

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071017\
 Data File : BF096606.D
 Acq On : 10 Jul 2017 17:39
 Operator : SJ/JU
 Sample : PB100517BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100517BL

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-BF070117.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	4.881	471	475	487	rBV	430021	636133	20.12%	2.492%
2	5.304	541	547	550	rBV	2872333	3006401	95.08%	11.780%
3	6.328	715	721	724	rBV	2662473	2922196	92.41%	11.450%
4	6.457	738	743	746	rBV	2817164	2917276	92.26%	11.430%
5	6.681	777	781	785	rBV	770926	632001	19.99%	2.476%
6	6.840	804	808	811	rBV	2441927	2228359	70.47%	8.731%
7	7.245	871	877	880	rBV	1805116	1882411	59.53%	7.376%
8	7.963	994	999	1003	rBV	957852	870466	27.53%	3.411%
9	9.039	1176	1182	1186	rBV	2730508	3162085	100.00%	12.390%
10	9.710	1291	1296	1300	rBV	871887	908784	28.74%	3.561%
11	10.498	1424	1430	1434	rBV	1623089	1783727	56.41%	6.989%
12	11.192	1543	1548	1551	rBV	917341	854723	27.03%	3.349%
13	12.780	1812	1818	1821	rBV	2532014	2566482	81.16%	10.056%
14	13.816	1990	1994	1998	rBV	756780	656434	20.76%	2.572%
15	15.210	2226	2231	2236	rBV	432158	494645	15.64%	1.938%

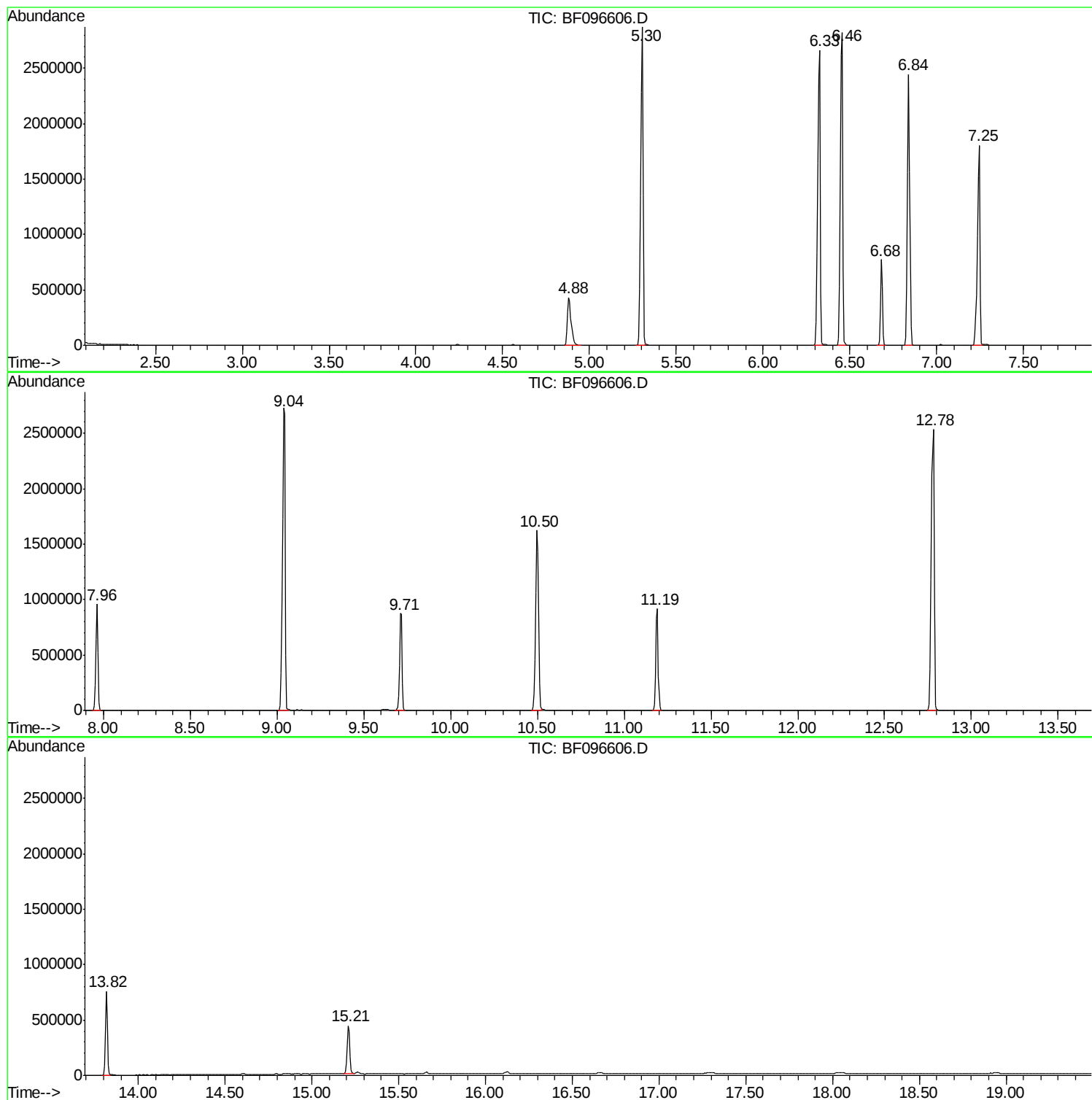
Sum of corrected areas: 25522123

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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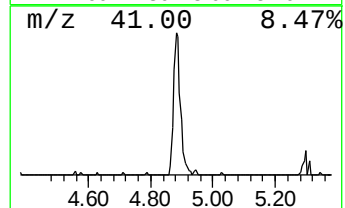
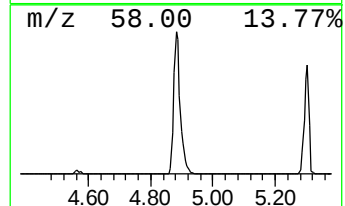
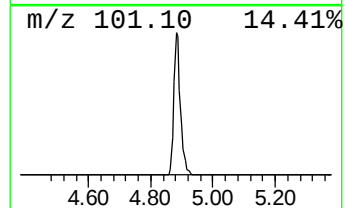
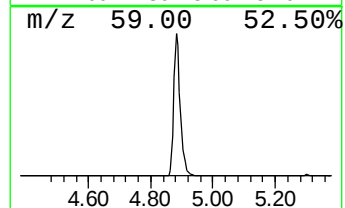
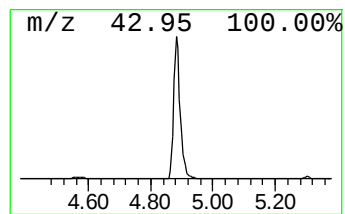
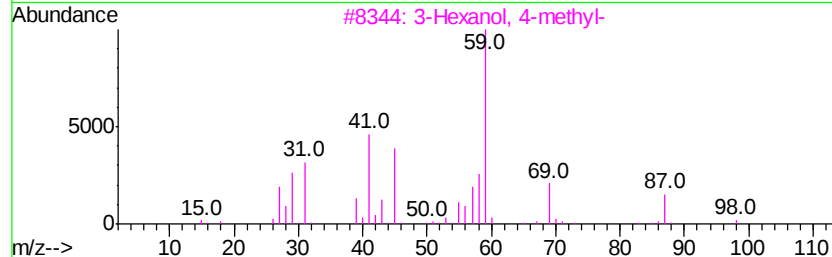
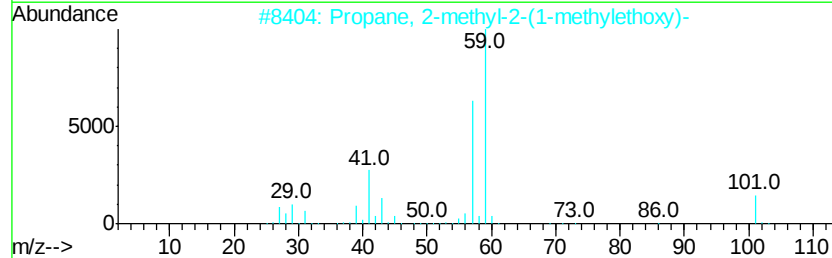
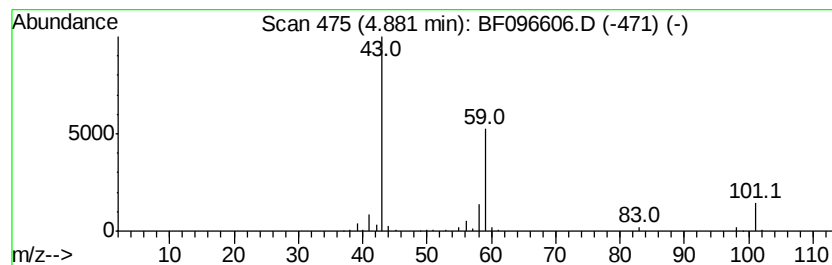
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	20.13 ng	636133	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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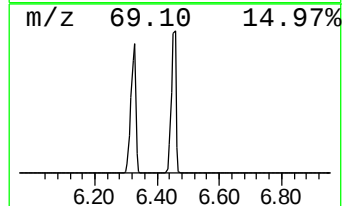
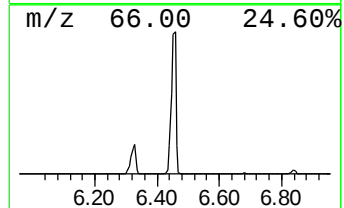
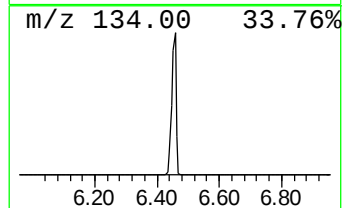
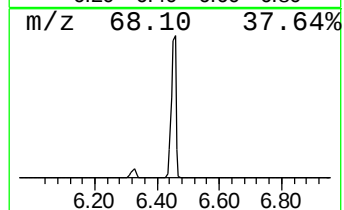
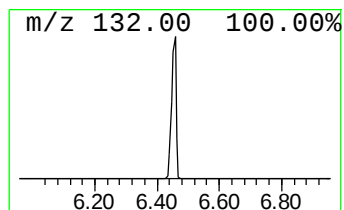
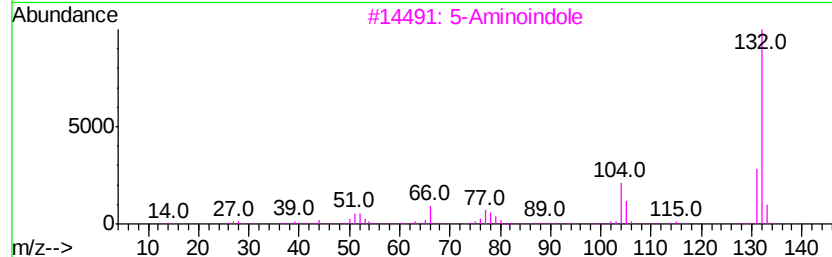
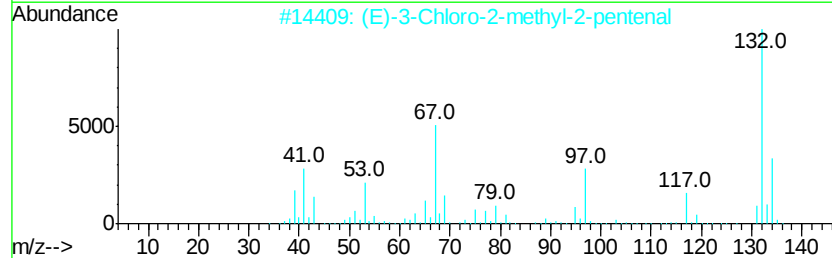
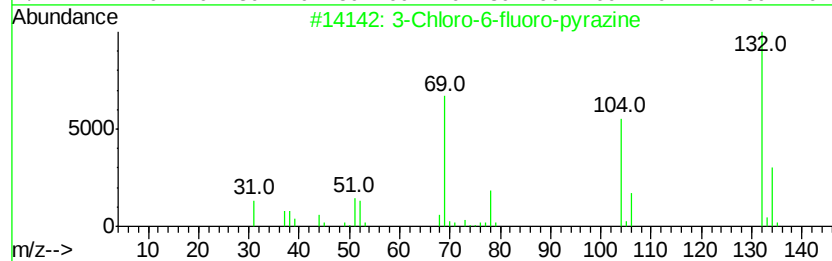
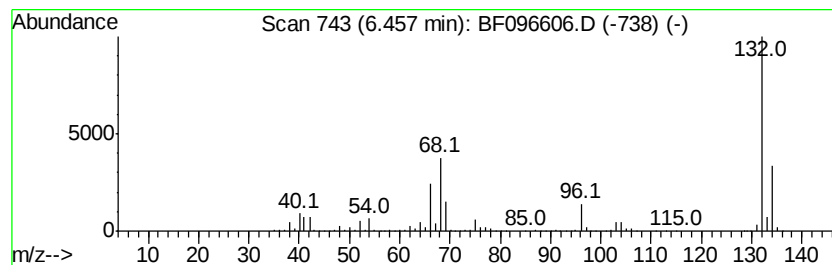
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Peak Number 2 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	92.32 ng	2917280	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
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	12
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.88	20.1	ng	636133	1	6.68	632001	20.0
unknown6.46	6.46	92.3	ng	2917280	1	6.68	632001	20.0