

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092518\
 Data File : BG036958.D
 Acq On : 25 Sep 2018 20:54
 Operator : JU/SJ
 Sample : J4774-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-03-092418

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-BG092018.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	309	317	329	rVB	84128	142463	4.12%	0.861%
2	5.615	390	396	414	rBV	654296	1156931	33.49%	6.996%
3	7.201	659	666	689	rBV	555175	1113301	32.22%	6.732%
4	7.577	722	730	745	rBV	668989	1291575	37.38%	7.810%
5	8.041	803	809	825	rBV	149474	285443	8.26%	1.726%
6	8.364	856	864	880	rBV	555467	1040985	30.13%	6.295%
7	9.205	999	1007	1024	rBV	426496	888682	25.72%	5.374%
8	10.844	1277	1286	1297	rBV	190128	380727	11.02%	2.302%
9	13.288	1695	1702	1719	rBV	1117385	1924157	55.69%	11.635%
10	14.669	1930	1937	1947	rBV2	322857	551013	15.95%	3.332%
11	16.155	2183	2190	2208	rBV	1032389	1712544	49.57%	10.356%
12	17.413	2397	2404	2425	rBV	490274	779237	22.55%	4.712%
13	20.021	2841	2848	2863	rBV	2538477	3455052	100.00%	20.892%
14	21.678	3123	3130	3141	rBV2	525249	921228	26.66%	5.571%
15	24.886	3665	3676	3691	rBV2	298830	894166	25.88%	5.407%

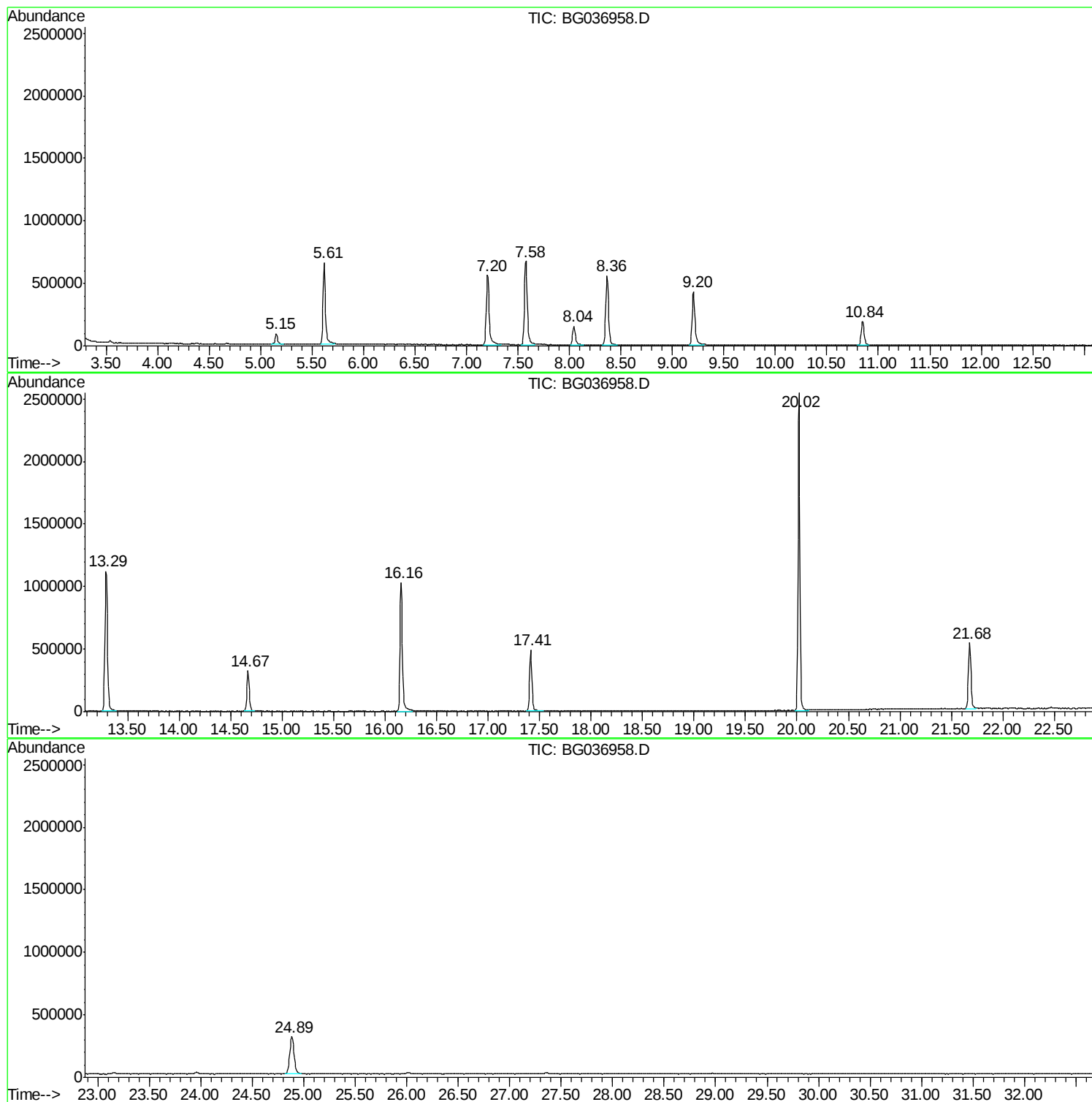
Sum of corrected areas: 16537504

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.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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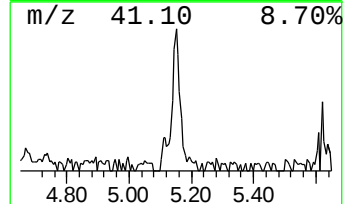
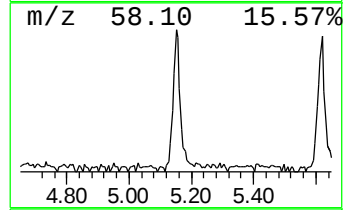
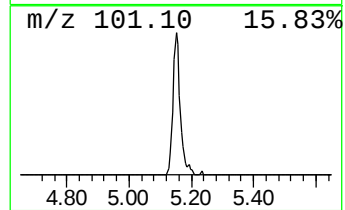
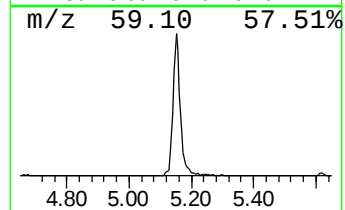
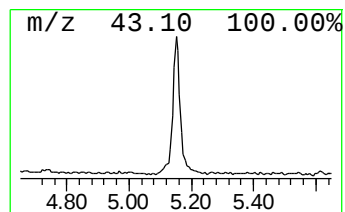
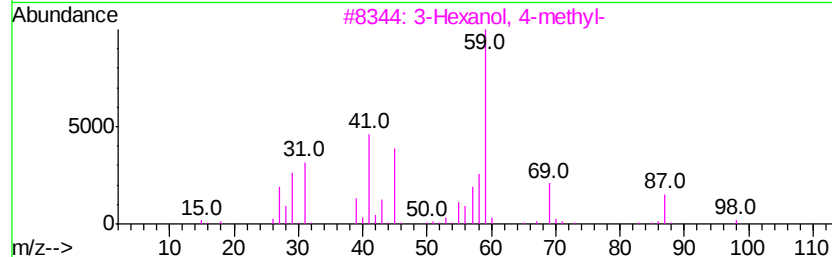
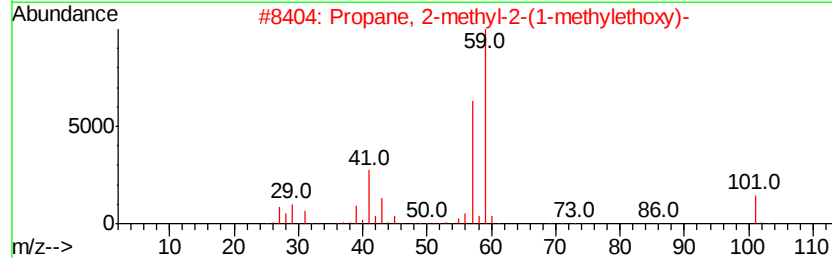
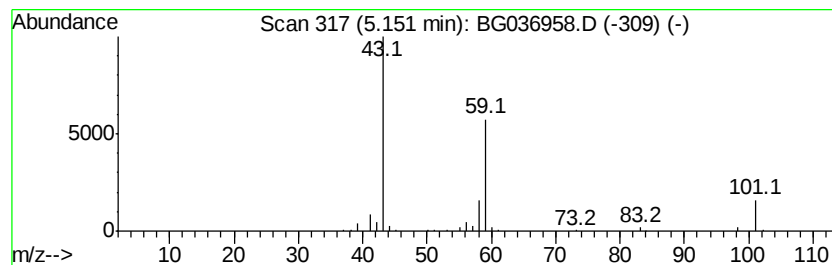
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	9.98 ng	142463	1,4-Dichlorobenzene-d4	8.04

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	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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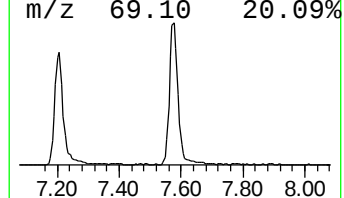
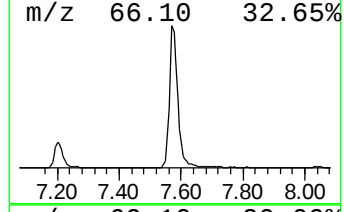
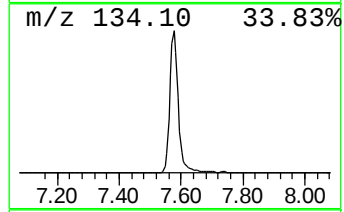
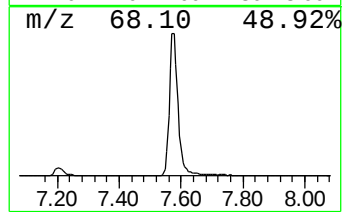
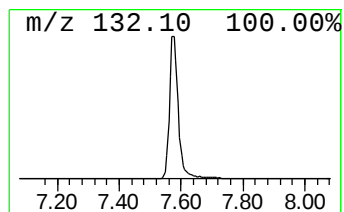
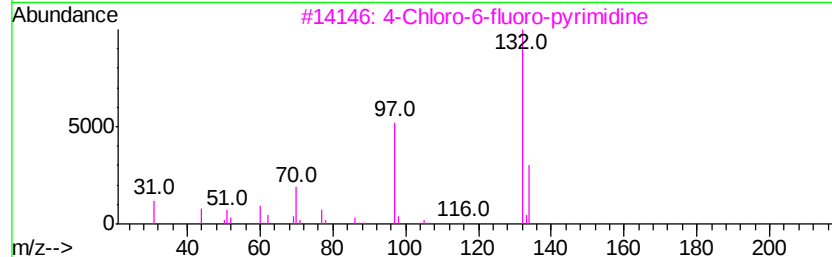
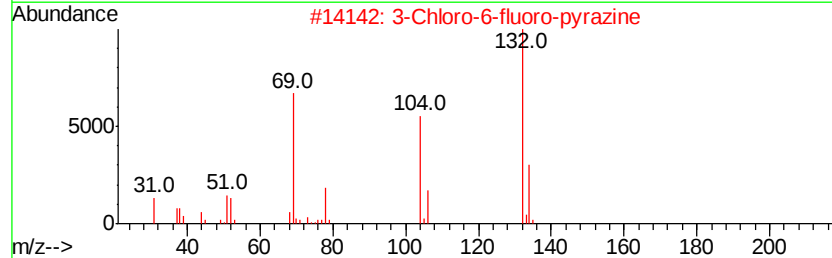
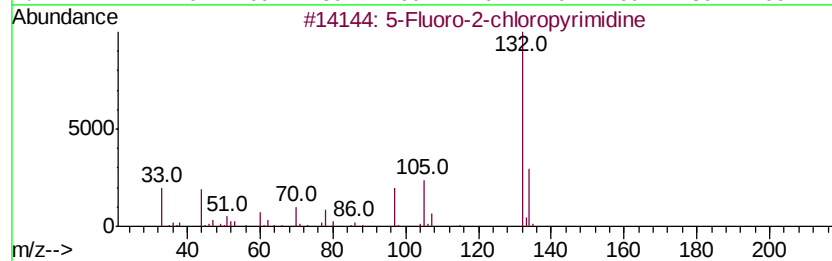
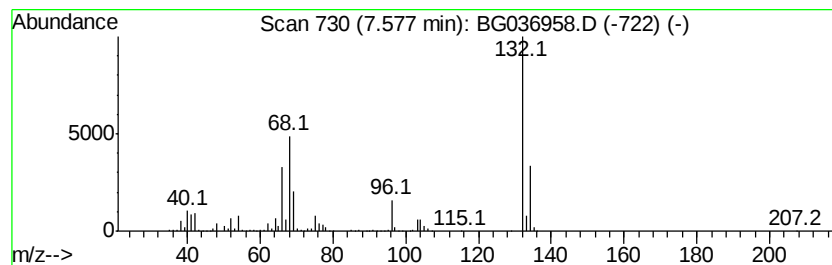
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	90.50 ng	1291580	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	10.0	ng	142463	1	8.04	285443	20.0
unknown7.58	7.58	90.5	ng	1291580	1	8.04	285443	20.0