

0000 000000 00000 00000 00000 00 000000 0000000 00000000 000000 0000 3 0000 00
00 000000 0000000 00000000 0000000 .0000 0000000 00000 0000000 00000 00 0000000
0000 00 0000 .0000 00000000 00000000 00000000 0000 00 00000 00000000 00 0000000000
0000000 0000000000 00 0000 0000000 0000000 0000000 000000 000000 00000 000000 00 0000
00 0000 00000000 00000 0000000 00000 00000 00000 0000 0000000 00 0000 00000000 .000000
000000000 000000000000 00000 0000 0000000 00 0000 000000000000 0000000 000000000
0 (As) 00000000 0000000 0000000 0000000 00000 0000000 0000000000 0 000000
0000000 00000000 000000000 00 (Pb) 00000 0 (Ni) 000000 0 (Cr) 000000 0 (Cd) 0000000000
00000 000000 0000000000 000000 00 000000000 000000 000000 00 0000000 0000000000 000000000
00000 00000 0 0000000 000000 00 0000000 0000 00 0000000 0000000000 00000000 000000 0
0000000000 000000 00 00000000000 00 0000000000 00000 00000 0000000000 .000000 0000000
0000000 0000000000 0000000 000000 00 000000 0000000000 0000000 00 0000000 0000000 00 0000

0000000 00 0000000000000 000000000 00000000 0000000 0000000000 00 0000 0000000 00 0000000
0000000 00 0000000000 .0000 0000000 0000 000000000 0000000000 00000 0000000 00 00000000
0000000 00000 0000 0000 00 .0000000 00000000 0000000 00 0000 0000000 -0000000 0000 00
0000000 00 0000 0000000 0000 00 0000000 0000000 00 0000000 0000000000000 00 00000000
00000000 00000 0000000 0000000) 0000000 0000 00000 0000000 0000 00 000000000 0000000
0000000 000000 4 00 000000 00 00000000 00000 000000000000 0000000000 (000000000000
1000 0 500 0 100 0000000 00 0000000 3 0000 00 00 0 0000 0000000 0000 0 0000 0000000
00 00 .00 0000000 0000000 (000000000000 15 00 0) 0000 0000000 00000 00 000000
0.2 00000 0000000 5 00000 00000000000 00 0000000 00 0000000000 0000000 00000000
.00 00000000 0000000 00000000 00 00000 00 00 0000 0000000 00000 00 00000000000
0000000 00 00 0000000 00 .000000 000000000 0000000 50 000000 0000000 00 0000000000-
0000000 0000000000000 00 0 0000 0000000 0000 0000 0000 00 00 0000000 0000000000
0000 000000 00 0000 00 0000000 0000 00 0000 0000 000000000000000000 00 .00 0000000
0000 0000 0000000 00 0000 0.5 00000000 0000000 00000 0000000 00000 0000000 00 0
-HCl-HF 0000 0000000 00. 00 4-HClO₃HNO 00000 00000 10 00 0000 0 00 0000000
00 0000000 0000000 000000 0000000000 00000000000 00000000 0000000 00 0000000
0000000 (MS-ICP) 000000000 0000 0000 0000000 00000 00000 0000 000000000 00 0000000000
0000000000 000000000 00 000000 00000000 00 0000000000 000000 0000000000 .00 0000000
(EPA US) 000000000 000000 000000 0000000 000000000 000000 00000000000 00000 00000000 0

00000000 0 0000000 0000000 0 00000 000000000 000000 00 0000000 0 00000000 00000000
0000000 0000 00000 0000000 0000000 0000 00 0000000 0000000 00000000000 00000
.0000 000000 0000000000 0000000000 (00000000000 000000000 000000 0000000 00000000)
0000000000 0000000000 00 000000 00000000 000000000 0000 00 0000 0000000000000
00 0 (Cr) 000000 0 (Co) 00000000 0 (Cd) 00000000000 0 (As) 00000000) 00000000 00000000

