

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072716\
 Data File : BF089331.D
 Acq On : 27 Jul 2016 16:23
 Operator : UM/SJ
 Sample : PB92412BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92412BL

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-BF072716.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	1.690	7	9	56	rVB	868279	1964474	100.00%	14.190%
2	4.673	267	270	282	rBV	100853	209740	10.68%	1.515%
3	5.130	308	310	329	rBV	901487	1281792	65.25%	9.259%
4	6.205	402	404	406	rBV	897447	1201342	61.15%	8.678%
5	6.307	411	413	432	rVV	1094402	1331119	67.76%	9.615%
6	6.547	432	434	445	rVV	153995	241957	12.32%	1.748%
7	6.696	445	447	461	rVB	1045859	1036468	52.76%	7.487%
8	7.119	482	484	510	rBV	513963	836223	42.57%	6.040%
9	7.828	544	546	557	rBV	298928	345264	17.58%	2.494%
10	8.913	638	641	643	rBV	971448	1525574	77.66%	11.020%
11	9.576	696	699	701	rBV	250652	374170	19.05%	2.703%
12	10.354	765	767	770	rBV2	669930	865882	44.08%	6.255%
13	11.039	825	827	844	rBV	284559	415177	21.13%	2.999%
14	12.628	964	966	969	rBV	1476918	1544174	78.60%	11.154%
15	13.657	1054	1056	1066	rBV	239024	374667	19.07%	2.706%
16	14.994	1171	1173	1181	rBV	184849	295607	15.05%	2.135%

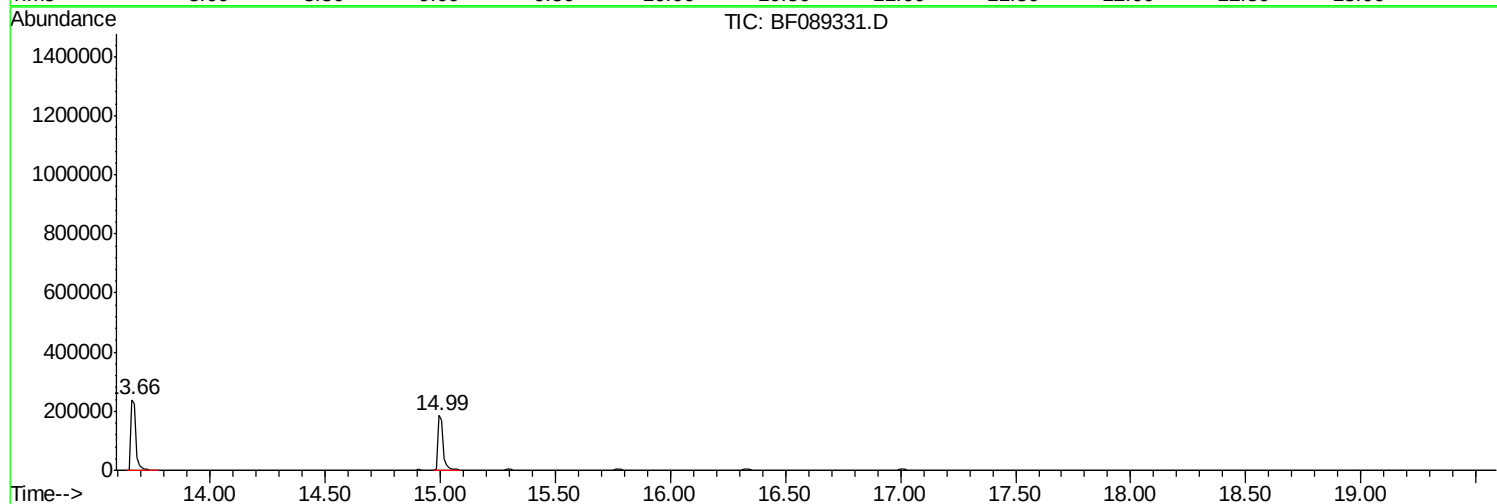
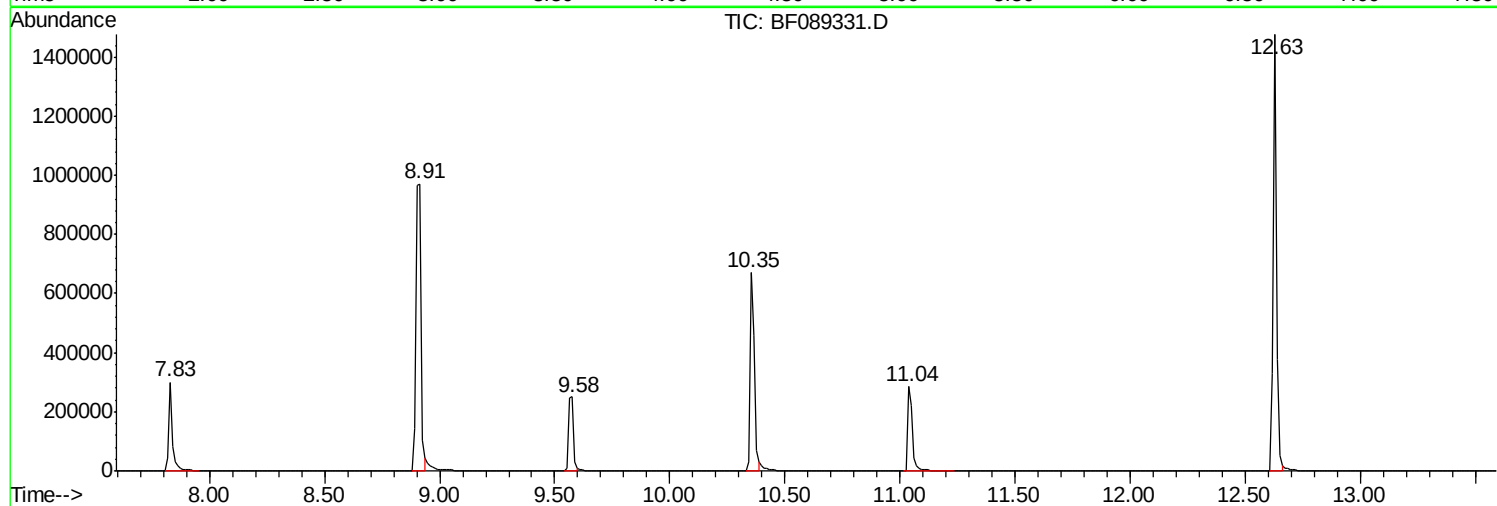
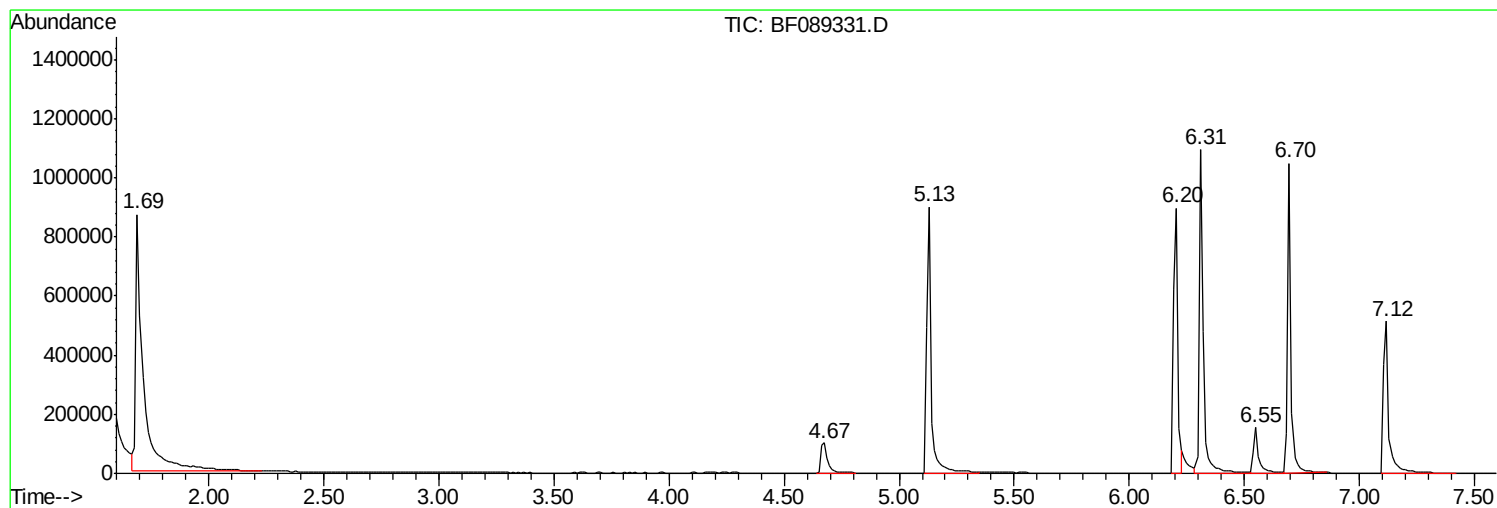
Sum of corrected areas: 13843630

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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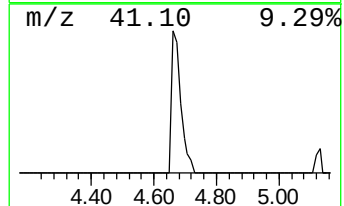
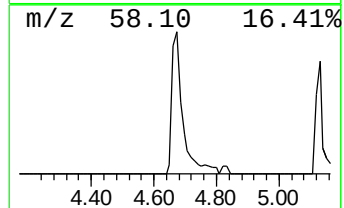
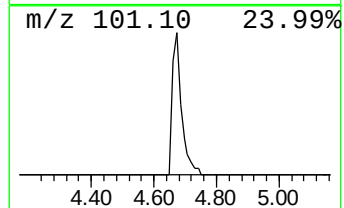
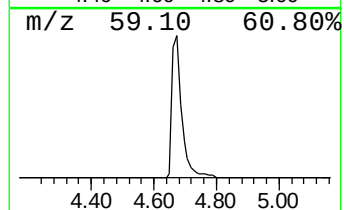
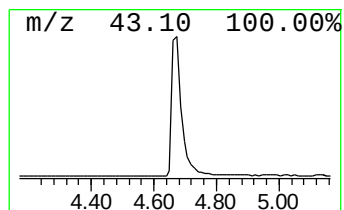
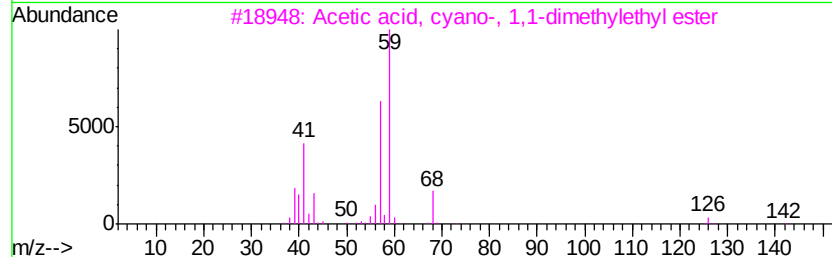
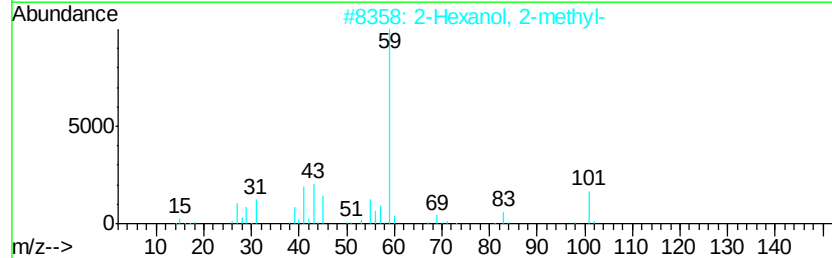
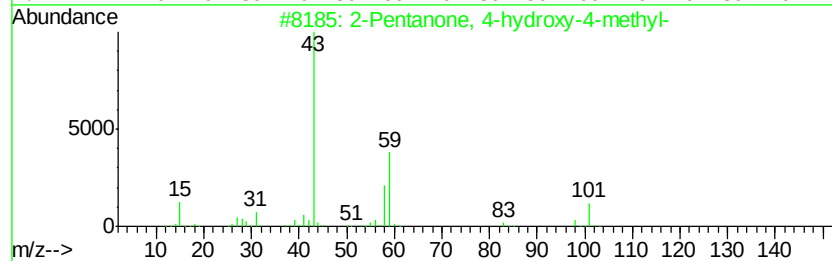
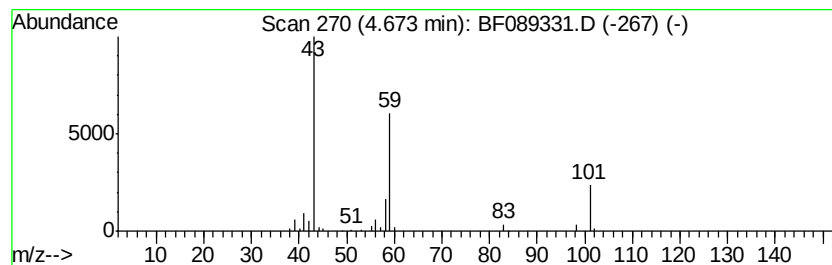
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.67	17.34 ng	209740	1,4-Dichlorobenzene-d4	6.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
5		Acetone	58	C3H6O	000067-64-1	10



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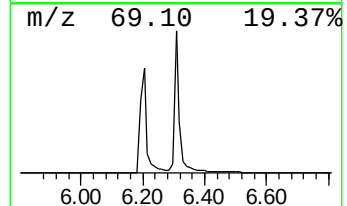
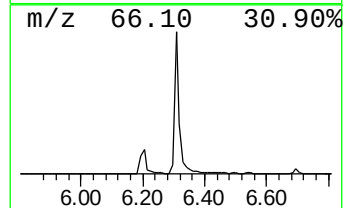
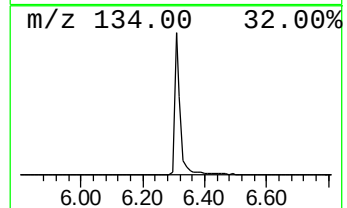
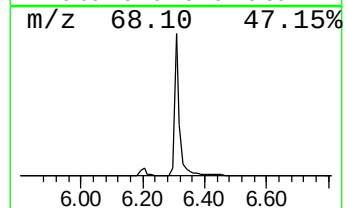
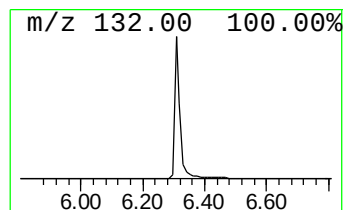
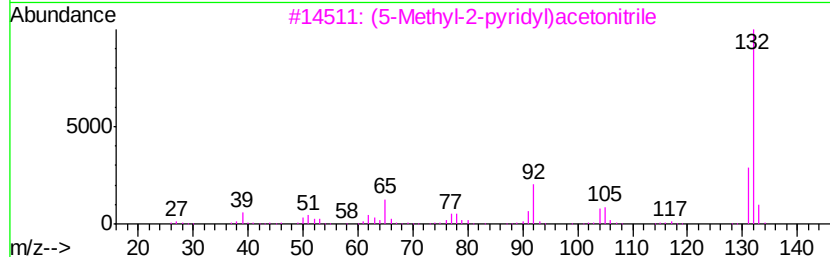
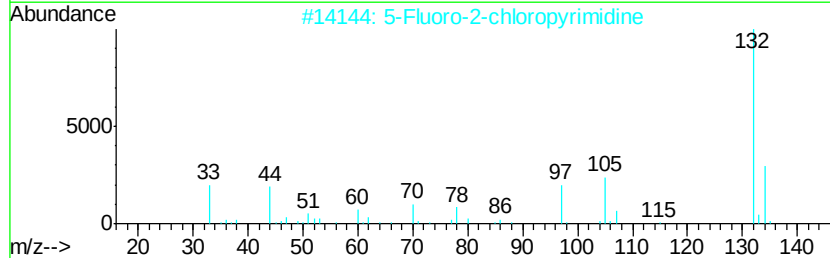
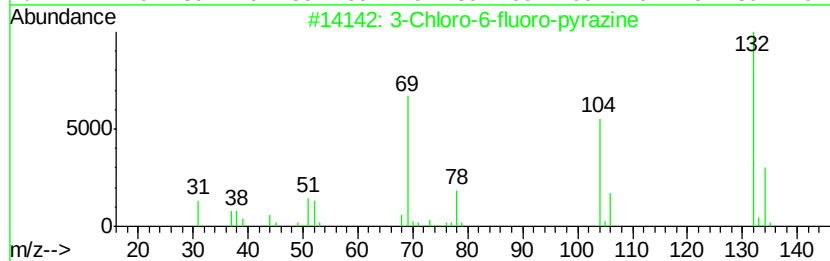
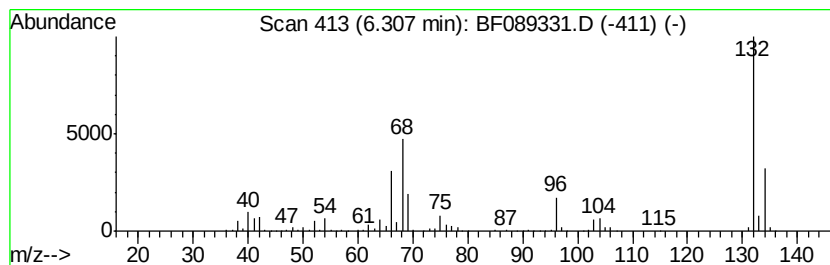
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Peak Number 3 unknown6.31 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	110.03 ng	1331120	1,4-Dichlorobenzene-d4	6.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.67	17.3	ng	209740	1	6.55	241957	20.0
unknown6.31	6.31	110.0	ng	1331120	1	6.55	241957	20.0