

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
 Data File : BF091393.D
 Acq On : 2 Dec 2016 17:06
 Operator : UM/SJ
 Sample : PB94987BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94987BL

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-BF112916.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	5.053	239	242	248	rBV	1329649	1944445	30.85%	3.817%
2	5.464	275	278	281	rBV	4003804	4526852	71.81%	8.886%
3	6.493	364	368	370	rBV	4099818	5348879	84.85%	10.499%
4	6.630	377	380	382	rBV	4332385	5081474	80.61%	9.974%
5	6.859	398	400	402	rBV	1550290	1333022	21.15%	2.617%
6	7.019	411	414	416	rVB	3599000	3863692	61.29%	7.584%
7	7.419	446	449	451	rBV	2863265	2903818	46.07%	5.700%
8	8.139	509	512	515	rBV	1588279	1682318	26.69%	3.302%
9	9.225	603	607	609	rBV	5247564	5762725	91.42%	11.312%
10	9.796	655	657	663	rVB	74507	122788	1.95%	0.241%
11	9.899	663	666	668	rVB	2136669	2113524	33.53%	4.149%
12	10.688	731	735	737	rBV	3223120	3894297	61.78%	7.644%
13	11.385	793	796	798	rBV	1545225	2149846	34.10%	4.220%
14	12.974	932	935	937	rBV	5667659	6303630	100.00%	12.373%
15	14.014	1023	1026	1028	rBV	2322377	2173839	34.49%	4.267%
16	15.442	1148	1151	1154	rBV	1254407	1739786	27.60%	3.415%

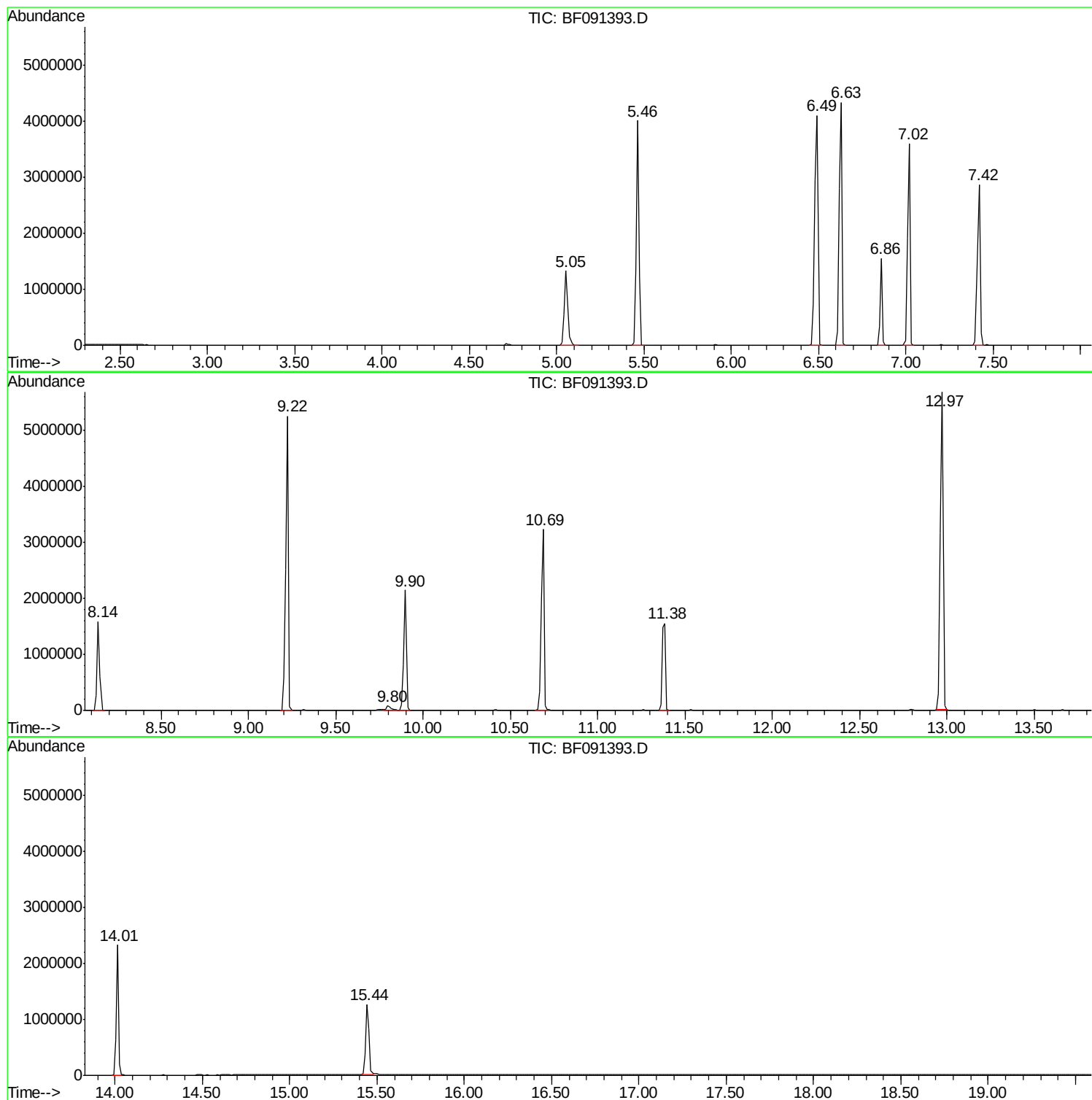
Sum of corrected areas: 50944935

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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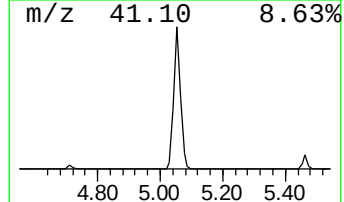
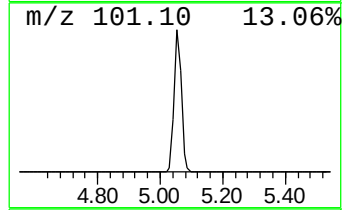
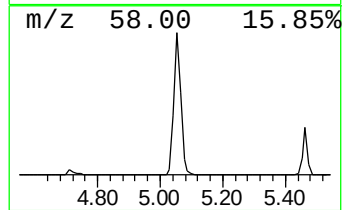
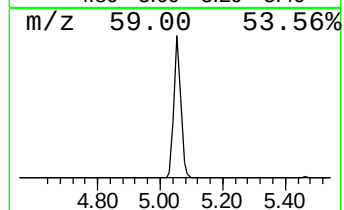
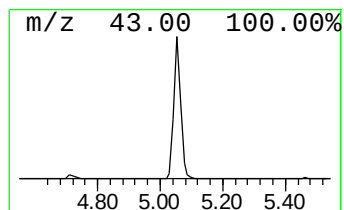
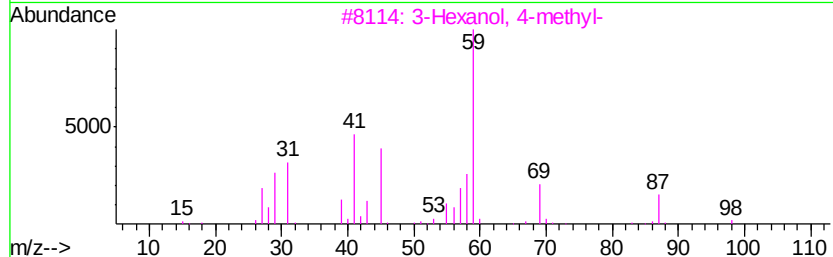
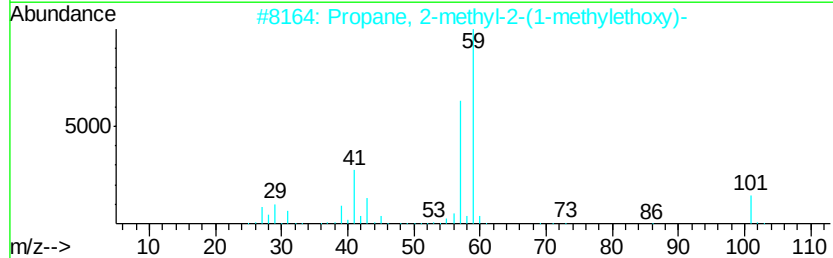
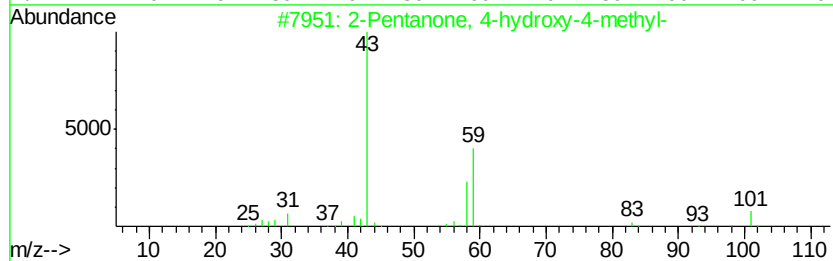
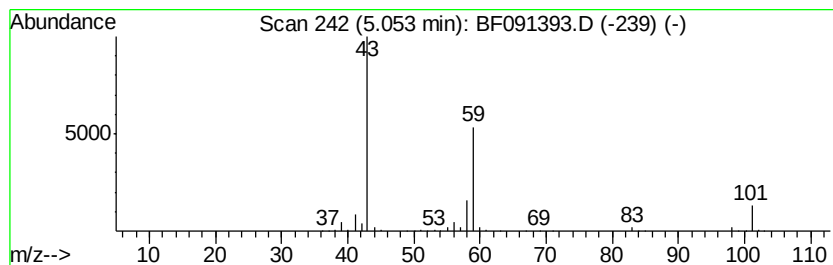
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	29.17 ng	1944450	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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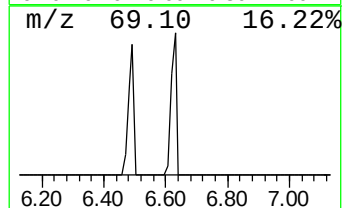
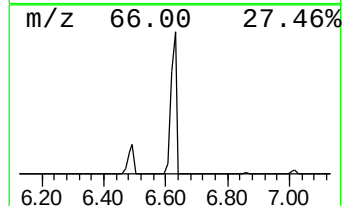
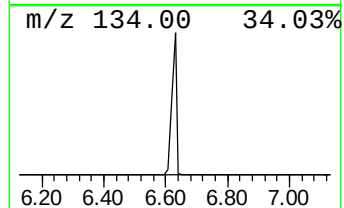
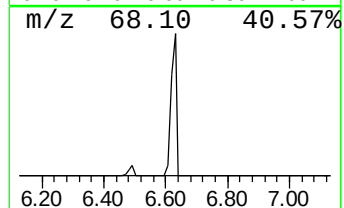
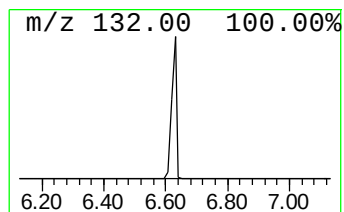
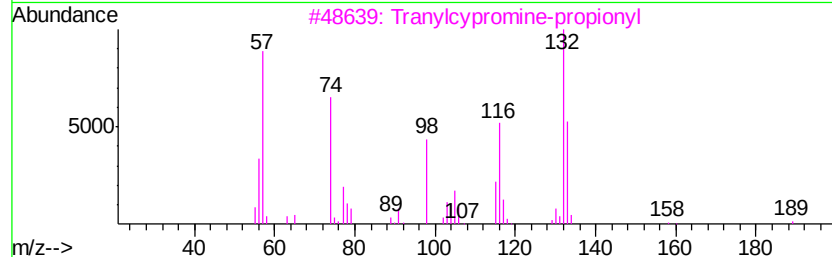
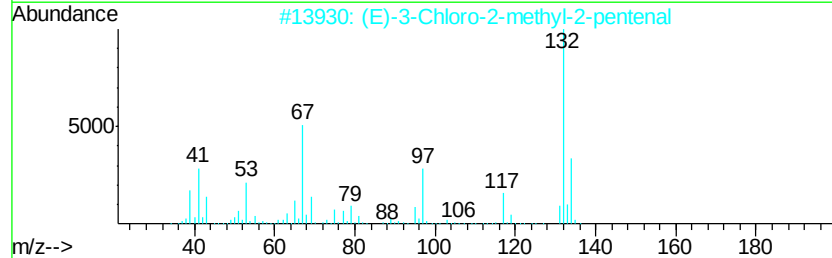
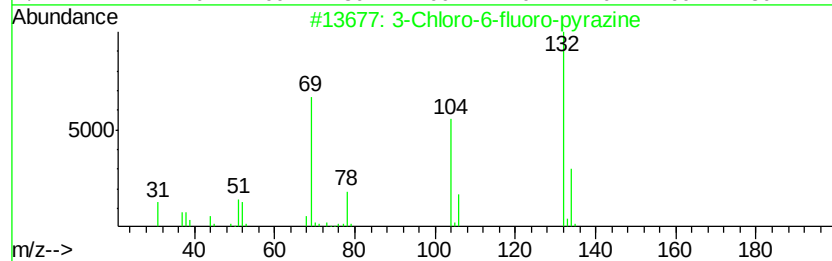
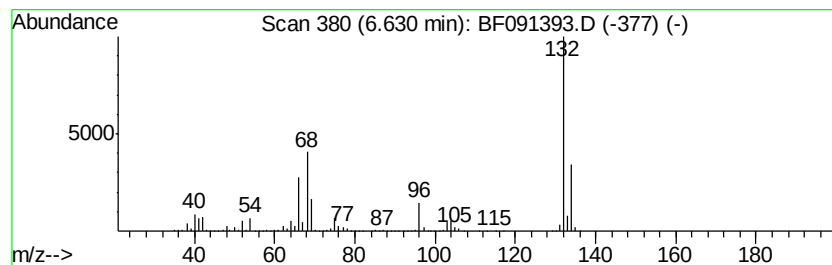
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	76.24 ng	5081470	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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...	5.05	29.2	ng	1944450	1	6.86	1333020	20.0
unknown6.63	6.63	76.2	ng	5081470	1	6.86	1333020	20.0