

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091817\
 Data File : BF098626.D
 Acq On : 19 Sep 2017 3:09
 Operator : SJ/JU
 Sample : I5262-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

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

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	474	rBV	128104	174633	5.33%	0.881%
2	5.239	533	536	557	rBV	855275	1031864	31.47%	5.205%
3	6.286	711	714	728	rBV	620752	699137	21.32%	3.527%
4	6.404	731	734	742	rVB	2075963	1852449	56.49%	9.344%
5	6.634	769	773	783	rBV	737652	625393	19.07%	3.155%
6	6.792	795	800	803	rBV	2351593	2252736	68.69%	11.363%
7	7.204	866	870	873	rBV	1756094	1793674	54.70%	9.047%
8	7.916	988	991	1007	rVB	945310	813919	24.82%	4.106%
9	8.998	1170	1175	1178	rBV	3657898	3279377	100.00%	16.542%
10	9.392	1239	1242	1250	rVB	132022	103916	3.17%	0.524%
11	9.669	1285	1289	1299	rVB	978786	854834	26.07%	4.312%
12	10.457	1419	1423	1435	rVB	1236341	1074546	32.77%	5.420%
13	10.574	1439	1443	1450	rBV	37450	39356	1.20%	0.199%
14	11.151	1537	1541	1547	rBV	1049643	946519	28.86%	4.774%
15	12.733	1805	1810	1813	rBV	3199973	2809996	85.69%	14.174%
16	13.633	1957	1963	1969	rBV2	38067	51211	1.56%	0.258%
17	13.780	1985	1988	1995	rVB	743418	711575	21.70%	3.589%
18	15.180	2221	2226	2237	rBV	594824	709944	21.65%	3.581%

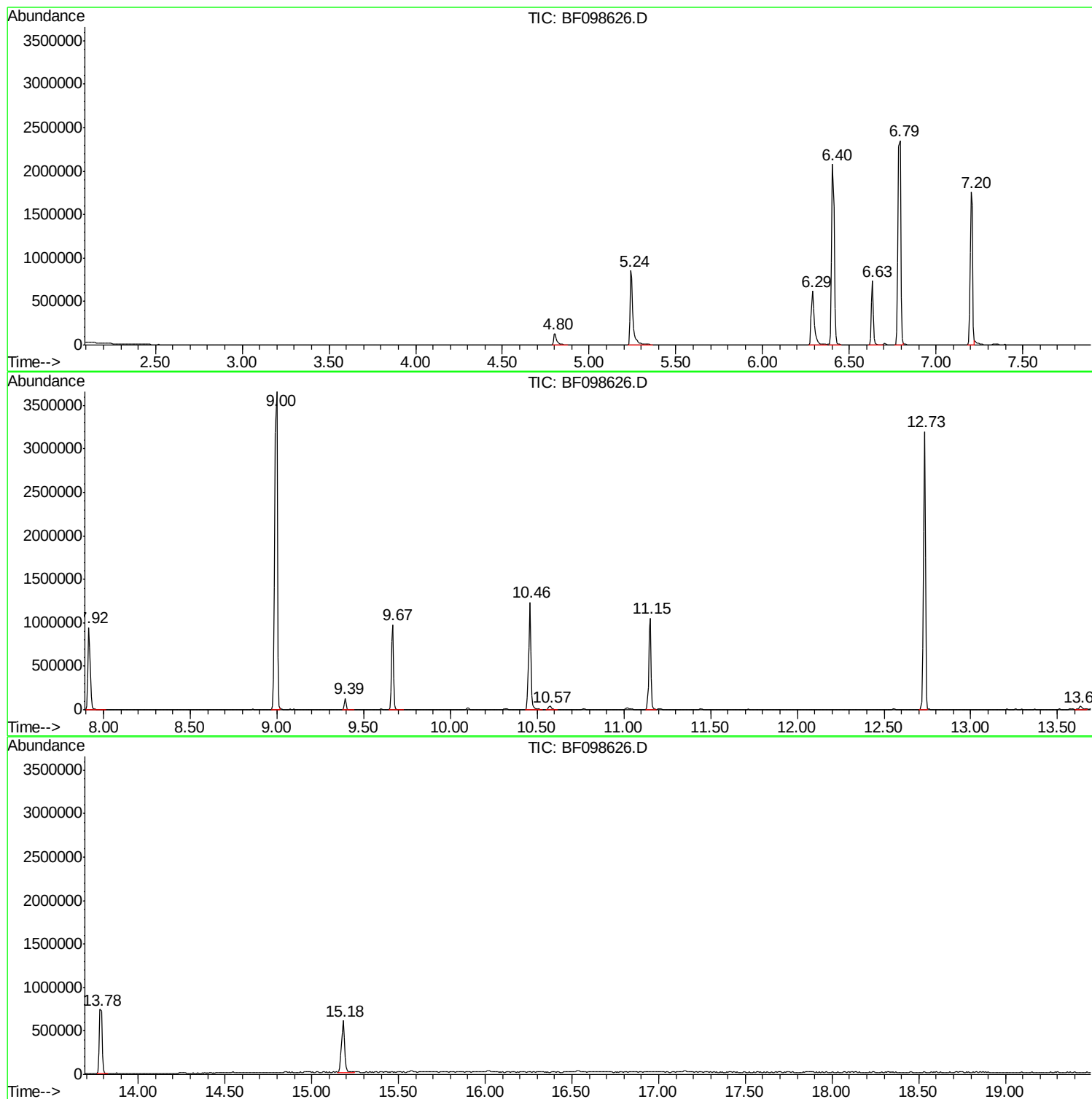
Sum of corrected areas: 19825079

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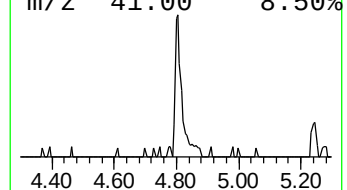
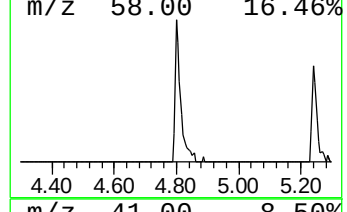
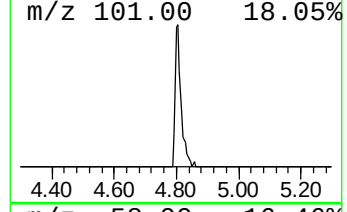
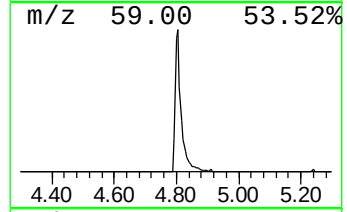
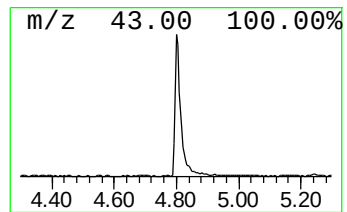
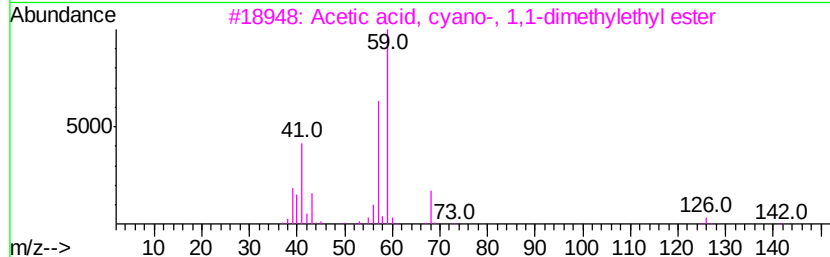
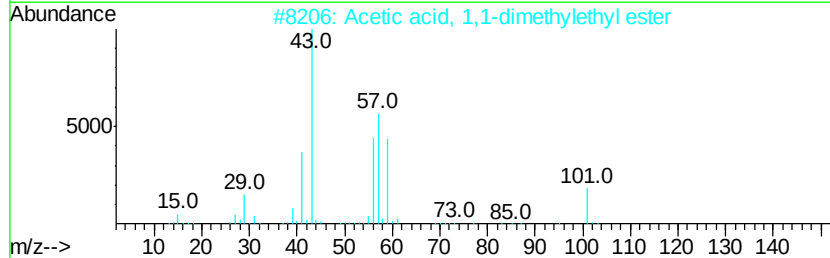
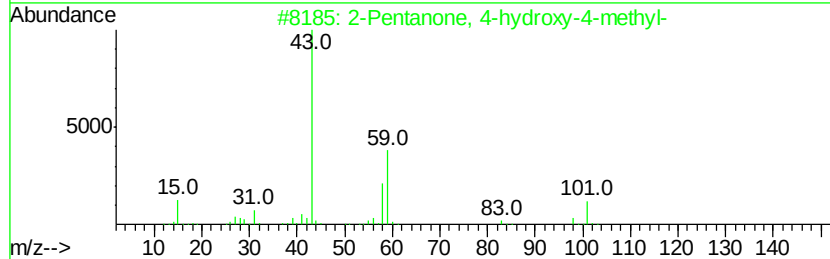
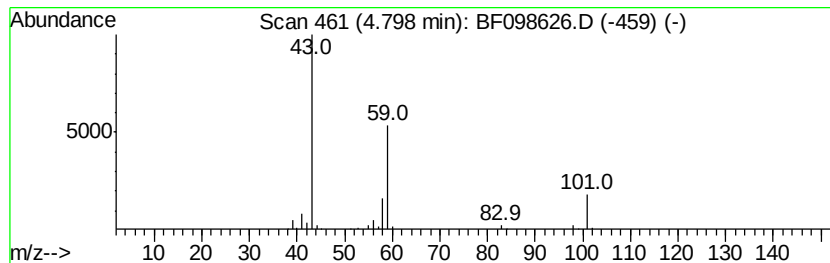
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	5.58 ng	174633	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	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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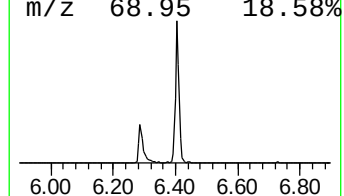
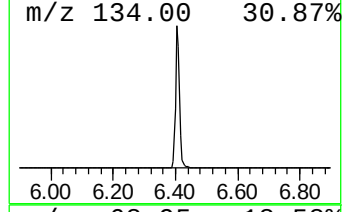
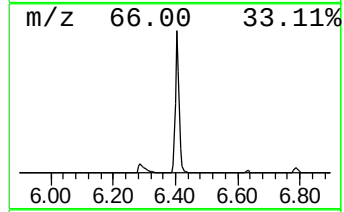
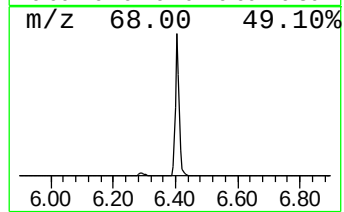
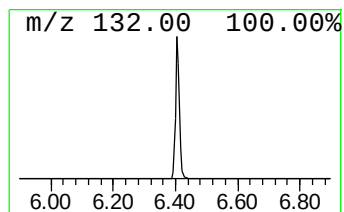
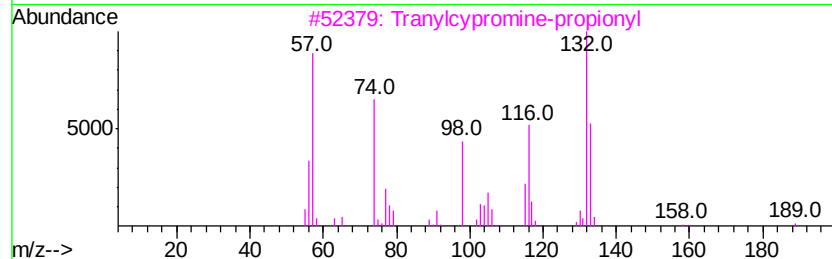
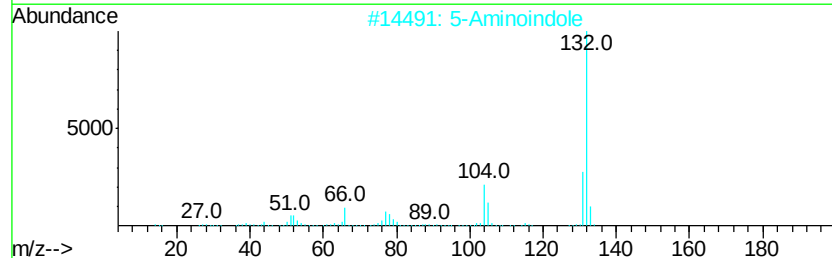
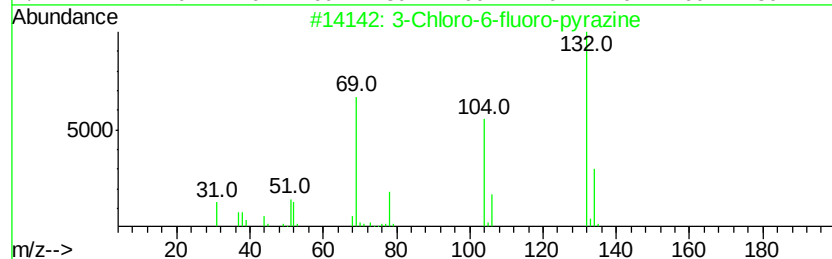
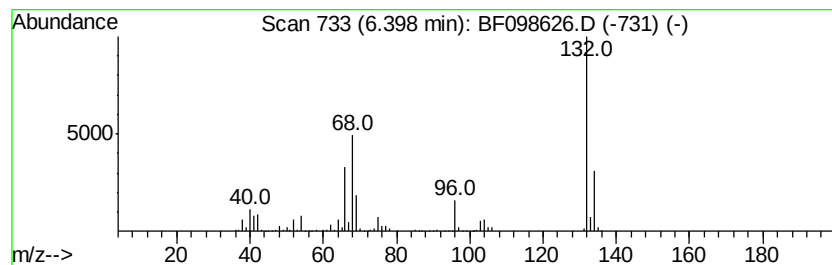
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Peak Number 2 unknown6.40 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	59.24 ng	1852450	1,4-Dichlorobenzene-d4	6.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
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...	4.80	5.6	ng	174633	1	6.63	625393	20.0
unknown6.40	6.40	59.2	ng	1852450	1	6.63	625393	20.0