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_publ_contact_author_name   'MATTEO ALVARO'
_publ_contact_author_address
;
'Dipartimento di Scienze della Terra,
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loop_
_publ_author_name
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'ALVARO, MATTEO'
;'Dipartimento di Scienze della Terra,
Universit\`a di Pavia, Via Ferrata 1, I-27100 Pavia (Italy)'
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'Nestola, Fabrizio'
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 Via Giotto 1, I-35137 Padova (Italy)'
;
'BOFFA BALLARAN, TIZIANA'
;'Bayerisches Geoinstitut, Universit\"at Bayreuth,
Universit\"atstrasse 37, D-95440 Bayreuth (Germany)'
;
'C\'AMARA, FERNANDO'
;'C.N.R. - Istituto di Geoscienze e Georisorse - Unit\`a di Pavia,
 Via Ferrata 1, I-27100 Pavia (Italy)'
;
'DOMENEGHETTI, CHIARA'
;'Dipartimento di Scienze della Terra, Universit\`a di Pavia,
 Via Ferrata 1, I-27100 Pavia (Italy)'
;
'TAZZOLI, VITTORIO'
;'Dipartimento di Scienze della Terra, Universit\`a di Pavia,
 Via Ferrata 1, I-27100 Pavia (Italy)'
;
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Angel, R.J. (2004) Absorption corrections for diamond-anvil pressure cells
implemented in the software package -- Absorb 6.0. Journal of Applied
Crystallography, 37, 486-492.
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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
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 F^2^ > 2sigma(F^2^) is used only for calculating R-factors(gt) etc. and is
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Fe2A FeP2 0.0000 0.2792(2) 0.2500 0.0146(7) Uani 0.75 2 d SP . .
FE2B MgP2 0.0000 0.2792(2) 0.2500 0.0146(7) Uani 0.12 2 d SP . .
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Fe2A 0.0202(17) 0.0134(12) 0.0085(9) 0.000 -0.0003(9) 0.000
FE2B 0.0202(17) 0.0134(12) 0.0085(9) 0.000 -0.0003(9) 0.000
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 used when they are defined by crystal symmetry. An approximate (isotropic)
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O3 Si 1.660(6) 1_556 ?
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