

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082415\
 Data File : BF081174.D
 Acq On : 24 Aug 2015 17:55
 Operator : UM/IZ
 Sample : PB85151BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85151BL

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	4.664	239	243	255	rBV	697127	1219410	61.04%	6.290%
2	5.110	280	282	285	rBV	1266702	1861240	93.17%	9.601%
3	6.196	374	377	379	rBV	1549189	1997730	100.00%	10.305%
4	6.310	384	387	389	rBV	1757258	1874090	93.81%	9.667%
5	6.539	405	407	410	rBV	741963	623230	31.20%	3.215%
6	6.699	418	421	423	rBV	1485893	1397470	69.95%	7.208%
7	7.110	454	457	459	rBV	1226204	1152880	57.71%	5.947%
8	7.830	517	520	522	rBV	693354	827495	41.42%	4.268%
9	8.905	611	614	616	rBV	2154013	1943637	97.29%	10.026%
10	9.568	669	672	675	rBV	1018621	917033	45.90%	4.730%
11	10.356	738	741	743	rBV2	814365	1150666	57.60%	5.935%
12	11.042	798	801	803	rBV	857469	986306	49.37%	5.088%
13	12.619	936	939	942	rBV	1595446	1899479	95.08%	9.798%
14	13.648	1027	1029	1032	rBV	836582	888332	44.47%	4.582%
15	14.985	1143	1146	1148	rBV	577305	647826	32.43%	3.342%

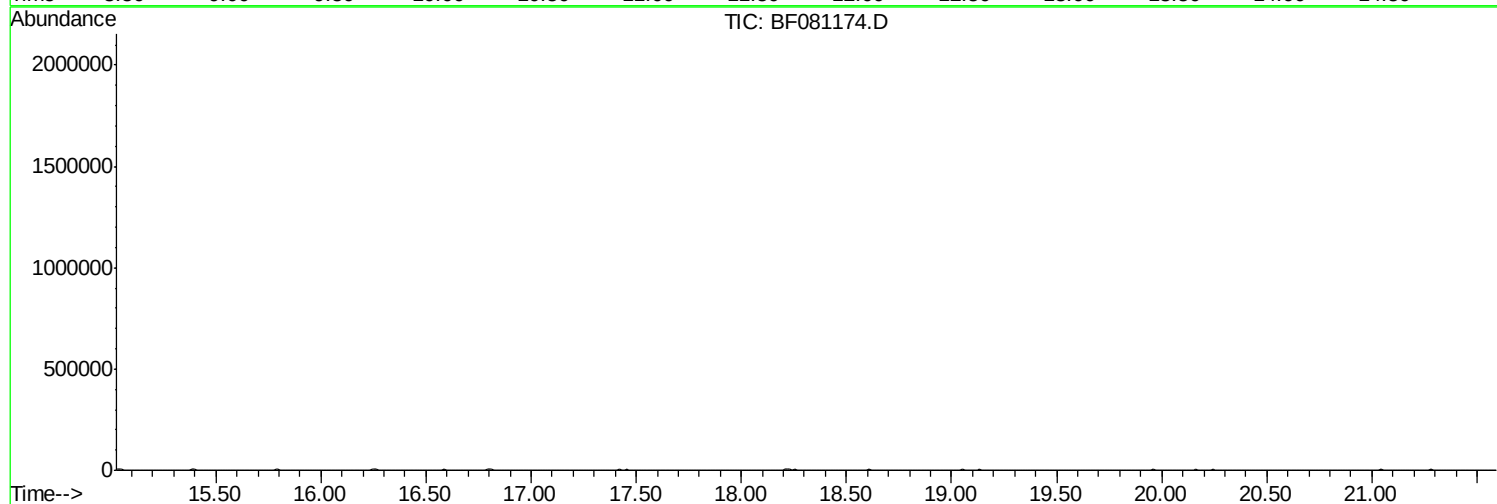
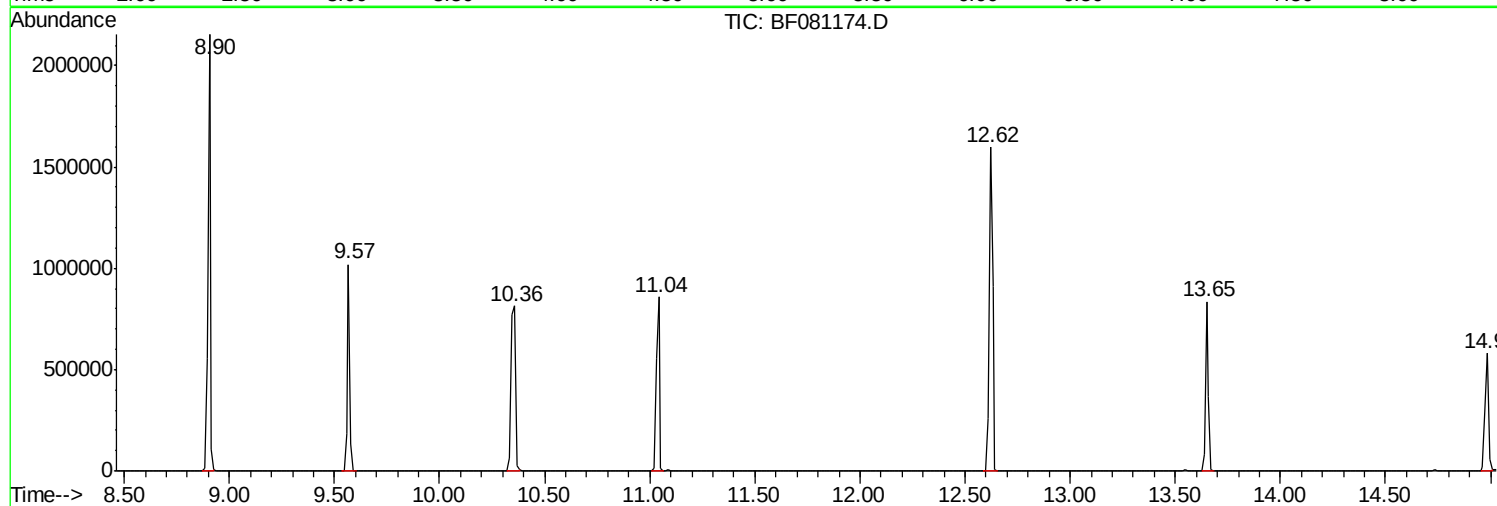
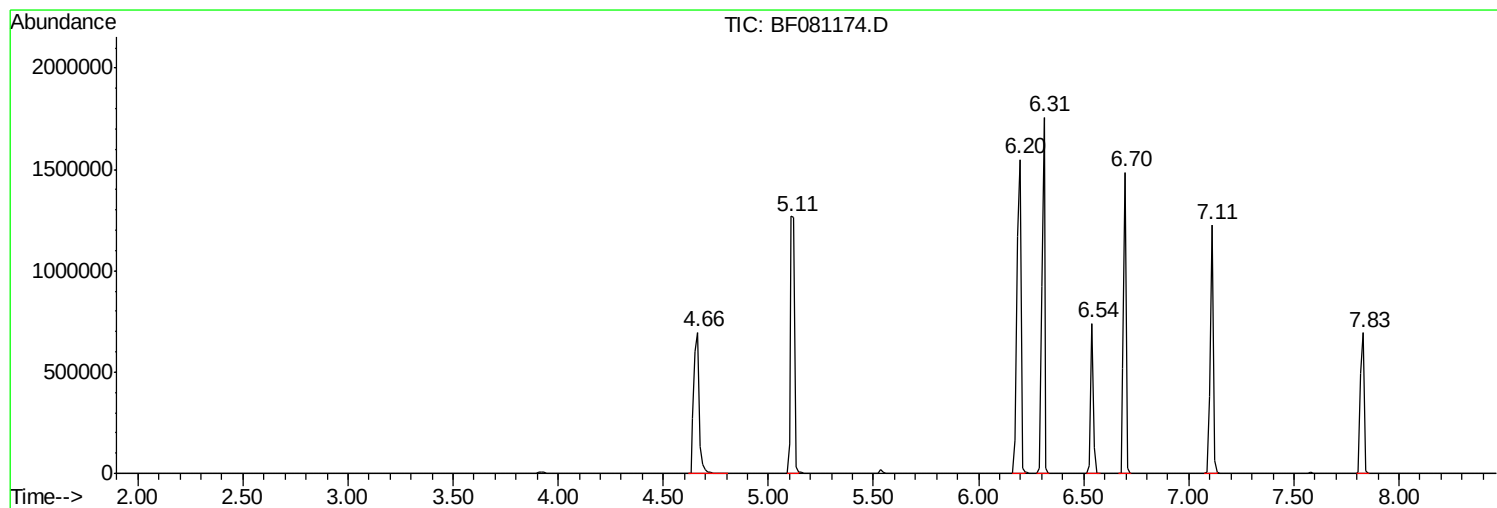
Sum of corrected areas: 19386824

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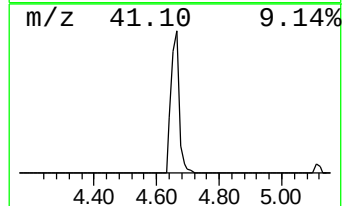
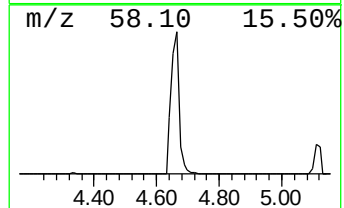
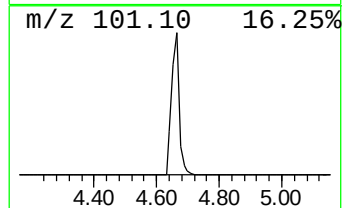
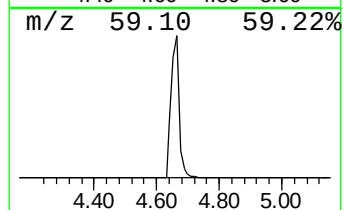
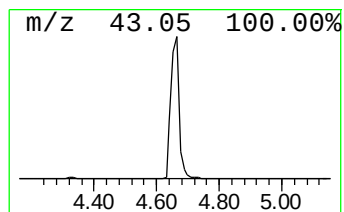
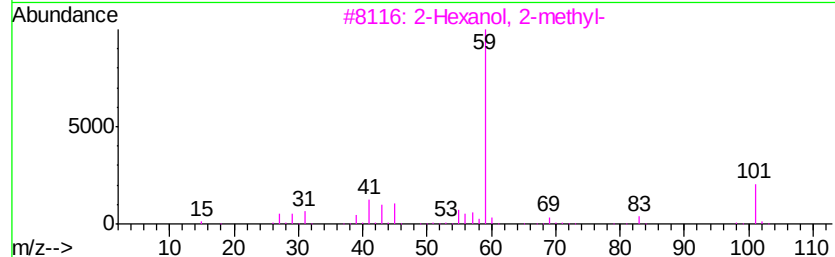
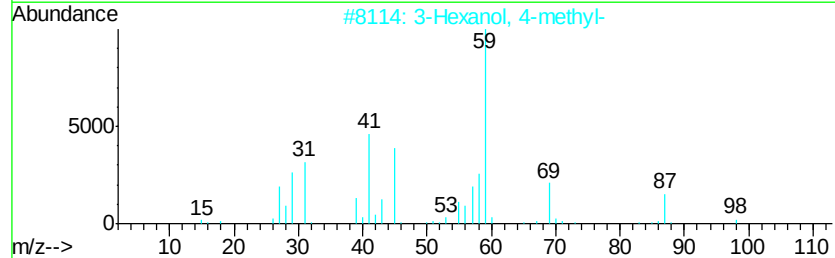
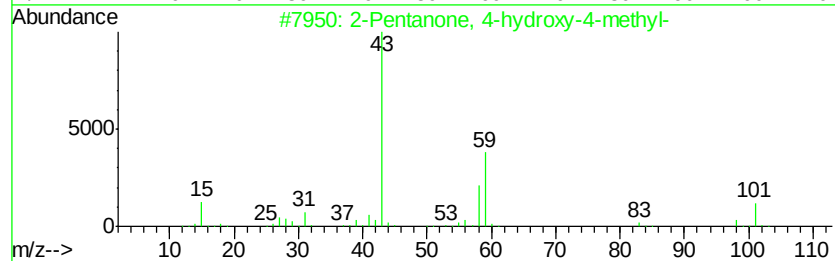
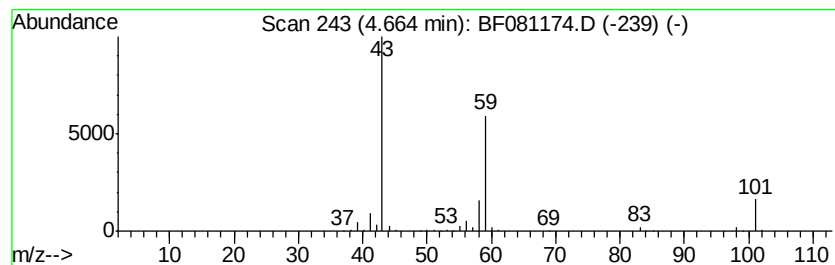
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	39.13 ng	1219410	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	45
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-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17



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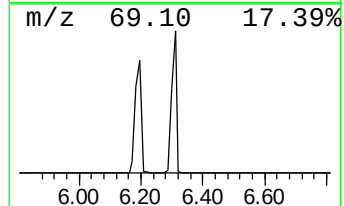
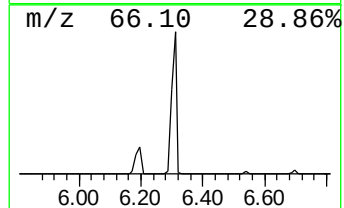
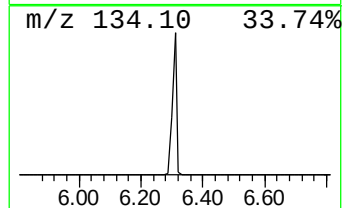
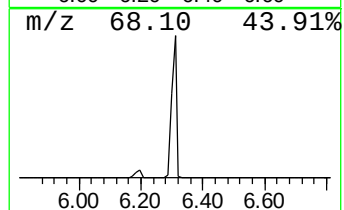
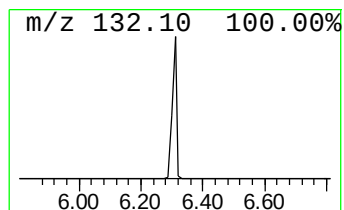
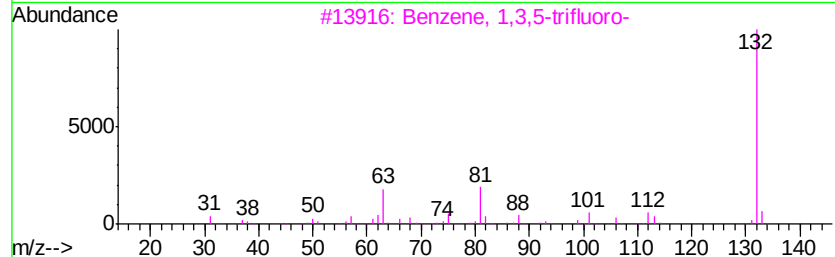
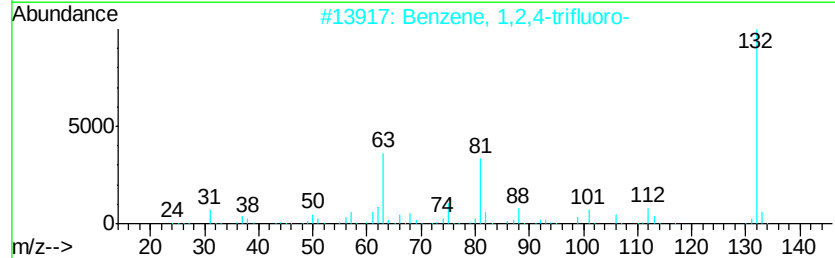
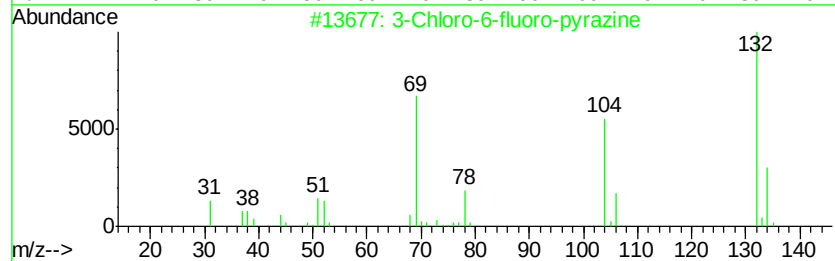
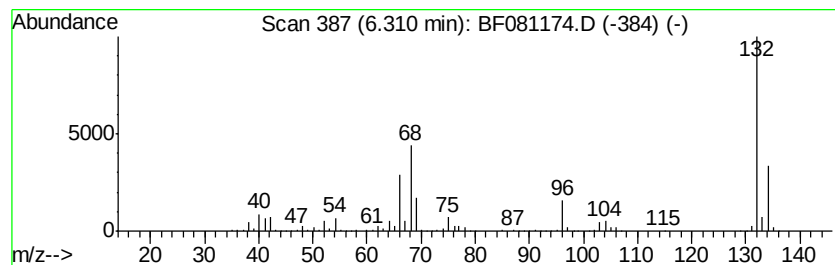
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Peak Number 2 unknown6.31 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	60.14 ng	1874090	1,4-Dichlorobenzene-d4	6.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
3			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	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...	4.66	39.1	ng	1219410	1	6.54	623230	20.0
unknown6.31	6.31	60.1	ng	1874090	1	6.54	623230	20.0