

Data Path : Z:\HPCHEM1\BNA G\DATA\BG072716\
 Data File : BG023289.D
 Acq On : 27 Jul 2016 19:01
 Operator : UM/SJ
 Sample : H4152-13
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-18-6.0-15.0

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-BG072616.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	5.190	394	400	415	rBV	1214263	1899565	16.07%	3.593%
2	5.672	475	482	492	rBV	1059476	1701300	14.39%	3.218%
3	7.287	750	757	769	rBV	816245	1479526	12.52%	2.799%
4	7.663	813	821	830	rBV	1461178	2491493	21.08%	4.713%
5	8.133	895	901	908	rBV	621680	1077762	9.12%	2.039%
6	8.462	949	957	966	rBV	2026811	3512170	29.71%	6.644%
7	9.314	1093	1102	1112	rBV	1797139	3307574	27.98%	6.257%
8	10.971	1377	1384	1392	rVB	872291	1559520	13.19%	2.950%
9	13.420	1793	1801	1811	rBV	4798841	7734608	65.43%	14.632%
10	14.249	1936	1942	1948	rBV	305181	463226	3.92%	0.876%
11	14.807	2029	2037	2046	rBV2	1456199	2362519	19.99%	4.469%
12	16.305	2285	2292	2299	rBV	1905070	2951954	24.97%	5.584%
13	16.498	2320	2325	2337	rBV4	93524	187818	1.59%	0.355%
14	17.568	2500	2507	2514	rBV2	2025202	3116415	26.36%	5.895%
15	20.176	2945	2951	2961	rBV	8313082	11820732	100.00%	22.362%
16	21.891	3236	3243	3254	rVB	2101282	3609932	30.54%	6.829%
17	23.648	3537	3542	3548	rVB2	68711	120536	1.02%	0.228%
18	25.298	3811	3823	3837	rBV2	1139511	3464998	29.31%	6.555%

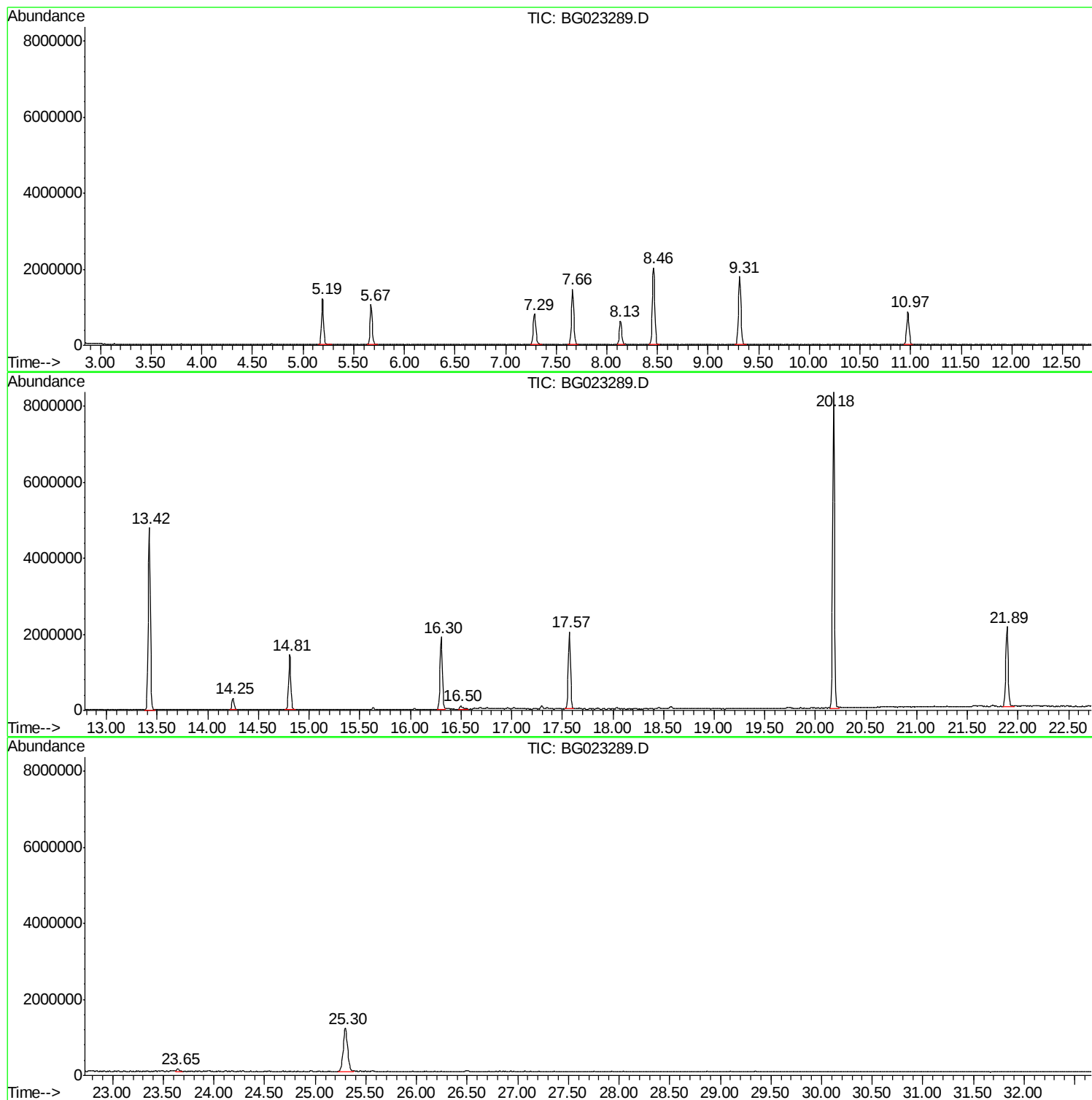
Sum of corrected areas: 52861648

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG072616.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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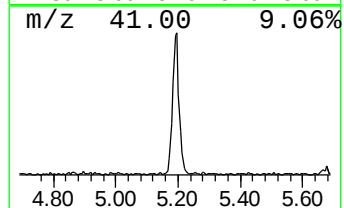
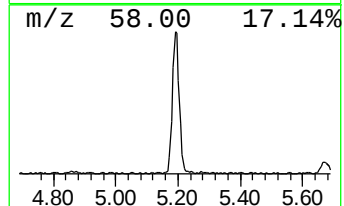
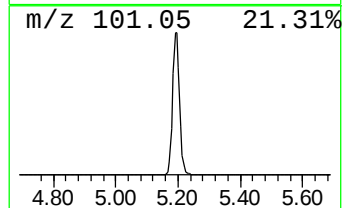
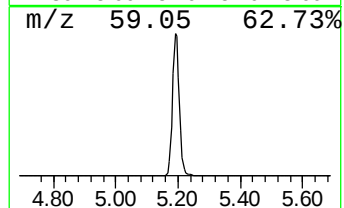
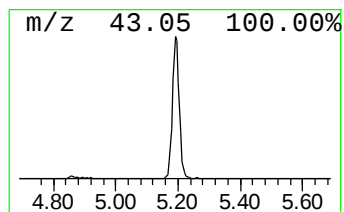
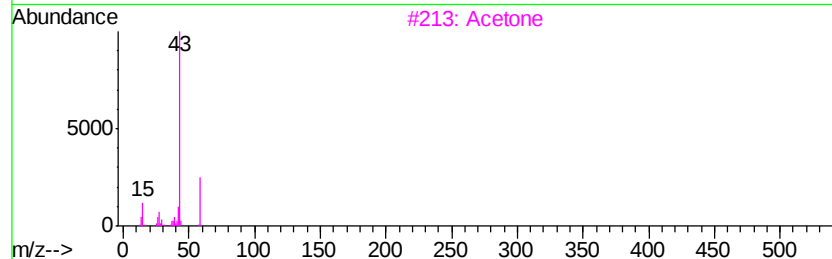
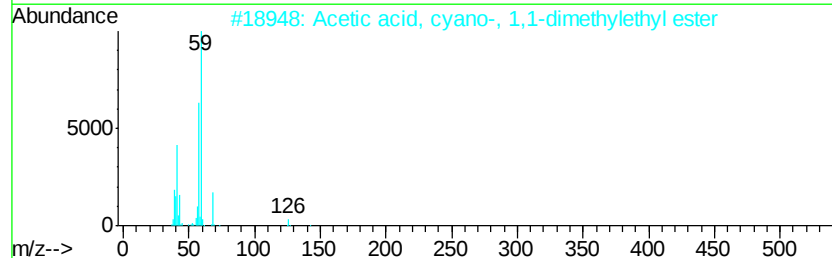
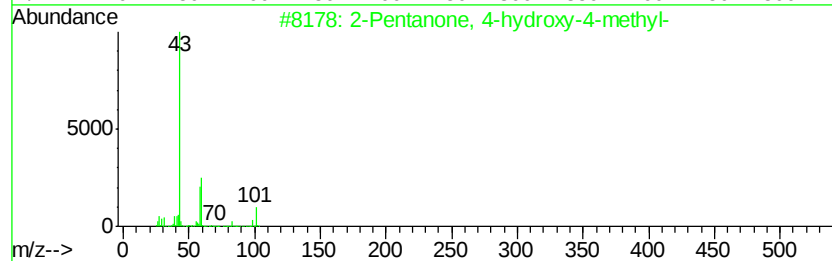
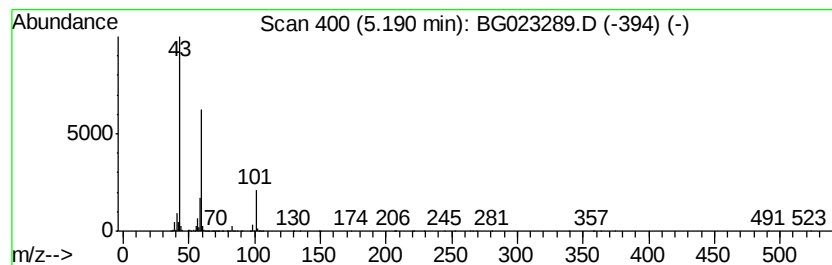
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	35.25 ng	1899570	1,4-Dichlorobenzene-d4	8.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
3	Acetone	58	C3H6O	000067-64-1	9	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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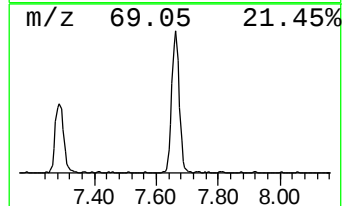
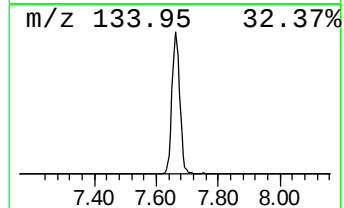
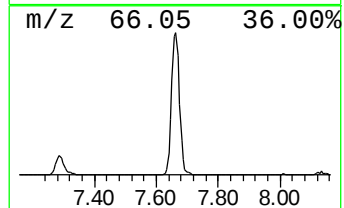
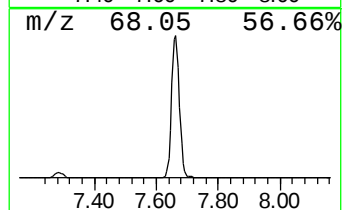
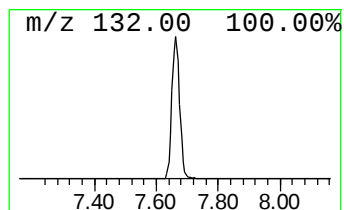
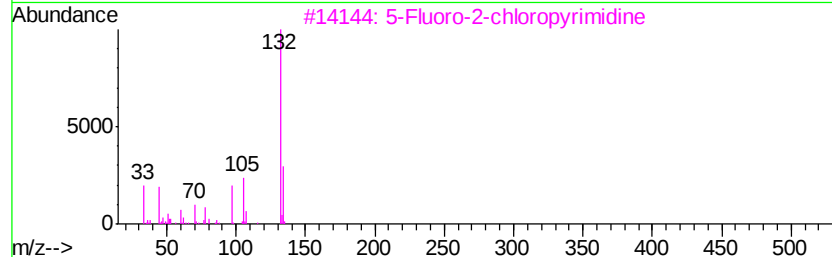
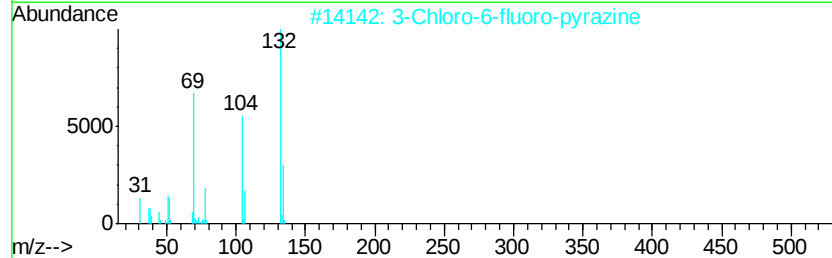
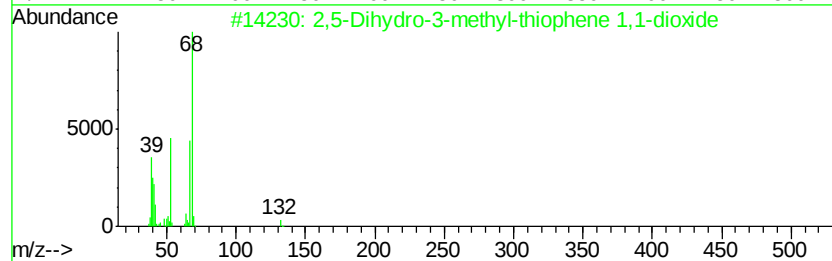
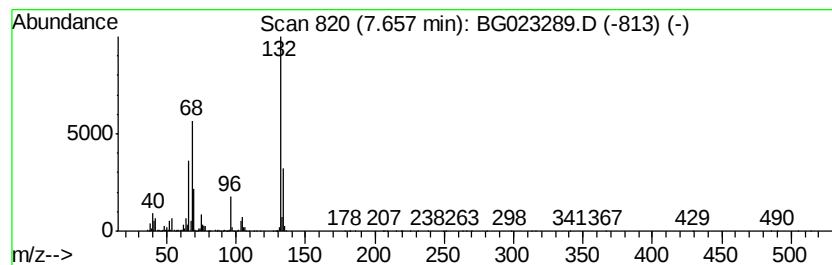
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	46.23 ng	2491490	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Dihydro-3-methyl-thiophene 1...	132	C5H8O2S	001193-10-8	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.19	35.3	ng	1899570	1	8.13	1077760	20.0
unknown7.66	7.66	46.2	ng	2491490	1	8.13	1077760	20.0