

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080416\
 Data File : BF089618.D
 Acq On : 5 Aug 2016 8:04
 Operator : UM/SJ
 Sample : H4288-24
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-DUP-04-072816

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-BF080416.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	1.701	5	10	12	rVB	3504344	7379792	100.00%	43.836%
2	4.650	265	268	278	rBV	40445	97009	1.31%	0.576%
3	5.324	326	327	348	rBV7	169619	447050	6.06%	2.655%
4	6.982	470	472	479	rBV	100311	257521	3.49%	1.530%
5	7.085	479	481	499	rVB	632257	1020614	13.83%	6.062%
6	7.450	511	513	530	rBV	103973	234653	3.18%	1.394%
7	7.679	530	533	549	rBV	698683	967418	13.11%	5.746%
8	8.353	590	592	614	rBV	473256	794474	10.77%	4.719%
9	9.473	687	690	716	rBV	202181	357627	4.85%	2.124%
10	11.256	842	846	848	rBV	1023833	1547792	20.97%	9.194%
11	12.296	934	937	943	rBV	259912	423114	5.73%	2.513%
12	13.577	1047	1049	1066	rVB	543823	873163	11.83%	5.187%
13	14.685	1144	1146	1154	rBV	287386	410484	5.56%	2.438%
14	17.051	1350	1353	1356	rBV	1113411	1460252	19.79%	8.674%
15	18.389	1468	1470	1480	rBV	188610	327503	4.44%	1.945%
16	20.046	1612	1615	1623	rBV	170414	236596	3.21%	1.405%

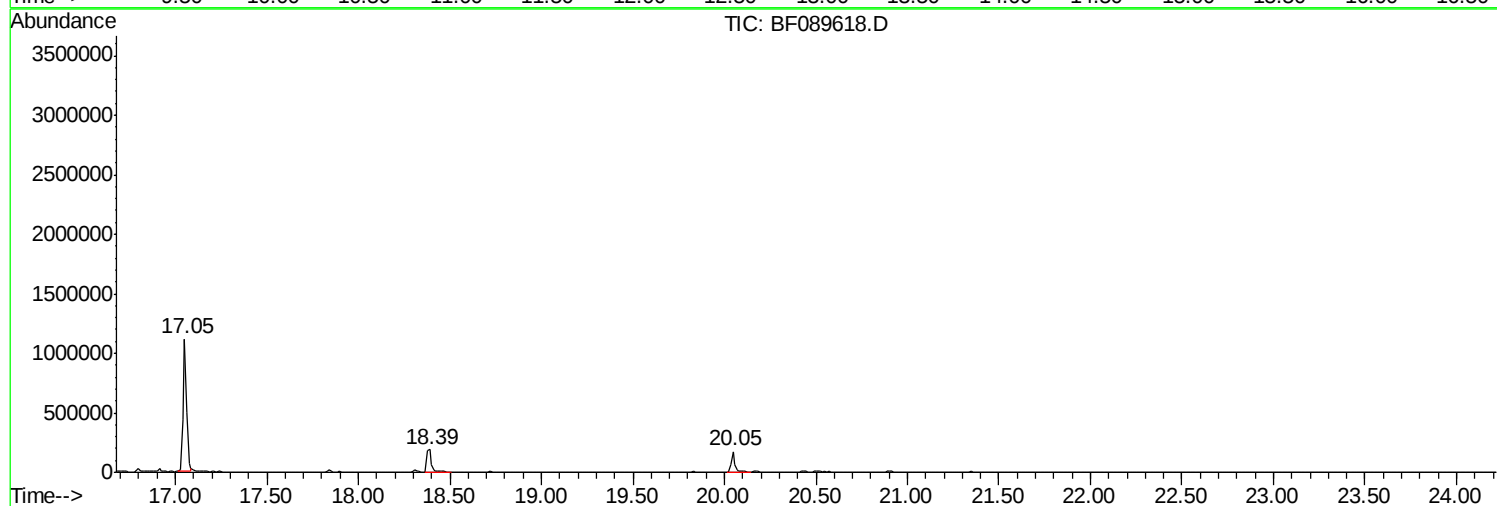
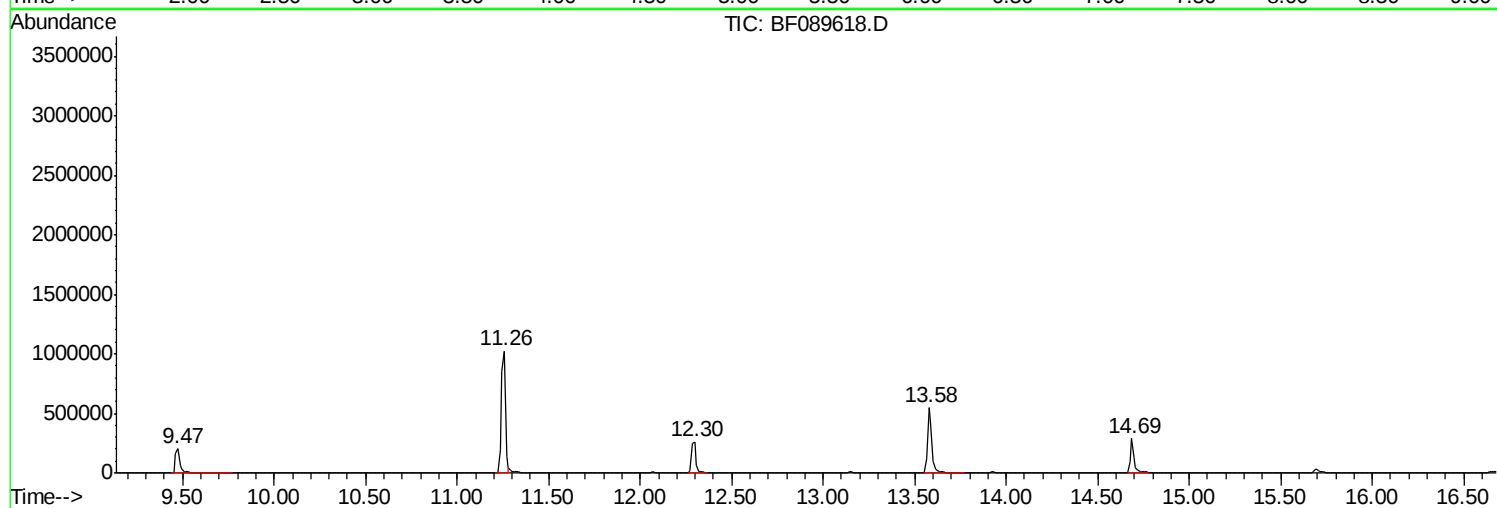
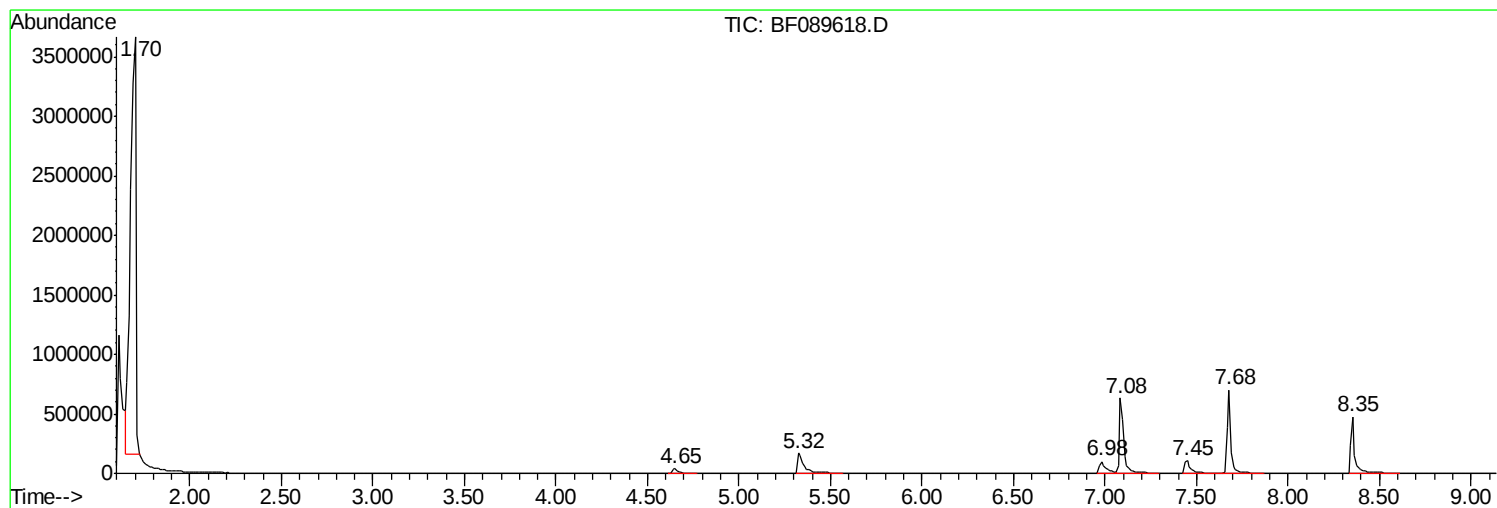
Sum of corrected areas: 16835062

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.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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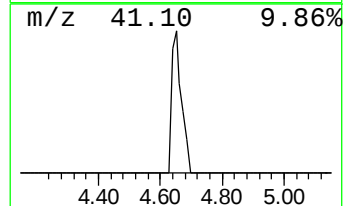
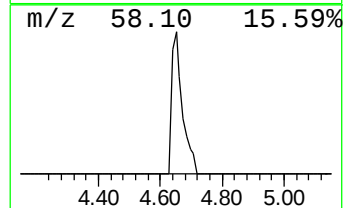
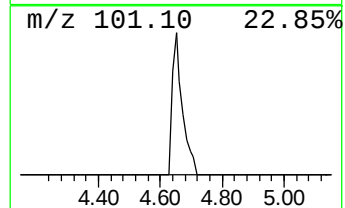
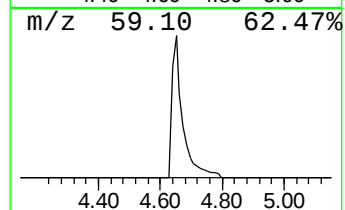
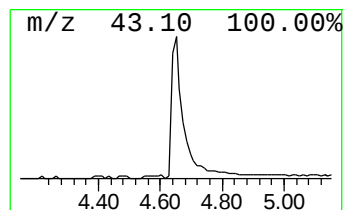
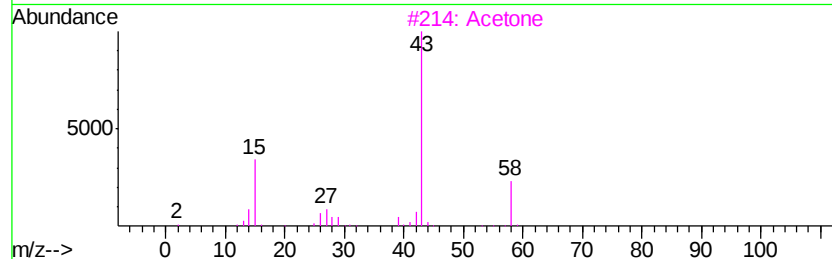
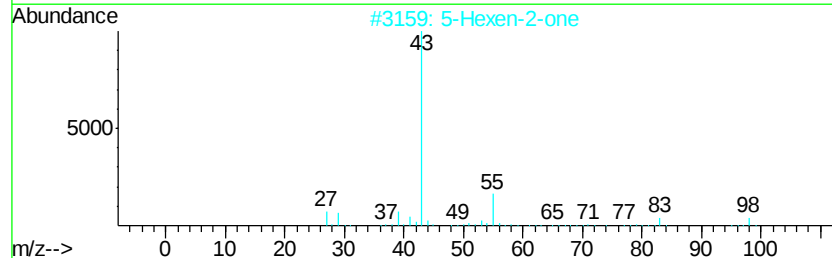
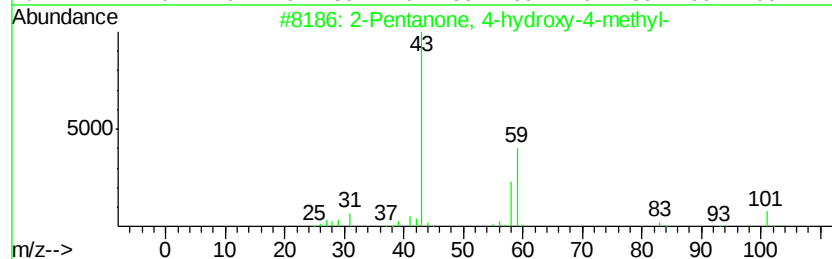
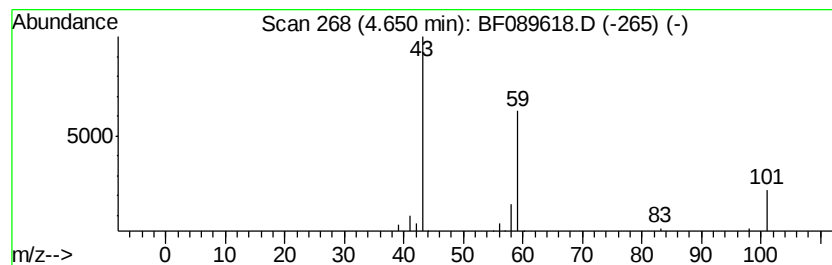
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.65	8.27 ng	97009	1,4-Dichlorobenzene-d4	7.45	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9
3	Acetone	58	C3H6O	000067-64-1	9
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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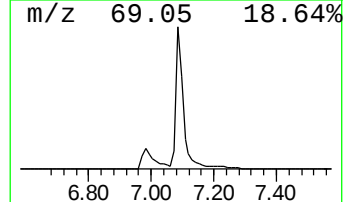
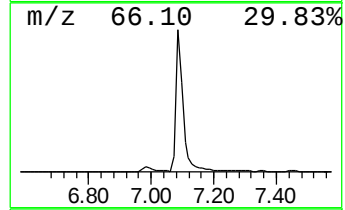
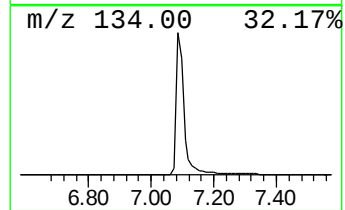
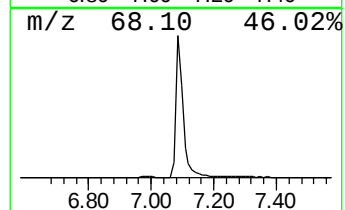
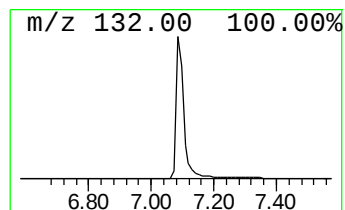
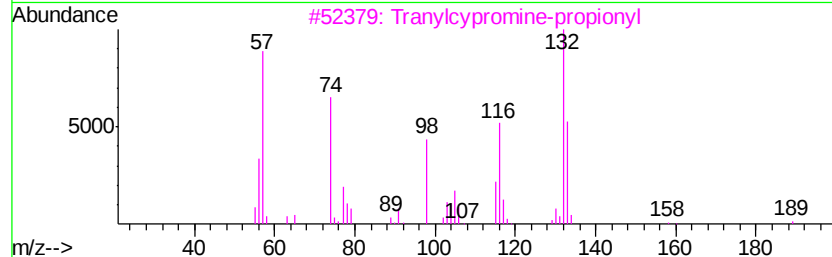
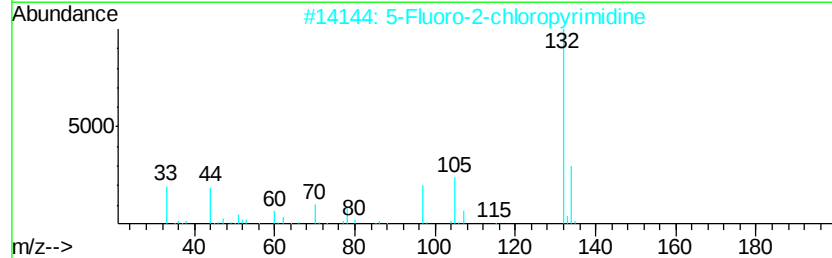
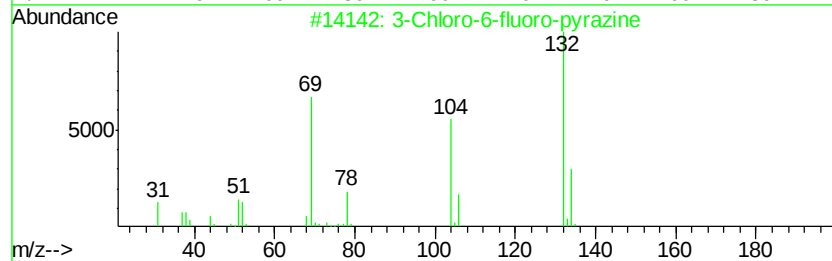
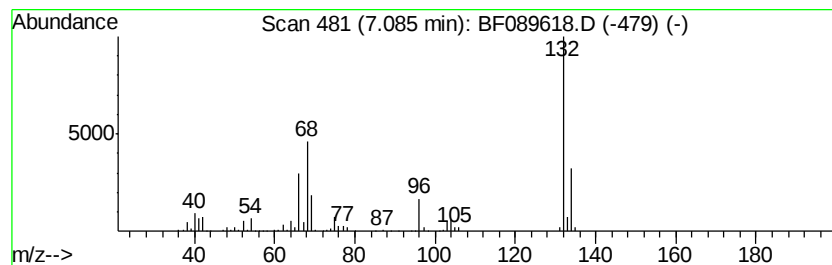
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Peak Number 3 unknown7.08 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	86.99 ng	1020610	1,4-Dichlorobenzene-d4	7.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.65	8.3	ng	97009	1	7.45	234653	20.0
unknown7.08	7.08	87.0	ng	1020610	1	7.45	234653	20.0