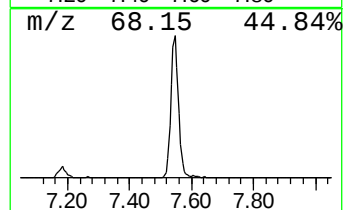
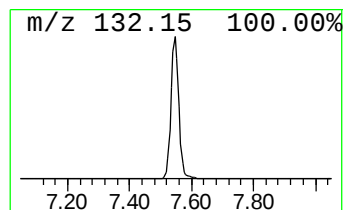
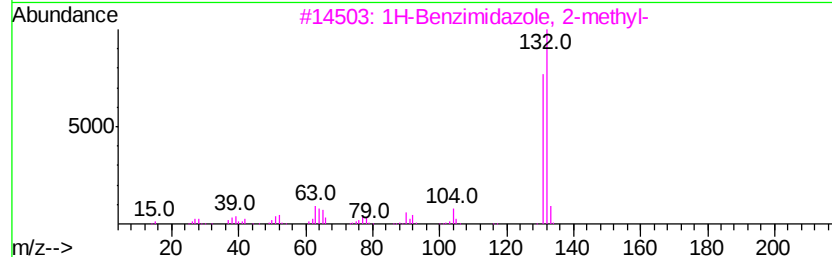
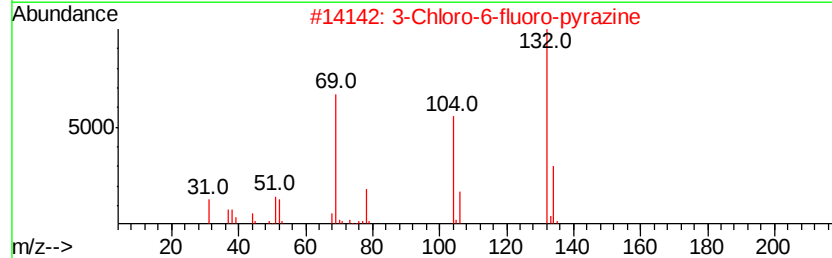
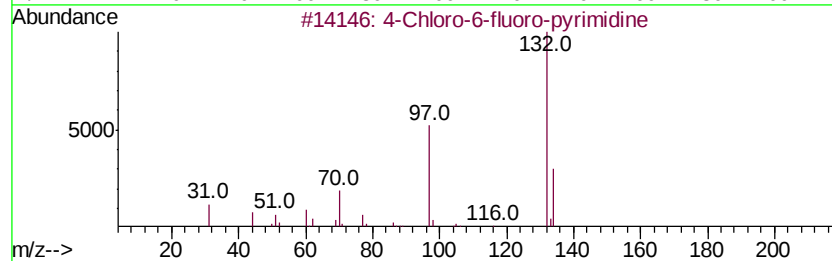
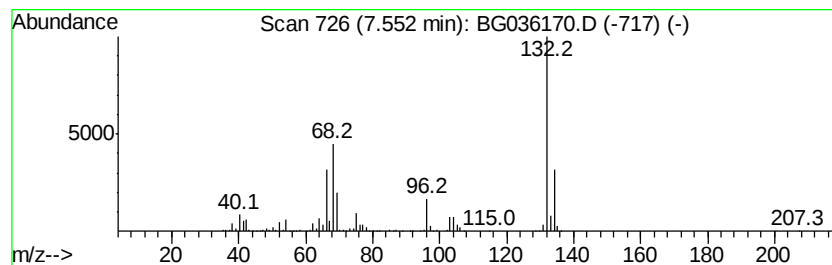
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Peak Number 2 unknown7.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	113.28 ng	1005860	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	15.1	ng	134167	1	8.00	177591	20.0
unknown7.55	7.55	113.3	ng	1005860	1	8.00	177591	20.0