

Data Path : Z:\HPCHEM1\BNA G\Data\BG042017\
 Data File : BG026679.D
 Acq On : 20 Apr 2017 17:30
 Operator : SJ/MA
 Sample : PB98380BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB98380BL

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-BG041817.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	5.126	341	347	356	rBV	241049	358905	2.72%	0.584%
2	5.620	423	431	446	rBV	2653176	4353008	32.93%	7.081%
3	7.218	696	703	714	rBV	2525308	4367011	33.04%	7.104%
4	7.571	756	763	776	rBV	2727770	4749497	35.93%	7.726%
5	8.029	832	841	849	rBV	533687	883827	6.69%	1.438%
6	8.352	887	896	907	rBV	2094431	3614701	27.35%	5.880%
7	9.186	1029	1038	1049	rBV	1487196	2645311	20.01%	4.303%
8	10.826	1308	1317	1327	rBV	733046	1343328	10.16%	2.185%
9	13.264	1725	1732	1743	rBV	4678379	7443705	56.31%	12.108%
10	14.639	1958	1966	1979	rBV2	1233915	1973313	14.93%	3.210%
11	16.125	2211	2219	2229	rBV	4390041	6657939	50.37%	10.830%
12	17.371	2424	2431	2441	rBV2	1944378	2868526	21.70%	4.666%
13	19.974	2867	2874	2886	rBV	9333390	13218195	100.00%	21.501%
14	21.619	3147	3154	3165	rBV	2178066	3480517	26.33%	5.662%
15	24.792	3683	3694	3707	rBV2	1254809	3518619	26.62%	5.724%

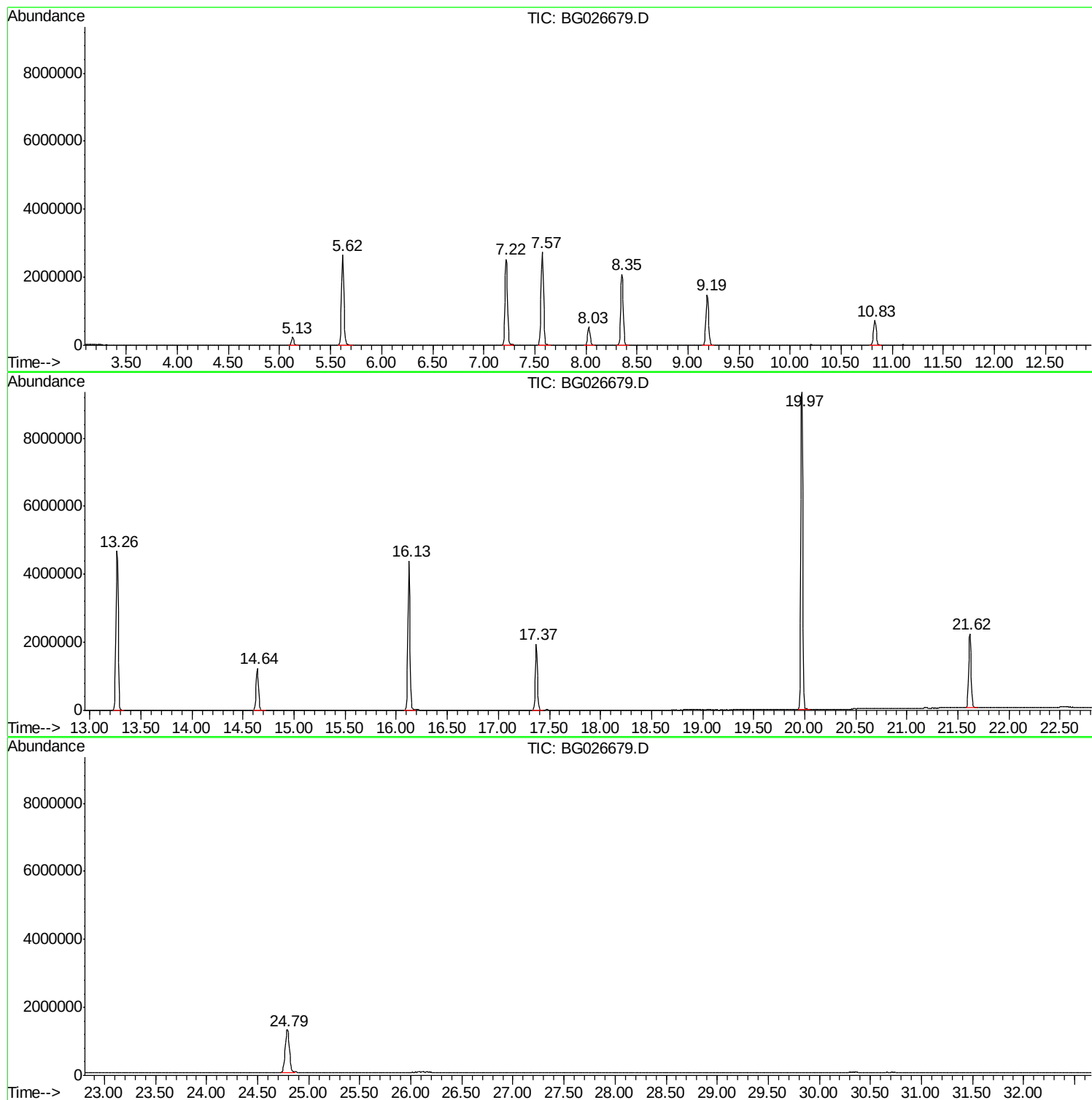
Sum of corrected areas: 61476402

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041817.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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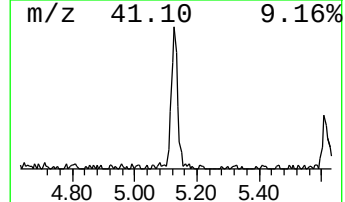
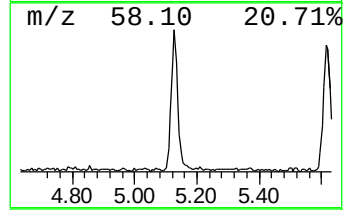
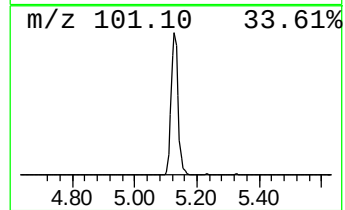
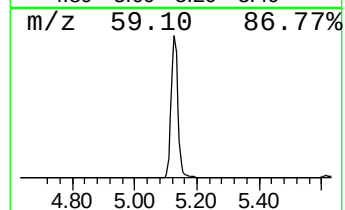
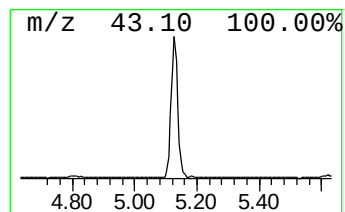
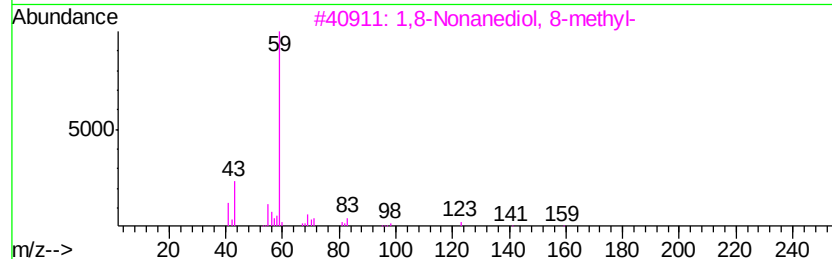
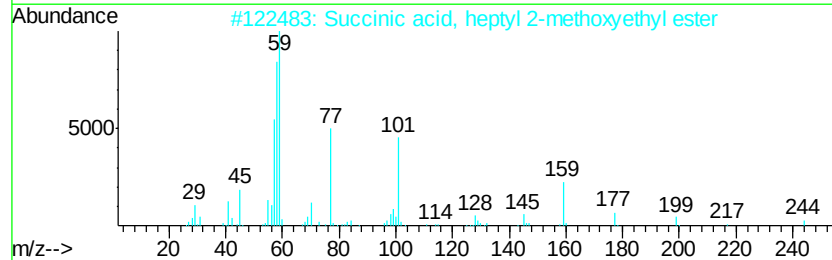
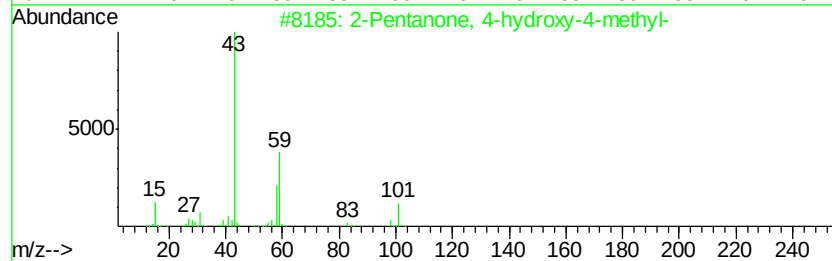
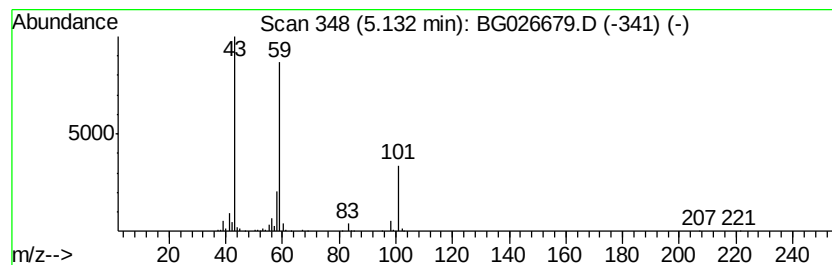
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.13	8.12 ng	358905	1,4-Dichlorobenzene-d4	8.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	36
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	28
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	12



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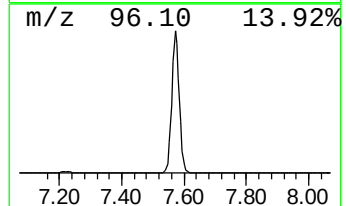
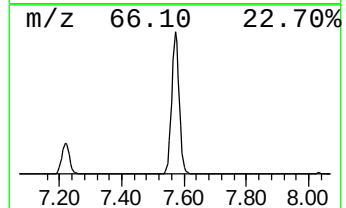
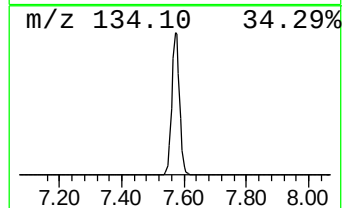
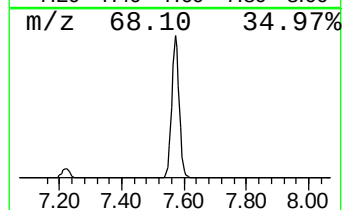
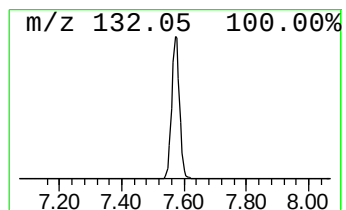
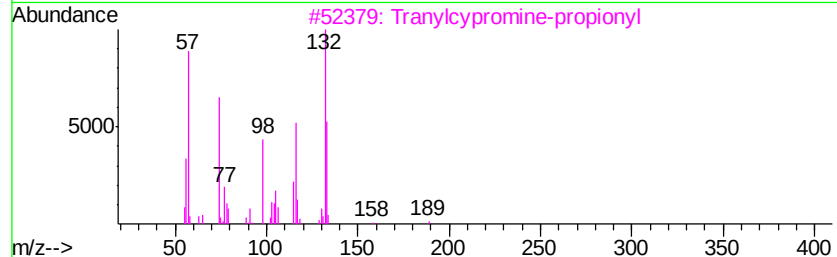
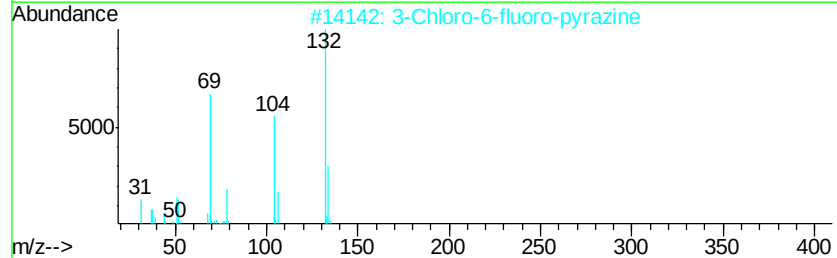
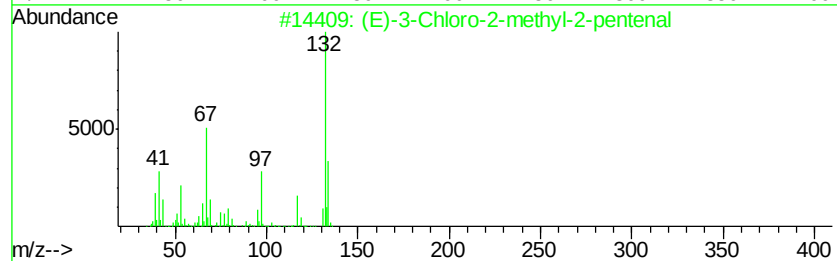
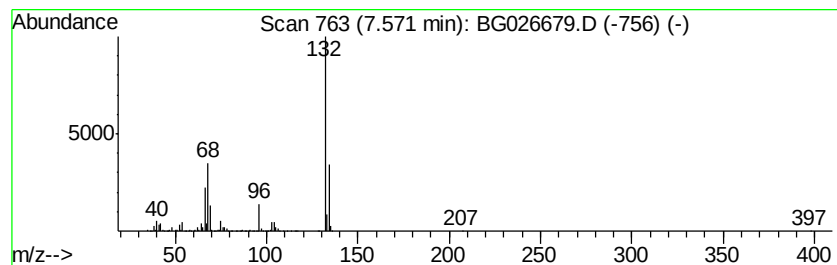
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	107.48 ng	4749500	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	8.1	ng	358905	1	8.03	883827	20.0
unknown7.57	7.57	107.5	ng	4749500	1	8.03	883827	20.0