

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091918\
 Data File : BG036842.D
 Acq On : 20 Sep 2018 10:09
 Operator : JU/SJ
 Sample : PB112998BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112998BL

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	5.151	311	317	324	rBV	40938	64705	1.81%	0.332%
2	5.621	390	397	410	rBV	988513	1564165	43.72%	8.014%
3	7.207	659	667	678	rBV	995666	1705642	47.68%	8.739%
4	7.577	723	730	741	rBV	946562	1610919	45.03%	8.254%
5	8.047	803	810	818	rVB	204895	357748	10.00%	1.833%
6	8.370	857	865	872	rBV	697683	1213292	33.91%	6.216%
7	9.205	999	1007	1018	rBV	584423	1032985	28.87%	5.293%
8	10.850	1277	1287	1296	rBV	287345	509675	14.25%	2.611%
9	13.294	1695	1703	1714	rBV	1567196	2450704	68.50%	12.557%
10	14.669	1927	1937	1947	rBV2	470640	750261	20.97%	3.844%
11	16.155	2182	2190	2202	rBV	1108636	1774784	49.61%	9.093%
12	17.413	2395	2404	2413	rBV2	660186	989133	27.65%	5.068%
13	20.021	2841	2848	2856	rBV	2719614	3577516	100.00%	18.330%
14	21.678	3123	3130	3142	rBV	613595	1019584	28.50%	5.224%
15	24.880	3666	3675	3693	rVB2	285929	896287	25.05%	4.592%

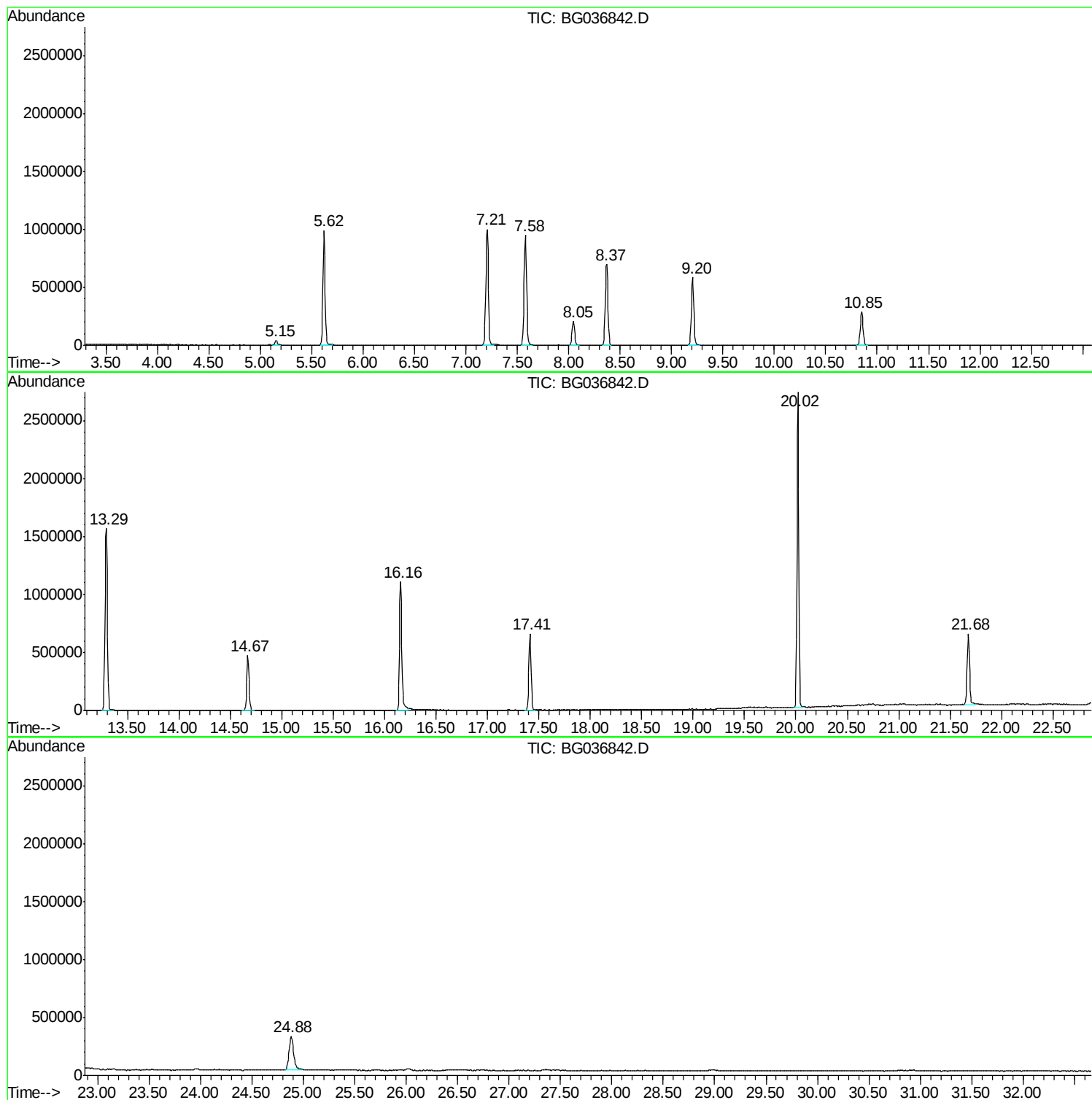
Sum of corrected areas: 19517400

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



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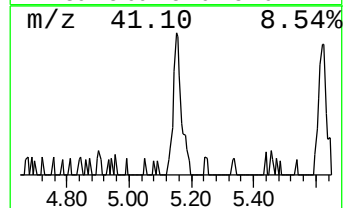
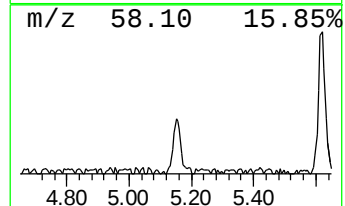
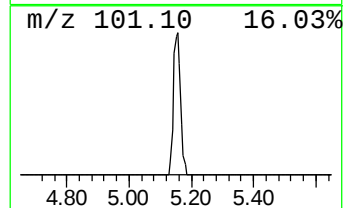
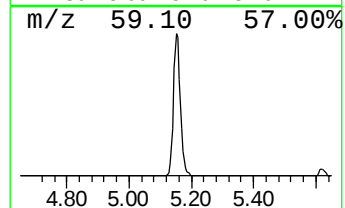
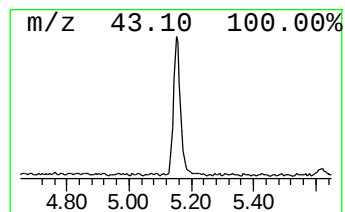
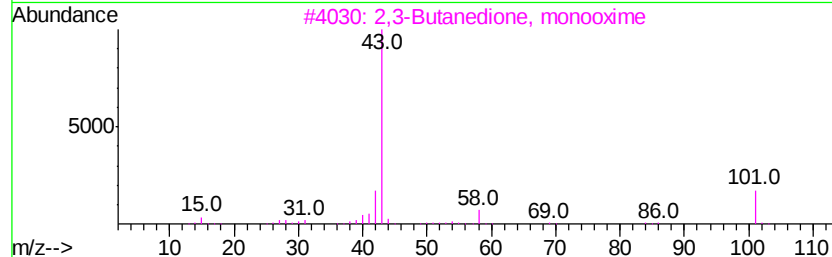
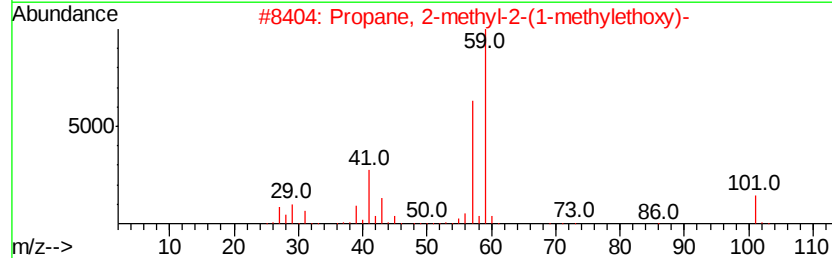
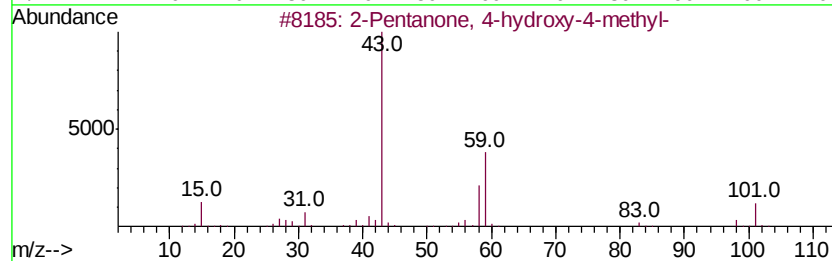
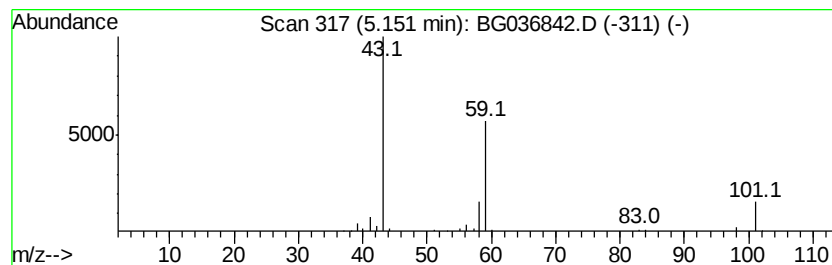
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.62 ng	64705	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	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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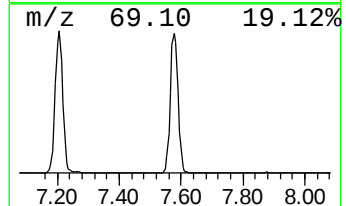
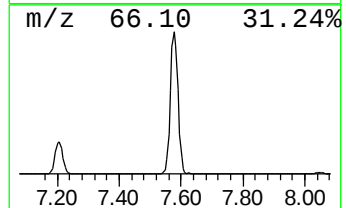
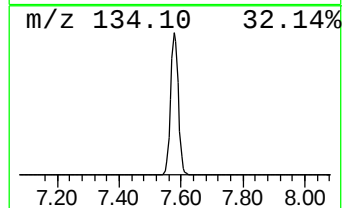
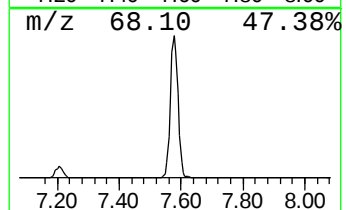
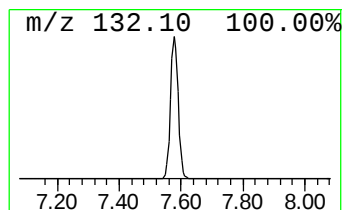
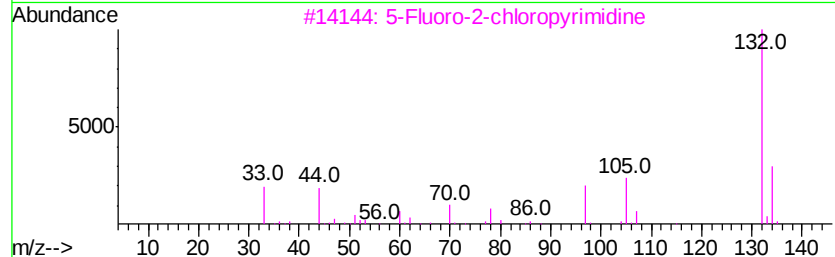
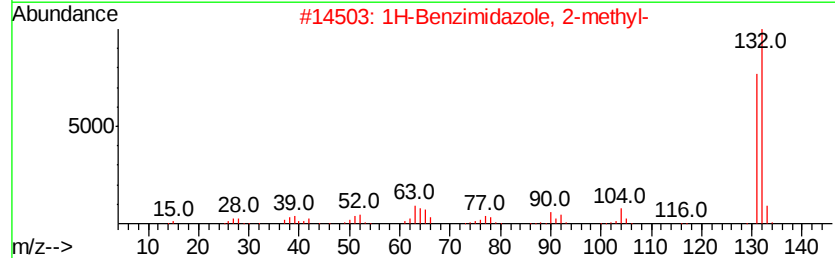
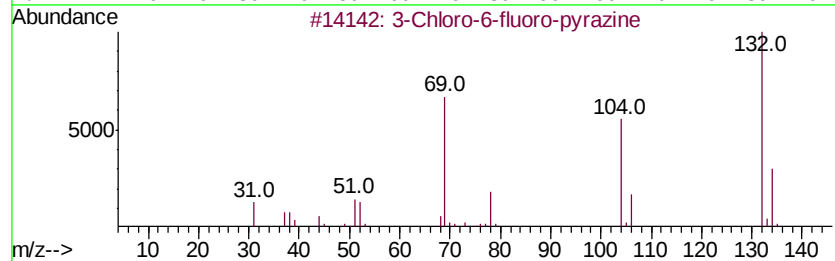
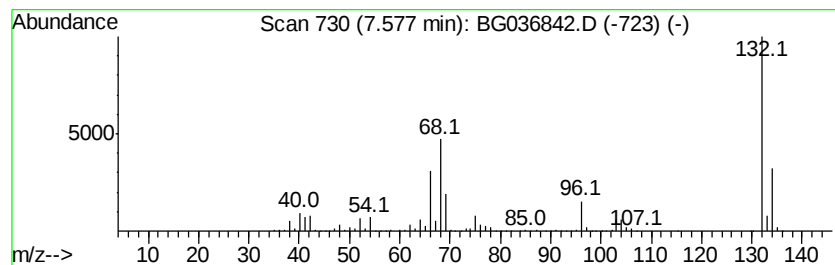
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	90.06 ng	1610920	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		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	3.6	ng	64705	1	8.05	357748	20.0
unknown7.58	7.58	90.1	ng	1610920	1	8.05	357748	20.0