

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030996.D
 Acq On : 14 Nov 2017 5:58
 Operator : SJ/JU
 Sample : PB104197BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104197BL

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-BG111017.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	5.216	323	328	337	rBV	84713	132311	3.70%	0.869%
2	5.704	403	411	424	rBV	547591	911879	25.51%	5.988%
3	7.308	676	684	694	rBV	514754	929028	25.99%	6.101%
4	7.678	738	747	759	rBV	542140	966023	27.03%	6.344%
5	8.142	816	826	836	rVB	151606	251871	7.05%	1.654%
6	8.465	873	881	894	rVB	447043	788535	22.06%	5.178%
7	9.306	1017	1024	1039	rBV	291970	563476	15.76%	3.700%
8	10.957	1296	1305	1316	rBV	172659	342490	9.58%	2.249%
9	13.383	1711	1718	1730	rBV	990714	1605267	44.91%	10.541%
10	14.758	1946	1952	1965	rBV2	299006	527659	14.76%	3.465%
11	16.244	2197	2205	2227	rBV	952346	1576970	44.12%	10.355%
12	17.502	2412	2419	2433	rBV	508922	814114	22.78%	5.346%
13	20.093	2853	2860	2873	rBV	2675176	3574509	100.00%	23.473%
14	21.773	3138	3146	3160	rBV2	648123	1157818	32.39%	7.603%
15	25.075	3697	3708	3723	rVB2	360332	1086505	30.40%	7.135%

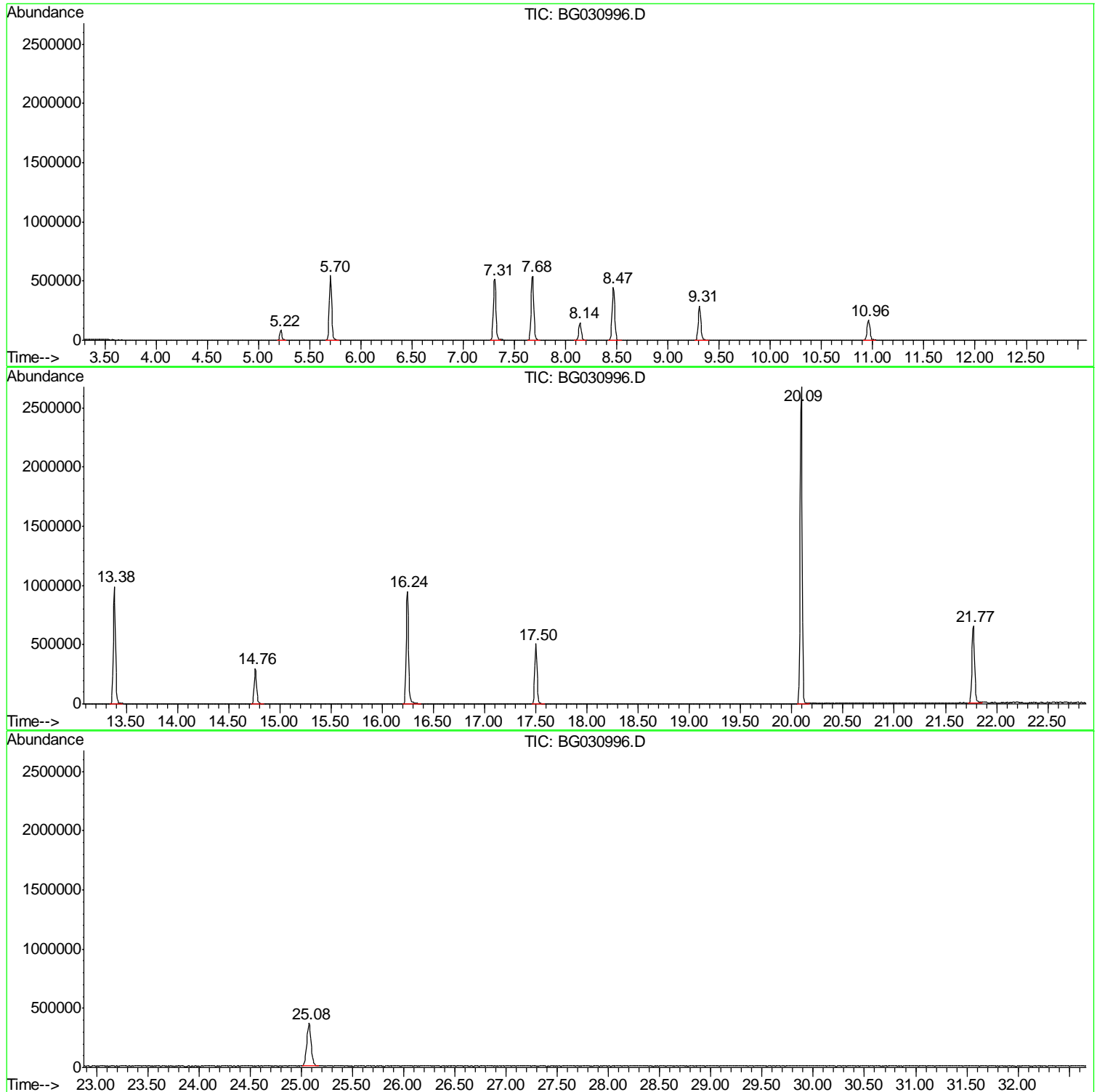
Sum of corrected areas: 15228455

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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 ClientSampled :
 PB104197BL

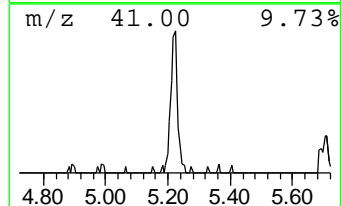
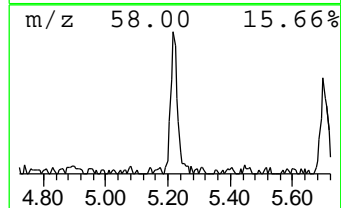
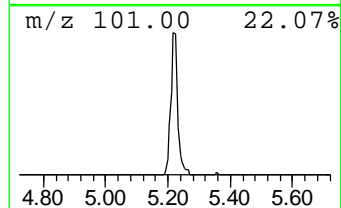
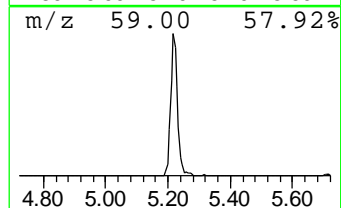
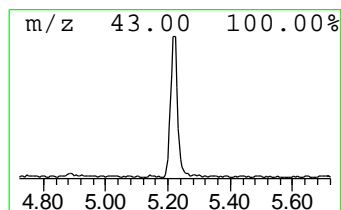
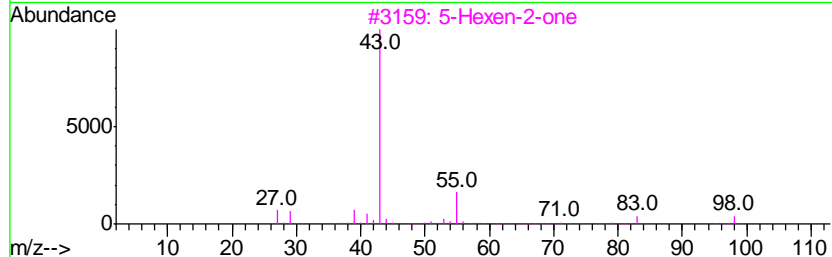
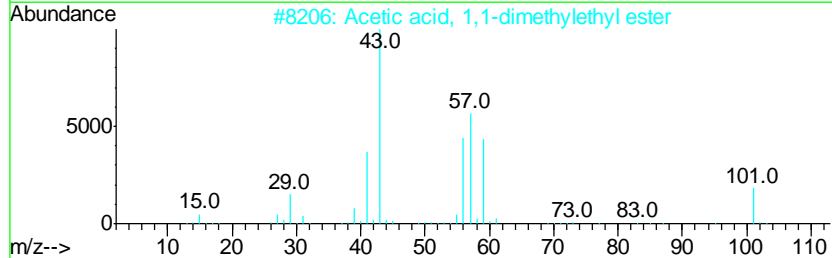
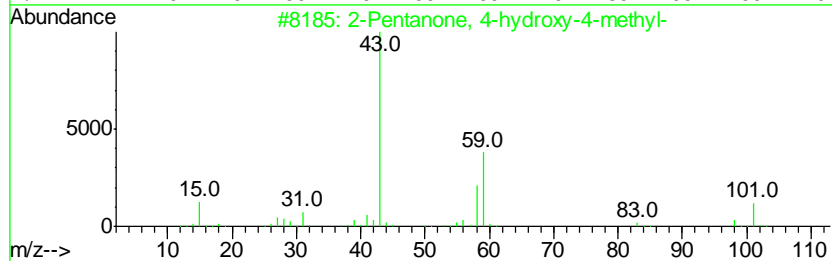
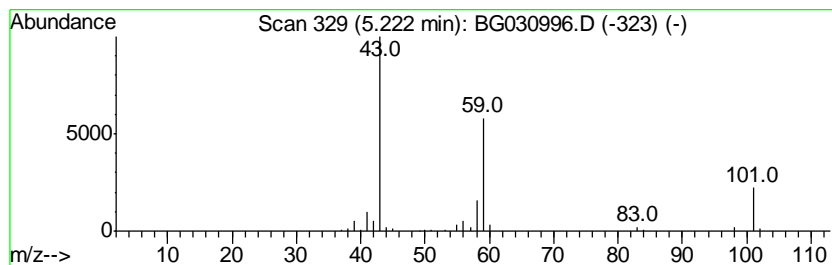
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 TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	10.51 ng	132311	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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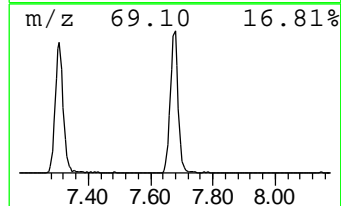
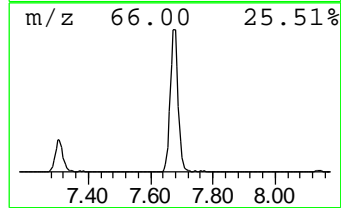
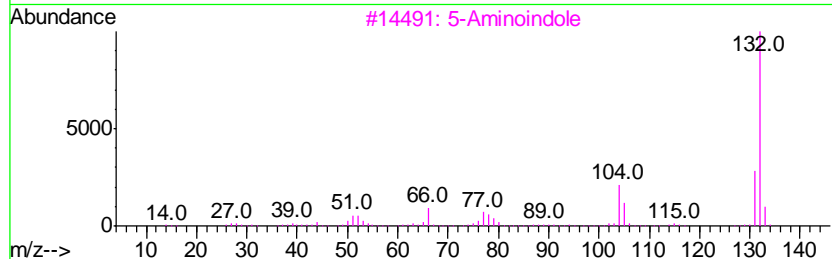
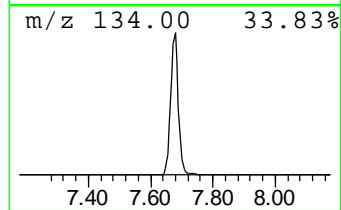
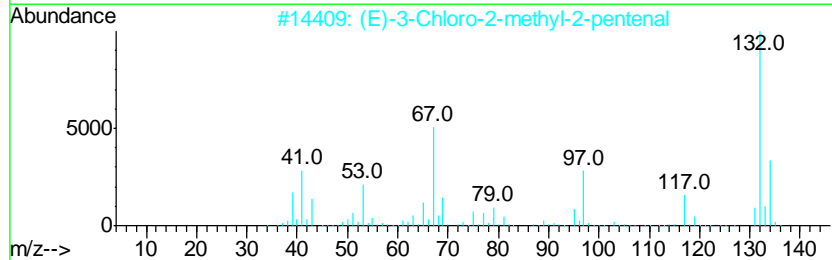
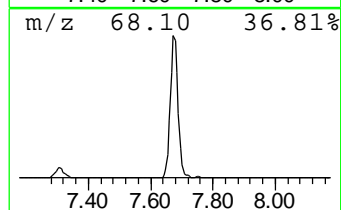
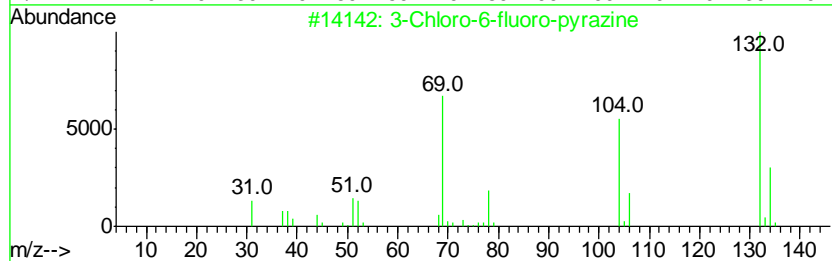
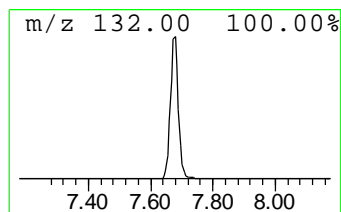
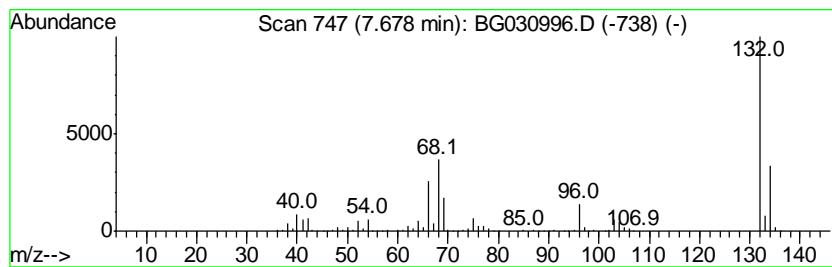
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 Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	76.71 ng	966023	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	10.5	ng	132311	1	8.14	251871	20.0
unknown7.68	7.68	76.7	ng	966023	1	8.14	251871	20.0