

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082415\
 Data File : BF081175.D
 Acq On : 24 Aug 2015 18:27
 Operator : UM/IZ
 Sample : PB84980BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84980BL

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-BF081715.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	3.910	176	177	182	rBV	11467	21331	1.18%	0.133%
2	4.664	239	243	257	rBV	765244	1196208	65.93%	7.471%
3	5.110	280	282	285	rBV	1148941	1694265	93.38%	10.581%
4	5.533	317	319	323	rBB	25454	25789	1.42%	0.161%
5	6.184	374	376	379	rBV	1234334	1716295	94.59%	10.719%
6	6.310	384	387	389	rBV	1497419	1738476	95.81%	10.858%
7	6.539	405	407	409	rBV	507005	412009	22.71%	2.573%
8	6.699	418	421	423	rBV	1314622	1309384	72.16%	8.178%
9	7.110	454	457	459	rBV	1156263	1110780	61.22%	6.937%
10	7.830	517	520	522	rBV	360185	496299	27.35%	3.100%
11	8.905	611	614	616	rBV	1971043	1814449	100.00%	11.332%
12	9.568	669	672	674	rBV	610002	540953	29.81%	3.378%
13	10.345	738	740	743	rBV2	825959	948580	52.28%	5.924%
14	11.031	798	800	803	rBV2	442811	530523	29.24%	3.313%
15	12.619	936	939	942	rBV	1660273	1588383	87.54%	9.920%
16	13.648	1027	1029	1032	rBV	552856	503455	27.75%	3.144%
17	14.985	1143	1146	1148	rBV	245671	364520	20.09%	2.277%

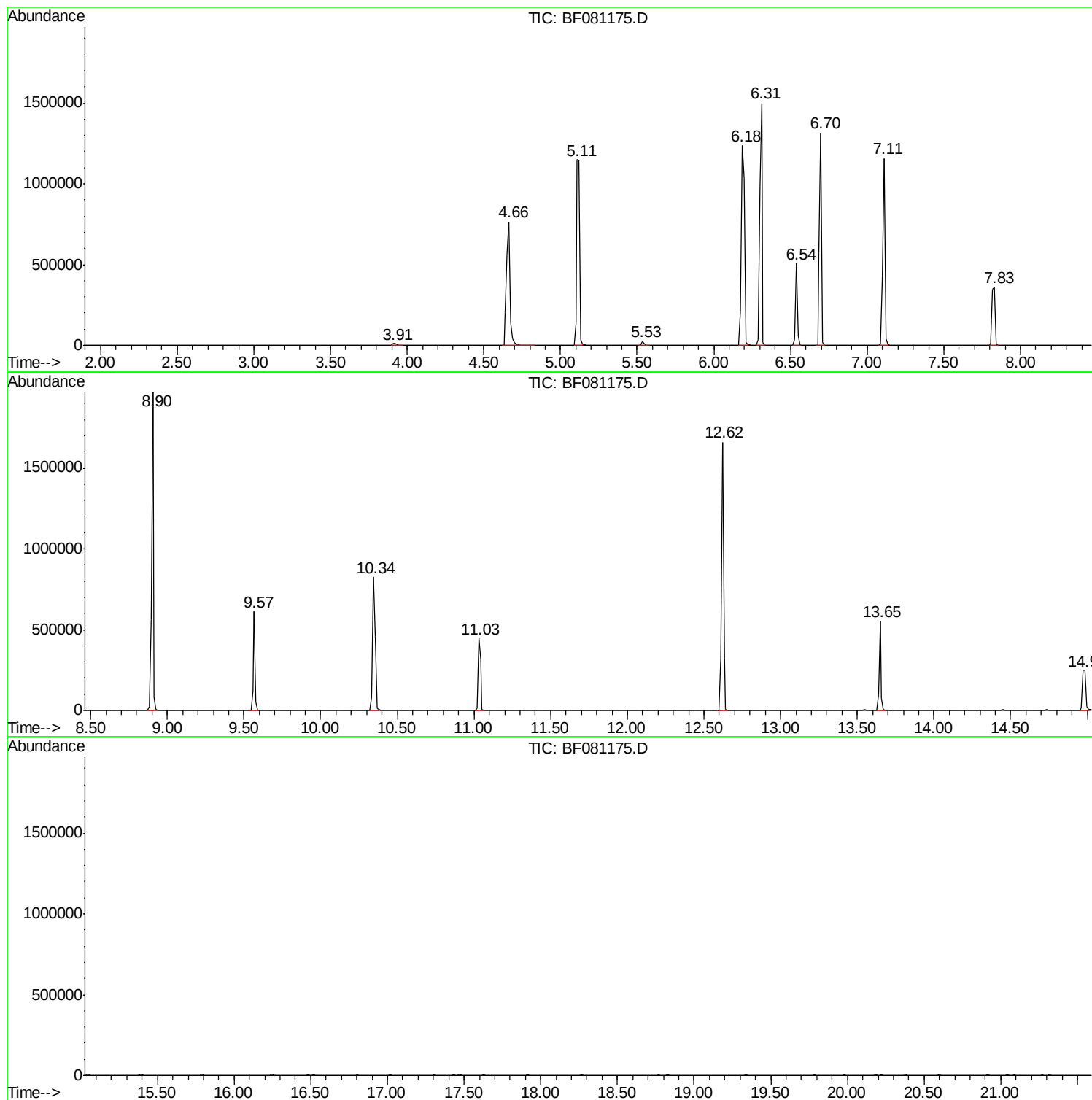
Sum of corrected areas: 16011699

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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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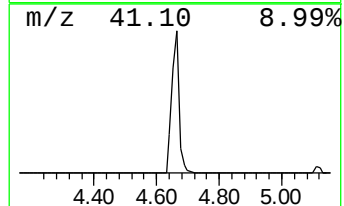
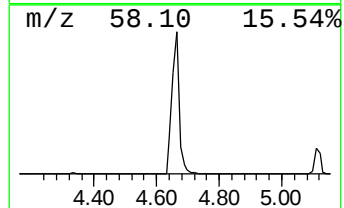
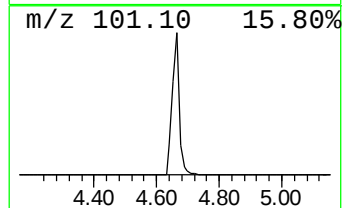
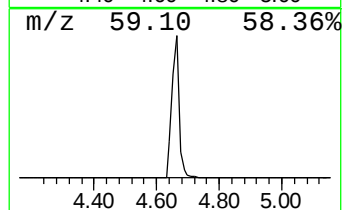
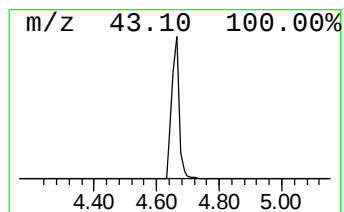
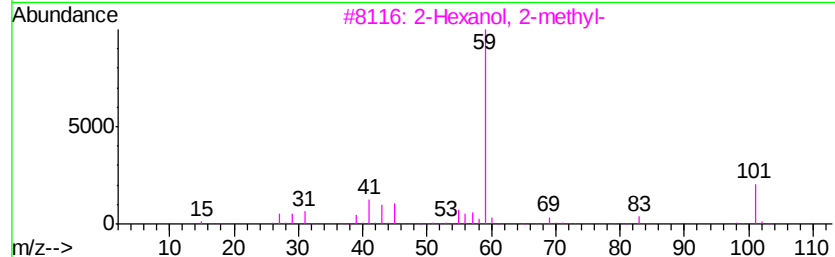
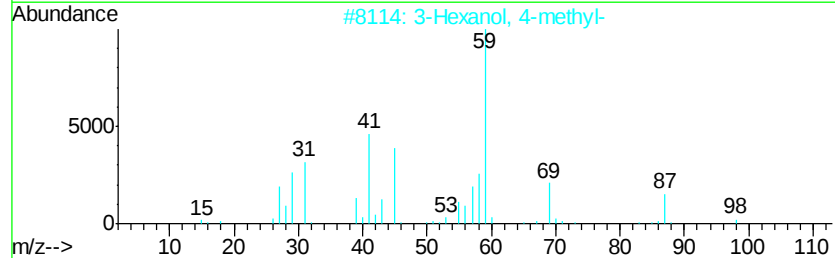
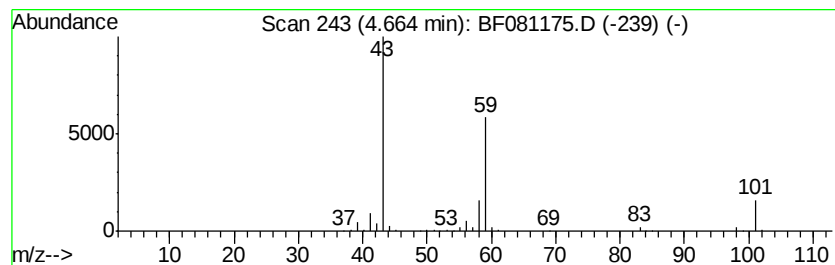
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	58.07 ng	1196210	1,4-Dichlorobenzene-d4	6.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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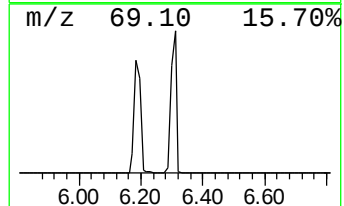
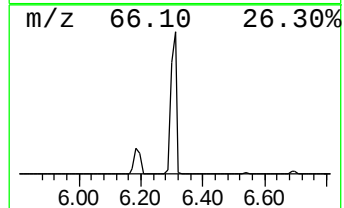
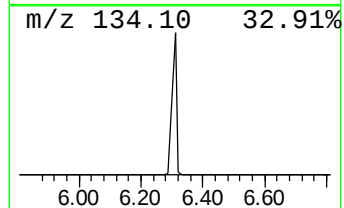
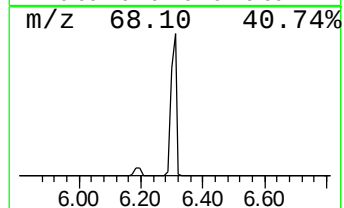
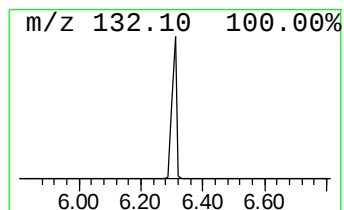
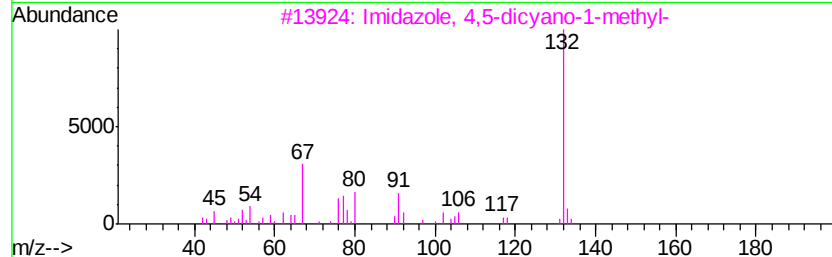
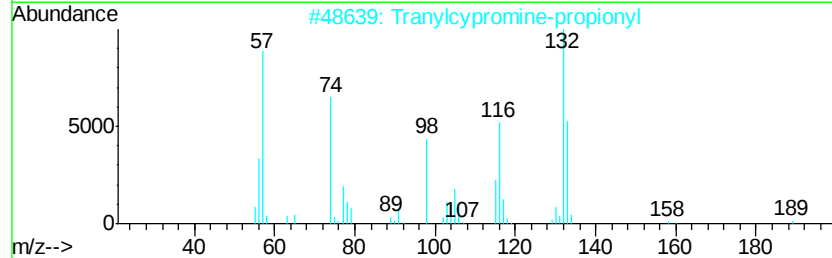
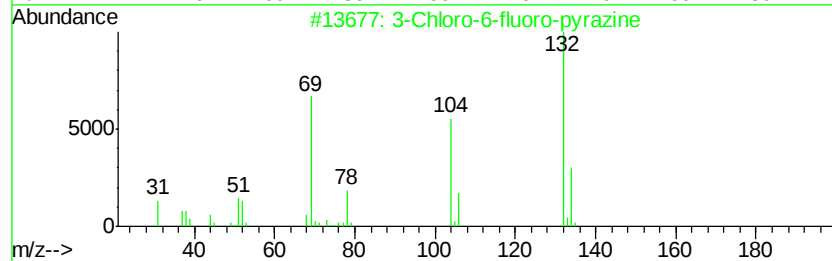
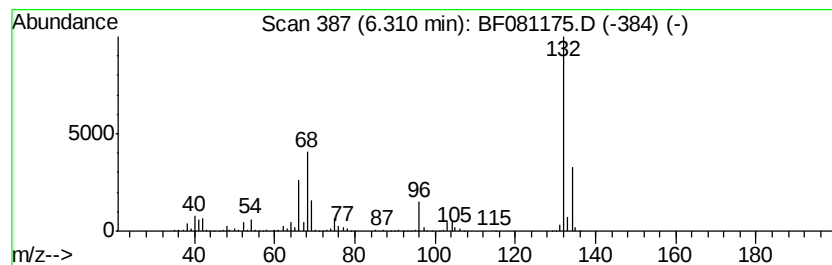
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Peak Number 2 unknown6.31 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	84.39 ng	1738480	1,4-Dichlorobenzene-d4	6.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
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...	4.66	58.1	ng	1196210	1	6.54	412009	20.0
unknown6.31	6.31	84.4	ng	1738480	1	6.54	412009	20.0