

data\_ctx1dc\_abs

\_audit\_creation\_method        SHELXL-97  
\_chemical\_name\_systematic  
;  
?  
;  
\_chemical\_name\_common        ?  
\_chemical\_melting\_point       ?  
\_chemical\_formula\_moiety       ?  
\_chemical\_formula\_sum  
'H B Ca O5 Si'  
\_chemical\_formula\_weight       159.99

loop\_  
\_atom\_type\_symbol  
\_atom\_type\_description  
\_atom\_type\_scatter\_dispersion\_real  
\_atom\_type\_scatter\_dispersion\_imag  
\_atom\_type\_scatter\_source  
'H' 'H' 0.0000 0.0000  
'International Tables Vol C Tables 4.2.6.8 and 6.1.1.4'  
'B' 'B' 0.0013 0.0007  
'International Tables Vol C Tables 4.2.6.8 and 6.1.1.4'  
'O' 'O' 0.0106 0.0060  
'International Tables Vol C Tables 4.2.6.8 and 6.1.1.4'  
'Si' 'Si' 0.0817 0.0704  
'International Tables Vol C Tables 4.2.6.8 and 6.1.1.4'  
'Ca' 'Ca' 0.2262 0.3064  
'International Tables Vol C Tables 4.2.6.8 and 6.1.1.4'

\_symmetry\_cell\_setting        ?  
\_symmetry\_space\_group\_name\_H-M ?

loop\_  
\_symmetry\_equiv\_pos\_as\_xyz  
'x, y, z'  
'-x, y+1/2, -z+1/2'  
'-x, -y, -z'  
'x, -y-1/2, z-1/2'

\_cell\_length\_a                4.83550(10)  
\_cell\_length\_b                7.61020(10)  
\_cell\_length\_c                9.63170(10)  
\_cell\_angle\_alpha              90.00  
\_cell\_angle\_beta               90.1500(10)  
\_cell\_angle\_gamma              90.00  
\_cell\_volume                   354.437(9)  
\_cell\_formula\_units\_Z         4  
\_cell\_measurement\_temperature 293(2)  
\_cell\_measurement\_reflns\_used ?  
\_cell\_measurement\_theta\_min   ?

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_cell_measurement_theta_max      ?

_exptl_crystal_description      ?
_exptl_crystal_colour           ?
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_exptl_crystal_size_mid         ?
_exptl_crystal_size_min         ?
_exptl_crystal_density_meas     ?
_exptl_crystal_density_diffn    2.998
_exptl_crystal_density_method   'not measured'
_exptl_crystal_F_000            320
_exptl_absorpt_coefficient_mu    1.997
_exptl_absorpt_correction_type  ?
_exptl_absorpt_correction_T_min ?
_exptl_absorpt_correction_T_max ?
_exptl_absorpt_process_details  ?

_exptl_special_details
;
?
;

_diffn_ambient_temperature      293(2)
_diffn_radiation_wavelength     0.71073
_diffn_radiation_type           MoK\alpha
_diffn_radiation_source         'fine-focus sealed tube'
_diffn_radiation_monochromator  graphite
_diffn_measurement_device_type  ?
_diffn_measurement_method       ?
_diffn_detector_area_resol_mean ?
_diffn_standards_number         ?
_diffn_standards_interval_count ?
_diffn_standards_interval_time  ?
_diffn_standards_decay_%        ?
_diffn_reflns_number            7893
_diffn_reflns_av_R_equivalents  0.0578
_diffn_reflns_av_sigmaI/netI    0.0329
_diffn_reflns_limit_h_min       -8
_diffn_reflns_limit_h_max       7
_diffn_reflns_limit_k_min       -13
_diffn_reflns_limit_k_max       13
_diffn_reflns_limit_l_min       -16
_diffn_reflns_limit_l_max       17
_diffn_reflns_theta_min         3.41
_diffn_reflns_theta_max         40.08
_reflns_number_total            2211
_reflns_number_gt               2081
_reflns_threshold_expression     >2sigma(I)

_computing_data_collection      ?
_computing_cell_refinement      ?
_computing_data_reduction       ?
_computing_structure_solution   ?
_computing_structure_refinement 'SHELXL-97 (Sheldrick, 1997)'

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\_computing\_molecular\_graphics ?  
\_computing\_publication\_material ?

\_refine\_special\_details

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Refinement of  $F^2$  against ALL reflections. The weighted R-factor  $wR$  and goodness of fit  $S$  are based on  $F^2$ , conventional R-factors  $R$  are based on  $F$ , with  $F$  set to zero for negative  $F^2$ . The threshold expression of  $F^2 > 2\sigma(F^2)$  is used only for calculating R-factors(gt) etc. and is not relevant to the choice of reflections for refinement. R-factors based on  $F^2$  are statistically about twice as large as those based on  $F$ , and R-factors based on ALL data will be even larger.

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\_refine\_ls\_structure\_factor\_coef Fsqd  
\_refine\_ls\_matrix\_type full  
\_refine\_ls\_weighting\_scheme calc  
\_refine\_ls\_weighting\_details  
'calc w=1/[\sigma^2(Fo^2)+(0.0300P)^2+0.0000P] where P=(Fo^2+2Fc^2)/3'  
\_atom\_sites\_solution\_primary direct  
\_atom\_sites\_solution\_secondary difmap  
\_atom\_sites\_solution\_hydrogens geom  
\_refine\_ls\_hydrogen\_treatment mixed  
\_refine\_ls\_extinction\_method SHELXL  
\_refine\_ls\_extinction\_coef 0.077(8)  
\_refine\_ls\_extinction\_expression  
'Fc^\*=kFc[1+0.001xFc^2\l^3/sin(2\q)]^-1/4^'  
\_refine\_ls\_number\_reflns 2211  
\_refine\_ls\_number\_parameters 78  
\_refine\_ls\_number\_restraints 0  
\_refine\_ls\_R\_factor\_all 0.0361  
\_refine\_ls\_R\_factor\_gt 0.0346  
\_refine\_ls\_wR\_factor\_ref 0.0890  
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loop\_

\_atom\_site\_label  
\_atom\_site\_type\_symbol  
\_atom\_site\_fract\_x  
\_atom\_site\_fract\_y  
\_atom\_site\_fract\_z  
\_atom\_site\_U\_iso\_or\_equiv  
\_atom\_site\_adp\_type  
\_atom\_site\_occupancy  
\_atom\_site\_symmetry\_multiplicity  
\_atom\_site\_calc\_flag  
\_atom\_site\_refinement\_flags  
\_atom\_site\_disorder\_assembly  
\_atom\_site\_disorder\_group

Ca Ca 0.99165(4) 0.10622(2) 0.336128(19) 0.00629(7) Uani 1 1 d . . .

Si Si 0.46903(5) 0.26609(3) 0.08442(3) 0.00424(7) Uani 1 1 d . . .  
 B B 0.5682(2) 0.41054(13) 0.34086(11) 0.00512(15) Uani 1 1 d . . .  
 O1 O 0.24040(16) 0.40095(9) 0.03741(8) 0.00767(13) Uani 1 1 d . . .  
 O2 O 0.66992(14) 0.30113(9) 0.45732(8) 0.00621(12) Uani 1 1 d . . .  
 O3 O 0.67513(14) 0.33452(10) 0.21048(7) 0.00607(12) Uani 1 1 d . . .  
 O4 O 0.31507(15) 0.08816(9) 0.14600(8) 0.00699(12) Uani 1 1 d . . .  
 O5 O 0.25878(15) 0.41350(10) 0.33644(8) 0.00724(12) Uani 1 1 d . . .  
 H H 0.199(4) 0.452(3) 0.412(3) 0.022(6) Uiso 1 1 d . . .

loop\_

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 \_atom\_site\_aniso\_U\_11  
 \_atom\_site\_aniso\_U\_22  
 \_atom\_site\_aniso\_U\_33  
 \_atom\_site\_aniso\_U\_23  
 \_atom\_site\_aniso\_U\_13  
 \_atom\_site\_aniso\_U\_12

Ca 0.00581(9) 0.00731(9) 0.00575(10) -0.00004(5) 0.00018(6) -0.00048(5)  
 Si 0.00429(11) 0.00456(11) 0.00388(12) -0.00020(7) -0.00016(8) 0.00034(7)  
 B 0.0048(4) 0.0058(3) 0.0048(4) 0.0002(3) 0.0001(3) 0.0002(3)  
 O1 0.0077(3) 0.0076(3) 0.0077(3) -0.0002(2) -0.0014(2) 0.0025(2)  
 O2 0.0058(2) 0.0077(3) 0.0051(3) 0.00230(19) 0.00108(18) 0.00122(19)  
 O3 0.0054(2) 0.0085(3) 0.0043(3) -0.0018(2) -0.00029(18) -0.00001(19)  
 O4 0.0070(3) 0.0046(2) 0.0094(3) 0.0007(2) 0.0025(2) 0.00049(19)  
 O5 0.0045(2) 0.0107(3) 0.0066(3) -0.0011(2) 0.0003(2) 0.0006(2)

\_geom\_special\_details

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All esds (except the esd in the dihedral angle between two l.s. planes)  
 are estimated using the full covariance matrix. The cell esds are taken  
 into account individually in the estimation of esds in distances, angles  
 and torsion angles; correlations between esds in cell parameters are only  
 used when they are defined by crystal symmetry. An approximate (isotropic)  
 treatment of cell esds is used for estimating esds involving l.s. planes.

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loop\_

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 \_geom\_bond\_atom\_site\_label\_2  
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 \_geom\_bond\_site\_symmetry\_2  
 \_geom\_bond\_publ\_flag  
 Ca O1 2.2782(8) 2\_645 ?  
 Ca O1 2.2795(8) 4\_666 ?  
 Ca O4 2.4152(8) 1\_655 ?  
 Ca O2 2.4480(7) . ?  
 Ca O5 2.5238(8) 2\_645 ?  
 Ca O3 2.6107(7) . ?  
 Ca O3 2.6603(8) 2\_745 ?  
 Ca O5 2.6716(8) 1\_655 ?  
 Si O1 1.5741(8) . ?  
 Si O2 1.6463(7) 4\_565 ?  
 Si O3 1.6529(7) . ?  
 Si O4 1.6559(7) . ?

B O4 1.4702(12) 2\_655 ?  
B O3 1.4773(13) . ?  
B O2 1.4800(12) . ?  
B O5 1.4969(13) . ?  
O1 Ca 2.2782(8) 2\_655 ?  
O1 Ca 2.2795(8) 4\_465 ?  
O2 Si 1.6463(7) 4\_566 ?  
O3 Ca 2.6603(8) 2\_755 ?  
O4 B 1.4702(12) 2\_645 ?  
O4 Ca 2.4152(8) 1\_455 ?  
O5 Ca 2.5238(8) 2\_655 ?  
O5 Ca 2.6715(8) 1\_455 ?

loop\_  
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O1 Ca O1 77.78(3) 2\_645 4\_666 ?  
O1 Ca O4 133.41(3) 2\_645 1\_655 ?  
O1 Ca O4 107.59(3) 4\_666 1\_655 ?  
O1 Ca O2 81.14(3) 2\_645 . ?  
O1 Ca O2 86.76(3) 4\_666 . ?  
O4 Ca O2 144.06(2) 1\_655 . ?  
O1 Ca O5 73.59(3) 2\_645 2\_645 ?  
O1 Ca O5 142.89(3) 4\_666 2\_645 ?  
O4 Ca O5 77.18(3) 1\_655 2\_645 ?  
O2 Ca O5 111.21(3) . 2\_645 ?  
O1 Ca O3 114.53(3) 2\_645 . ?  
O1 Ca O3 135.85(3) 4\_666 . ?  
O4 Ca O3 93.81(2) 1\_655 . ?  
O2 Ca O3 56.31(2) . . ?  
O5 Ca O3 78.57(2) 2\_645 . ?  
O1 Ca O3 81.76(3) 2\_645 2\_745 ?  
O1 Ca O3 78.84(2) 4\_666 2\_745 ?  
O4 Ca O3 55.54(2) 1\_655 2\_745 ?  
O2 Ca O3 159.62(2) . 2\_745 ?  
O5 Ca O3 74.16(2) 2\_645 2\_745 ?  
O3 Ca O3 142.526(11) . 2\_745 ?  
O1 Ca O5 146.97(3) 2\_645 1\_655 ?  
O1 Ca O5 76.46(2) 4\_666 1\_655 ?  
O4 Ca O5 74.73(2) 1\_655 1\_655 ?  
O2 Ca O5 77.11(2) . 1\_655 ?  
O5 Ca O5 137.781(17) 2\_645 1\_655 ?  
O3 Ca O5 72.59(2) . 1\_655 ?  
O3 Ca O5 112.80(2) 2\_745 1\_655 ?  
O1 Si O2 113.83(4) . 4\_565 ?  
O1 Si O3 115.31(4) . . ?  
O2 Si O3 106.76(4) 4\_565 . ?  
O1 Si O4 108.67(4) . . ?  
O2 Si O4 106.20(4) 4\_565 . ?

O3 Si O4 105.38(4) . . ?  
 O4 B O3 107.34(8) 2\_655 . ?  
 O4 B O2 108.98(8) 2\_655 . ?  
 O3 B O2 107.91(7) . . ?  
 O4 B O5 111.84(7) 2\_655 . ?  
 O3 B O5 109.50(8) . . ?  
 O2 B O5 111.11(8) . . ?  
 Si O1 Ca 129.77(5) . 2\_655 ?  
 Si O1 Ca 126.71(4) . 4\_465 ?  
 Ca O1 Ca 102.22(3) 2\_655 4\_465 ?  
 B O2 Si 122.88(6) . 4\_566 ?  
 B O2 Ca 100.94(5) . . ?  
 Si O2 Ca 122.90(4) 4\_566 . ?  
 B O3 Si 122.42(6) . . ?  
 B O3 Ca 94.17(5) . . ?  
 Si O3 Ca 118.84(4) . . ?  
 B O3 Ca 93.01(5) . 2\_755 ?  
 Si O3 Ca 119.03(4) . 2\_755 ?  
 Ca O3 Ca 103.94(2) . 2\_755 ?  
 B O4 Si 127.58(6) 2\_645 . ?  
 B O4 Ca 103.67(5) 2\_645 1\_455 ?  
 Si O4 Ca 121.18(4) . 1\_455 ?  
 B O5 Ca 120.35(6) . 2\_655 ?  
 B O5 Ca 118.04(5) . 1\_455 ?  
 Ca O5 Ca 106.07(3) 2\_655 1\_455 ?

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