

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG041919\
 Data File : BG040574.D
 Acq On : 19 Apr 2019 22:57
 Operator : JU/SJ
 Sample : PB119049BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119049BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	5.120	341	346	355	rBV	79872	127120	3.88%	0.743%
2	5.590	419	426	441	rBV	619728	1024879	31.31%	5.992%
3	7.188	688	698	712	rBV	607292	1067904	32.62%	6.243%
4	7.558	753	761	774	rBV	597812	1056914	32.29%	6.179%
5	8.028	834	841	850	rBV	243225	395316	12.08%	2.311%
6	8.351	889	896	905	rVB	967323	1661860	50.77%	9.716%
7	9.203	1033	1041	1059	rBV	525525	976186	29.82%	5.707%
8	10.842	1312	1320	1333	rBV	227891	432359	13.21%	2.528%
9	13.287	1728	1736	1752	rBV	1331872	2179669	66.58%	12.743%
10	14.667	1963	1971	1983	rBV2	427886	760467	23.23%	4.446%
11	16.154	2217	2224	2241	rBV	602097	1098316	33.55%	6.421%
12	17.411	2431	2438	2450	rBV	523226	821789	25.10%	4.805%
13	19.826	2838	2849	2853	rBV9	26985	87651	2.68%	0.512%
14	19.861	2853	2855	2866	rVB5	26081	63387	1.94%	0.371%
15	20.014	2874	2881	2889	rBV	2341699	3273613	100.00%	19.139%
16	21.683	3158	3165	3178	rBV2	551462	952330	29.09%	5.568%
17	22.429	3287	3292	3298	rBV4	32045	58758	1.79%	0.344%
18	23.075	3391	3402	3403	rBV9	27223	74760	2.28%	0.437%
19	23.921	3541	3546	3553	rVB6	18080	39274	1.20%	0.230%
20	24.955	3712	3722	3738	rVB2	304140	951747	29.07%	5.564%

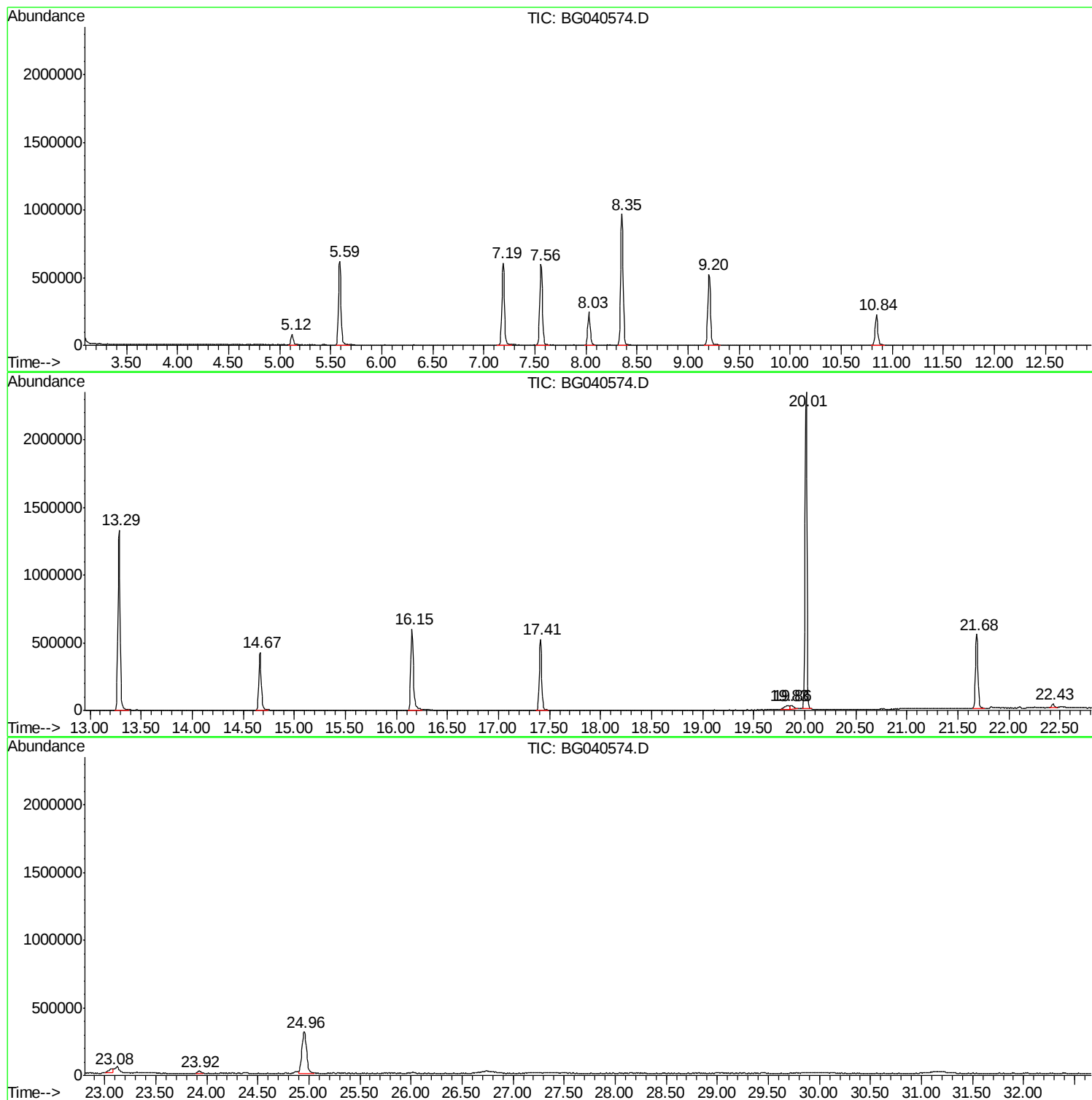
Sum of corrected areas: 17104299

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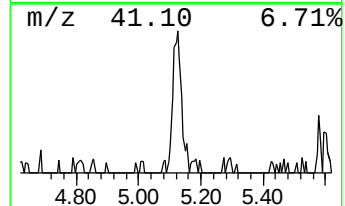
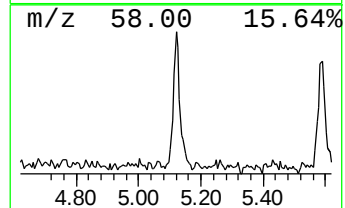
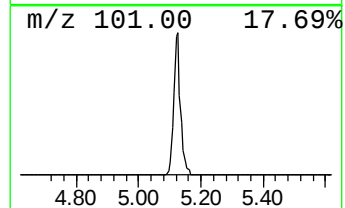
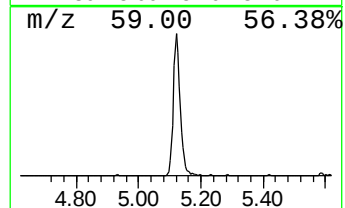
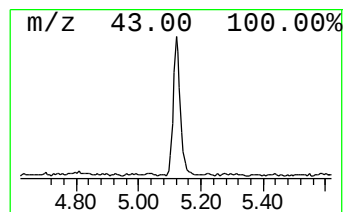
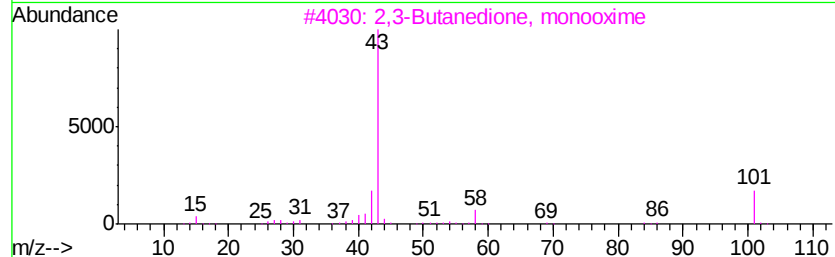
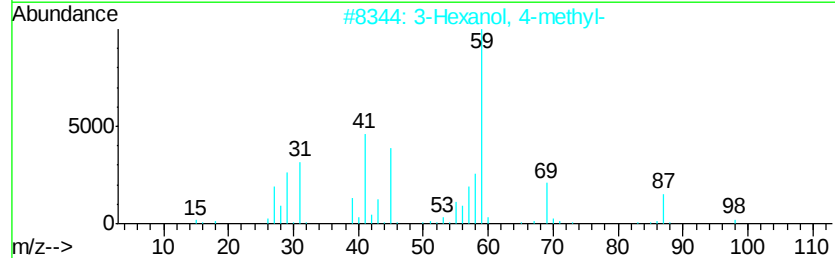
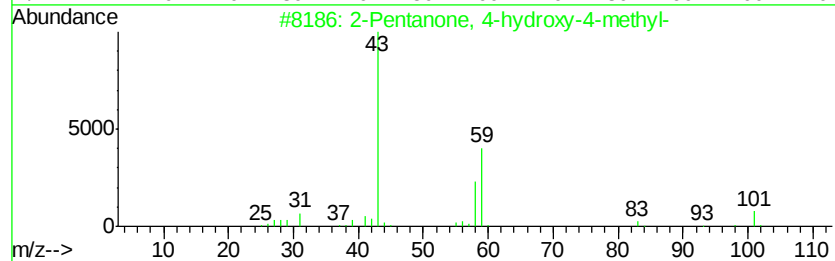
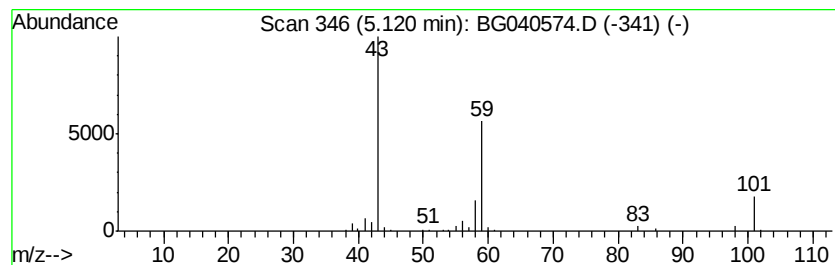
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	6.43 ng	127120	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			Guanidine	59	CH5N3	000113-00-8	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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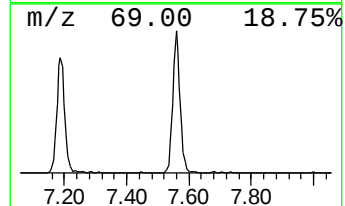
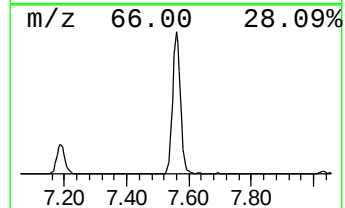
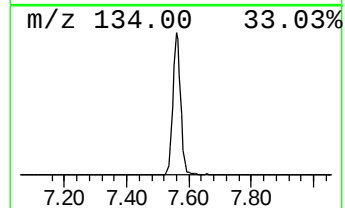
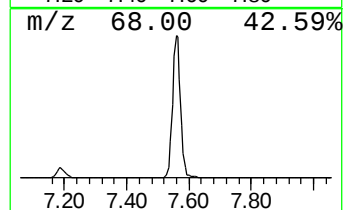
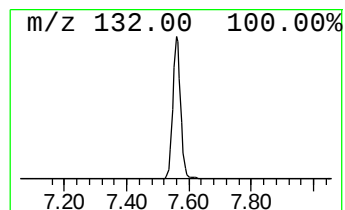
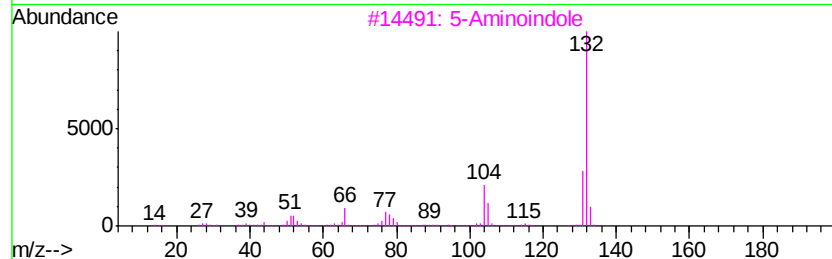
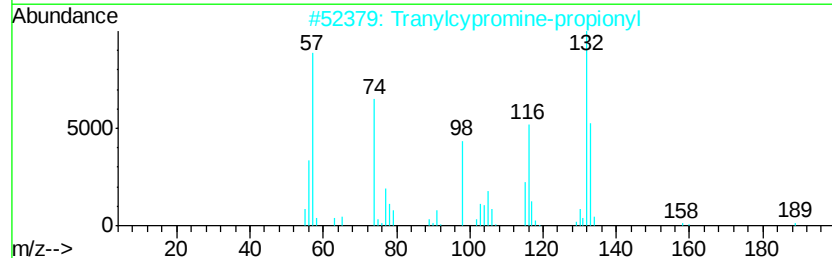
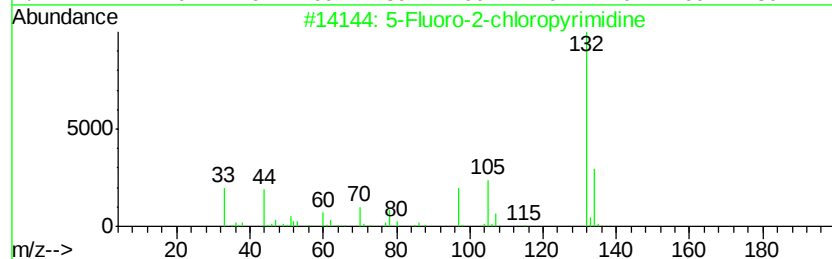
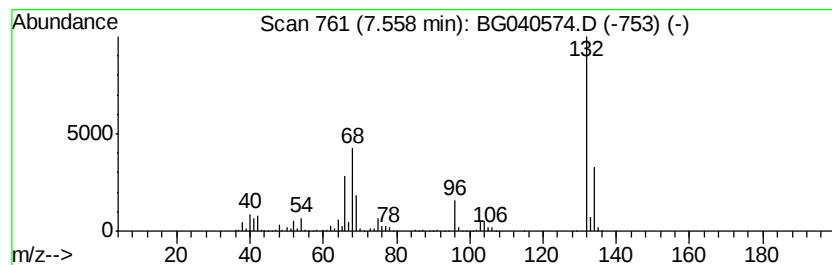
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Peak Number 2 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	53.47 ng	1056910	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	6.4	ng	127120	1	8.03	395316	20.0
unknown7.56	7.56	53.5	ng	1056910	1	8.03	395316	20.0