

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011719\
 Data File : BG039186.D
 Acq On : 17 Jan 2019 14:46
 Operator : JU/SJ
 Sample : PB116407BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116407BL

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-BG010919.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	5.097	337	342	349	rVB	15586	24734	1.30%	0.338%
2	5.567	415	422	435	rVB	283167	467193	24.55%	6.387%
3	7.171	688	695	707	rBV	261186	466136	24.50%	6.373%
4	7.542	751	758	769	rBV	264903	478591	25.15%	6.543%
5	8.018	831	839	846	rVB2	66931	114571	6.02%	1.566%
6	8.335	886	893	904	rVB	229141	396753	20.85%	5.424%
7	9.193	1032	1039	1051	rBV	150707	286216	15.04%	3.913%
8	10.832	1310	1318	1327	rBV	72983	139206	7.32%	1.903%
9	13.276	1727	1734	1746	rBV	478429	780527	41.02%	10.671%
10	14.657	1962	1969	1980	rBB2	134489	220300	11.58%	3.012%
11	16.149	2216	2223	2242	rBV	415069	665396	34.97%	9.097%
12	17.406	2430	2437	2447	rBV	205759	329669	17.33%	4.507%
13	20.009	2874	2880	2888	rBV	1447963	1902648	100.00%	26.012%
14	21.678	3157	3164	3172	rBV	262166	446284	23.46%	6.101%
15	22.418	3285	3290	3295	rVB2	13229	19438	1.02%	0.266%
16	23.106	3400	3407	3414	rBV3	18328	38401	2.02%	0.525%
17	23.911	3537	3544	3550	rVB4	11842	24175	1.27%	0.331%
18	24.851	3697	3704	3710	rBV7	9593	25477	1.34%	0.348%
19	24.945	3710	3720	3731	rVB	164654	488672	25.68%	6.681%

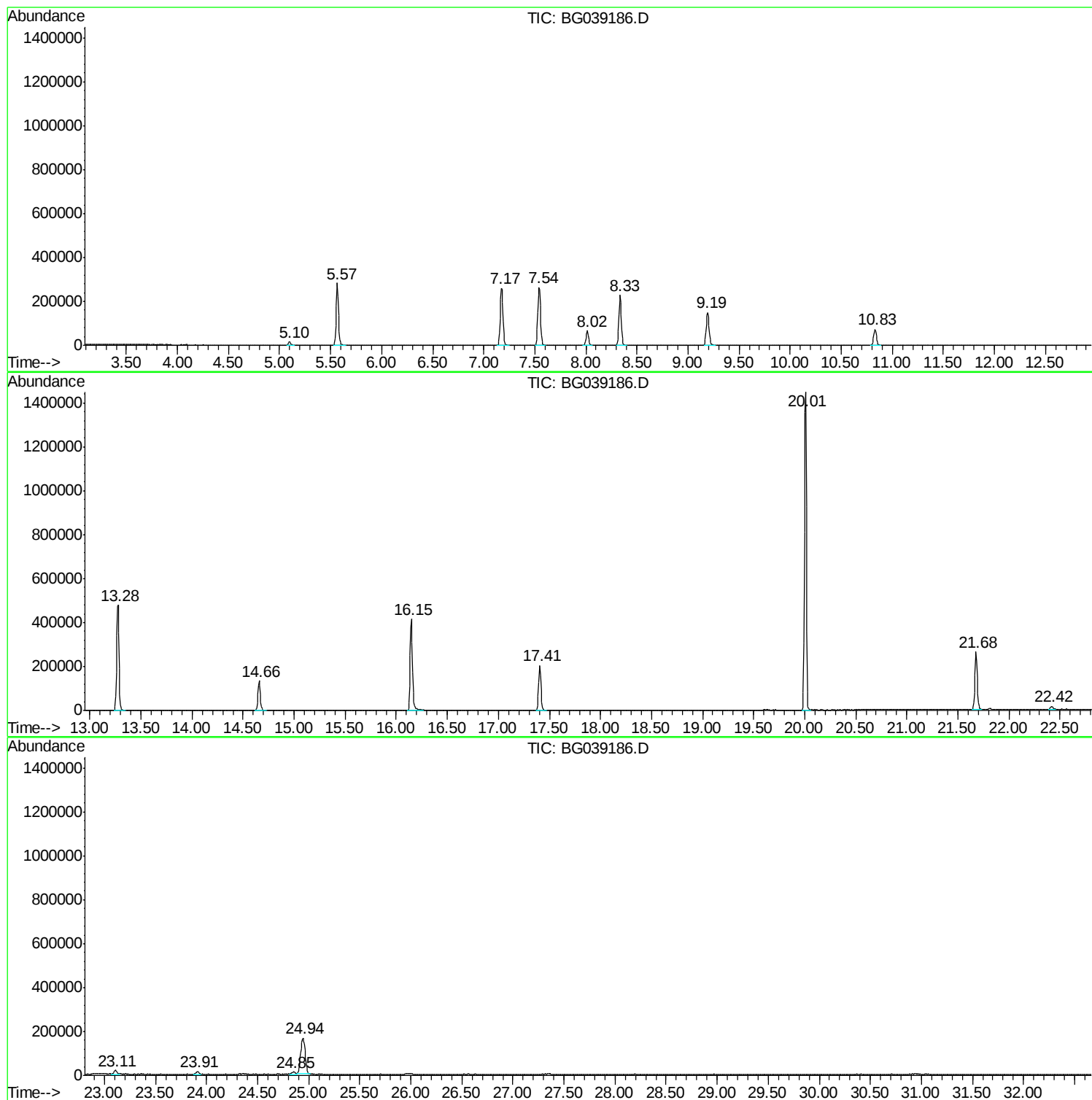
Sum of corrected areas: 7314387

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.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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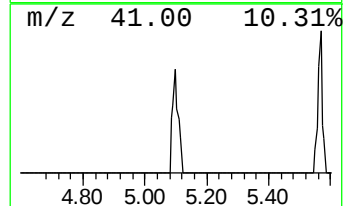
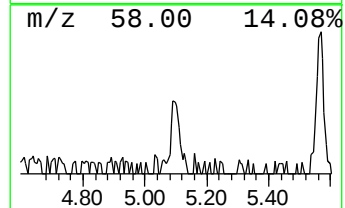
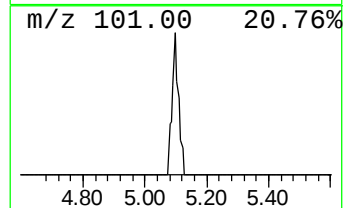
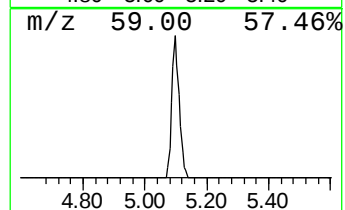
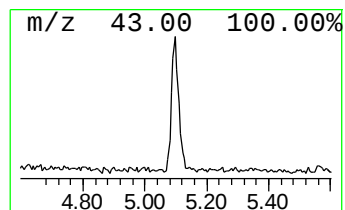
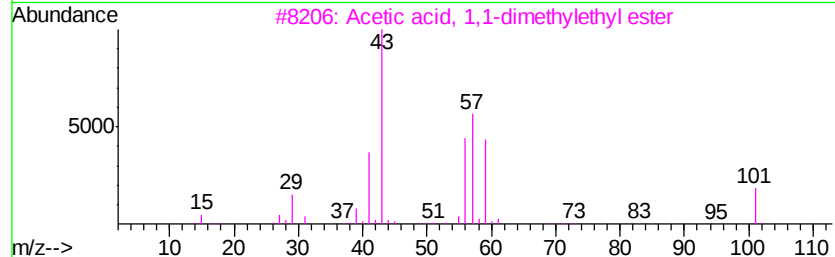
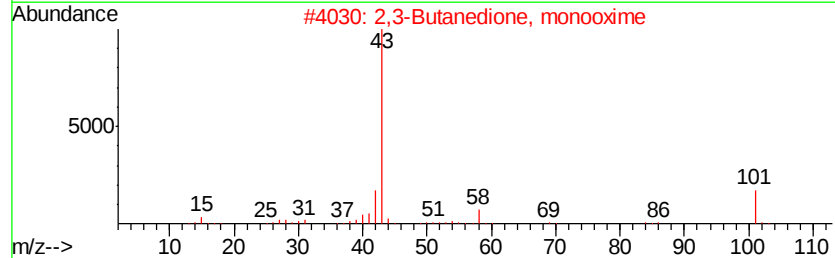
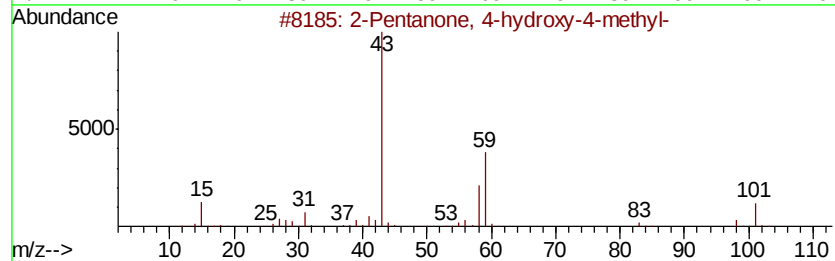
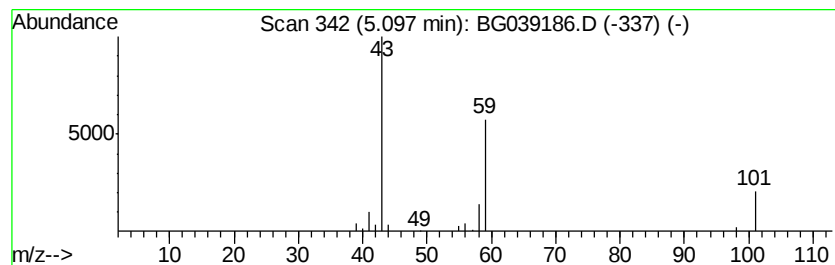
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	4.32 ng	24734	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4			1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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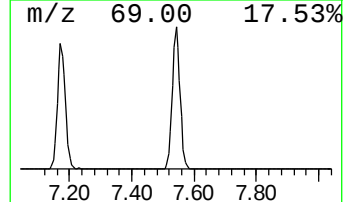
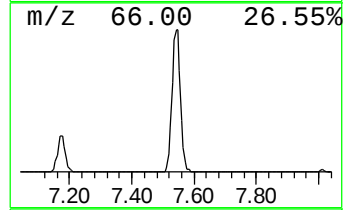
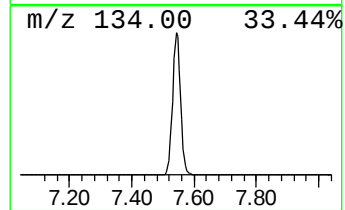
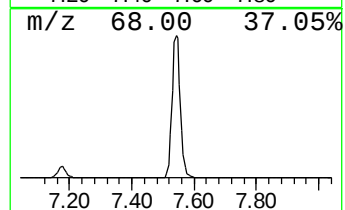
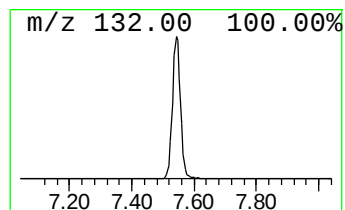
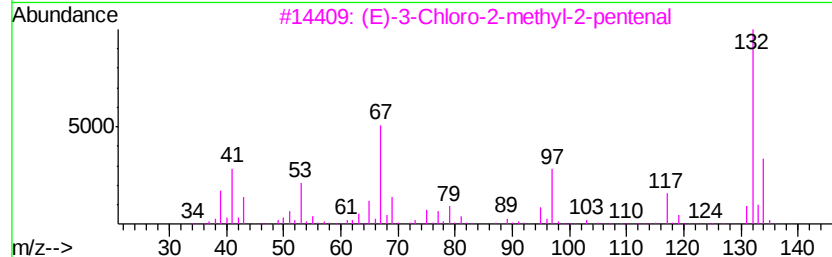
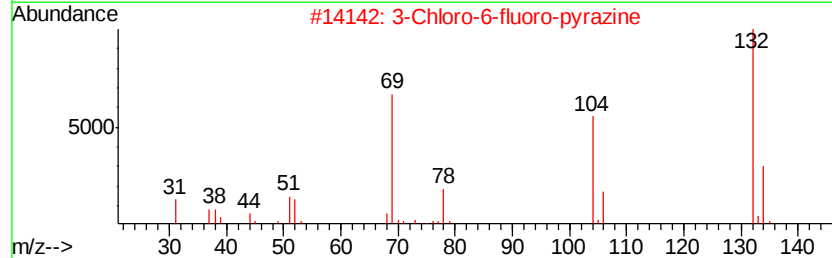
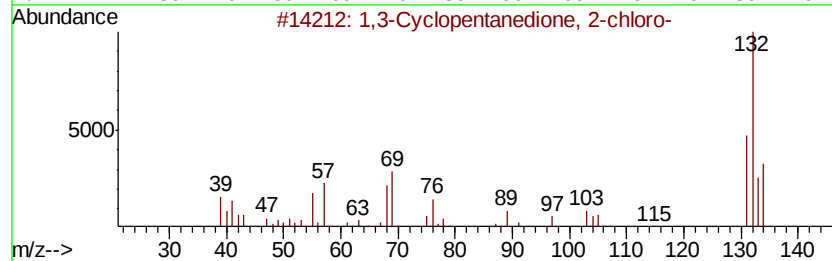
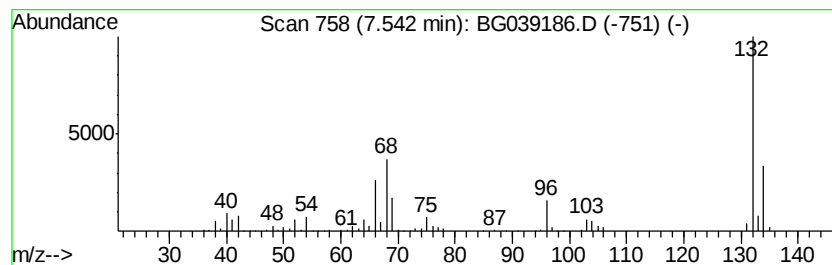
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Peak Number 2 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	83.54 ng	478591	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	37
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	4.3	ng	24734	1	8.02	114571	20.0
unknown7.54	7.54	83.5	ng	478591	1	8.02	114571	20.0