

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024093.D
 Acq On : 23 Sep 2016 18:47
 Operator : UM/SJ
 Sample : H4803-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WP

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-BG092316.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	2.902	7	11	31	rVB	4848058	7404051	100.00%	11.943%
2	5.105	381	386	395	rBV	166344	274496	3.71%	0.443%
3	5.587	462	468	481	rBV	879352	1480759	20.00%	2.388%
4	7.214	738	745	746	rBV	1083107	1566536	21.16%	2.527%
5	7.237	747	749	762	rVB	1362366	1988283	26.85%	3.207%
6	7.572	799	806	809	rBV	971158	1784290	24.10%	2.878%
7	7.607	809	812	823	rVB	1183434	1845748	24.93%	2.977%
8	8.042	879	886	898	rVB	164998	292651	3.95%	0.472%
9	8.365	933	941	949	rBV	651811	1136103	15.34%	1.833%
10	8.500	957	964	981	rBV	1356261	2349557	31.73%	3.790%
11	8.829	1011	1020	1032	rBV	1618765	2886550	38.99%	4.656%
12	9.223	1080	1087	1098	rBV	647825	1151365	15.55%	1.857%
13	9.987	1209	1217	1221	rBV	1003780	1826779	24.67%	2.947%
14	10.045	1221	1227	1239	rVB	1655414	2884877	38.96%	4.653%
15	10.280	1246	1267	1275	rBV2	380205	1708045	23.07%	2.755%
16	10.533	1303	1310	1332	rVB2	438383	1025323	13.85%	1.654%
17	10.874	1361	1368	1377	rBV	241365	437169	5.90%	0.705%
18	12.178	1583	1590	1608	rBV	901037	1648827	22.27%	2.660%
19	13.153	1749	1756	1764	rBV	2024177	3300799	44.58%	5.324%
20	13.235	1764	1770	1780	rVV	814304	1686888	22.78%	2.721%
21	13.329	1780	1786	1794	rVB	1735259	2643460	35.70%	4.264%
22	14.716	2015	2022	2032	rBV2	443153	710861	9.60%	1.147%
23	14.856	2037	2046	2060	rBV	833946	1608254	21.72%	2.594%
24	14.974	2060	2066	2092	rVB	1014512	2295214	31.00%	3.702%
25	15.873	2210	2219	2232	rBV	1618764	2601301	35.13%	4.196%
26	16.219	2271	2278	2288	rBV	1721022	2591705	35.00%	4.180%
27	17.142	2428	2435	2453	rBV	1477617	2636055	35.60%	4.252%
28	17.482	2486	2493	2507	rVB	661958	972546	13.14%	1.569%
29	20.096	2931	2938	2953	rBV	3878733	5088678	68.73%	8.208%
30	21.794	3220	3227	3237	rBV	600737	1146389	15.48%	1.849%
31	25.131	3787	3795	3808	rVB2	330405	1023212	13.82%	1.650%

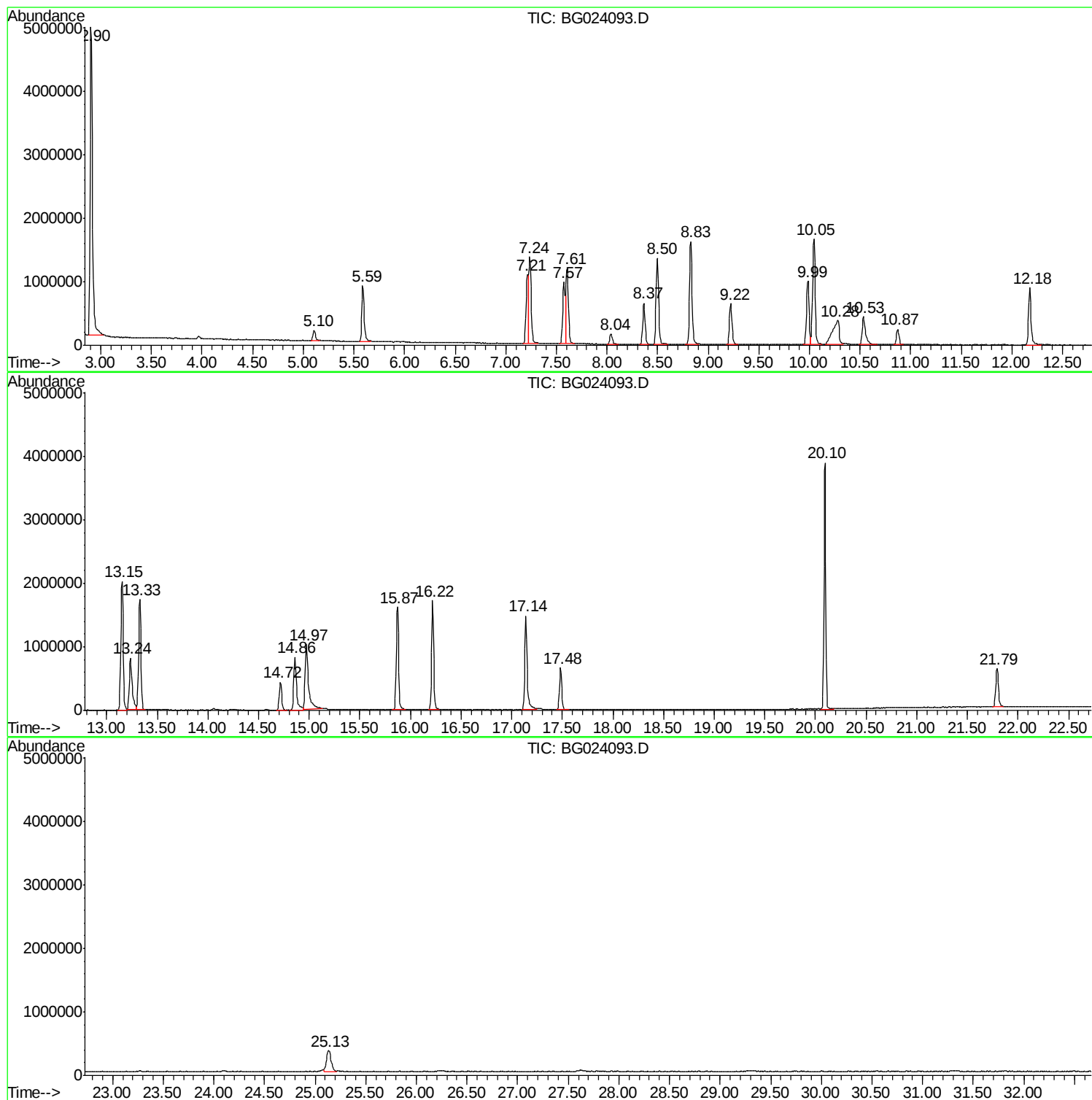
Sum of corrected areas: 61996771

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.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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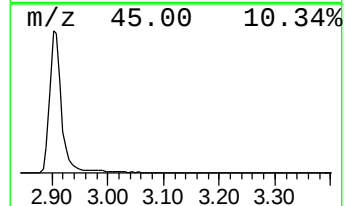
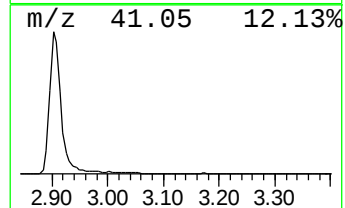
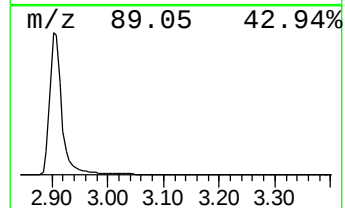
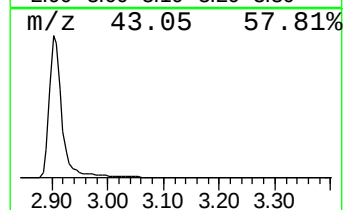
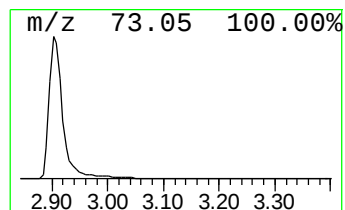
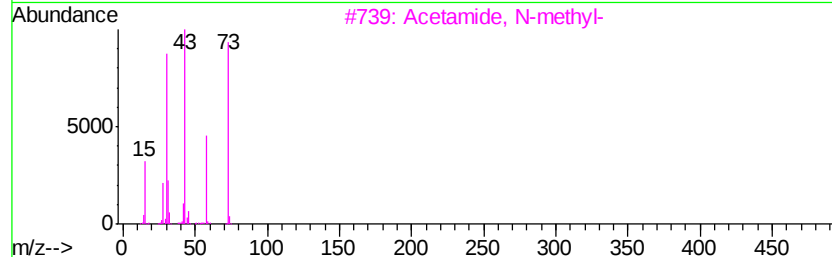
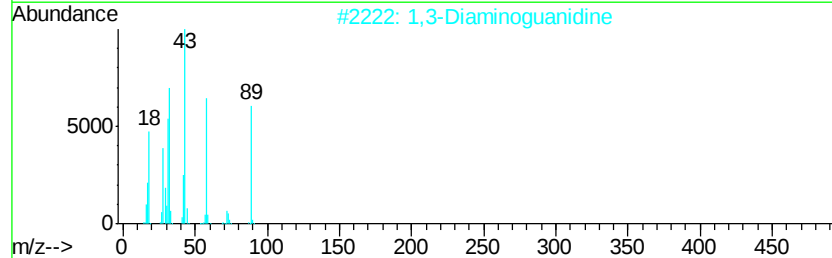
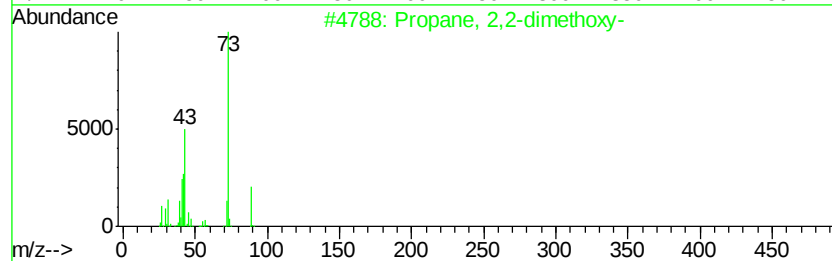
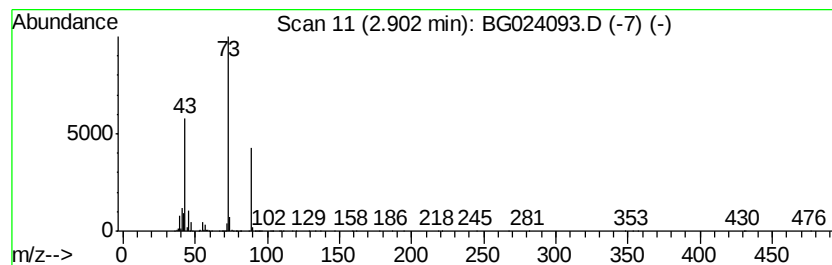
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.90 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.90	506.00 ng	7404050	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	40
2		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	23
3		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	9



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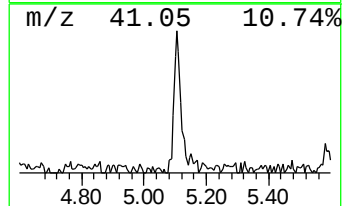
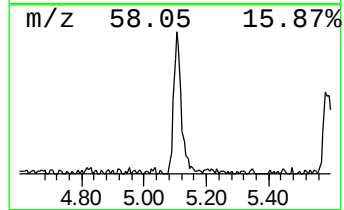
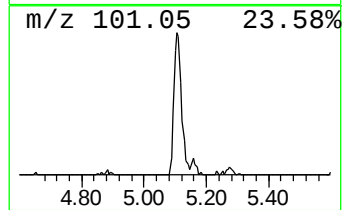
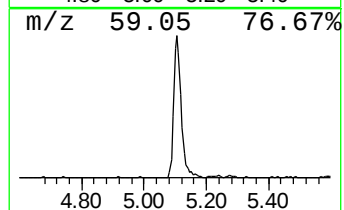
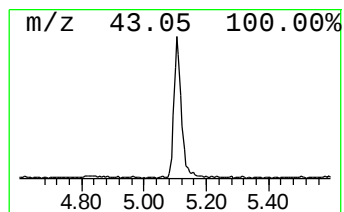
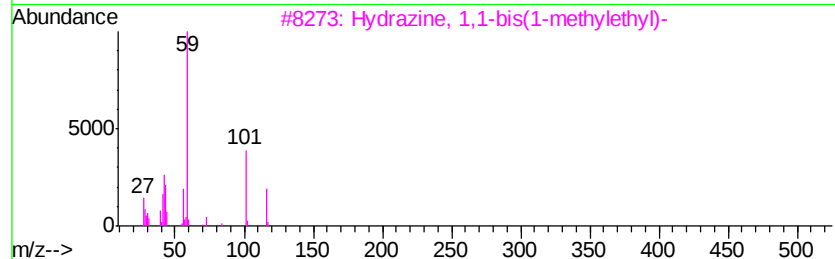
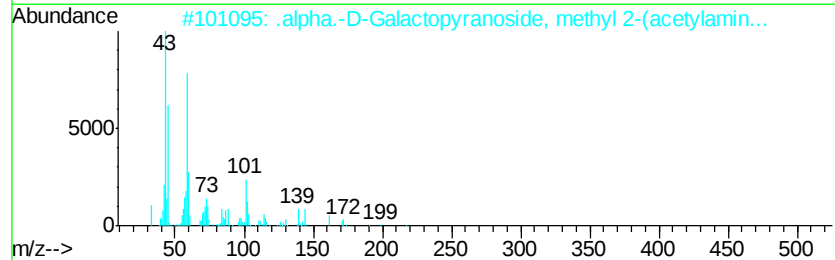
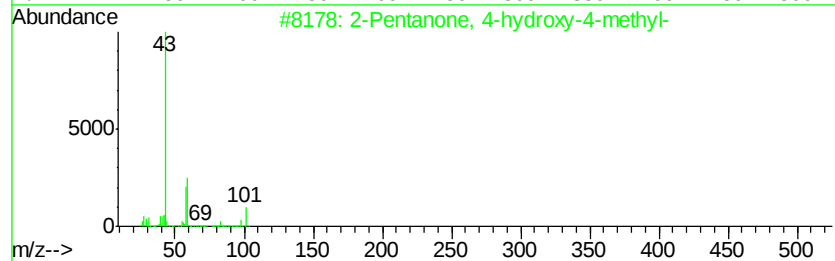
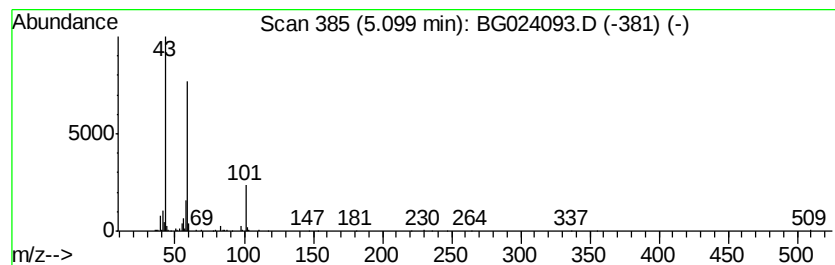
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	18.76 ng	274496	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		.alpha.-D-Galactopyranoside, met...	249	C10H19NO6	017296-07-0	38
3		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	12
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown2.90	2.90	506.0	ng	7404050	1	8.04	292651	20.0
2-Pentanone, 4-hy...	5.10	18.8	ng	274496	1	8.04	292651	20.0