

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091018\
 Data File : BF108911.D
 Acq On : 10 Sep 2018 23:25
 Operator : JU/SJ
 Sample : PB112743BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB112743BL

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-BF090518.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	4.954	438	445	454	rBV	115355	159890	4.02%	0.488%
2	5.360	509	514	517	rBV	3943986	3741740	94.04%	11.426%
3	6.378	681	687	690	rBV	3725029	3833715	96.35%	11.707%
4	6.507	705	709	713	rBV	3904792	3814179	95.86%	11.647%
5	6.737	745	748	752	rBV	1229758	1077970	27.09%	3.292%
6	6.896	771	775	779	rBV	3414397	3111215	78.19%	9.500%
7	7.295	838	843	846	rBV	2275450	2346662	58.98%	7.166%
8	8.019	962	966	970	rBV	1511985	1249684	31.41%	3.816%
9	9.095	1144	1149	1152	rBV	4460221	3978815	100.00%	12.150%
10	9.772	1260	1264	1268	rBV	1319550	1168245	29.36%	3.567%
11	10.554	1393	1397	1414	rBV	2075547	1866300	46.91%	5.699%
12	11.248	1511	1515	1518	rBV	1301169	1015877	25.53%	3.102%
13	12.830	1780	1784	1787	rBV	3777015	3361456	84.48%	10.264%
14	13.872	1957	1961	1965	rBV	1292391	1191681	29.95%	3.639%
15	15.283	2196	2201	2206	rBV	700784	831145	20.89%	2.538%

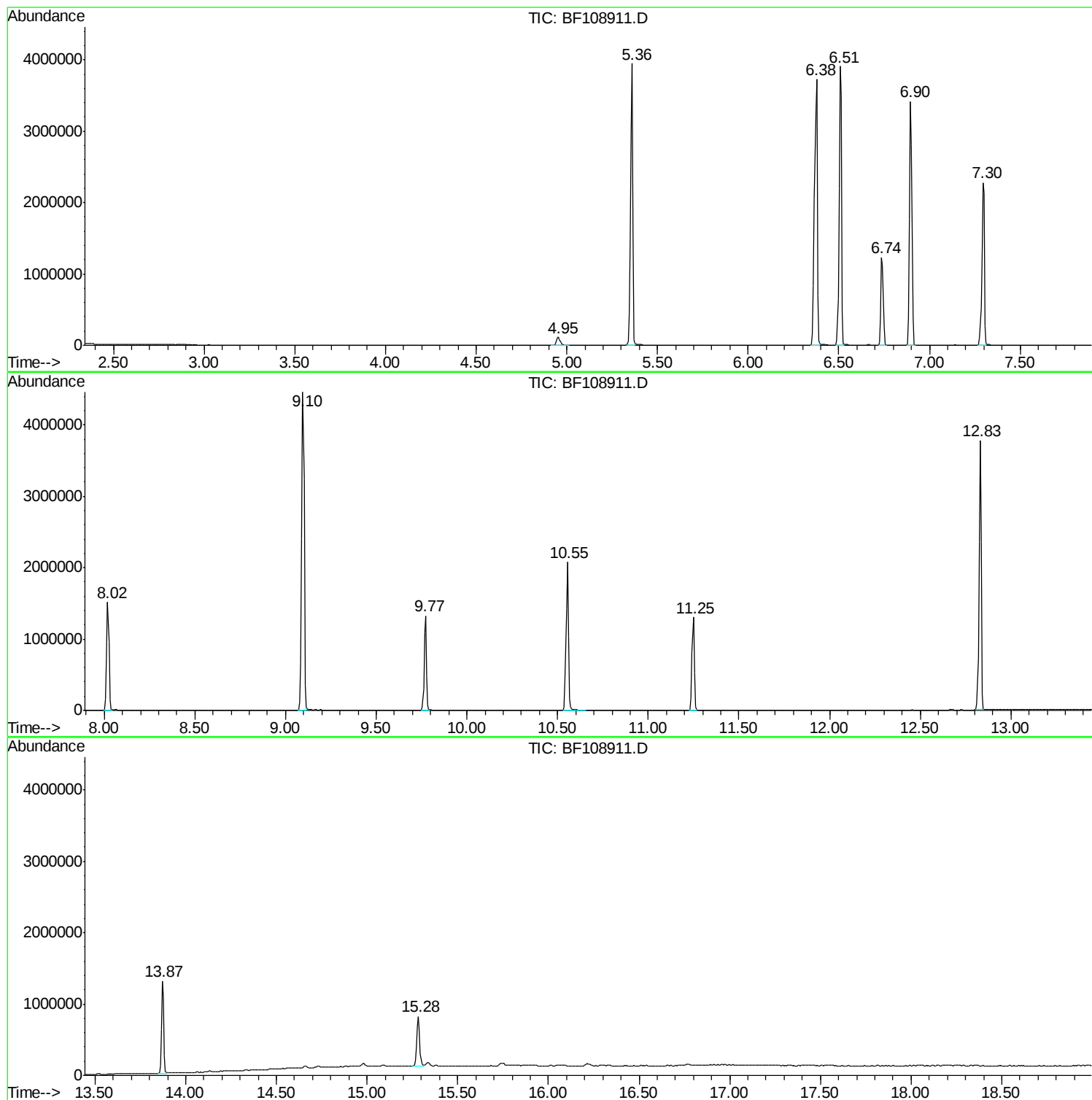
Sum of corrected areas: 32748574

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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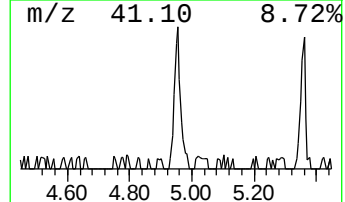
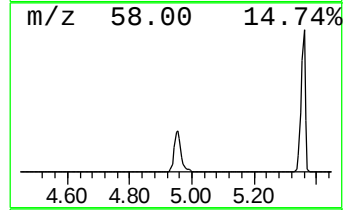
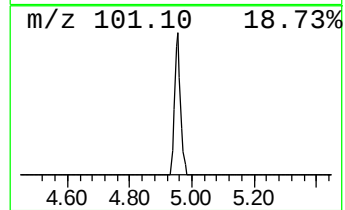
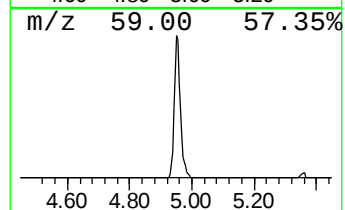
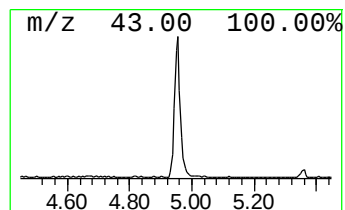
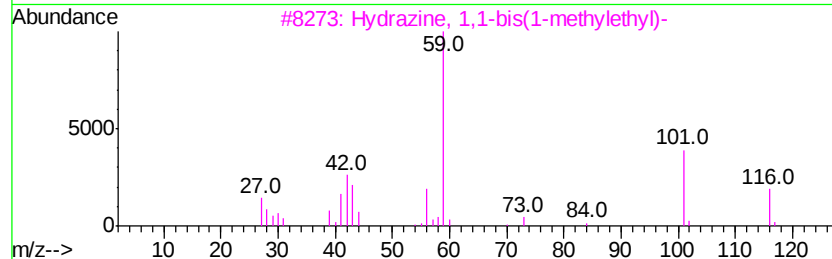
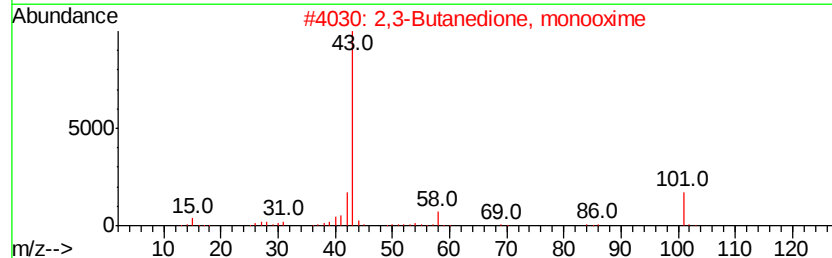
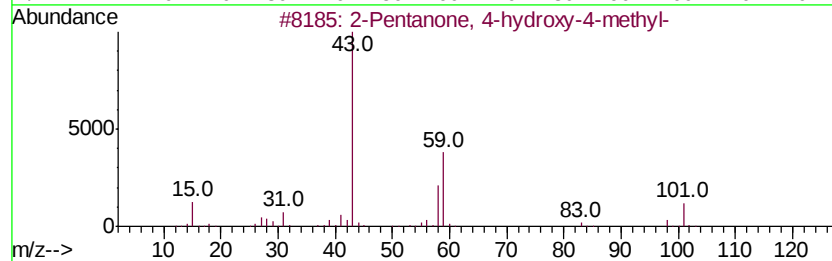
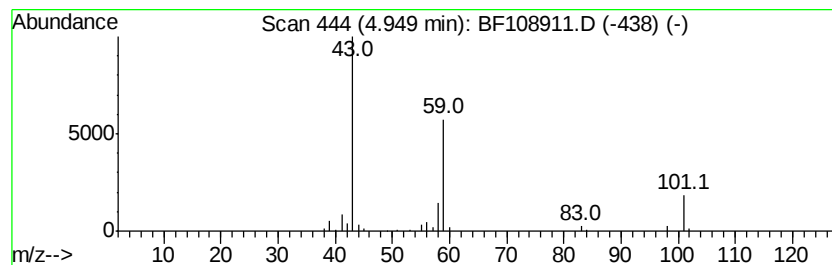
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	2.97 ng	159890	1,4-Dichlorobenzene-d4	6.74

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	27
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9



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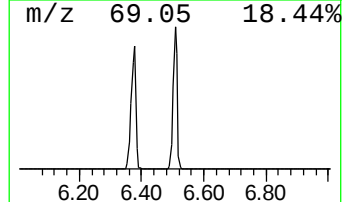
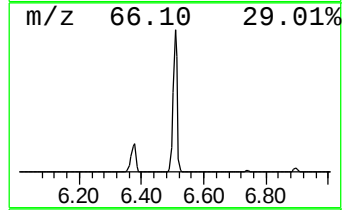
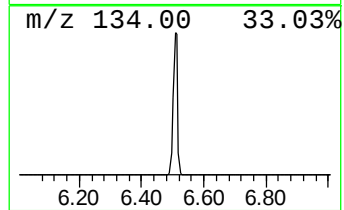
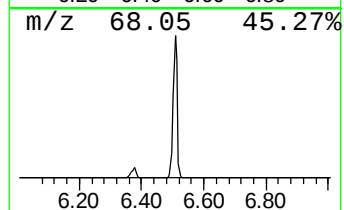
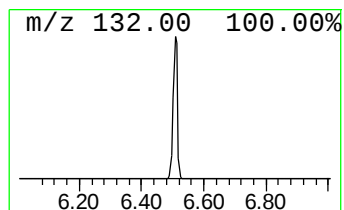
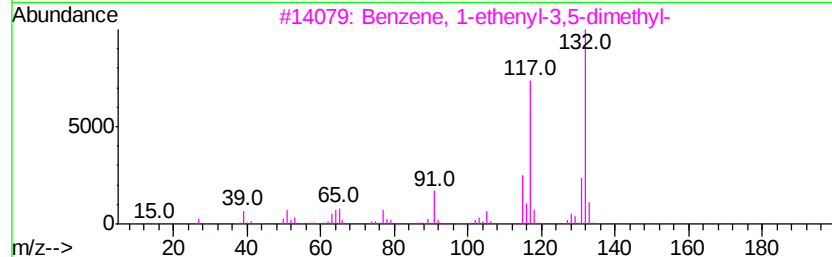
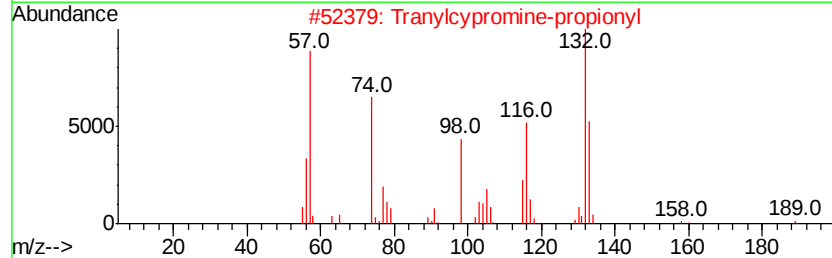
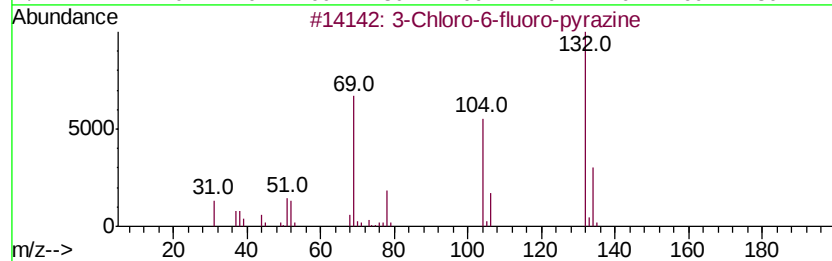
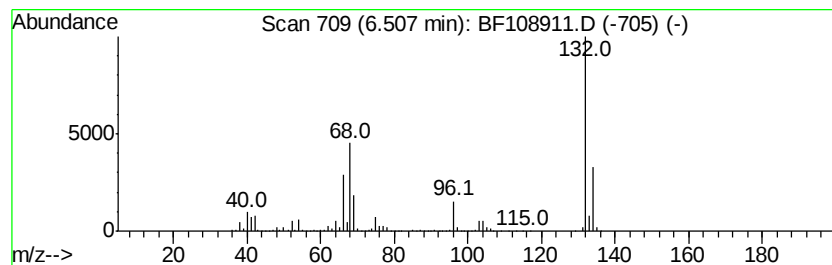
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Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	70.77 ng	3814180	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.95	3.0	ng	159890	1	6.74	1077970	20.0
unknown6.51	6.51	70.8	ng	3814180	1	6.74	1077970	20.0