

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036646.D
 Acq On : 11 Sep 2018 4:46
 Operator : JU/SJ
 Sample : J4829-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OILY-STONES-PADS

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	311	318	327	rVB	112030	174613	4.27%	0.873%
2	5.622	390	397	410	rBV	891420	1403322	34.33%	7.018%
3	7.209	659	667	677	rBV	834049	1428805	34.95%	7.146%
4	7.585	724	731	741	rBV	942269	1599382	39.12%	7.999%
5	8.055	804	811	819	rVB	197547	328944	8.05%	1.645%
6	8.378	859	866	873	rBV	686384	1189001	29.09%	5.947%
7	9.212	1000	1008	1023	rVB	607678	1096914	26.83%	5.486%
8	10.858	1281	1288	1296	rBV	269415	486283	11.90%	2.432%
9	13.302	1696	1704	1714	rBV	1723199	2639102	64.56%	13.199%
10	14.677	1931	1938	1947	rBV2	458991	750871	18.37%	3.755%
11	16.163	2184	2191	2211	rBV	1216216	1953894	47.80%	9.772%
12	17.420	2399	2405	2413	rBV2	635328	976650	23.89%	4.885%
13	20.029	2843	2849	2859	rBV	2948517	4087970	100.00%	20.445%
14	21.686	3124	3131	3146	rBV2	598850	1056648	25.85%	5.285%
15	24.894	3668	3677	3694	rBV	236072	822436	20.12%	4.113%

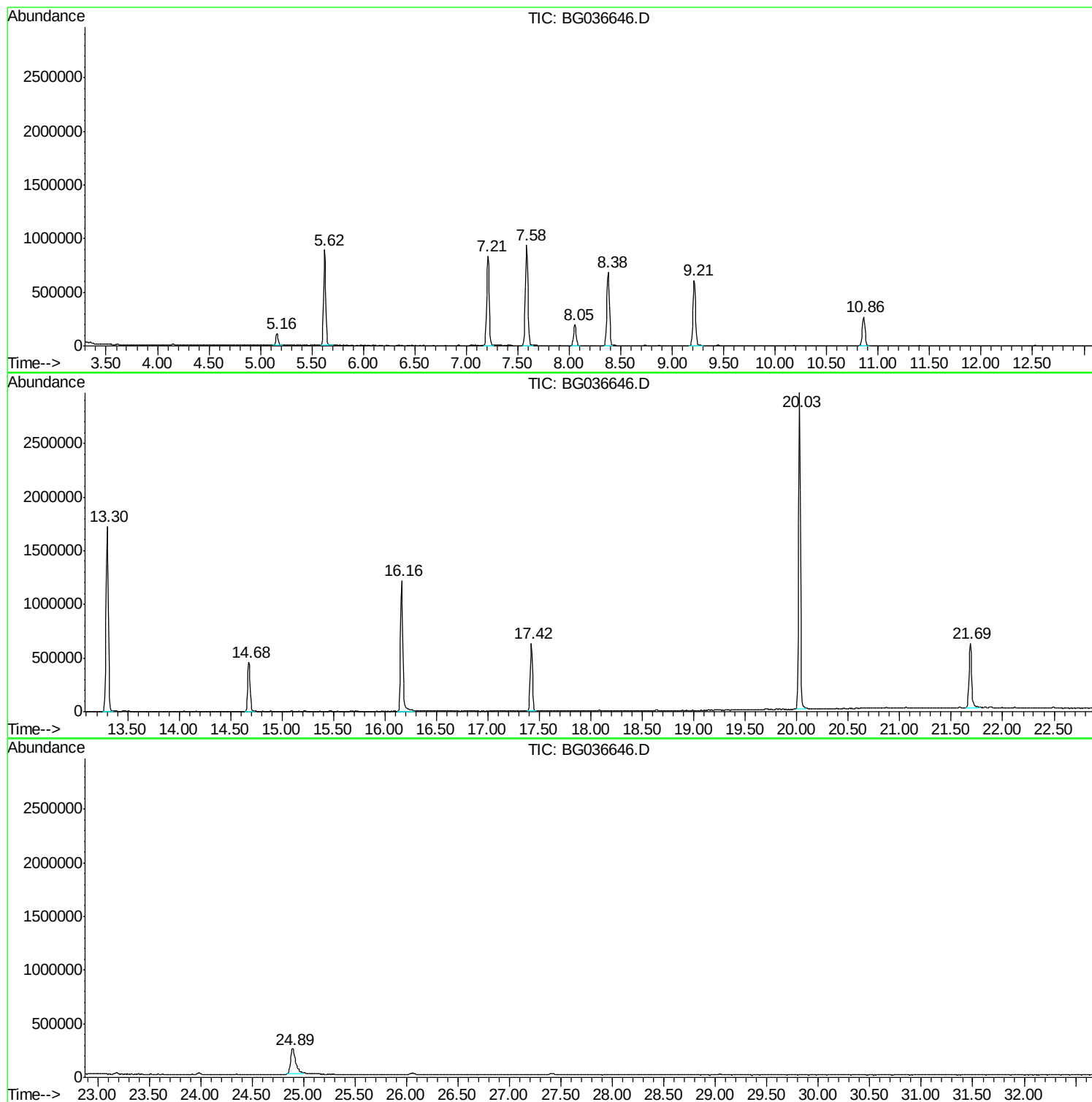
Sum of corrected areas: 19994835

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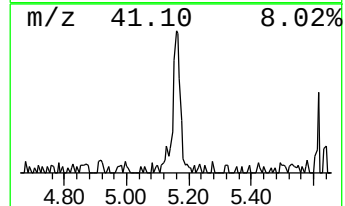
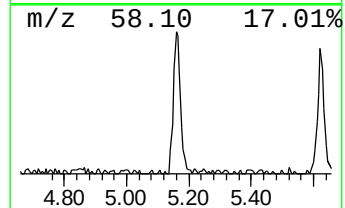
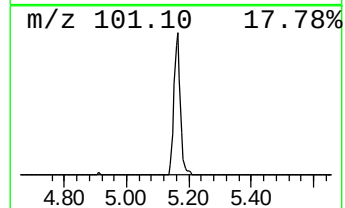
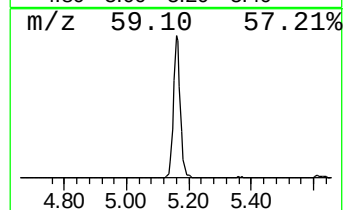
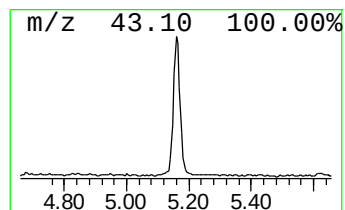
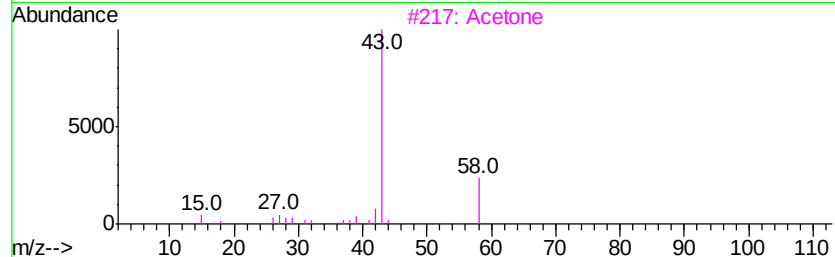
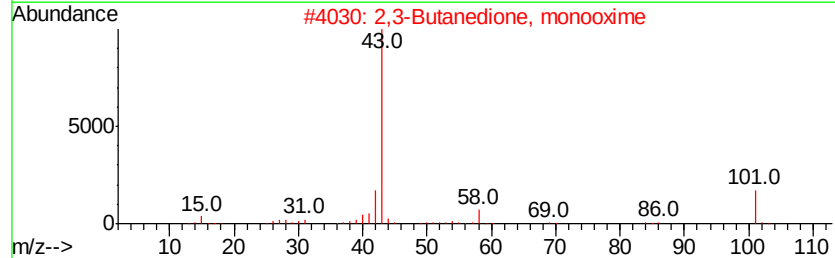
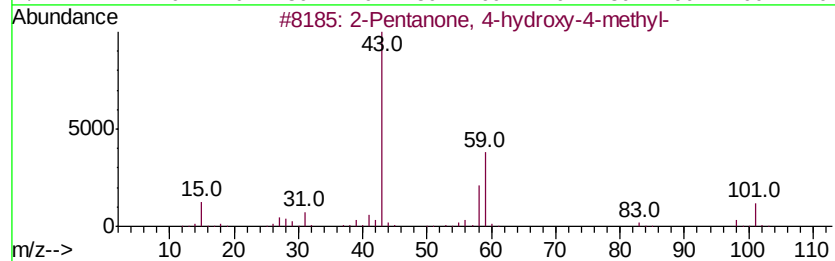
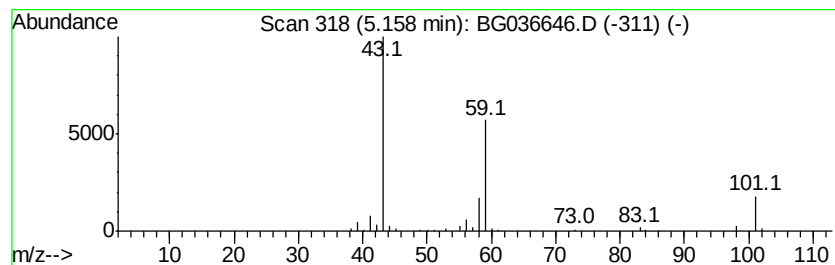
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	10.62 ng	174613	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	53
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
3			Acetone	58	C3H6O	000067-64-1	9
4			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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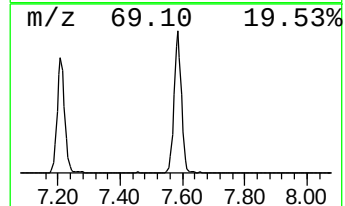
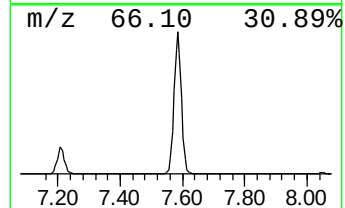
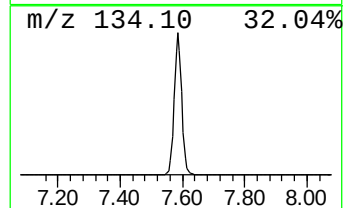
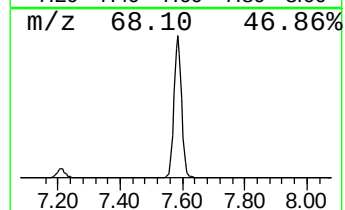
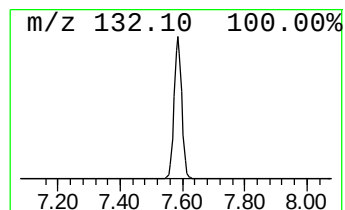
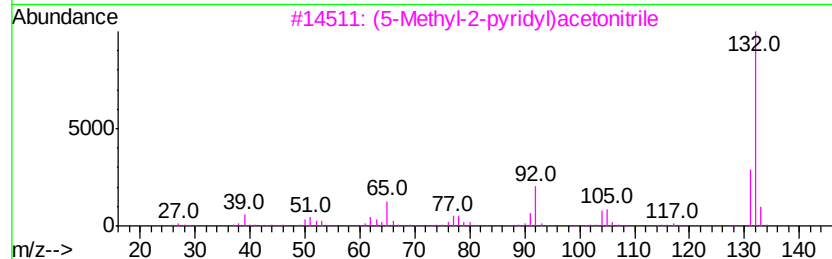
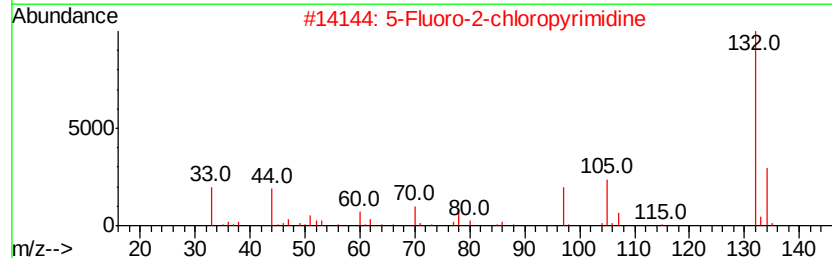
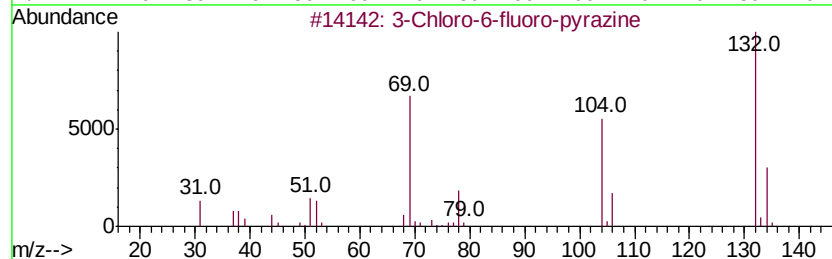
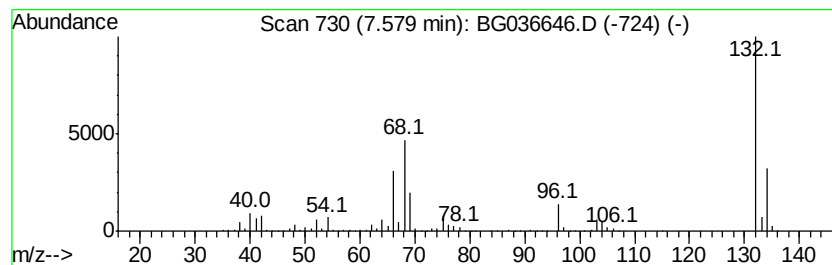
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	97.24 ng	1599380	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	10.6	ng	174613	1	8.05	328944	20.0
unknown7.58	7.58	97.2	ng	1599380	1	8.05	328944	20.0