

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
 Data File : BF097393.D
 Acq On : 5 Aug 2017 13:21
 Operator : SJ/JU
 Sample : PB101266BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101266BL

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-BF072517.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	4.575	470	474	495	rBV	163387	309533	10.78%	1.279%
2	5.034	548	552	569	rBV	2254880	2872525	100.00%	11.867%
3	6.098	729	733	746	rBV	2435342	2827075	98.42%	11.679%
4	6.198	746	750	760	rVB	2689873	2668378	92.89%	11.023%
5	6.422	785	788	797	rBV	666528	604370	21.04%	2.497%
6	6.581	810	815	832	rVB	2003791	2132692	74.24%	8.810%
7	6.992	881	885	900	rBV	1555907	1772349	61.70%	7.322%
8	7.704	1002	1006	1020	rBV	892159	882974	30.74%	3.648%
9	8.786	1184	1190	1193	rBV	2466697	2848167	99.15%	11.766%
10	9.445	1297	1302	1310	rVB	896296	928858	32.34%	3.837%
11	10.239	1432	1437	1449	rBV	1534294	1634297	56.89%	6.751%
12	10.921	1549	1553	1560	rVB	905308	849832	29.58%	3.511%
13	12.504	1817	1822	1825	rBV	2228216	2449043	85.26%	10.117%
14	13.539	1994	1998	2006	rBV	779233	801176	27.89%	3.310%
15	14.880	2221	2226	2230	rBV	525364	593347	20.66%	2.451%
16	15.262	2287	2291	2295	rBV3	25821	31994	1.11%	0.132%

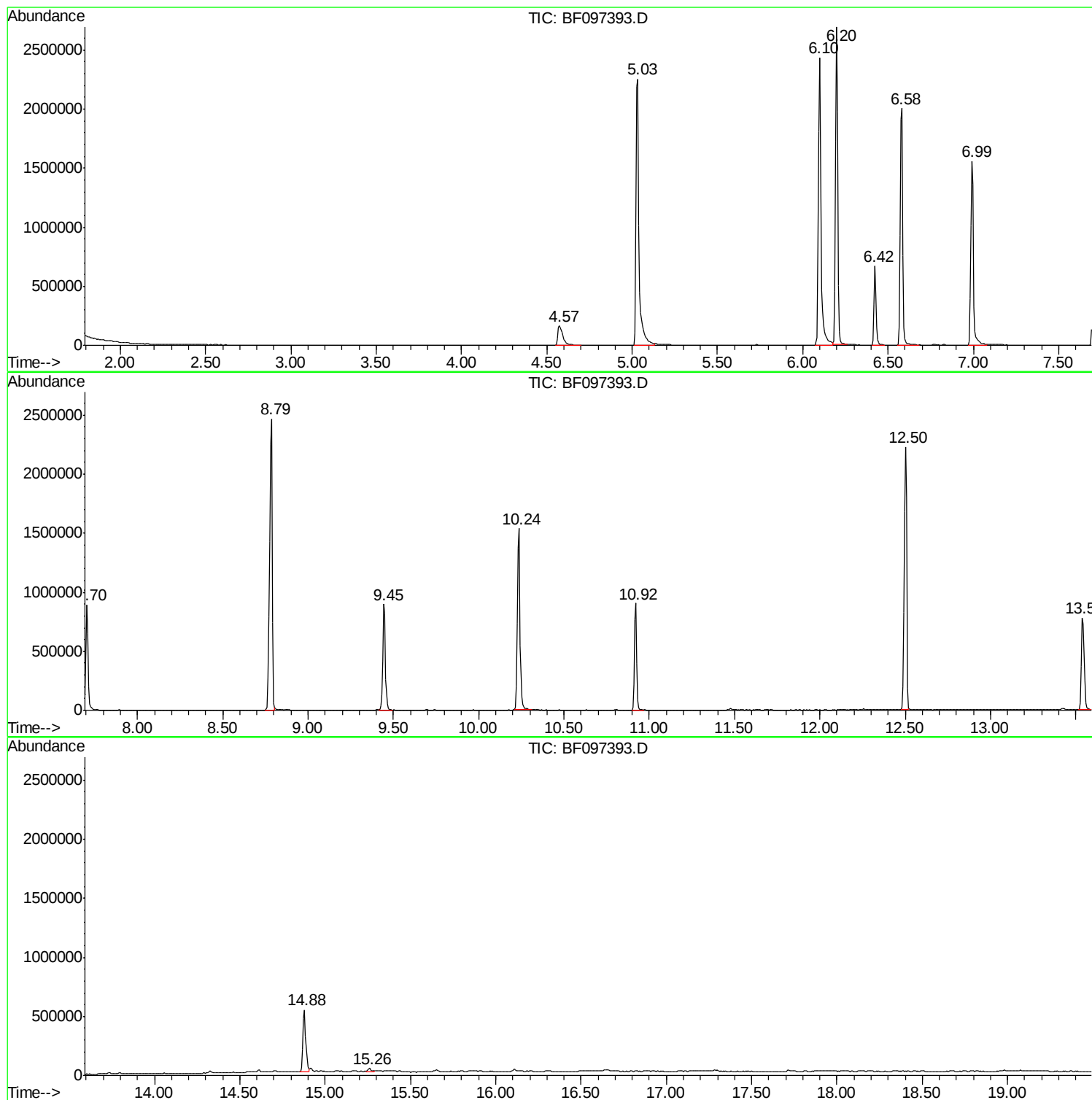
Sum of corrected areas: 24206610

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.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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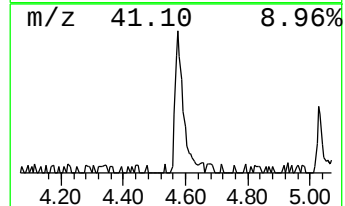
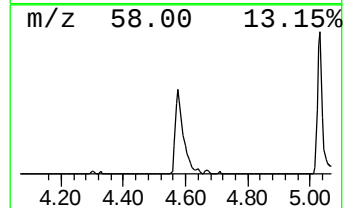
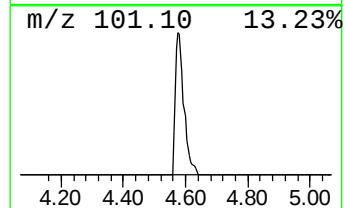
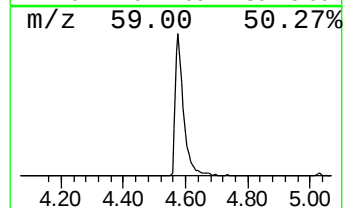
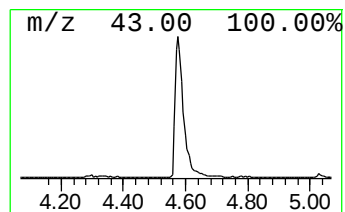
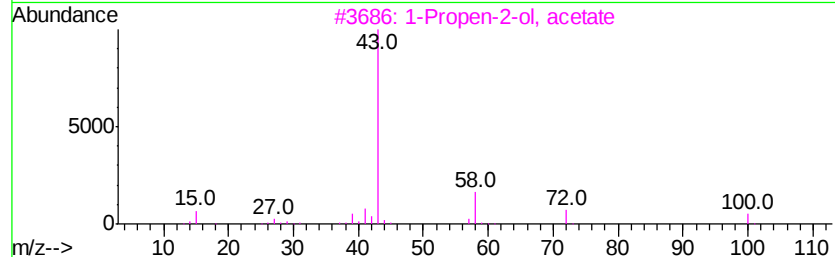
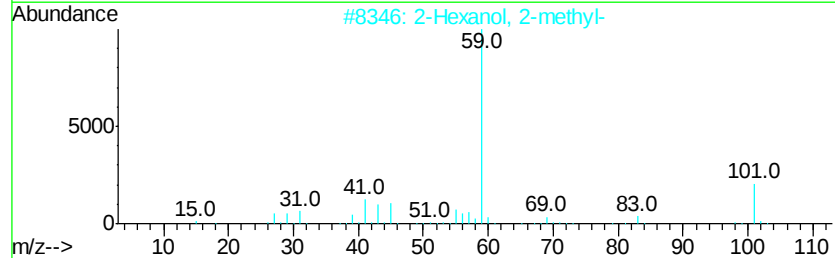
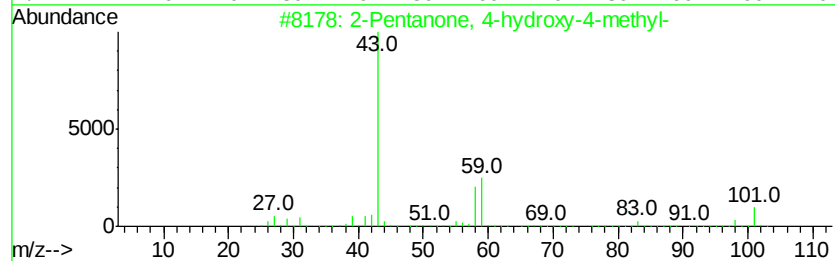
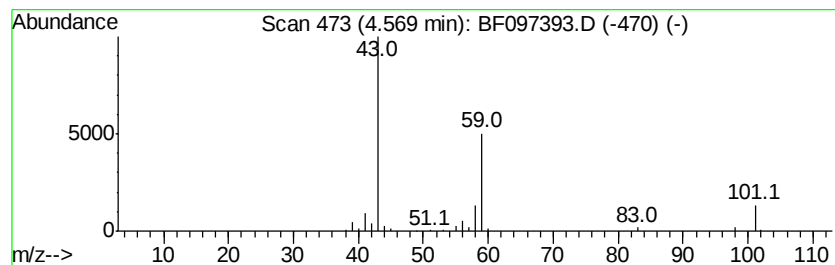
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.57	10.24 ng	309533	1,4-Dichlorobenzene-d4	6.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
5		Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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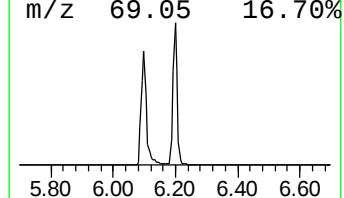
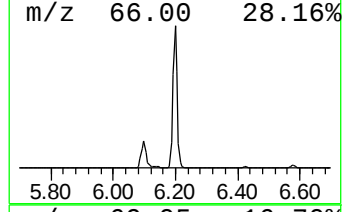
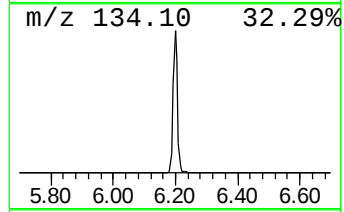
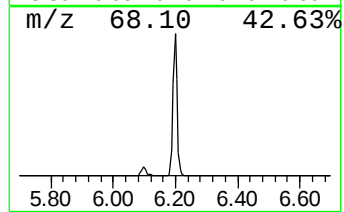
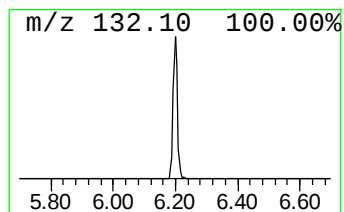
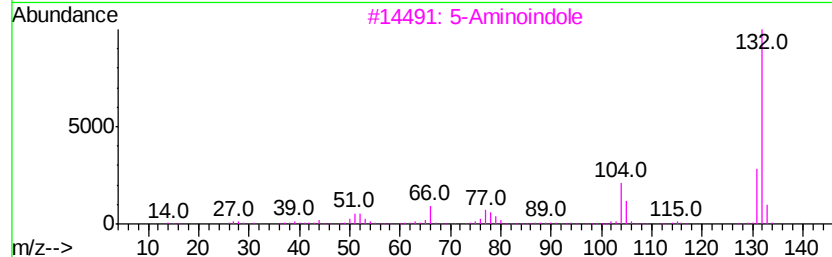
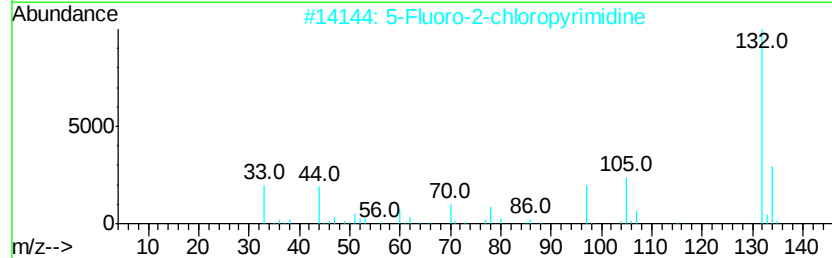
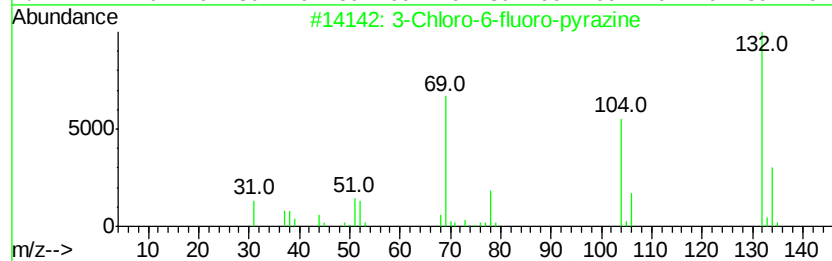
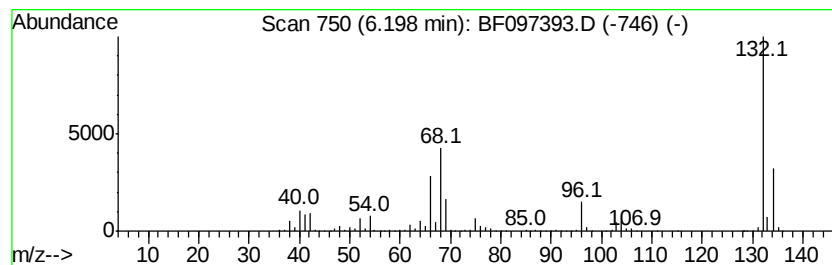
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Peak Number 2 unknown6.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	88.30 ng	2668380	1,4-Dichlorobenzene-d4	6.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.57	10.2	ng	309533	1	6.42	604370	20.0
unknown6.20	6.20	88.3	ng	2668380	1	6.42	604370	20.0