

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101416\
 Data File : BF090594.D
 Acq On : 15 Oct 2016 8:29
 Operator : UM/SJ
 Sample : H5227-04 2X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4

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-BF101316.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	5.091	239	241	250	rBV	475082	687675	25.18%	2.233%
2	5.514	276	278	281	rBV	2095220	2561499	93.79%	8.319%
3	6.543	366	368	378	rBV	1947770	2702198	98.94%	8.775%
4	6.680	378	380	383	rBV	2677928	2731173	100.00%	8.870%
5	6.920	399	401	412	rBV	1075757	1418852	51.95%	4.608%
6	7.069	412	414	417	rBV	1962053	1991364	72.91%	6.467%
7	7.480	447	450	457	rBV	1329624	1660321	60.79%	5.392%
8	8.200	511	513	516	rBV	1668922	1840610	67.39%	5.977%
9	9.286	605	608	610	rBV	1671509	2482033	90.88%	8.060%
10	9.400	616	618	620	rVB	27068	30390	1.11%	0.099%
11	9.675	640	642	644	rBV	599101	526458	19.28%	1.710%
12	9.892	660	661	663	rVB	42579	31314	1.15%	0.102%
13	9.960	664	667	669	rBV	2480068	2111928	77.33%	6.859%
14	10.338	691	700	701	rBV3	34851	155240	5.68%	0.504%
15	10.395	703	705	706	rVB	69943	66565	2.44%	0.216%
16	10.612	722	724	727	rBV	29985	46700	1.71%	0.152%
17	10.749	733	736	745	rBV	1358292	1527491	55.93%	4.961%
18	10.863	745	746	751	rVB	88481	126503	4.63%	0.411%
19	11.309	783	785	787	rBV	34237	52524	1.92%	0.171%
20	11.446	795	797	806	rBV	1547557	1839855	67.37%	5.975%
21	11.675	814	817	821	rBV2	72906	104881	3.84%	0.341%
22	11.743	821	823	825	rVB	28267	28127	1.03%	0.091%
23	11.972	840	843	845	rBV2	12619	31093	1.14%	0.101%
24	12.063	849	851	855	rVB2	23343	37706	1.38%	0.122%
25	12.498	887	889	891	rBV2	21832	35300	1.29%	0.115%
26	12.669	901	904	906	rBV	149502	200353	7.34%	0.651%
27	12.886	921	923	931	rBV	169894	257935	9.44%	0.838%
28	13.035	933	936	947	rBV	1843653	2100159	76.90%	6.820%
29	13.927	1012	1014	1021	rVB2	71623	132861	4.86%	0.431%
30	14.087	1025	1028	1036	rBV	1107718	1599849	58.58%	5.196%
31	15.127	1117	1119	1123	rBV	81723	196019	7.18%	0.637%
32	15.550	1154	1156	1167	rVB	766836	1411947	51.70%	4.585%
33	16.018	1195	1197	1199	rBV	44035	65812	2.41%	0.214%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

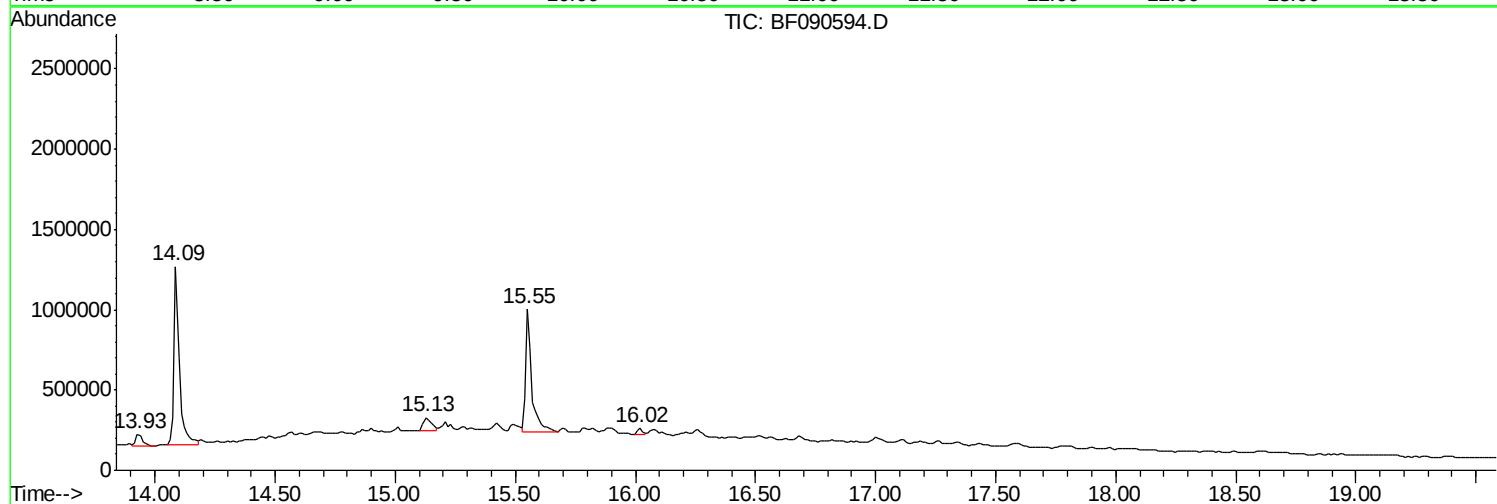
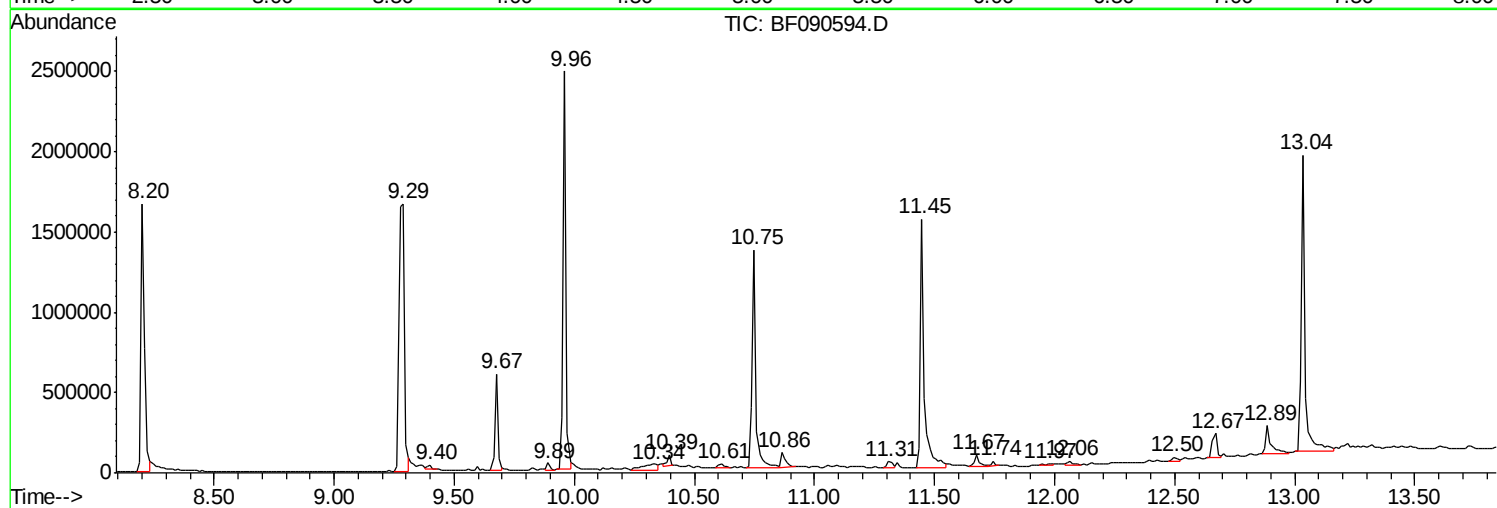
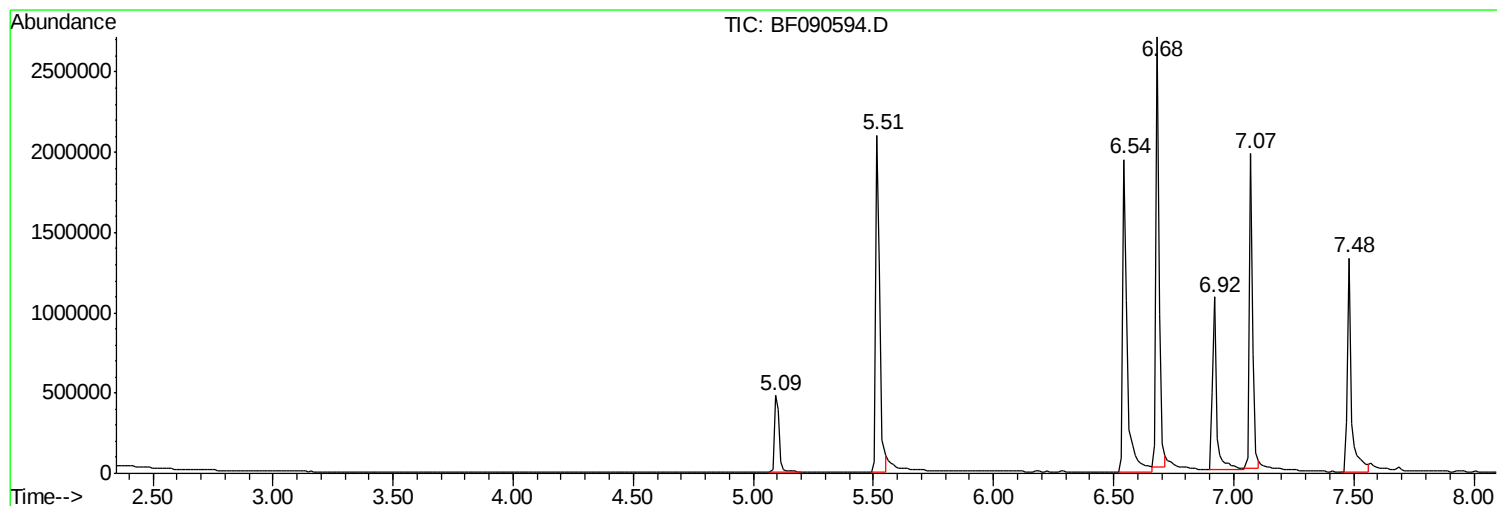
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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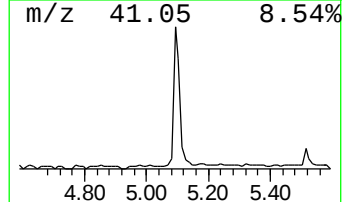
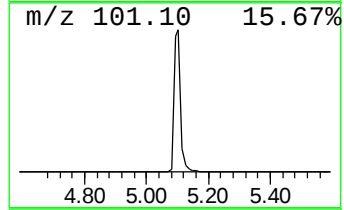
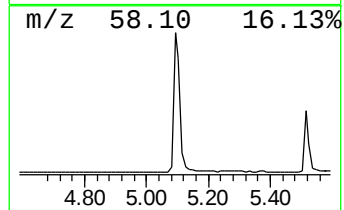
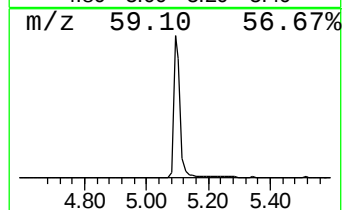
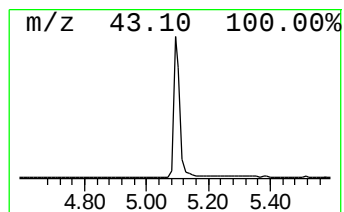
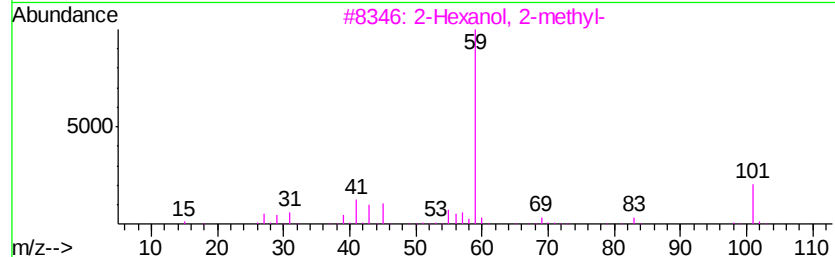
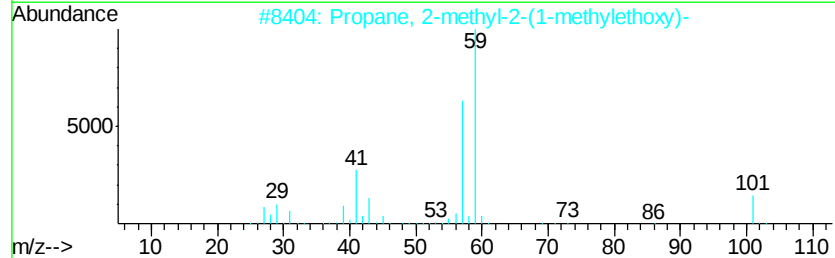
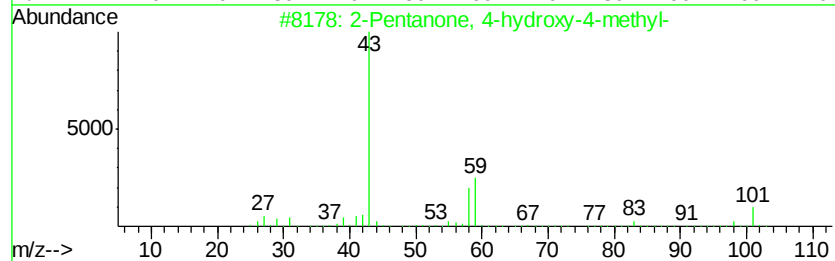
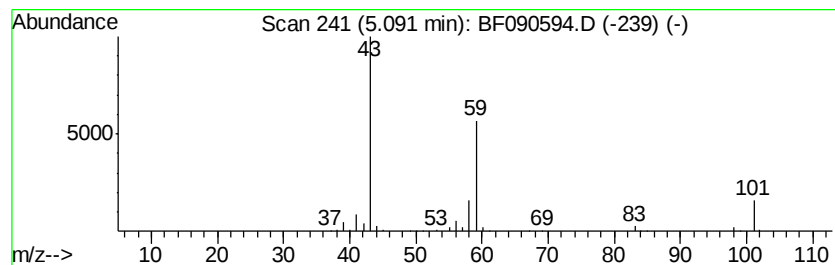
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	9.69 ng	687675	1,4-Dichlorobenzene-d4	6.92

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	42	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40	
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	



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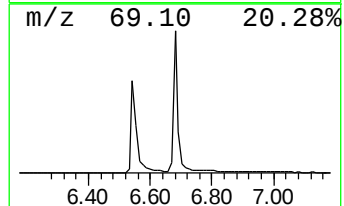
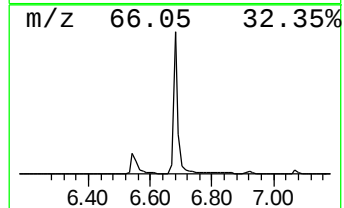
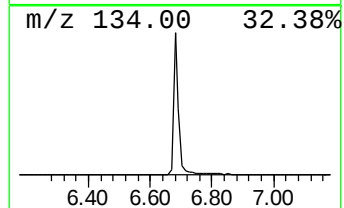
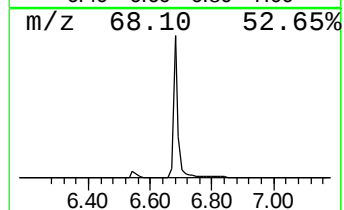
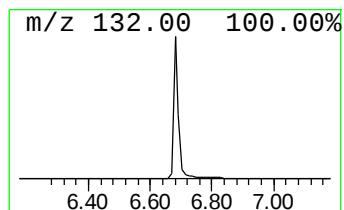
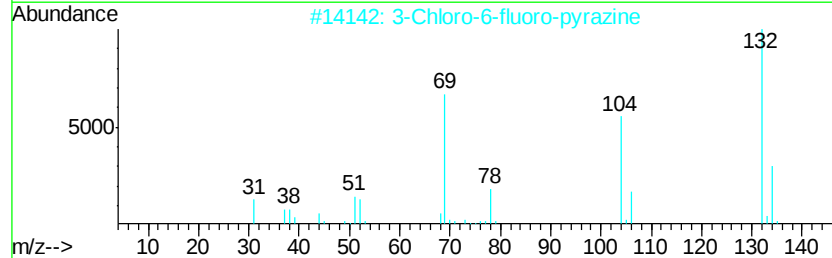
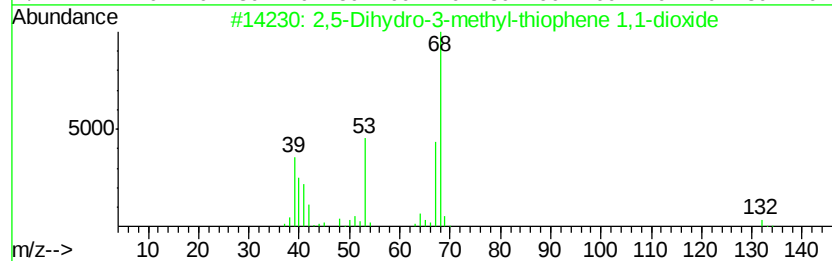
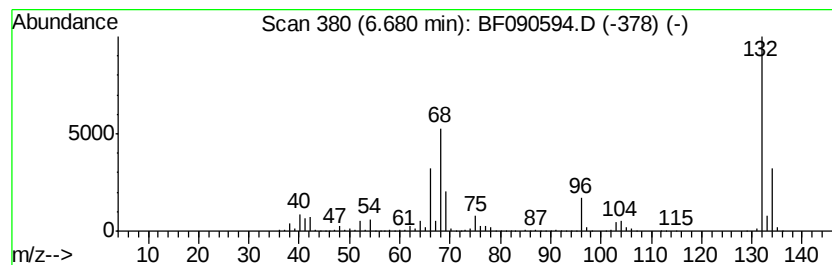
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Peak Number 2 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	38.50 ng	2731170	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Dihydro-3-methyl-thiophene 1...	132	C5H8O2S	001193-10-8	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.09	9.7	ng	687675	1	6.92	1418850	20.0
unknown6.68	6.68	38.5	ng	2731170	1	6.92	1418850	20.0