

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102919\
 Data File : BG043373.D
 Acq On : 29 Oct 2019 14:27
 Operator : JU
 Sample : PB124338BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124338BL

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-BG102119.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	5.127	341	347	358	rBV	154098	251174	5.60%	1.312%
2	5.614	422	430	447	rBV	788718	1393231	31.09%	7.280%
3	7.213	694	702	723	rBV	839186	1547828	34.54%	8.087%
4	7.565	753	762	773	rBV	877270	1558508	34.78%	8.143%
5	8.023	832	840	851	rVB	133005	234175	5.23%	1.224%
6	8.347	888	895	905	rBV	675330	1183683	26.41%	6.185%
7	9.187	1028	1038	1048	rBV	459673	903552	20.16%	4.721%
8	10.832	1308	1318	1328	rBV	133141	292673	6.53%	1.529%
9	13.276	1726	1734	1749	rBV	1496723	2458811	54.87%	12.847%
10	14.651	1961	1968	1986	rBV	275512	478493	10.68%	2.500%
11	16.137	2214	2221	2243	rBV	1249664	2217987	49.49%	11.589%
12	17.395	2428	2435	2454	rBV	324360	601152	13.41%	3.141%
13	20.003	2873	2879	2894	rBV	3149254	4481454	100.00%	23.416%
14	21.660	3155	3161	3172	rBV2	362520	707996	15.80%	3.699%
15	23.934	3541	3548	3556	rBV4	20364	49000	1.09%	0.256%
16	24.862	3695	3706	3720	rBV	244103	778996	17.38%	4.070%

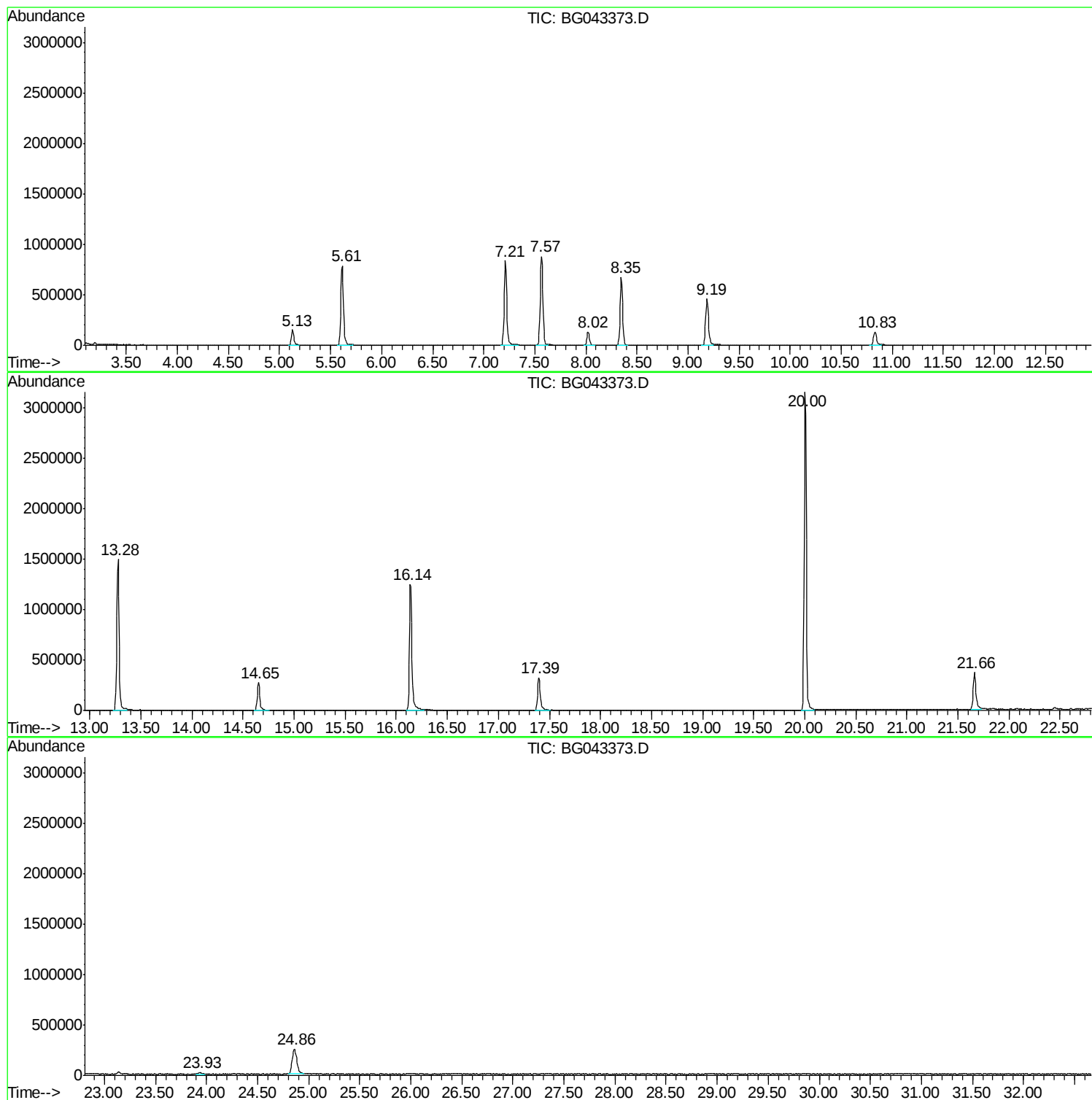
Sum of corrected areas: 19138713

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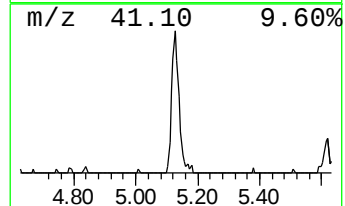
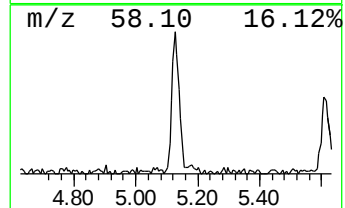
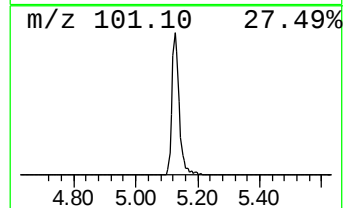
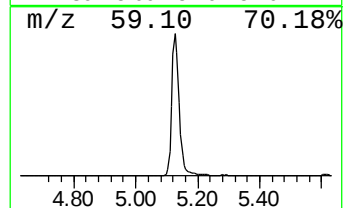
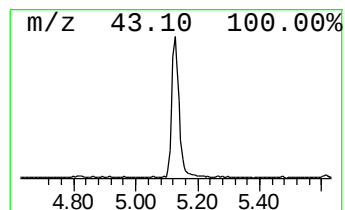
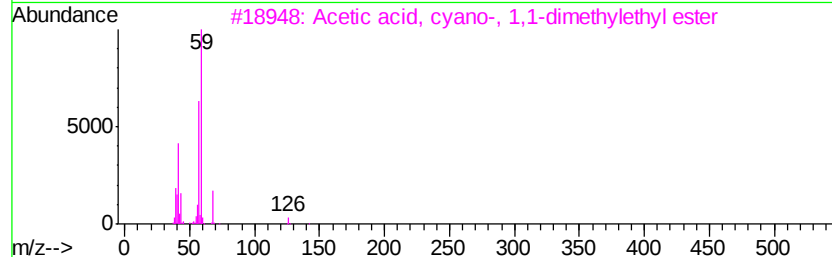
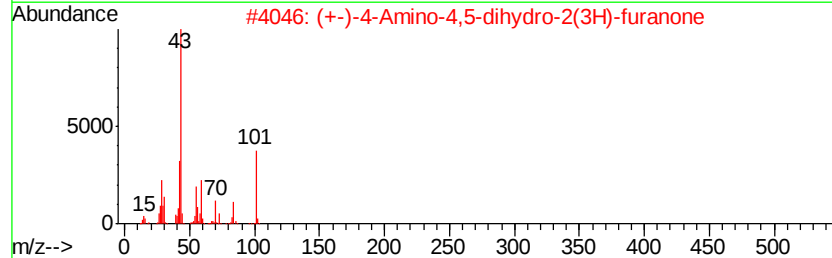
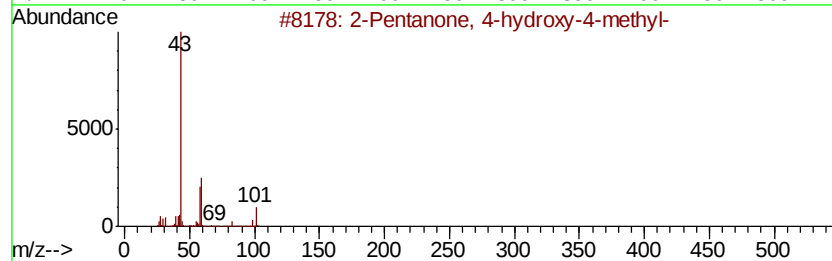
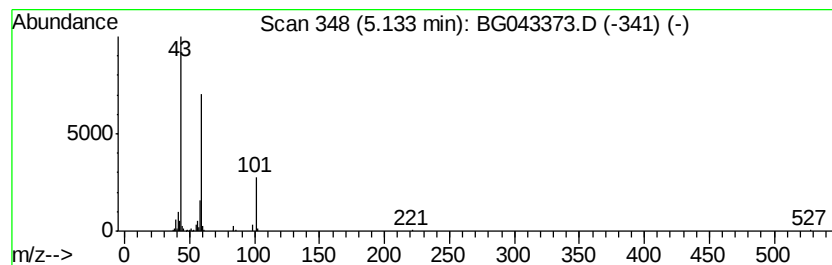
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	21.45 ng	251174	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	37
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Acetone	58	C3H6O	000067-64-1	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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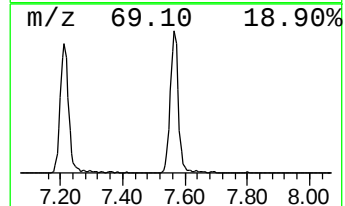
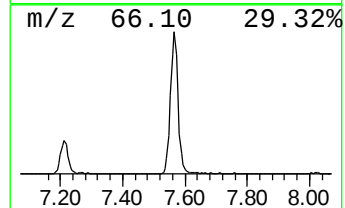
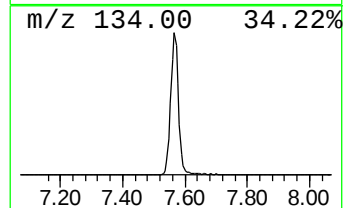
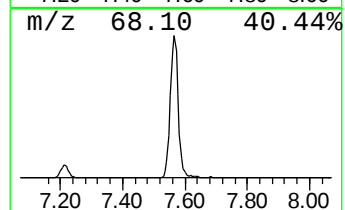
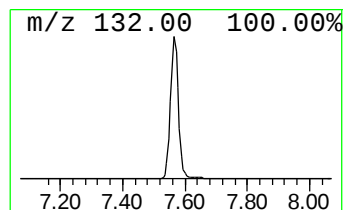
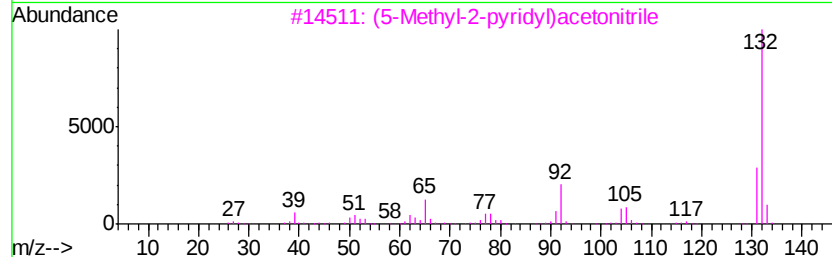
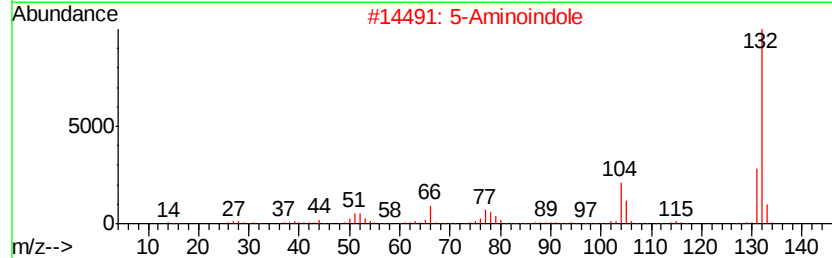
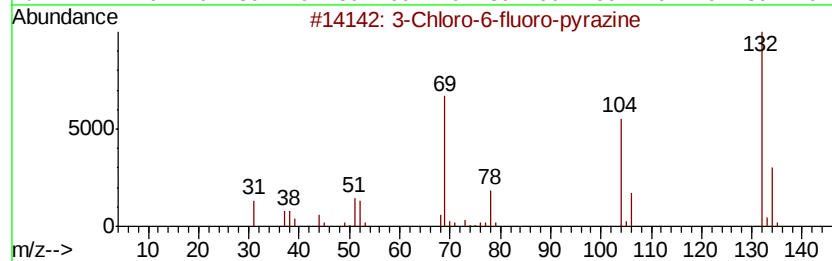
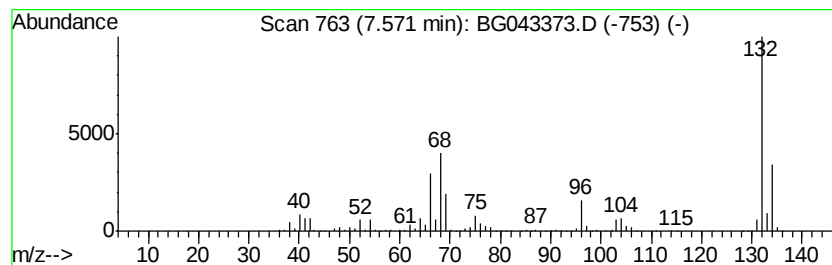
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	133.11 ng	1558510	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	21.4	ng	251174	1	8.02	234175	20.0
unknown7.57	7.57	133.1	ng	1558510	1	8.02	234175	20.0