

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100617\
Data File : BF099153.D
Acq On : 6 Oct 2017 17:56
Operator : SJ/JU
Sample : PB102940BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB102940BL

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-BF092317.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	5.104	509	513	523	rBV	413714	573256	16.03%	2.032%
2	5.492	574	579	582	rBV	2810296	3045962	85.18%	10.797%
3	6.498	744	750	753	rBV	2760354	2921673	81.71%	10.356%
4	6.634	768	773	777	rBV	2985618	3094614	86.54%	10.969%
5	6.863	808	812	815	rBV	976181	771103	21.56%	2.733%
6	7.022	834	839	842	rBV	2639755	2468805	69.04%	8.751%
7	7.428	902	908	911	rBV	1901464	1952988	54.62%	6.922%
8	8.145	1025	1030	1033	rBV	1108323	1051737	29.41%	3.728%
9	9.222	1207	1213	1216	rBV	3289790	3575873	100.00%	12.675%
10	9.898	1323	1328	1332	rBV	1274863	1134138	31.72%	4.020%
11	10.686	1457	1462	1475	rBV	1821874	1871046	52.32%	6.632%
12	11.386	1577	1581	1584	rBV	1281711	1106472	30.94%	3.922%
13	12.974	1845	1851	1854	rBV	2921240	3156255	88.27%	11.188%
14	14.021	2025	2029	2033	rBV	925777	837800	23.43%	2.970%
15	15.498	2274	2280	2292	rBV	501797	650518	18.19%	2.306%

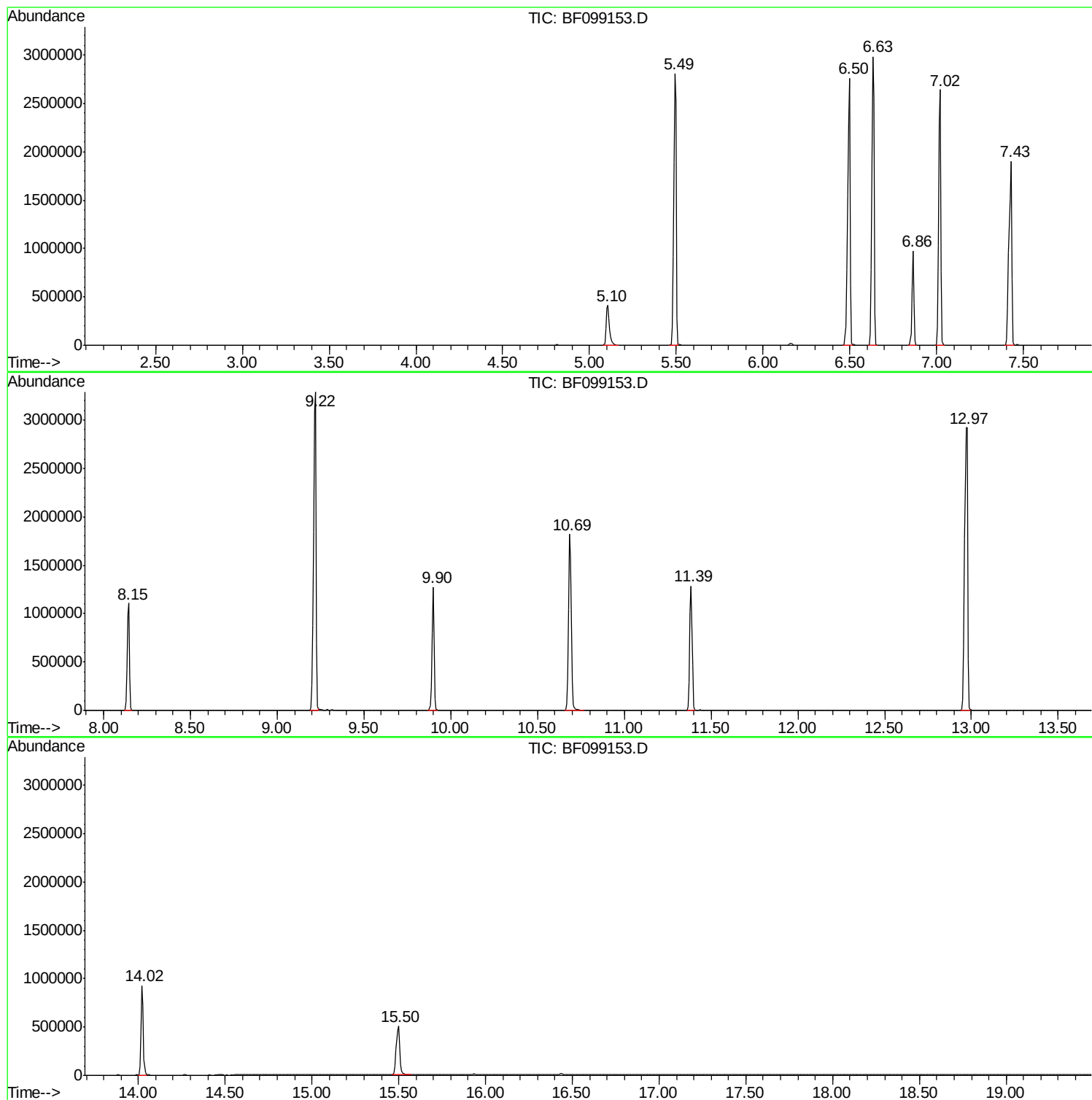
Sum of corrected areas: 28212240

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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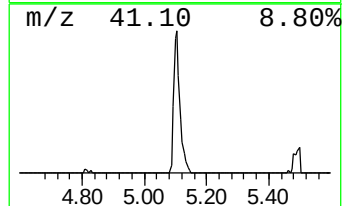
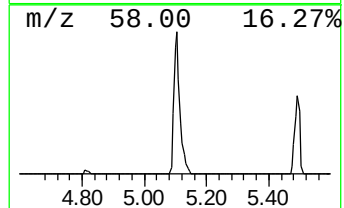
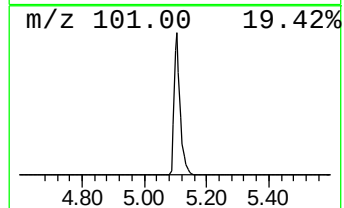
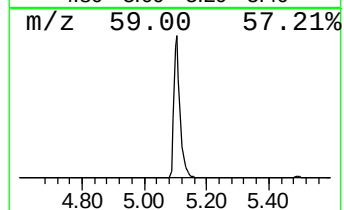
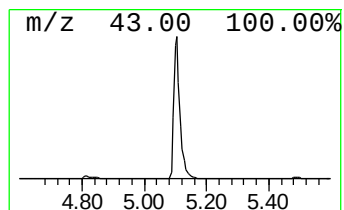
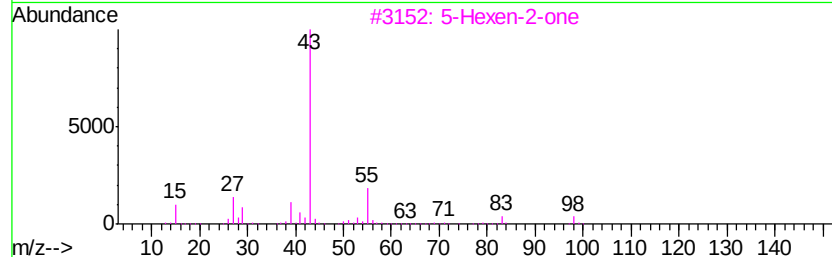
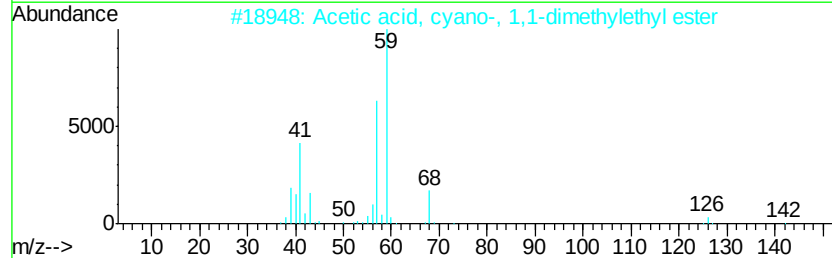
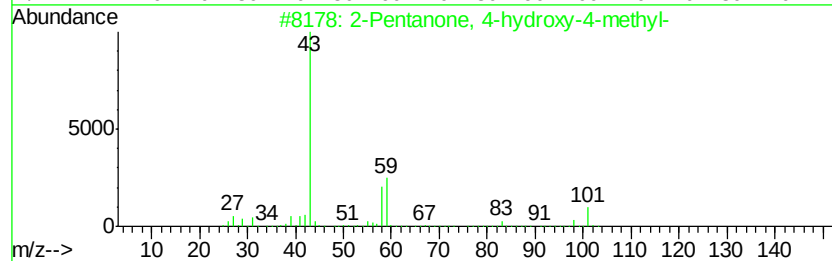
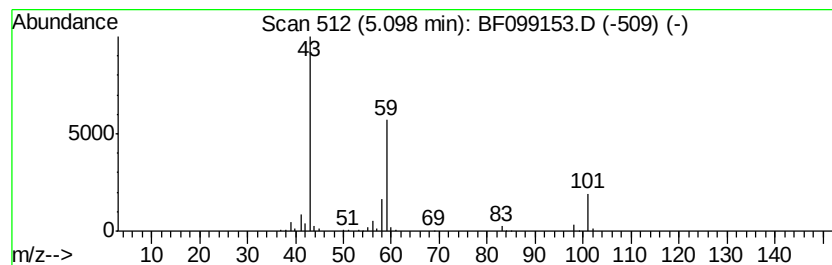
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	14.87 ng	573256	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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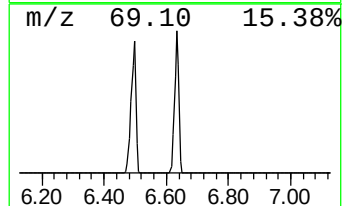
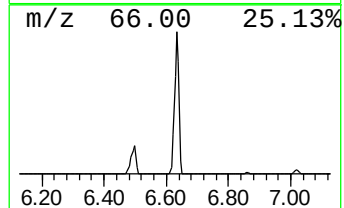
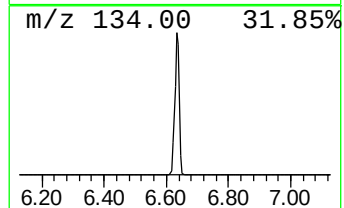
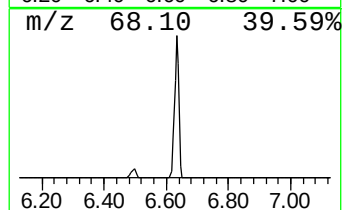
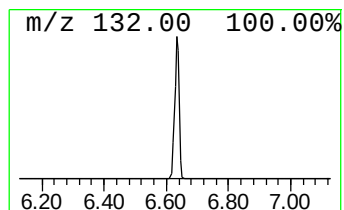
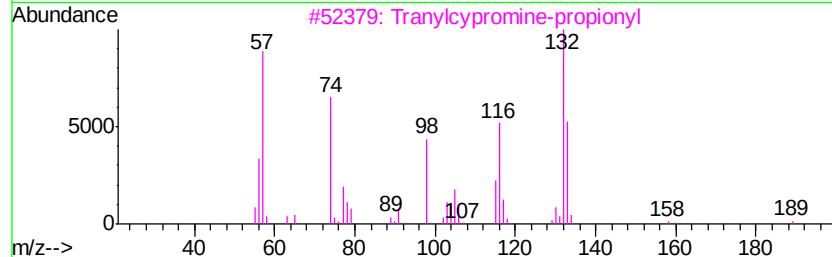
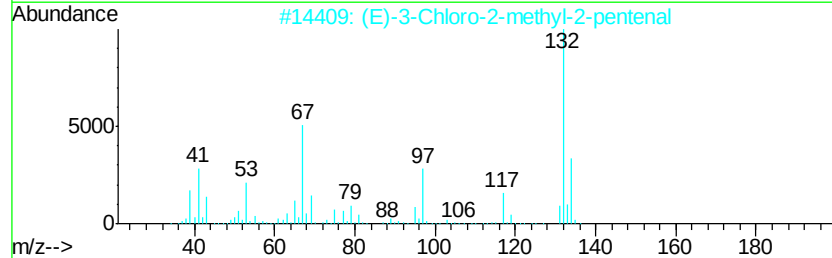
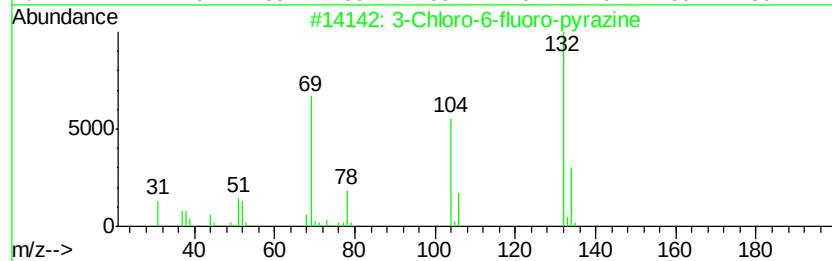
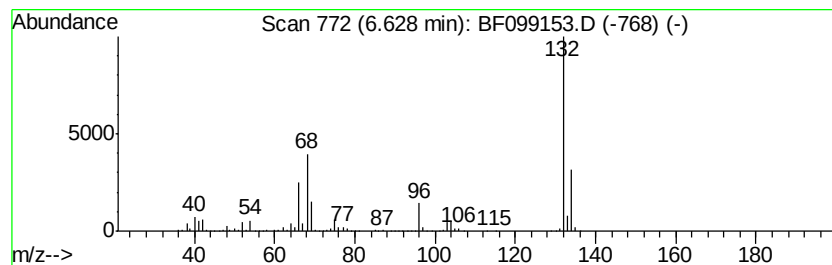
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	80.26 ng	3094610	1,4-Dichlorobenzene-d4	6.86

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	25
3		Tranylcypromine-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...	5.10	14.9	ng	573256	1	6.86	771103	20.0
unknown6.63	6.63	80.3	ng	3094610	1	6.86	771103	20.0