

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110617\
 Data File : BF100250.D
 Acq On : 6 Nov 2017 14:28
 Operator : SJ/JU
 Sample : PB103849BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103849BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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	4.992	492	494	523	rBV	115003	246952	11.32%	1.412%
2	5.392	559	562	609	rBV	1106876	1734493	79.48%	9.914%
3	6.410	732	735	753	rBV	1324460	1632434	74.80%	9.331%
4	6.534	753	756	771	rVB	1824390	1825731	83.66%	10.436%
5	6.757	792	794	810	rVB	472166	489242	22.42%	2.797%
6	6.910	817	820	839	rBV	1464855	1452659	66.56%	8.303%
7	7.333	889	892	919	rBV	861115	1040539	47.68%	5.948%
8	8.039	1009	1012	1035	rBV	596972	681604	31.23%	3.896%
9	9.110	1190	1194	1211	rBV	2202957	2182396	100.00%	12.475%
10	9.792	1307	1310	1323	rBV2	796559	779790	35.73%	4.457%
11	10.592	1442	1446	1466	rBV	1124359	1184262	54.26%	6.769%
12	11.286	1561	1564	1578	rBV	766039	819328	37.54%	4.683%
13	12.868	1829	1833	1836	rBV	2312570	2157356	98.85%	12.332%
14	13.945	2012	2016	2030	rBV	607826	722649	33.11%	4.131%
15	15.404	2260	2264	2281	rBV	324309	545217	24.98%	3.116%

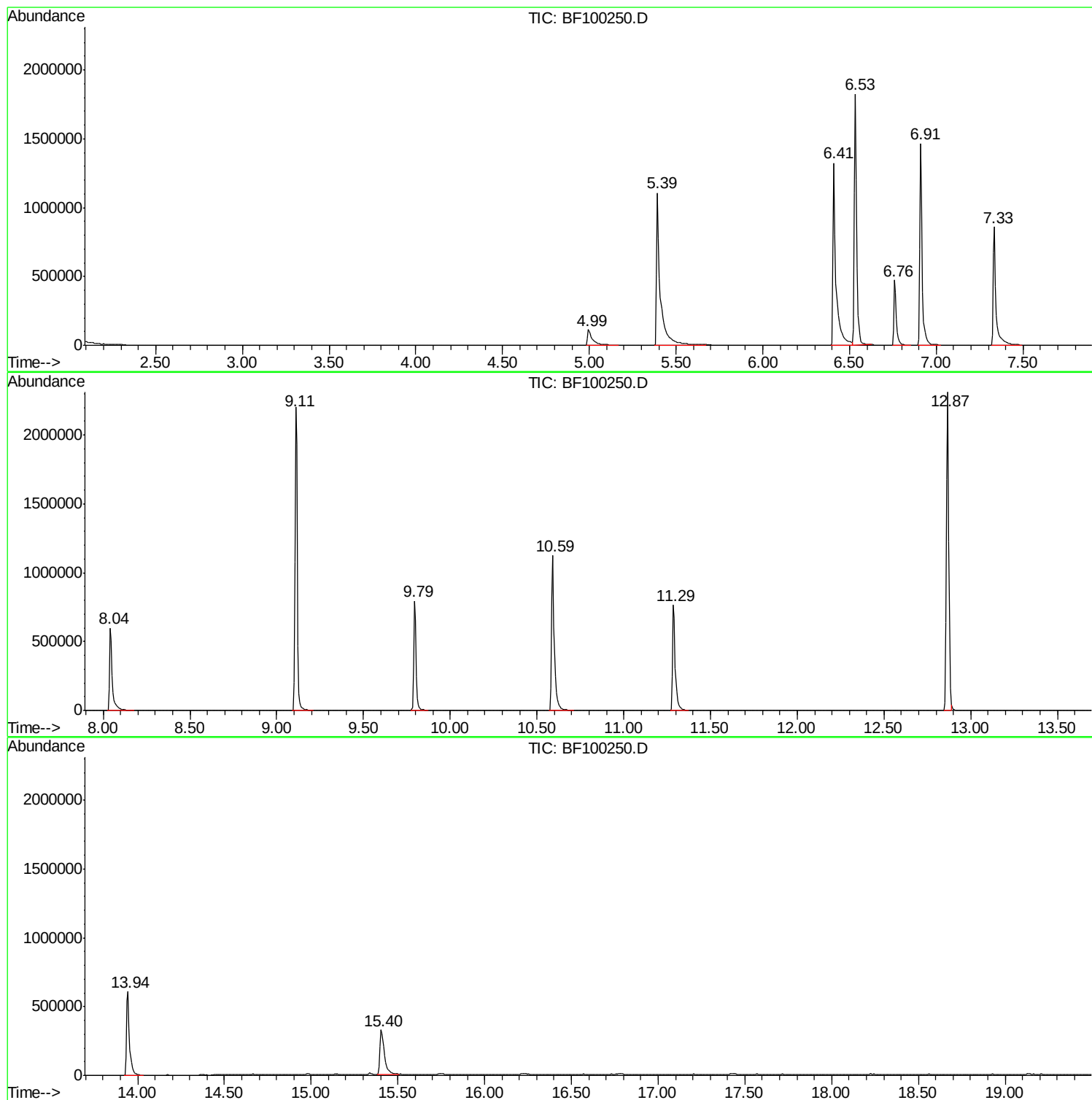
Sum of corrected areas: 17494652

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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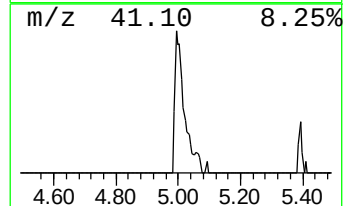
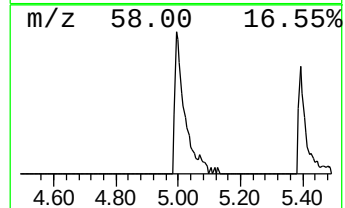
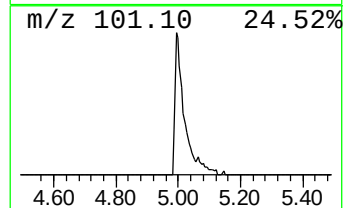
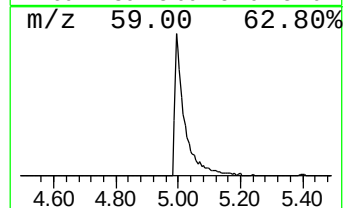
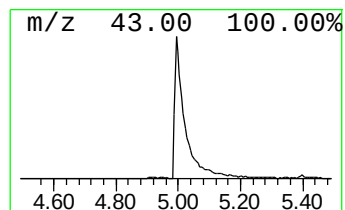
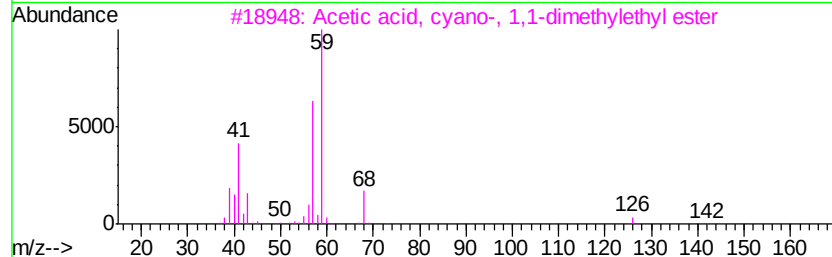
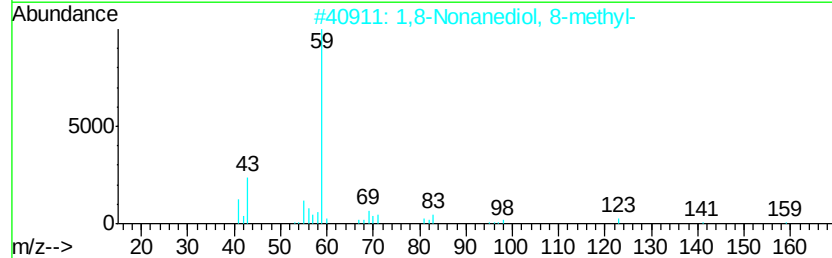
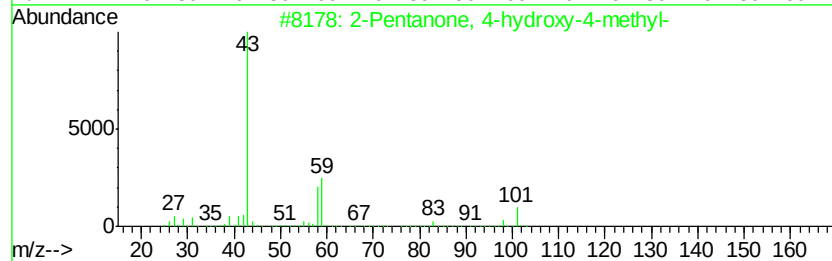
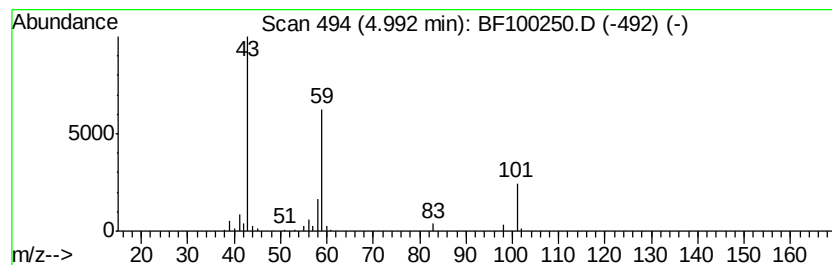
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	10.10 ng	246952	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
3			Acetic acid, cyano-, 1,1-dimethy...	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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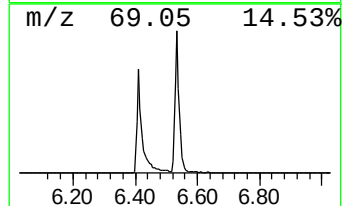
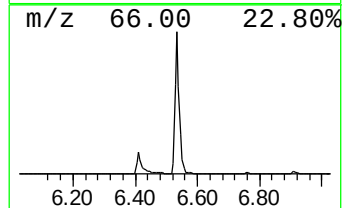
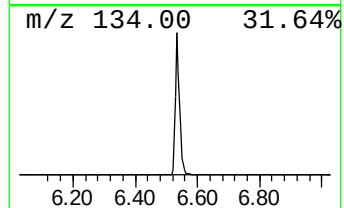
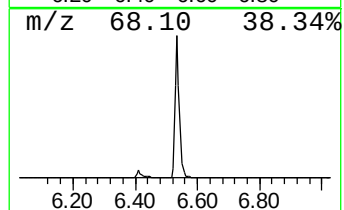
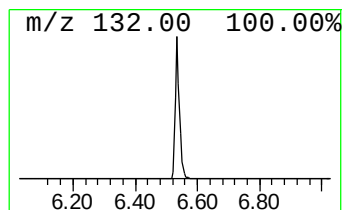
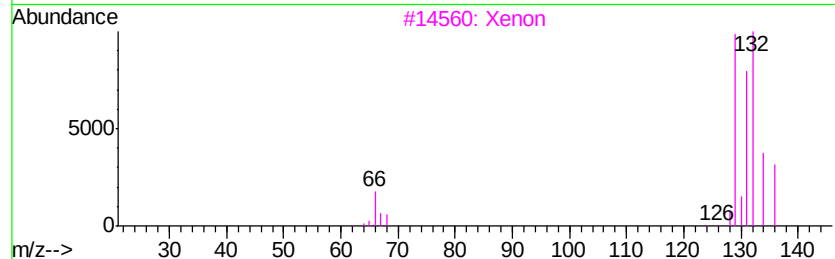
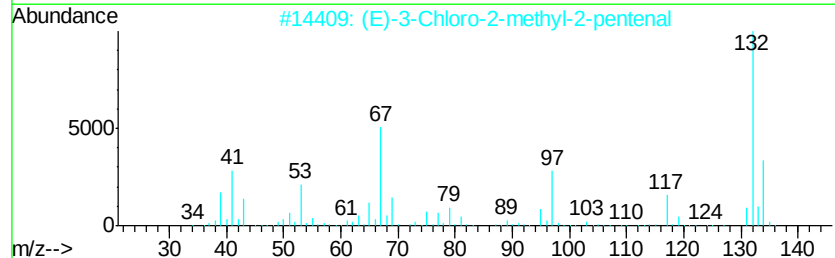
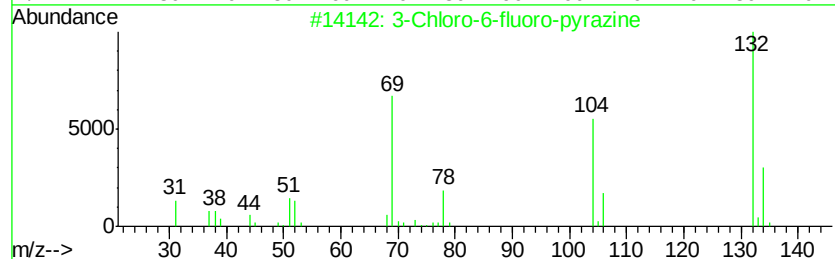
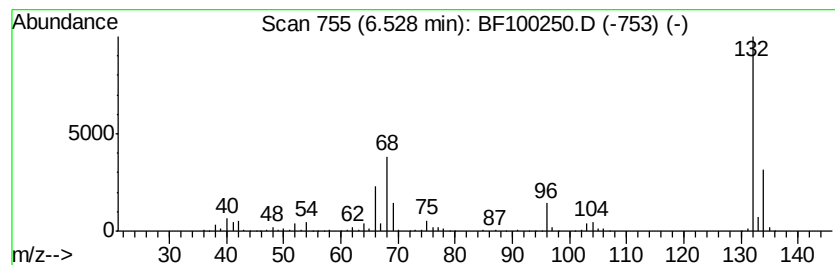
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	74.64 ng	1825730	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3			Xenon	132	Xe	007440-63-3	9
4			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5			5-Aminoindole	132	C8H8N2	005192-03-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	10.1	ng	246952	1	6.76	489242	20.0
unknown6.53	6.53	74.6	ng	1825730	1	6.76	489242	20.0