

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072017\
 Data File : BF096908.D
 Acq On : 20 Jul 2017 16:22
 Operator : SJ/JU
 Sample : PB100793BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100793BL

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-BF071317.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	4.746	464	469	485	rBV	233703	359395	11.45%	1.473%
2	5.181	538	543	546	rBV	2592338	2750206	87.59%	11.269%
3	6.222	715	720	723	rBV	2455056	2819782	89.80%	11.554%
4	6.340	735	740	743	rBV	2602878	2725693	86.81%	11.169%
5	6.569	775	779	785	rBV	622307	544012	17.33%	2.229%
6	6.728	801	806	809	rBV	2234113	2216189	70.58%	9.081%
7	7.134	869	875	878	rBV	1864473	1891241	60.23%	7.749%
8	7.851	992	997	1000	rBV	857511	823589	26.23%	3.375%
9	8.928	1175	1180	1184	rBV	2760301	3139942	100.00%	12.866%
10	9.598	1289	1294	1298	rBV	807962	753508	24.00%	3.088%
11	10.386	1422	1428	1431	rBV	1708994	1745302	55.58%	7.151%
12	11.075	1540	1545	1548	rBV	909952	827624	26.36%	3.391%
13	12.663	1810	1815	1818	rBV2	2488878	2725199	86.79%	11.167%
14	13.698	1987	1991	1995	rBV	739371	664650	21.17%	2.723%
15	15.063	2219	2223	2229	rVB	353583	418391	13.32%	1.714%

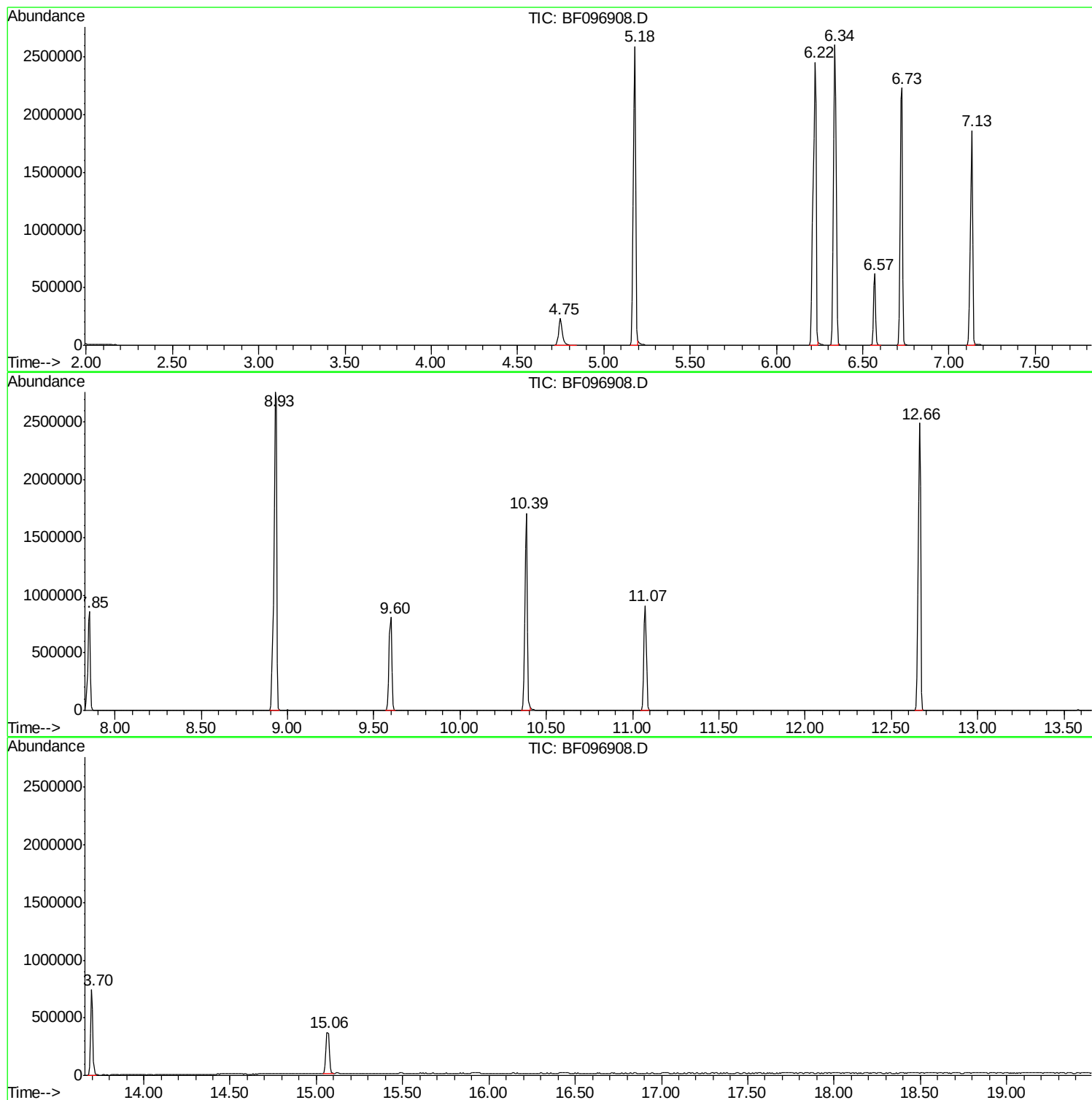
Sum of corrected areas: 24404723

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.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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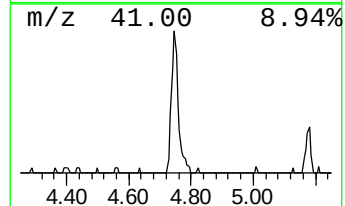
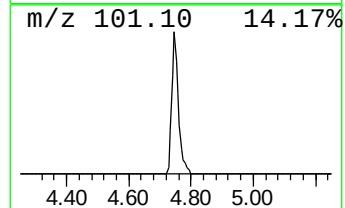
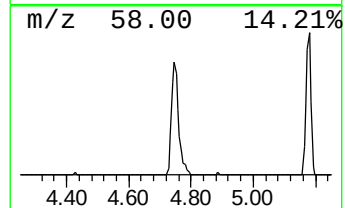
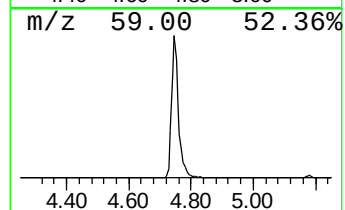
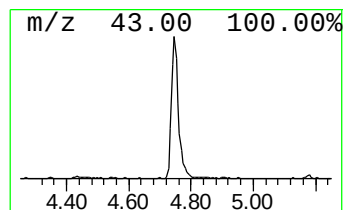
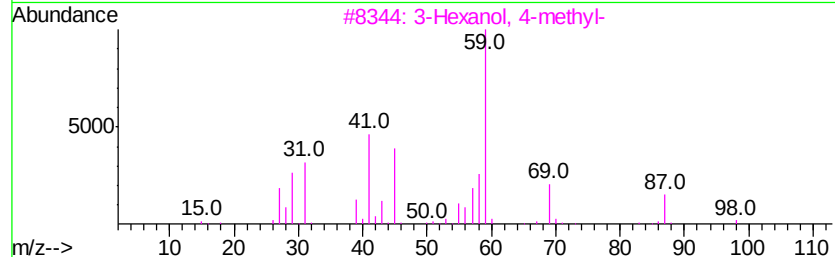
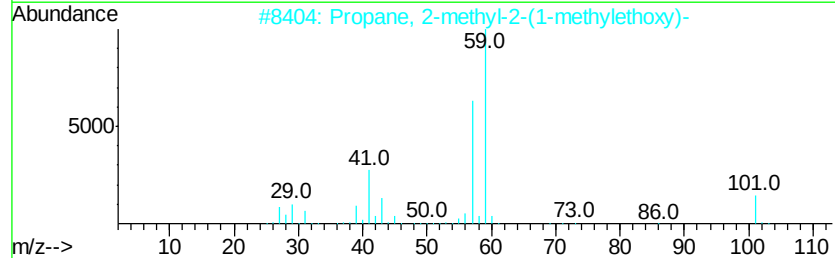
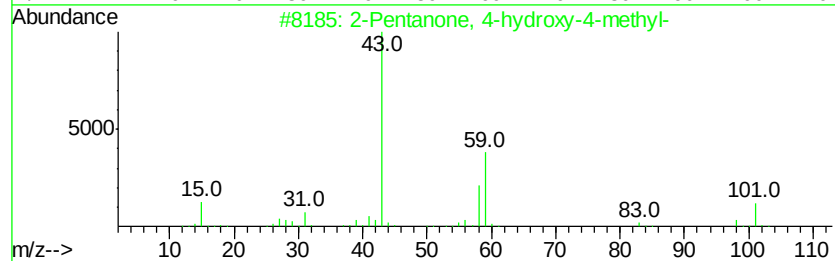
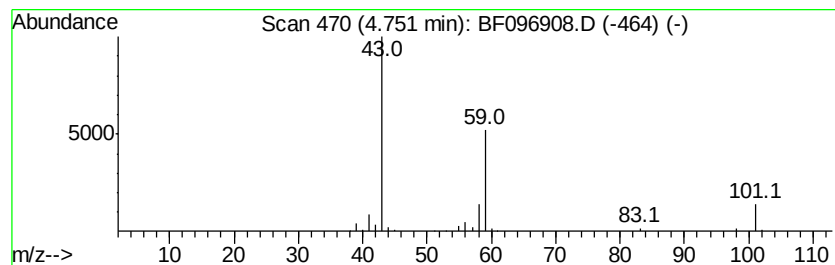
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	13.21 ng	359395	1,4-Dichlorobenzene-d4	6.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		3-Hexanol	102	C6H14O	000623-37-0	33
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33



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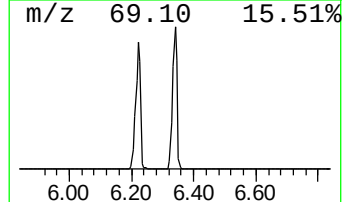
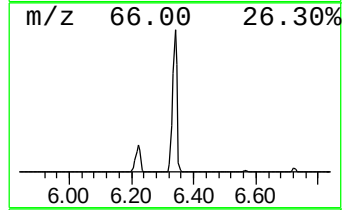
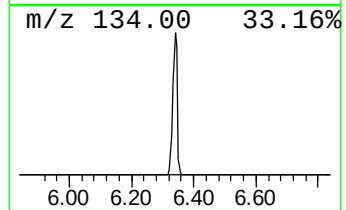
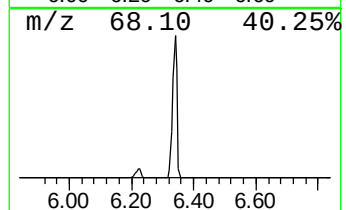
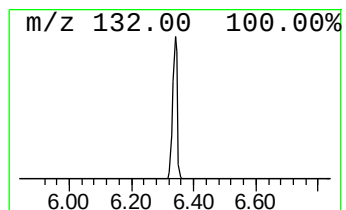
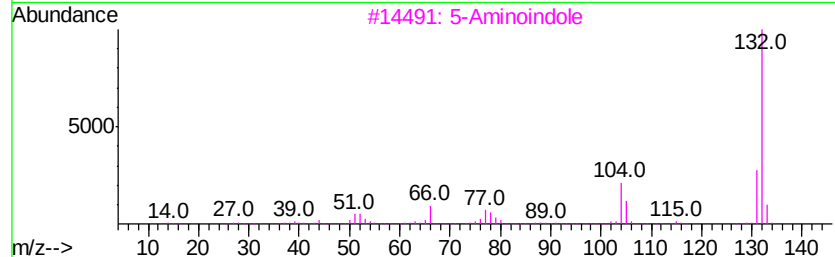
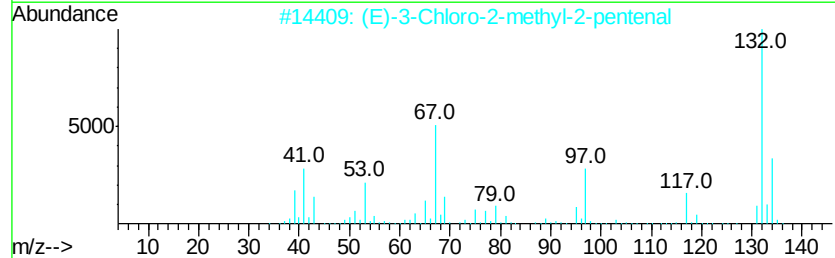
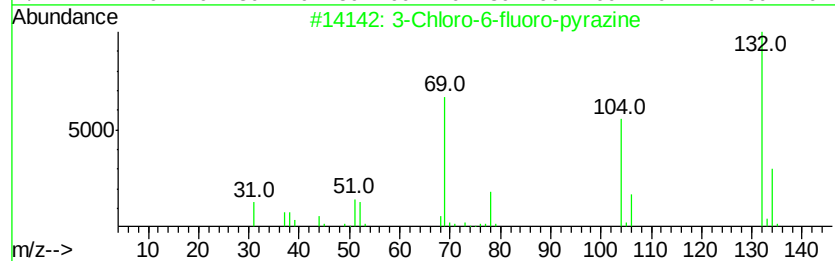
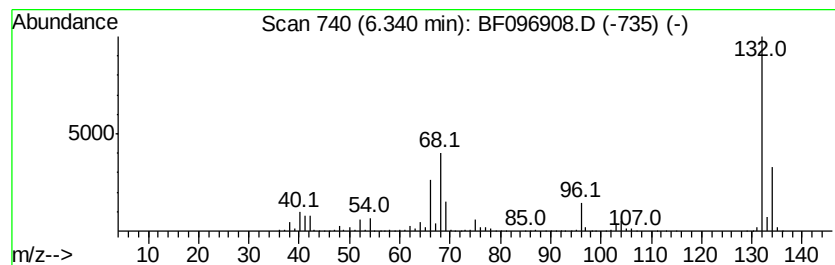
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Peak Number 2 unknown6.34 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	100.21 ng	2725690	1,4-Dichlorobenzene-d4	6.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.75	13.2	ng	359395	1	6.57	544012	20.0
unknown6.34	6.34	100.2	ng	2725690	1	6.57	544012	20.0