

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
 Data File : BF081274.D
 Acq On : 29 Aug 2015 16:49
 Operator : UM/IZ
 Sample : PB85281BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85281BL

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	4.583	246	249	263	rBV	565568	931488	58.38%	6.664%
2	5.051	287	290	292	rBV	1150073	1489032	93.32%	10.652%
3	6.137	382	385	387	rBV	1448705	1595658	100.00%	11.415%
4	6.240	392	394	397	rBV	1092444	1317472	82.57%	9.425%
5	6.480	413	415	417	rBV	428547	366173	22.95%	2.620%
6	6.640	426	429	431	rBV	763444	1033350	64.76%	7.392%
7	7.052	462	465	467	rBV	925552	890740	55.82%	6.372%
8	7.772	525	528	530	rBV	341710	458211	28.72%	3.278%
9	8.846	619	622	625	rBV	1362826	1425581	89.34%	10.198%
10	9.509	677	680	683	rBB	603635	525543	32.94%	3.760%
11	10.298	746	749	751	rBV	723873	933113	58.48%	6.675%
12	10.972	806	808	811	rBV	397807	541648	33.95%	3.875%
13	12.561	945	947	950	rBV	1252197	1502021	94.13%	10.745%
14	13.589	1035	1037	1040	rBV	589111	528733	33.14%	3.783%
15	14.915	1150	1153	1155	rBV	413472	439606	27.55%	3.145%

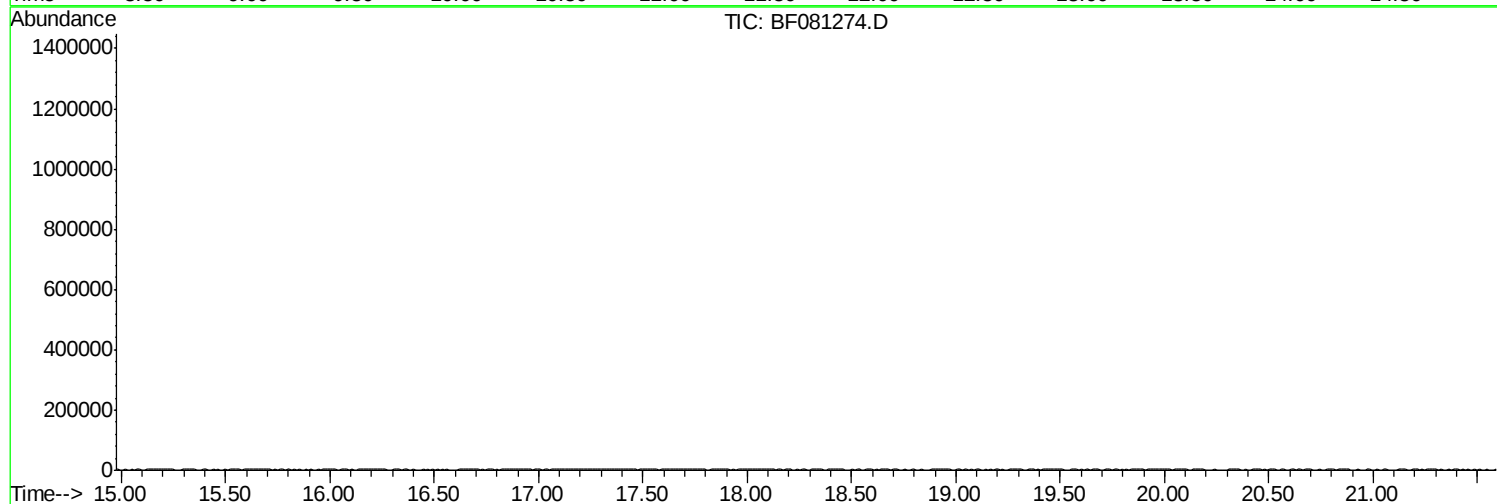
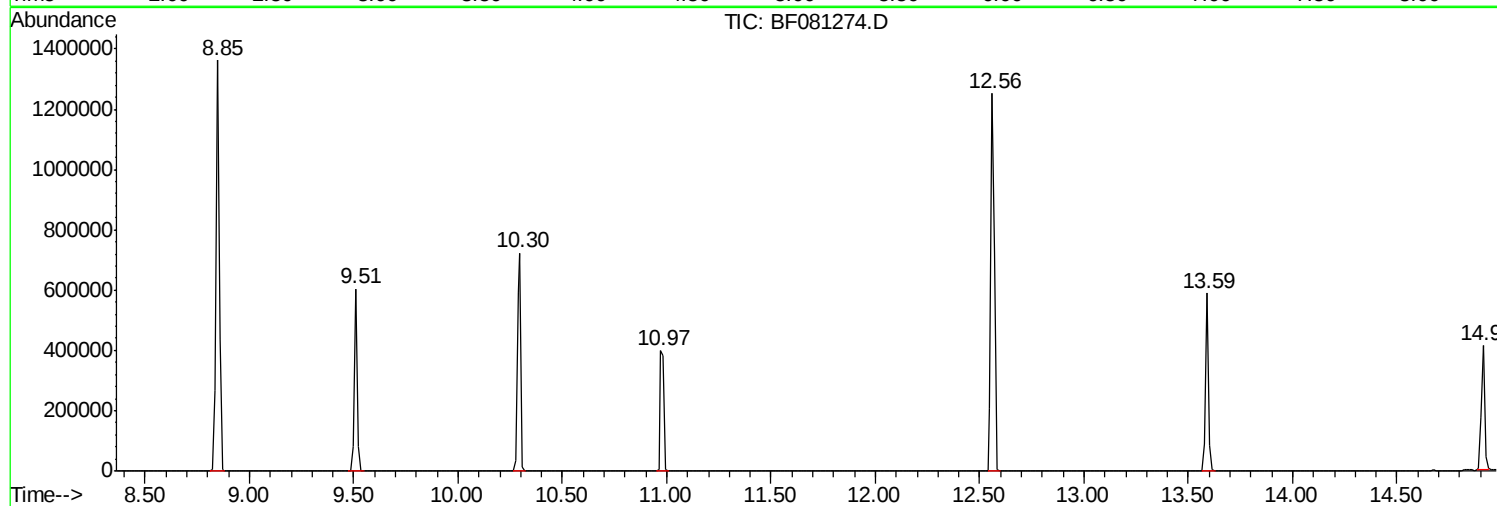
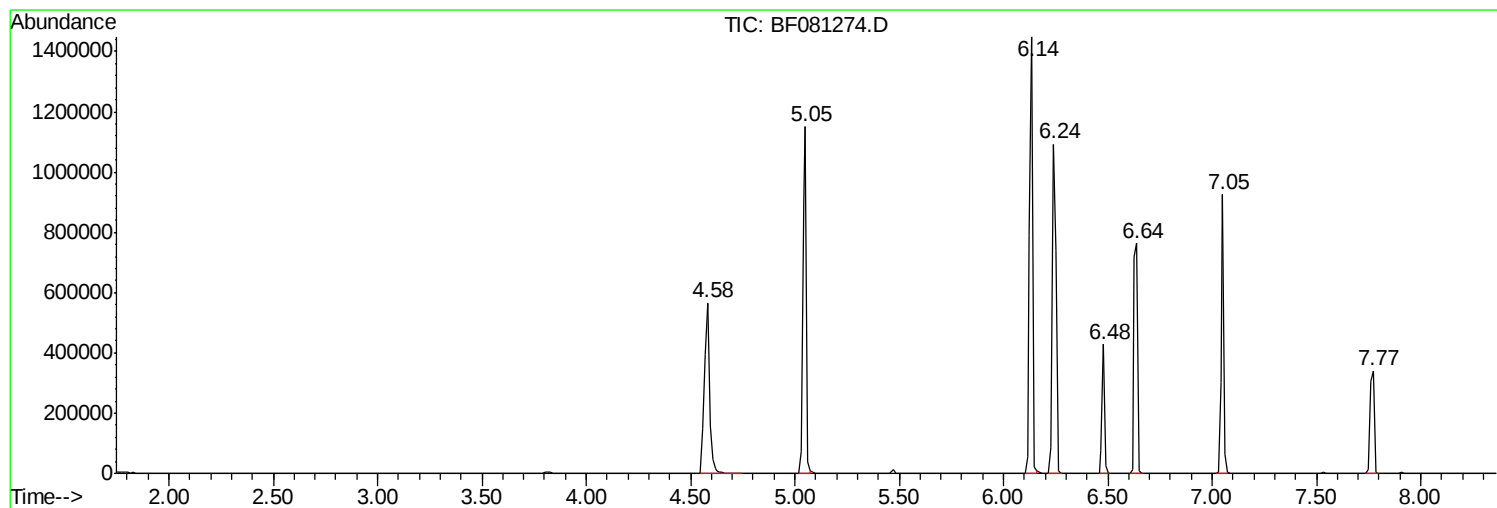
Sum of corrected areas: 13978369

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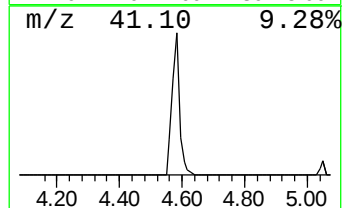
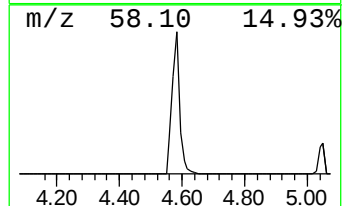
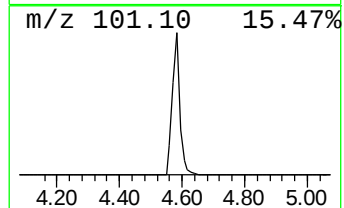
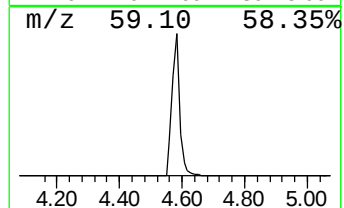
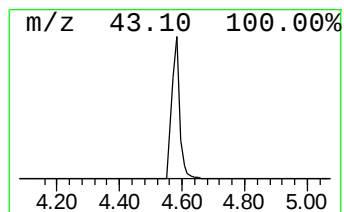
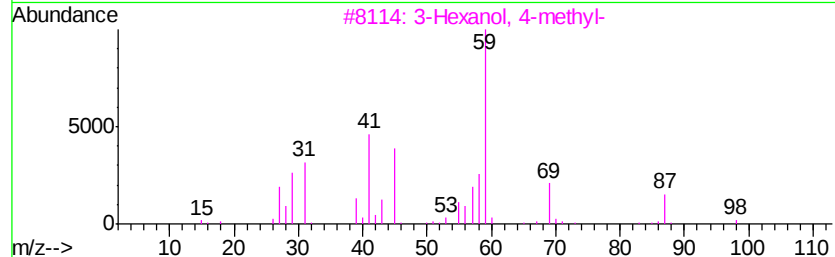
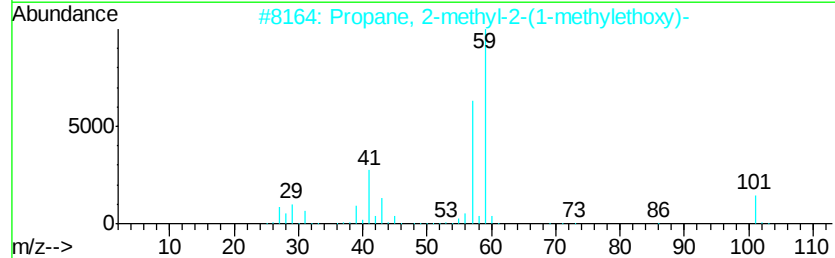
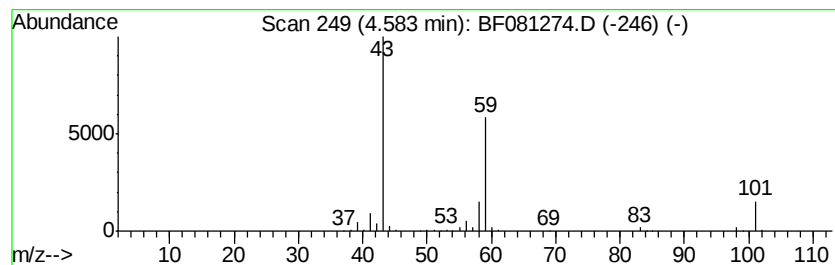
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.58	50.88 ng	931488	1,4-Dichlorobenzene-d4	6.48

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



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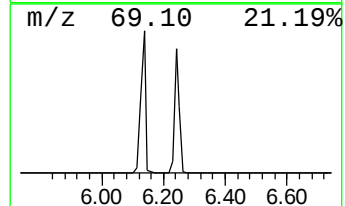
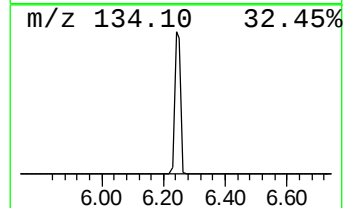
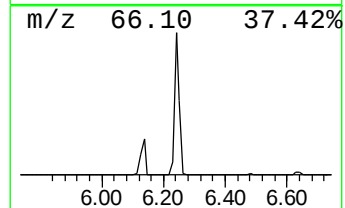
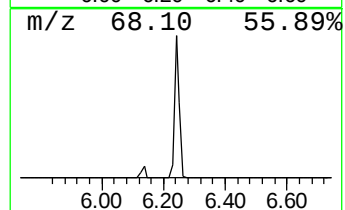
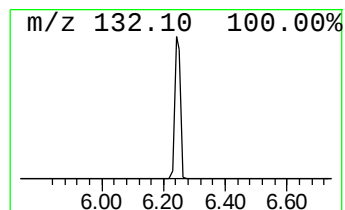
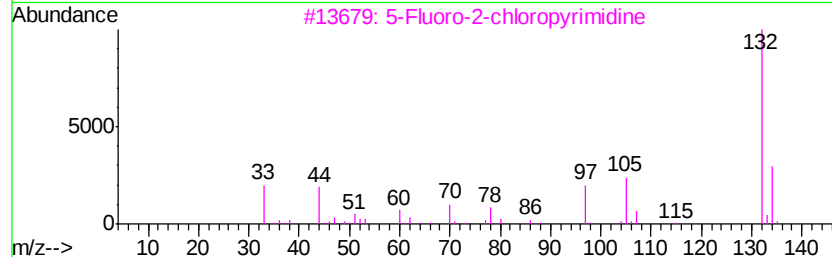
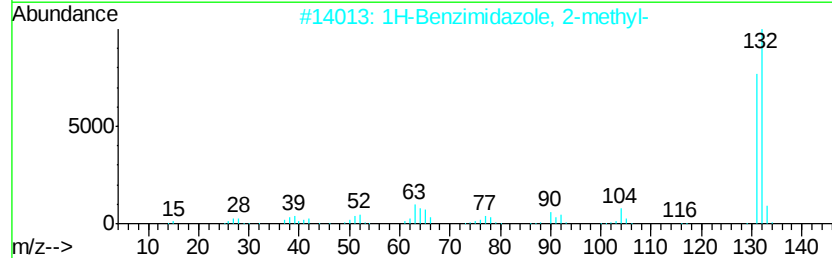
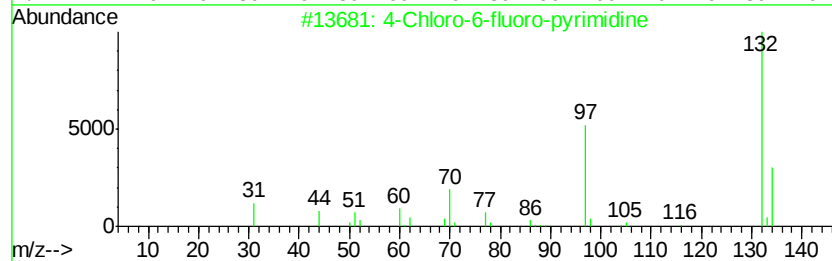
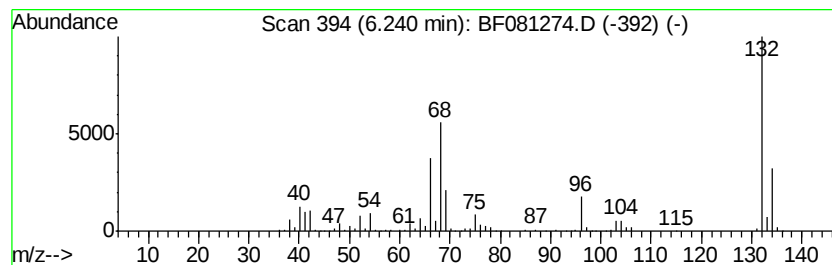
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Peak Number 2 unknown6.24 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.24	71.96 ng	1317470	1,4-Dichlorobenzene-d4	6.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.58	50.9	ng	931488	1	6.48	366173	20.0
unknown6.24	6.24	72.0	ng	1317470	1	6.48	366173	20.0