

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060617\
 Data File : BG027254.D
 Acq On : 6 Jun 2017 18:12
 Operator : SJ/MA
 Sample : PB99498BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99498BL

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-BG060517.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	5.706	337	343	356	rVB	103989	180958	3.69%	0.724%
2	6.147	411	418	431	rVB	1186550	2065398	42.12%	8.261%
3	7.698	674	682	694	rBV	1199964	2216882	45.21%	8.867%
4	8.098	741	750	760	rBV	1158848	2125398	43.34%	8.501%
5	8.568	819	830	838	rBV	179096	329066	6.71%	1.316%
6	8.891	874	885	895	rVB	785689	1452832	29.63%	5.811%
7	9.731	1019	1028	1037	rBV	645342	1235607	25.20%	4.942%
8	11.388	1301	1310	1317	rBV	271400	530243	10.81%	2.121%
9	13.773	1708	1716	1726	rBV	2015320	3321662	67.74%	13.285%
10	15.148	1942	1950	1959	rBV2	465724	795677	16.23%	3.182%
11	16.629	2195	2202	2214	rBV	1723470	2675037	54.55%	10.699%
12	17.898	2411	2418	2425	rBV	633928	995597	20.30%	3.982%
13	20.477	2850	2857	2863	rBV	3511374	4903627	100.00%	19.612%
14	22.287	3158	3165	3172	rVB	582451	1100575	22.44%	4.402%
15	25.983	3784	3794	3809	rVB2	325539	1074108	21.90%	4.296%

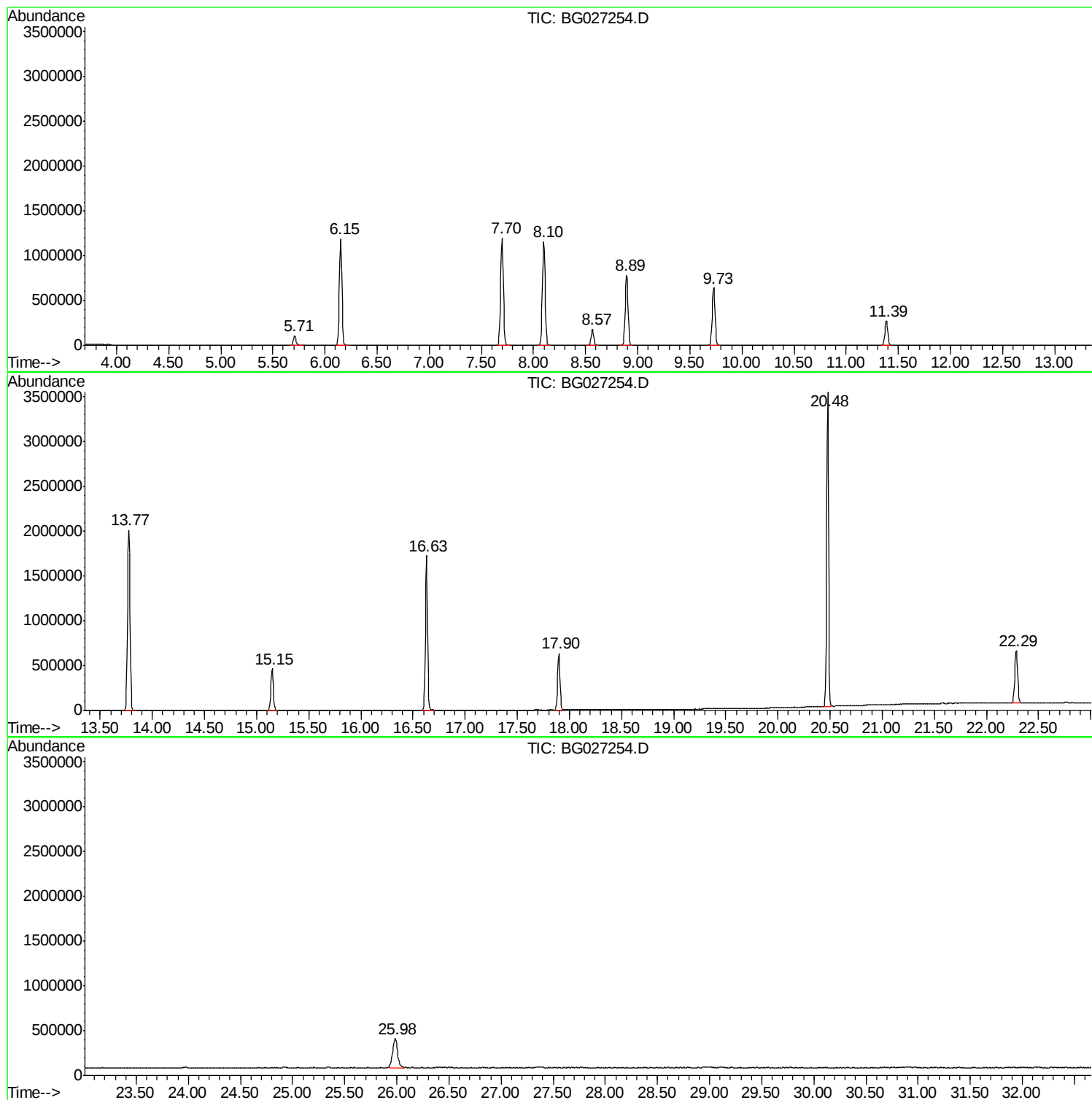
Sum of corrected areas: 25002667

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060517.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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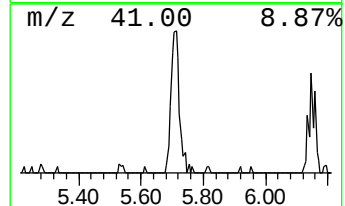
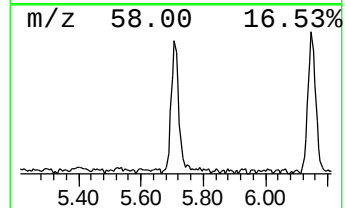
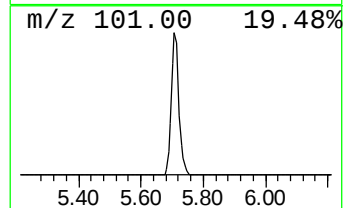
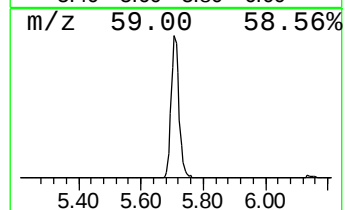
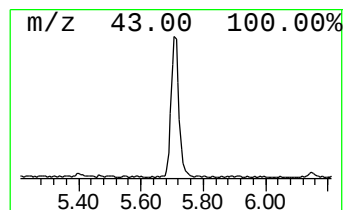
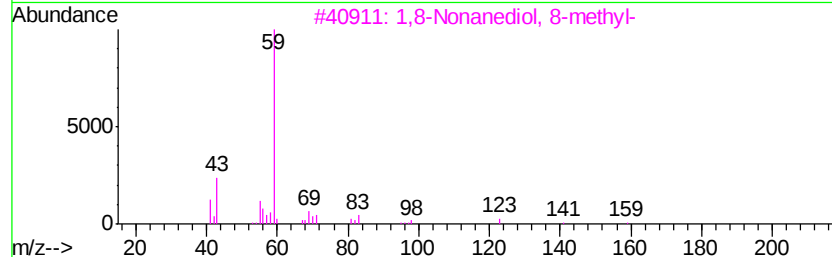
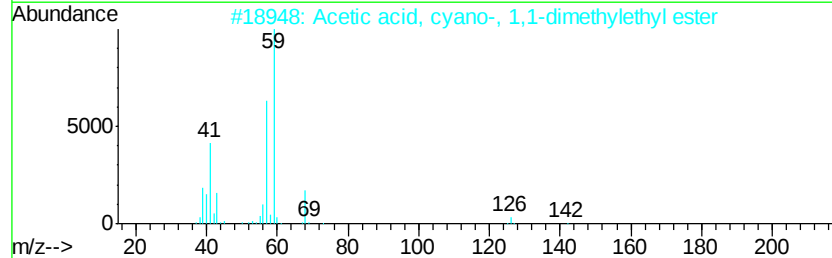
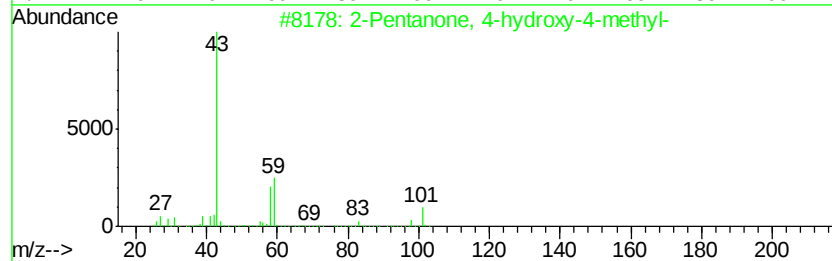
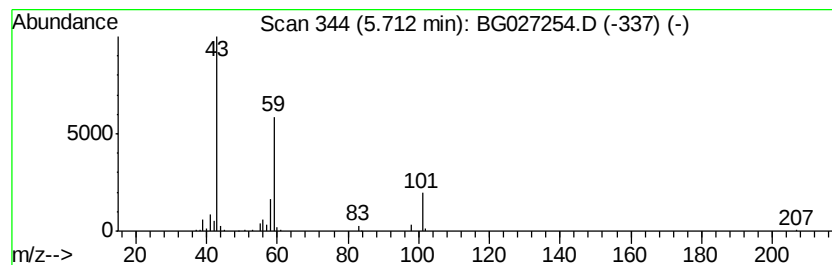
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.71	11.00 ng	180958	1,4-Dichlorobenzene-d4	8.57

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-dimethy...	141	C7H11NO2	001116-98-9	32		
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25		



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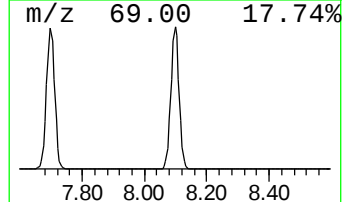
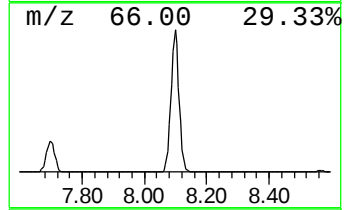
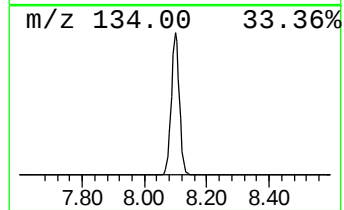
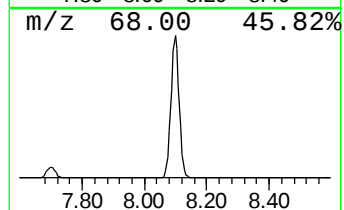
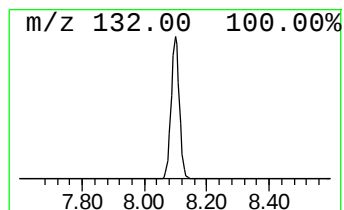
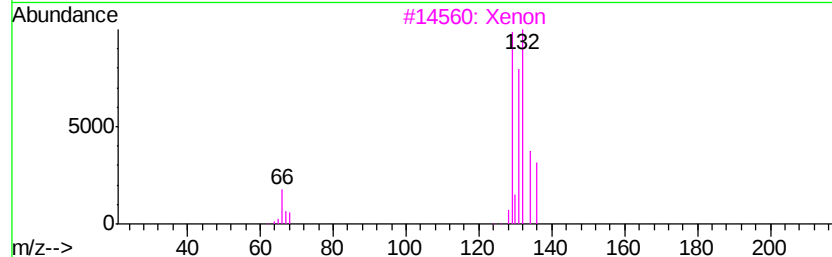
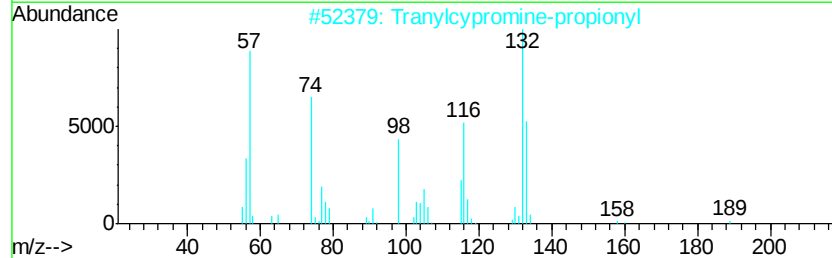
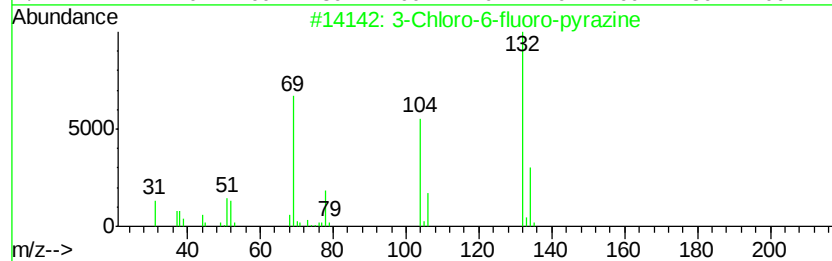
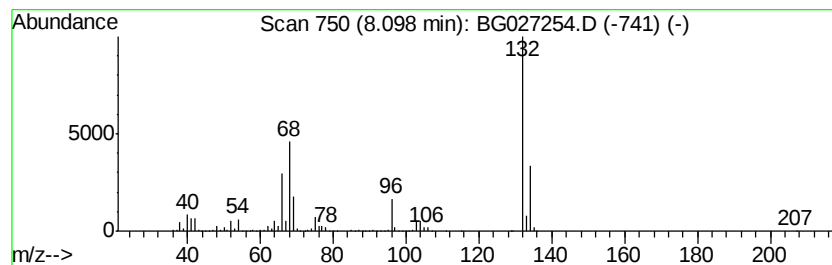
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Peak Number 2 unknown8.10 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.10	129.18 ng	2125400	1,4-Dichlorobenzene-d4	8.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Xenon	132	Xe	007440-63-3	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.71	11.0	ng	180958	1	8.57	329066	20.0
unknown8.10	8.10	129.2	ng	2125400	1	8.57	329066	20.0