

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112117\
 Data File : BG031174.D
 Acq On : 22 Nov 2017 5:21
 Operator : SJ/JU
 Sample : I6539-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 111517-TIDO-F-D-S

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	5.211	322	327	337	rBV	91896	144918	2.56%	0.722%
2	5.699	401	410	422	rBV	272812	476072	8.39%	2.373%
3	7.303	676	683	700	rBV	164086	345161	6.09%	1.721%
4	7.673	736	746	757	rBV	519497	939426	16.56%	4.684%
5	8.137	817	825	834	rBV	183956	320017	5.64%	1.595%
6	8.460	873	880	892	rBV	743022	1328259	23.42%	6.622%
7	9.306	1014	1024	1037	rBV	548614	1061247	18.71%	5.291%
8	10.952	1296	1304	1321	rBV	228303	473254	8.34%	2.359%
9	13.378	1710	1717	1737	rBV	1818836	3051484	53.80%	15.213%
10	14.759	1944	1952	1962	rBV2	429252	763792	13.47%	3.808%
11	16.240	2198	2204	2224	rBV	913996	1569477	27.67%	7.825%
12	17.497	2411	2418	2431	rBV	668795	1041630	18.37%	5.193%
13	18.361	2560	2565	2575	rBV2	34094	66046	1.16%	0.329%
14	19.624	2776	2780	2789	rBV3	41592	70099	1.24%	0.349%
15	20.088	2853	2859	2873	rBV	4165926	5671492	100.00%	28.276%
16	21.381	3075	3079	3087	rBV4	38412	80311	1.42%	0.400%
17	21.768	3138	3145	3158	rBV	789445	1375459	24.25%	6.857%
18	25.070	3696	3707	3719	rBV2	424658	1279717	22.56%	6.380%

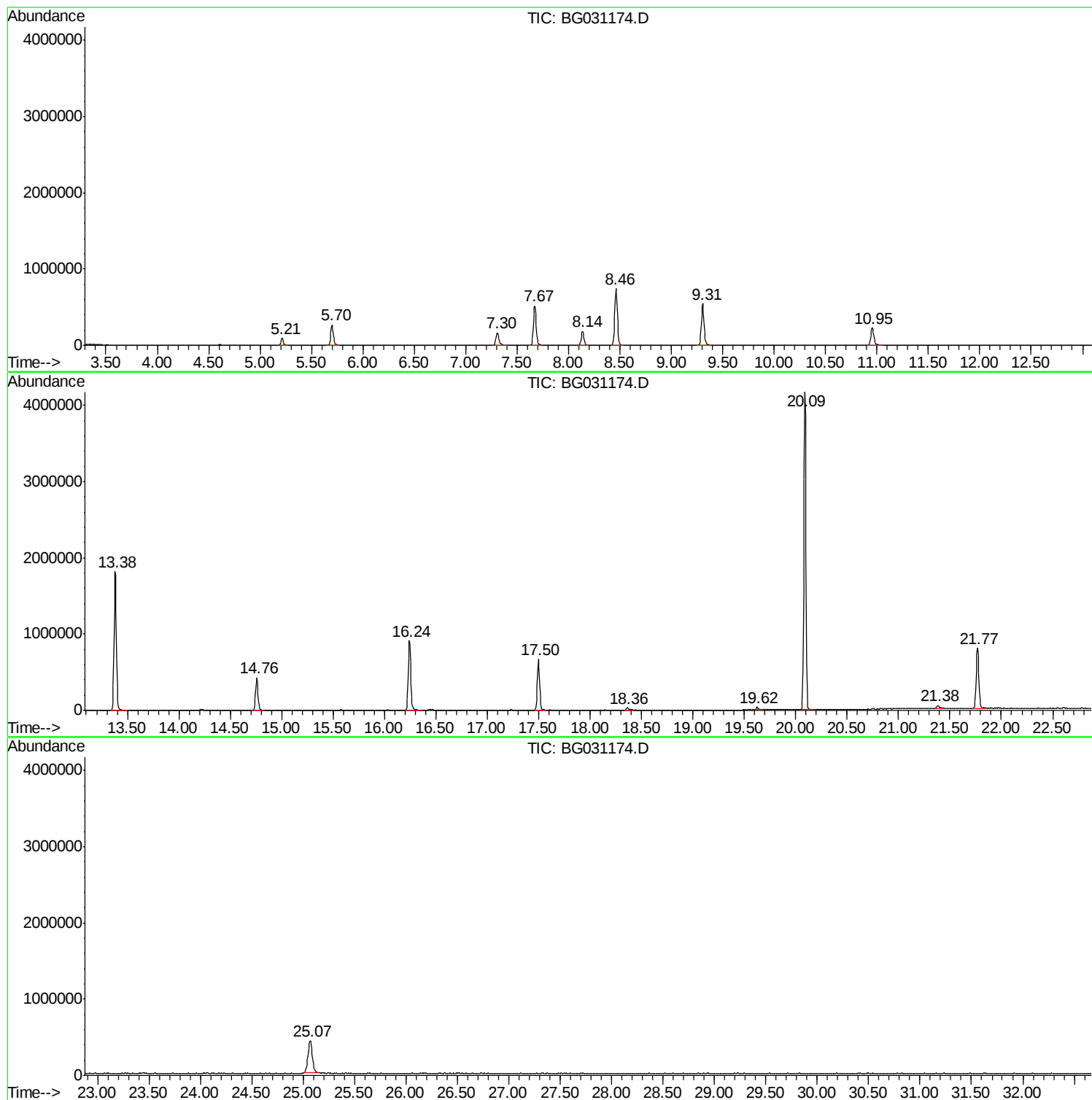
Sum of corrected areas: 20057861

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111517-TIDO-F-D-S

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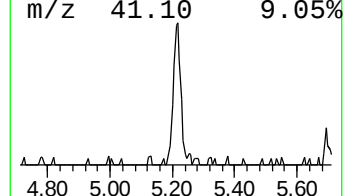
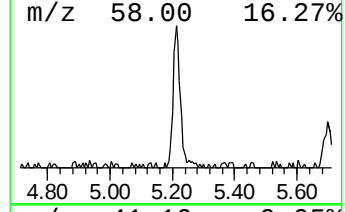
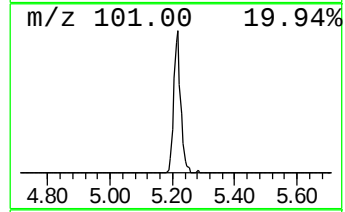
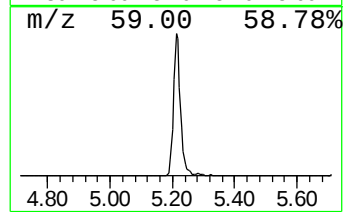
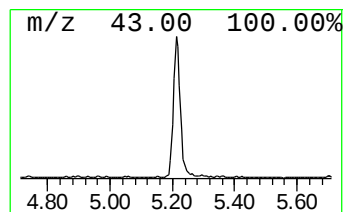
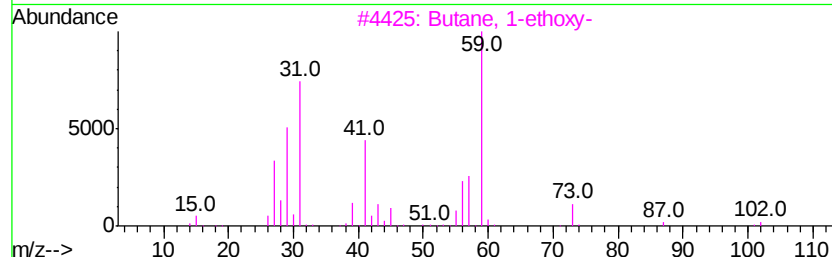
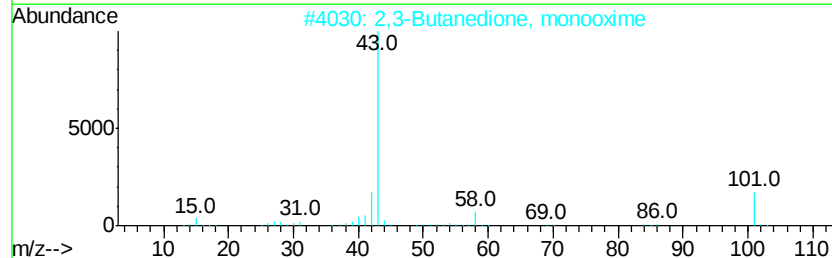
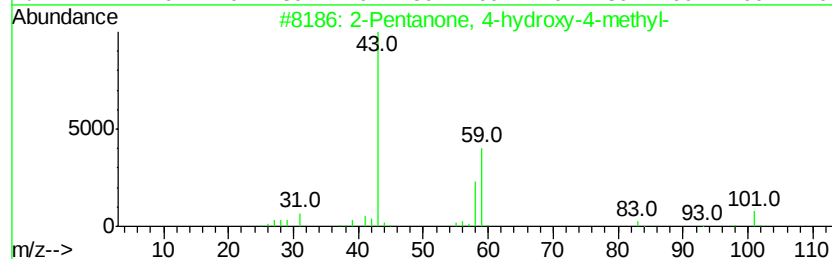
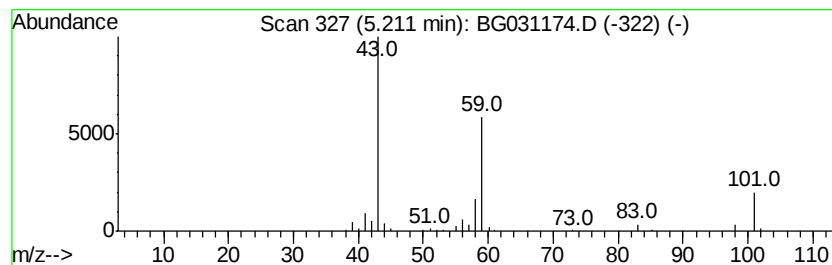
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	9.06 ng	144918	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	9
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetone	58	C3H6O	000067-64-1	9



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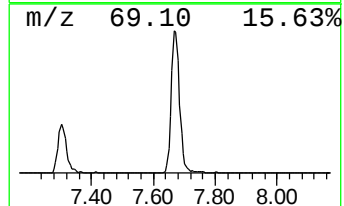
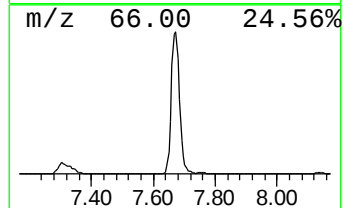
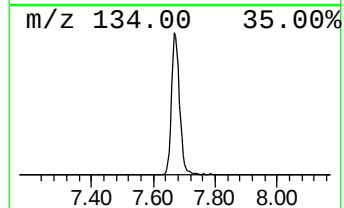
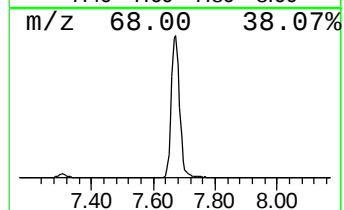
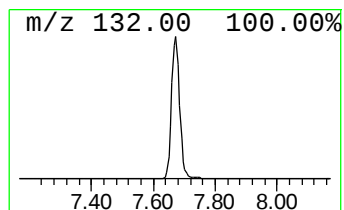
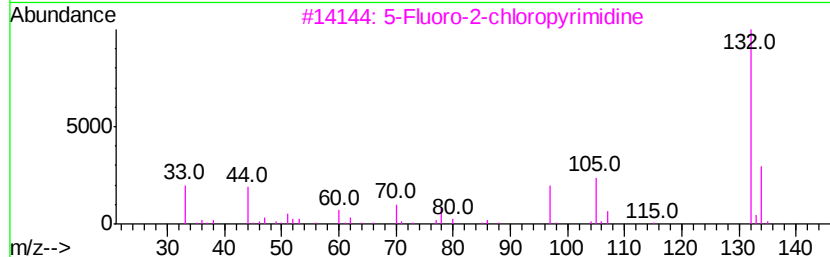
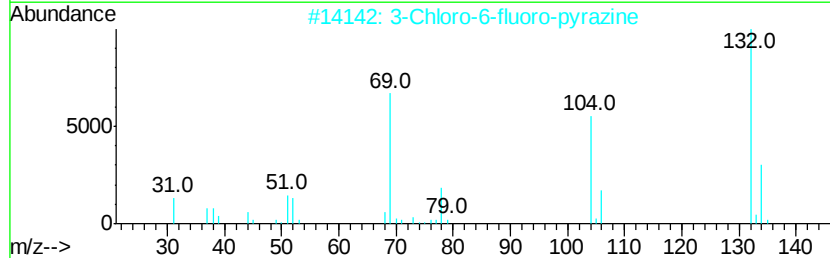
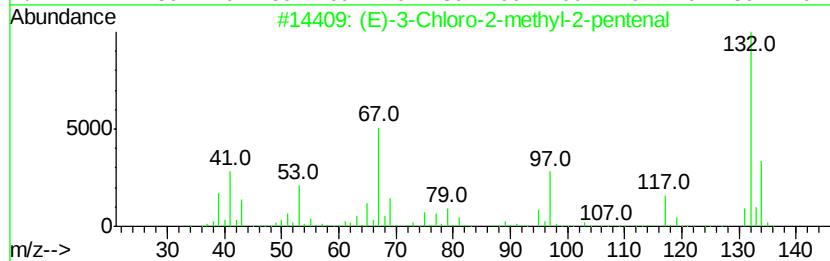
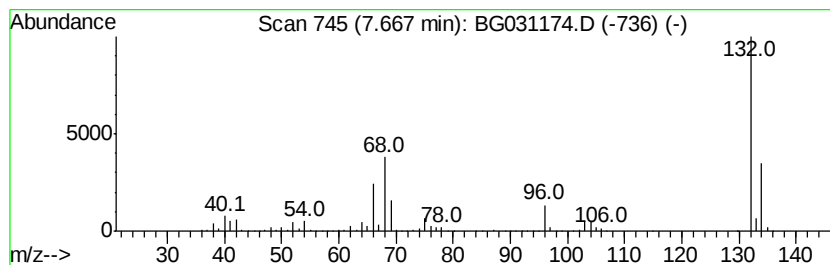
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Peak Number 2 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	58.71 ng	939426	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	25
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	9.1	ng	144918	1	8.14	320017	20.0
unknown7.67	7.67	58.7	ng	939426	1	8.14	320017	20.0