

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050517\
 Data File : BF094925.D
 Acq On : 5 May 2017 19:37
 Operator : SJ/MA
 Sample : PB98685BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98685BL

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-BF050217.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	5.054	526	530	548	rBV	182510	360971	11.55%	1.484%
2	5.678	632	636	661	rBV	1788762	2649802	84.79%	10.893%
3	7.266	901	906	921	rBV	1680862	2441840	78.13%	10.038%
4	7.390	921	927	944	rVB	2102404	2778050	88.89%	11.420%
5	7.725	980	984	994	rVB	482770	548257	17.54%	2.254%
6	7.960	1019	1024	1036	rBV	1812541	2223508	71.15%	9.141%
7	8.625	1131	1137	1161	rBV	1240220	1678089	53.69%	6.899%
8	9.742	1322	1327	1342	rBV	539861	730334	23.37%	3.002%
9	11.519	1623	1629	1646	rBV	2376187	3125277	100.00%	12.848%
10	12.572	1803	1808	1821	rBV2	558510	841767	26.93%	3.460%
11	13.866	2022	2028	2050	rBV	1232447	1835241	58.72%	7.545%
12	14.971	2211	2216	2230	rBV	564004	829952	26.56%	3.412%
13	17.301	2606	2612	2630	rBV	2530404	3119928	99.83%	12.826%
14	18.636	2835	2839	2854	rBV	489185	688297	22.02%	2.830%
15	20.306	3119	3123	3137	rBV	281791	473884	15.16%	1.948%

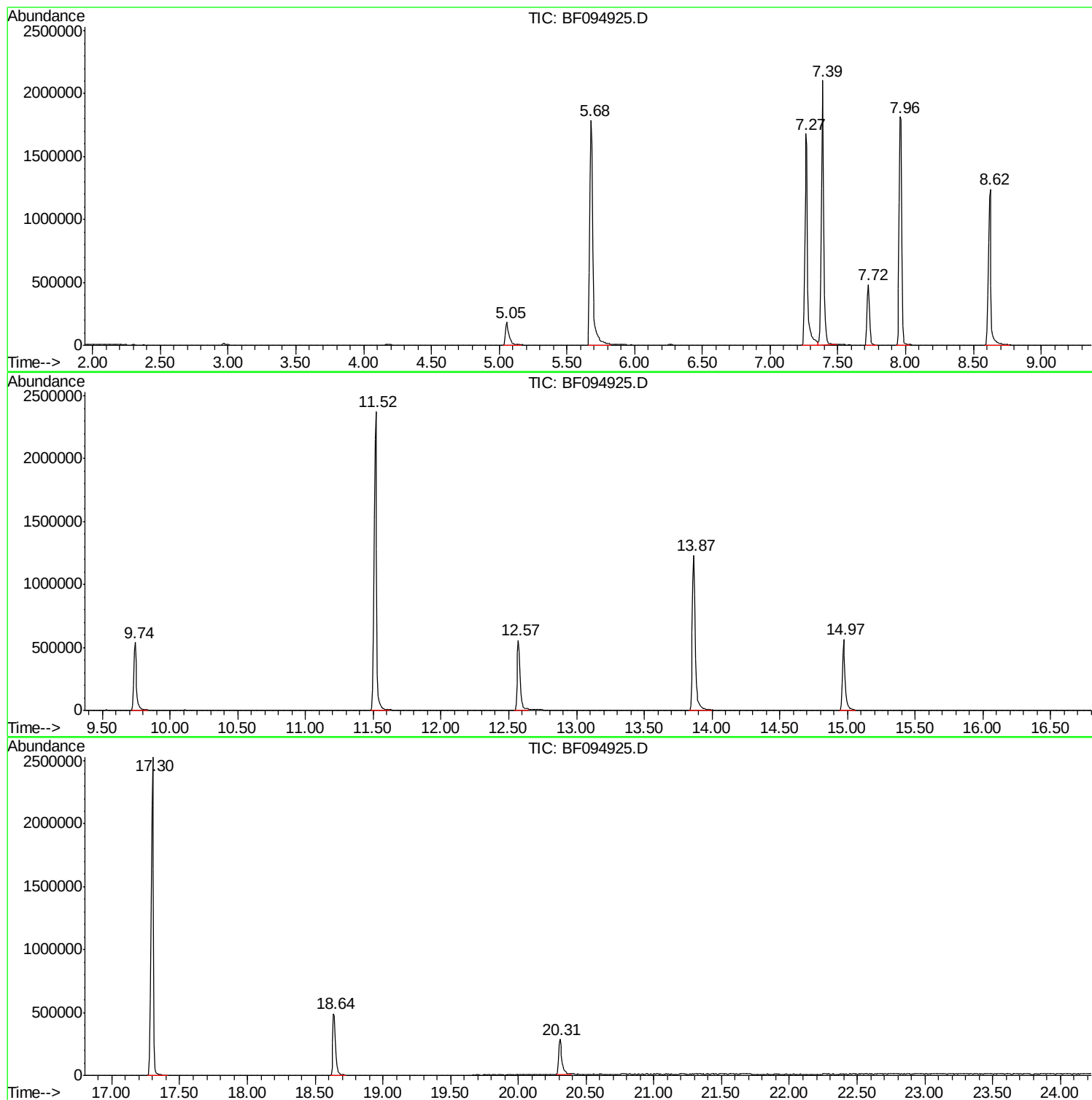
Sum of corrected areas: 24325197

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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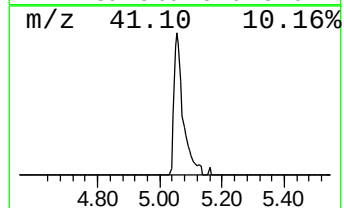
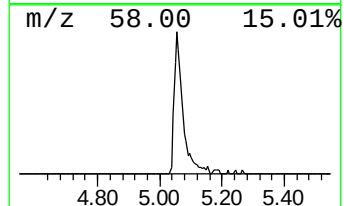
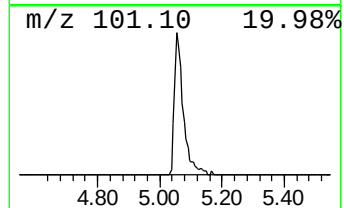
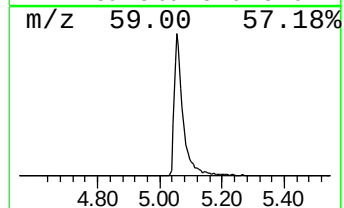
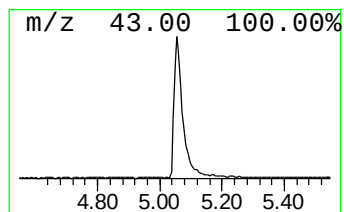
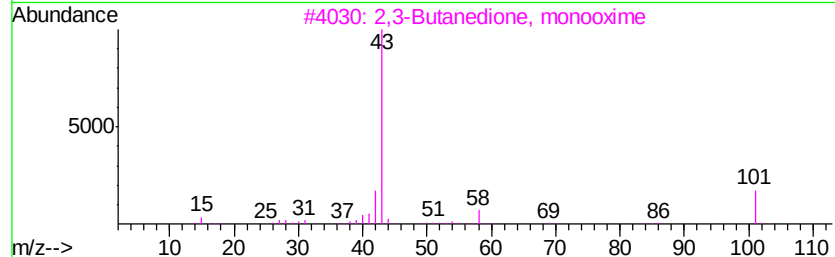
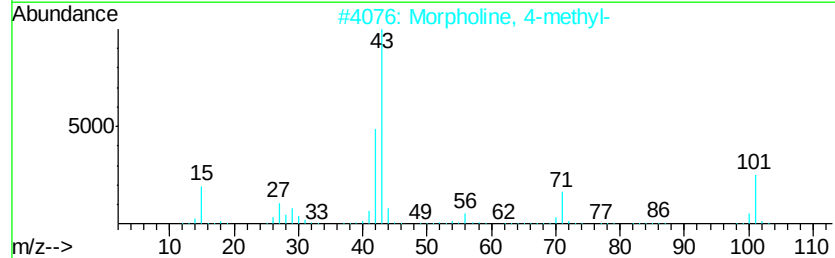
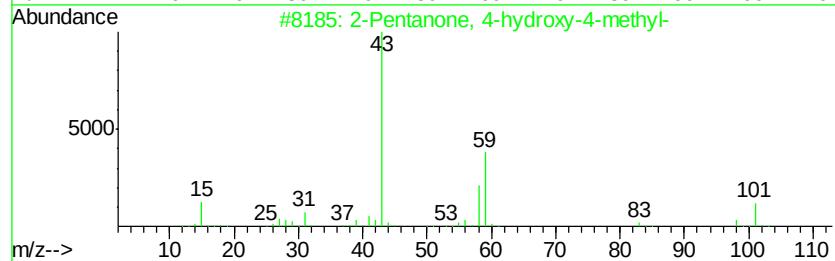
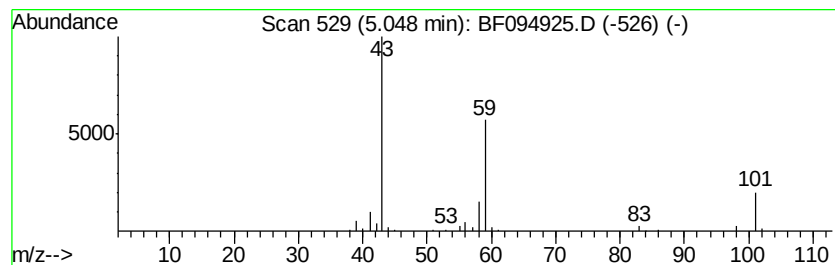
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	13.17 ng	360971	1,4-Dichlorobenzene-d4	7.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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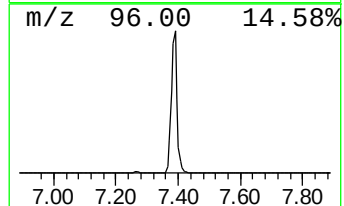
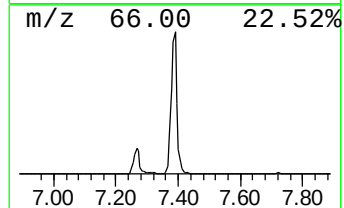
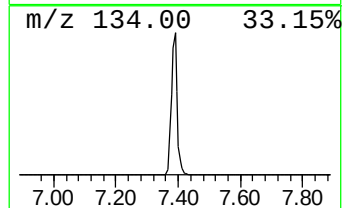
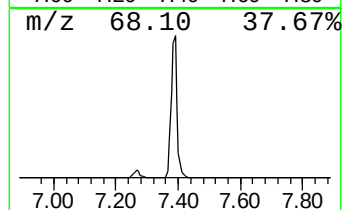
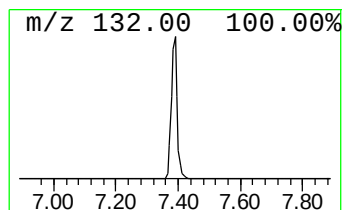
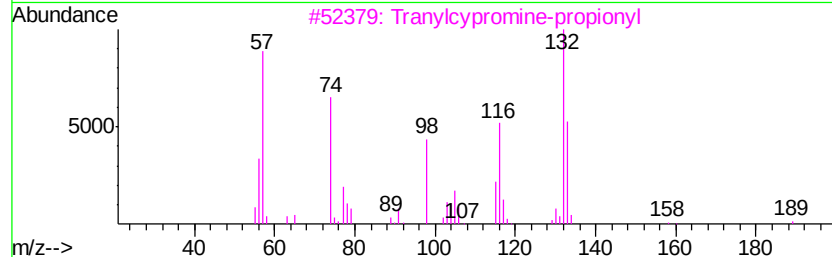
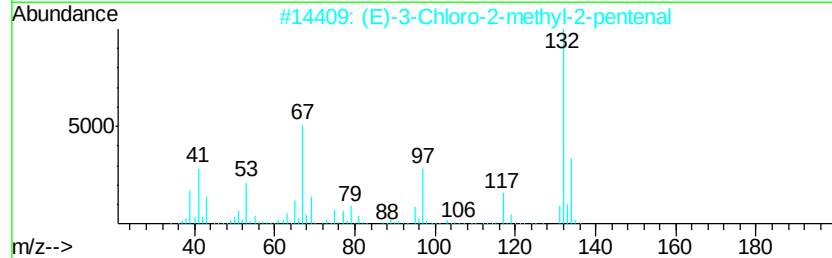
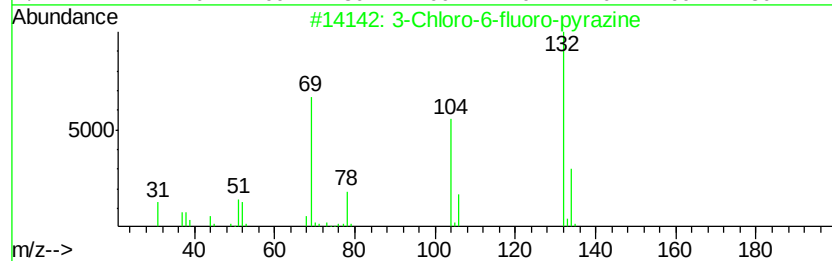
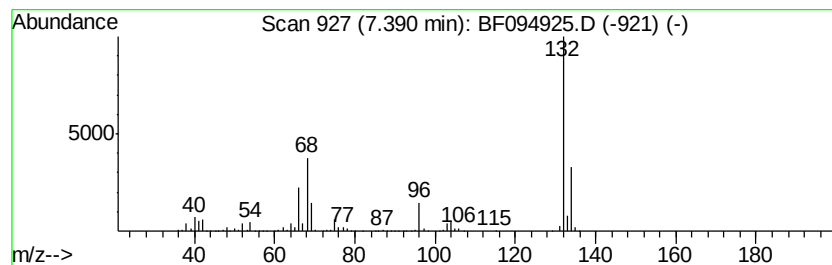
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Peak Number 2 unknown7.39 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.39	101.34 ng	2778050	1,4-Dichlorobenzene-d4	7.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.05	13.2	ng	360971	1	7.72	548257	20.0
unknown7.39	7.39	101.3	ng	2778050	1	7.72	548257	20.0