

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111017\
 Data File : BG030916.D
 Acq On : 10 Nov 2017 21:06
 Operator : SJ/JU
 Sample : PB104116BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104116BL

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	5.218	323	328	336	rBV	113801	172683	4.08%	0.939%
2	5.699	403	410	422	rBV	671779	1132987	26.76%	6.158%
3	7.309	676	684	694	rBV	680045	1197508	28.28%	6.508%
4	7.679	739	747	757	rBV	687518	1211418	28.61%	6.584%
5	8.143	819	826	834	rBV	166710	283560	6.70%	1.541%
6	8.467	872	881	893	rBV	541536	954854	22.55%	5.190%
7	9.307	1016	1024	1036	rBV	374375	713463	16.85%	3.878%
8	10.958	1297	1305	1316	rBV	209539	407588	9.63%	2.215%
9	13.384	1711	1718	1730	rBV	1271011	2021627	47.74%	10.988%
10	14.765	1946	1953	1963	rBV2	373715	647528	15.29%	3.519%
11	16.246	2198	2205	2221	rBV	1159808	1857783	43.88%	10.097%
12	17.503	2412	2419	2430	rBV2	583718	948682	22.41%	5.156%
13	20.094	2853	2860	2869	rBV	3080379	4234219	100.00%	23.013%
14	21.769	3138	3145	3158	rBV2	743604	1332079	31.46%	7.240%
15	25.071	3695	3707	3719	rBV2	443001	1283180	30.30%	6.974%

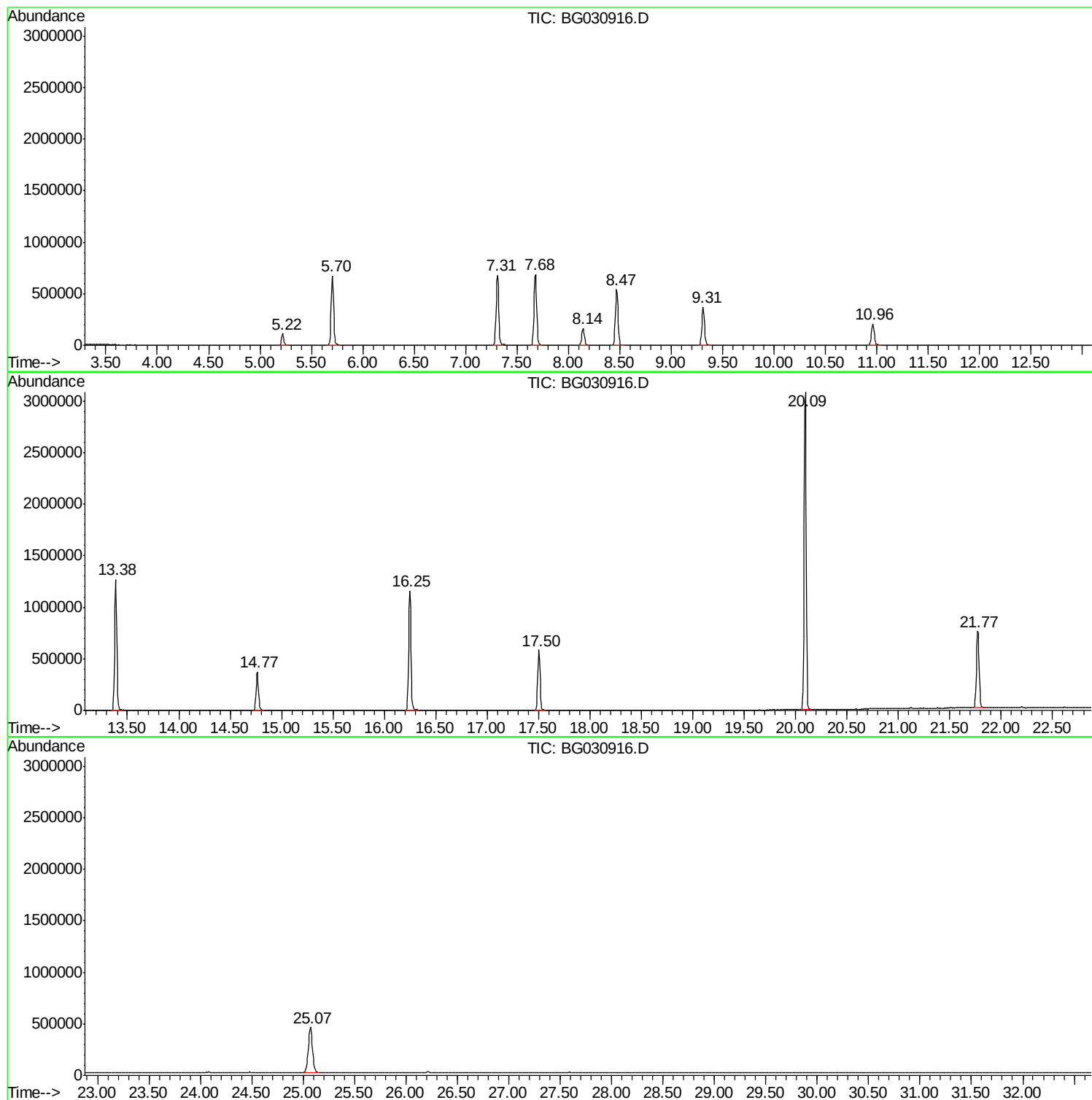
Sum of corrected areas: 18399159

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



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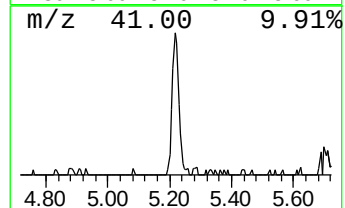
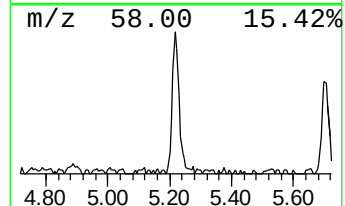
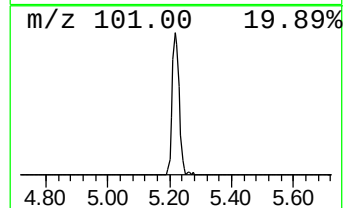
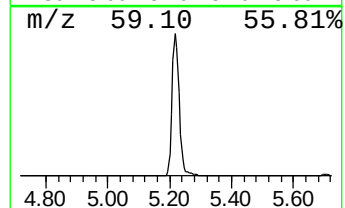
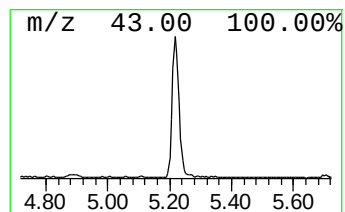
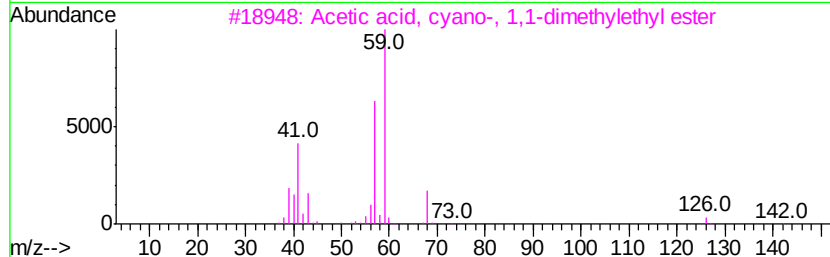
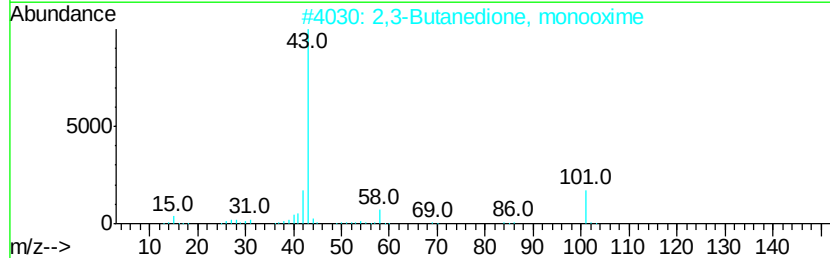
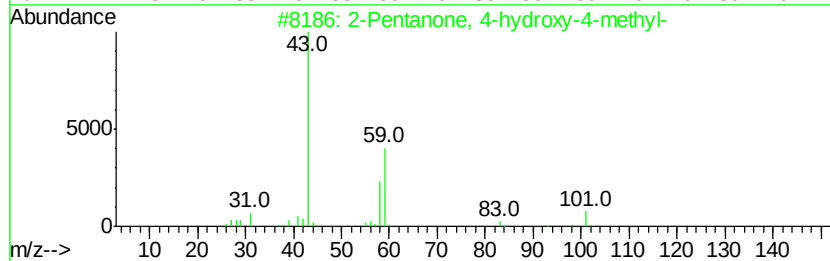
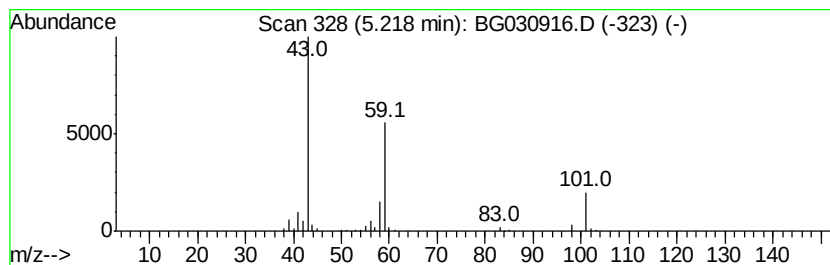
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	12.18 ng	172683	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	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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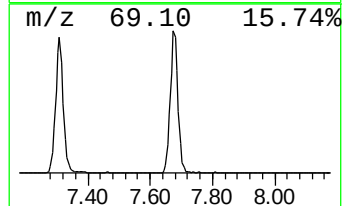
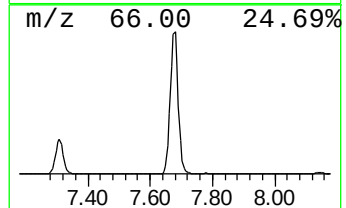
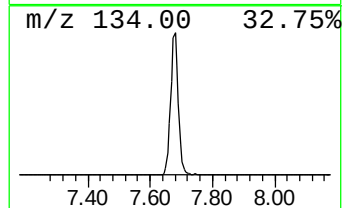
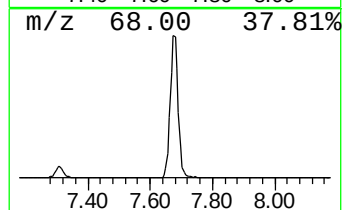
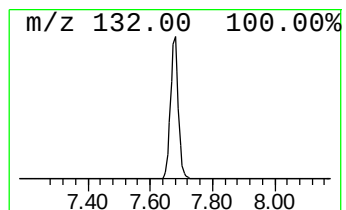
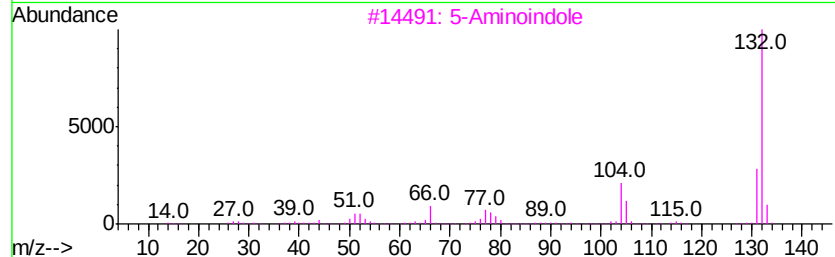
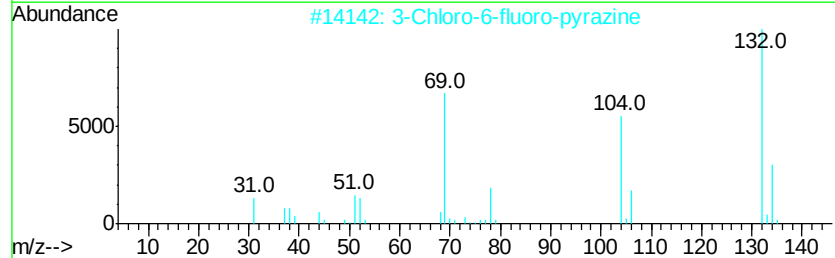
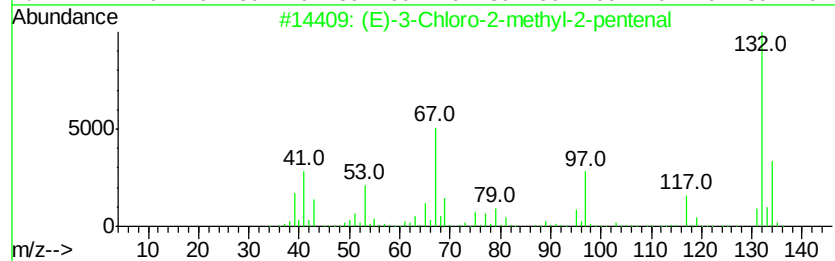
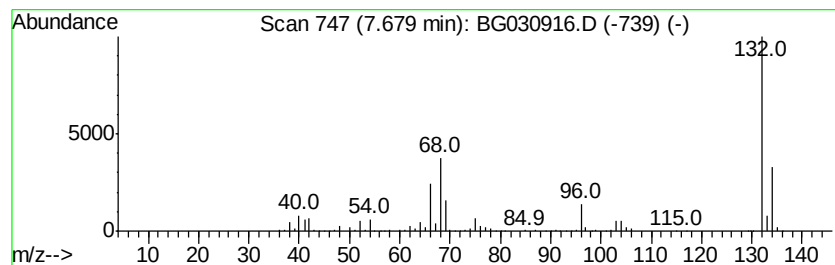
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	85.44 ng	1211420	1,4-Dichlorobenzene-d4	8.14

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	35
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	12.2	ng	172683	1	8.14	283560	20.0
unknown7.68	7.68	85.4	ng	1211420	1	8.14	283560	20.0