

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102716\
 Data File : BF090875.D
 Acq On : 27 Oct 2016 19:14
 Operator : UM/SJ
 Sample : PB94343BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94343BL

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-BF102616.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	4.944	247	250	260	rVB	410768	610573	13.45%	2.029%
2	5.367	285	287	290	rBV	2435427	2629686	57.95%	8.739%
3	6.407	376	378	381	rBV	2196538	2499387	55.08%	8.306%
4	6.533	387	389	392	rBV	2266352	2996039	66.02%	9.957%
5	6.773	408	410	421	rBV	723722	804053	17.72%	2.672%
6	6.922	421	423	426	rBV	1950705	2479646	54.64%	8.241%
7	7.333	457	459	462	rBV	1420206	1610022	35.48%	5.351%
8	8.053	520	522	525	rBV	796748	1057484	23.30%	3.514%
9	9.139	614	617	619	rBV	4220440	4088081	90.09%	13.586%
10	9.813	673	676	678	rBV	1335672	1292298	28.48%	4.295%
11	10.602	742	745	747	rBV	2464282	2443012	53.84%	8.119%
12	11.299	803	806	808	rBV	848978	1231657	27.14%	4.093%
13	12.888	942	945	947	rBV	3885202	4537924	100.00%	15.081%
14	13.928	1034	1036	1047	rBV	1055822	1047768	23.09%	3.482%
15	15.334	1156	1159	1162	rBV	570006	708736	15.62%	2.355%
16	15.700	1187	1191	1197	rBV2	12956	54125	1.19%	0.180%

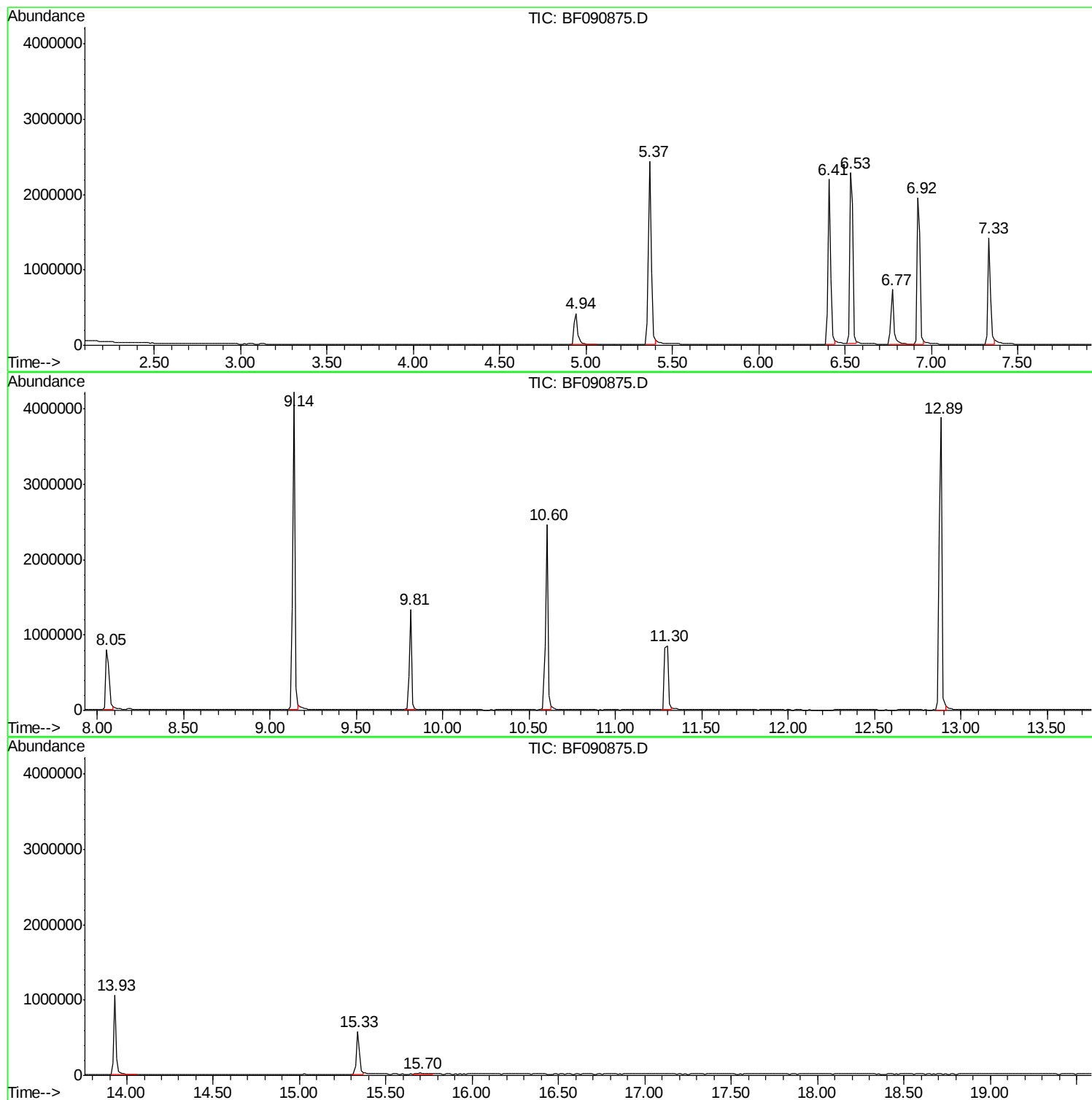
Sum of corrected areas: 30090491

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102616.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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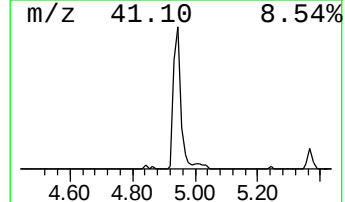
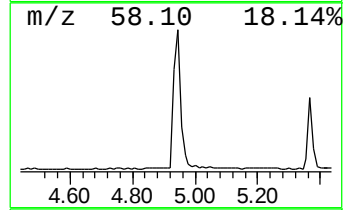
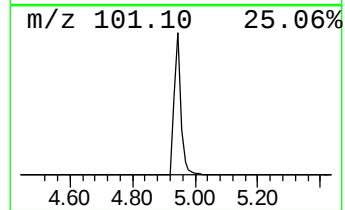
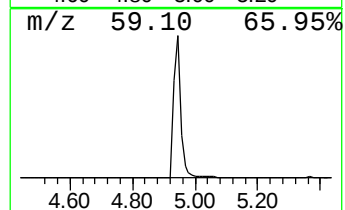
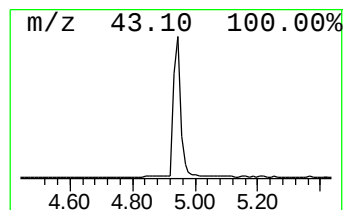
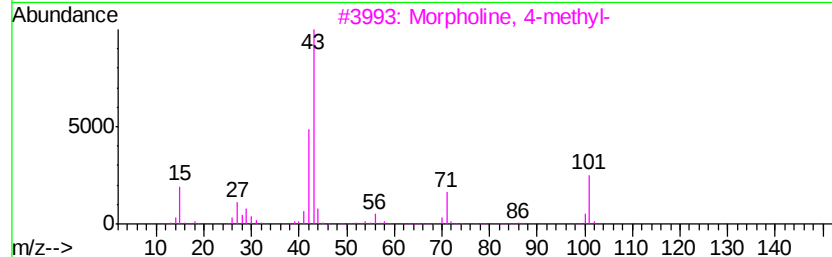
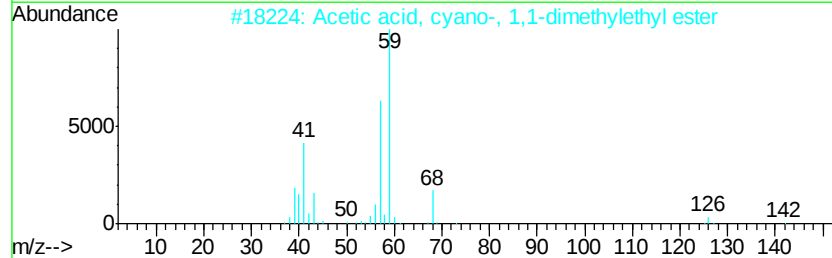
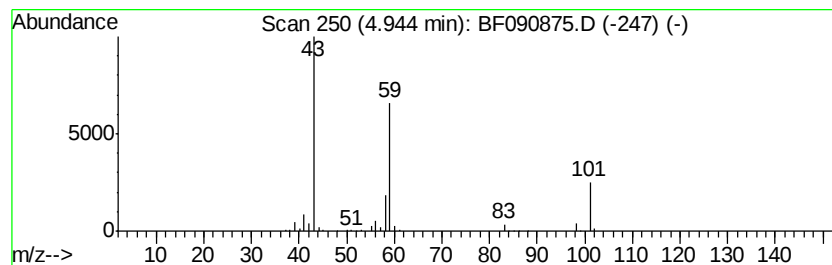
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	15.19 ng	610573	1,4-Dichlorobenzene-d4	6.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17	
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
4	Acetone	58	C3H6O	000067-64-1	9	
5	Guanidine	59	CH5N3	000113-00-8	9	



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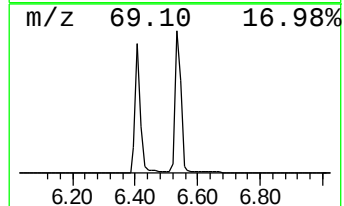
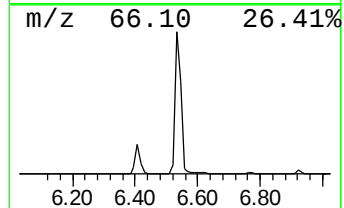
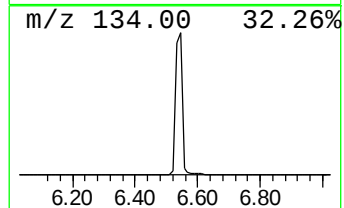
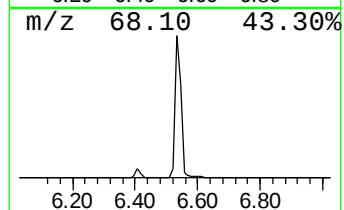
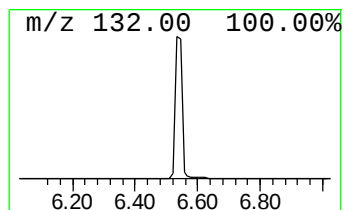
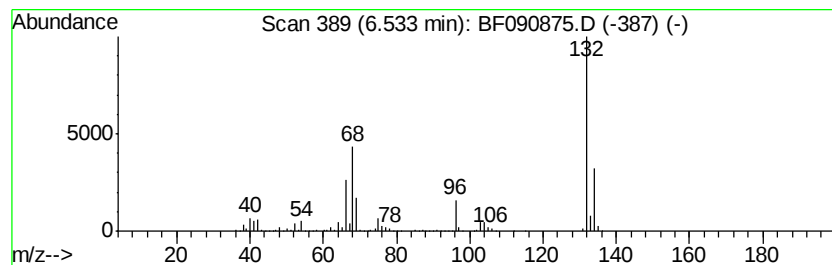
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	74.52 ng	2996040	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.94	15.2	ng	610573	1	6.77	804053	20.0
unknown6.53	6.53	74.5	ng	2996040	1	6.77	804053	20.0