

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102918\
 Data File : BM017404.D
 Acq On : 29 Oct 2018 12:35
 Operator : MJ/SJ
 Sample : PB114216BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB114216BL

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_M\METHODS\8270-BM101718.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	4.822	306	312	322	rBV	148779	222810	1.38%	0.266%
2	5.269	381	388	404	rBV	4275909	6075639	37.76%	7.249%
3	6.834	647	654	676	rBV	4216385	6605402	41.05%	7.881%
4	7.186	706	714	725	rBV	4321251	6714003	41.73%	8.010%
5	7.645	785	792	800	rBV	1036291	1614615	10.03%	1.926%
6	7.963	838	846	858	rBV	3205842	5048671	31.38%	6.023%
7	8.798	978	988	1008	rBV	2384603	4083456	25.38%	4.872%
8	10.422	1253	1264	1278	rBV	1353409	2320582	14.42%	2.769%
9	12.904	1679	1686	1706	rBV	6666777	10105433	62.80%	12.056%
10	14.292	1914	1922	1939	rBV2	2039967	3256342	20.24%	3.885%
11	15.786	2169	2176	2193	rBV	4829303	7878396	48.96%	9.399%
12	17.045	2383	2390	2403	rBV	2722074	4178470	25.97%	4.985%
13	19.386	2784	2788	2793	rBV2	205469	233230	1.45%	0.278%
14	19.686	2832	2839	2849	rBV	12587897	16090402	100.00%	19.197%
15	20.439	2963	2967	2974	rBV2	178660	217291	1.35%	0.259%
16	21.238	3098	3103	3113	rBV	3831478	5078216	31.56%	6.059%
17	23.450	3472	3479	3495	rVB	2004605	4094733	25.45%	4.885%

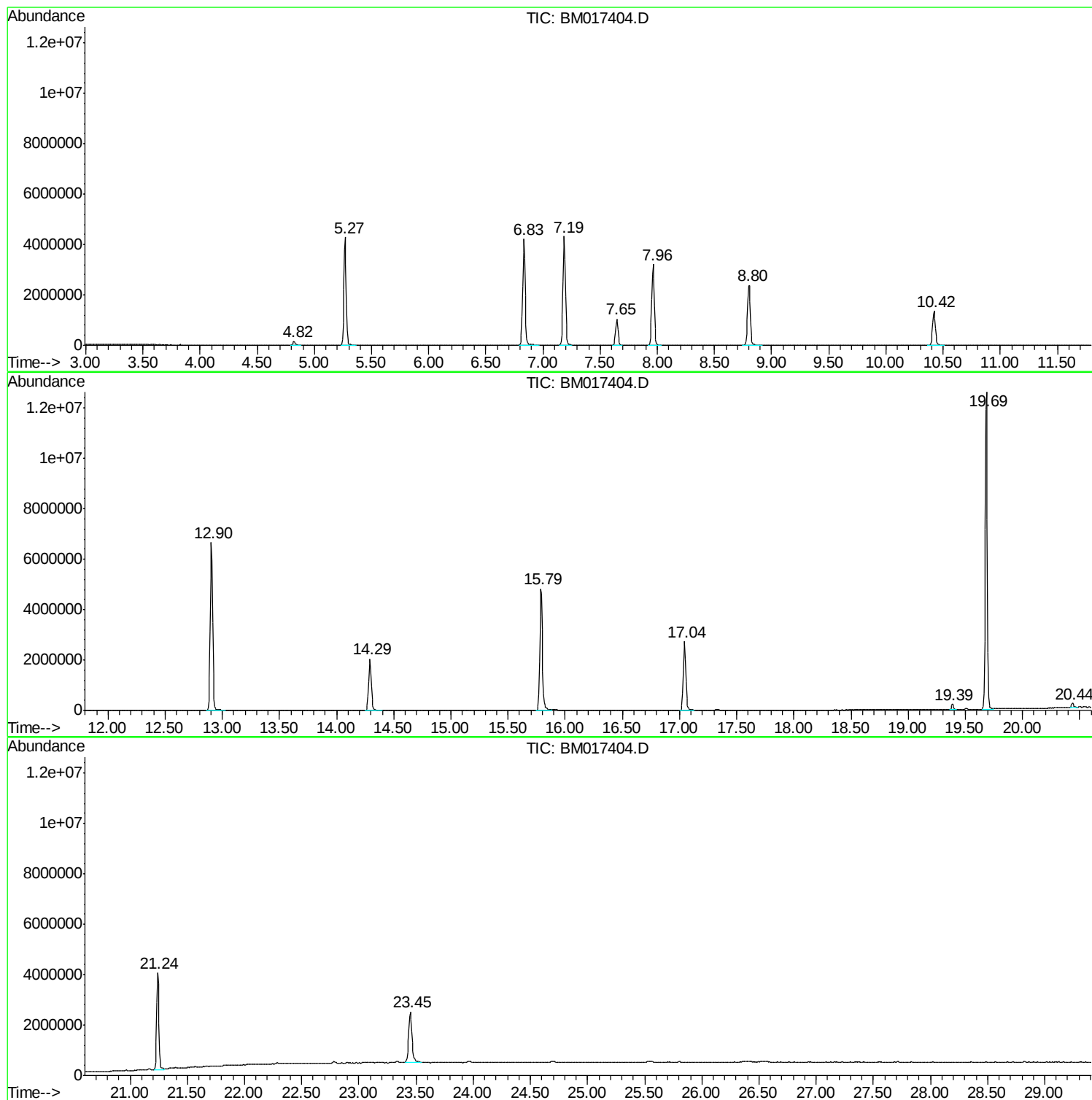
Sum of corrected areas: 83817691

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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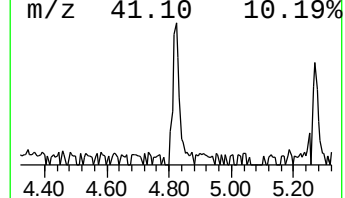
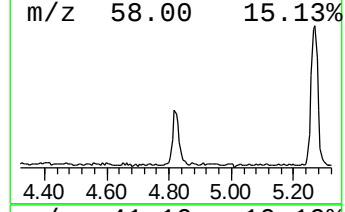
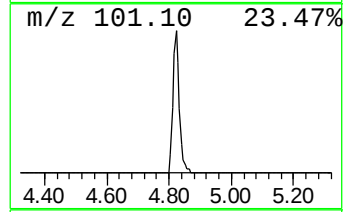
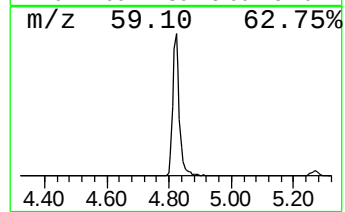
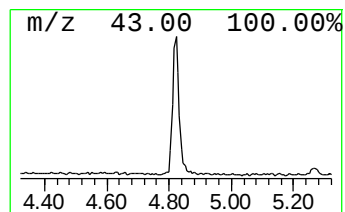
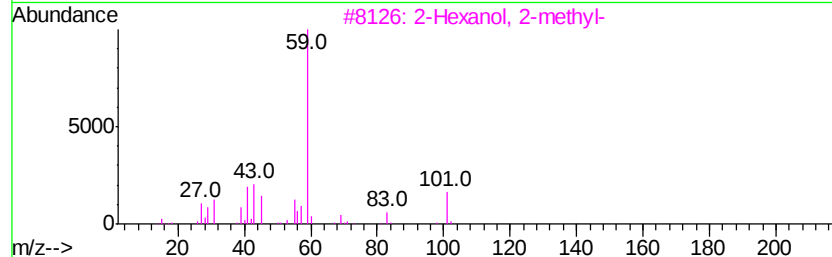
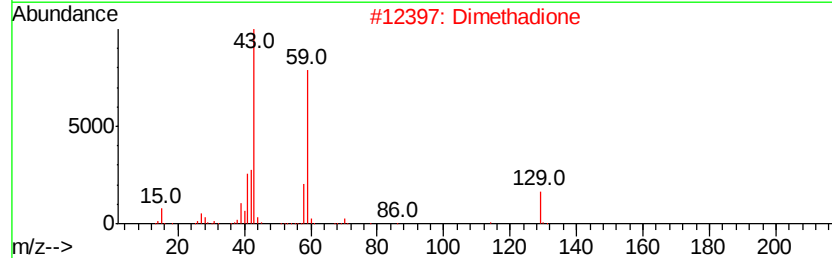
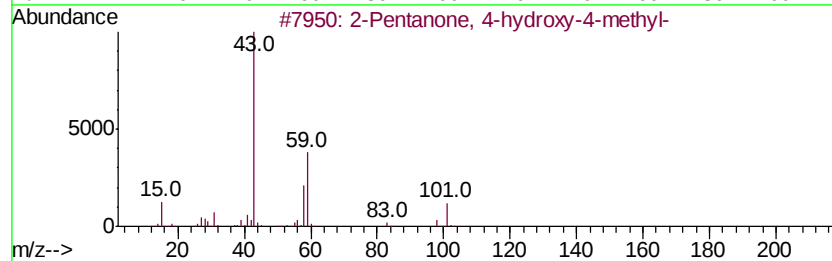
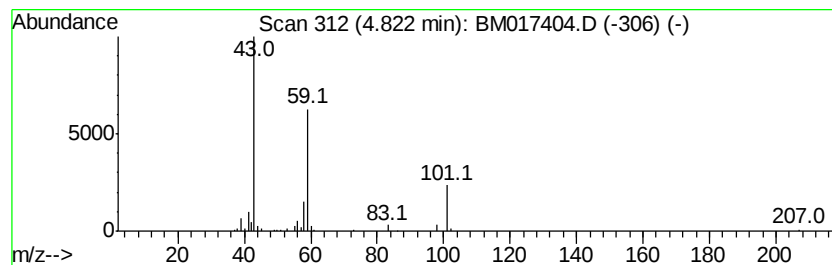
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	2.76 ng	222810	1,4-Dichlorobenzene-d4	7.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Dimethadione	129	C5H7NO3	000695-53-4	40
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	23
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12



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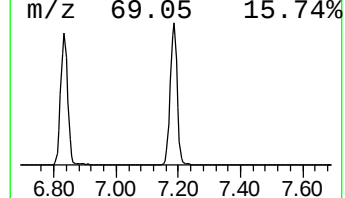
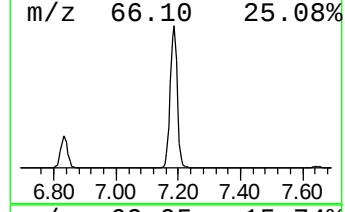
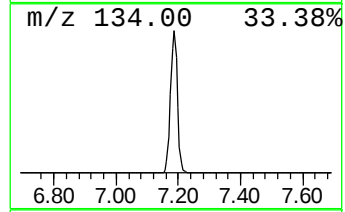
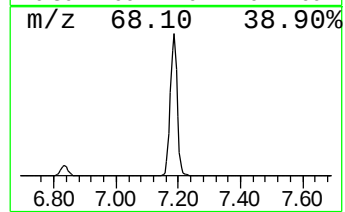
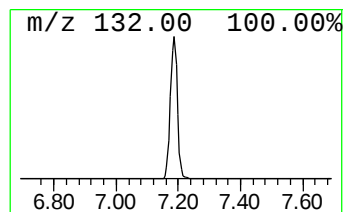
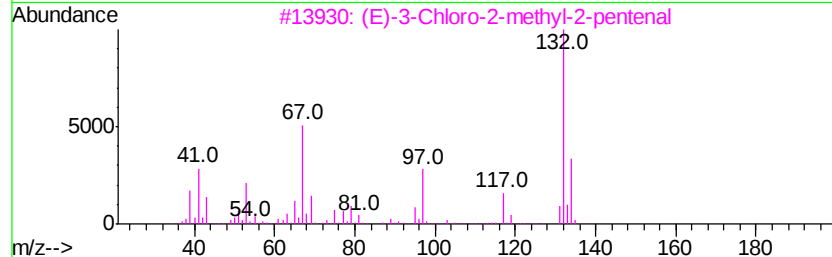
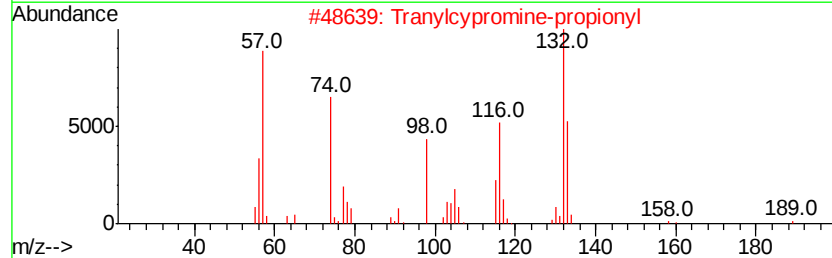
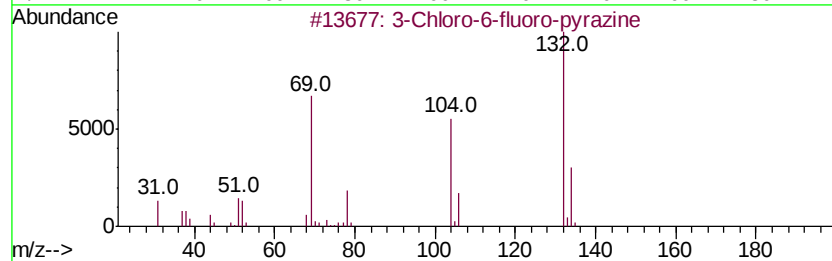
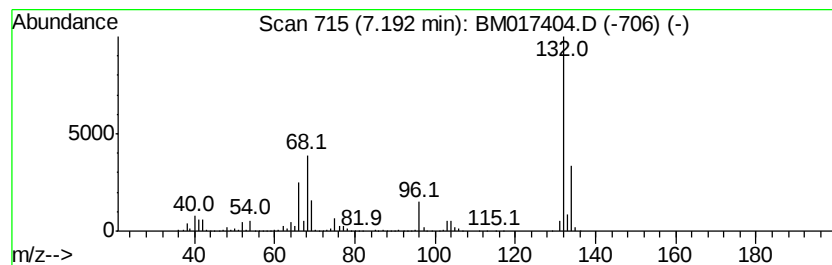
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Peak Number 2 unknown7.19 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	83.17 ng	6714000	1,4-Dichlorobenzene-d4	7.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.82	2.8	ng	222810	1	7.65	1614620	20.0
unknown7.19	7.19	83.2	ng	6714000	1	7.65	1614620	20.0