

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101216\
 Data File : BF090530.D
 Acq On : 12 Oct 2016 12:51
 Operator : UM/SJ
 Sample : PB93973BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93973BL

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-BF101116.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	5.160	244	247	263	rBV	845616	1294105	19.29%	2.482%
2	5.583	281	284	301	rBV	2722921	5124584	76.40%	9.828%
3	6.589	369	372	381	rBV	3390342	5111812	76.21%	9.804%
4	6.726	381	384	386	rBV	4054851	5330959	79.47%	10.224%
5	6.943	401	403	413	rVB	914494	1242812	18.53%	2.384%
6	7.103	414	417	431	rVB	4232935	4253254	63.41%	8.157%
7	7.515	450	453	455	rBV	2532577	3386703	50.49%	6.495%
8	8.235	513	516	524	rBV	1639220	1774184	26.45%	3.403%
9	9.309	607	610	613	rBV	5109447	6244197	93.09%	11.975%
10	9.995	667	670	672	rBV	1746840	2017814	30.08%	3.870%
11	10.783	736	739	742	rBV	3821375	3996876	59.59%	7.665%
12	11.481	797	800	802	rBV	2372005	2088843	31.14%	4.006%
13	13.069	936	939	942	rBV	6265409	6707730	100.00%	12.864%
14	14.121	1029	1031	1034	rBV	2022622	1923456	28.68%	3.689%
15	15.595	1157	1160	1163	rBV	1185204	1644436	24.52%	3.154%

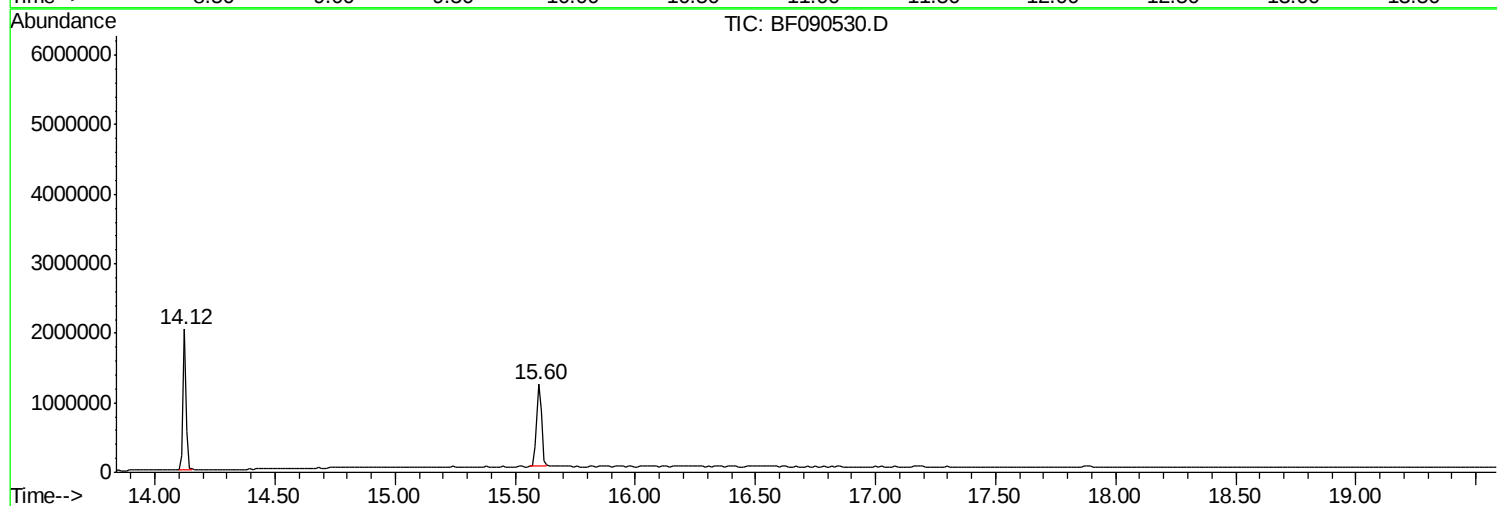
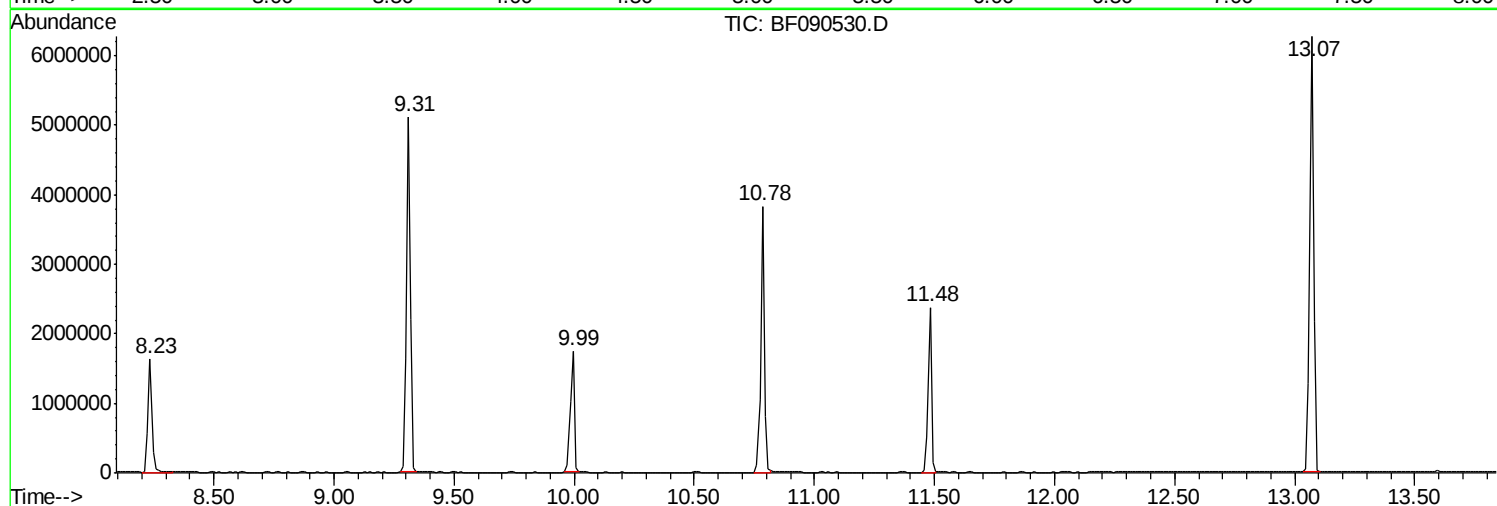
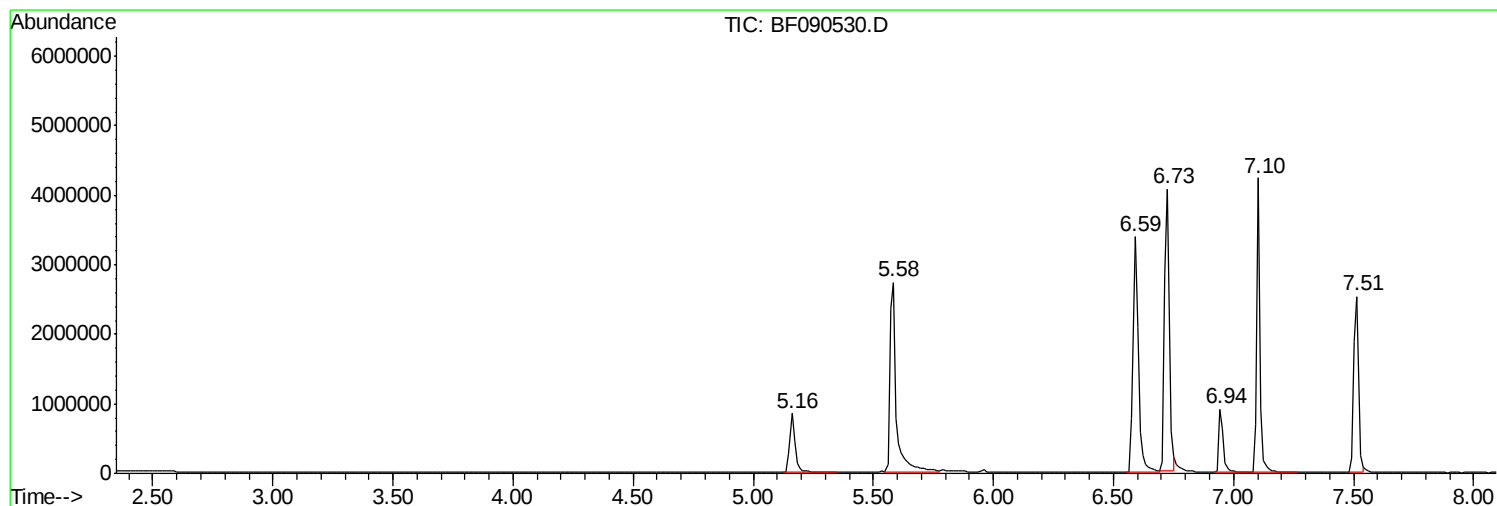
Sum of corrected areas: 52141765

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101116.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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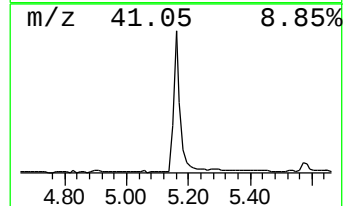
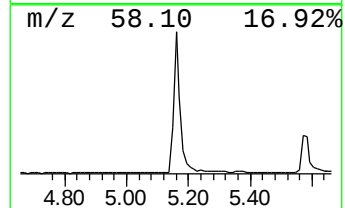
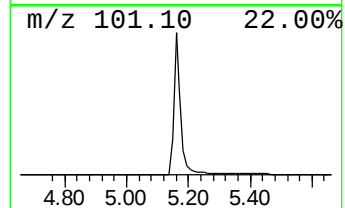
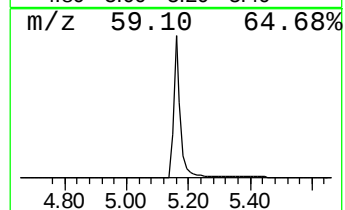
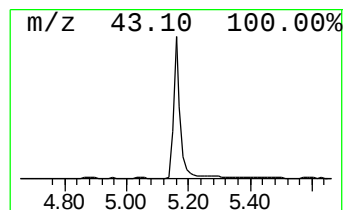
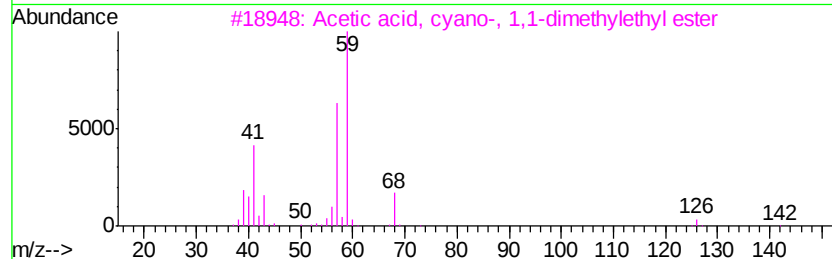
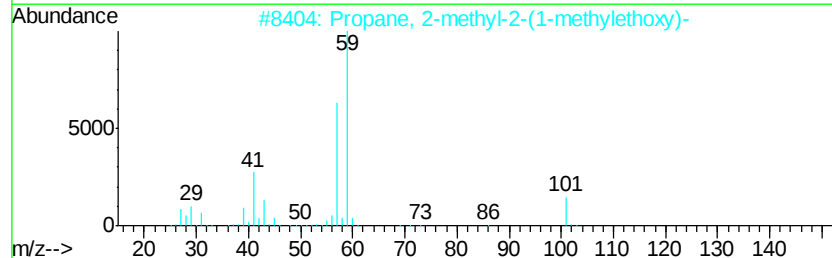
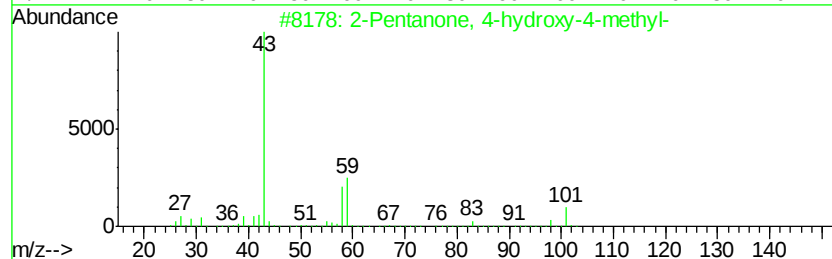
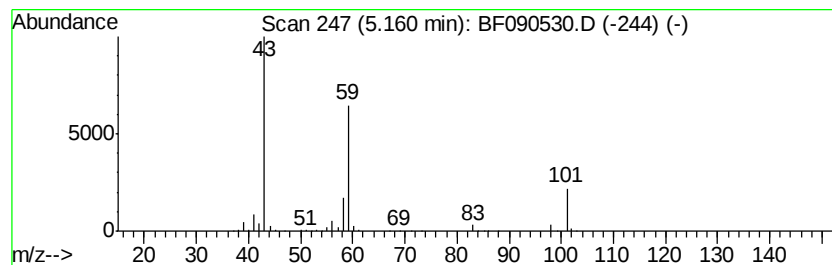
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101116.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	20.83 ng	1294110	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
4		Guanidine	59	CH5N3	000113-00-8	9
5		Acetone	58	C3H6O	000067-64-1	9



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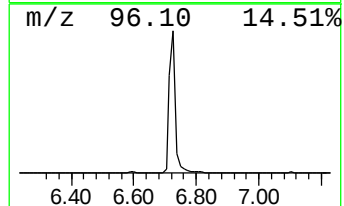
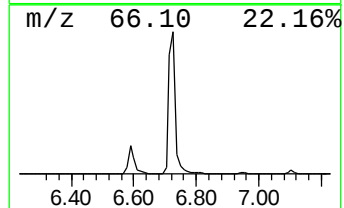
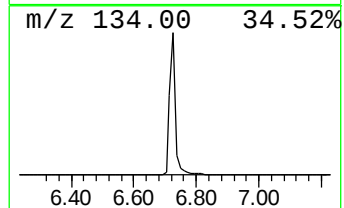
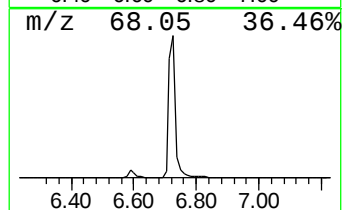
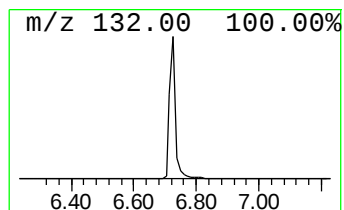
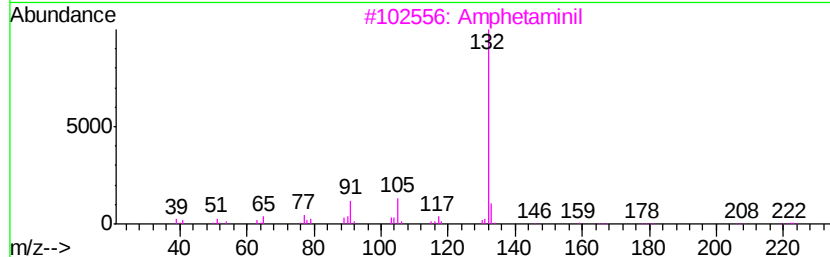
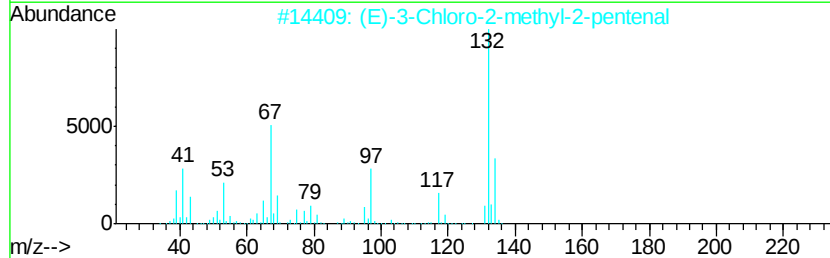
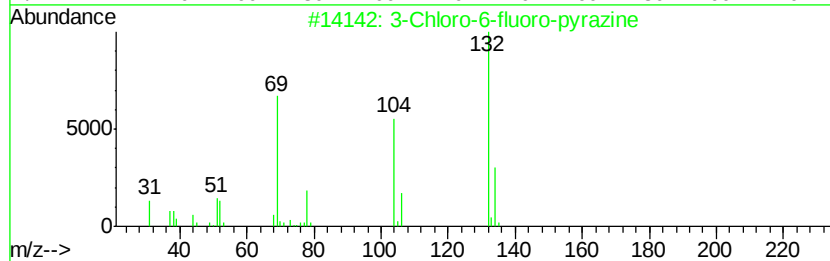
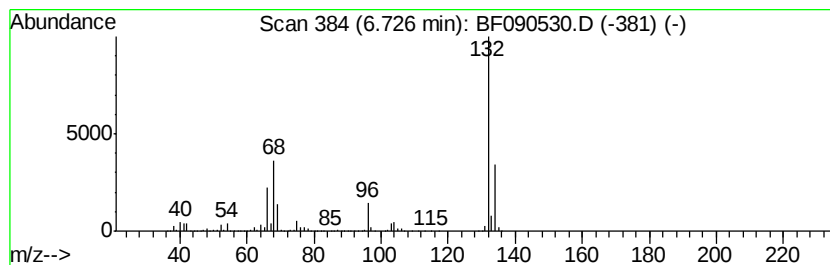
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	85.79 ng	5330960	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	20.8	ng	1294110	1	6.94	1242810	20.0
unknown6.73	6.73	85.8	ng	5330960	1	6.94	1242810	20.0