

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100715\
 Data File : BF082114.D
 Acq On : 7 Oct 2015 20:04
 Operator : UM/IZ
 Sample : G3897-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RB

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-BF092815.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	5.040	251	254	264	rVB	285322	331467	10.43%	1.749%
2	5.463	288	291	293	rBV	785903	881105	27.74%	4.650%
3	6.492	379	381	391	rBV	525461	566496	17.83%	2.990%
4	6.640	391	394	396	rBV	1699511	1927195	60.67%	10.171%
5	6.869	412	414	423	rBV	431296	499457	15.72%	2.636%
6	7.029	426	428	430	rBV	2015992	1869622	58.86%	9.867%
7	7.440	461	464	476	rBV	1272811	1436986	45.24%	7.584%
8	8.160	525	527	536	rBV	699757	682151	21.47%	3.600%
9	9.235	619	621	624	rBV	2416655	2847458	89.64%	15.028%
10	9.921	678	681	683	rBV	611907	765875	24.11%	4.042%
11	10.709	747	750	752	rBV	1759100	1819969	57.29%	9.605%
12	11.406	809	811	813	rBV	674490	744218	23.43%	3.928%
13	12.995	947	950	952	rBV	3064497	3176513	100.00%	16.765%
14	14.047	1039	1042	1054	rVB	627939	770857	24.27%	4.068%
15	15.487	1165	1168	1183	rBV	421921	628449	19.78%	3.317%

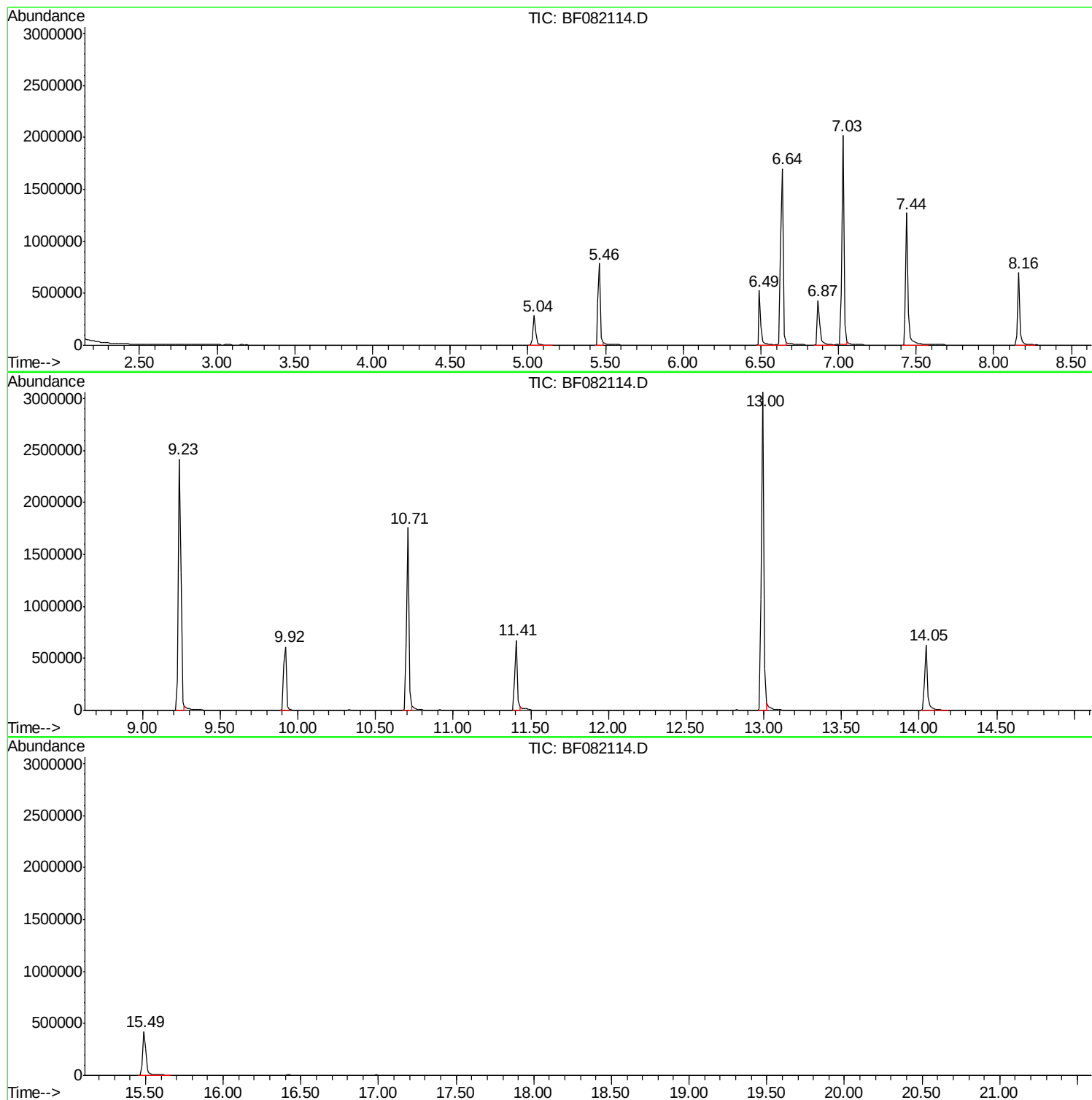
Sum of corrected areas: 18947818

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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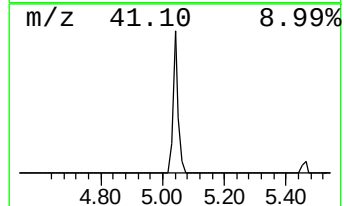
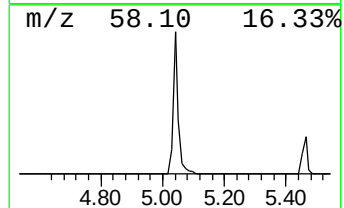
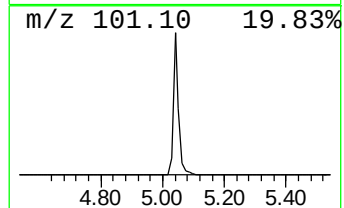
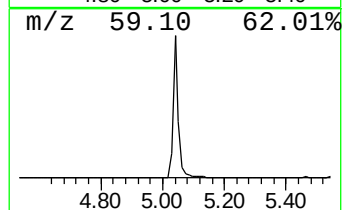
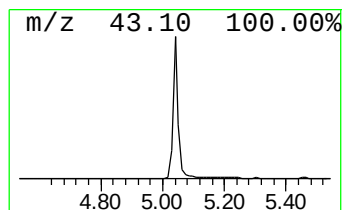
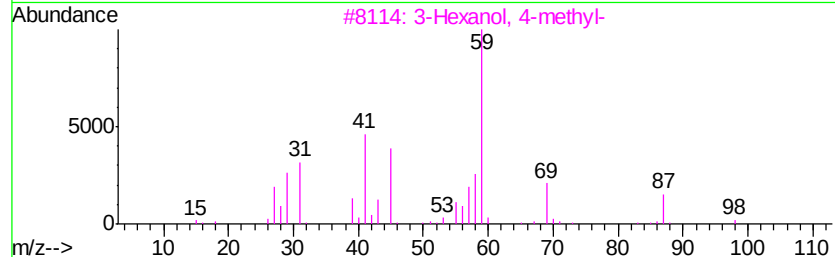
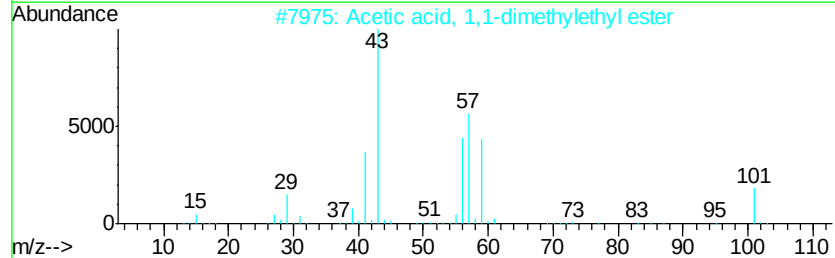
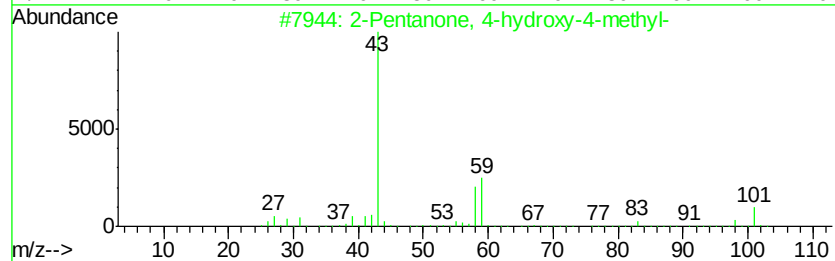
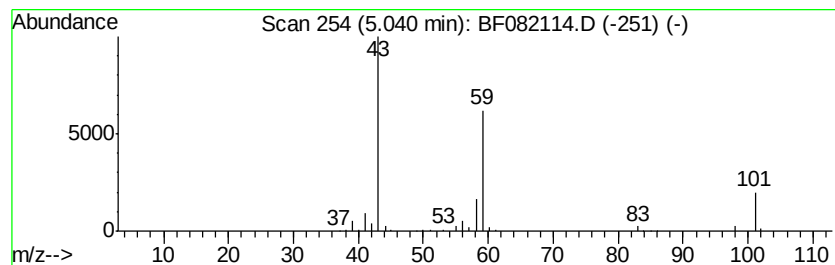
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	13.27 ng	331467	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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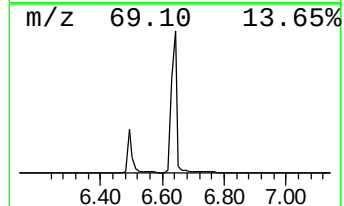
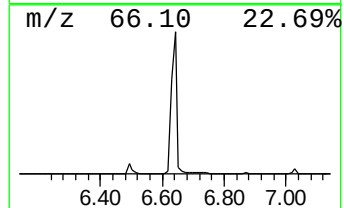
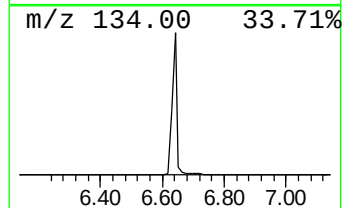
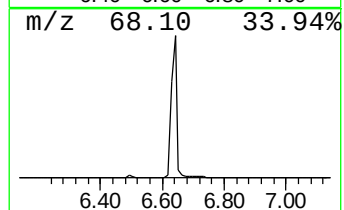
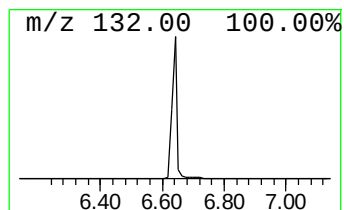
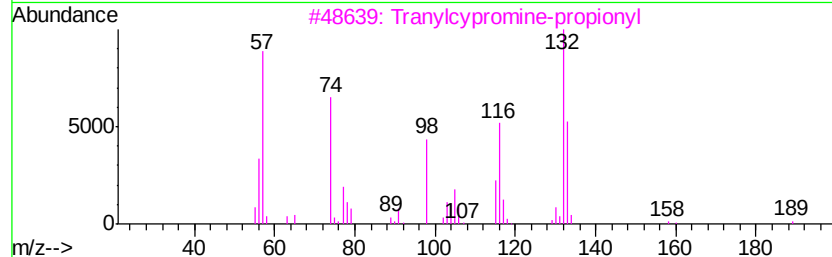
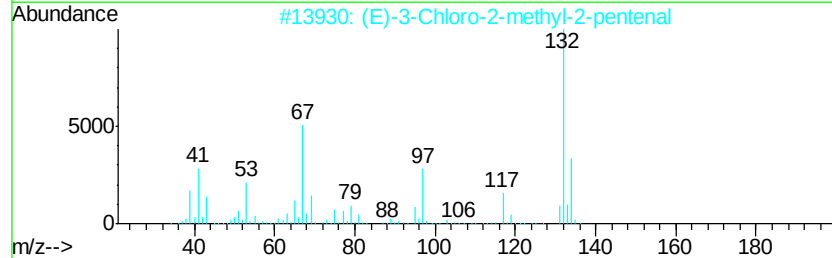
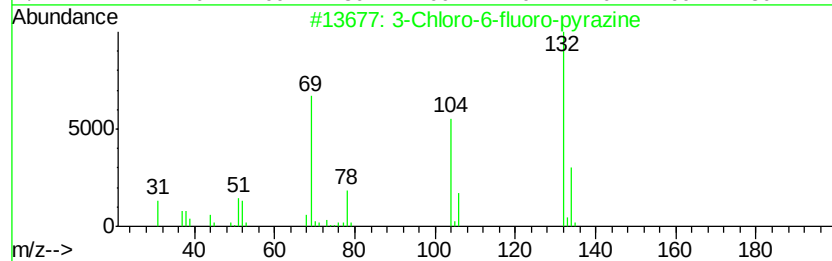
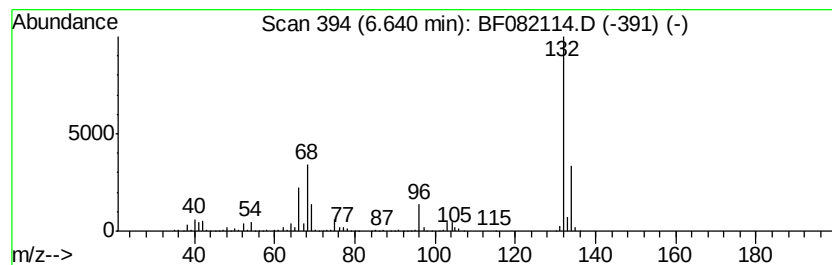
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Peak Number 2 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	77.17 ng	1927200	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Amphetaminil	250	C17H18N2	017590-01-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.04	13.3	ng	331467	1	6.87	499457	20.0
unknown6.64	6.64	77.2	ng	1927200	1	6.87	499457	20.0