

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091418\
 Data File : BF109012.D
 Acq On : 14 Sep 2018 23:08
 Operator : JU/SJ
 Sample : J4906-24
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-PIT-10

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_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	2.596	40	44	53	rVB	33670	44137	1.02%	0.134%
2	4.925	433	440	449	rBV	237939	337148	7.81%	1.024%
3	5.337	505	510	513	rBV	3023777	3016089	69.86%	9.159%
4	6.354	678	683	686	rBV	2719789	2963257	68.63%	8.999%
5	6.489	701	706	709	rBV	3483046	3341634	77.40%	10.148%
6	6.719	742	745	749	rBV	1128252	886138	20.52%	2.691%
7	6.878	768	772	775	rBV	3254085	2702470	62.59%	8.207%
8	7.278	835	840	843	rBV	2262972	2243249	51.96%	6.812%
9	8.001	959	963	966	rBV	1328785	1151131	26.66%	3.496%
10	9.078	1141	1146	1149	rBV	4448625	4074763	94.38%	12.374%
11	9.748	1256	1260	1264	rBV2	1547342	1324067	30.67%	4.021%
12	10.536	1389	1394	1397	rBV	2634710	2660610	61.62%	8.079%
13	11.224	1507	1511	1515	rBV	1666189	1417883	32.84%	4.306%
14	12.813	1776	1781	1784	rBV	4360060	4317610	100.00%	13.111%
15	13.848	1953	1957	1961	rBV	1367922	1234486	28.59%	3.749%
16	14.954	2141	2145	2149	rBV	42468	45755	1.06%	0.139%
17	15.254	2190	2196	2202	rBV	837366	1067635	24.73%	3.242%
18	15.307	2202	2205	2209	rVB	44446	53798	1.25%	0.163%
19	15.712	2269	2274	2278	rBV2	36058	48556	1.12%	0.147%

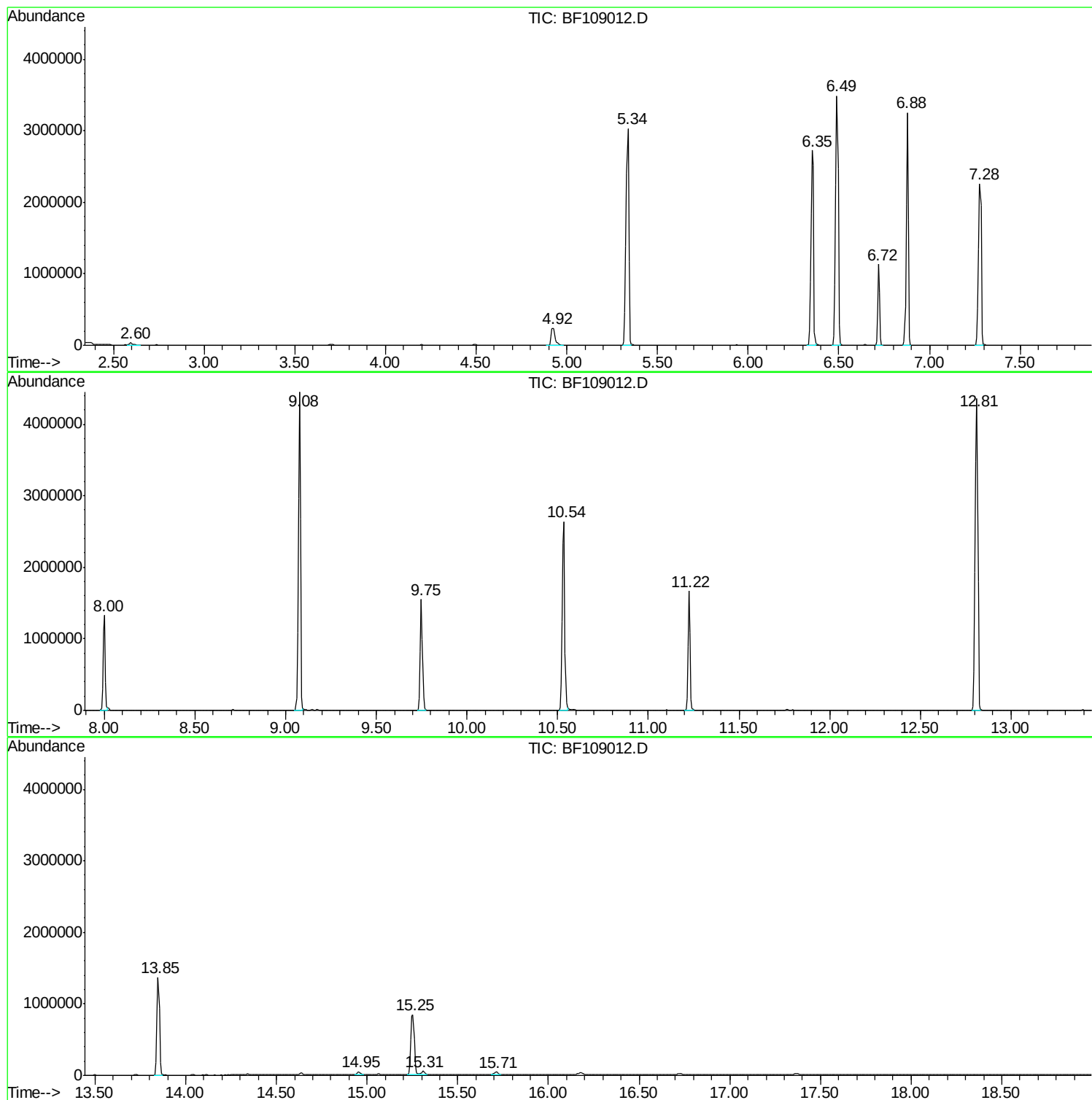
Sum of corrected areas: 32930416

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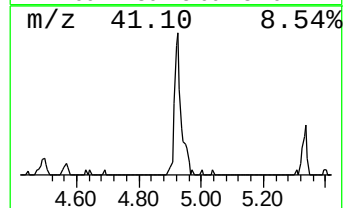
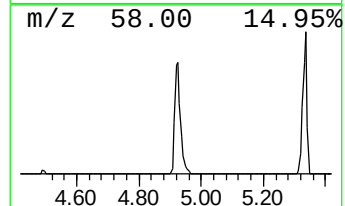
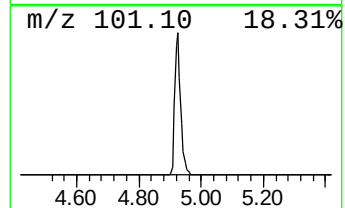
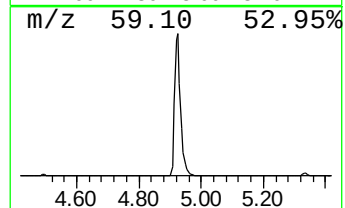
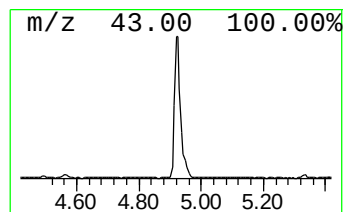
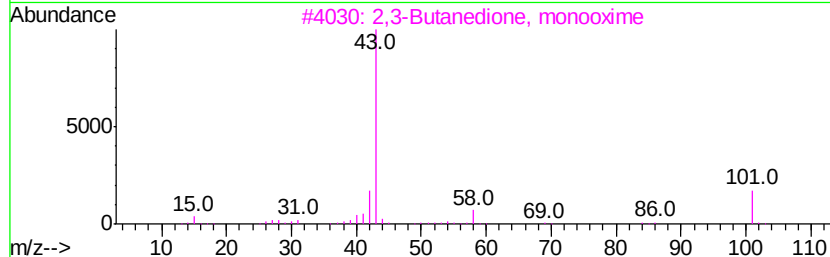
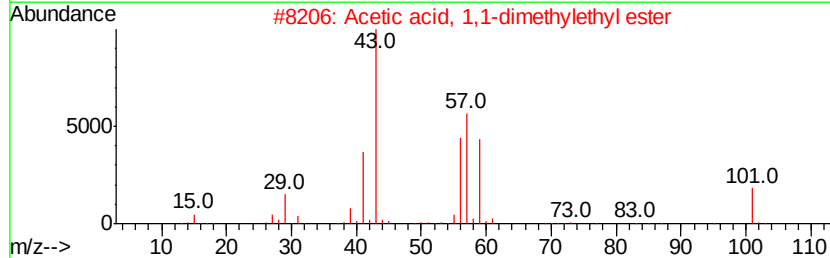
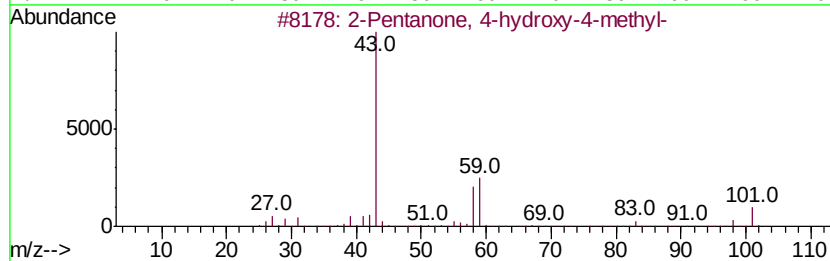
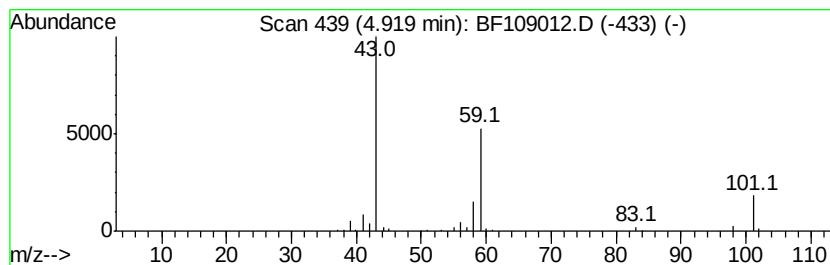
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	7.61 ng	337148	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
4			3-Hexanol	102	C6H14O	000623-37-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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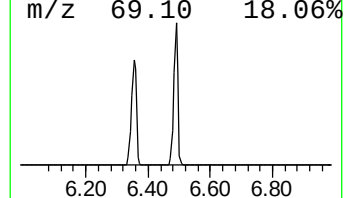
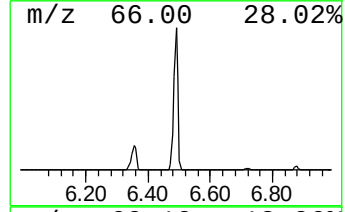
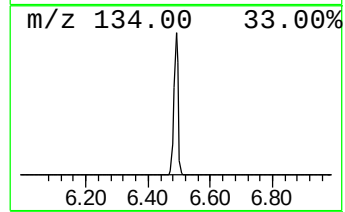
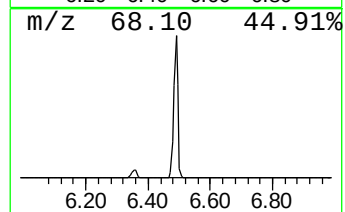
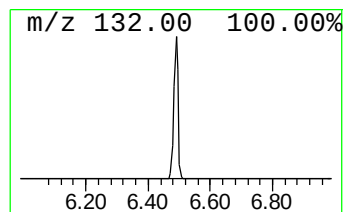
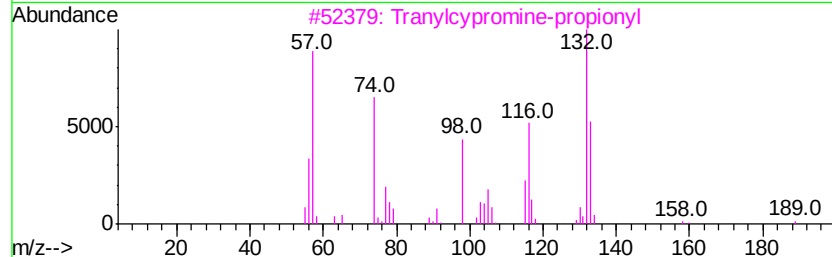
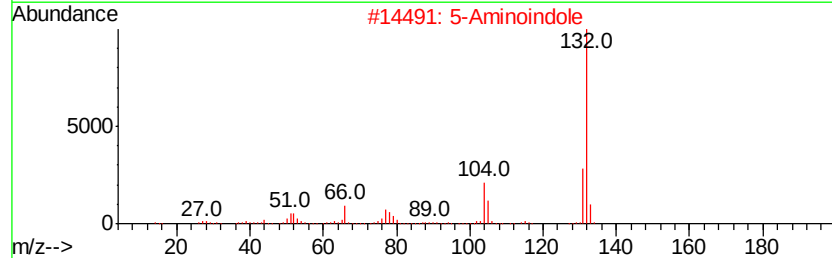
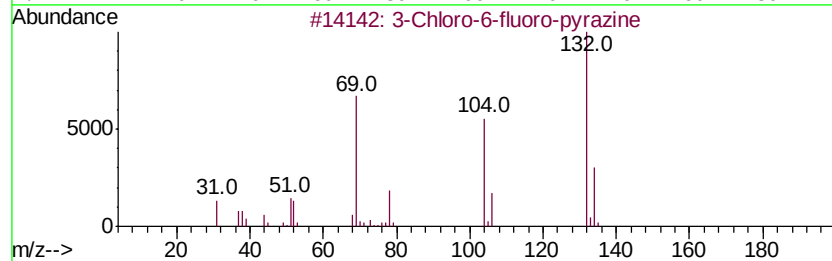
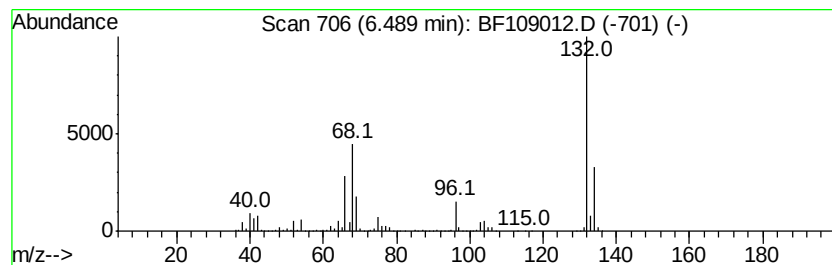
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Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	75.42 ng	3341630	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	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
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...	4.92	7.6	ng	337148	1	6.72	886138	20.0
unknown6.49	6.49	75.4	ng	3341630	1	6.72	886138	20.0