

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG091018\
 Data File : BG036647.D
 Acq On : 11 Sep 2018 5:25
 Operator : JU/SJ
 Sample : J4829-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OILY-STONES-3-4FT

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	5.163	308	319	329	rVB	134771	208269	4.79%	1.026%
2	5.627	390	398	409	rBV	865484	1414669	32.55%	6.969%
3	7.207	659	667	676	rBV	803984	1385952	31.89%	6.828%
4	7.583	724	731	741	rBV	939930	1613972	37.14%	7.951%
5	8.053	803	811	819	rVB	200300	343407	7.90%	1.692%
6	8.376	857	866	880	rBV	711150	1214888	27.95%	5.985%
7	9.217	1000	1009	1019	rVB	605304	1106094	25.45%	5.449%
8	10.856	1281	1288	1294	rBV	276526	512373	11.79%	2.524%
9	10.909	1294	1297	1305	rVB2	35352	58172	1.34%	0.287%
10	13.300	1697	1704	1717	rBV	1686714	2650579	60.99%	13.058%
11	14.675	1931	1938	1946	rBV2	474001	779501	17.94%	3.840%
12	16.161	2184	2191	2210	rBV	1060690	1749784	40.26%	8.620%
13	17.419	2399	2405	2419	rBV2	662544	1010693	23.26%	4.979%
14	20.027	2843	2849	2865	rBV	3237517	4346105	100.00%	21.411%
15	21.684	3124	3131	3153	rBV2	568071	1109227	25.52%	5.464%
16	24.898	3668	3678	3702	rBV2	195002	795204	18.30%	3.917%

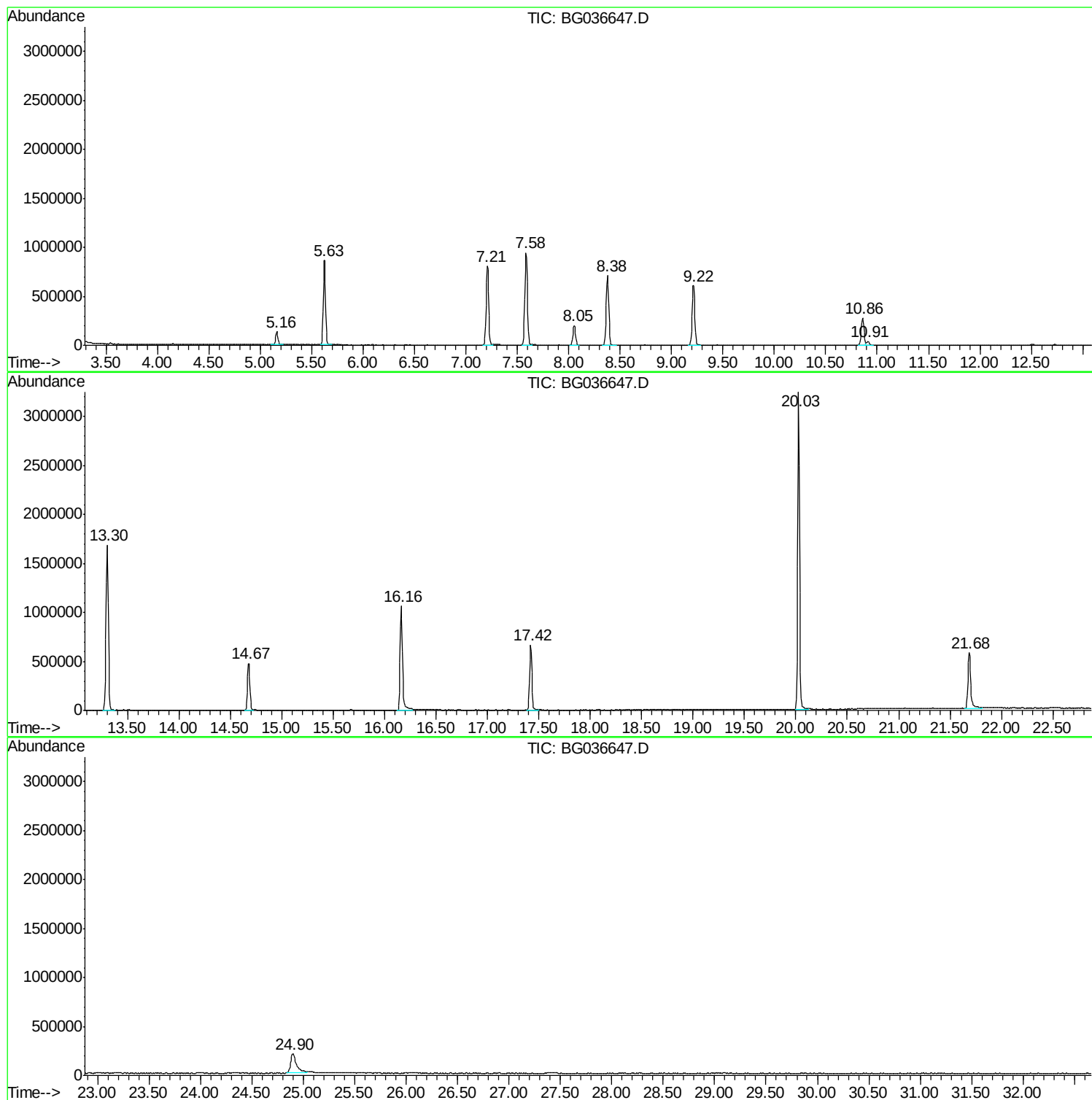
Sum of corrected areas: 20298889

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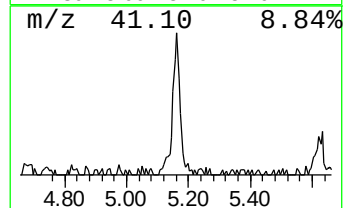
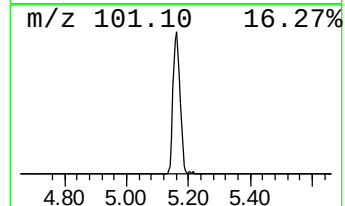
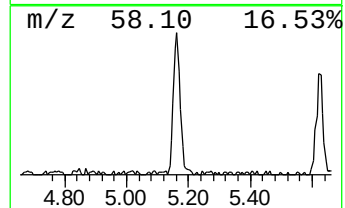
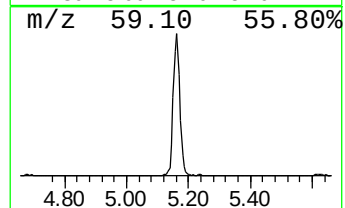
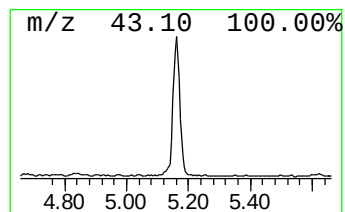
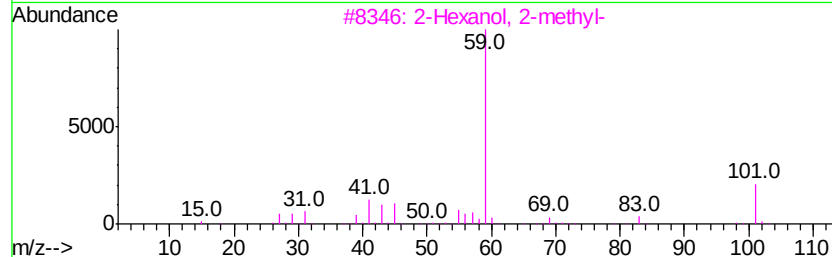
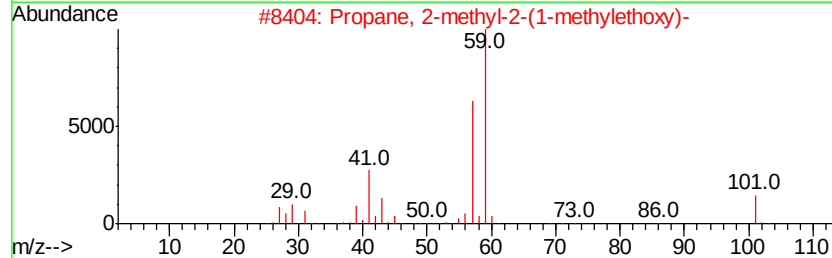
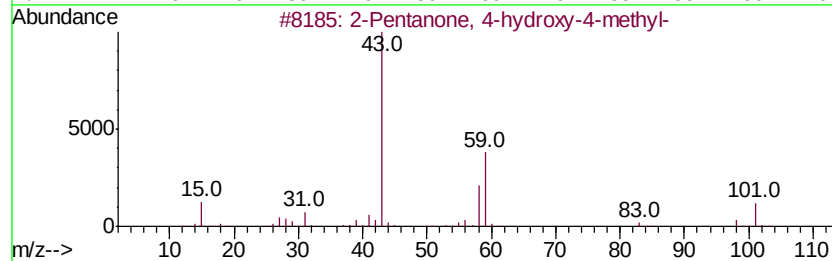
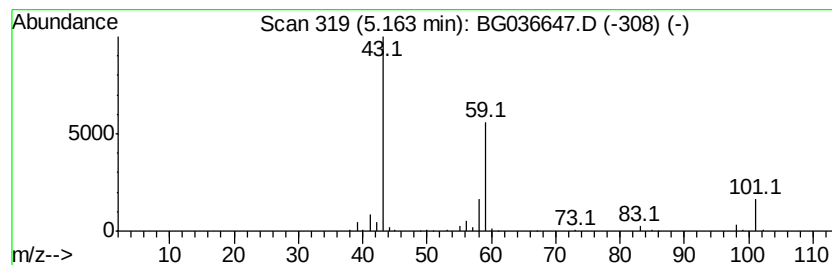
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	12.13 ng	208269	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	59
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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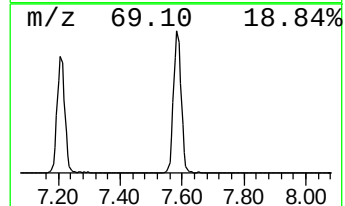
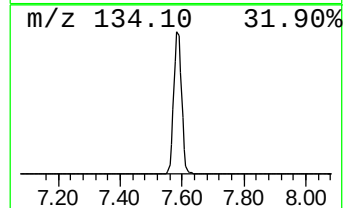
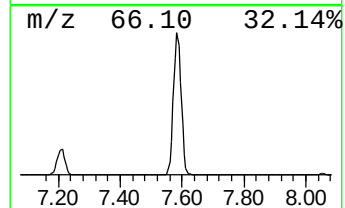
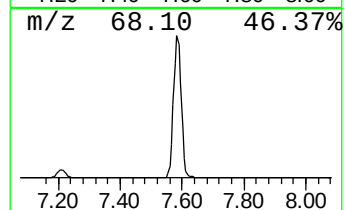
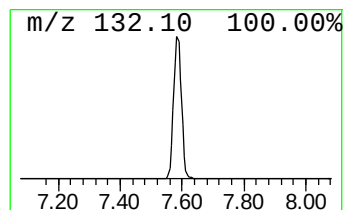
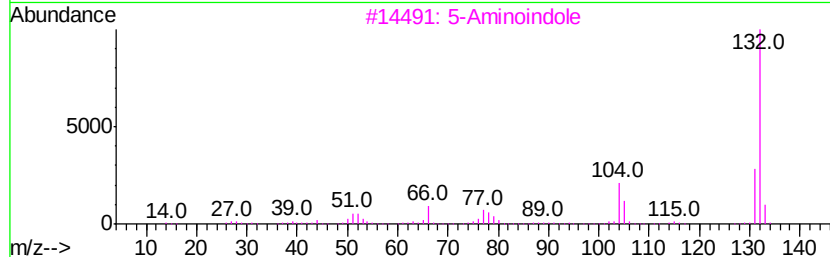
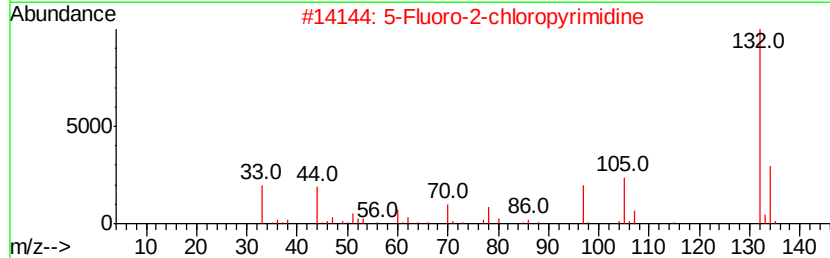
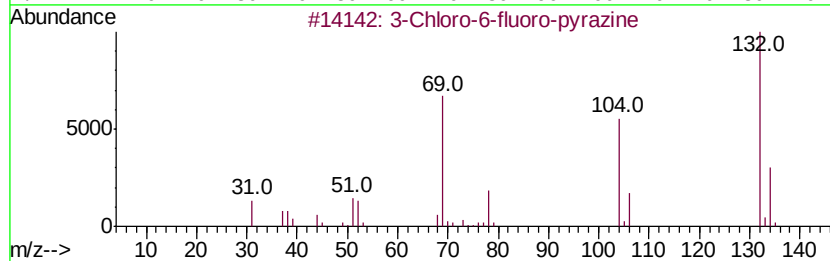
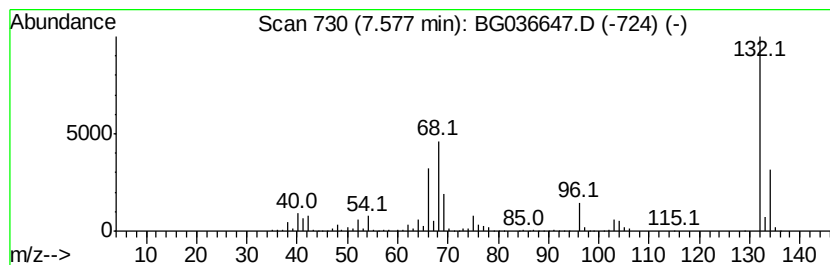
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	94.00 ng	1613970	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	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
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...	5.16	12.1	ng	208269	1	8.05	343407	20.0
unknown7.58	7.58	94.0	ng	1613970	1	8.05	343407	20.0