

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
 Data File : BF089395.D
 Acq On : 29 Jul 2016 6:55
 Operator : UM/SJ
 Sample : PB92528BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92528BL

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-BF072716.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	1.678	6	8	52	rVB	705034	1604534	100.00%	12.709%
2	4.661	266	269	279	rVB	35955	75743	4.72%	0.600%
3	5.107	305	308	322	rBV	1150998	1353399	84.35%	10.720%
4	6.193	400	403	405	rBV	796902	1257583	78.38%	9.961%
5	6.296	410	412	414	rBV	1407267	1454371	90.64%	11.519%
6	6.524	430	432	437	rBV	246882	255607	15.93%	2.025%
7	6.673	443	445	448	rBV	745690	1081559	67.41%	8.566%
8	7.096	480	482	500	rBV	632876	836445	52.13%	6.625%
9	7.816	542	545	547	rBV	218556	323041	20.13%	2.559%
10	8.890	636	639	642	rBV	1541259	1474572	91.90%	11.679%
11	9.553	695	697	700	rBV	405122	361420	22.52%	2.863%
12	10.342	764	766	778	rBB	692347	699130	43.57%	5.537%
13	11.028	824	826	839	rBV	302325	297115	18.52%	2.353%
14	12.605	962	964	967	rBV	749684	1079828	67.30%	8.553%
15	13.645	1053	1055	1069	rBV	262222	258960	16.14%	2.051%
16	14.982	1169	1172	1175	rBV	176875	212186	13.22%	1.681%

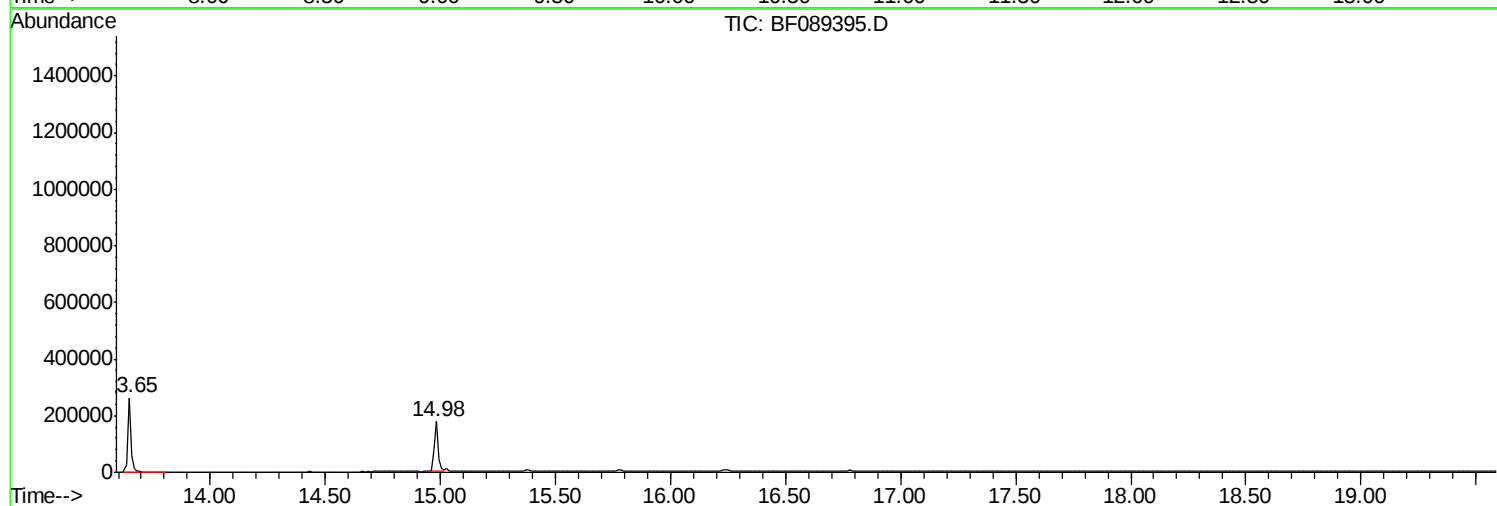
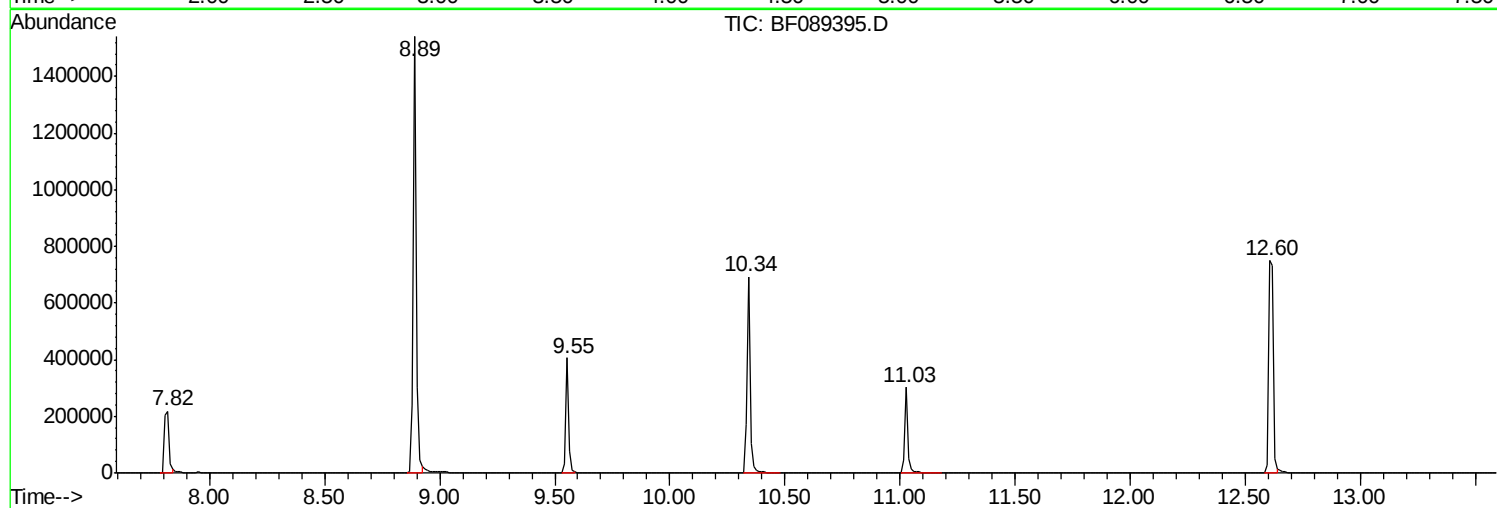
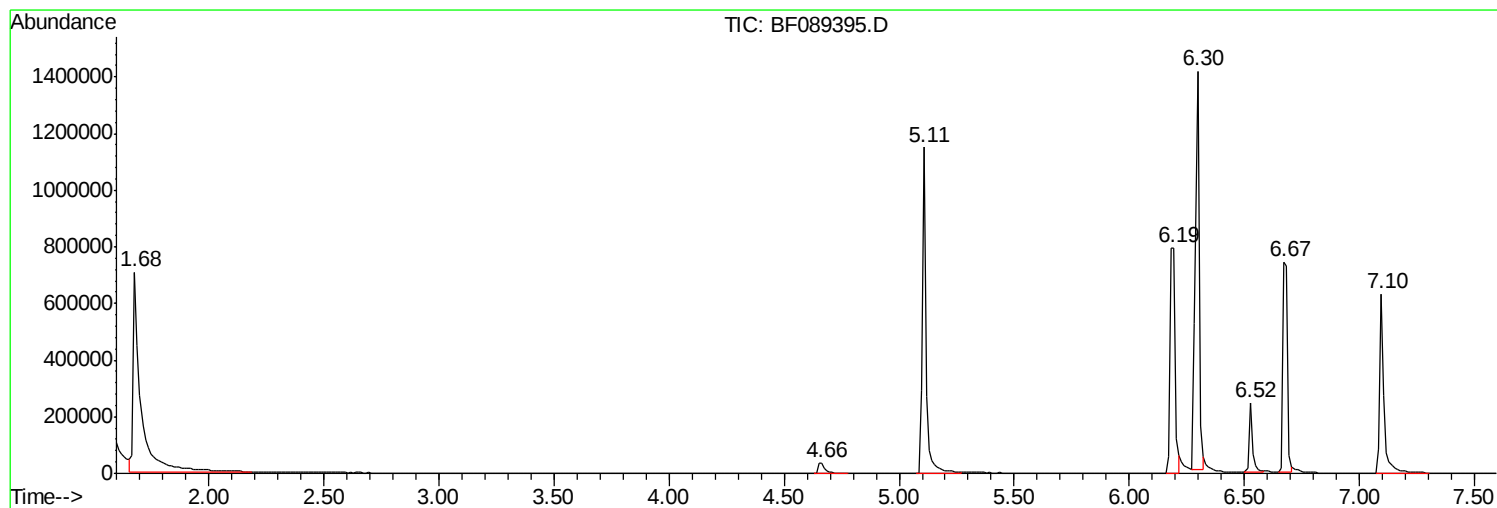
Sum of corrected areas: 12625493

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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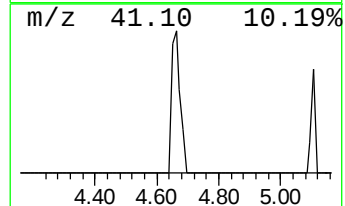
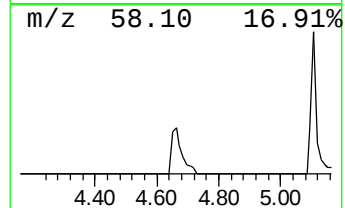
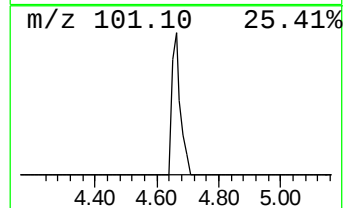
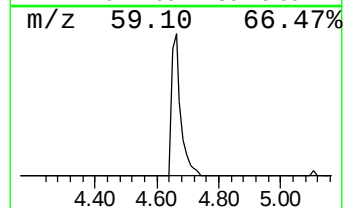
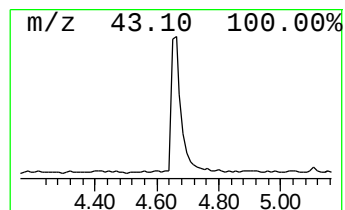
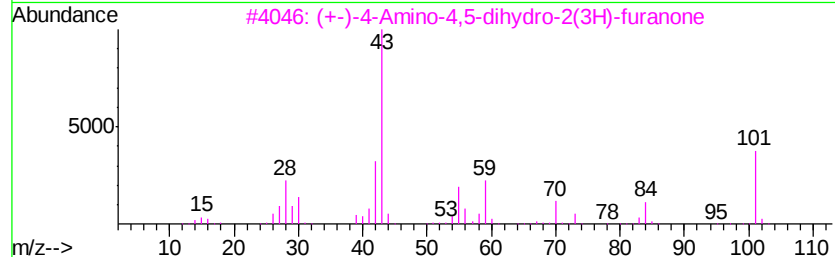
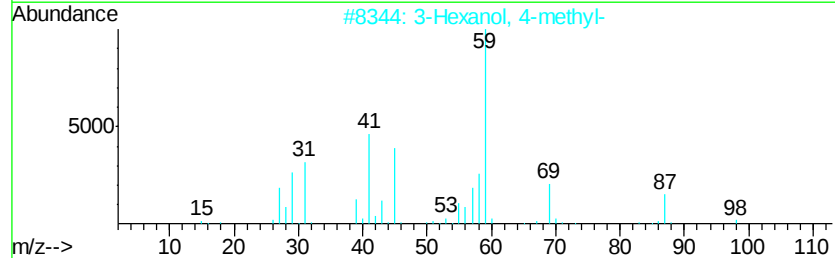
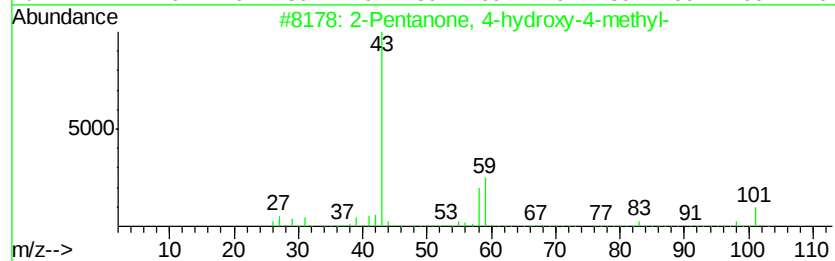
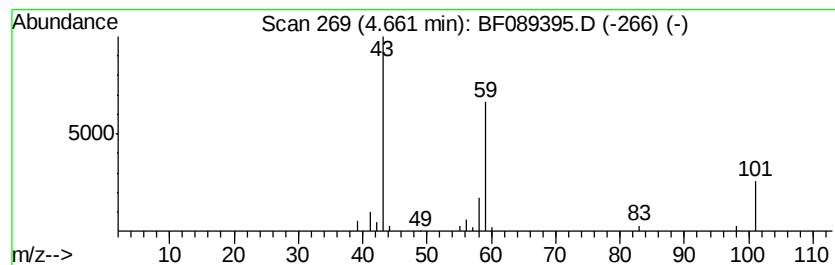
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	5.93 ng	75743	1,4-Dichlorobenzene-d4	6.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	28
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			3-Octanol	130	C8H18O	000589-98-0	9



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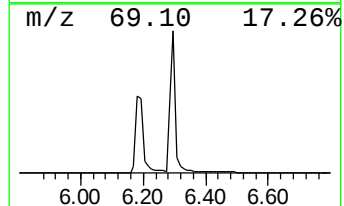
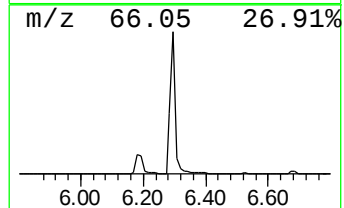
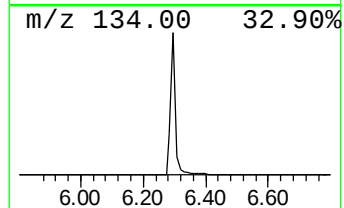
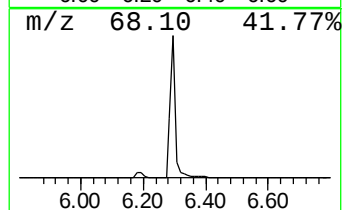
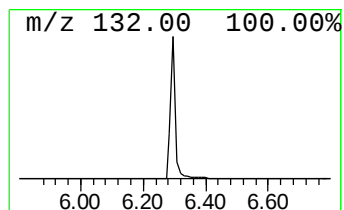
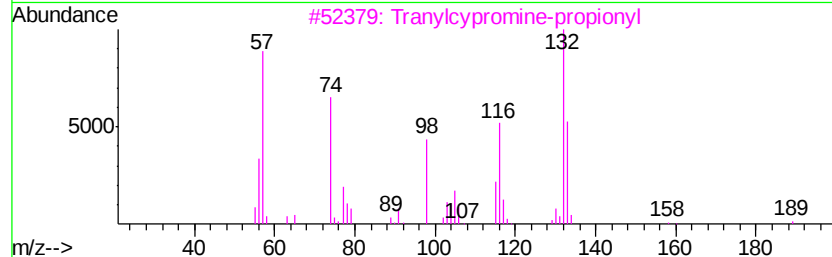
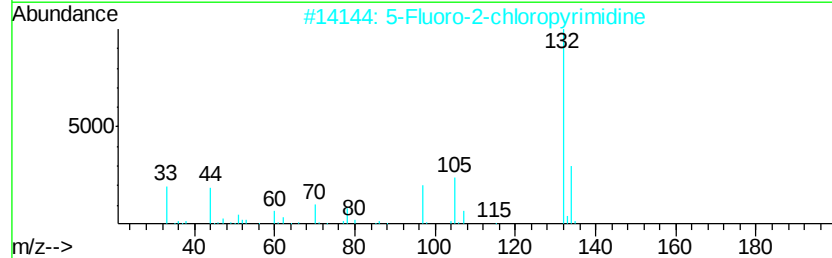
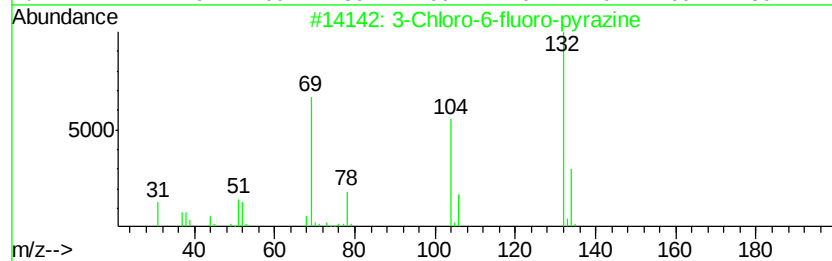
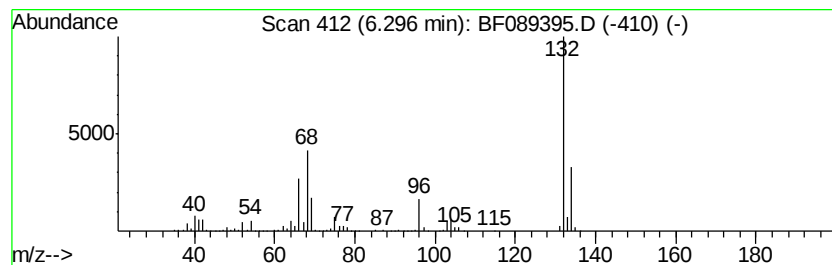
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Peak Number 3 unknown6.30 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.30	113.80 ng	1454370	1,4-Dichlorobenzene-d4	6.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.66	5.9	ng	75743	1	6.52	255607	20.0
unknown6.30	6.30	113.8	ng	1454370	1	6.52	255607	20.0