

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060116\
 Data File : BF087623.D
 Acq On : 1 Jun 2016 18:33
 Operator : UM/SJ
 Sample : PB91016BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91016BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.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	4.684	269	271	292	rBV	94706	351213	8.68%	1.077%
2	5.336	325	328	368	rBV	1671321	3694174	91.30%	11.325%
3	6.982	469	472	478	rBV	2100540	3303881	81.66%	10.128%
4	7.073	478	480	508	rVV	2541559	4046097	100.00%	12.404%
5	7.416	508	510	525	rVV	468279	769519	19.02%	2.359%
6	7.645	528	530	552	rVB	1631105	2732887	67.54%	8.378%
7	8.342	589	591	617	rBV	1255673	2447259	60.48%	7.502%
8	9.451	686	688	710	rVB	539151	935117	23.11%	2.867%
9	11.222	840	843	846	rBV	2747458	3734879	92.31%	11.450%
10	12.274	933	935	947	rBV	716546	1033037	25.53%	3.167%
11	13.565	1046	1048	1068	rBV	1449348	2586100	63.92%	7.928%
12	14.685	1143	1146	1160	rBV	543577	1076581	26.61%	3.300%
13	17.040	1349	1352	1355	rBV	3329531	3943004	97.45%	12.088%
14	18.400	1469	1471	1485	rBV	605495	1012286	25.02%	3.103%
15	20.080	1616	1618	1634	rBV	405793	953830	23.57%	2.924%

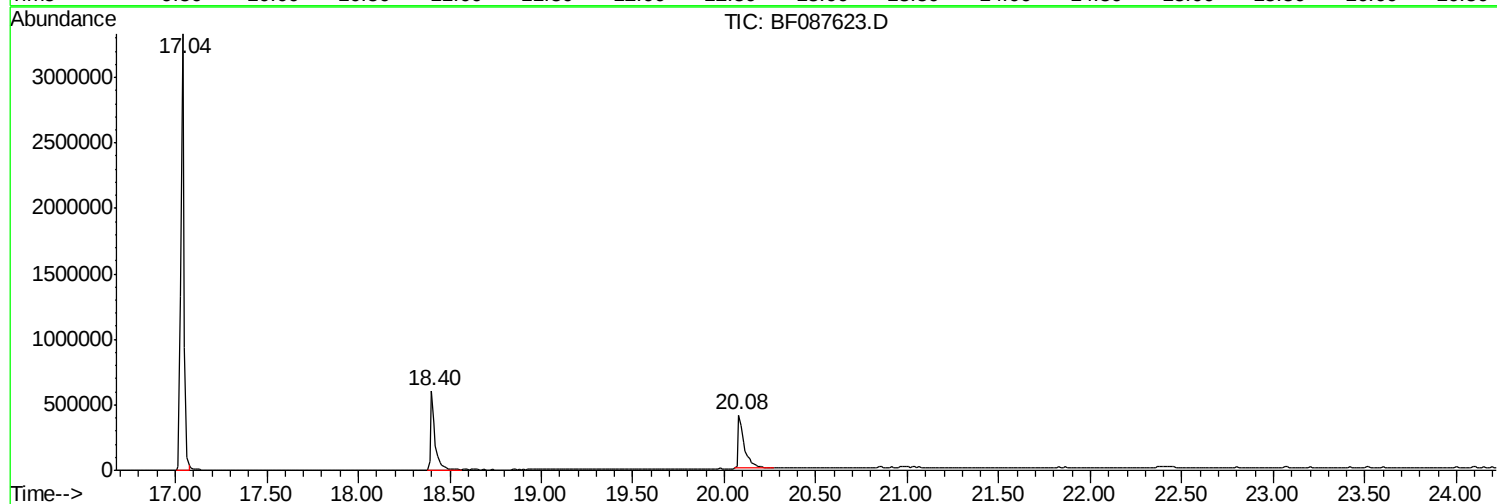
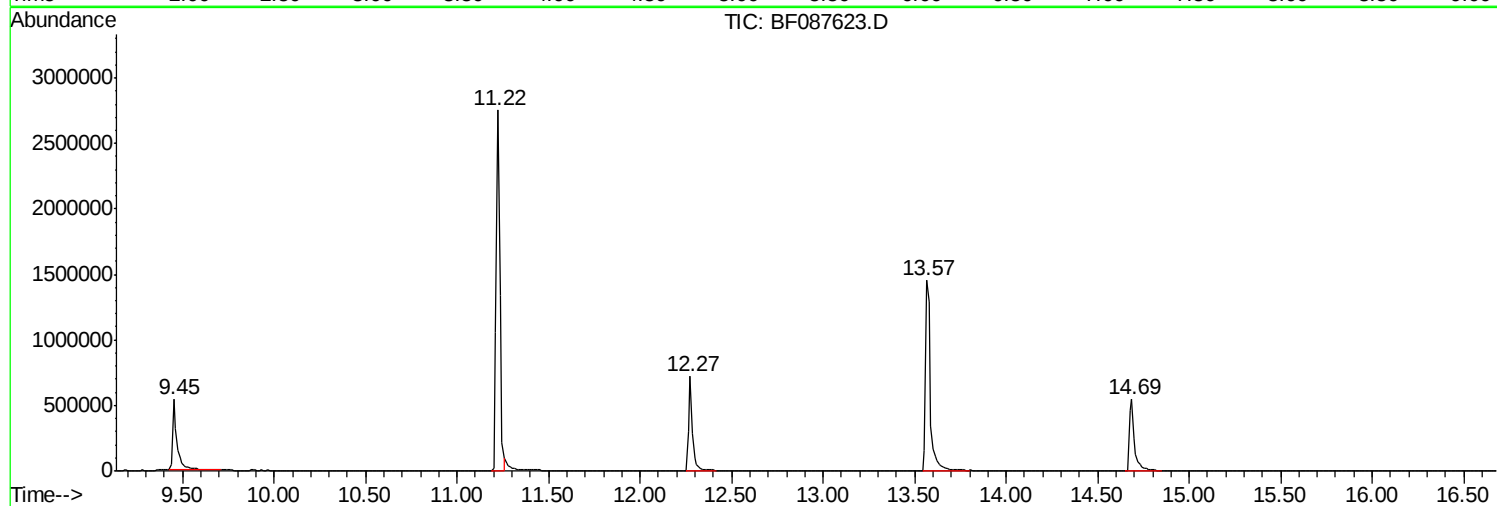
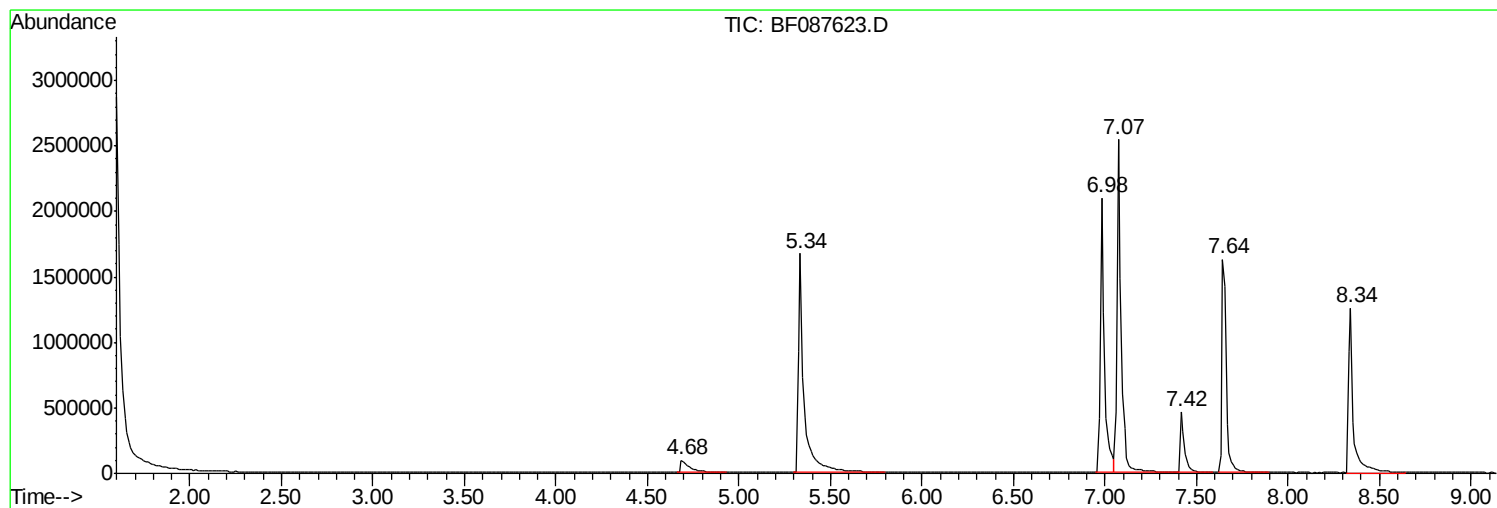
Sum of corrected areas: 32619864

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.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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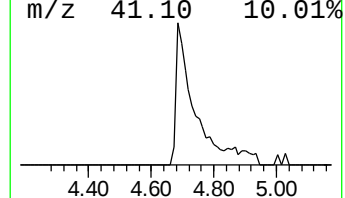
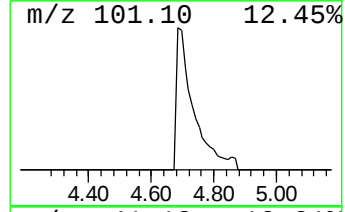
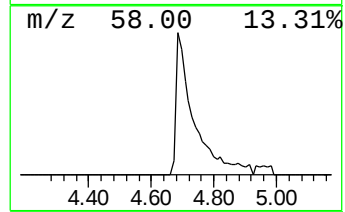
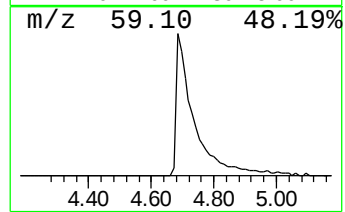
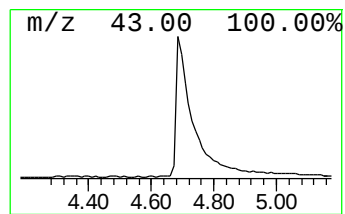
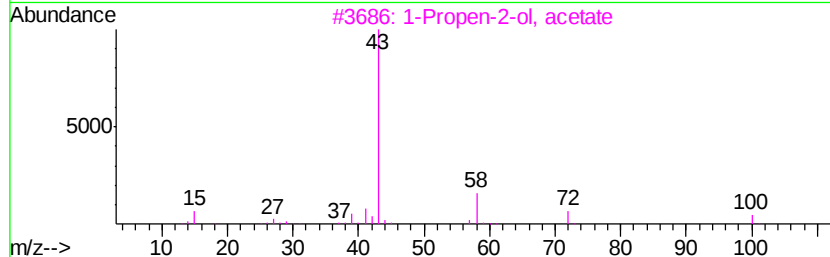
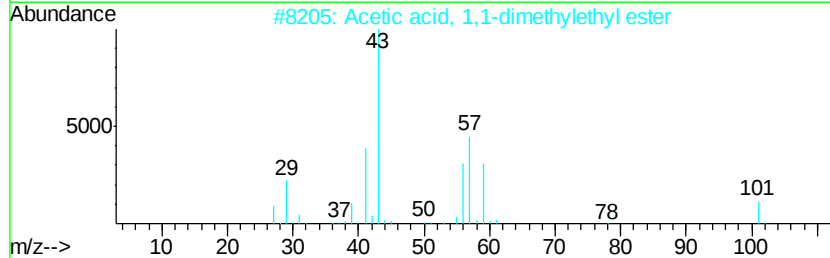
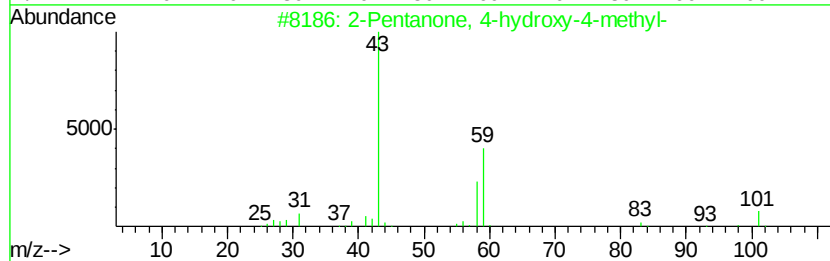
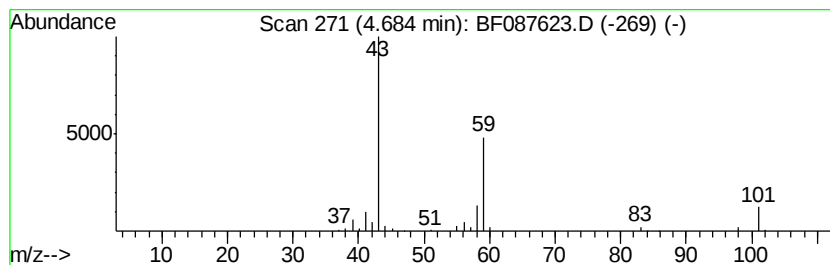
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.68	9.13 ng	351213	1,4-Dichlorobenzene-d4	7.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10
5	Butane	58	C4H10	000106-97-8	9



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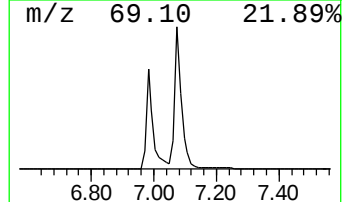
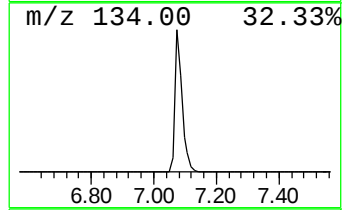
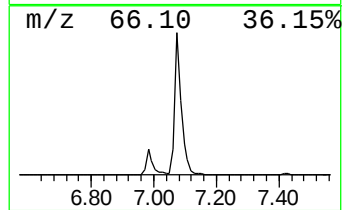
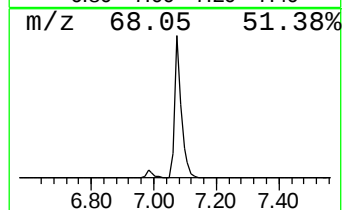
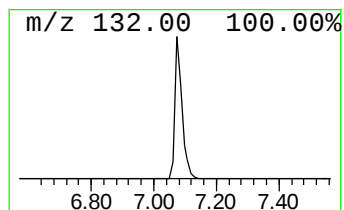
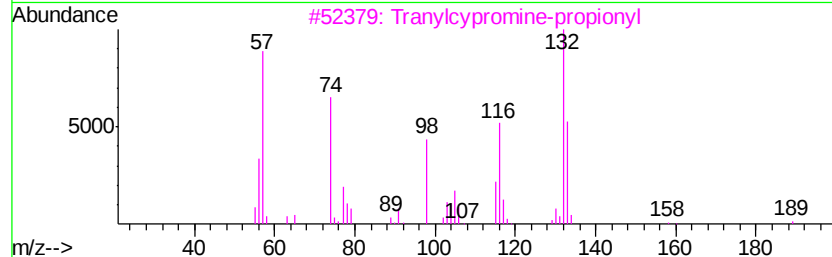
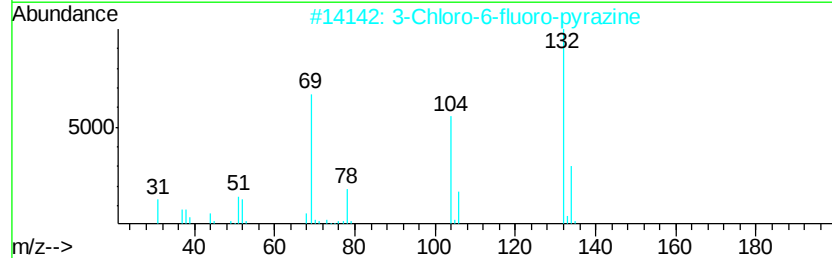
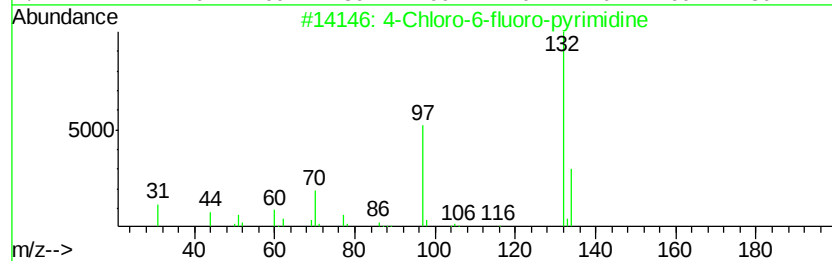
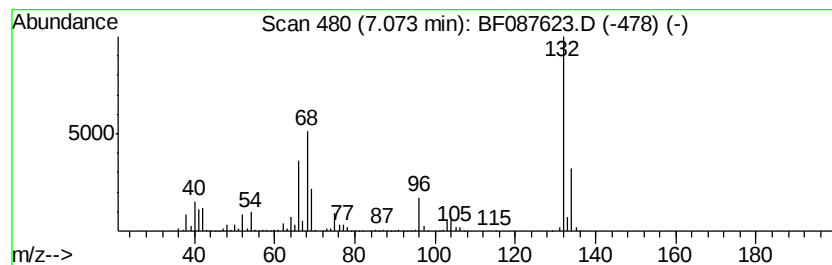
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Peak Number 2 unknown7.07 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	105.16 ng	4046100	1,4-Dichlorobenzene-d4	7.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.68	9.1	ng	351213	1	7.42	769519	20.0
unknown7.07	7.07	105.2	ng	4046100	1	7.42	769519	20.0