

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111017\
 Data File : BG030920.D
 Acq On : 10 Nov 2017 23:37
 Operator : SJ/JU
 Sample : PB104039BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104039BL

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	338	rBV	117088	189409	4.17%	1.037%
2	5.703	404	411	423	rBV	768568	1245786	27.43%	6.818%
3	7.307	676	684	696	rBV	763495	1322864	29.13%	7.240%
4	7.678	739	747	757	rBV	758428	1327755	29.24%	7.267%
5	8.142	820	826	834	rVB	124778	209191	4.61%	1.145%
6	8.471	874	882	892	rVB	584956	1049136	23.10%	5.742%
7	9.311	1014	1025	1037	rBV	414437	782889	17.24%	4.285%
8	10.962	1297	1306	1325	rBV	155467	305301	6.72%	1.671%
9	13.389	1710	1719	1730	rBV	1424157	2230582	49.12%	12.208%
10	14.763	1946	1953	1963	rBV2	275759	477026	10.50%	2.611%
11	16.244	2199	2205	2229	rBV	1192771	1956702	43.09%	10.709%
12	17.501	2412	2419	2430	rBV2	445559	707999	15.59%	3.875%
13	20.092	2854	2860	2870	rBV	3445180	4541076	100.00%	24.853%
14	21.773	3138	3146	3157	rBV2	560650	972845	21.42%	5.324%
15	25.075	3697	3708	3720	rBV2	316468	953004	20.99%	5.216%

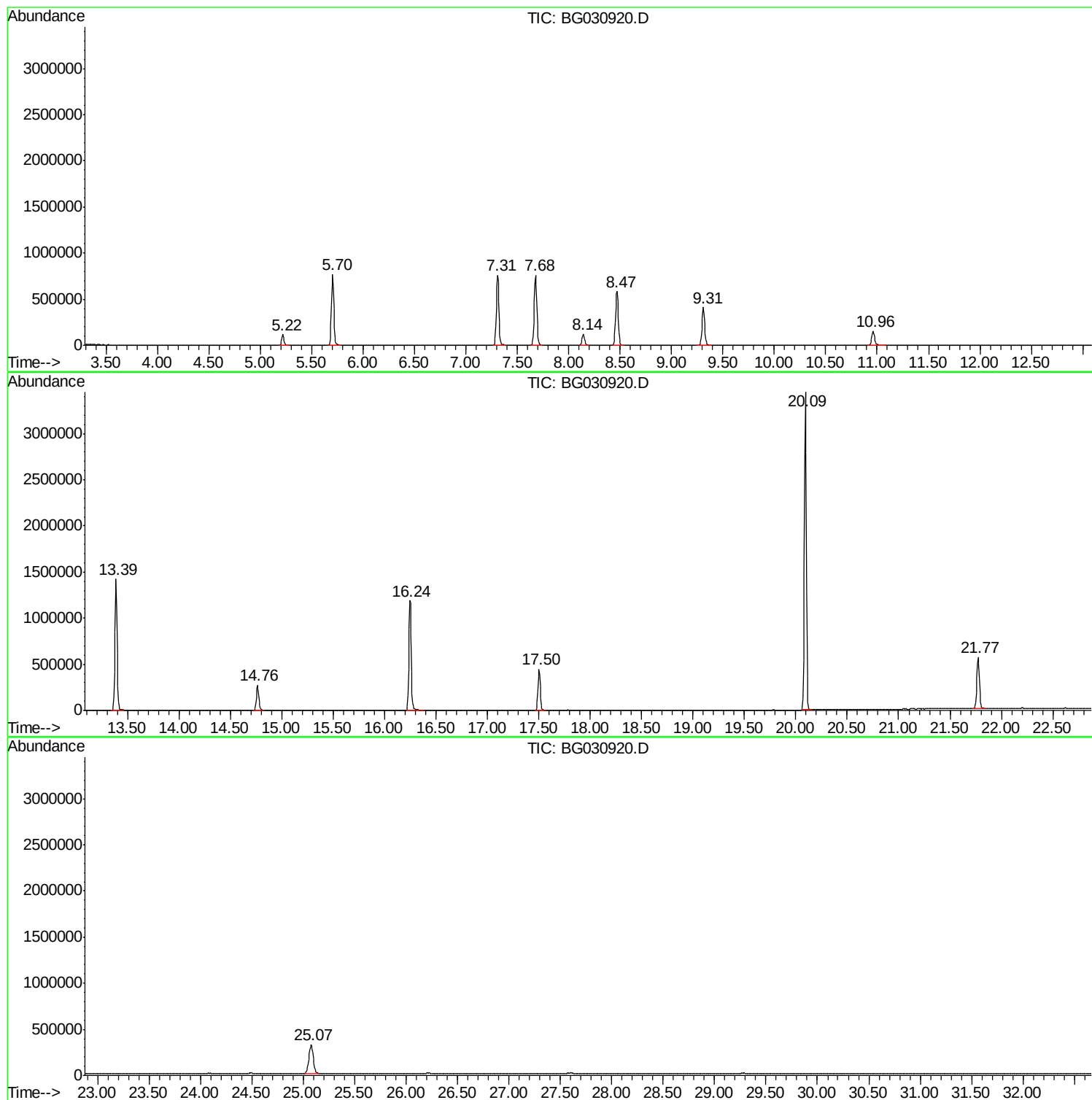
Sum of corrected areas: 18271565

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



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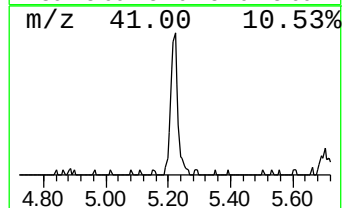
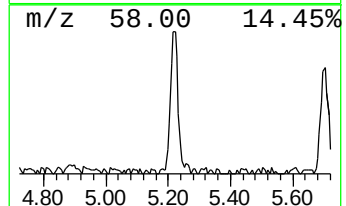
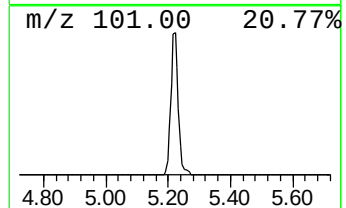
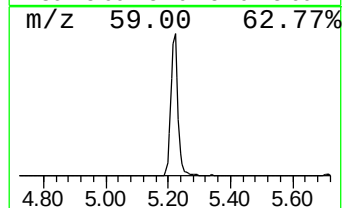
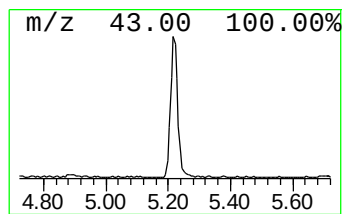
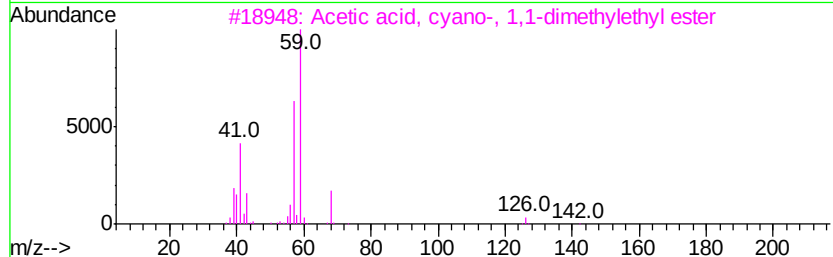
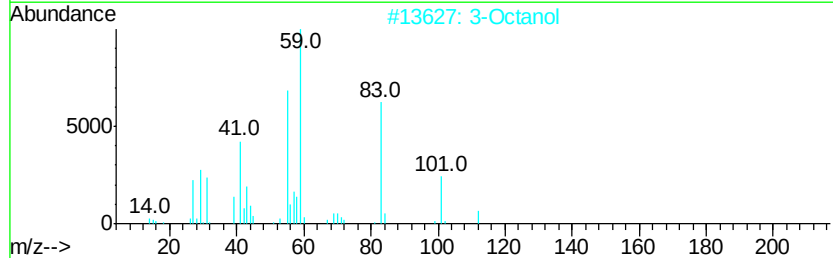
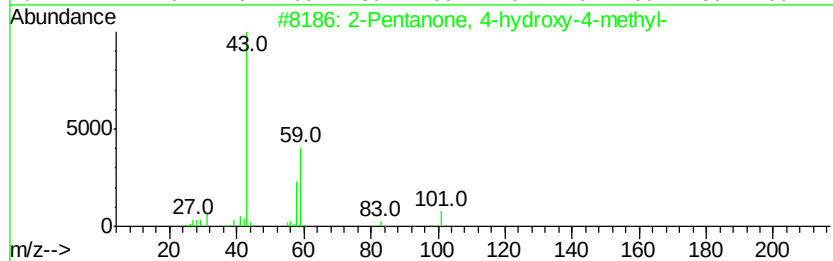
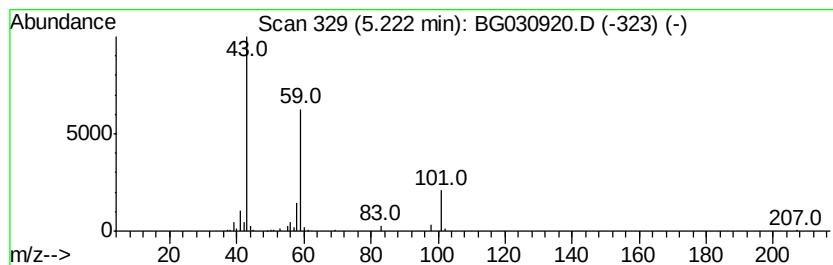
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	18.11 ng	189409	1,4-Dichlorobenzene-d4	8.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			3-Octanol	130	C8H18O	000589-98-0	36
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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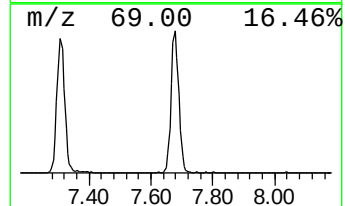
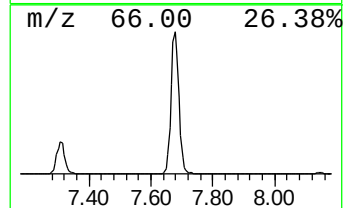
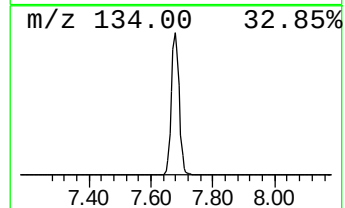
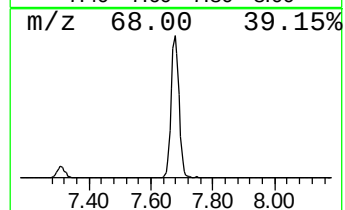
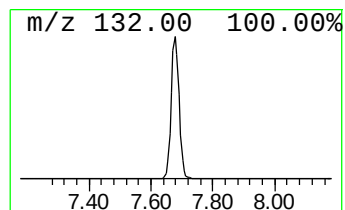
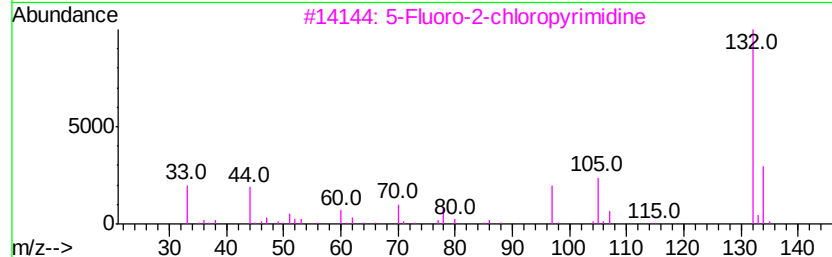
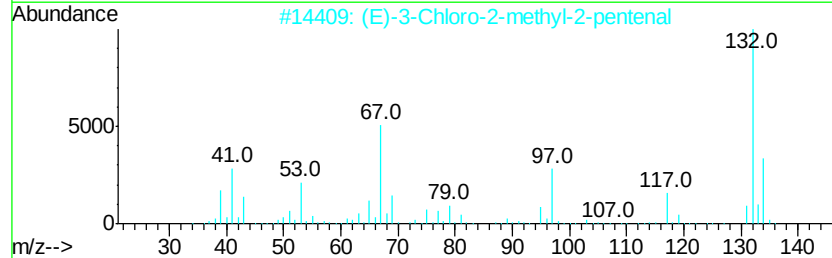
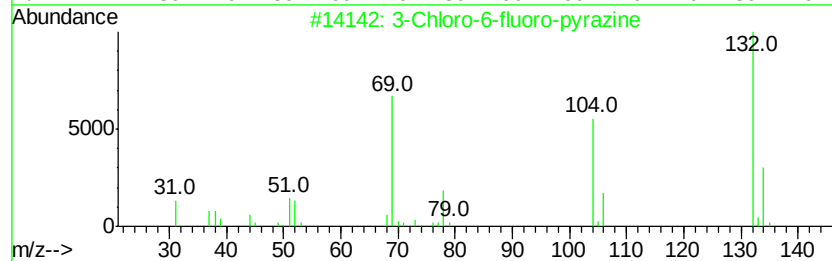
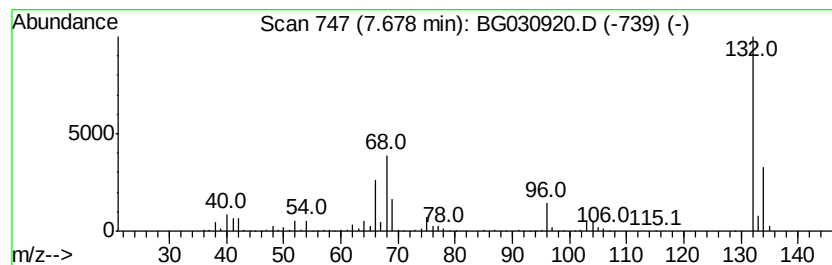
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	126.94 ng	1327760	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	18.1	ng	189409	1	8.14	209191	20.0
unknown7.68	7.68	126.9	ng	1327760	1	8.14	209191	20.0