

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112817\
 Data File : BG031260.D
 Acq On : 28 Nov 2017 20:29
 Operator : SJ/JU
 Sample : I6610-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-2(0-2)-20171127

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	5.203	320	326	337	rVB	56093	92430	3.01%	0.716%
2	5.691	402	409	421	rBV	407856	703583	22.93%	5.447%
3	7.295	673	682	701	rBV	411076	788662	25.71%	6.105%
4	7.665	736	745	761	rBV	427473	776704	25.32%	6.013%
5	8.129	817	824	832	rBV	123177	218973	7.14%	1.695%
6	8.452	872	879	892	rBV	330266	615121	20.05%	4.762%
7	9.298	1016	1023	1034	rBV	219832	450526	14.68%	3.488%
8	10.949	1296	1304	1326	rBV	133866	298253	9.72%	2.309%
9	13.376	1710	1717	1733	rBV	796490	1347011	43.90%	10.428%
10	14.210	1852	1859	1867	rBV	35355	78064	2.54%	0.604%
11	14.756	1945	1952	1964	rVB2	269396	469143	15.29%	3.632%
12	15.573	2085	2091	2097	rBV2	25882	35666	1.16%	0.276%
13	16.237	2198	2204	2222	rVB	672323	1130444	36.84%	8.751%
14	16.431	2232	2237	2247	rVB2	21383	37555	1.22%	0.291%
15	17.494	2411	2418	2429	rBV2	442109	747616	24.37%	5.788%
16	20.086	2852	2859	2868	rBV	2303696	3068118	100.00%	23.752%
17	21.378	3073	3079	3088	rBV	22387	61066	1.99%	0.473%
18	21.766	3138	3145	3156	rBV	585862	1016178	33.12%	7.867%
19	25.062	3695	3706	3719	rBV3	314153	982263	32.02%	7.604%

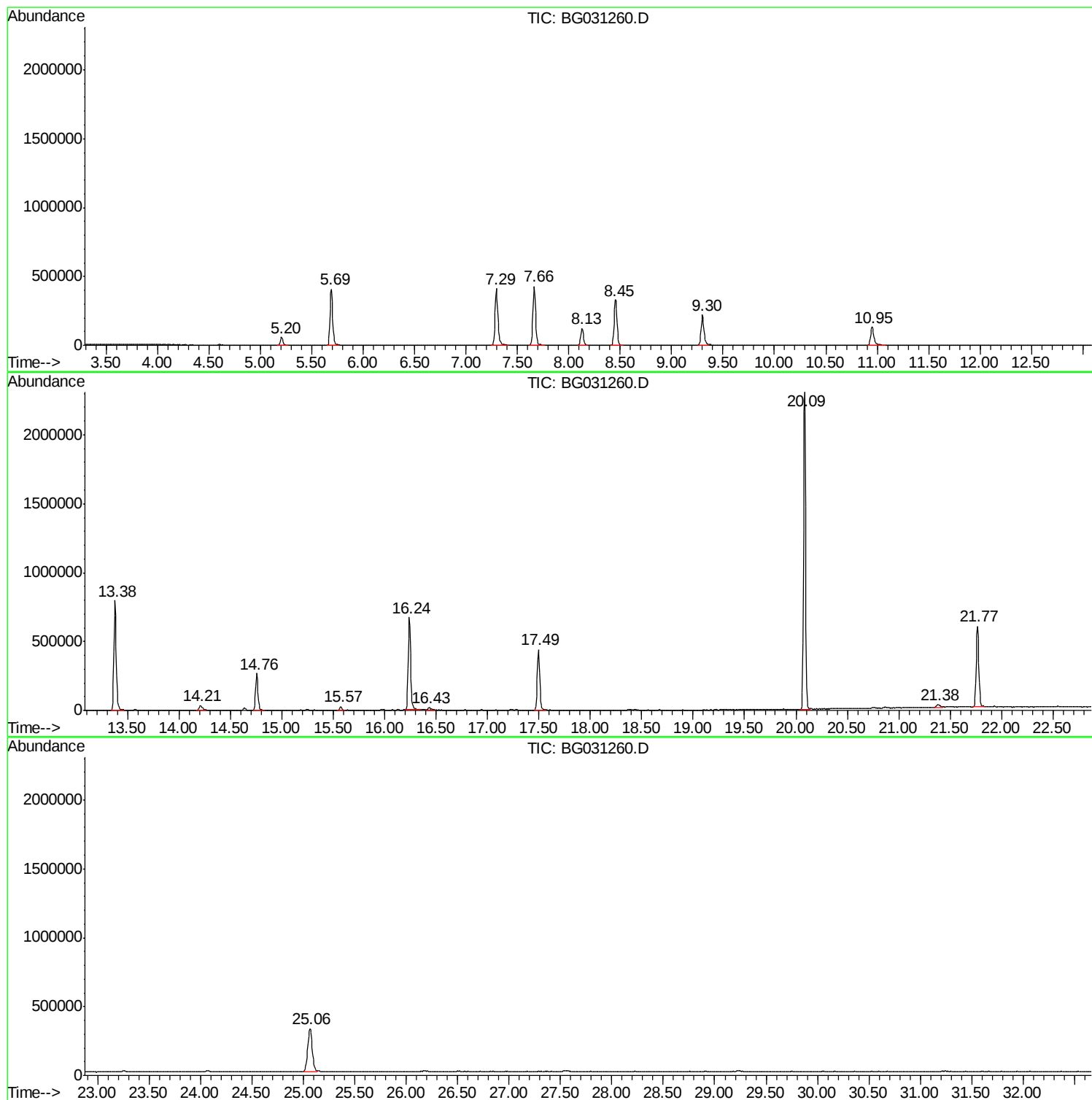
Sum of corrected areas: 12917376

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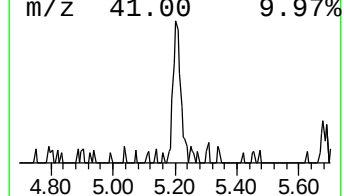
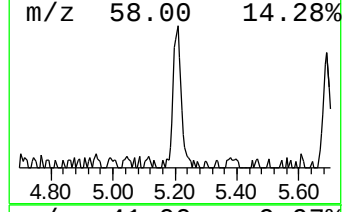
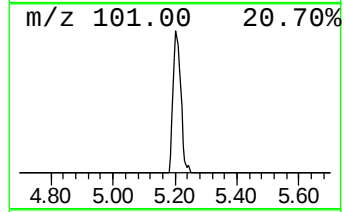
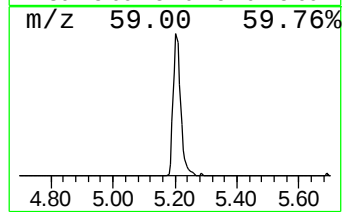
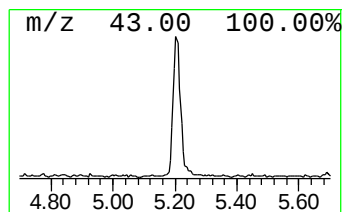
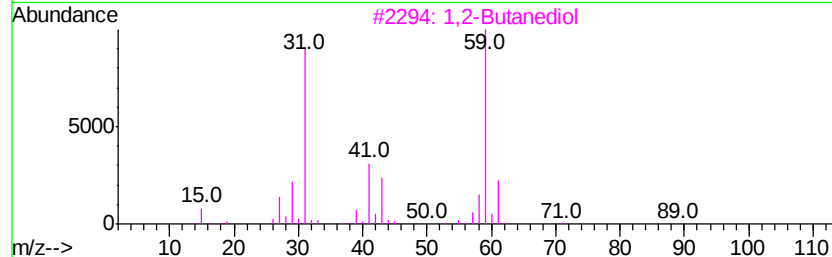
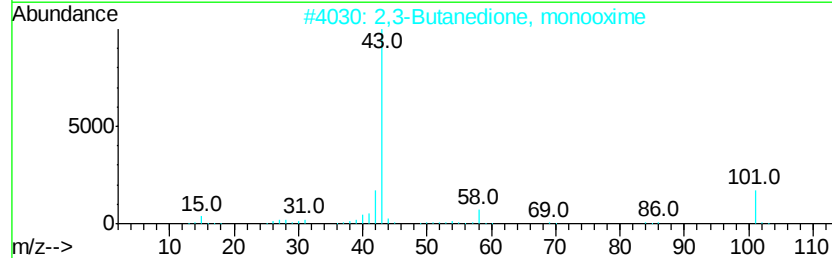
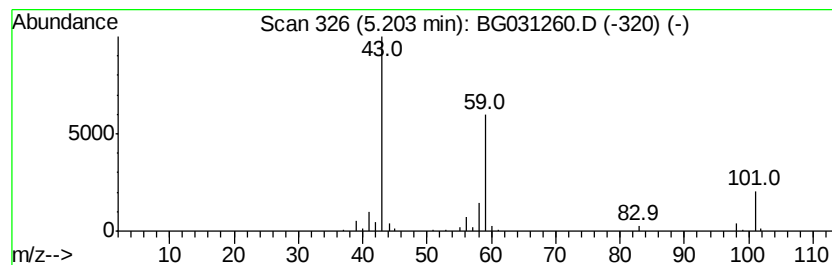
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	8.44 ng	92430	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	72
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3		1,2-Butanediol	90	C4H10O2	000584-03-2	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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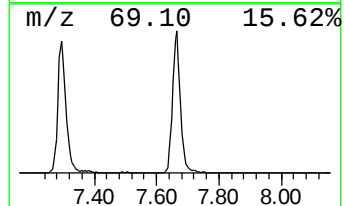
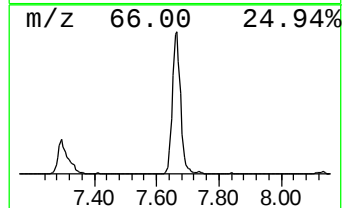
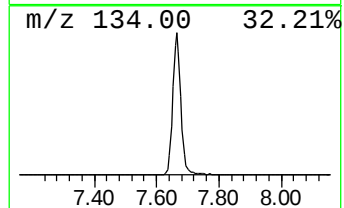
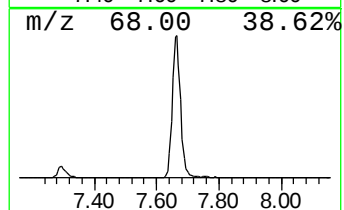
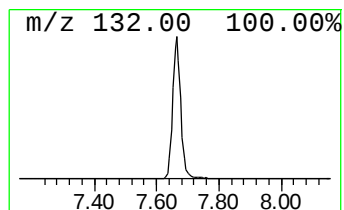
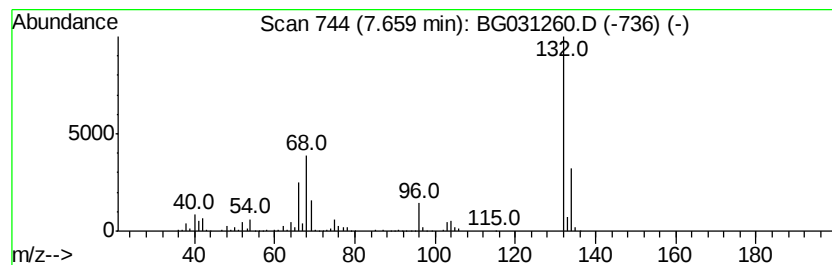
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Peak Number 2 unknown7.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	70.94 ng	776704	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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	8.4	ng	92430	1	8.13	218973	20.0
unknown7.66	7.66	70.9	ng	776704	1	8.13	218973	20.0