

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080817\
 Data File : BF097480.D
 Acq On : 8 Aug 2017 20:44
 Operator : SJ/JU
 Sample : PB101311BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101311BL

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	4.557	467	471	483	rBV	106184	210607	8.37%	1.009%
2	5.004	544	547	577	rBV	1866477	2372391	94.30%	11.367%
3	6.075	725	729	743	rBV	2043334	2277203	90.52%	10.910%
4	6.181	743	747	762	rVB	2156691	2271878	90.31%	10.885%
5	6.404	782	785	797	rVB	572889	526254	20.92%	2.521%
6	6.557	807	811	839	rVB	1826662	1782953	70.87%	8.542%
7	6.975	879	882	908	rBV	1462178	1571136	62.45%	7.528%
8	7.687	1000	1003	1019	rBV	791823	791754	31.47%	3.793%
9	8.763	1181	1186	1201	rBV	2247451	2515752	100.00%	12.053%
10	9.428	1295	1299	1312	rBV2	822530	819107	32.56%	3.924%
11	10.222	1430	1434	1452	rBV	1259086	1397990	55.57%	6.698%
12	10.904	1546	1550	1571	rBB	849957	784152	31.17%	3.757%
13	12.486	1815	1819	1823	rBV	1973227	2124619	84.45%	10.179%
14	13.527	1992	1996	2011	rBV	758003	766192	30.46%	3.671%
15	14.868	2219	2224	2235	rBV	494191	659680	26.22%	3.161%

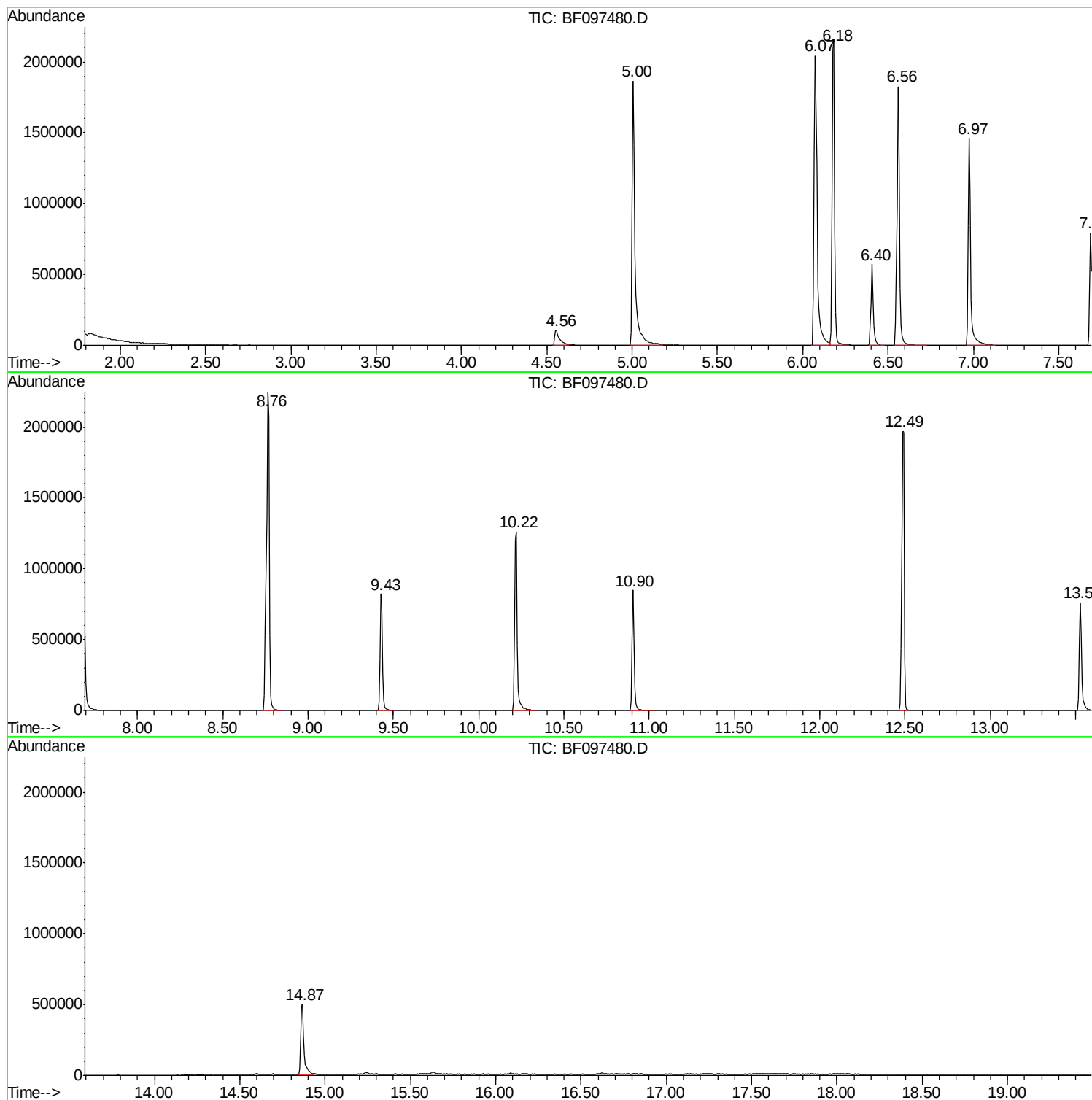
Sum of corrected areas: 20871668

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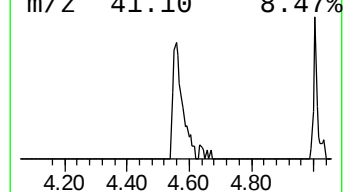
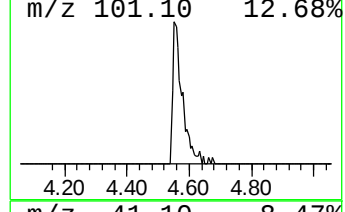
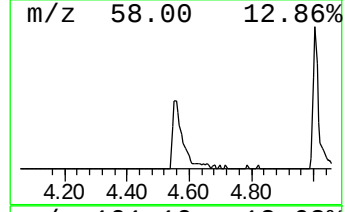
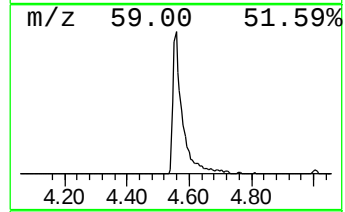
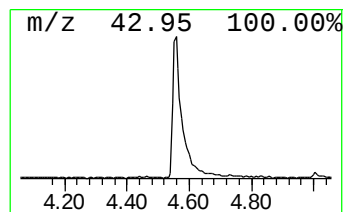
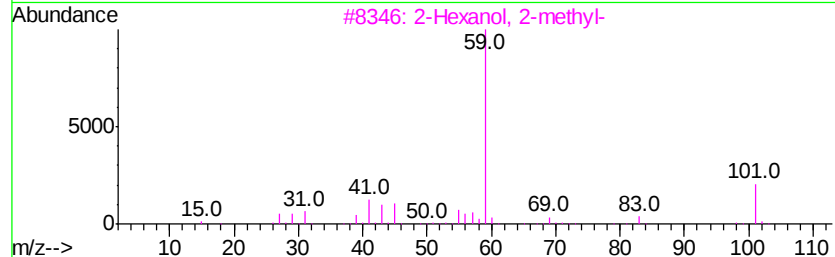
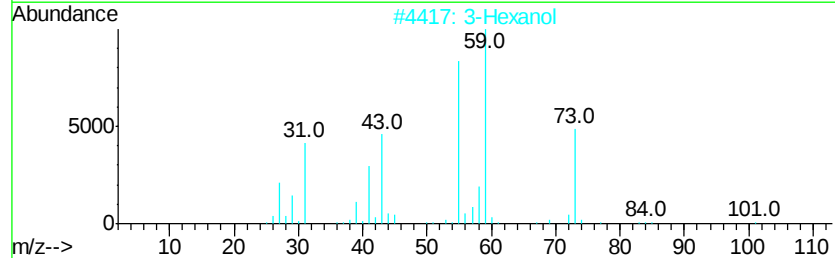
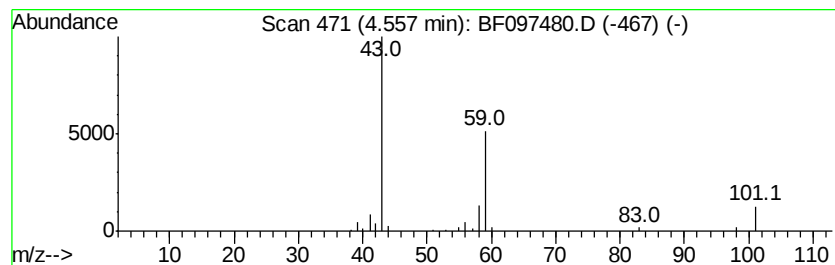
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	8.00 ng	210607	1,4-Dichlorobenzene-d4	6.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Hexanol	102	C6H14O	000623-37-0	33
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		3-Pentanol	88	C5H12O	000584-02-1	9
5		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9



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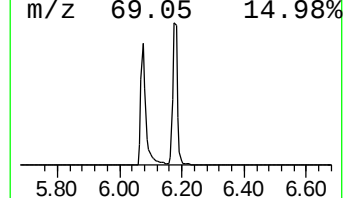
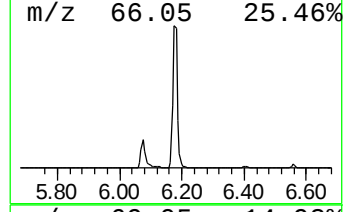
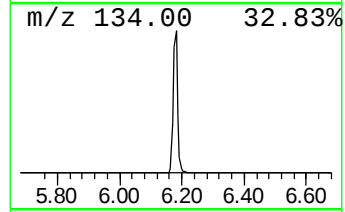
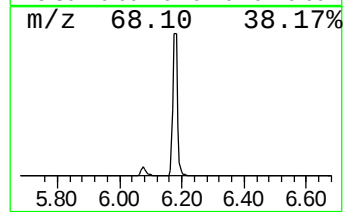
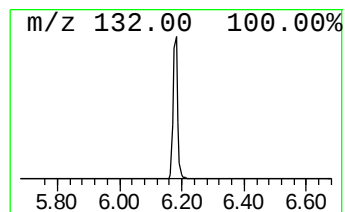
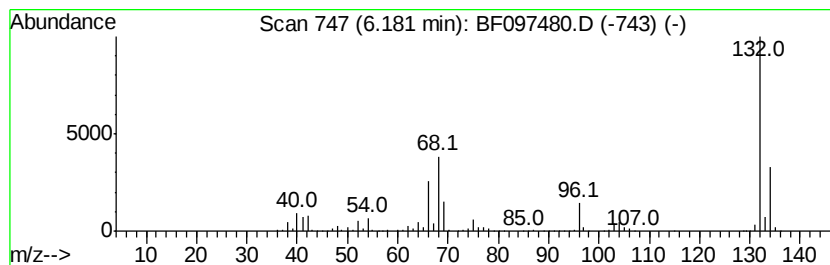
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Peak Number 2 unknown6.18 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.18	86.34 ng	2271880	1,4-Dichlorobenzene-d4	6.40

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		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...	4.56	8.0	ng	210607	1	6.40	526254	20.0
unknown6.18	6.18	86.3	ng	2271880	1	6.40	526254	20.0