

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051117\
 Data File : BF095084.D
 Acq On : 11 May 2017 17:27
 Operator : SJ/MA
 Sample : PB98865BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98865BL

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	5.013	519	523	540	rBV	128207	284993	9.79%	1.297%
2	5.648	627	631	665	rBV	1200451	2302659	79.10%	10.481%
3	7.248	898	903	918	rBV	1378198	2133083	73.28%	9.709%
4	7.366	918	923	938	rVB	1753004	2538644	87.21%	11.556%
5	7.701	976	980	997	rVB	427775	513363	17.64%	2.337%
6	7.942	1016	1021	1032	rBV	1515994	1899969	65.27%	8.648%
7	8.601	1129	1133	1165	rBV	937934	1444298	49.62%	6.574%
8	9.725	1320	1324	1345	rBV	383623	681862	23.42%	3.104%
9	11.495	1620	1625	1643	rBV	2119625	2910890	100.00%	13.250%
10	12.554	1801	1805	1820	rBV2	408086	701673	24.11%	3.194%
11	13.848	2020	2025	2049	rBV	905301	1566707	53.82%	7.131%
12	14.960	2209	2214	2233	rBV	485071	814122	27.97%	3.706%
13	17.283	2604	2609	2622	rBV	2246431	2825436	97.06%	12.861%
14	18.630	2834	2838	2854	rBV	497569	788517	27.09%	3.589%
15	20.300	3117	3122	3134	rBV	324344	562924	19.34%	2.562%

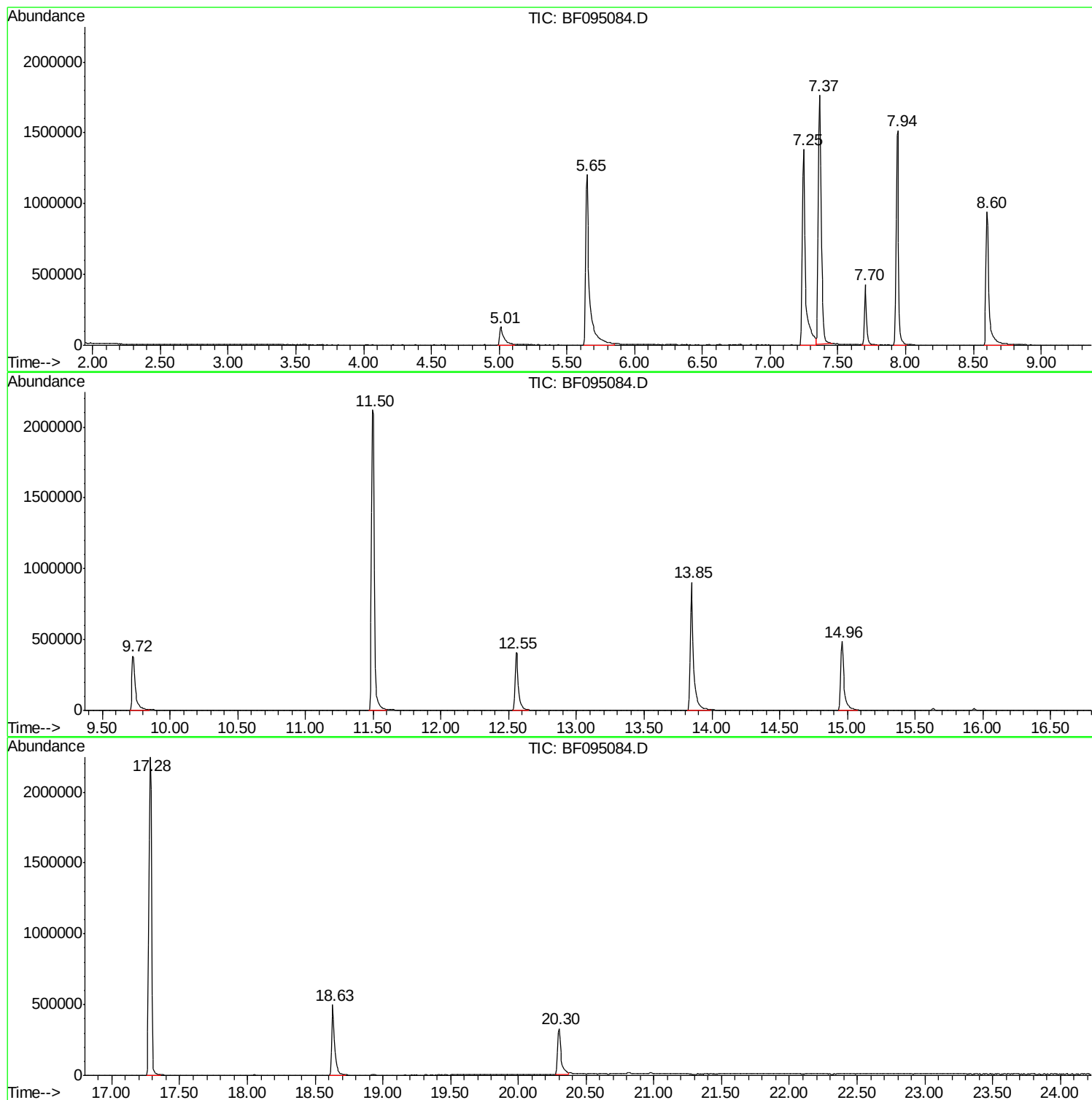
Sum of corrected areas: 21969140

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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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Instrument :
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ClientSampled :
PB98865BL

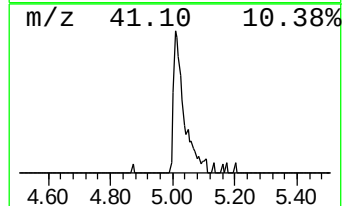
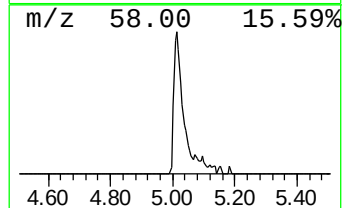
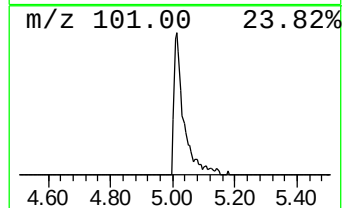
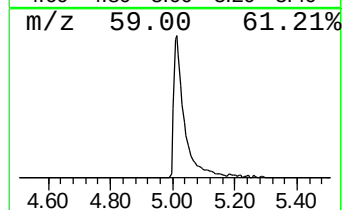
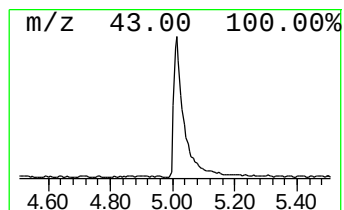
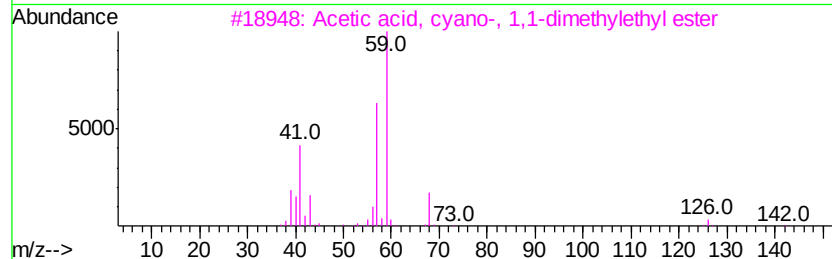
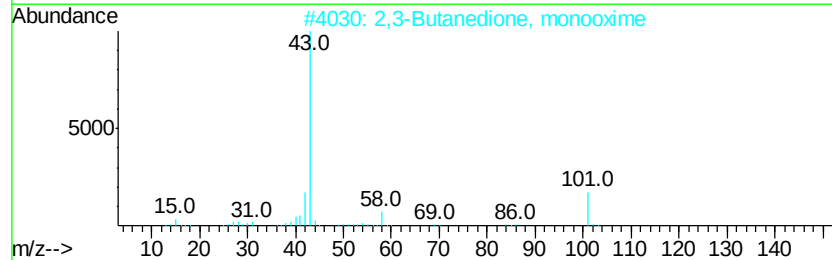
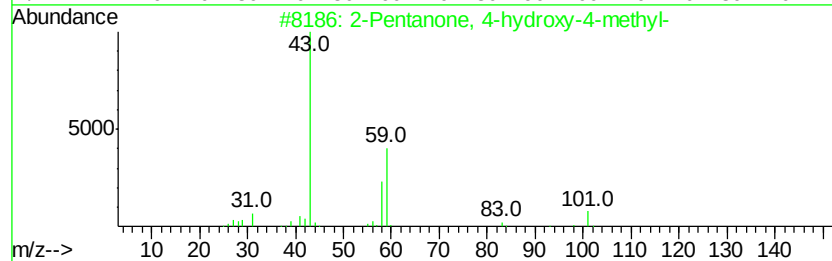
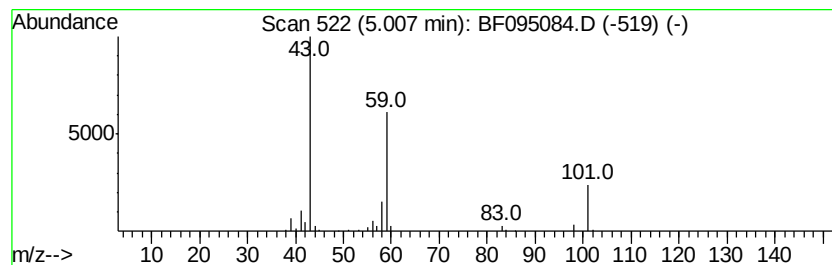
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	11.10 ng	284993	1,4-Dichlorobenzene-d4	7.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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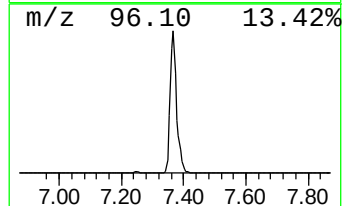
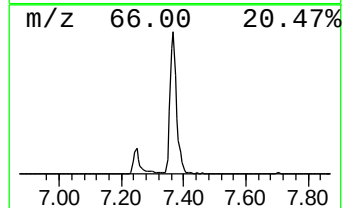
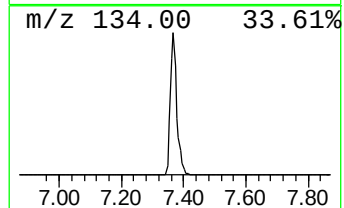
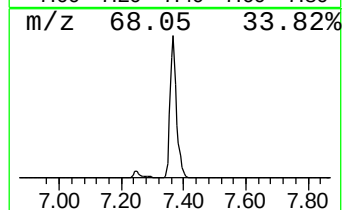
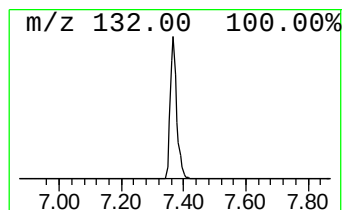
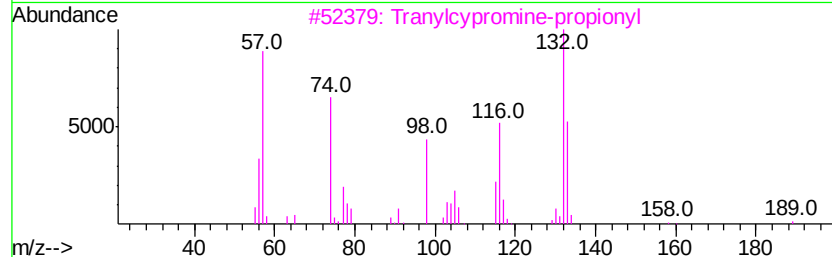
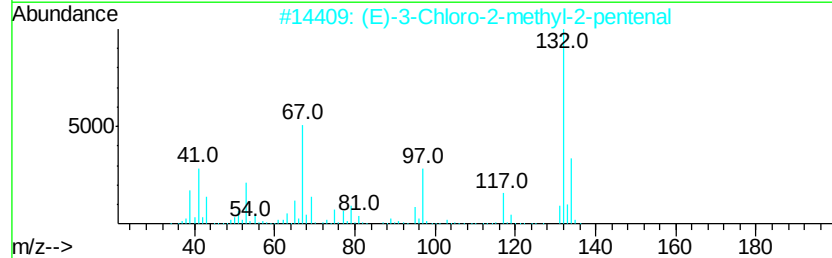
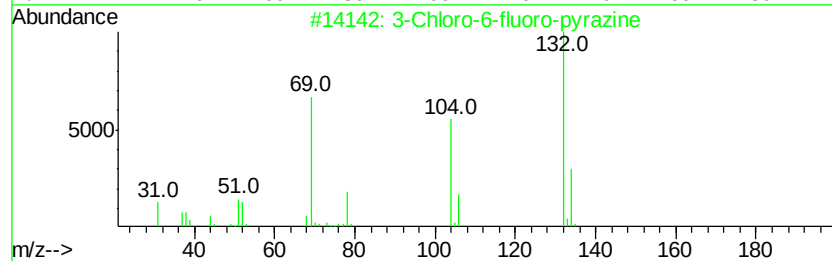
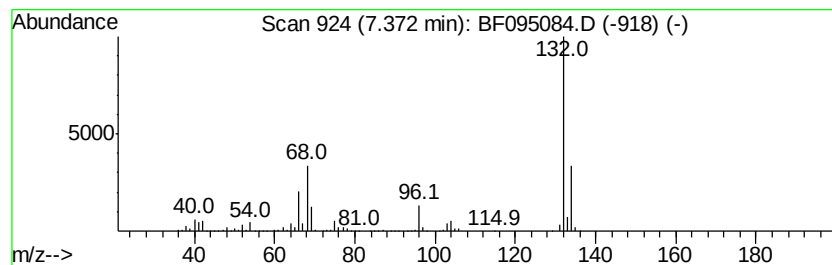
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Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	98.90 ng	2538640	1,4-Dichlorobenzene-d4	7.70

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
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...	5.01	11.1	ng	284993	1	7.70	513363	20.0
unknown7.37	7.37	98.9	ng	2538640	1	7.70	513363	20.0