

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF113016\
 Data File : BF091345.D
 Acq On : 30 Nov 2016 10:59
 Operator : UM/SJ
 Sample : PB94951BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94951BL

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-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	244	rBV	1168706	1335508	27.27%	4.042%
2	5.464	275	278	280	rBV	2668969	2394744	48.89%	7.249%
3	6.482	364	367	370	rBV	2296510	2153475	43.97%	6.518%
4	6.630	377	380	382	rBV	2401022	2493297	50.90%	7.547%
5	6.859	398	400	403	rBV	1079643	1027557	20.98%	3.110%
6	7.019	411	414	416	rBV	3584500	3125176	63.80%	9.460%
7	7.419	446	449	452	rBV	1997261	2574997	52.57%	7.794%
8	8.150	510	513	515	rBV	1328168	1415932	28.91%	4.286%
9	9.225	604	607	610	rBV	4570429	4448767	90.83%	13.466%
10	9.899	663	666	669	rBV	1758743	1543399	31.51%	4.672%
11	10.688	732	735	737	rBV	1534943	1710655	34.92%	5.178%
12	11.385	793	796	798	rBV	1576610	1610352	32.88%	4.874%
13	12.802	918	920	922	rBV	58168	65364	1.33%	0.198%
14	12.974	932	935	937	rBV	4105134	4898143	100.00%	14.826%
15	14.014	1023	1026	1028	rBV	1294063	1322382	27.00%	4.003%
16	15.443	1148	1151	1154	rBV	746298	917518	18.73%	2.777%

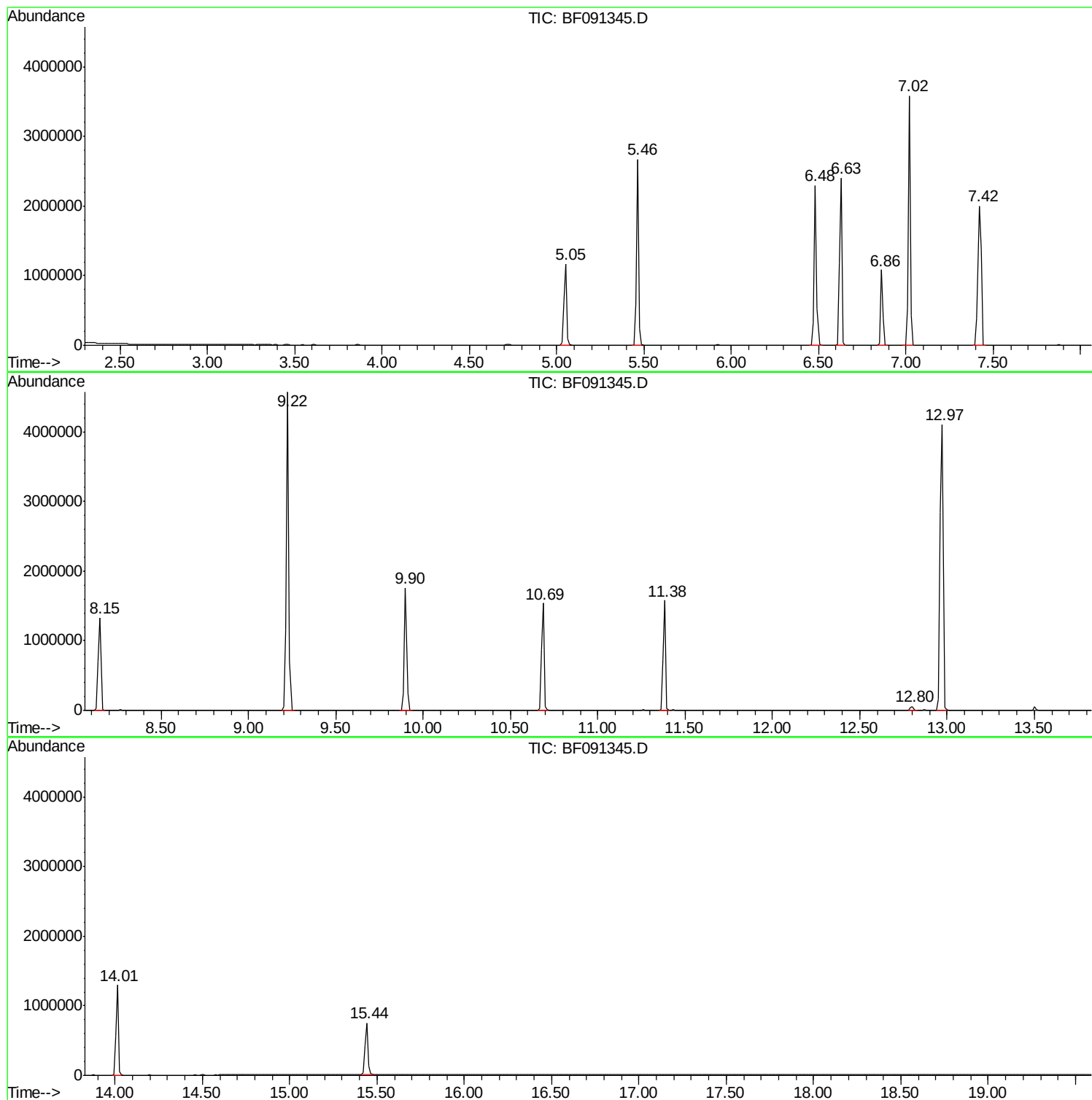
Sum of corrected areas: 33037266

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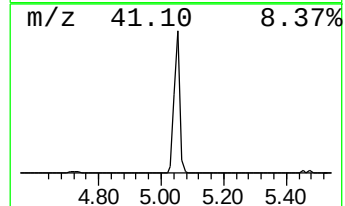
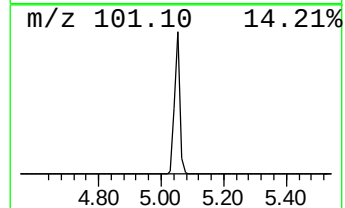
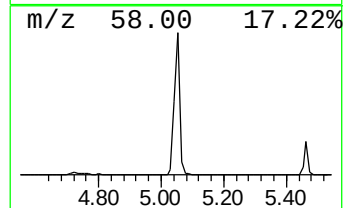
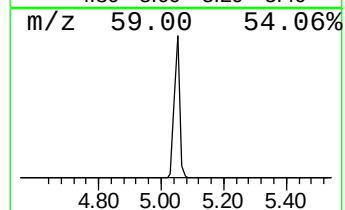
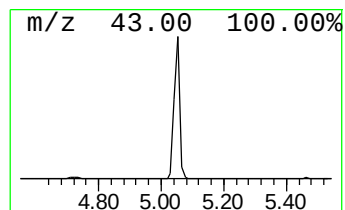
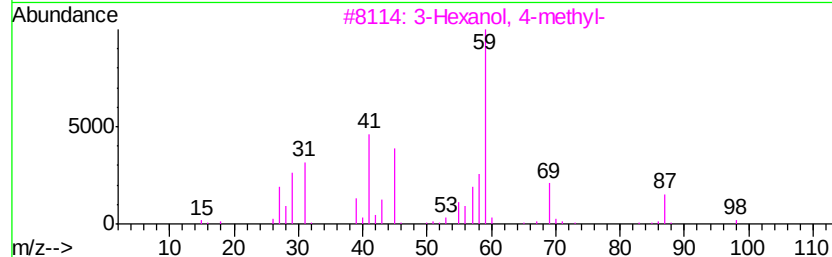
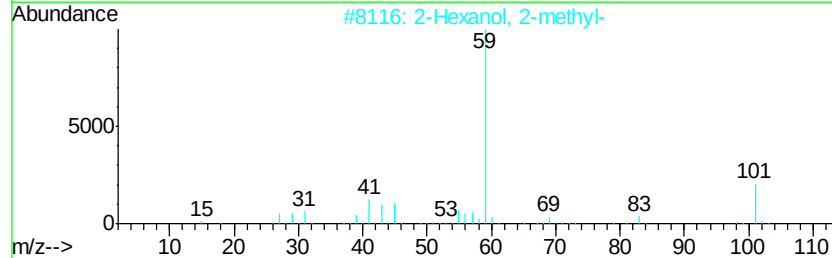
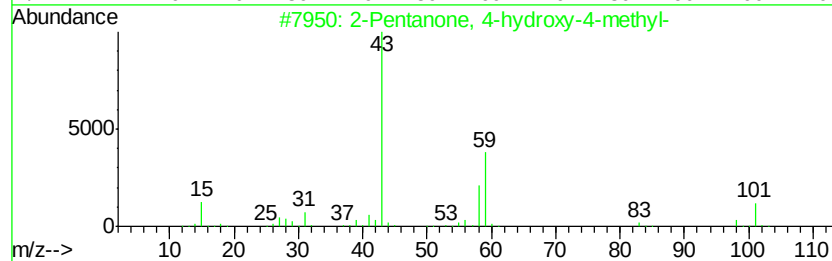
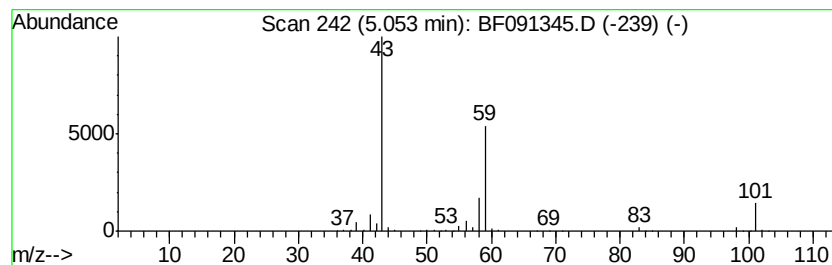
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	25.99 ng	1335510	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	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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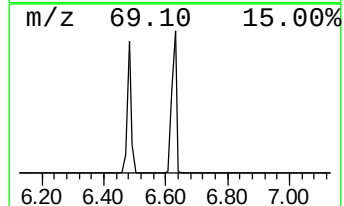
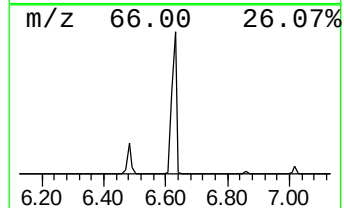
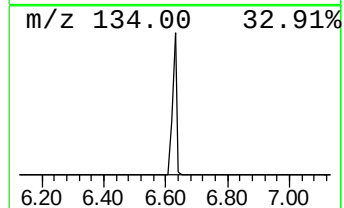
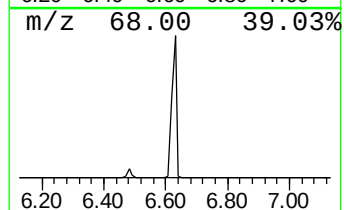
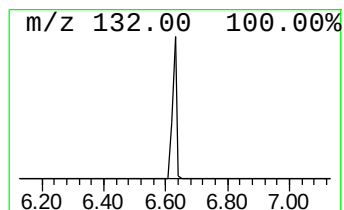
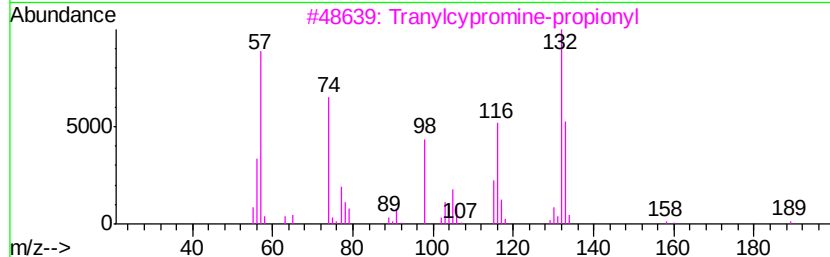
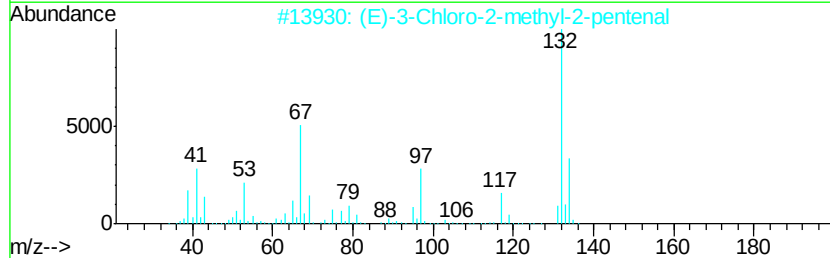
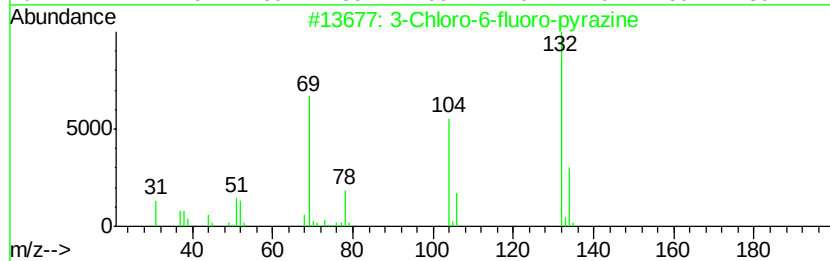
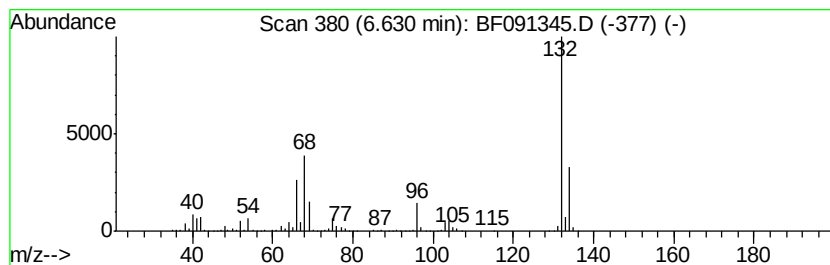
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	48.53 ng	2493300	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	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.05	26.0	ng	1335510	1	6.86	1027560	20.0
unknown6.63	6.63	48.5	ng	2493300	1	6.86	1027560	20.0