

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072116\
 Data File : BF089166.D
 Acq On : 21 Jul 2016 8:17
 Operator : UM/SJ
 Sample : H4089-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOCATION-33A-9-11

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-BF072016.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	1.724	10	12	50	rVB	1270619	2034846	100.00%	13.131%
2	4.741	273	276	289	rBV	30098	66073	3.25%	0.426%
3	5.187	313	315	341	rBV	768020	1088378	53.49%	7.023%
4	6.261	406	409	416	rBV	744742	1172819	57.64%	7.568%
5	6.364	416	418	437	rVV	894389	1251620	61.51%	8.077%
6	6.604	437	439	449	rVV	190349	308837	15.18%	1.993%
7	6.753	449	452	468	rVB	940975	956920	47.03%	6.175%
8	7.164	486	488	504	rBV	506943	749500	36.83%	4.836%
9	7.885	549	551	562	rBV	403651	430369	21.15%	2.777%
10	8.959	642	645	648	rBV	1284111	1414950	69.54%	9.130%
11	9.359	678	680	692	rBB	253243	239577	11.77%	1.546%
12	9.622	701	703	706	rBV	327779	483744	23.77%	3.122%
13	10.410	770	772	775	rBV2	688292	801926	39.41%	5.175%
14	10.548	781	784	790	rVB	18568	23310	1.15%	0.150%
15	11.096	830	832	835	rBV	422738	510071	25.07%	3.291%
16	12.525	955	957	959	rBB	59354	54650	2.69%	0.353%
17	12.685	968	971	973	rBV	1534008	1501728	73.80%	9.690%
18	13.222	1016	1018	1020	rBV	45804	39031	1.92%	0.252%
19	13.634	1046	1054	1058	rBV4	18668	70792	3.48%	0.457%
20	13.714	1058	1061	1064	rBV	1144263	1965866	96.61%	12.685%
21	15.062	1176	1179	1182	rBV	294487	331971	16.31%	2.142%

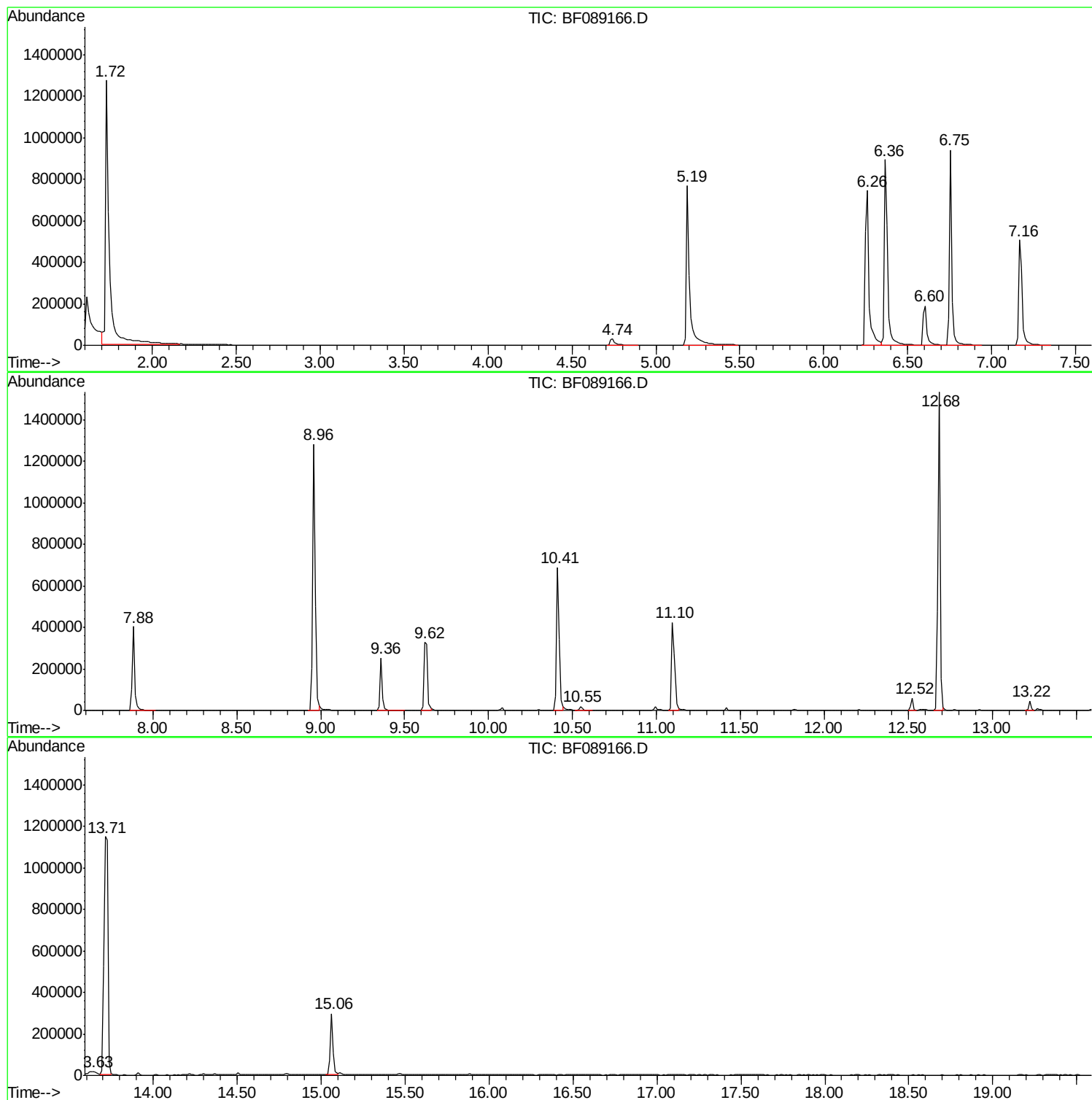
Sum of corrected areas: 15496978

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072016.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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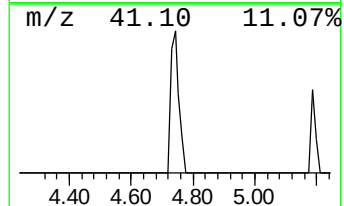
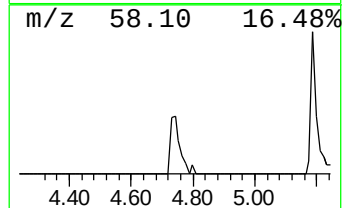
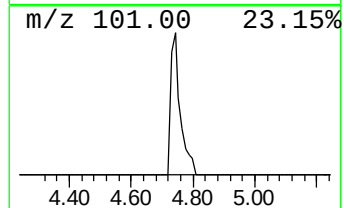
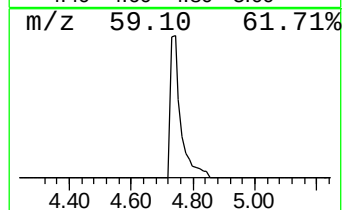
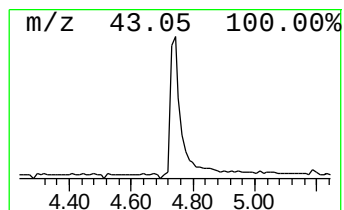
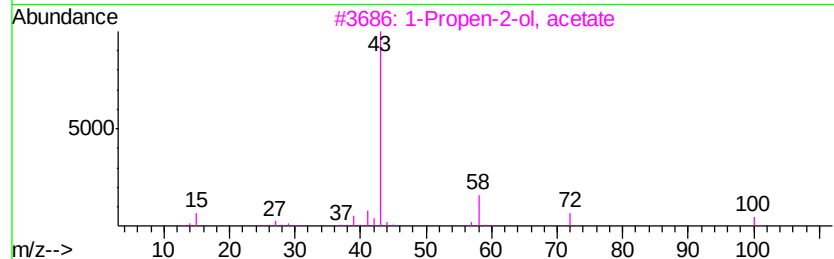
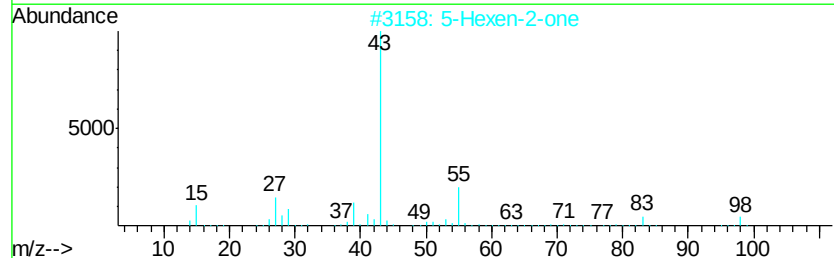
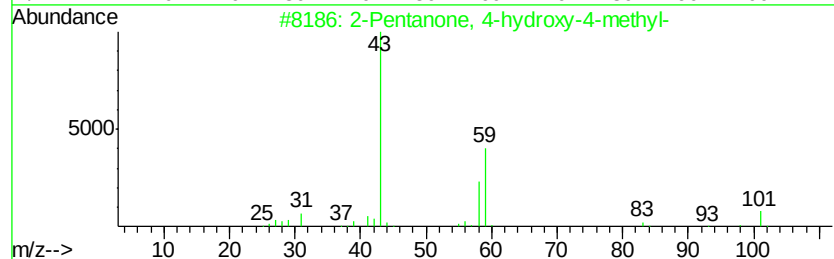
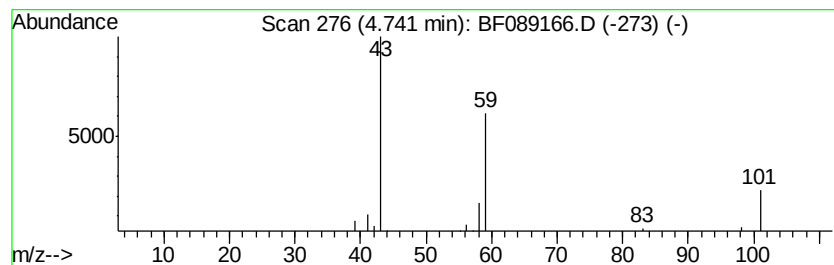
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	4.28 ng	66073	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		Acetone	58	C3H6O	000067-64-1	7
5		Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	5



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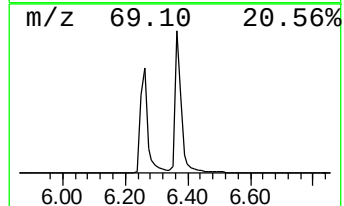
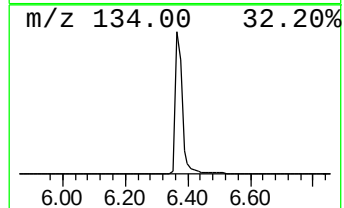
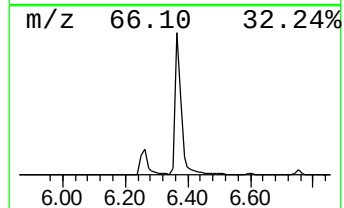
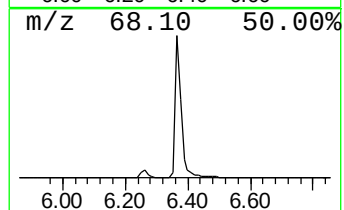
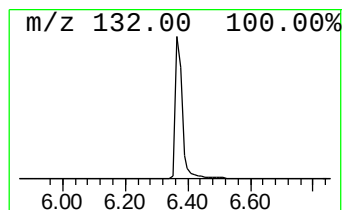
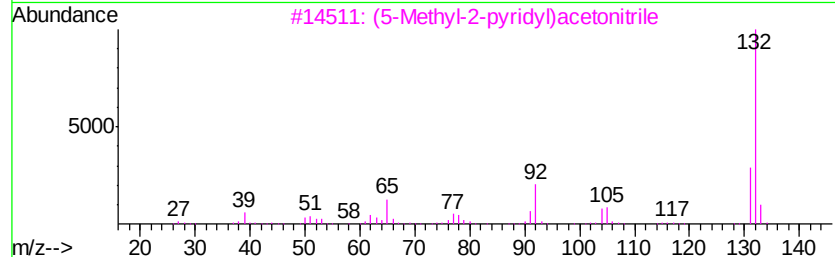
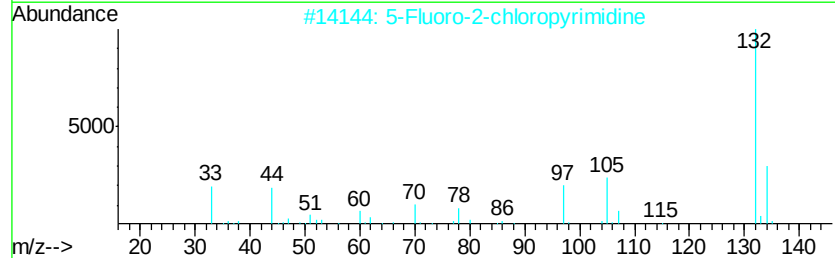
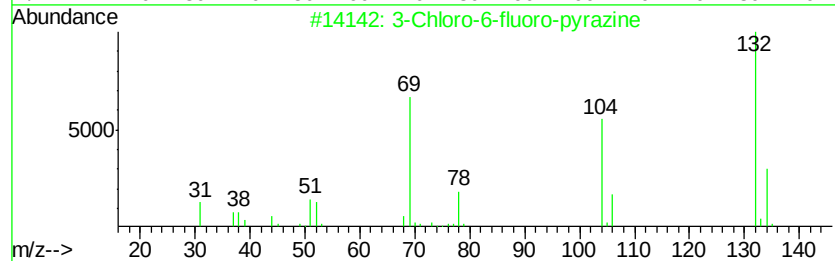
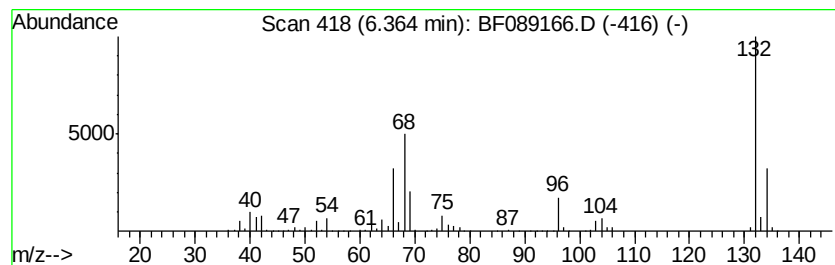
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Peak Number 3 unknown6.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	81.05 ng	1251620	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Tranvlcyproamine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.74	4.3	ng	66073	1	6.60	308837	20.0
unknown6.36	6.36	81.0	ng	1251620	1	6.60	308837	20.0