

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062317\
 Data File : BG027636.D
 Acq On : 23 Jun 2017 18:32
 Operator : SJ/JU
 Sample : PB100074BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100074BL

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:\HPCHEM1\BNA_G\METHODS\8270-BG062117.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	5.642	292	298	306	rBV	42652	66885	3.10%	0.769%
2	6.088	366	374	388	rBV	374340	631335	29.23%	7.259%
3	7.645	629	639	651	rBV	356463	656076	30.38%	7.544%
4	8.039	697	706	714	rBV	358373	649161	30.06%	7.464%
5	8.503	779	785	792	rVB2	61164	108363	5.02%	1.246%
6	8.832	833	841	849	rBB	251181	458240	21.22%	5.269%
7	9.666	975	983	996	rBV	204125	383335	17.75%	4.408%
8	11.323	1258	1265	1275	rVB	83996	158434	7.34%	1.822%
9	13.715	1664	1672	1682	rBV	590384	964140	44.64%	11.086%
10	15.089	1899	1906	1915	rBB2	137167	230207	10.66%	2.647%
11	16.570	2151	2158	2168	rBV	630503	986662	45.69%	11.345%
12	17.833	2366	2373	2381	rBV2	227074	363069	16.81%	4.175%
13	20.413	2806	2812	2818	rBV	1629828	2159634	100.00%	24.833%
14	22.205	3108	3117	3125	rBV	246177	456421	21.13%	5.248%
15	25.842	3723	3736	3754	rBV3	122836	424762	19.67%	4.884%

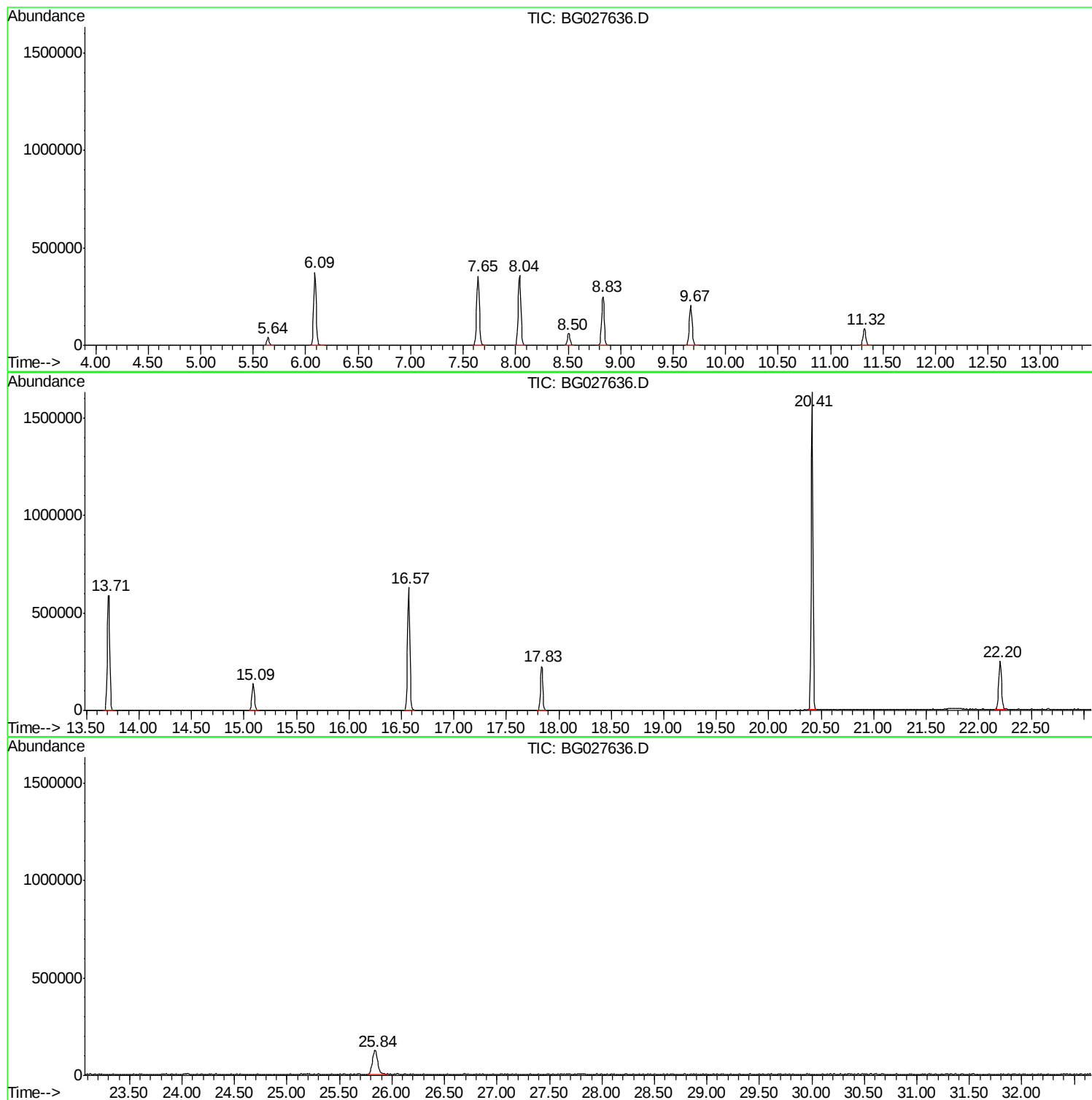
Sum of corrected areas: 8696724

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062117.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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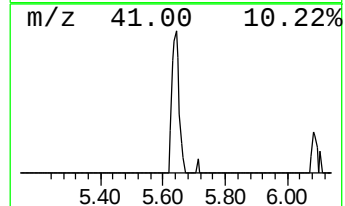
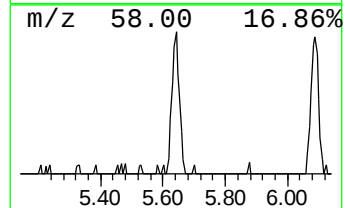
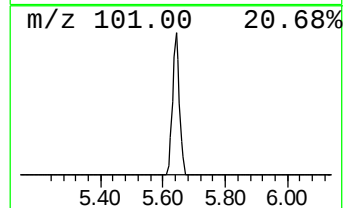
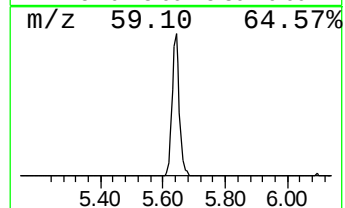
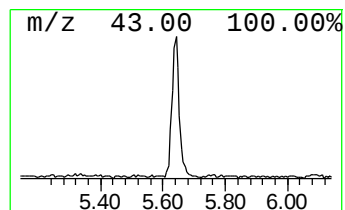
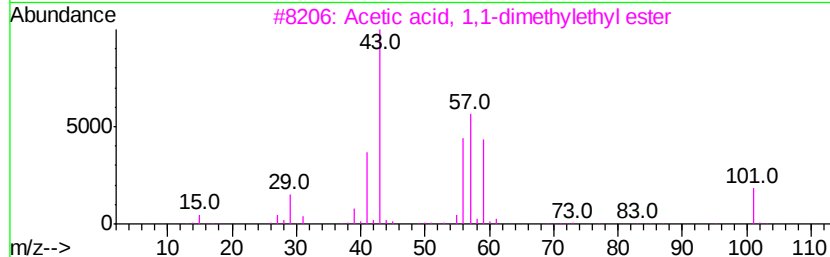
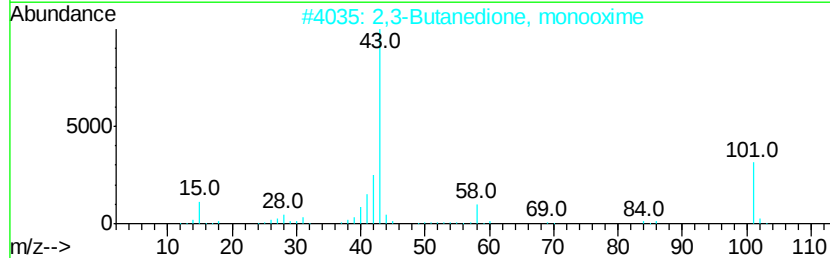
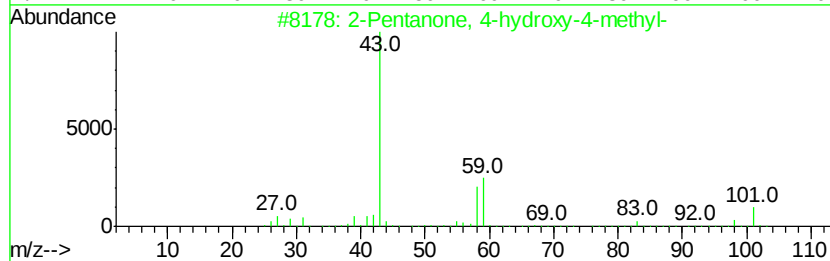
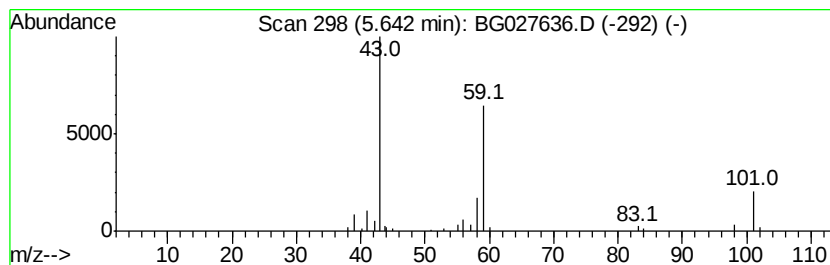
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	12.34 ng	66885	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	43
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25



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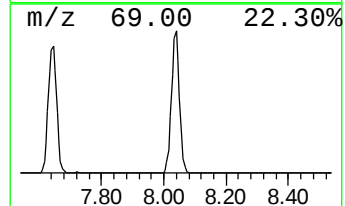
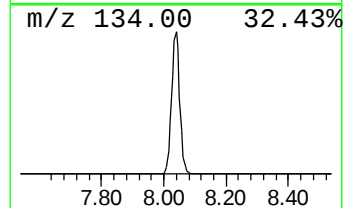
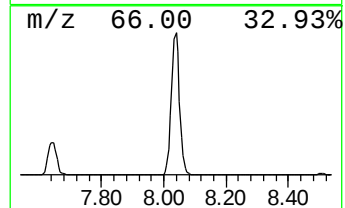
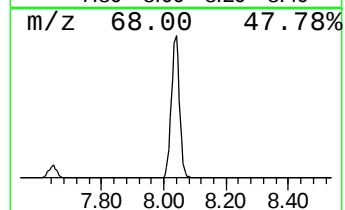
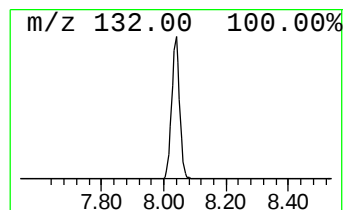
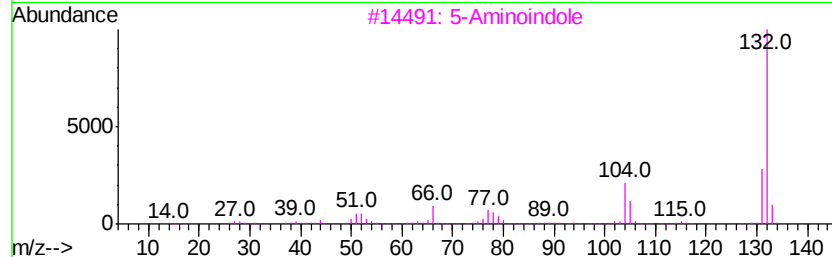
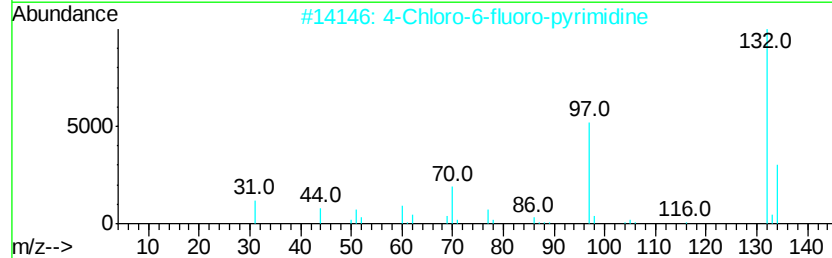
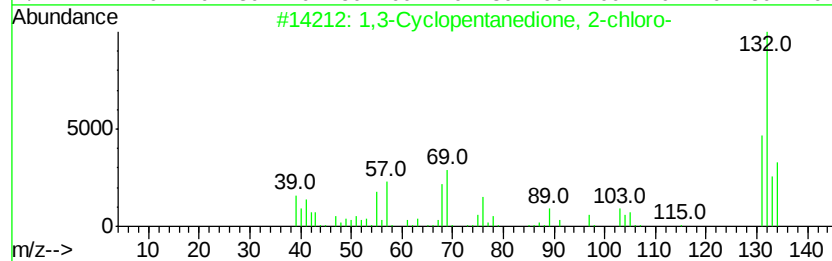
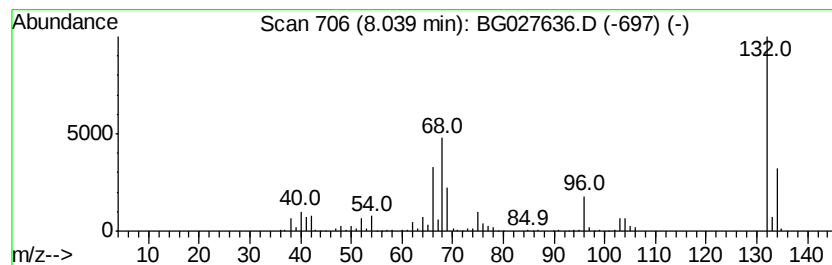
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Peak Number 2 unknown8.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	119.81 ng	649161	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.64	12.3	ng	66885	1	8.50	108363	20.0
unknown8.04	8.04	119.8	ng	649161	1	8.50	108363	20.0