

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031301.D
 Acq On : 30 Nov 2017 23:30
 Operator : SJ/JU
 Sample : I6635-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112817-JETT-E-D-B

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	5.209	321	327	335	rBV	65131	106606	1.61%	0.470%
2	5.690	402	409	429	rBV	319811	573339	8.67%	2.526%
3	7.294	675	682	694	rBV	182714	386594	5.85%	1.703%
4	7.664	737	745	758	rBV	634670	1134498	17.16%	4.998%
5	8.129	818	824	832	rBV	206309	359586	5.44%	1.584%
6	8.452	871	879	892	rBV	906104	1682237	25.45%	7.412%
7	9.298	1014	1023	1037	rBV	648988	1293365	19.56%	5.698%
8	10.949	1295	1304	1322	rBV	234621	512172	7.75%	2.257%
9	13.375	1709	1717	1731	rBV	2247363	3632060	54.94%	16.002%
10	14.756	1944	1952	1963	rBV3	435668	812675	12.29%	3.580%
11	16.237	2195	2204	2217	rBV	1116120	1834743	27.75%	8.084%
12	17.494	2411	2418	2431	rBV2	667462	1088370	16.46%	4.795%
13	20.085	2852	2859	2872	rBV	5034055	6610845	100.00%	29.126%
14	21.766	3138	3145	3161	rBV2	793626	1366844	20.68%	6.022%
15	25.062	3695	3706	3722	rBV2	435263	1303387	19.72%	5.742%

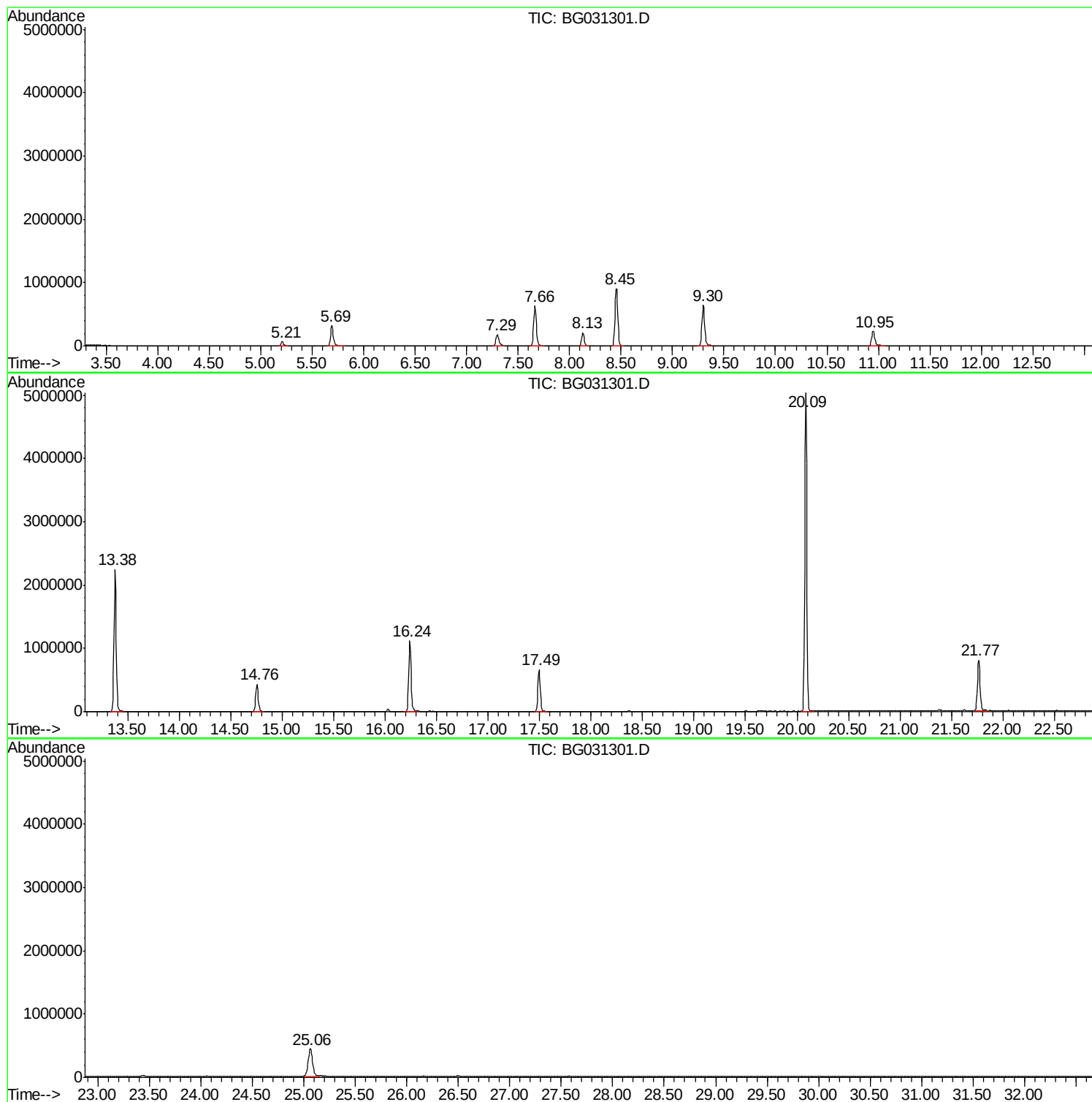
Sum of corrected areas: 22697321

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.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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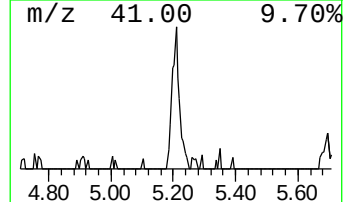
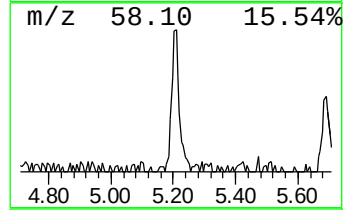
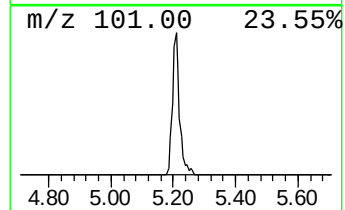
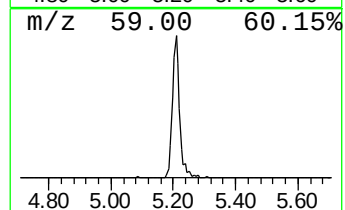
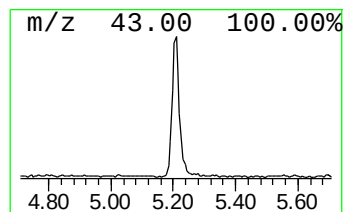
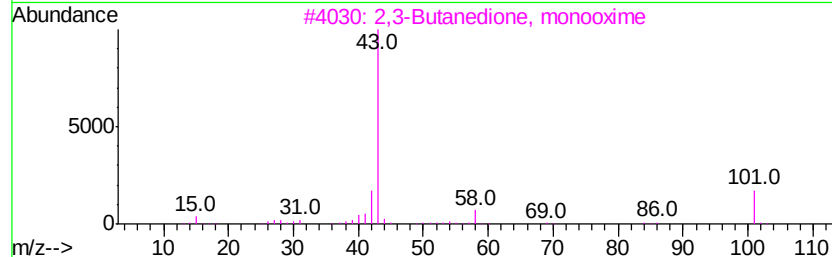
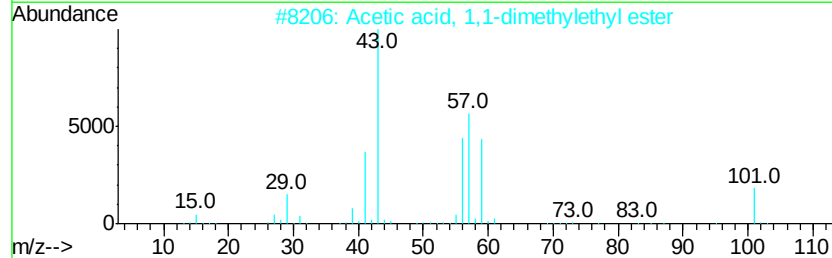
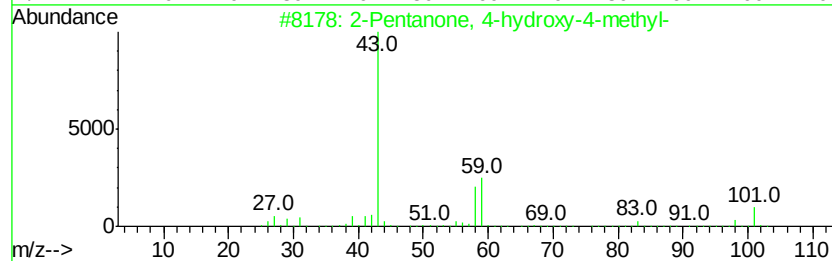
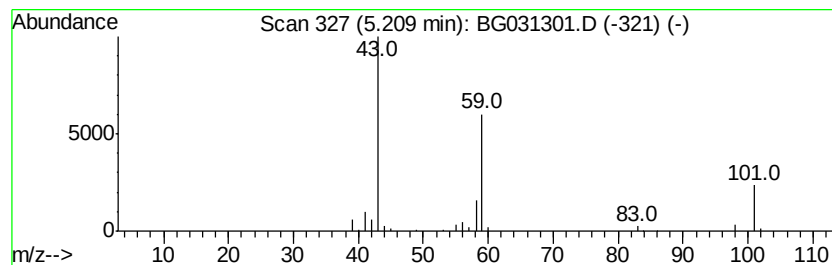
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.93 ng	106606	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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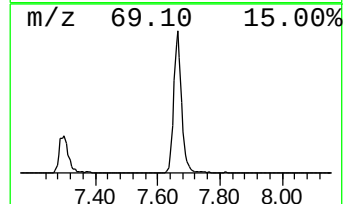
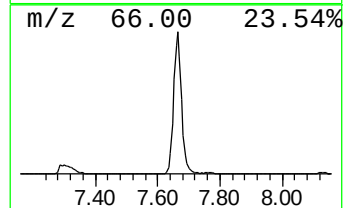
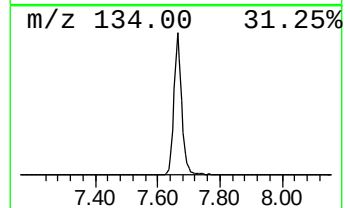
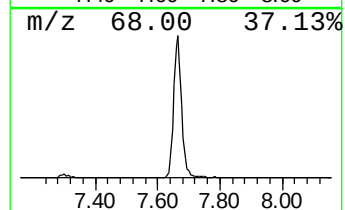
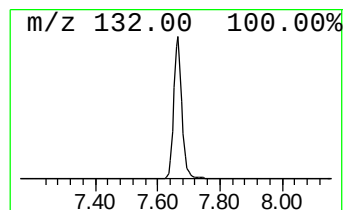
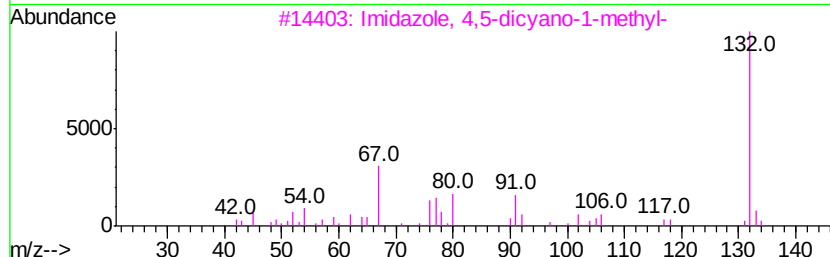
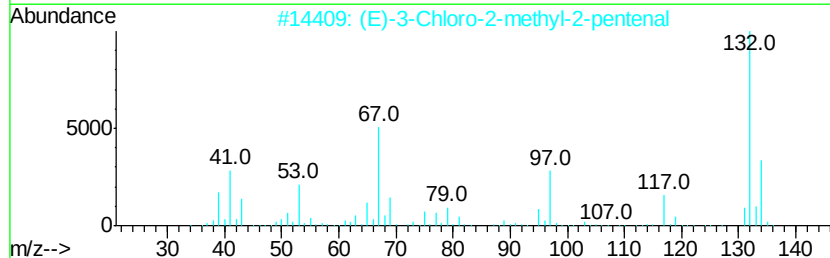
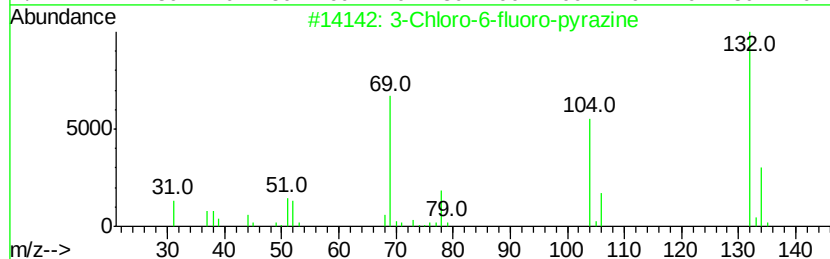
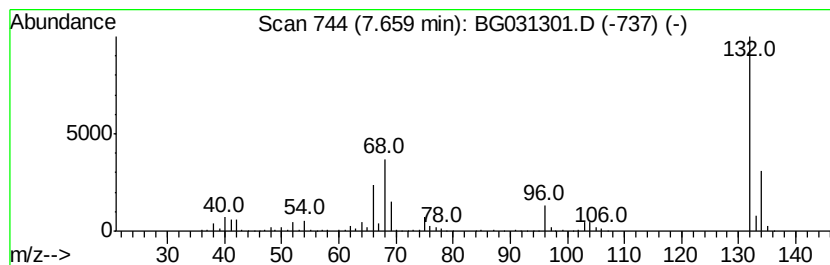
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	63.10 ng	1134500	1,4-Dichlorobenzene-d4	8.13

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	16
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.9	ng	106606	1	8.13	359586	20.0
unknown7.66	7.66	63.1	ng	1134500	1	8.13	359586	20.0