

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112217\
 Data File : BG031198.D
 Acq On : 22 Nov 2017 20:33
 Operator : SJ/JU
 Sample : I6525-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUP111617

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	5.214	322	328	337	rVB	36681	60209	1.14%	0.286%
2	5.696	404	410	426	rBV	432198	744846	14.08%	3.534%
3	7.300	676	683	698	rBV	262934	495543	9.37%	2.351%
4	7.670	738	746	757	rBV	781143	1424235	26.92%	6.757%
5	8.134	816	825	834	rBV	177958	311077	5.88%	1.476%
6	8.458	872	880	893	rBV	739271	1343229	25.39%	6.373%
7	9.304	1015	1024	1038	rBV	540868	1055749	19.95%	5.009%
8	10.949	1295	1304	1324	rBV	223972	477604	9.03%	2.266%
9	13.381	1710	1718	1728	rBV	1837357	3066032	57.95%	14.546%
10	14.756	1942	1952	1968	rBV2	422279	743330	14.05%	3.527%
11	16.237	2198	2204	2228	rBV	1391343	2378554	44.95%	11.285%
12	17.494	2411	2418	2433	rBV2	608402	1029503	19.46%	4.884%
13	20.085	2853	2859	2873	rBV	3936894	5291094	100.00%	25.103%
14	21.766	3138	3145	3159	rBV	786384	1360571	25.71%	6.455%
15	25.068	3695	3707	3722	rBV2	428653	1295962	24.49%	6.149%

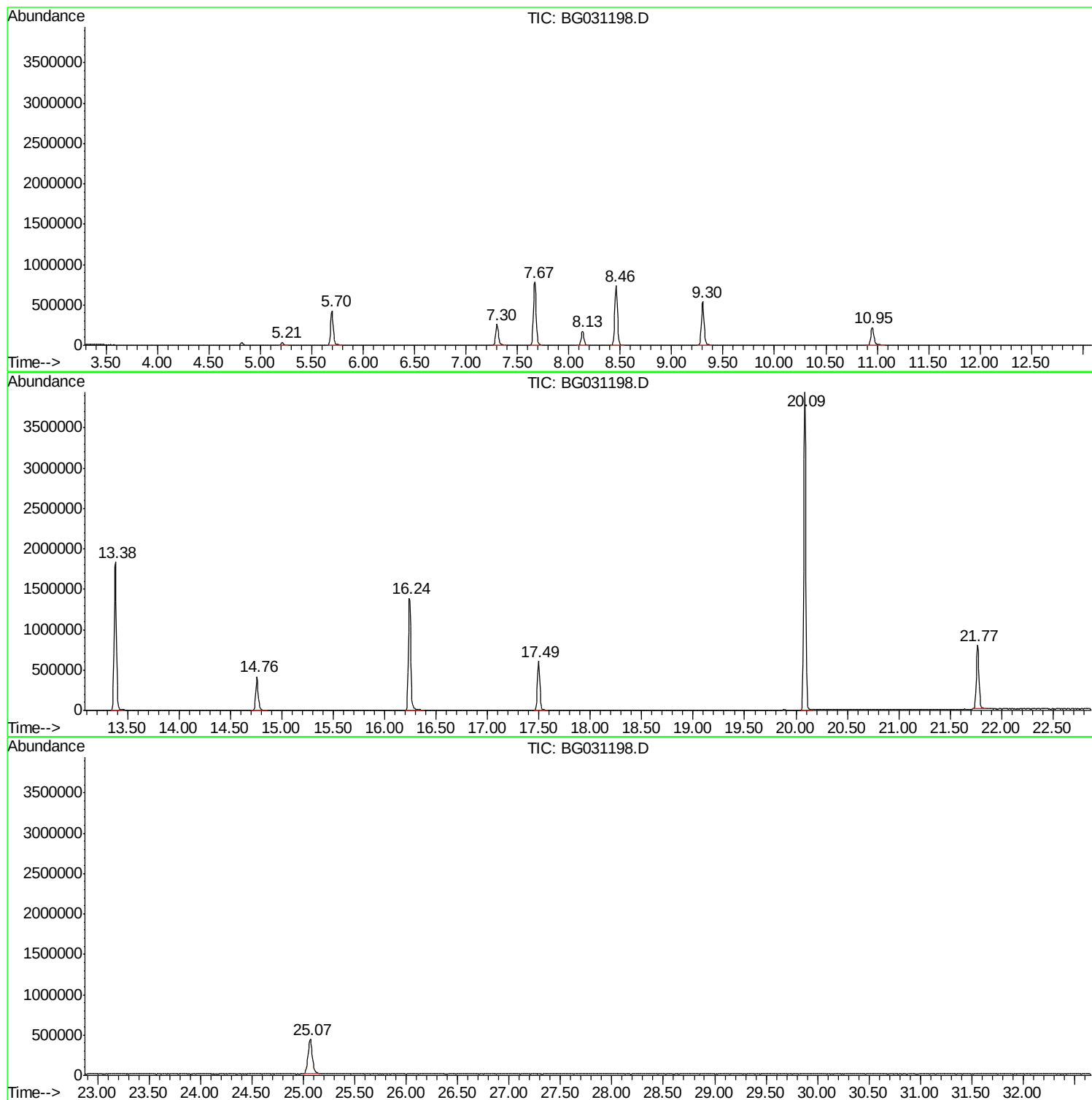
Sum of corrected areas: 21077538

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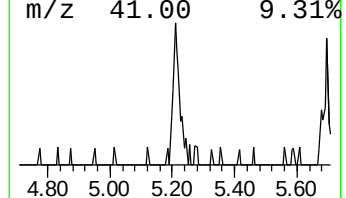
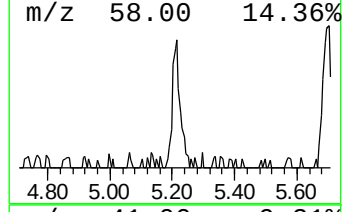
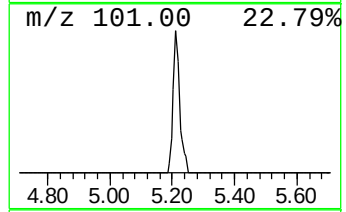
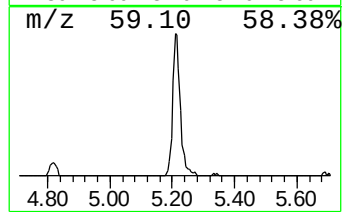
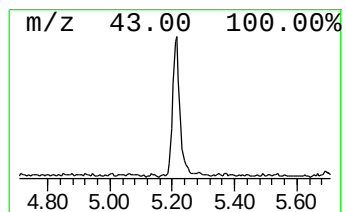
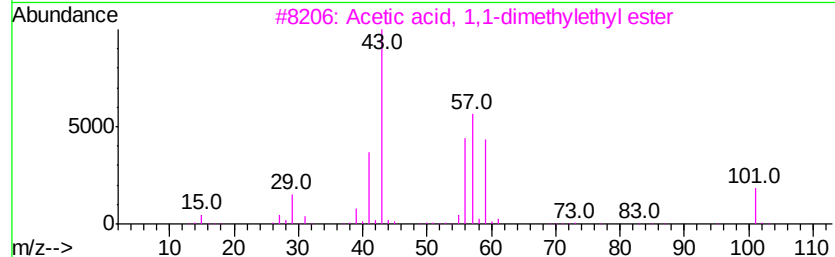
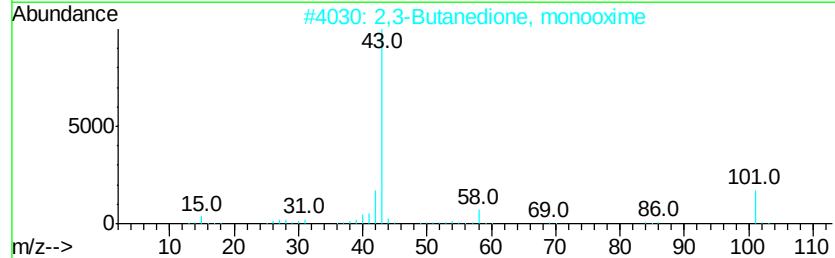
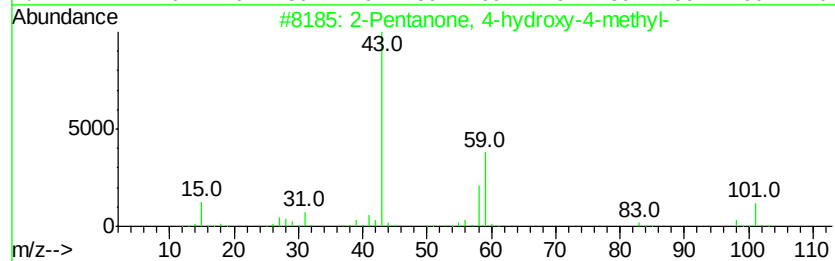
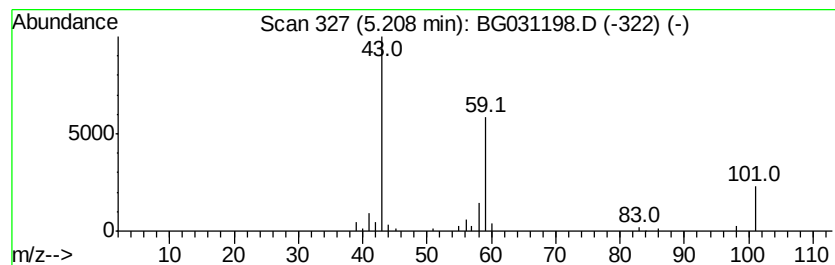
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	3.87 ng	60209	1,4-Dichlorobenzene-d4	8.13

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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25



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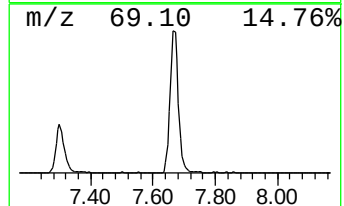
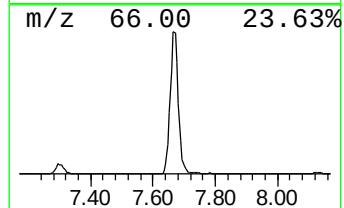
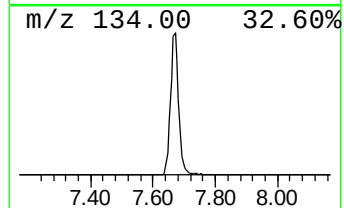
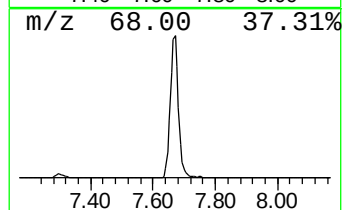
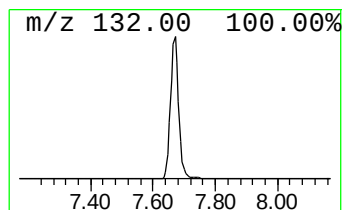
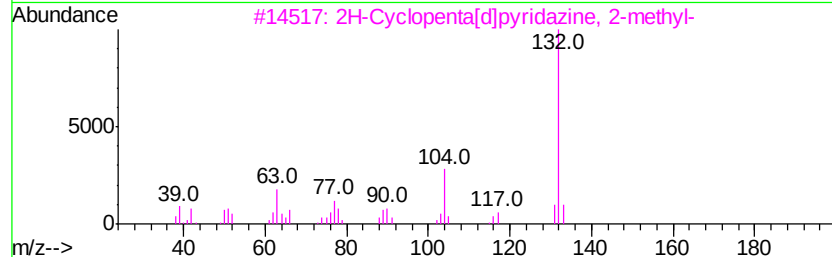
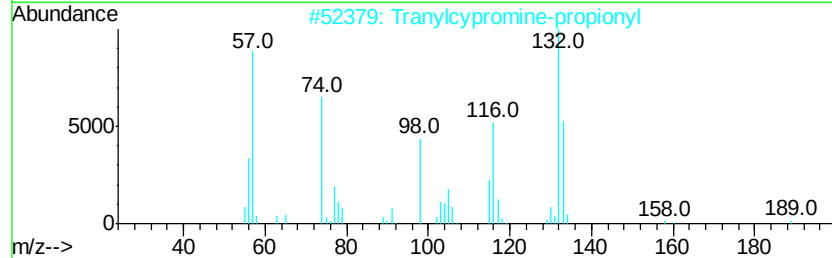
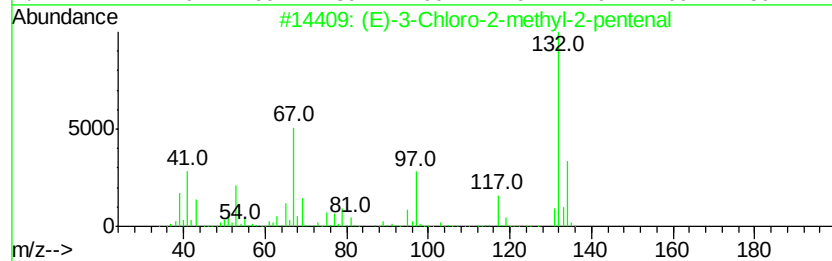
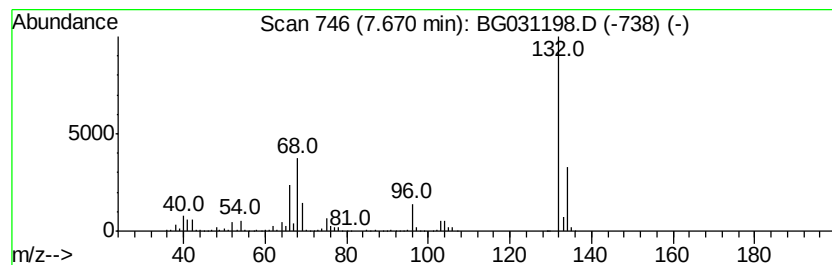
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Peak Number 2 unknown7.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	91.57 ng	1424240	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	3.9	ng	60209	1	8.13	311077	20.0
unknown7.67	7.67	91.6	ng	1424240	1	8.13	311077	20.0