

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060116\
 Data File : BF087635.D
 Acq On : 2 Jun 2016 1:28
 Operator : UM/SJ
 Sample : H3349-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-33

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	4.684	269	271	285	rBV	30226	133456	3.01%	0.460%
2	5.347	326	329	347	rBV	458781	1650925	37.27%	5.685%
3	7.005	472	474	479	rBV	368948	832578	18.80%	2.867%
4	7.085	479	481	500	rVB	1753586	3691387	83.34%	12.710%
5	7.416	508	510	521	rBV	467992	805828	18.19%	2.775%
6	7.656	528	531	546	rVB	1823892	3039822	68.63%	10.467%
7	8.342	589	591	615	rBV	1810847	2845360	64.24%	9.797%
8	9.451	686	688	704	rBV	564746	1006538	22.73%	3.466%
9	11.222	840	843	846	rBV	2731112	4429133	100.00%	15.251%
10	11.931	903	905	913	rBV	88981	181702	4.10%	0.626%
11	12.274	933	935	947	rVB	793465	1066201	24.07%	3.671%
12	13.108	1003	1008	1010	rVB2	48164	64247	1.45%	0.221%
13	13.577	1046	1049	1064	rBV	1766384	2809424	63.43%	9.673%
14	13.874	1073	1075	1081	rBV	52681	85554	1.93%	0.295%
15	14.605	1136	1139	1141	rBV	33565	45218	1.02%	0.156%
16	14.685	1144	1146	1159	rVB	565051	1015742	22.93%	3.497%
17	15.188	1188	1190	1196	rBV	27256	56920	1.29%	0.196%
18	17.040	1349	1352	1355	rBV	3210867	3879741	87.60%	13.359%
19	18.286	1458	1461	1466	rBV2	24880	51175	1.16%	0.176%
20	18.400	1470	1471	1487	rVV	382529	791995	17.88%	2.727%
21	20.092	1617	1619	1627	rBV	242908	559587	12.63%	1.927%

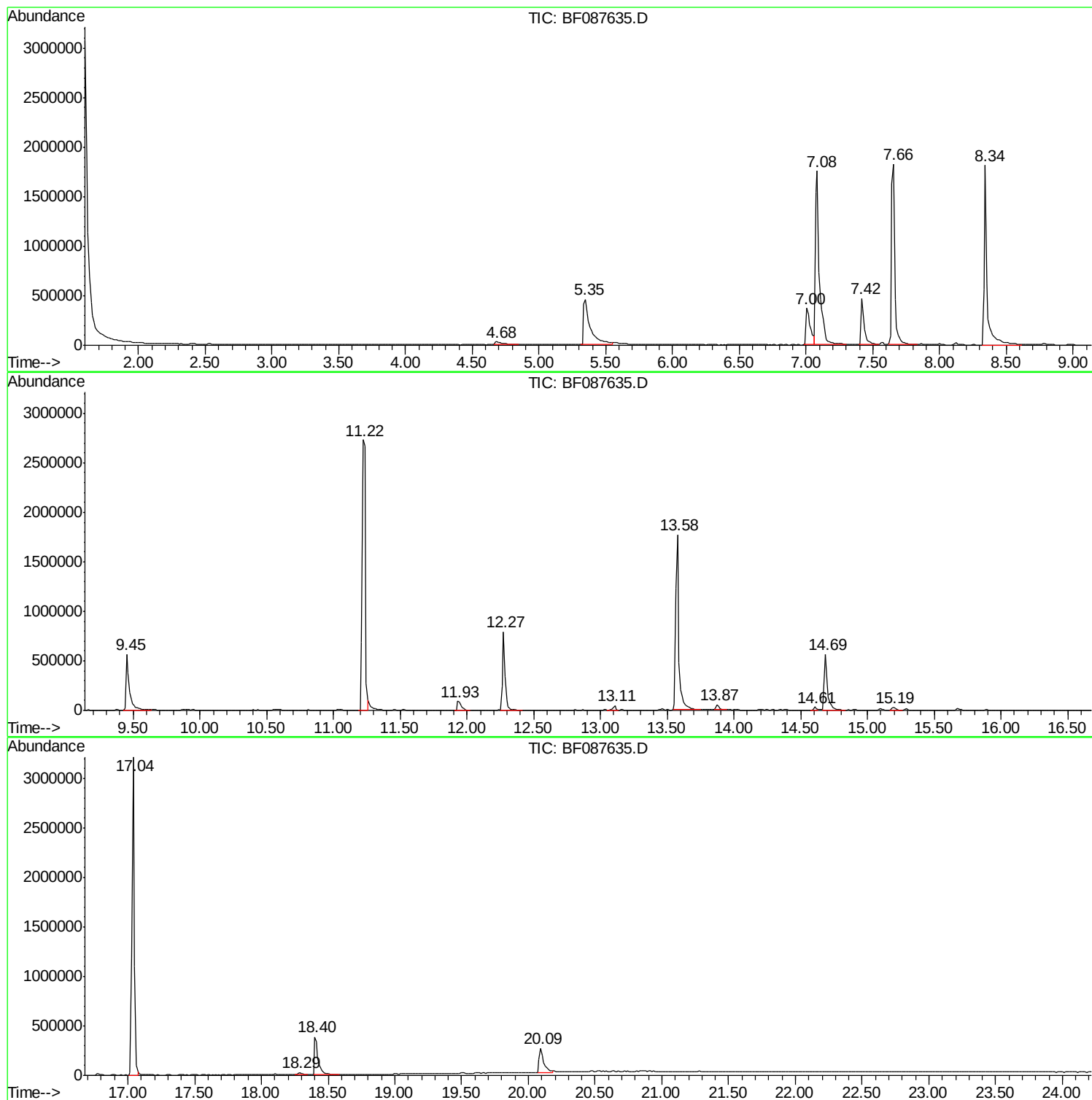
Sum of corrected areas: 29042533

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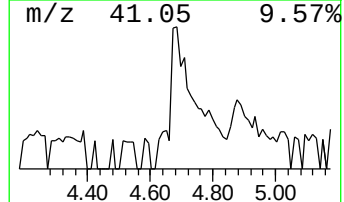
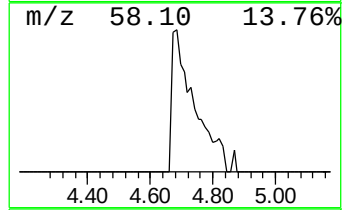
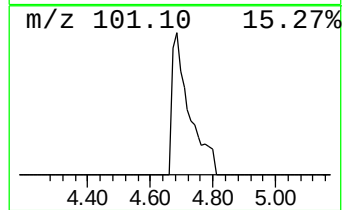
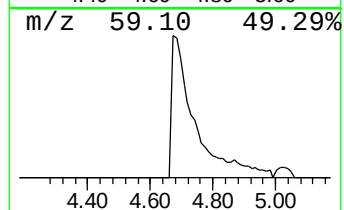
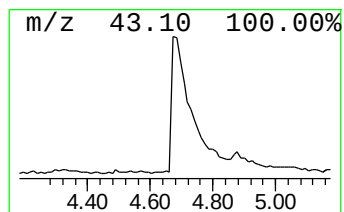
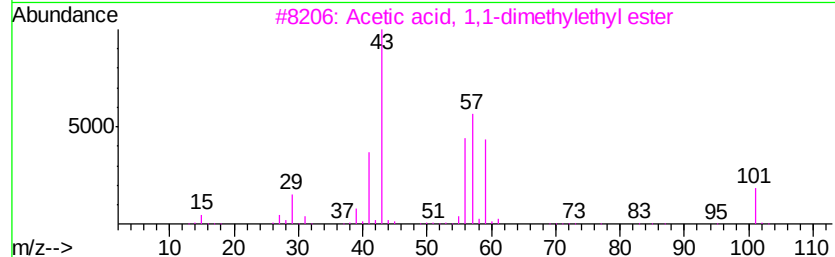
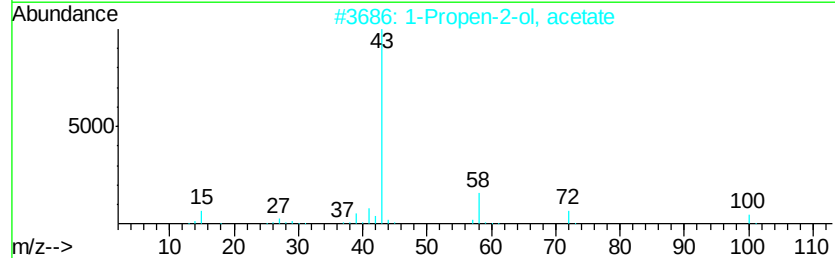
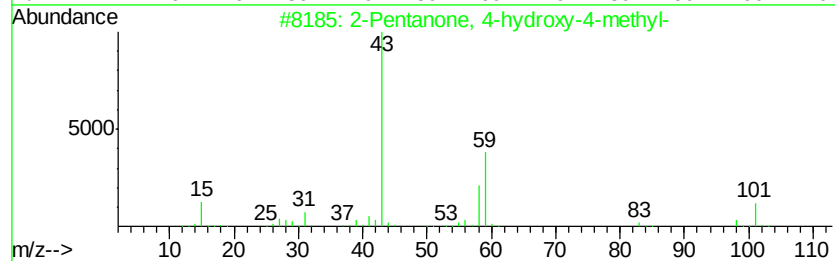
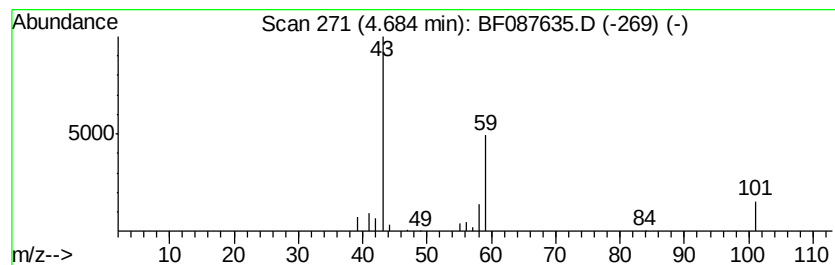
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.68	3.31 ng	133456	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	72	
2	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9	
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9	
4	Butane	58	C4H10	000106-97-8	7	
5	Diacetamide	101	C4H7NO2	000625-77-4	7	



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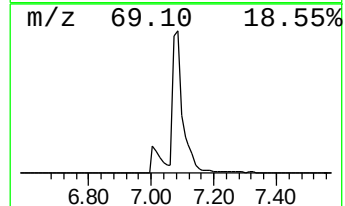
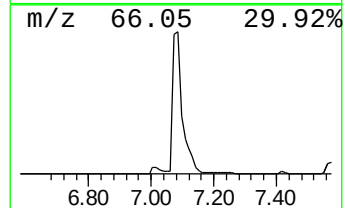
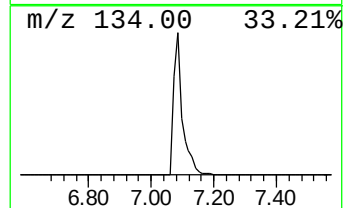
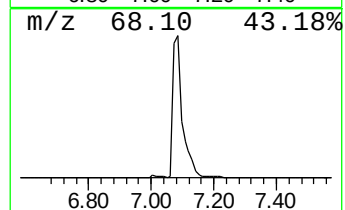
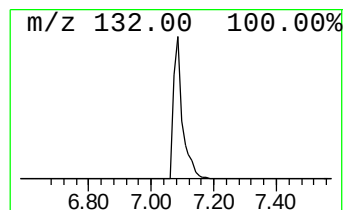
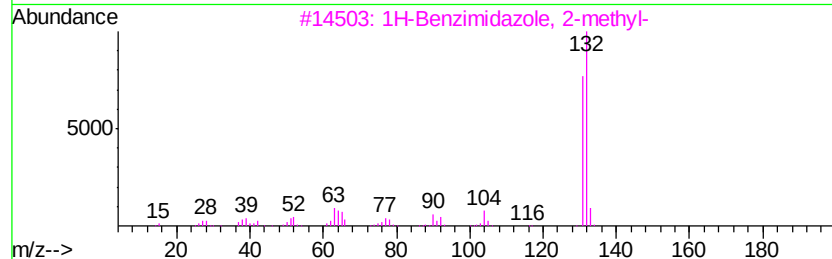
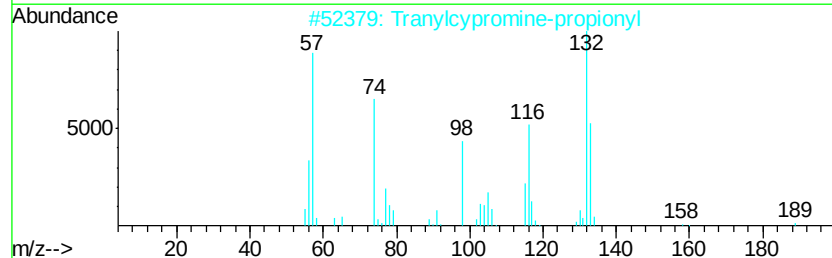
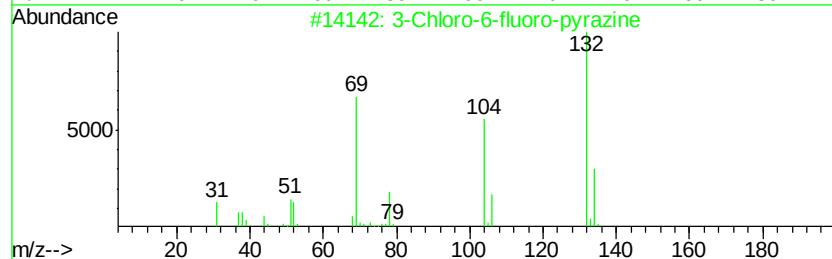
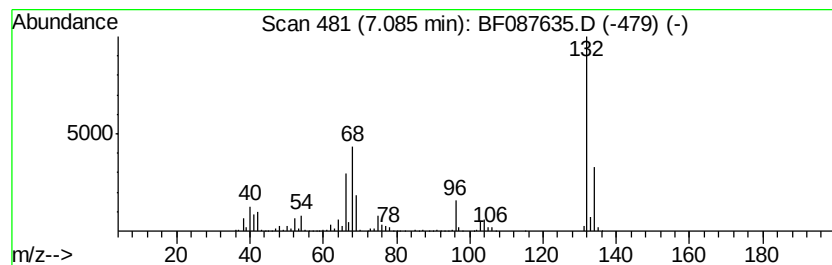
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Peak Number 2 unknown7.08 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	91.62 ng	3691390	1,4-Dichlorobenzene-d4	7.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.68	3.3	ng	133456	1	7.42	805828	20.0
unknown7.08	7.08	91.6	ng	3691390	1	7.42	805828	20.0