

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
 Data File : BF087510.D
 Acq On : 29 May 2016 8:52
 Operator : UM/SJ
 Sample : H3272-11
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB-201605

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-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	4.707	271	273	298	rBV	65252	268444	6.75%	1.031%
2	5.370	329	331	351	rBV	458434	1393795	35.07%	5.352%
3	5.976	382	384	392	rBV	17364	47398	1.19%	0.182%
4	7.027	475	476	482	rBV	313274	756306	19.03%	2.904%
5	7.119	482	484	498	rVB	1986414	3028820	76.20%	11.631%
6	7.473	513	515	527	rBV	387003	737903	18.57%	2.834%
7	7.702	532	535	551	rVB	1802438	2690376	67.69%	10.332%
8	8.387	592	595	627	rBV	1471123	2544628	64.02%	9.772%
9	9.496	690	692	710	rBV	533860	909712	22.89%	3.493%
10	11.279	844	848	850	rBV	2958977	3974692	100.00%	15.264%
11	12.319	937	939	958	rVB	608502	963348	24.24%	3.699%
12	13.622	1050	1053	1074	rBV	1495383	2471291	62.18%	9.490%
13	14.731	1147	1150	1163	rBV	597441	956804	24.07%	3.674%
14	17.085	1353	1356	1359	rBV	3368949	3970280	99.89%	15.247%
15	18.434	1472	1474	1486	rBV	421932	765170	19.25%	2.938%
16	20.114	1619	1621	1634	rBV	241567	561440	14.13%	2.156%

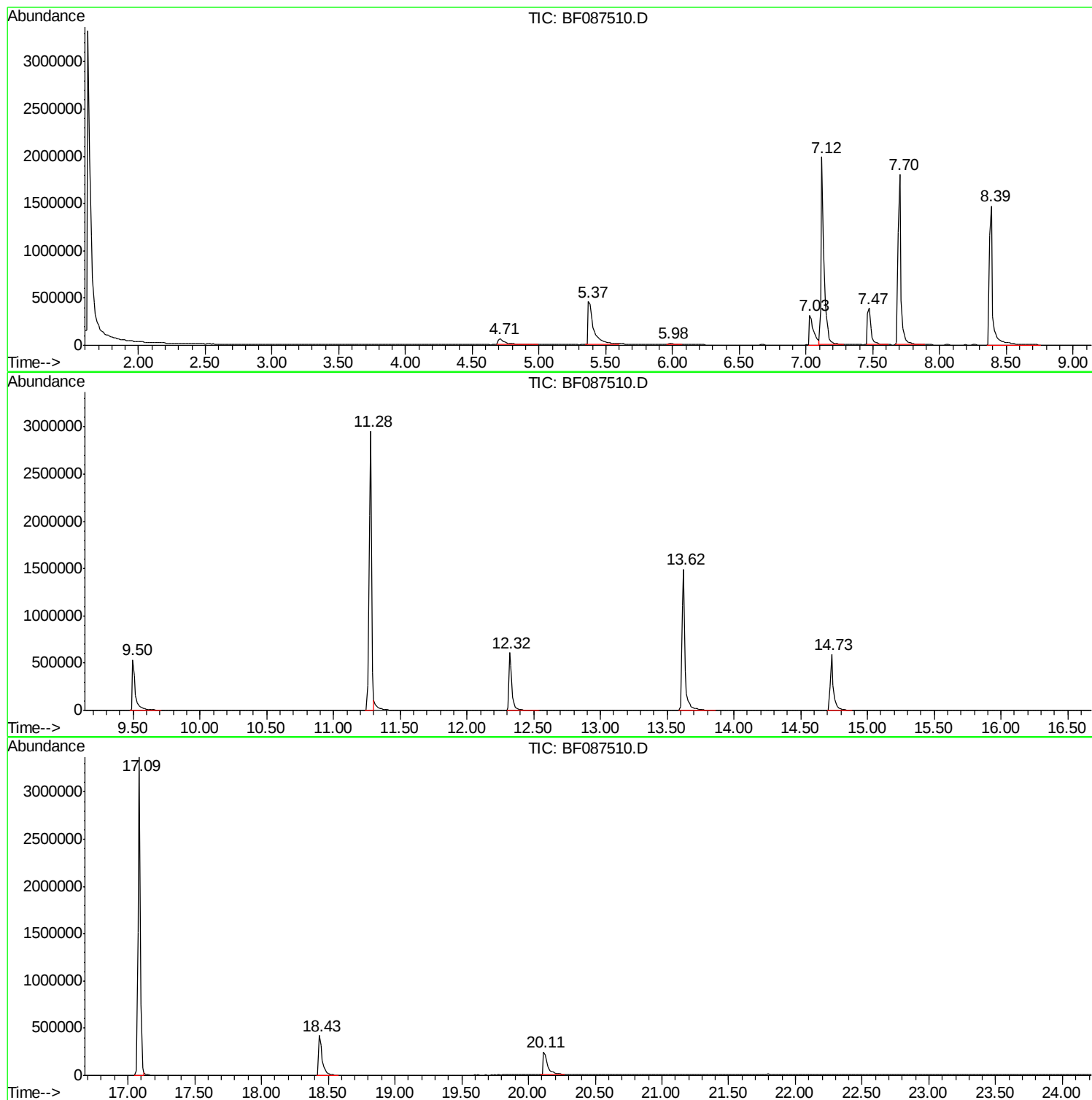
Sum of corrected areas: 26040407

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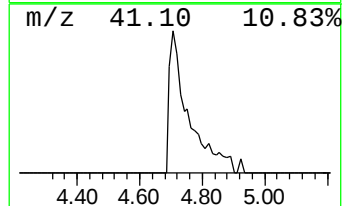
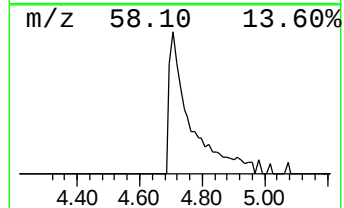
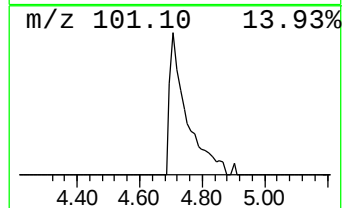
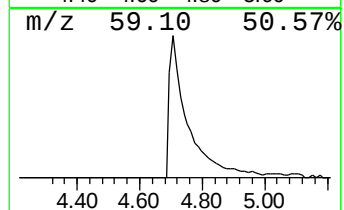
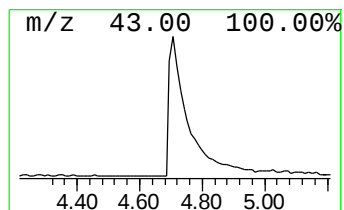
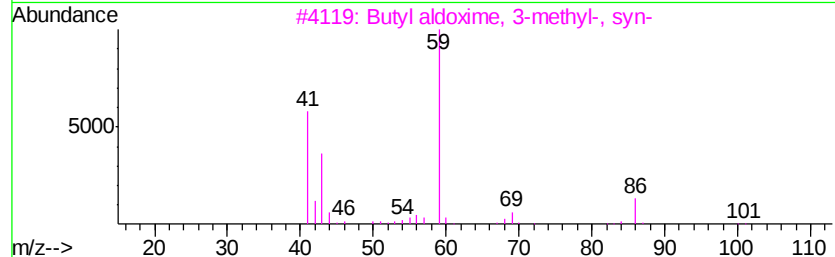
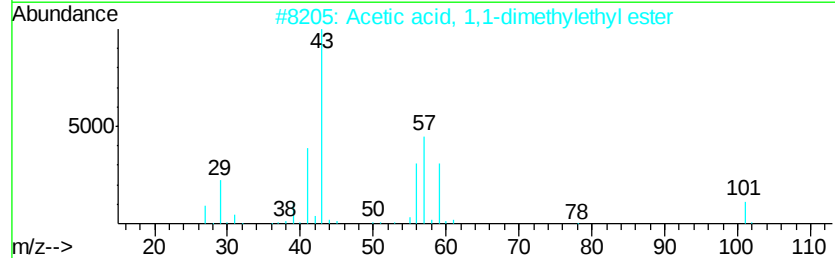
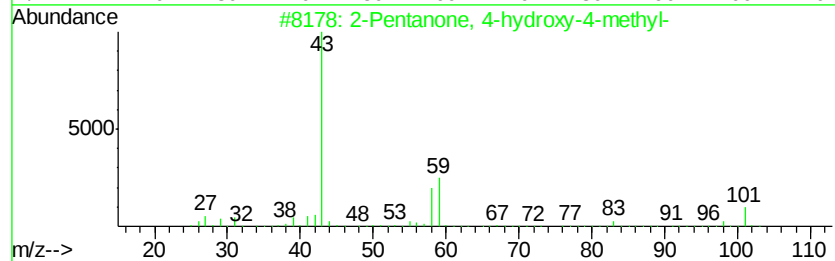
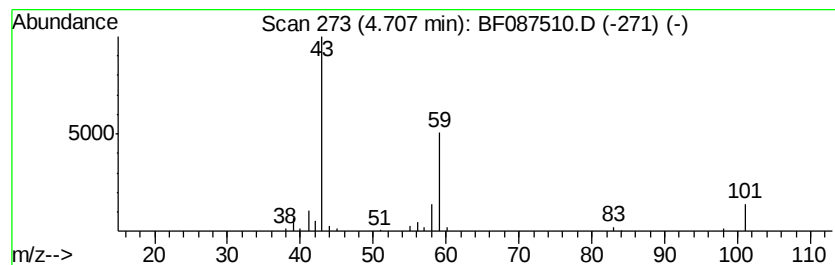
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.71	7.28 ng	268444	1,4-Dichlorobenzene-d4	7.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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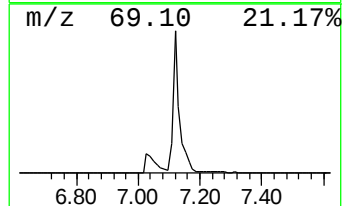
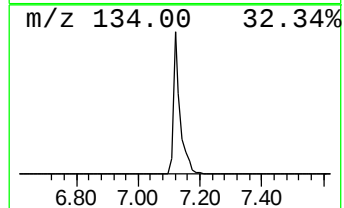
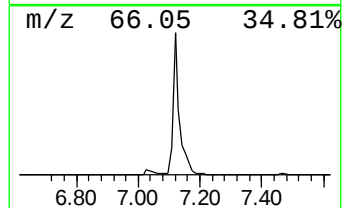
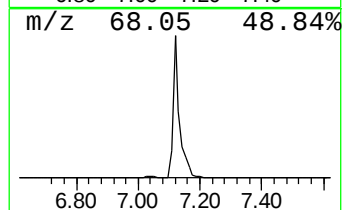
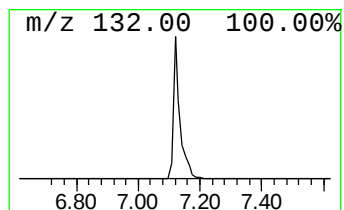
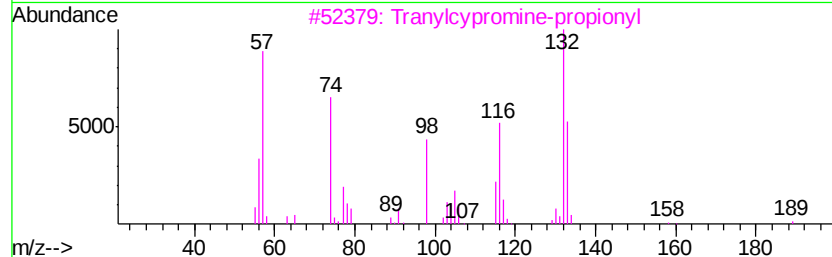
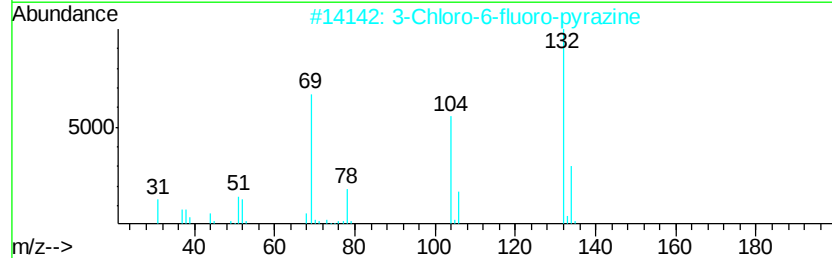
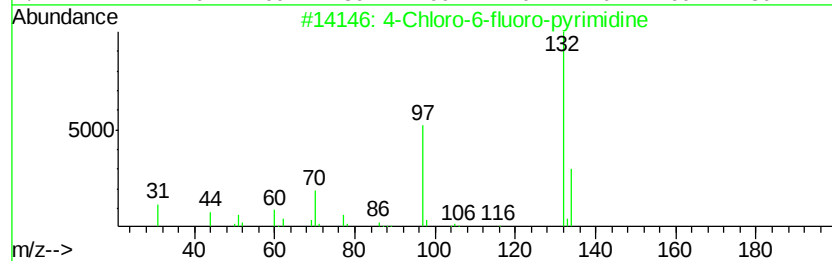
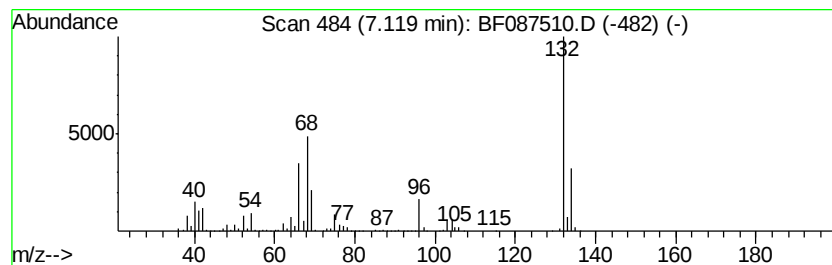
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Peak Number 2 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	82.09 ng	3028820	1,4-Dichlorobenzene-d4	7.47

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		Tranylcypromine-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...	4.71	7.3	ng	268444	1	7.47	737903	20.0
unknown7.12	7.12	82.1	ng	3028820	1	7.47	737903	20.0