

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050117\
 Data File : BF094756.D
 Acq On : 1 May 2017 18:38
 Operator : SJ/MA
 Sample : PB98574BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98574BL

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-BF041717.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	3.937	336	340	351	rBV	290181	353391	12.34%	1.635%
2	4.331	403	407	417	rBV	2421527	2461227	85.93%	11.385%
3	5.401	584	589	592	rBV	2235959	2268743	79.21%	10.494%
4	5.525	605	610	613	rBV	2497027	2535084	88.51%	11.726%
5	5.766	647	651	655	rBV	674468	612992	21.40%	2.835%
6	5.942	676	681	684	rBV	2327794	2098990	73.29%	9.709%
7	6.413	756	761	764	rBV	1507147	1559063	54.43%	7.212%
8	7.254	900	904	918	rBV	880480	820064	28.63%	3.793%
9	8.578	1124	1129	1132	rBV	2965678	2864146	100.00%	13.248%
10	9.389	1263	1267	1283	rBV2	815460	831267	29.02%	3.845%
11	10.366	1429	1433	1449	rBV	1121024	1267223	44.24%	5.862%
12	11.213	1573	1577	1587	rBV	712897	705810	24.64%	3.265%
13	13.189	1908	1913	1917	rBV	1943736	2137810	74.64%	9.889%
14	14.465	2125	2130	2137	rBV	527568	568628	19.85%	2.630%
15	16.095	2402	2407	2419	rVB	443034	534312	18.66%	2.472%

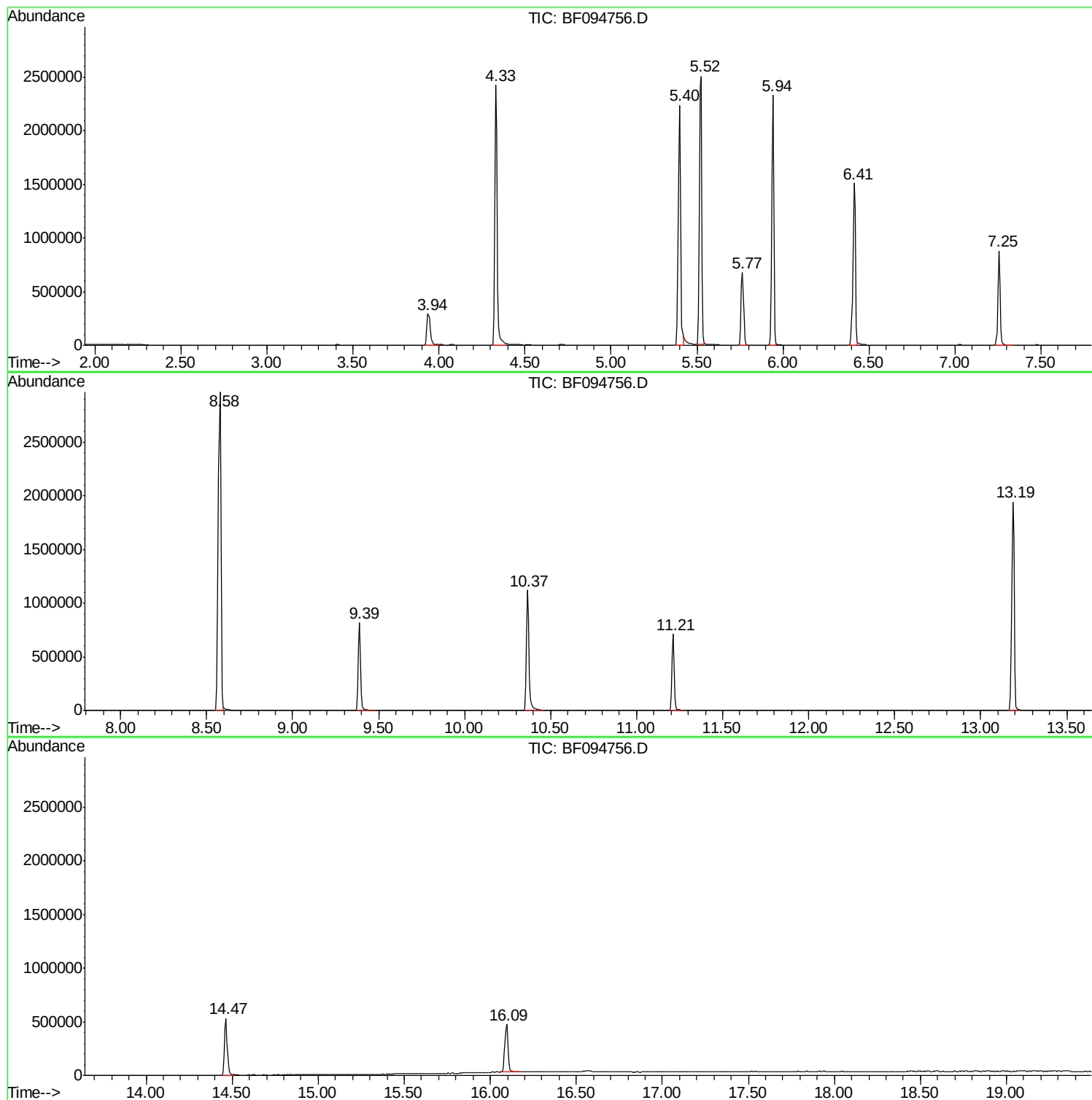
Sum of corrected areas: 21618750

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.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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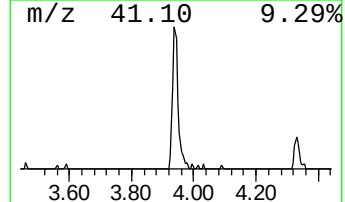
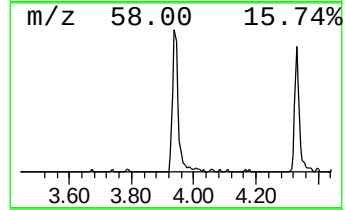
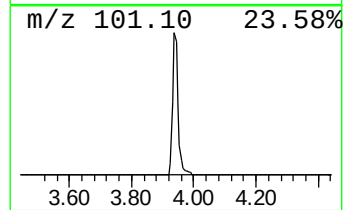
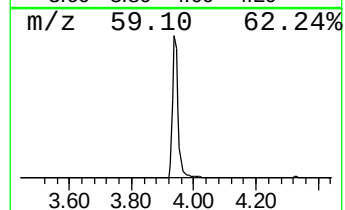
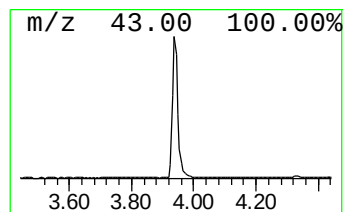
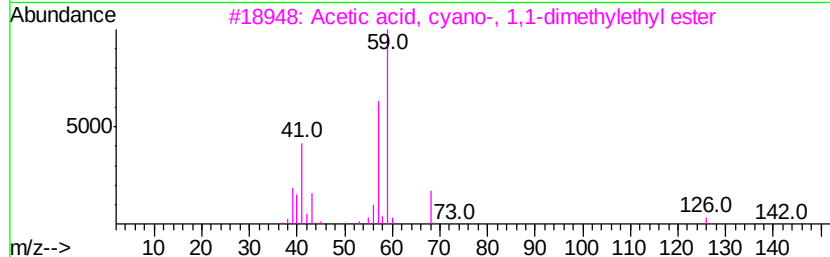
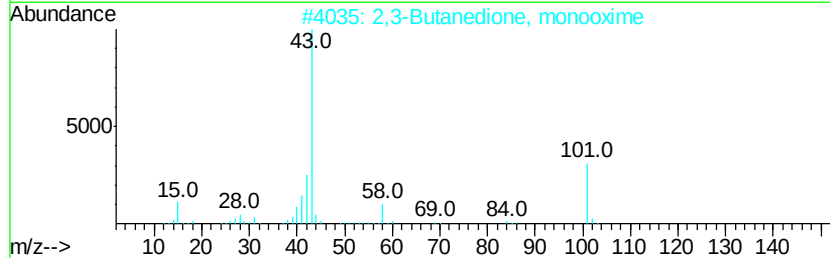
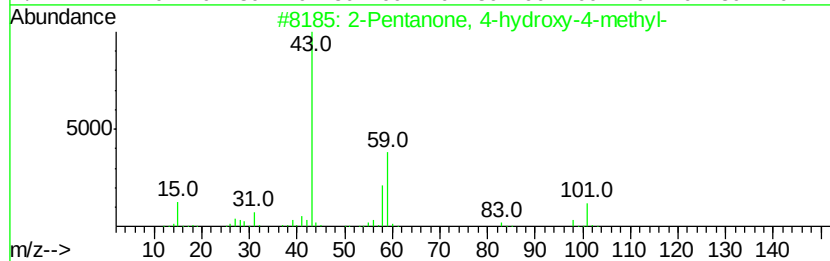
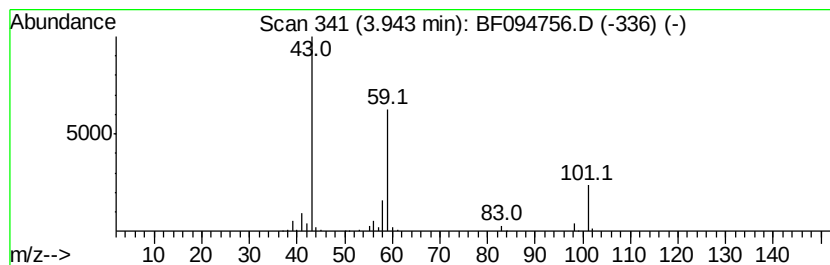
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	11.53 ng	353391	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
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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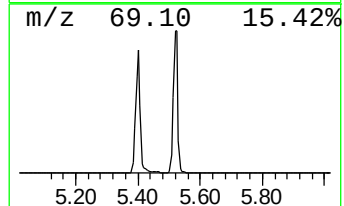
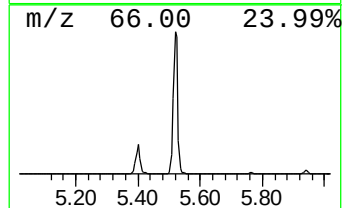
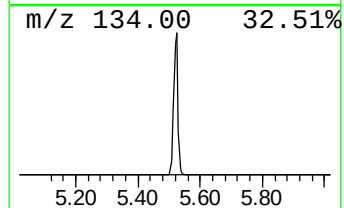
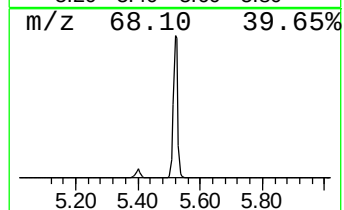
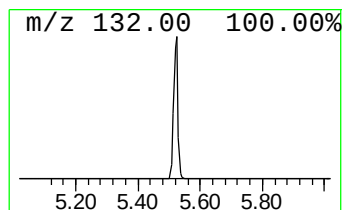
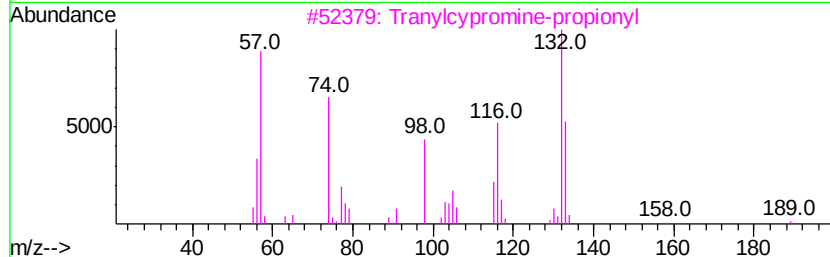
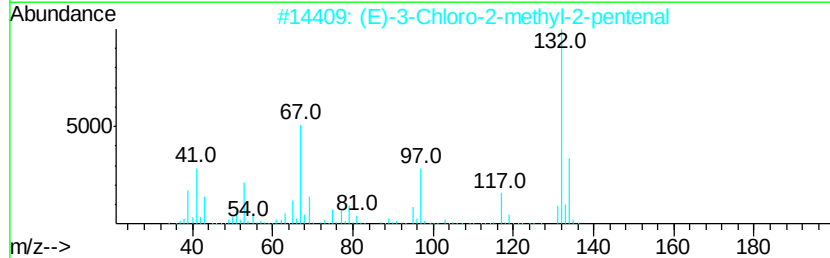
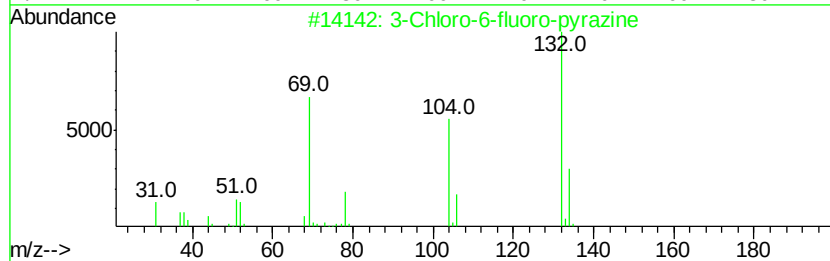
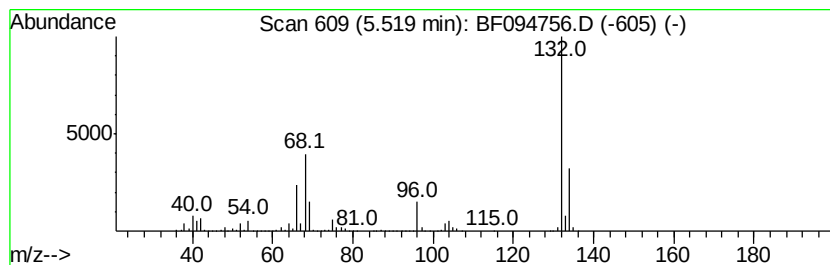
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Peak Number 2 unknown5.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	82.71 ng	2535080	1,4-Dichlorobenzene-d4	5.77

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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	3.94	11.5	ng	353391	1	5.77	612992	20.0
unknown5.52	5.52	82.7	ng	2535080	1	5.77	612992	20.0