

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091817\
 Data File : BF098627.D
 Acq On : 19 Sep 2017 3:37
 Operator : SJ/JU
 Sample : I5262-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 091417-SIDI-E-D-S

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	4.798	459	461	473	rBV	139761	187422	5.12%	0.902%
2	5.240	533	536	551	rBV	965852	1035190	28.25%	4.984%
3	6.287	711	714	729	rBV	708633	771102	21.04%	3.713%
4	6.404	731	734	749	rVB	2211207	1902522	51.92%	9.160%
5	6.634	769	773	782	rBV	814809	683193	18.65%	3.289%
6	6.793	795	800	803	rBV	2536116	2459151	67.11%	11.840%
7	7.204	866	870	873	rBV	1942953	1901289	51.89%	9.154%
8	7.916	988	991	999	rBV	1164256	956631	26.11%	4.606%
9	8.998	1169	1175	1178	rBV	3964012	3664138	100.00%	17.642%
10	9.392	1239	1242	1249	rBV	122151	96350	2.63%	0.464%
11	9.669	1285	1289	1295	rVB	969435	851246	23.23%	4.098%
12	10.457	1419	1423	1433	rVB	1159348	1004364	27.41%	4.836%
13	11.151	1537	1541	1544	rBV	990580	900875	24.59%	4.337%
14	12.733	1805	1810	1813	rBV	3106065	2728529	74.47%	13.137%
15	13.633	1957	1963	1971	rBV2	43203	60121	1.64%	0.289%
16	13.780	1984	1988	1995	rVB	694101	675320	18.43%	3.251%
17	15.180	2219	2226	2236	rVB	581307	722027	19.71%	3.476%
18	15.574	2289	2293	2298	rVB2	35176	49514	1.35%	0.238%
19	16.021	2364	2369	2376	rBV2	36136	67135	1.83%	0.323%
20	16.539	2453	2457	2463	rVB6	25583	53706	1.47%	0.259%

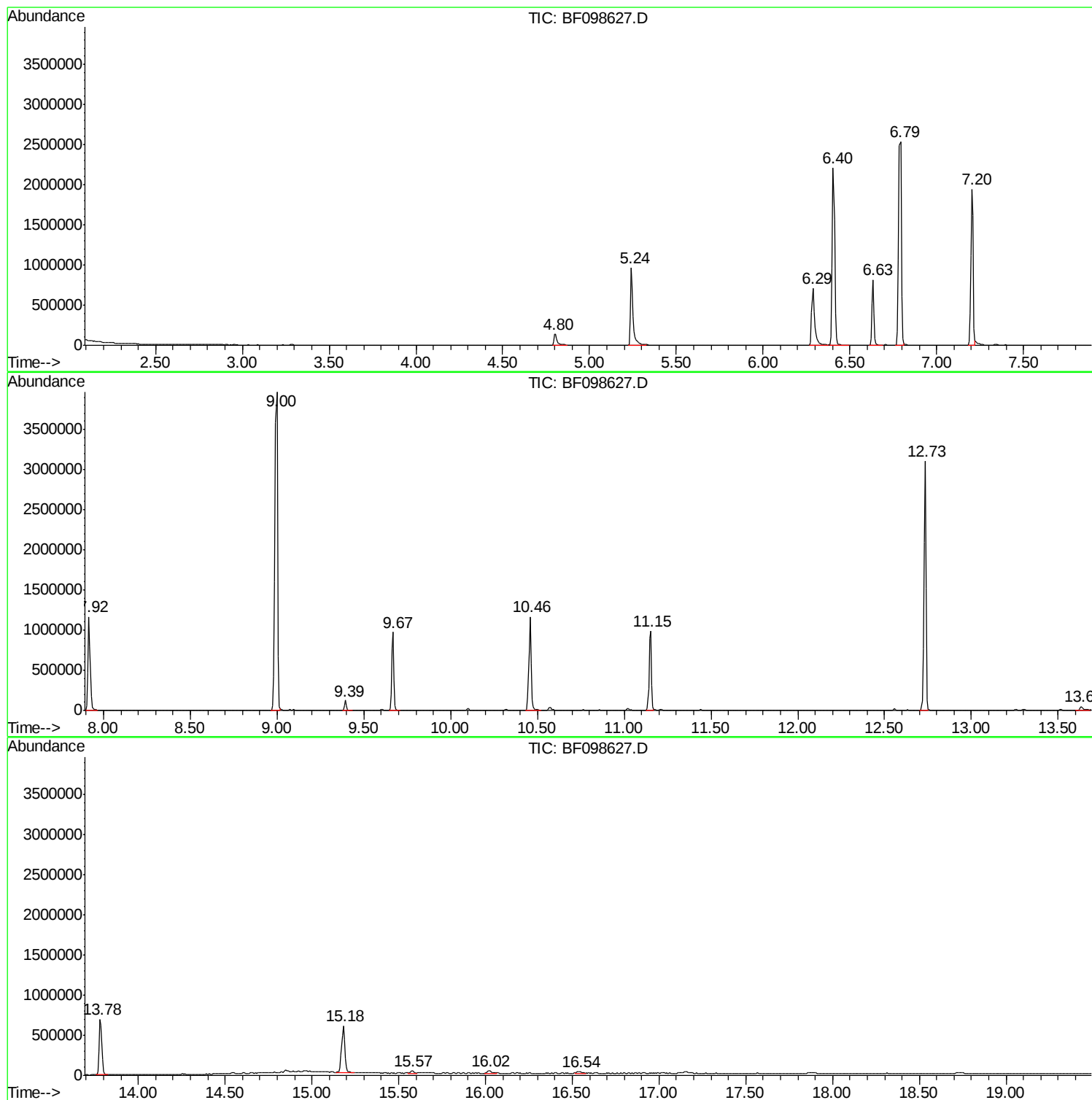
Sum of corrected areas: 20769825

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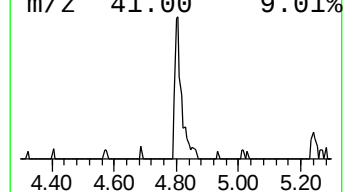
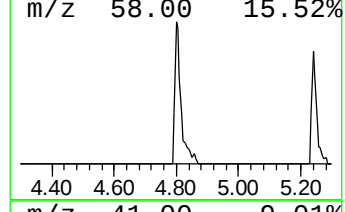
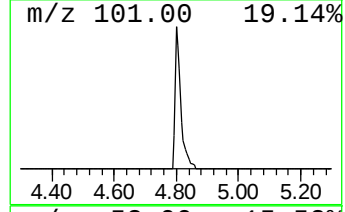
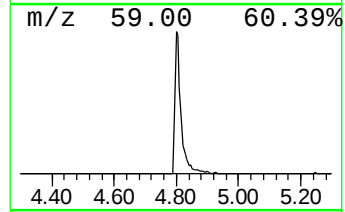
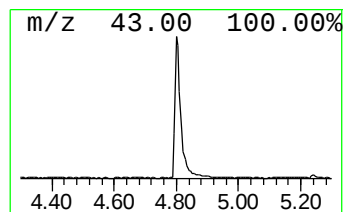
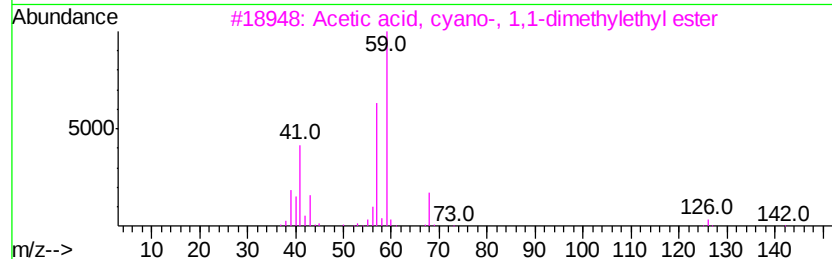
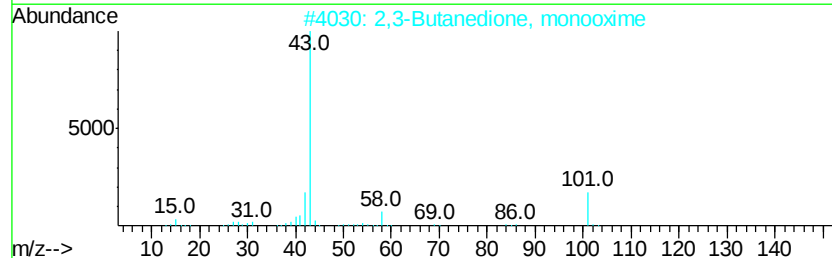
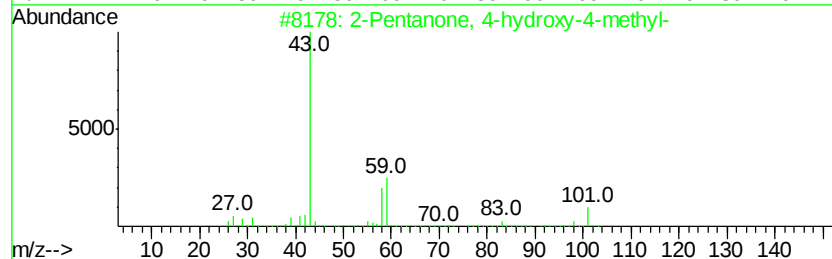
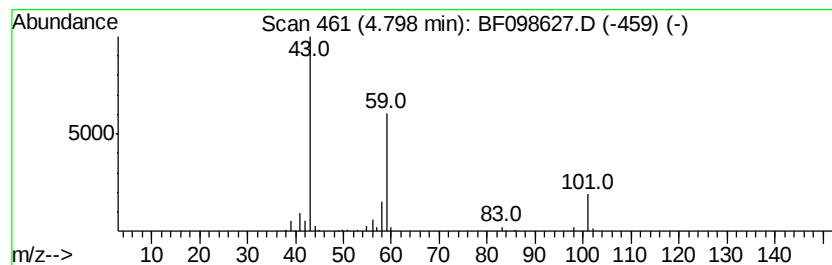
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	5.49 ng	187422	1,4-Dichlorobenzene-d4	6.63

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Guanidine	59	CH5N3	000113-00-8	9



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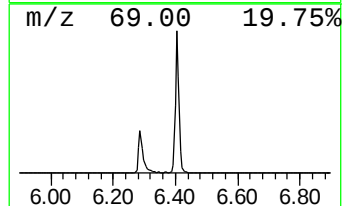
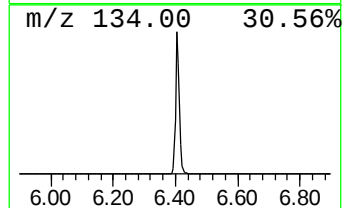
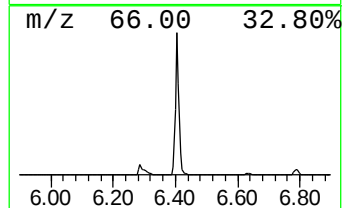
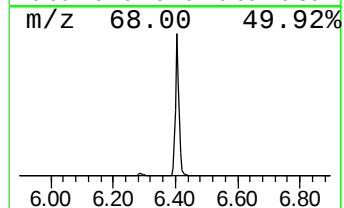
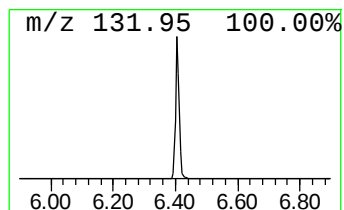
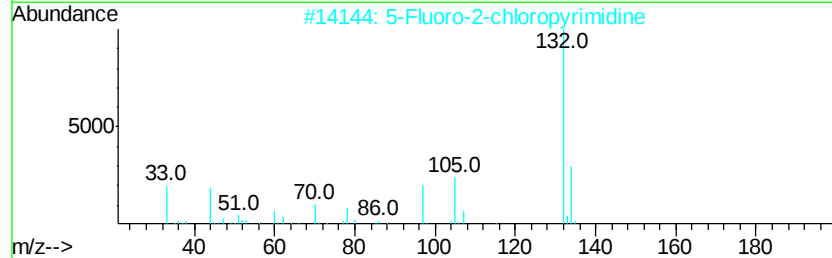
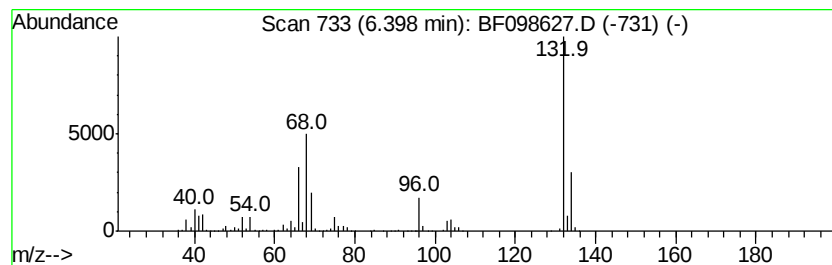
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Peak Number 2 unknown6.40 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	55.70 ng	1902520	1,4-Dichlorobenzene-d4	6.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16	
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10	
4	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9	
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.80	5.5	ng	187422	1	6.63	683193	20.0
unknown6.40	6.40	55.7	ng	1902520	1	6.63	683193	20.0