

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091418\
 Data File : BG036736.D
 Acq On : 15 Sep 2018 11:42
 Operator : JU/SJ
 Sample : J4883-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 292

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	5.158	307	318	326	rBV2	130182	224697	5.68%	1.203%
2	5.622	390	397	406	rBV	788486	1244958	31.47%	6.663%
3	7.208	659	667	677	rBV	720022	1250250	31.61%	6.692%
4	7.578	721	730	739	rBV	817608	1406199	35.55%	7.526%
5	8.048	803	810	819	rBV	179744	305471	7.72%	1.635%
6	8.371	856	865	873	rBV	578322	990239	25.03%	5.300%
7	9.212	999	1008	1019	rVB	536754	940283	23.77%	5.033%
8	10.851	1280	1287	1294	rBV	241515	440115	11.13%	2.356%
9	13.295	1696	1703	1711	rBV	1370411	2162911	54.68%	11.576%
10	14.676	1929	1938	1945	rBV2	408191	670749	16.96%	3.590%
11	16.156	2184	2190	2212	rBV	1144044	1839823	46.51%	9.847%
12	17.414	2398	2404	2411	rBV2	596921	927950	23.46%	4.967%
13	20.022	2842	2848	2860	rBV	3013790	3955659	100.00%	21.172%
14	21.679	3123	3130	3138	rBV2	652896	1053324	26.63%	5.638%
15	23.172	3378	3384	3390	rVB3	44361	77446	1.96%	0.415%
16	23.965	3514	3519	3526	rVB2	41773	86428	2.18%	0.463%
17	24.887	3665	3676	3689	rVB	358228	1066776	26.97%	5.710%
18	26.033	3867	3871	3879	rVB7	17857	40389	1.02%	0.216%

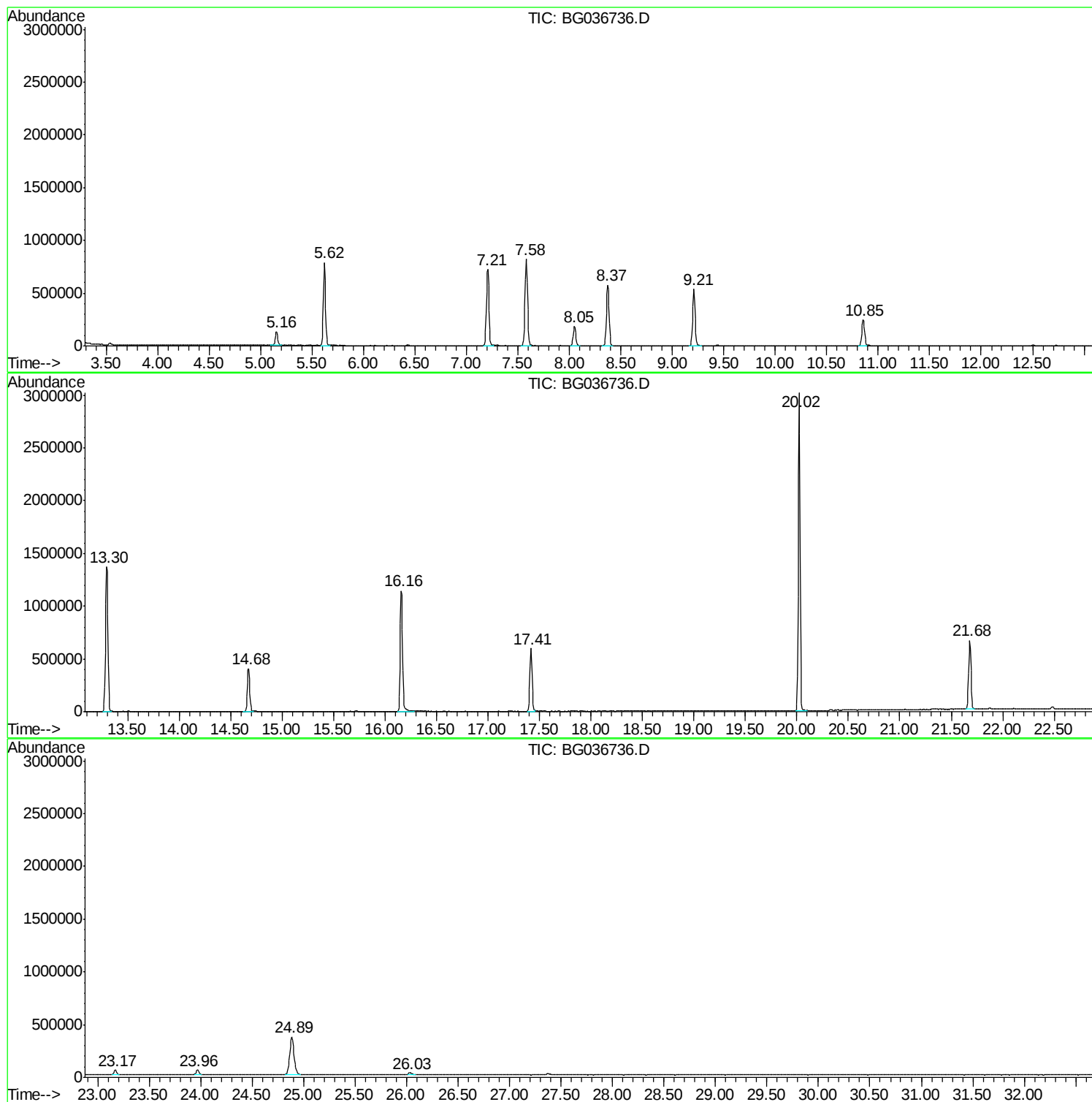
Sum of corrected areas: 18683667

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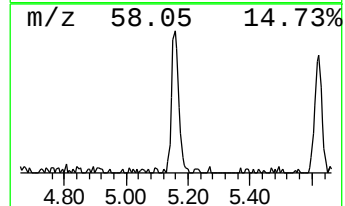
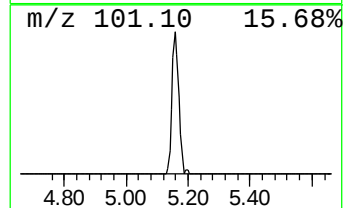
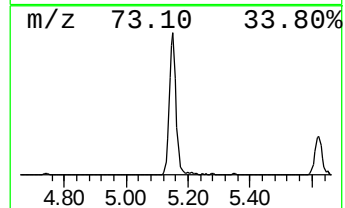
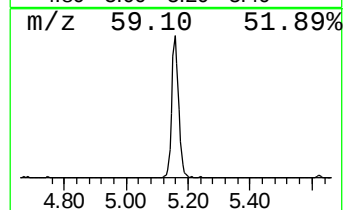
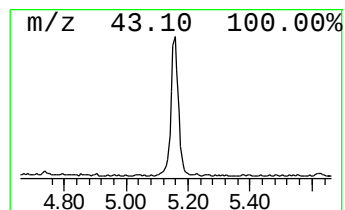
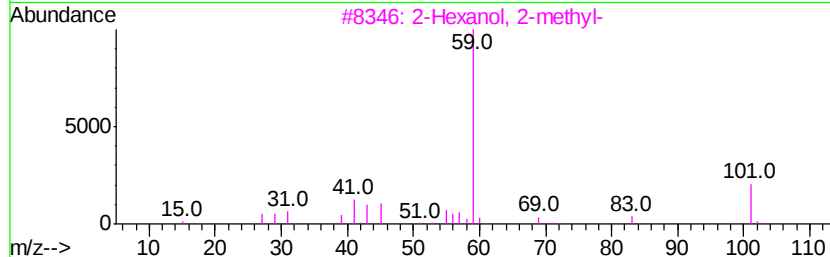
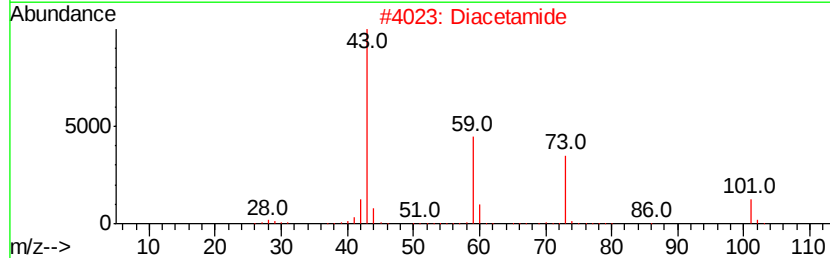
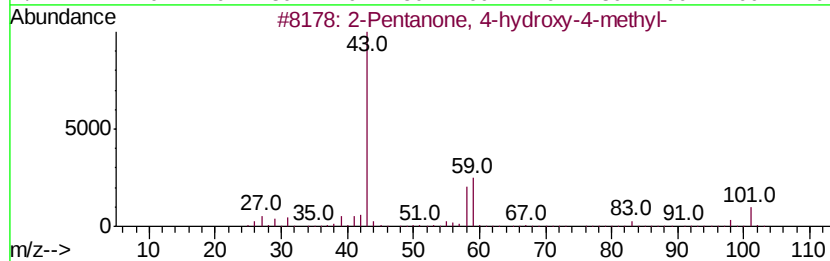
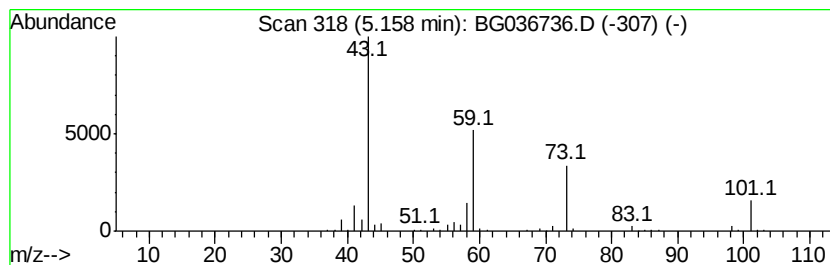
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	14.71 ng	224697	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Diacetamide	101	C4H7NO2	000625-77-4	45
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	43
4		3-Hexanol	102	C6H14O	000623-37-0	40
5		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	28



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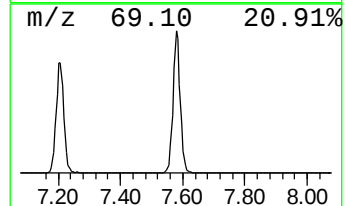
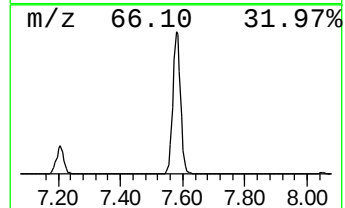
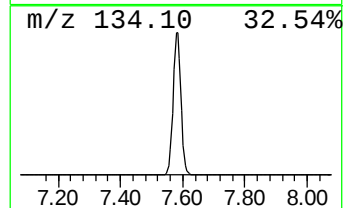
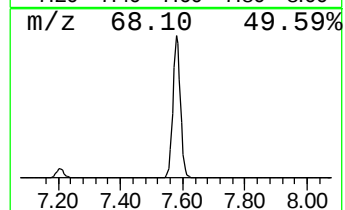
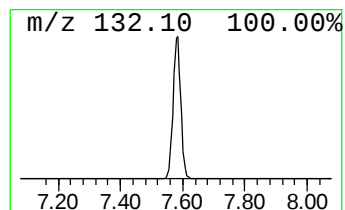
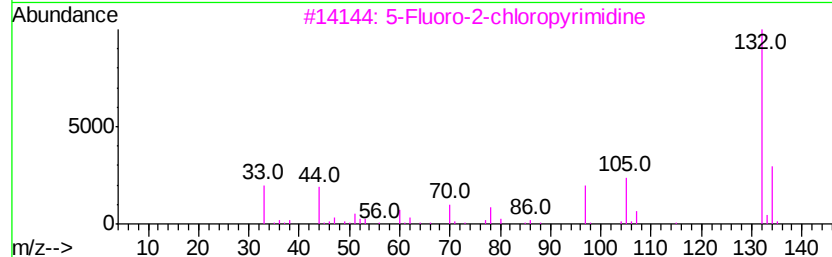
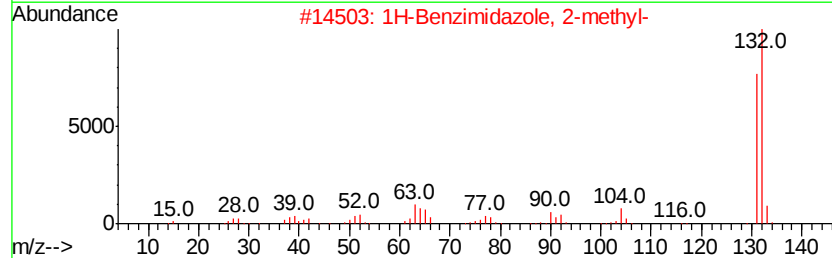
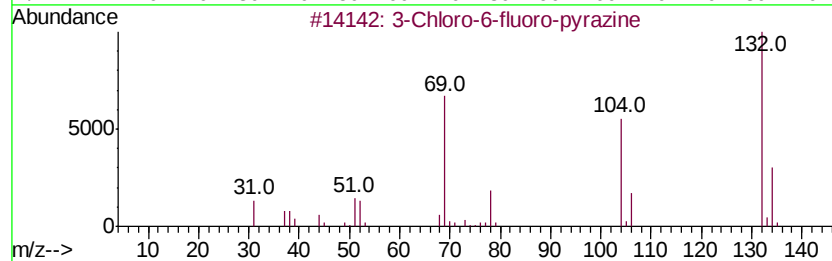
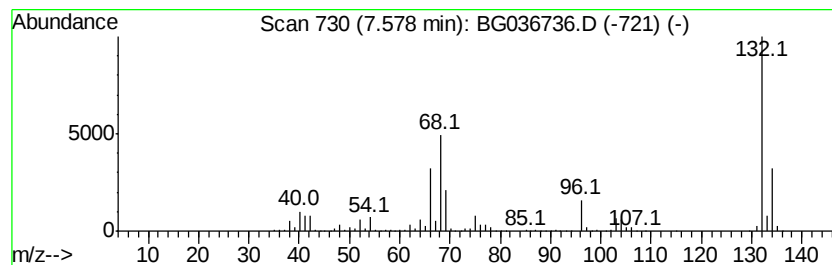
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	92.07 ng	1406200	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	25
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...	5.16	14.7	ng	224697	1	8.05	305471	20.0
unknown7.58	7.58	92.1	ng	1406200	1	8.05	305471	20.0