

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024116.D
 Acq On : 24 Sep 2016 9:28
 Operator : UM/SJ
 Sample : H4893-07
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LIP-12

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-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	5.102	382	385	390	rVB2	60424	87496	1.92%	0.412%
2	5.590	463	468	481	rVB	423361	754867	16.59%	3.556%
3	7.217	739	745	757	rBV	262741	510552	11.22%	2.405%
4	7.569	799	805	820	rBV	908020	1598760	35.13%	7.532%
5	8.039	879	885	894	rVB	170148	293186	6.44%	1.381%
6	8.362	931	940	954	rBV	806504	1422151	31.25%	6.700%
7	9.220	1079	1086	1100	rBV	754043	1402115	30.81%	6.605%
8	10.877	1359	1368	1379	rBV	232115	456440	10.03%	2.150%
9	13.326	1778	1785	1802	rBV	2077426	3287138	72.24%	15.485%
10	14.719	2015	2022	2033	rBV	442775	710250	15.61%	3.346%
11	16.222	2270	2278	2289	rBV	1850696	3006350	66.07%	14.163%
12	17.480	2486	2492	2503	rBV2	577037	942440	20.71%	4.440%
13	18.355	2637	2641	2649	rBV5	43711	77702	1.71%	0.366%
14	20.094	2931	2937	2948	rBV	3352333	4550359	100.00%	21.436%
15	21.791	3220	3226	3237	rVB	613294	1085952	23.87%	5.116%
16	25.122	3783	3793	3806	rBV	339772	1041633	22.89%	4.907%

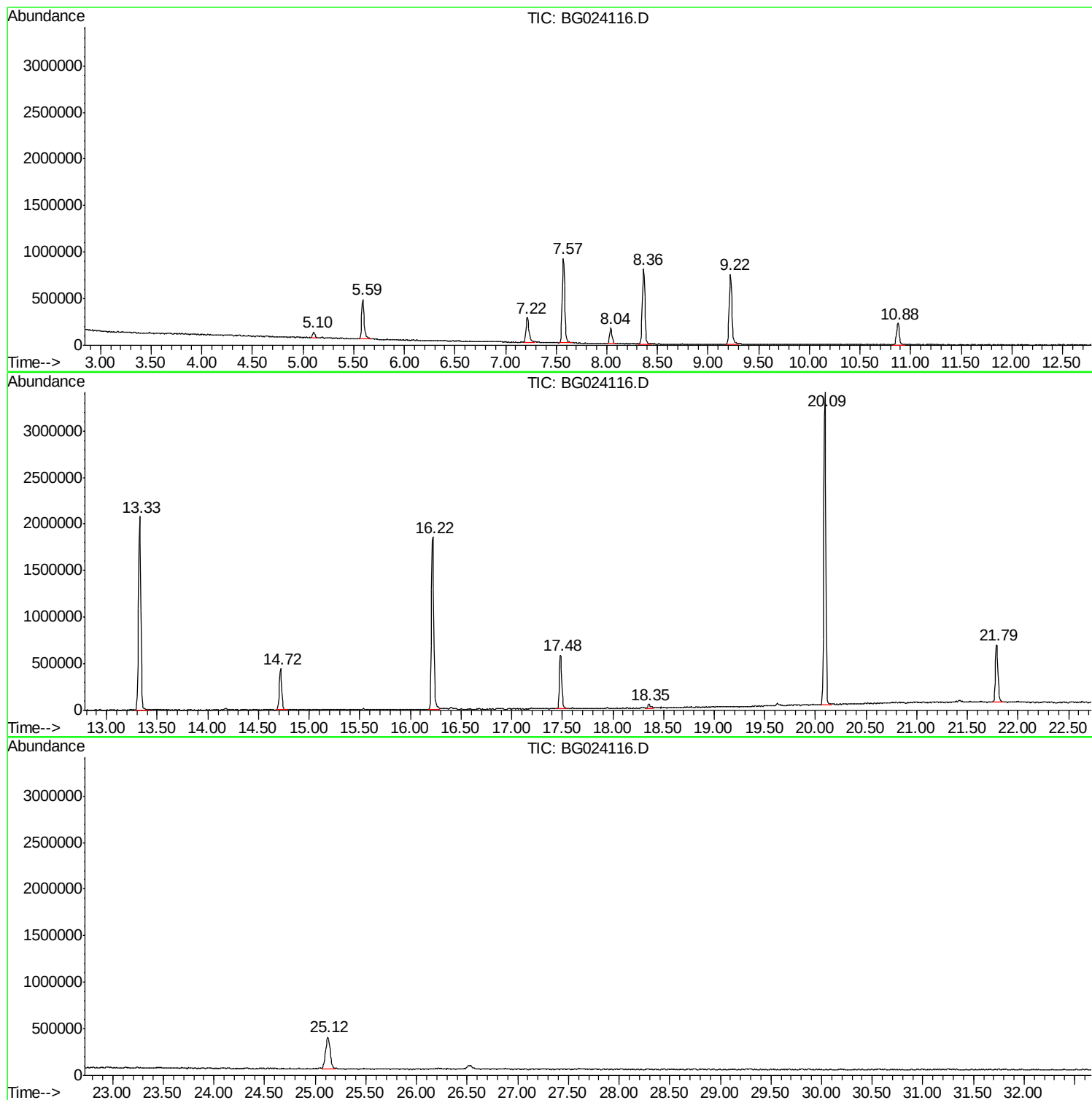
Sum of corrected areas: 21227391

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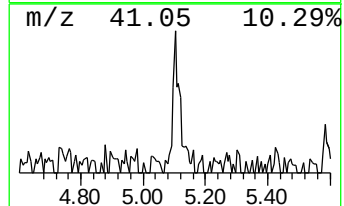
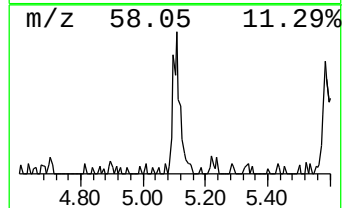
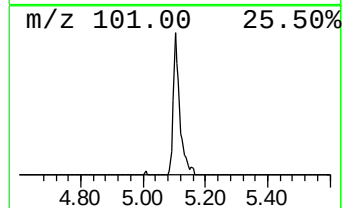
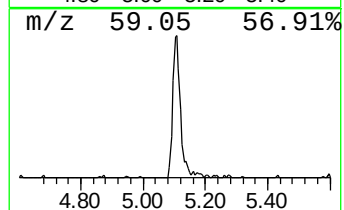
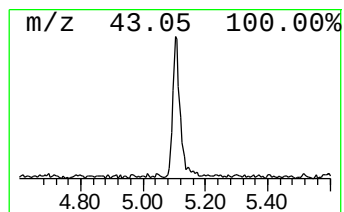
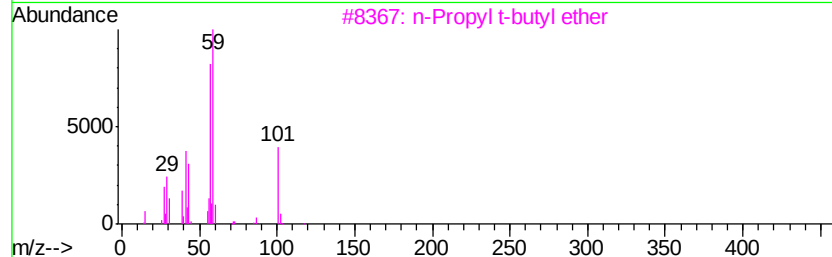
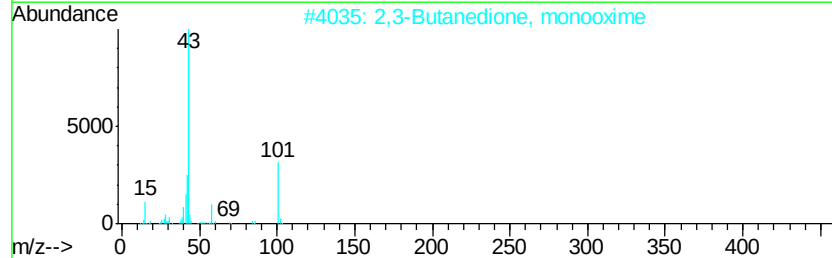
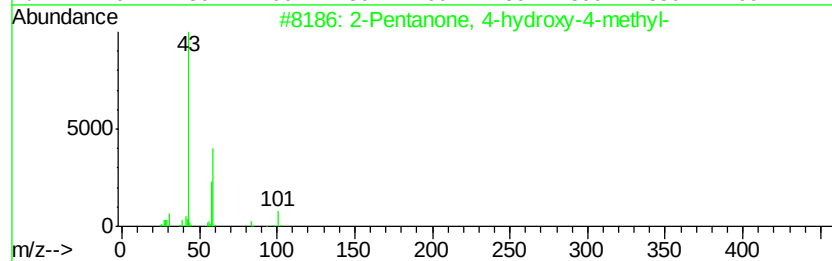
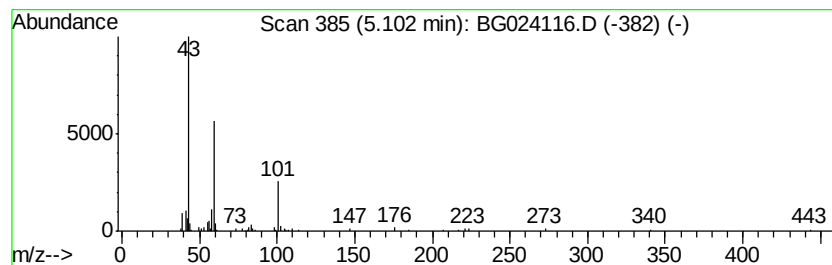
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	5.97 ng	87496	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	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	25
4			3-Heptadecanol	256	C17H36O	084534-30-5	23
5			3-Hexanol	102	C6H14O	000623-37-0	23



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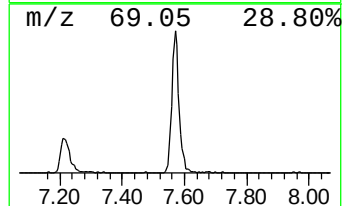
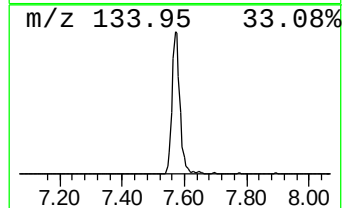
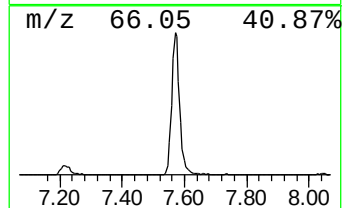
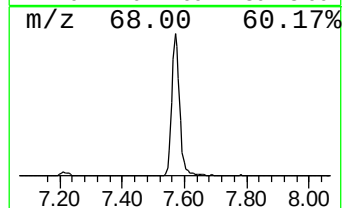
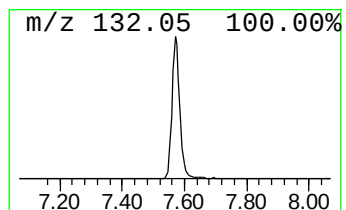
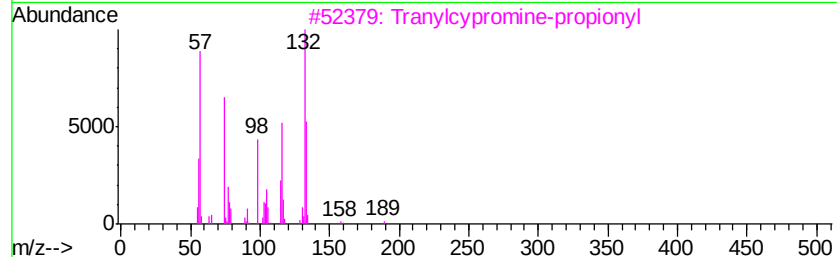
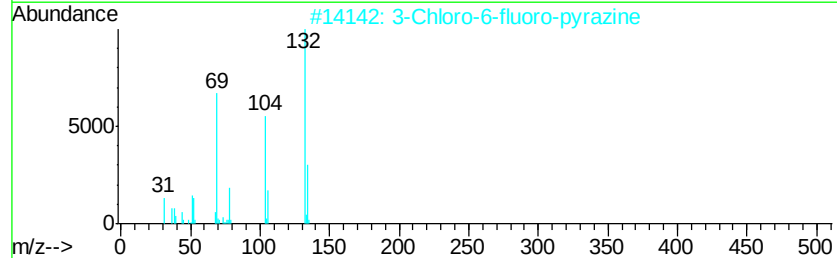
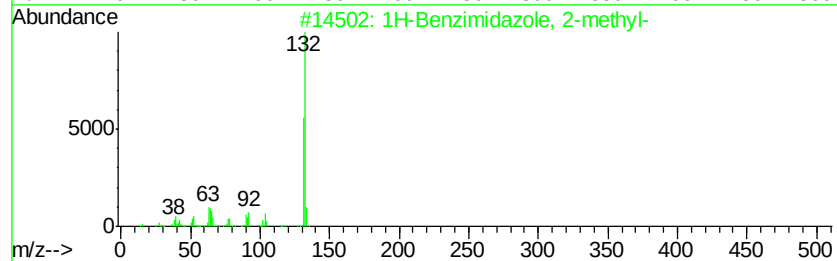
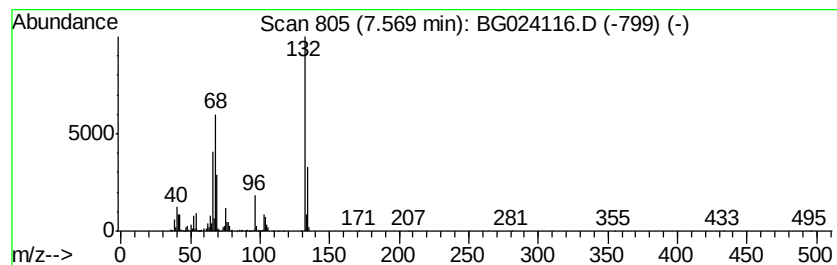
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	109.06 ng	1598760	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	6.0	ng	87496	1	8.04	293186	20.0
unknown7.57	7.57	109.1	ng	1598760	1	8.04	293186	20.0