

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062617\
 Data File : BG027685.D
 Acq On : 26 Jun 2017 21:48
 Operator : SJ/JU
 Sample : I3884-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CAN-GW-05

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-BG062117.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	5.624	290	295	304	rBV2	15858	24645	2.99%	0.648%
2	6.071	365	371	381	rBV	125968	211025	25.64%	5.546%
3	7.628	629	636	647	rBB	120377	216561	26.31%	5.692%
4	8.021	696	703	712	rBV	127021	229046	27.83%	6.020%
5	8.486	776	782	789	rBB	47710	85859	10.43%	2.257%
6	8.815	831	838	845	rBB	87441	161683	19.64%	4.250%
7	9.649	973	980	993	rVB	71110	135857	16.50%	3.571%
8	11.306	1254	1262	1270	rBB	66901	129001	15.67%	3.391%
9	13.697	1661	1669	1676	rBB	219792	356220	43.28%	9.363%
10	15.072	1897	1903	1911	rBB2	118192	199932	24.29%	5.255%
11	16.558	2150	2156	2166	rBV2	71937	129665	15.75%	3.408%
12	17.822	2364	2371	2378	rBV2	209614	327753	39.82%	8.614%
13	20.395	2803	2809	2816	rBV	617147	823145	100.00%	21.635%
14	22.181	3105	3113	3121	rBV	212146	404986	49.20%	10.644%
15	25.801	3717	3729	3747	rVB2	114620	369370	44.87%	9.708%

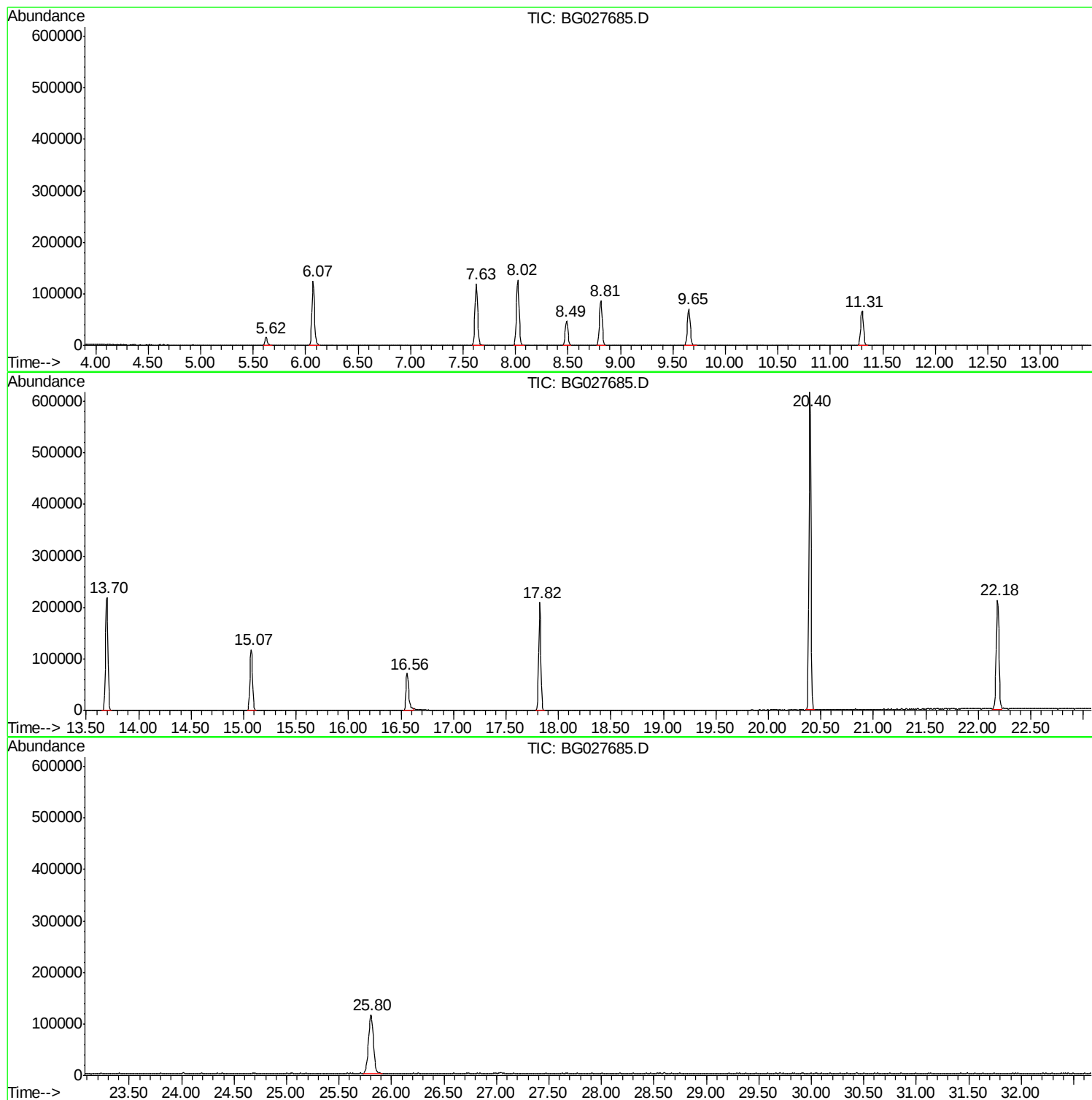
Sum of corrected areas: 3804748

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062117.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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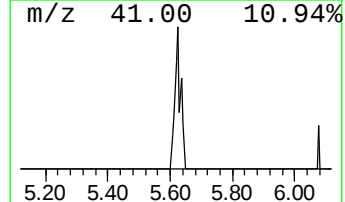
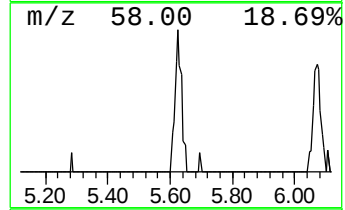
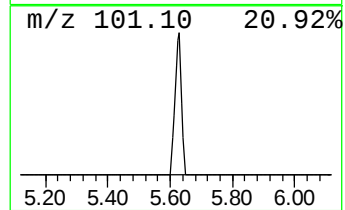
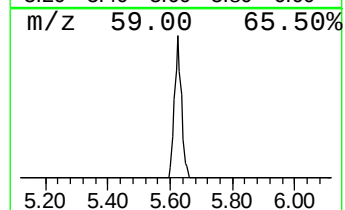
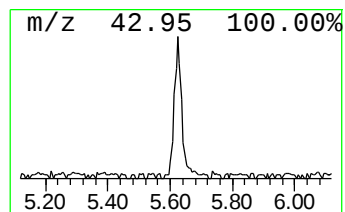
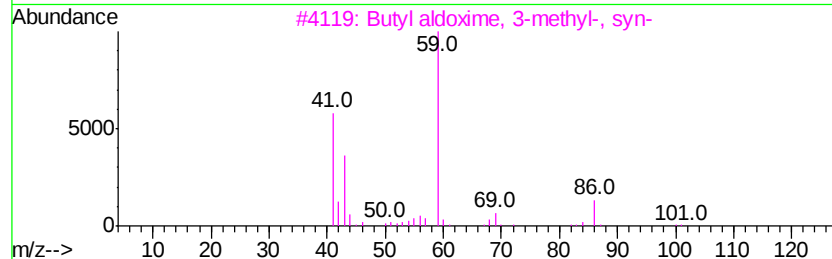
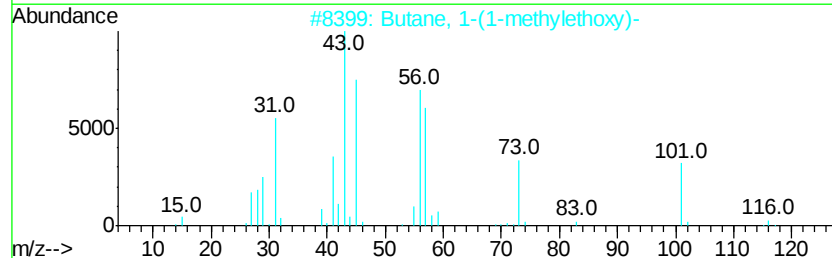
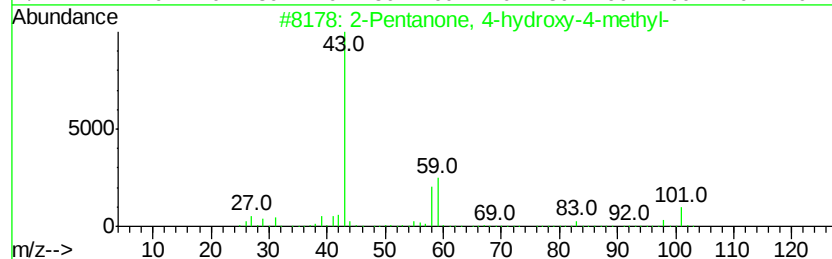
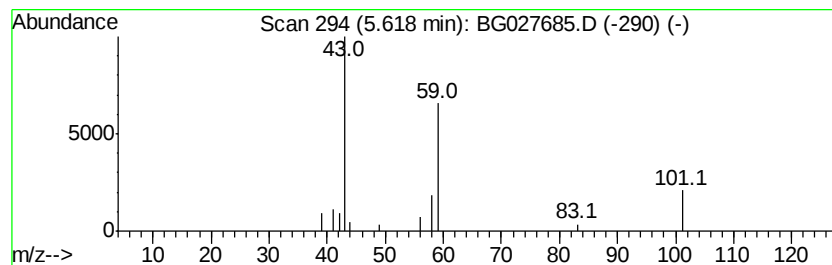
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.62	5.74 ng	24645	1,4-Dichlorobenzene-d4	8.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	32
3		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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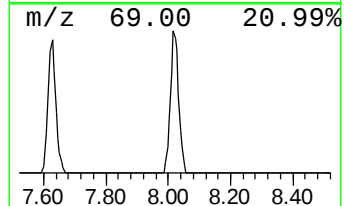
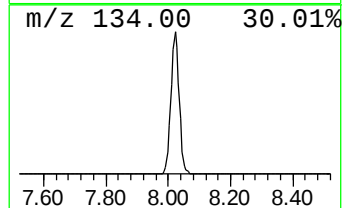
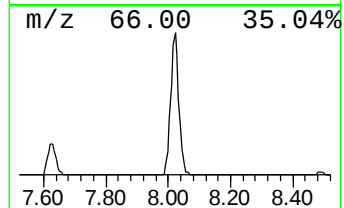
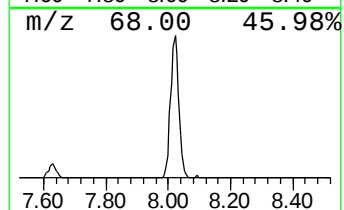
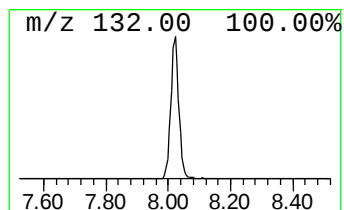
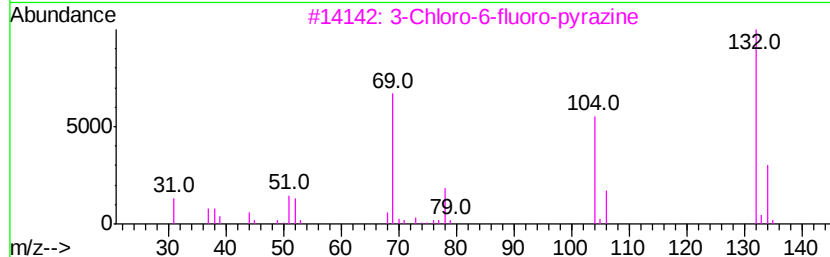
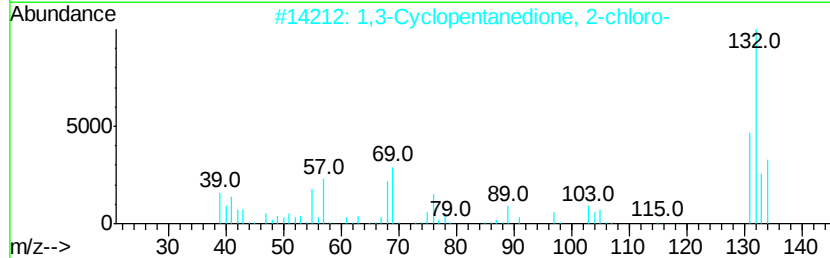
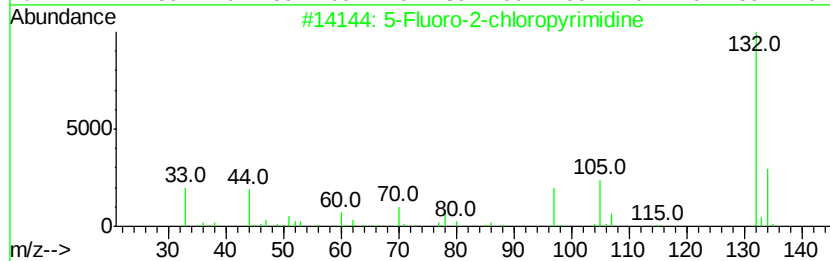
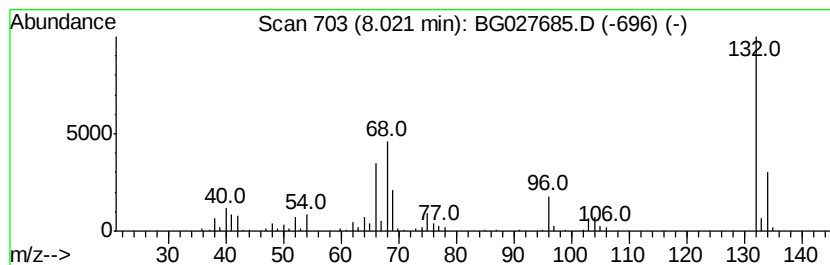
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Peak Number 2 unknown8.02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	53.35 ng	229046	1,4-Dichlorobenzene-d4	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	12
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.62	5.7	ng	24645	1	8.49	85859	20.0
unknown8.02	8.02	53.4	ng	229046	1	8.49	85859	20.0