



00000000 00000000 00000000 00000000 000000 000000 00000000 000000 000000
00 00000000 0000 00000000 00 00000000 00000000 00 00000000 00 00000000
:0000 00000000 000000 000000 0000

000000 0000 00000000 «00000000» 0000 .0
0000 00000000 00000000 00 0000 0000 00 000000 0 000000 00000000 0000
000000 00 00 (BMR) 000000 00000000 00 000000 00 00 00000000
00 00000000 000000 00000000 00 0000 0000 00000000 000000 0000000000
000000 000000 00.0000 00000000 0000 0000000000 000000 00000000 0000
000000:0000 00000000 00000000 0000 00 00000000 00000000 00000000 0000
000000 0000000000 00000000 00000000 0000000000 :000000 00000000 0 00000000
0000 00 000000 0000000000 000000 0000 000000 .00000000 00000000 00 000000
00000000 00000000000000 000000 0 0000 0000000000 00000000 00 00000000
00000000 00000000 00 00000000 00000000 00 (000000000000 0 000000000000)
00 0000000000 000000000000 0 000000 00000000 000000 0000000000 0000 00
.0000 0000 00000000 00000000

:0000000 000000

0000000 0 000000 00000000 00000 00 00000000 00 00000000000000
0 000000 000000 00 00000000000 0000000000 00 000000000 00000
0 00000 0000 00000000000000 000000 000000000-00000000 00000000
.0000000 00000000 00 00000 0000000 00000

:00000000 00000000 00 00000 000000

000000 00000 00 000000000 + (0000 00000) 0000000 00000000000000
+ 00 00 0000 0000 + 0000 00000 0000 :00000 00000 00000 + (BV) 0000
00000+00000+0000000000 00000000000 0000 0000

00000000 0000 0000 0 000000000 00 000000 00000000000000 00.0
0 00000 00000 0000 00000000000 00000 0000 00 000000 0000000000
:000000 : (Iron) 0000 :00000000000 0000 000000 00 00000000
0000000 0 000000000 00000000 00000 00 000000 0000000000
000000 00 0000) 0000 0000 00000000 .00000 000000000 00000000
0000000 0000 0000 00 (000000 000000 00000 00000000 000000 000000
.0000000 000000 00000 0000000 0 00000000

0000000 0000000000 0000000000 00000000 00000000 00000 00000 :000000
0000 0000000000000000 00000000 000000 C 000000000 00
:000000 : (EPA 0 DHA 00000000) 0-00000 0000 000000000 .0000000
DHA 0000 000000000 00 00000000000000000000 00000 0000000000 00000000
00000000 0 0000 00000000 0 0000000 0000000 00 0000000 0 0000
.0000000 0000000 00 000000 0000000000

0 0000 00000 0(00000000) 00000 0000 000000000 :000000
:000000 : (0000000000000000 00000) D 000000000 0 0000000 .00000
00000 0000» 00 0000000000 00000 0000000 0000000000 0 0000000 00000
00000 0 00 0 :000000 .0000 (Mass Bone Peak) «0000000000
000000 00 0000 00 000000000000/0000000 000000000000 00000000
00000000 0000 00000000 00000000000 0 000000 0000 00000
0000000 000000 0000 :000000 : (Iodine) 00 0 (Zinc) 0000 .000000
00000 000000 00 0 000000000 00 0 0000000 0000000 00000000

.000000 000000 00000000 000000 0 0000000000 0000000000
00000000 00000000 0 0000 00 0000 0 00 00000 0 000000 :000000

0000 000000 000000 00 00000000000000000000 0000 000000 00 000000 .0
00000 0000000000000000 000000 00000000 00000000 0 00000000 0000 00 0000
00 000000 0 0000.00000000 000000 000000000000000000 000000 00000000
0 0000000000000000 00000000000000 00000000 0000 00000000 00 000000 :0000

.000000 000000 000000 000000
00 0000000000 00000000000000 00 0000000000 00000000 00 000000 :000000
0 00000 0000 0000 00000000 0000) 0000000 000000 000000000000000000
0000000 0000 00 00 000000 00000000 00000000 0 (0000 000000 000000
.000000 00000000 00000000 00 000000