

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
 Data File : BG040407.D
 Acq On : 9 Apr 2019 20:26
 Operator : JU/SJ
 Sample : PB118717BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118717BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.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	5.138	343	349	357	rBV	81748	129155	3.32%	0.734%
2	5.602	422	428	444	rBV	658912	1062283	27.32%	6.037%
3	7.200	691	700	714	rBV	627346	1103393	28.38%	6.271%
4	7.577	757	764	776	rBV	621748	1087386	27.97%	6.180%
5	8.047	837	844	852	rBV	218003	362032	9.31%	2.058%
6	8.370	891	899	908	rBV	704439	1242228	31.95%	7.060%
7	9.222	1036	1044	1056	rBV	541695	1002092	25.77%	5.695%
8	10.861	1313	1323	1337	rBV	248394	484341	12.46%	2.753%
9	13.299	1731	1738	1756	rBV	1417130	2308919	59.39%	13.122%
10	14.680	1966	1973	1986	rBV2	395881	675699	17.38%	3.840%
11	16.166	2219	2226	2248	rBV	449683	931169	23.95%	5.292%
12	17.424	2433	2440	2452	rBV2	573783	949071	24.41%	5.394%
13	20.021	2876	2882	2901	rBV	2790461	3888036	100.00%	22.097%
14	21.695	3160	3167	3182	rBV2	714375	1226611	31.55%	6.971%
15	24.980	3716	3726	3746	rVB3	329542	1142742	29.39%	6.495%

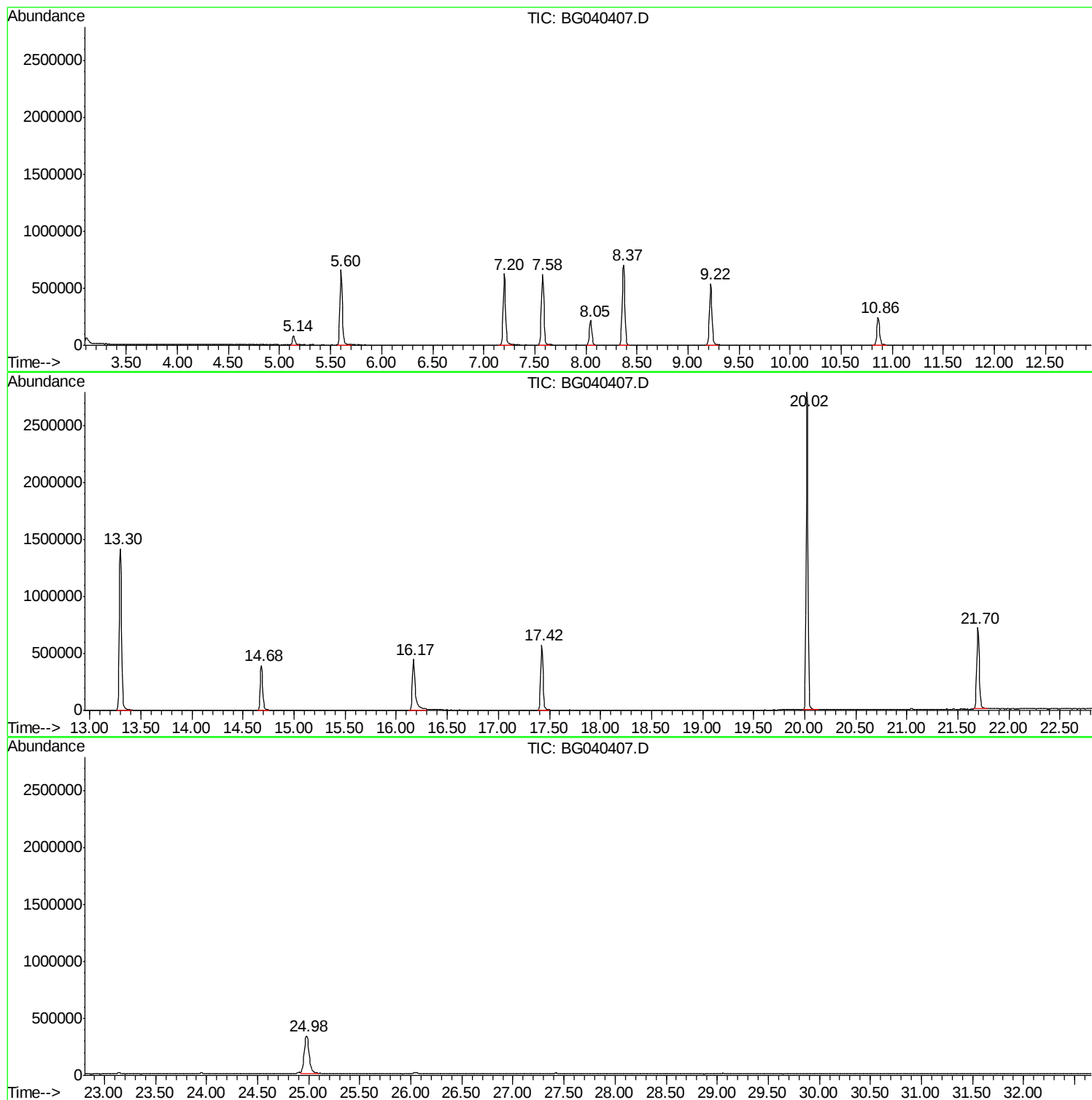
Sum of corrected areas: 17595157

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.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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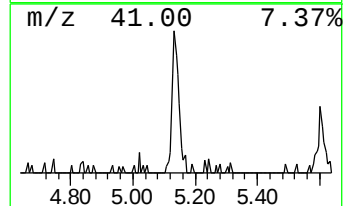
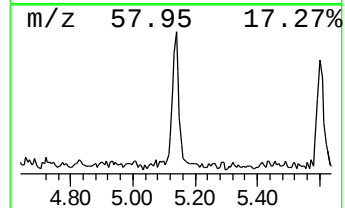
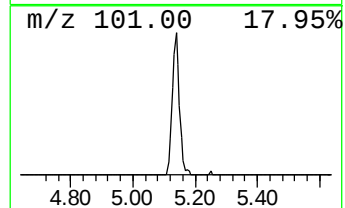
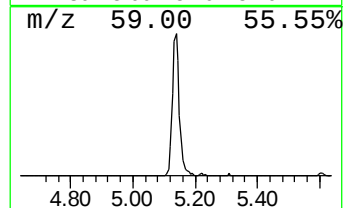
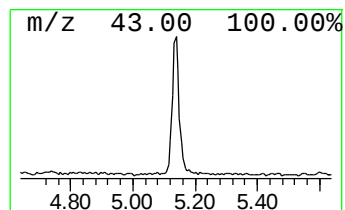
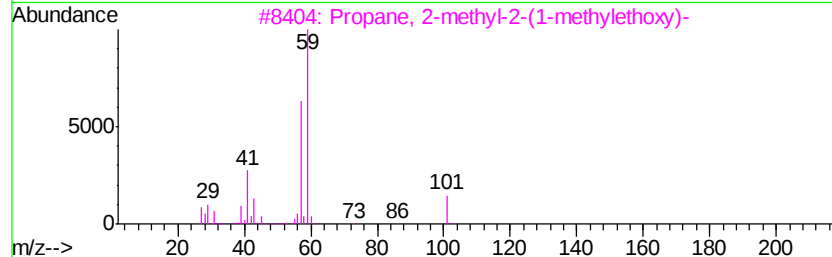
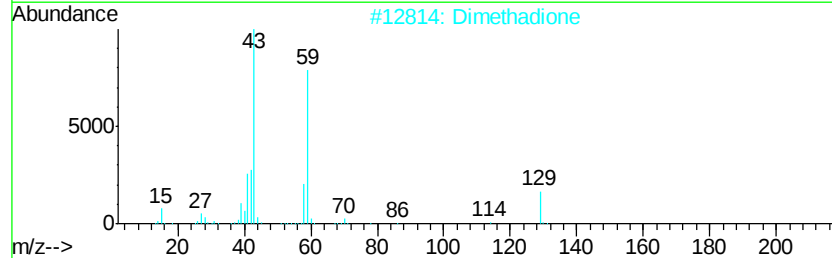
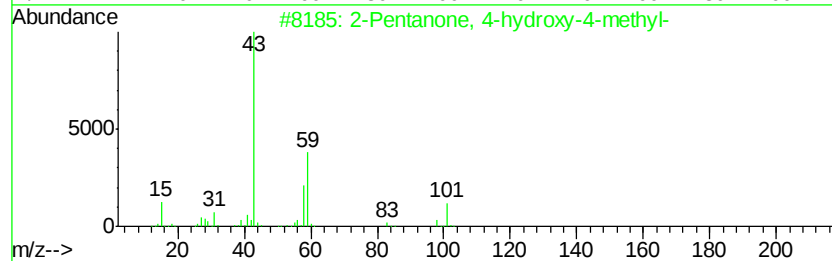
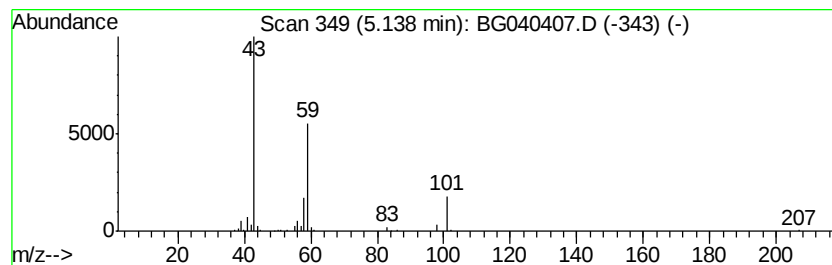
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	7.14 ng	129155	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Dimethadione	129	C5H7NO3	000695-53-4	42
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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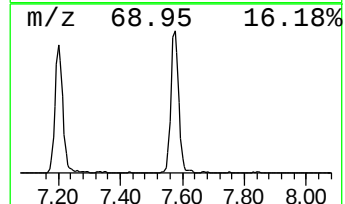
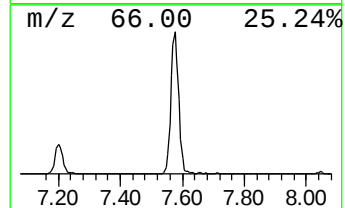
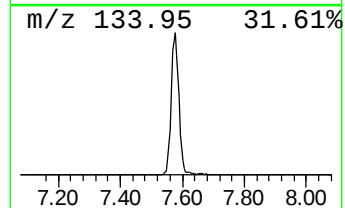
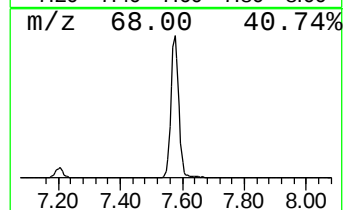
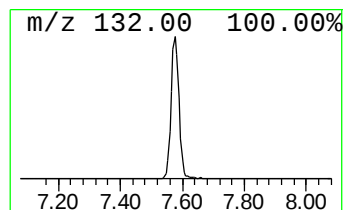
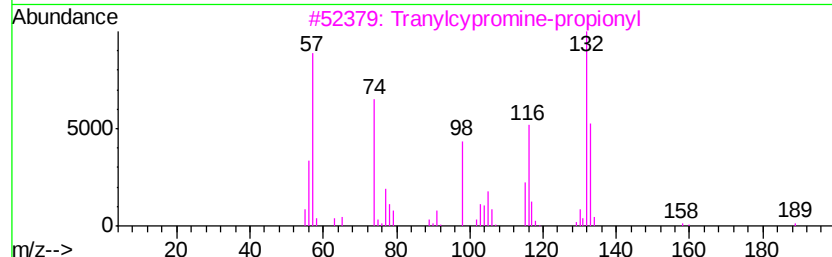
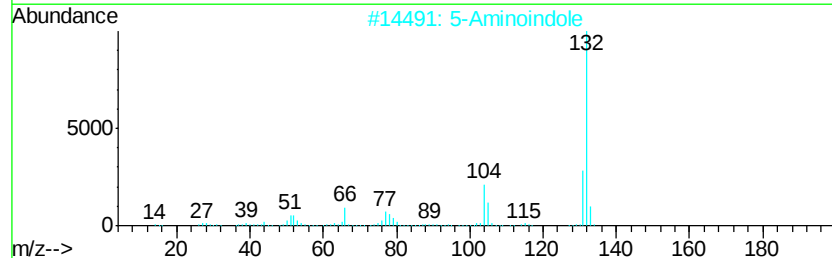
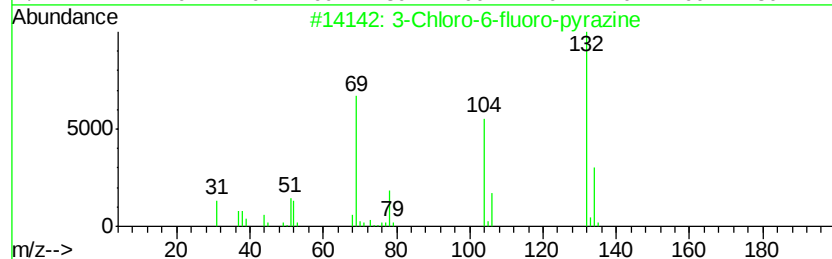
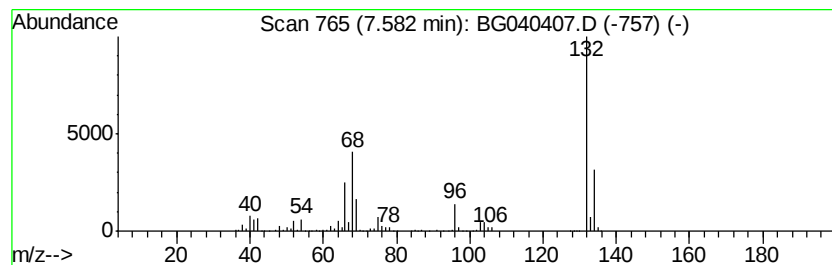
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	60.07 ng	1087390	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	7.1	ng	129155	1	8.05	362032	20.0
unknown7.58	7.58	60.1	ng	1087390	1	8.05	362032	20.0