

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091817\
 Data File : BF098625.D
 Acq On : 19 Sep 2017 2:40
 Operator : SJ/JU
 Sample : I5260-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 091317-TIDO-E-U-M

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:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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	4.804	458	462	475	rBV	136678	182219	4.62%	0.837%
2	5.240	533	536	551	rBV	936697	1096711	27.80%	5.037%
3	6.287	711	714	728	rBV	621111	710200	18.00%	3.262%
4	6.404	731	734	744	rVB	2081066	1897478	48.09%	8.714%
5	6.634	769	773	782	rBV	844040	699523	17.73%	3.213%
6	6.792	795	800	803	rBV	2735644	2568399	65.09%	11.795%
7	7.204	866	870	873	rBV	1923178	1937356	49.10%	8.897%
8	7.916	988	991	1006	rVB	1149791	969417	24.57%	4.452%
9	8.998	1169	1175	1178	rBV	4288266	3945623	100.00%	18.120%
10	9.392	1239	1242	1250	rVB	146498	118839	3.01%	0.546%
11	9.669	1285	1289	1298	rVB	1236158	1072747	27.19%	4.927%
12	10.457	1419	1423	1435	rVB	1365510	1201724	30.46%	5.519%
13	10.574	1440	1443	1449	rBV	38856	41233	1.05%	0.189%
14	11.151	1537	1541	1548	rBV	1031525	936724	23.74%	4.302%
15	12.733	1805	1810	1813	rBV	3267711	2930200	74.26%	13.457%
16	13.633	1960	1963	1967	rBV	45641	55199	1.40%	0.253%
17	13.786	1985	1989	1995	rVB	736439	712081	18.05%	3.270%
18	15.180	2221	2226	2235	rBV	571562	699240	17.72%	3.211%

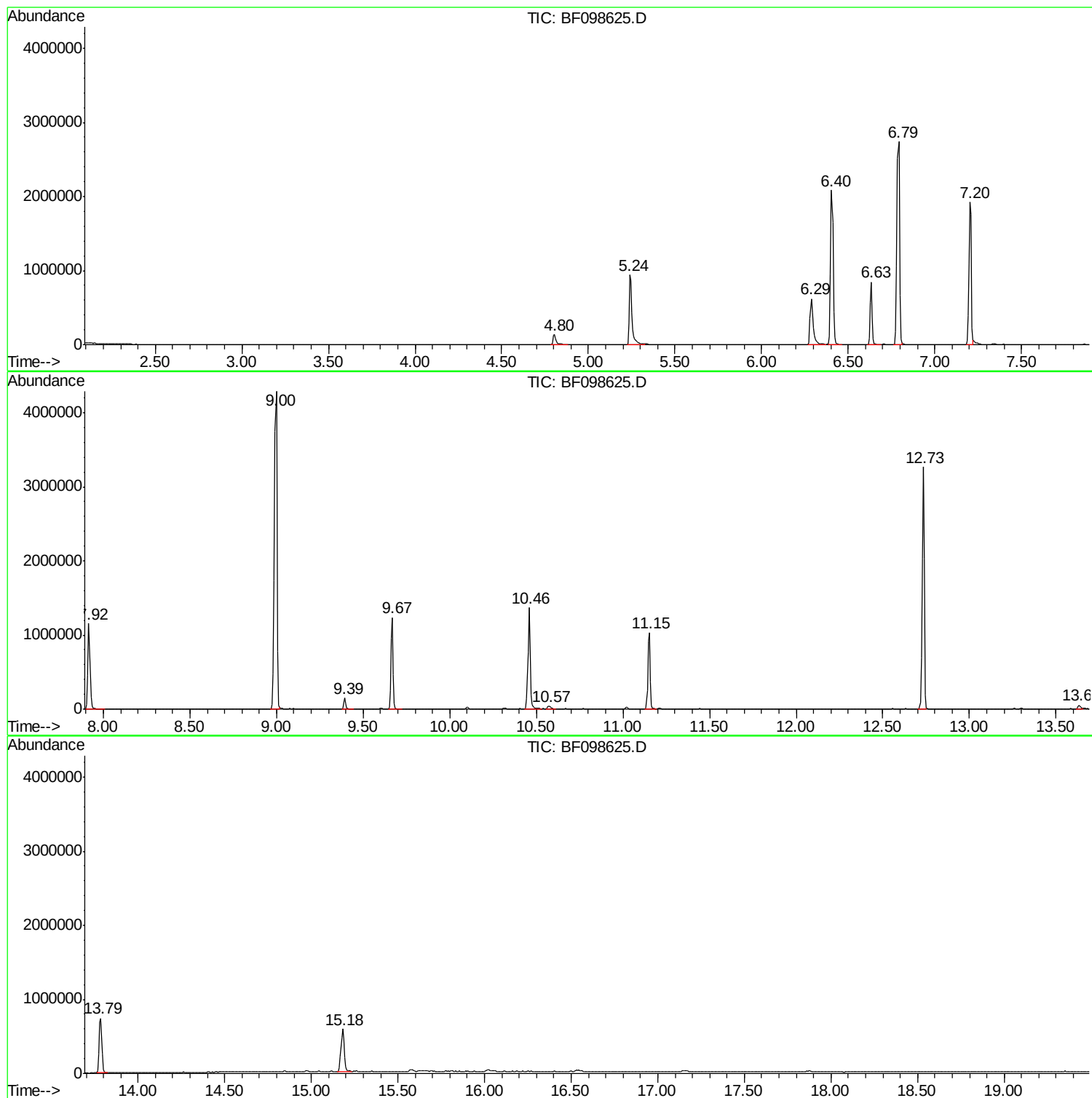
Sum of corrected areas: 21774913

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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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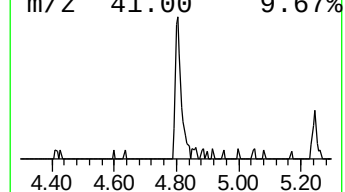
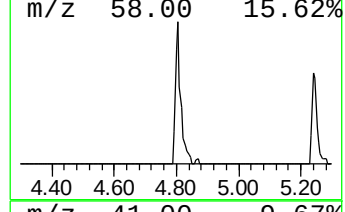
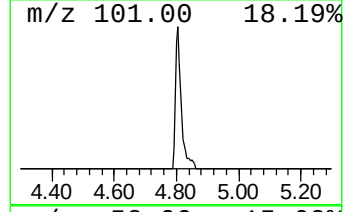
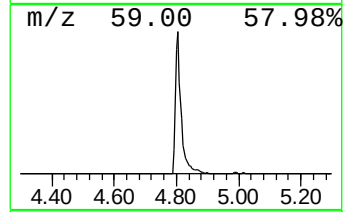
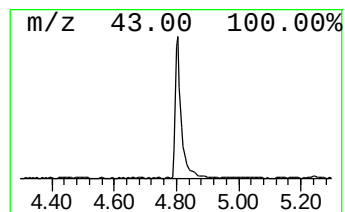
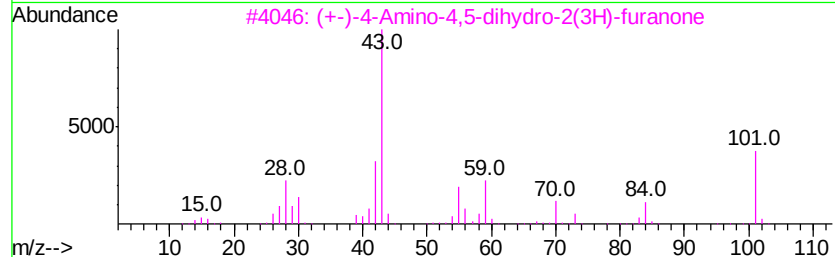
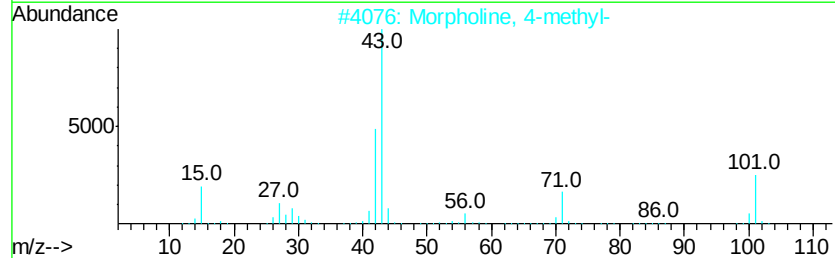
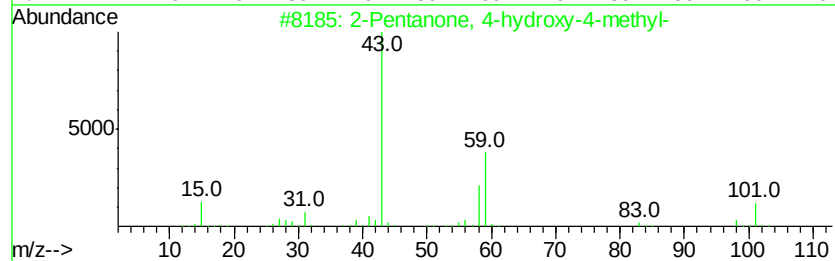
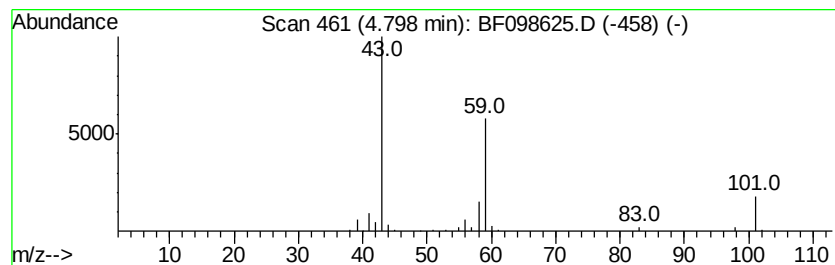
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	5.21 ng	182219	1,4-Dichlorobenzene-d4	6.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Guanidine	59	CH5N3	000113-00-8	9



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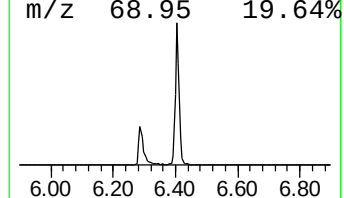
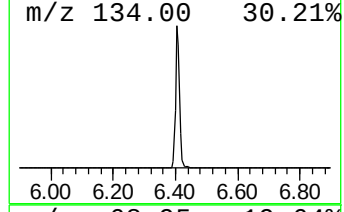
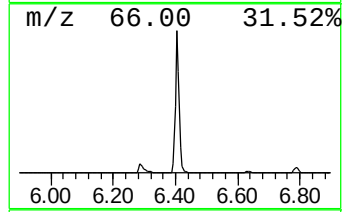
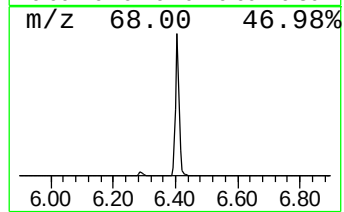
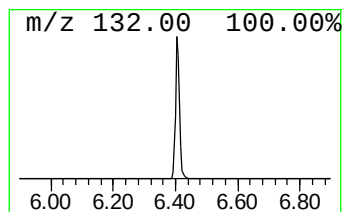
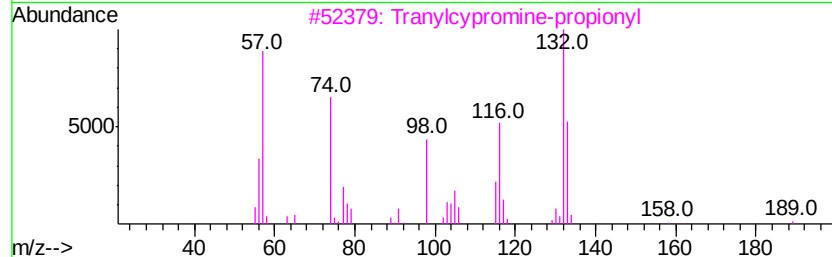
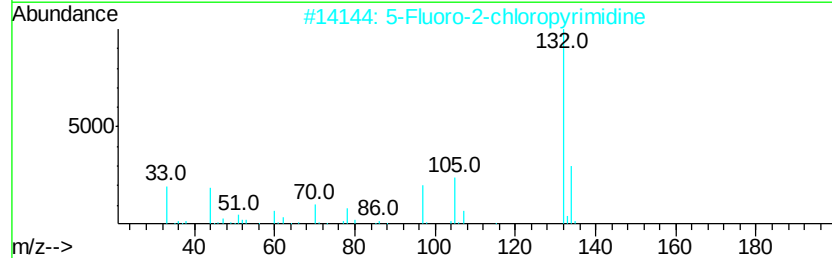
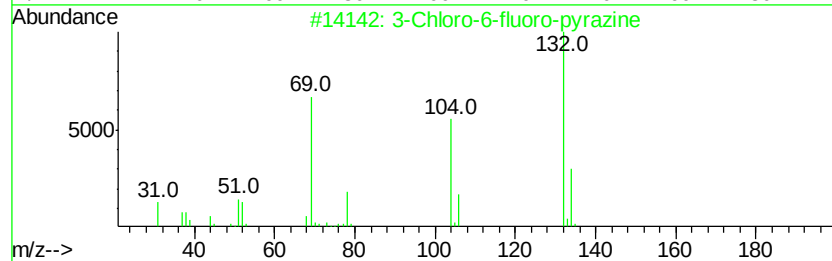
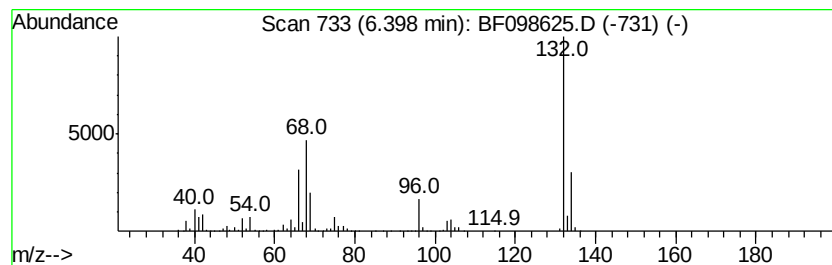
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Peak Number 2 unknown6.40 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	54.25 ng	1897480	1,4-Dichlorobenzene-d4	6.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.80	5.2	ng	182219	1	6.63	699523	20.0
unknown6.40	6.40	54.3	ng	1897480	1	6.63	699523	20.0