

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091418\
 Data File : BF109014.D
 Acq On : 15 Sep 2018 00:02
 Operator : JU/SJ
 Sample : J4906-32
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-PIT-2

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_F\METHODS\8270-BF091418.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	2.607	42	46	54	rVB	46844	62514	1.54%	0.198%
2	4.919	435	439	443	rBV	186490	227261	5.61%	0.719%
3	5.337	505	510	513	rBV	2982105	2907136	71.73%	9.200%
4	6.354	678	683	686	rBV	2795407	2744293	67.71%	8.684%
5	6.489	701	706	709	rBV	3414278	3304362	81.53%	10.457%
6	6.719	742	745	749	rVB	1085200	873252	21.55%	2.763%
7	6.878	768	772	775	rBV	3251580	2733444	67.45%	8.650%
8	7.278	835	840	843	rBV	2205959	2244887	55.39%	7.104%
9	8.001	959	963	965	rBV	1364719	1135269	28.01%	3.593%
10	8.019	965	966	970	rVB	218394	133554	3.30%	0.423%
11	8.713	1080	1084	1087	rBV	63375	49833	1.23%	0.158%
12	9.078	1141	1146	1149	rBV	4547163	4052781	100.00%	12.825%
13	9.748	1256	1260	1264	rBV2	1551284	1270884	31.36%	4.022%
14	10.530	1389	1393	1397	rBV	2436325	2504715	61.80%	7.926%
15	11.224	1507	1511	1514	rBV	1591206	1306569	32.24%	4.135%
16	12.813	1776	1781	1784	rBV	3910385	3875936	95.64%	12.265%
17	13.848	1953	1957	1961	rBV	1204145	1033529	25.50%	3.271%
18	14.954	2141	2145	2150	rBV	38827	44106	1.09%	0.140%
19	15.248	2190	2195	2201	rBV	812222	980843	24.20%	3.104%
20	15.307	2201	2205	2210	rVB	44657	55472	1.37%	0.176%
21	15.712	2268	2274	2282	rVB2	36714	59857	1.48%	0.189%

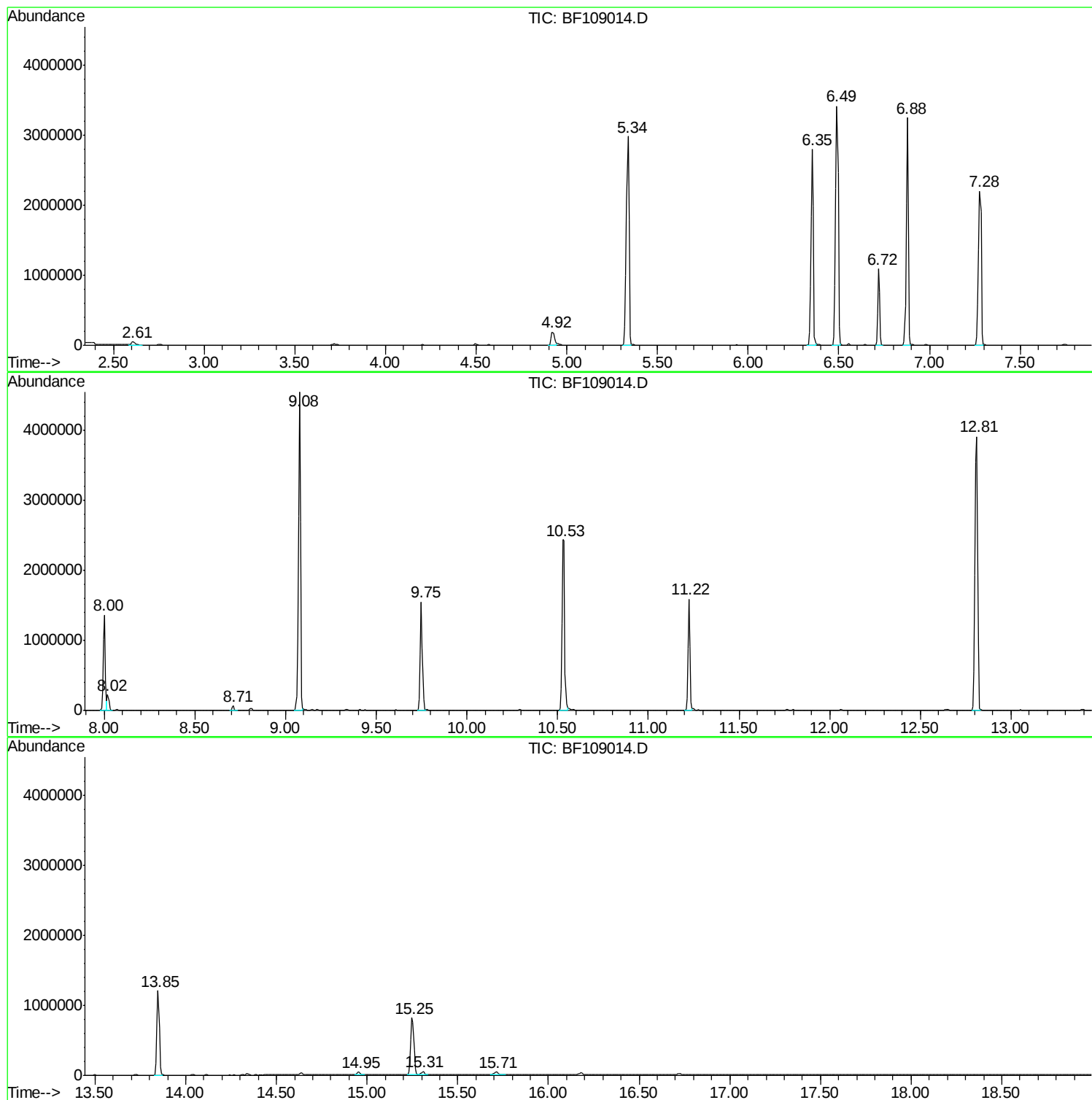
Sum of corrected areas: 31600497

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.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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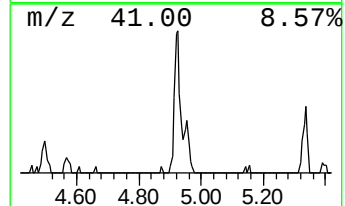
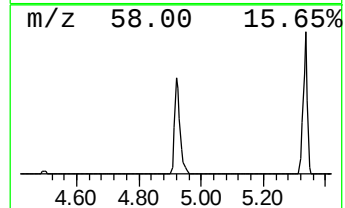
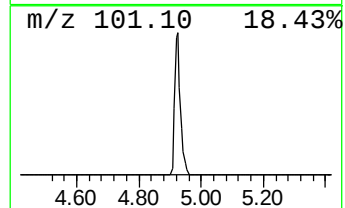
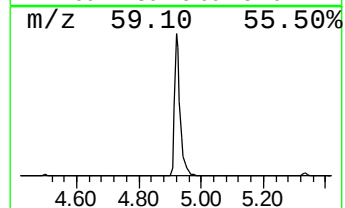
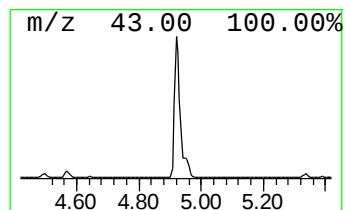
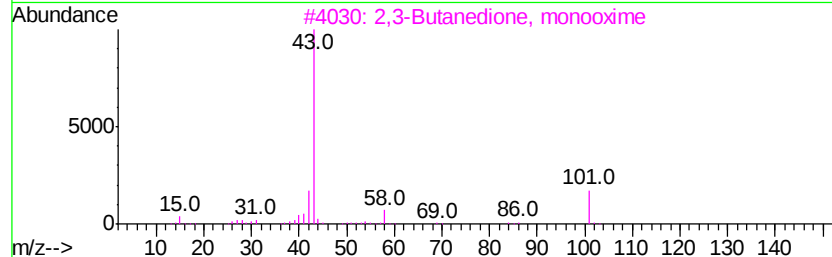
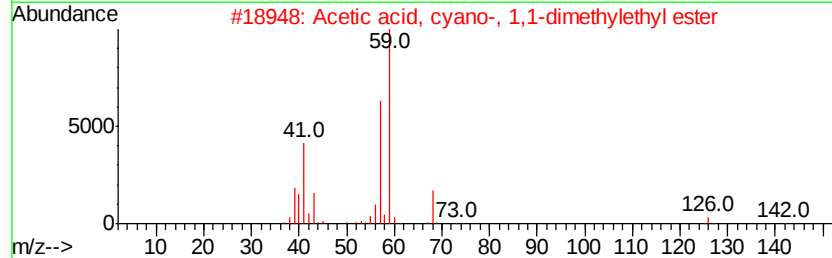
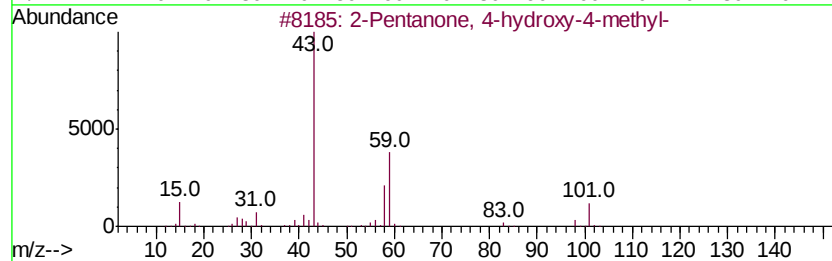
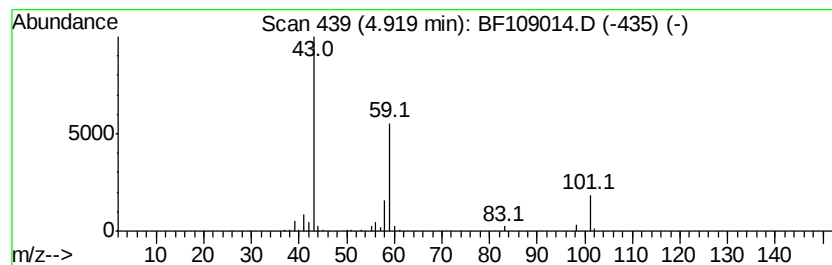
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	5.20 ng	227261	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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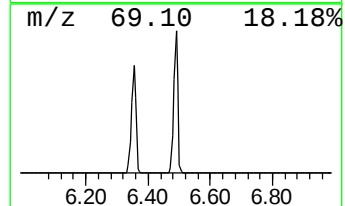
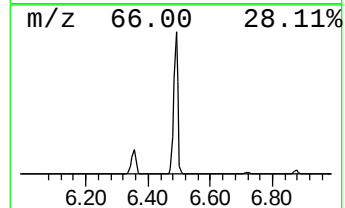
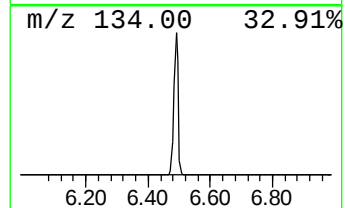
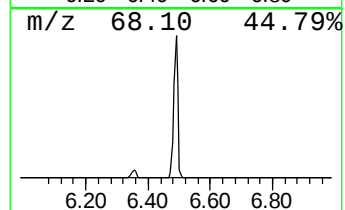
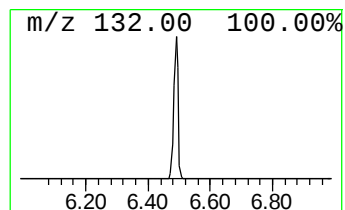
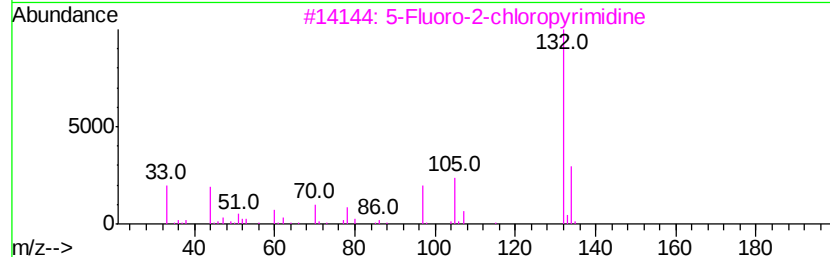
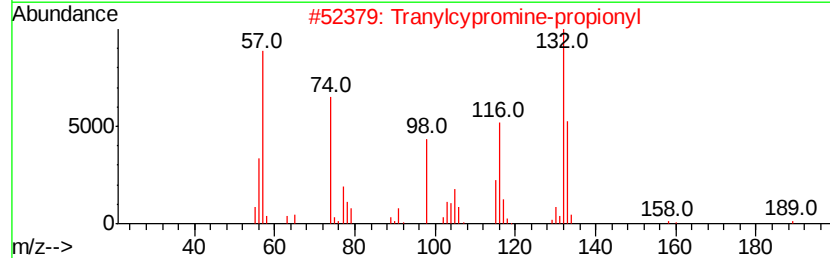
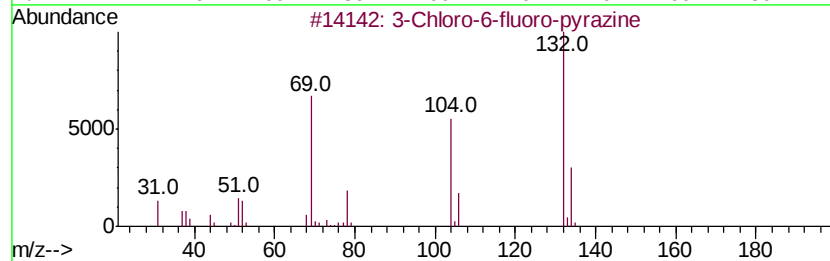
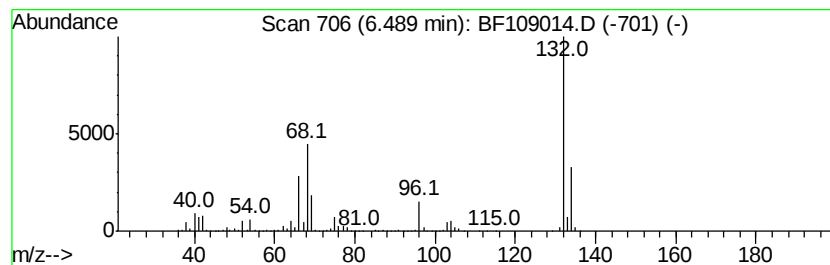
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Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	75.68 ng	3304360	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.92	5.2	ng	227261	1	6.72	873252	20.0
unknown6.49	6.49	75.7	ng	3304360	1	6.72	873252	20.0