

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024119.D
 Acq On : 24 Sep 2016 11:23
 Operator : UM/SJ
 Sample : H4902-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-1-092116

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	5.099	382	385	394	rVB2	77383	139430	2.52%	0.614%
2	5.593	463	469	479	rBV	449799	789121	14.26%	3.473%
3	7.214	740	745	757	rBV	308175	587882	10.63%	2.587%
4	7.572	799	806	815	rBV	917852	1627022	29.41%	7.161%
5	8.037	879	885	892	rBV	171966	312044	5.64%	1.373%
6	8.366	933	941	952	rVB	781484	1401123	25.33%	6.167%
7	9.223	1079	1087	1106	rBV	746225	1439357	26.02%	6.335%
8	10.874	1360	1368	1380	rBV	232715	453971	8.21%	1.998%
9	13.324	1779	1785	1795	rBV	2003442	3230286	58.39%	14.217%
10	14.164	1922	1928	1934	rVB	60031	91885	1.66%	0.404%
11	14.716	2016	2022	2029	rVB	445110	694568	12.56%	3.057%
12	16.220	2270	2278	2290	rBV	2002773	3086351	55.79%	13.583%
13	17.483	2487	2493	2503	rVB	612452	957768	17.31%	4.215%
14	18.352	2637	2641	2646	rBV5	45105	67994	1.23%	0.299%
15	19.627	2854	2858	2865	rBV7	45184	74582	1.35%	0.328%
16	20.091	2931	2937	2944	rBV	4367042	5531940	100.00%	24.347%
17	21.789	3219	3226	3234	rBV	674944	1133857	20.50%	4.990%
18	25.119	3784	3793	3809	rVB2	359217	1102163	19.92%	4.851%

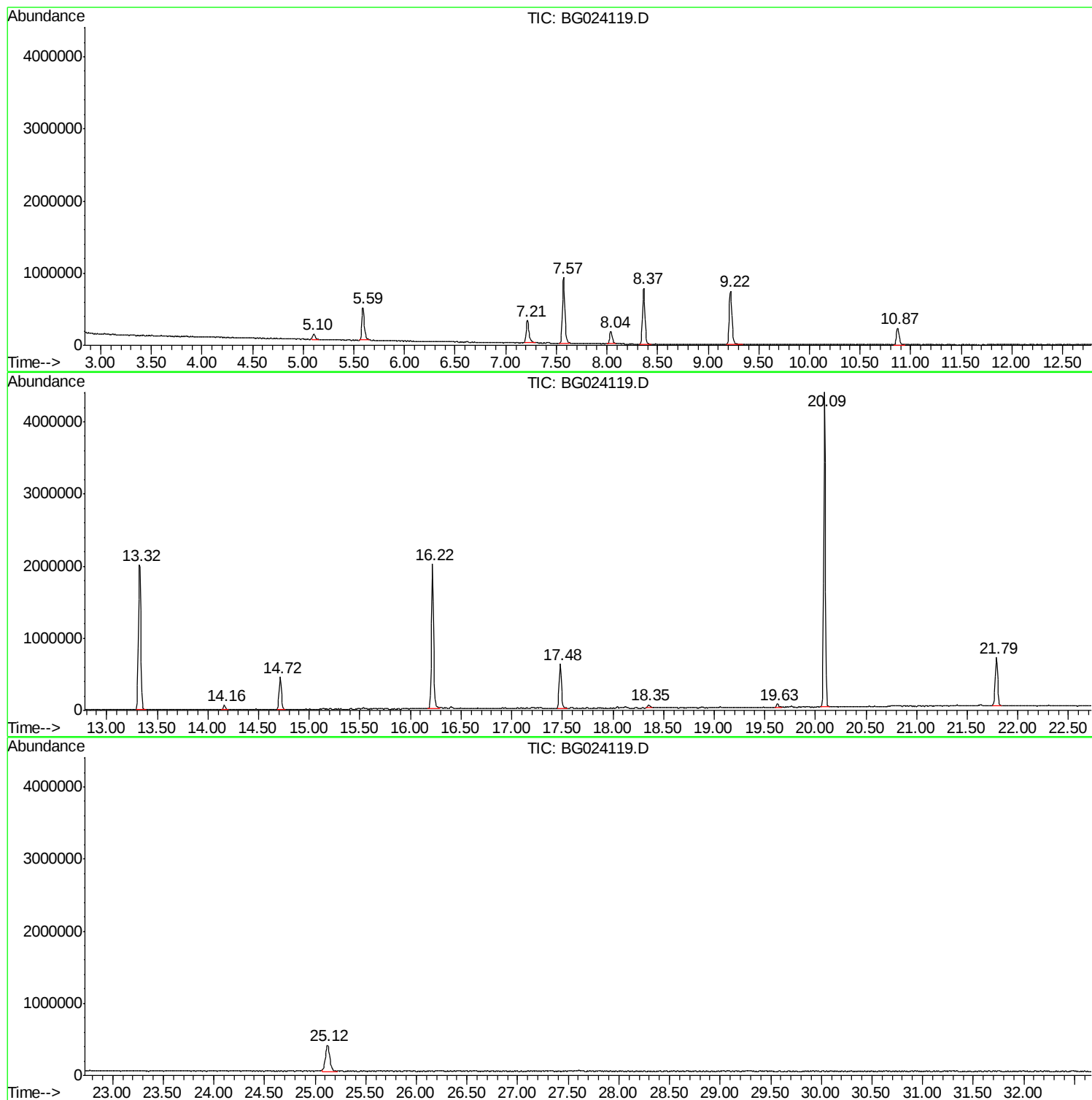
Sum of corrected areas: 22721344

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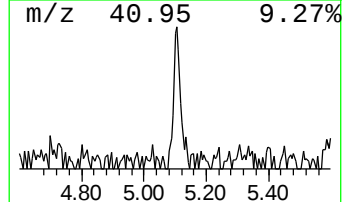
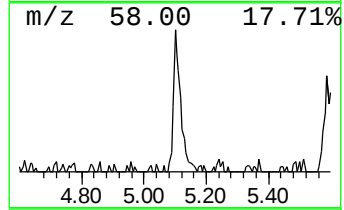
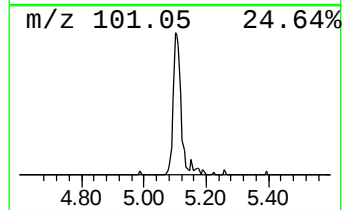
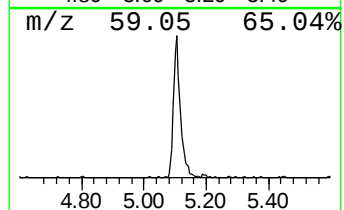
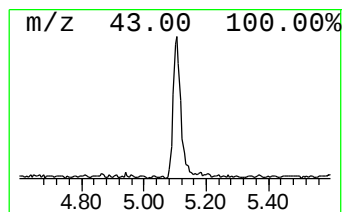
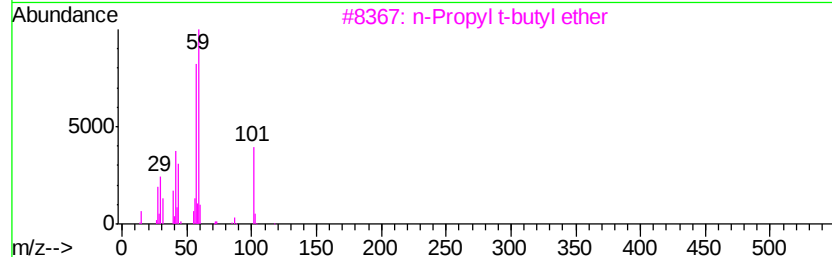
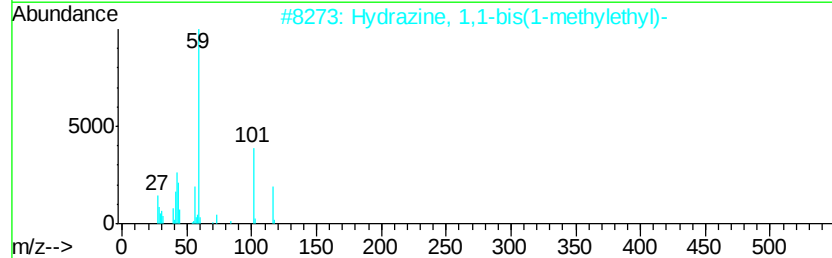
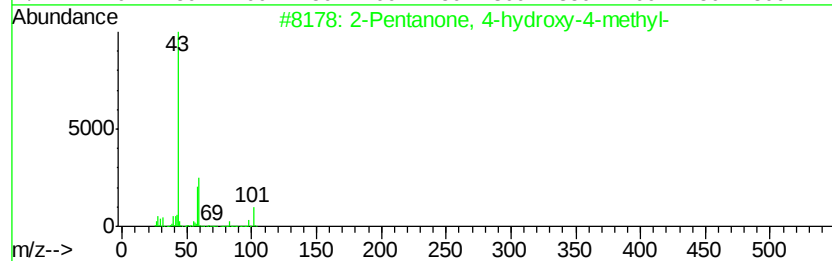
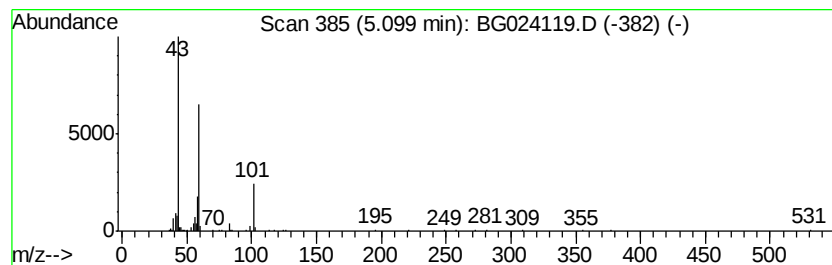
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	8.94 ng	139430	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			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
4			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	25
5			Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	23



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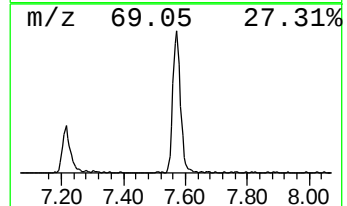
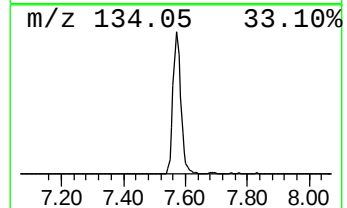
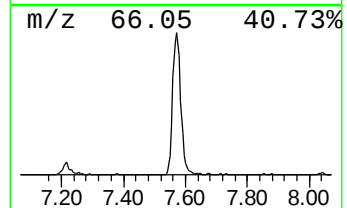
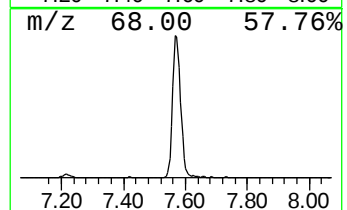
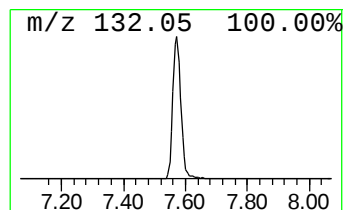
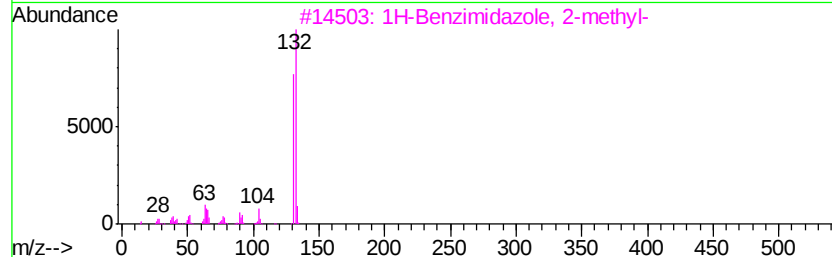
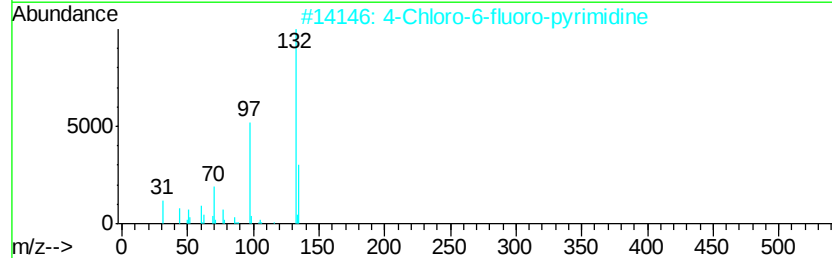
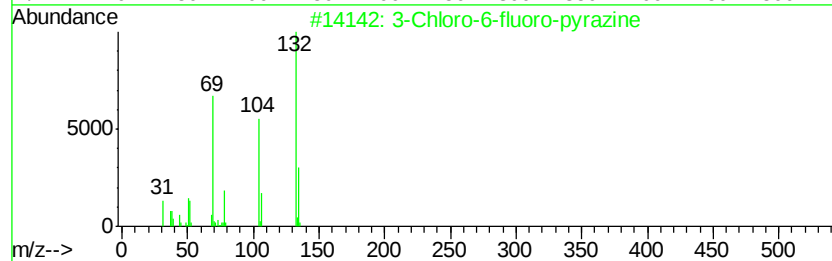
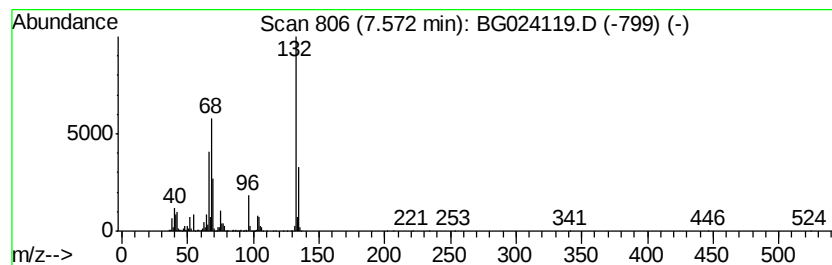
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	104.28 ng	1627020	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	18
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	8.9	ng	139430	1	8.04	312044	20.0
unknown7.57	7.57	104.3	ng	1627020	1	8.04	312044	20.0