

Data Path : Z:\HPCHEM1\BNA G\DATA\BG091916\
 Data File : BG024027.D
 Acq On : 19 Sep 2016 20:48
 Operator : UM/SJ
 Sample : H4819-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-MW11-04

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.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	5.123	385	389	395	rBV3	55792	89418	1.62%	0.441%
2	5.605	466	471	485	rVB	399858	677815	12.29%	3.343%
3	7.226	742	747	758	rBV2	244610	473225	8.58%	2.334%
4	7.585	802	808	819	rBV	743975	1344129	24.37%	6.629%
5	8.054	881	888	895	rBV	155752	261572	4.74%	1.290%
6	8.378	937	943	957	rVB	693300	1218523	22.10%	6.009%
7	9.235	1083	1089	1108	rBV	630235	1186622	21.52%	5.852%
8	10.892	1363	1371	1384	rBV	188438	371005	6.73%	1.830%
9	13.341	1780	1788	1801	rBV	1792140	2769285	50.22%	13.657%
10	14.181	1926	1931	1939	rBV	87259	147955	2.68%	0.730%
11	14.728	2017	2024	2034	rBV	356930	597449	10.83%	2.946%
12	16.232	2272	2280	2294	rBV	1734309	2817922	51.10%	13.897%
13	17.495	2489	2495	2506	rBV2	500815	842127	15.27%	4.153%
14	20.103	2933	2939	2948	rBV	4423349	5514580	100.00%	27.196%
15	21.801	3222	3228	3239	rVB	572269	1023774	18.56%	5.049%
16	25.149	3789	3798	3811	rVB3	290987	942066	17.08%	4.646%

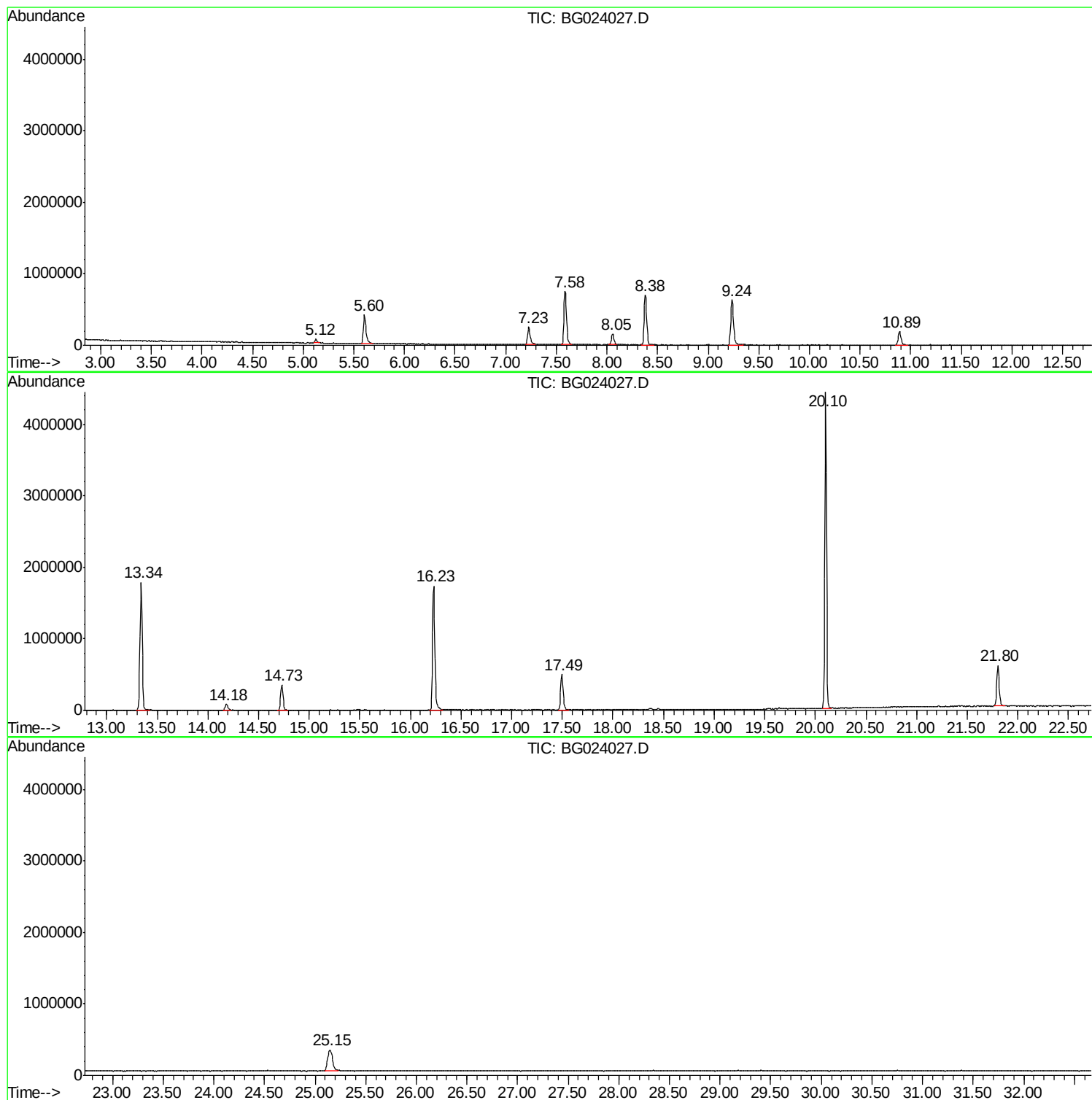
Sum of corrected areas: 20277467

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.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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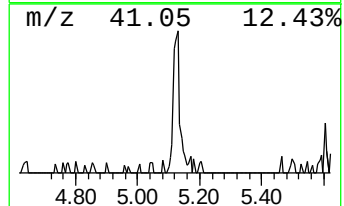
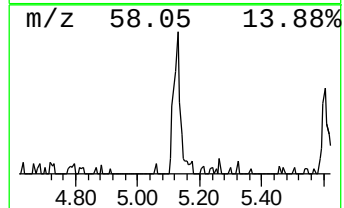
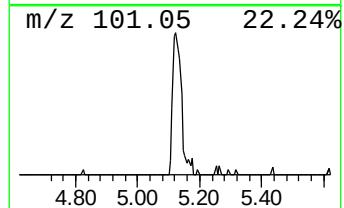
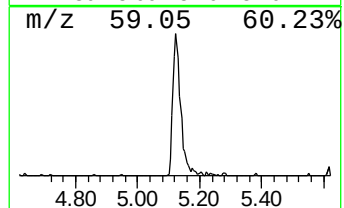
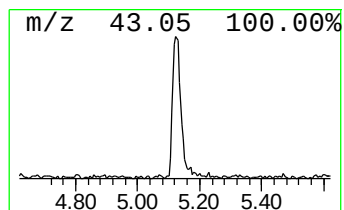
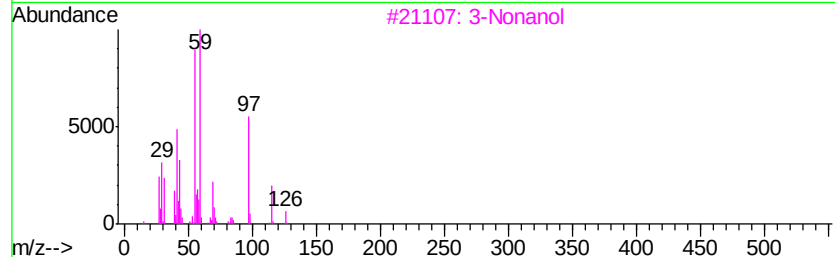
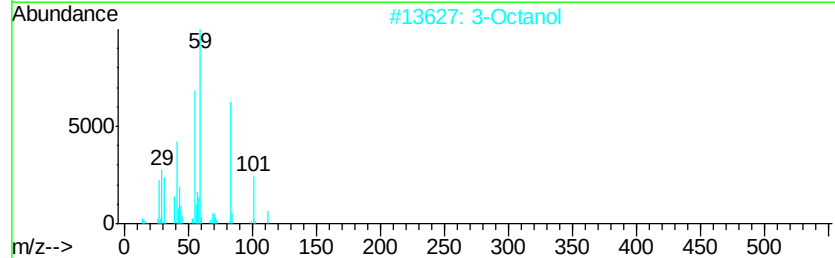
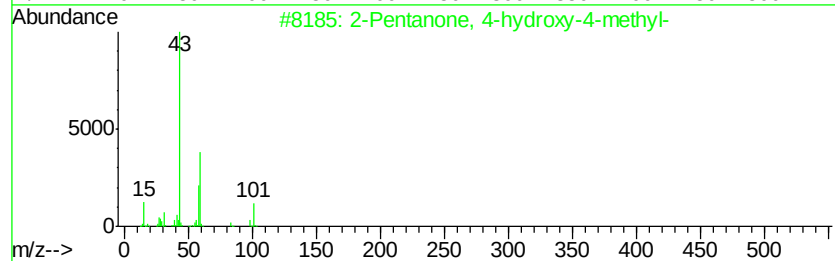
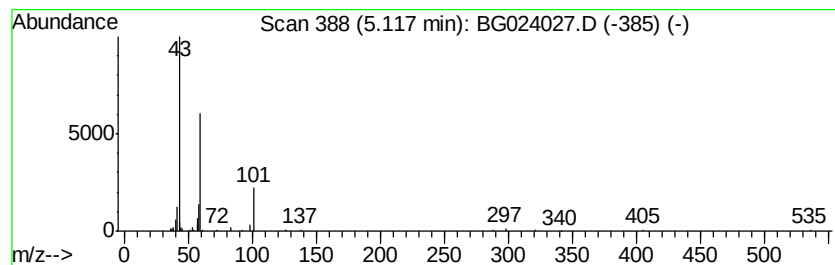
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	6.84 ng	89418	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	45
2		3-Octanol	130	C8H18O	000589-98-0	33
3		3-Nonanol	144	C9H20O	000624-51-1	23
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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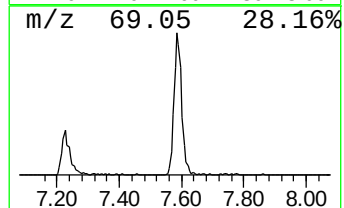
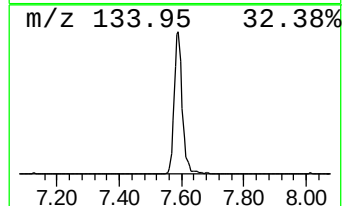
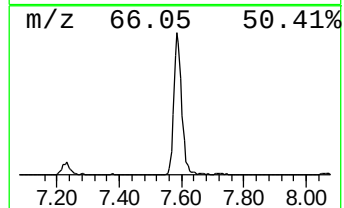
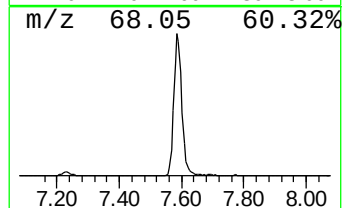
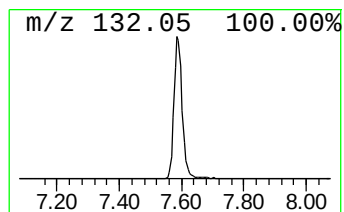
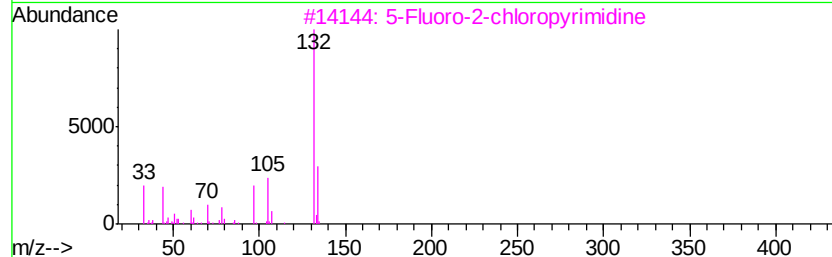
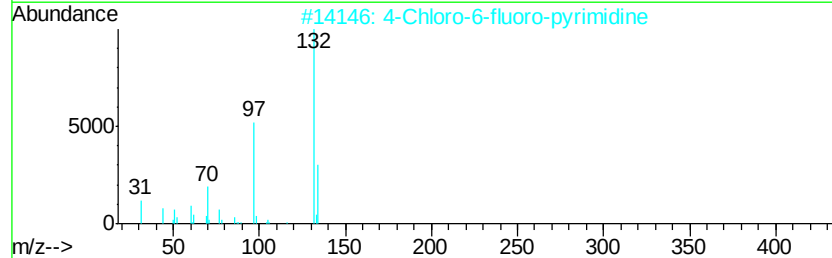
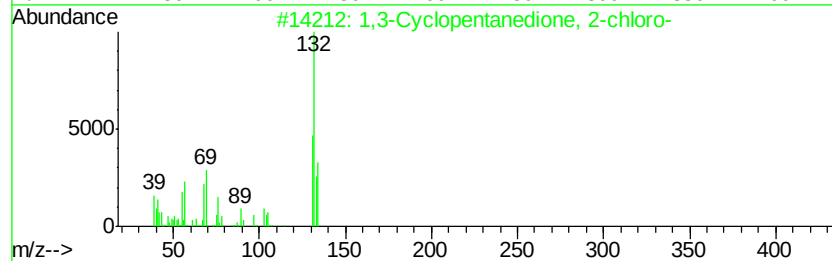
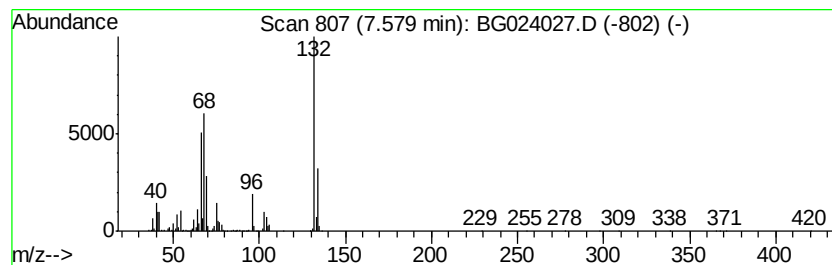
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	102.77 ng	1344130	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	6.8	ng	89418	1	8.05	261572	20.0
unknown7.58	7.58	102.8	ng	1344130	1	8.05	261572	20.0