

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112217\
 Data File : BG031192.D
 Acq On : 22 Nov 2017 16:48
 Operator : SJ/JU
 Sample : I6525-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-24

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	336	rBV	38178	64609	1.17%	0.303%
2	5.693	403	409	422	rBV	444683	792039	14.39%	3.720%
3	7.302	676	683	695	rBV	295098	545119	9.90%	2.560%
4	7.667	733	745	764	rBV	777804	1442787	26.21%	6.776%
5	8.137	815	825	834	rBV	177500	304739	5.54%	1.431%
6	8.460	872	880	892	rBV	751970	1359043	24.69%	6.383%
7	9.306	1016	1024	1038	rBV	533512	1051550	19.11%	4.938%
8	10.951	1297	1304	1323	rBV	212448	460376	8.36%	2.162%
9	13.378	1710	1717	1739	rBV	1871199	3103999	56.40%	14.577%
10	14.758	1945	1952	1964	rBV	411151	711360	12.92%	3.341%
11	16.239	2197	2204	2222	rBV	1470262	2392282	43.47%	11.235%
12	17.496	2411	2418	2434	rBV2	619947	996913	18.11%	4.682%
13	20.087	2852	2859	2870	rBV	4067595	5503855	100.00%	25.848%
14	21.768	3138	3145	3155	rBV	776193	1326337	24.10%	6.229%
15	25.064	3695	3706	3719	rBV2	404397	1238219	22.50%	5.815%

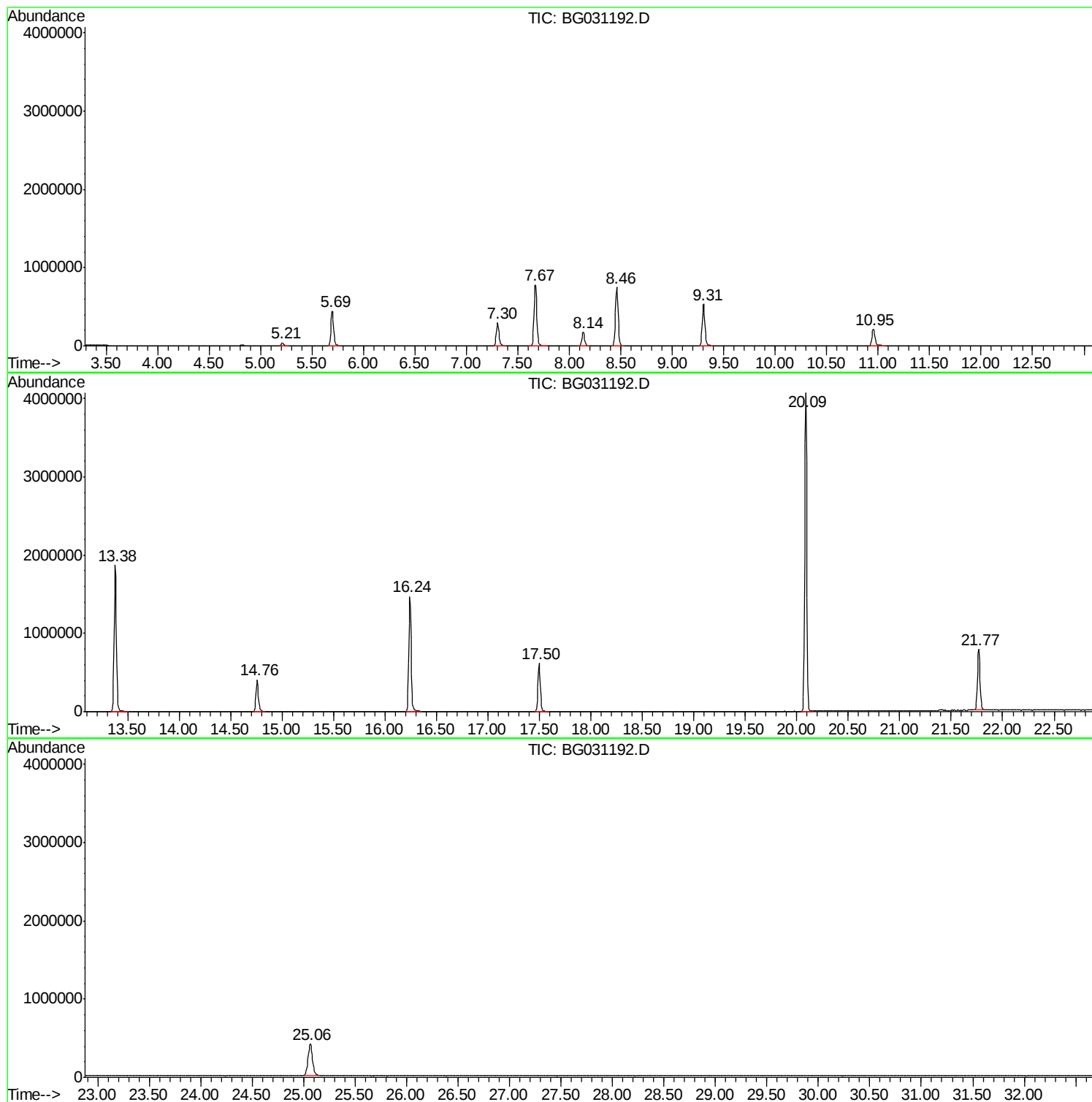
Sum of corrected areas: 21293227

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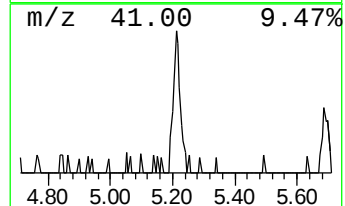
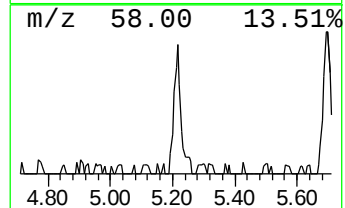
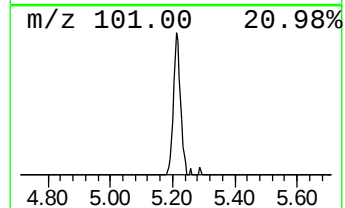
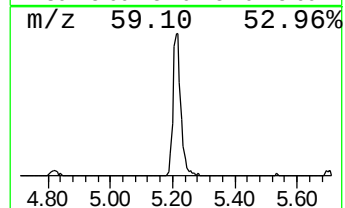
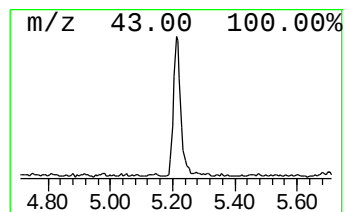
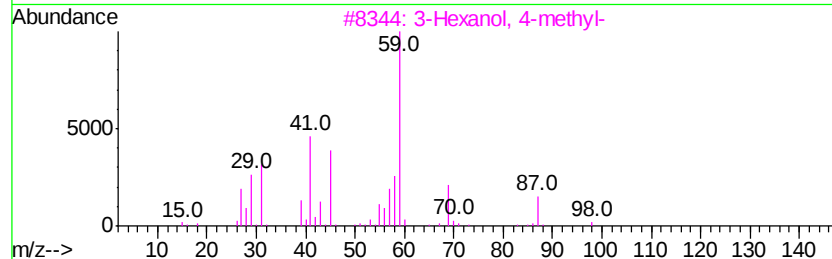
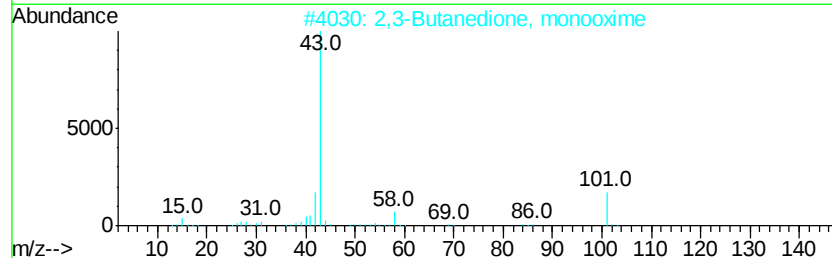
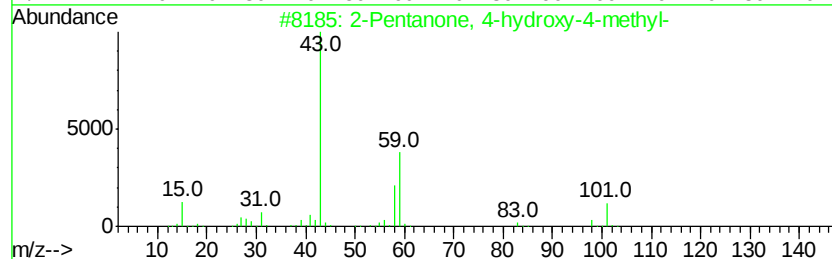
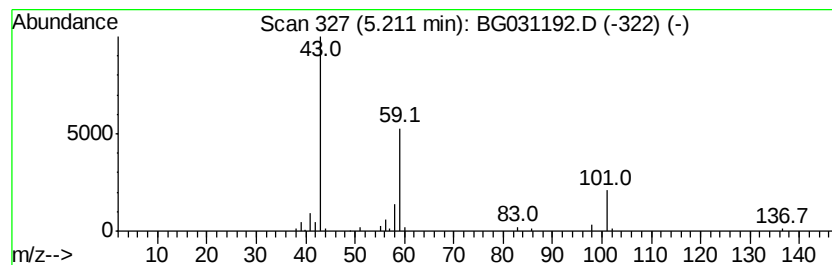
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	4.24 ng	64609	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	64		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	43		
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25		
4	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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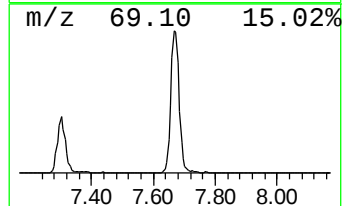
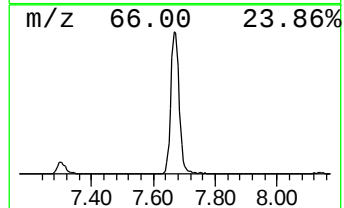
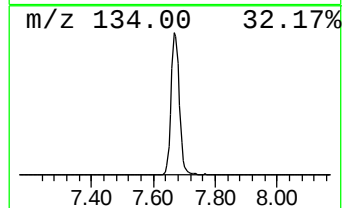
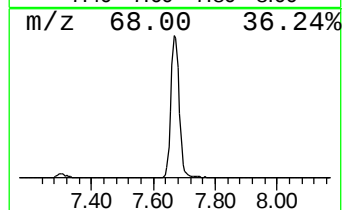
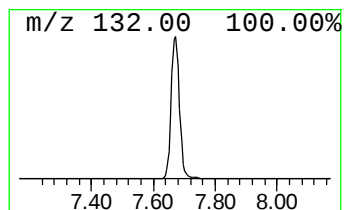
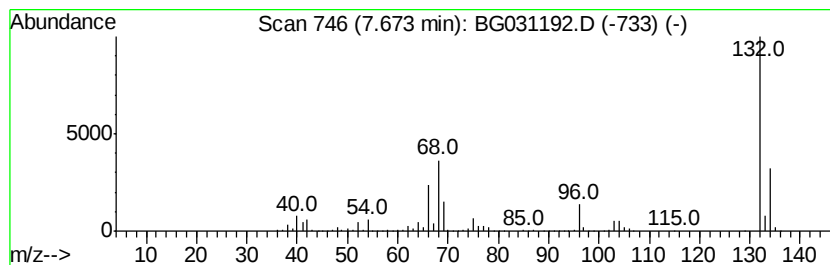
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Peak Number 2 unknown7.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	94.69 ng	1442790	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	4.2	ng	64609	1	8.14	304739	20.0
unknown7.67	7.67	94.7	ng	1442790	1	8.14	304739	20.0