



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 6, 2022 – 02:27 pm BST

PDB ID : 7R0G
Title : STRUCTURAL BASIS OF ION UPTAKE IN COPPER-TRANSPORTING P1B-TYPE ATPASES
Authors : Salustros, N.; Groenberg, C.; Wang, K.; Gourdon, P.
Deposited on : 2022-02-02
Resolution : 4.01 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.30
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.30

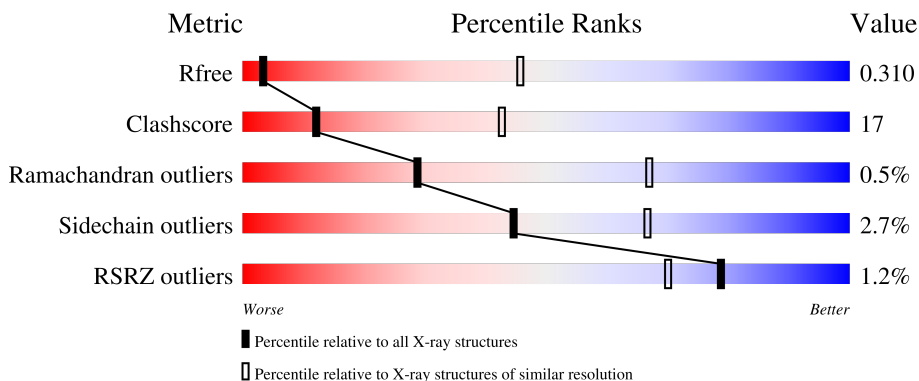
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.01 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1098 (4.34-3.70)
Clashscore	141614	1159 (4.34-3.70)
Ramachandran outliers	138981	1118 (4.34-3.70)
Sidechain outliers	138945	1108 (4.34-3.70)
RSRZ outliers	127900	1034 (4.38-3.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	658	 63% 35% ..
1	B	658	 62% 35% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9841 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Putative copper-exporting P-type ATPase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	654	4929	3176	829	909	15	0	0	0
1	B	652	4912	3167	827	903	15	0	0	0

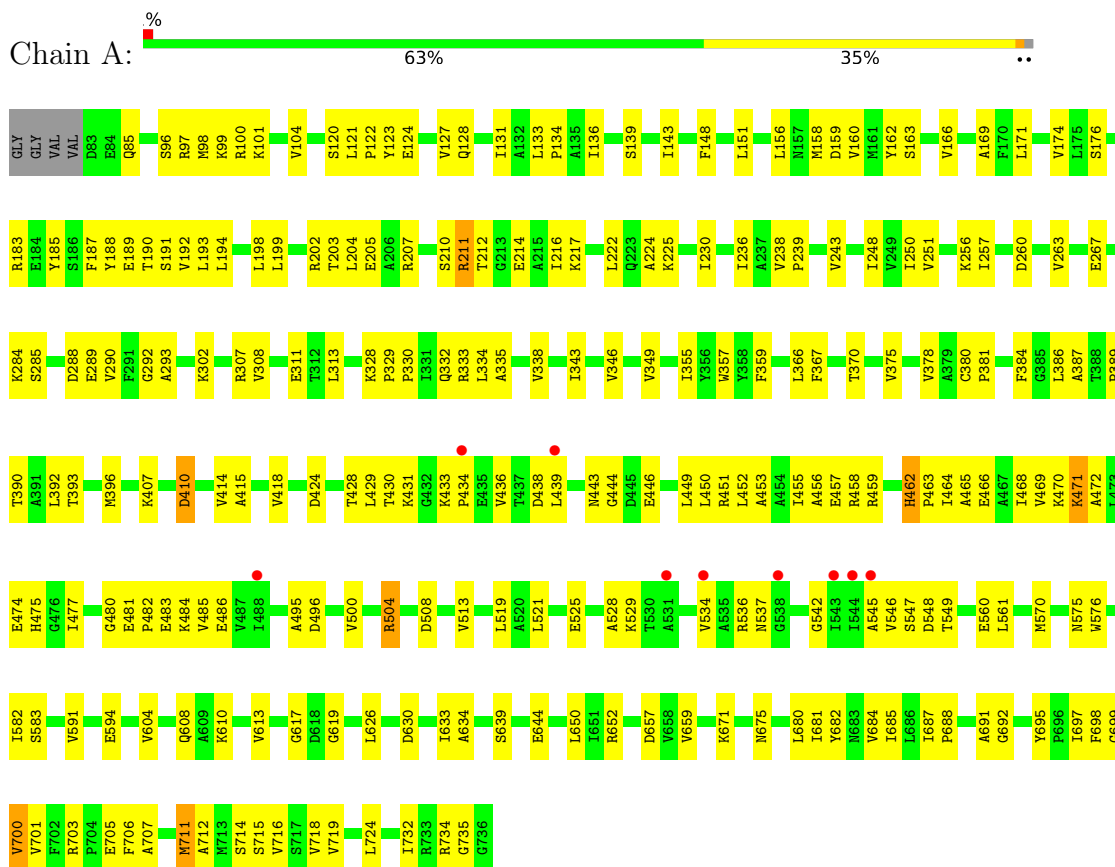
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	GLY	-	expression tag	UNP A0A117KM49
B	79	GLY	-	expression tag	UNP A0A117KM49

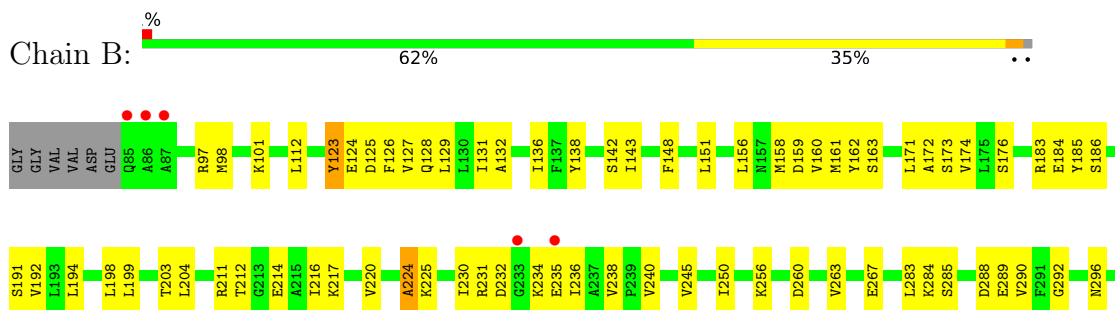
3 Residue-property plots [i](#)

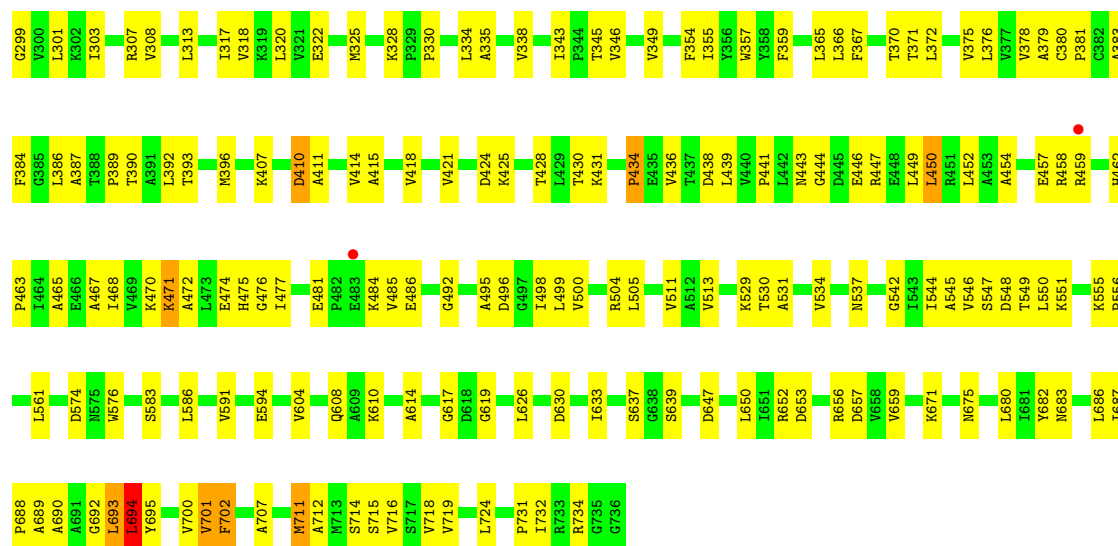
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Putative copper-exporting P-type ATPase A



- Molecule 1: Putative copper-exporting P-type ATPase A





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.19Å 226.06Å 111.11Å 90.00° 134.06° 90.00°	Depositor
Resolution (Å)	48.75 – 4.01 48.75 – 4.01	Depositor EDS
% Data completeness (in resolution range)	94.0 (48.75-4.01) 94.0 (48.75-4.01)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 4.00Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.263 , 0.312 0.261 , 0.310	Depositor DCC
R_{free} test set	1788 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	191.1	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9841	wwPDB-VP
Average B, all atoms (Å ²)	203.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.31	0/5003	0.63	1/6785 (0.0%)
1	B	0.33	0/4986	0.66	2/6762 (0.0%)
All	All	0.32	0/9989	0.65	3/13547 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	694	LEU	CB-CG-CD2	-12.81	89.22	111.00
1	A	519	LEU	CA-CB-CG	6.48	130.21	115.30
1	B	450	LEU	CA-CB-CG	5.21	127.28	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4929	0	5172	170	0
1	B	4912	0	5162	175	0
All	All	9841	0	10334	345	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (345) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:529:LYS:HG2	1:A:547:SER:HB2	1.51	0.92
1:B:438:ASP:HB2	1:B:545:ALA:HB3	1.52	0.89
1:A:313:LEU:HD12	1:A:639:SER:HA	1.58	0.83
1:B:457:GLU:HB2	1:B:468:ILE:HD11	1.60	0.83
1:B:692:GLY:O	1:B:694:LEU:N	2.12	0.82
1:B:424:ASP:O	1:B:428:THR:OG1	1.97	0.81
1:B:176:SER:HB2	1:B:185:TYR:HB2	1.63	0.81
1:A:267:GLU:HG3	1:A:285:SER:HB3	1.68	0.76
1:A:560:GLU:HG3	1:A:732:ILE:HD13	1.65	0.75
1:B:357:TRP:HZ2	1:B:692:GLY:HA2	1.52	0.75
1:B:355:ILE:O	1:B:359:PHE:HB2	1.87	0.74
1:B:212:THR:HG22	1:B:393:THR:HG21	1.71	0.73
1:A:525:GLU:OE2	1:A:575:ASN:ND2	2.23	0.72
1:A:459:ARG:NH1	1:A:483:GLU:OE2	2.23	0.72
1:A:101:LYS:HG2	1:A:139:SER:HA	1.72	0.71
1:B:158:MET:HA	1:B:380:CYS:SG	2.30	0.71
1:B:320:LEU:HD11	1:B:407:LYS:HG2	1.73	0.70
1:B:143:ILE:HG21	1:B:198:LEU:HD21	1.72	0.70
1:A:158:MET:HA	1:A:380:CYS:SG	2.31	0.70
1:A:212:THR:HG22	1:A:393:THR:HG21	1.72	0.70
1:A:159:ASP:O	1:A:163:SER:OG	2.05	0.70
1:B:450:LEU:HD21	1:B:472:ALA:HA	1.73	0.69
1:B:452:LEU:HD22	1:B:498:ILE:HG12	1.72	0.69
1:A:143:ILE:HG21	1:A:198:LEU:HD21	1.76	0.68
1:B:191:SER:HA	1:B:194:LEU:HB2	1.76	0.68
1:B:731:PRO:HA	1:B:734:ARG:HE	1.59	0.68
1:B:425:LYS:NZ	1:B:574:ASP:OD2	2.21	0.67
1:A:329:PRO:HD2	1:A:332:GLN:HB2	1.77	0.67
1:A:429:LEU:HD12	1:A:570:MET:HG3	1.76	0.67
1:B:136:ILE:HD11	1:B:192:VAL:HG22	1.76	0.67
1:A:251:VAL:HG21	1:A:257:ILE:HG12	1.76	0.66
1:B:296:ASN:ND2	1:B:299:GLY:O	2.27	0.66
1:B:357:TRP:CD2	1:B:367:PHE:HE1	2.13	0.66
1:A:366:LEU:O	1:A:370:THR:HG23	1.95	0.66
1:B:366:LEU:O	1:B:370:THR:HG23	1.95	0.66
1:B:688:PRO:HB2	1:B:693:LEU:HD22	1.77	0.66
1:B:325:MET:HE3	1:B:393:THR:HA	1.79	0.65
1:A:734:ARG:NE	1:A:735:GLY:H	1.96	0.64
1:B:550:LEU:HD11	1:B:586:LEU:HD21	1.78	0.64
1:B:583:SER:HB2	1:B:591:VAL:HG21	1.80	0.64
1:A:438:ASP:HB2	1:A:545:ALA:HB3	1.80	0.64
1:B:101:LYS:NZ	1:B:138:TYR:O	2.31	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ILE:HD11	1:A:192:VAL:HG22	1.78	0.64
1:B:439:LEU:HD13	1:B:446:GLU:HB3	1.80	0.63
1:B:499:LEU:HD22	1:B:505:LEU:HD11	1.80	0.63
1:A:176:SER:OG	1:A:183:ARG:NH2	2.30	0.63
1:A:482:PRO:HB3	1:A:496:ASP:HB2	1.79	0.63
1:A:187:PHE:HZ	1:A:370:THR:HG22	1.64	0.63
1:B:225:LYS:O	1:B:240:VAL:N	2.31	0.63
1:B:434:PRO:HA	1:B:548:ASP:OD1	1.99	0.63
1:A:250:ILE:HG22	1:A:302:LYS:HG2	1.81	0.62
1:A:355:ILE:O	1:A:359:PHE:HB2	1.99	0.62
1:B:125:ASP:OD2	1:B:183:ARG:NH2	2.25	0.62
1:A:381:PRO:HA	1:A:711:MET:HE3	1.81	0.62
1:B:230:ILE:HG12	1:B:235:GLU:HG2	1.80	0.62
1:A:456:ALA:HA	1:A:495:ALA:HB3	1.81	0.62
1:A:171:LEU:HA	1:A:174:VAL:HG12	1.81	0.61
1:A:289:GLU:OE2	1:A:307:ARG:NH2	2.30	0.61
1:A:205:GLU:HA	1:A:386:LEU:HD21	1.82	0.61
1:A:96:SER:HA	1:A:99:LYS:NZ	2.15	0.61
1:A:561:LEU:HG	1:A:659:VAL:HG22	1.83	0.61
1:B:267:GLU:HG2	1:B:283:LEU:HD11	1.82	0.61
1:B:731:PRO:HG3	1:B:734:ARG:HH21	1.66	0.60
1:B:267:GLU:HG3	1:B:285:SER:HB3	1.82	0.60
1:A:687:ILE:HG23	1:A:688:PRO:HD3	1.82	0.60
1:B:313:LEU:HD13	1:B:639:SER:HA	1.83	0.60
1:A:583:SER:HB2	1:A:591:VAL:HG21	1.84	0.59
1:A:500:VAL:HG22	1:A:534:VAL:HG12	1.84	0.59
1:A:191:SER:H	1:A:194:LEU:HD23	1.67	0.59
1:A:452:LEU:HD13	1:A:536:ARG:HB2	1.85	0.59
1:B:159:ASP:O	1:B:163:SER:OG	2.16	0.59
1:B:263:VAL:HG12	1:B:288:ASP:O	2.02	0.59
1:B:551:LYS:NZ	1:B:653:ASP:HB3	2.17	0.59
1:A:484:LYS:HB3	1:A:496:ASP:H	1.66	0.59
1:B:230:ILE:HD11	1:B:250:ILE:HB	1.84	0.59
1:B:284:LYS:HE3	1:B:290:VAL:HG12	1.85	0.59
1:A:212:THR:O	1:A:216:ILE:HD12	2.03	0.59
1:B:313:LEU:O	1:B:317:ILE:HG12	2.03	0.59
1:B:392:LEU:O	1:B:396:MET:HG3	2.02	0.59
1:A:191:SER:HA	1:A:194:LEU:HB2	1.84	0.58
1:A:263:VAL:HG11	1:A:285:SER:O	2.03	0.58
1:B:98:MET:HB3	1:B:203:THR:HG22	1.85	0.58
1:B:224:ALA:HB2	1:B:256:LYS:HD2	1.83	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:LEU:HD23	1:A:343:ILE:HG22	1.87	0.57
1:B:434:PRO:HG2	1:B:463:PRO:O	2.03	0.57
1:A:346:VAL:HA	1:A:349:VAL:HG22	1.85	0.57
1:A:671:LYS:HE3	1:A:724:LEU:HD22	1.85	0.57
1:B:561:LEU:HG	1:B:659:VAL:HG22	1.87	0.57
1:B:604:VAL:O	1:B:608:GLN:HG3	2.04	0.57
1:B:367:PHE:HE2	1:B:695:TYR:CE1	2.23	0.57
1:B:212:THR:O	1:B:216:ILE:HD12	2.05	0.56
1:A:230:ILE:HD11	1:A:250:ILE:HG23	1.87	0.56
1:A:604:VAL:O	1:A:608:GLN:HG3	2.05	0.56
1:B:459:ARG:HD2	1:B:485:VAL:HG21	1.87	0.56
1:B:199:LEU:O	1:B:203:THR:HG23	2.05	0.56
1:B:381:PRO:HB3	1:B:683:ASN:HD21	1.71	0.56
1:B:555:LYS:HB2	1:B:556:PRO:HD3	1.87	0.56
1:A:436:VAL:HA	1:A:546:VAL:HG12	1.87	0.56
1:B:457:GLU:HB2	1:B:468:ILE:CD1	2.34	0.56
1:B:389:PRO:O	1:B:393:THR:HG23	2.06	0.56
1:B:438:ASP:OD2	1:B:529:LYS:NZ	2.39	0.55
1:A:433:LYS:NZ	1:A:466:GLU:OE1	2.40	0.55
1:A:424:ASP:O	1:A:428:THR:HB	2.07	0.55
1:B:260:ASP:HB2	1:B:307:ARG:HB2	1.89	0.55
1:A:407:LYS:NZ	1:A:644:GLU:O	2.38	0.55
1:A:504:ARG:O	1:A:508:ASP:HB2	2.07	0.55
1:B:411:ALA:HB2	1:B:647:ASP:HB3	1.88	0.55
1:A:330:PRO:HD3	1:A:410:ASP:OD1	2.07	0.54
1:A:468:ILE:HD12	1:A:469:VAL:N	2.23	0.54
1:B:204:LEU:HD11	1:B:716:VAL:HG12	1.89	0.54
1:A:619:GLY:HA2	1:A:626:LEU:CD1	2.38	0.54
1:B:712:ALA:O	1:B:716:VAL:HG13	2.08	0.54
1:B:156:LEU:HD23	1:B:343:ILE:HG22	1.90	0.54
1:A:162:TYR:HE1	1:A:380:CYS:HB3	1.73	0.54
1:A:466:GLU:O	1:A:470:LYS:HB2	2.08	0.54
1:A:471:LYS:NZ	1:A:474:GLU:OE2	2.33	0.53
1:A:430:THR:HB	1:A:548:ASP:OD2	2.08	0.53
1:B:173:SER:HA	1:B:185:TYR:O	2.08	0.53
1:B:112:LEU:HD11	1:B:131:ILE:HB	1.90	0.53
1:A:462:HIS:CG	1:A:463:PRO:HD2	2.43	0.53
1:B:414:VAL:HG11	1:B:630:ASP:O	2.08	0.53
1:B:450:LEU:HD22	1:B:477:ILE:HG13	1.90	0.53
1:A:484:LYS:HB3	1:A:495:ALA:HA	1.91	0.53
1:A:458:ARG:HA	1:A:469:VAL:HG11	1.91	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:LYS:HE2	1:A:433:LYS:HA	1.90	0.52
1:A:449:LEU:HD13	1:A:542:GLY:HA3	1.91	0.52
1:B:576:TRP:HD1	1:B:594:GLU:OE2	1.93	0.52
1:A:263:VAL:HG12	1:A:288:ASP:O	2.10	0.52
1:B:434:PRO:HG3	1:B:463:PRO:HB2	1.91	0.52
1:B:330:PRO:HD3	1:B:410:ASP:OD1	2.11	0.51
1:B:447:ARG:NH2	1:B:476:GLY:O	2.43	0.51
1:B:686:LEU:HA	1:B:689:ALA:HB3	1.93	0.51
1:B:689:ALA:O	1:B:694:LEU:HD21	2.10	0.51
1:A:521:LEU:O	1:A:525:GLU:HG3	2.11	0.51
1:B:471:LYS:HD2	1:B:475:HIS:ND1	2.25	0.51
1:B:694:LEU:HD22	1:B:702:PHE:CE1	2.44	0.51
1:A:187:PHE:HB3	1:A:190:THR:OG1	2.10	0.51
1:A:211:ARG:HD2	1:A:390:THR:HG23	1.93	0.51
1:A:230:ILE:HD11	1:A:250:ILE:CG2	2.41	0.51
1:B:245:VAL:HG23	1:B:308:VAL:HG12	1.92	0.51
1:B:449:LEU:HD13	1:B:542:GLY:HA3	1.93	0.51
1:A:127:VAL:O	1:A:131:ILE:HG13	2.11	0.50
1:B:367:PHE:HE2	1:B:695:TYR:HE1	1.59	0.50
1:A:712:ALA:O	1:A:716:VAL:HG13	2.10	0.50
1:A:464:ILE:O	1:A:468:ILE:HG13	2.09	0.50
1:B:161:MET:HE1	1:B:376:LEU:HD22	1.92	0.50
1:B:633:ILE:HG23	1:B:650:LEU:HD23	1.93	0.50
1:B:171:LEU:HA	1:B:174:VAL:HG12	1.93	0.50
1:B:431:LYS:HG2	1:B:549:THR:O	2.12	0.50
1:B:652:ARG:HD3	1:B:657:ASP:OD2	2.11	0.50
1:B:492:GLY:HA3	1:B:505:LEU:HG	1.93	0.50
1:B:211:ARG:HD2	1:B:390:THR:HG23	1.93	0.49
1:A:96:SER:HA	1:A:99:LYS:HZ2	1.75	0.49
1:A:100:ARG:O	1:A:104:VAL:HG23	2.12	0.49
1:A:85:GLN:H	1:A:85:GLN:CD	2.15	0.49
1:B:434:PRO:HG3	1:B:463:PRO:CB	2.42	0.49
1:B:671:LYS:HE3	1:B:724:LEU:HD22	1.93	0.49
1:B:172:ALA:HB1	1:B:185:TYR:HD2	1.78	0.49
1:A:128:GLN:HB3	1:A:188:TYR:CD2	2.47	0.49
1:B:124:GLU:O	1:B:128:GLN:HG2	2.12	0.49
1:B:162:TYR:CD2	1:B:198:LEU:HB2	2.48	0.49
1:B:290:VAL:HG21	1:B:303:ILE:HG12	1.95	0.49
1:B:367:PHE:CE2	1:B:701:VAL:HG21	2.48	0.49
1:A:734:ARG:HE	1:A:735:GLY:H	1.58	0.48
1:B:465:ALA:O	1:B:468:ILE:HG13	2.12	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:389:PRO:O	1:A:393:THR:HG23	2.12	0.48
1:A:335:ALA:HA	1:A:338:VAL:HG22	1.96	0.48
1:A:378:VAL:HG23	1:A:707:ALA:HB1	1.95	0.48
1:A:392:LEU:O	1:A:396:MET:HG3	2.12	0.48
1:A:190:THR:HG22	1:A:705:GLU:OE1	2.14	0.48
1:A:652:ARG:HD3	1:A:657:ASP:OD2	2.13	0.48
1:B:439:LEU:HD23	1:B:439:LEU:HA	1.73	0.48
1:B:289:GLU:OE2	1:B:307:ARG:NH2	2.45	0.48
1:A:199:LEU:O	1:A:203:THR:HG23	2.13	0.48
1:A:343:ILE:HA	1:A:346:VAL:HG22	1.96	0.48
1:A:482:PRO:CB	1:A:496:ASP:HB2	2.44	0.48
1:A:633:ILE:HG23	1:A:650:LEU:HD23	1.95	0.48
1:B:112:LEU:HD21	1:B:132:ALA:HB2	1.96	0.48
1:B:530:THR:HB	1:B:546:VAL:CG1	2.44	0.47
1:A:334:LEU:O	1:A:338:VAL:HG13	2.13	0.47
1:A:224:ALA:HB2	1:A:256:LYS:CD	2.44	0.47
1:A:230:ILE:HB	1:A:248:ILE:HB	1.96	0.47
1:A:387:ALA:HB1	1:A:675:ASN:HB3	1.95	0.47
1:B:381:PRO:HA	1:B:711:MET:HE3	1.96	0.47
1:A:450:LEU:HD21	1:A:472:ALA:HA	1.95	0.47
1:A:681:ILE:HA	1:A:684:VAL:HG22	1.94	0.47
1:B:260:ASP:OD1	1:B:292:GLY:N	2.39	0.47
1:B:434:PRO:HG2	1:B:467:ALA:HB2	1.96	0.47
1:B:458:ARG:NH1	1:B:481:GLU:OE2	2.47	0.47
1:A:459:ARG:HD2	1:A:485:VAL:HG21	1.96	0.47
1:B:216:ILE:O	1:B:220:VAL:HG23	2.14	0.47
1:B:386:LEU:HD12	1:B:715:SER:HB3	1.97	0.47
1:A:671:LYS:CE	1:A:724:LEU:HD22	2.44	0.47
1:B:367:PHE:CE2	1:B:695:TYR:HE1	2.32	0.47
1:B:421:VAL:HG22	1:B:614:ALA:HB3	1.96	0.47
1:B:714:SER:O	1:B:718:VAL:HG13	2.14	0.47
1:B:260:ASP:OD2	1:B:313:LEU:N	2.48	0.47
1:B:439:LEU:CD1	1:B:446:GLU:HB3	2.44	0.47
1:B:454:ALA:HA	1:B:468:ILE:CD1	2.45	0.47
1:A:453:ALA:HB2	1:A:534:VAL:HG11	1.97	0.47
1:B:537:ASN:OD1	1:B:537:ASN:N	2.48	0.47
1:A:176:SER:HB2	1:A:185:TYR:HB2	1.97	0.46
1:A:732:ILE:HG13	1:A:732:ILE:O	2.15	0.46
1:B:367:PHE:CE2	1:B:695:TYR:CE1	3.02	0.46
1:B:499:LEU:HD23	1:B:499:LEU:HA	1.78	0.46
1:A:97:ARG:O	1:A:101:LYS:HB2	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:451:ARG:HG2	1:A:477:ILE:HG23	1.97	0.46
1:A:367:PHE:CE2	1:A:695:TYR:HE1	2.34	0.46
1:A:455:ILE:HD12	1:A:480:GLY:H	1.79	0.46
1:B:217:LYS:NZ	1:B:220:VAL:HG21	2.30	0.46
1:A:357:TRP:CE2	1:A:367:PHE:HE1	2.33	0.46
1:A:224:ALA:HB2	1:A:256:LYS:HD2	1.97	0.46
1:B:231:ARG:O	1:B:234:LYS:HG2	2.16	0.46
1:A:121:LEU:HB3	1:A:122:PRO:HD2	1.98	0.46
1:A:458:ARG:HD3	1:A:481:GLU:OE1	2.16	0.46
1:A:243:VAL:O	1:A:308:VAL:HG21	2.16	0.46
1:B:354:PHE:HB2	1:B:372:LEU:HB2	1.97	0.46
1:A:431:LYS:HG3	1:A:549:THR:O	2.16	0.46
1:A:617:GLY:O	1:A:634:ALA:HA	2.16	0.46
1:B:263:VAL:HG13	1:B:285:SER:O	2.17	0.45
1:B:694:LEU:HD22	1:B:702:PHE:CD1	2.52	0.45
1:A:151:LEU:HD11	1:A:160:VAL:HG13	1.99	0.45
1:A:189:GLU:O	1:A:193:LEU:HG	2.16	0.45
1:A:687:ILE:HG23	1:A:688:PRO:CD	2.47	0.45
1:B:346:VAL:HG23	1:B:379:ALA:HB1	1.99	0.45
1:A:211:ARG:HA	1:A:214:GLU:CD	2.37	0.45
1:A:484:LYS:CB	1:A:496:ASP:H	2.28	0.45
1:A:698:PHE:HE1	1:A:701:VAL:HG22	1.82	0.45
1:A:169:ALA:HB3	1:A:191:SER:HB3	1.98	0.45
1:A:415:ALA:HA	1:A:418:VAL:HG23	1.98	0.45
1:B:484:LYS:HB3	1:B:496:ASP:H	1.81	0.45
1:B:436:VAL:HG21	1:B:471:LYS:HG2	1.99	0.45
1:A:328:LYS:HG2	1:A:332:GLN:OE1	2.16	0.45
1:A:465:ALA:O	1:A:469:VAL:HG22	2.17	0.45
1:B:357:TRP:CG	1:B:367:PHE:HE1	2.34	0.45
1:B:529:LYS:HD3	1:B:547:SER:HB2	1.98	0.45
1:A:163:SER:HA	1:A:166:VAL:HG12	1.98	0.44
1:A:537:ASN:OD1	1:A:537:ASN:N	2.48	0.44
1:A:613:VAL:O	1:A:630:ASP:HB2	2.17	0.44
1:A:222:LEU:HB3	1:A:293:ALA:HB1	1.99	0.44
1:B:551:LYS:HZ2	1:B:653:ASP:HB3	1.81	0.44
1:A:685:ILE:HD13	1:A:685:ILE:HA	1.83	0.44
1:B:97:ARG:NH2	1:B:142:SER:OG	2.50	0.44
1:B:357:TRP:CD2	1:B:367:PHE:CE1	2.99	0.44
1:A:459:ARG:CD	1:A:485:VAL:HG21	2.47	0.44
1:B:334:LEU:O	1:B:338:VAL:HG13	2.17	0.44
1:A:697:ILE:HG13	1:A:697:ILE:O	2.17	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345:THR:O	1:B:349:VAL:HG13	2.17	0.44
1:A:210:SER:O	1:A:214:GLU:HG3	2.18	0.44
1:A:528:ALA:HB1	1:A:582:ILE:HD11	1.99	0.44
1:B:428:THR:HG21	1:B:617:GLY:HA2	2.00	0.44
1:A:307:ARG:HD2	1:A:311:GLU:O	2.18	0.44
1:B:500:VAL:HG22	1:B:534:VAL:HG12	1.99	0.44
1:B:687:ILE:HG13	1:B:688:PRO:HD3	2.00	0.44
1:A:230:ILE:HD13	1:A:230:ILE:HG21	1.74	0.44
1:B:216:ILE:HG23	1:B:318:VAL:HG13	1.98	0.43
1:A:207:ARG:HH11	1:A:211:ARG:HH22	1.66	0.43
1:A:375:VAL:HG21	1:A:691:ALA:HB2	1.99	0.43
1:B:462:HIS:CG	1:B:463:PRO:HD2	2.53	0.43
1:B:484:LYS:HB3	1:B:495:ALA:HA	2.00	0.43
1:A:162:TYR:CE1	1:A:380:CYS:HB3	2.53	0.43
1:A:217:LYS:HA	1:A:217:LYS:HD3	1.77	0.43
1:A:705:GLU:CD	1:A:705:GLU:H	2.21	0.43
1:B:387:ALA:HB1	1:B:675:ASN:HB3	1.99	0.43
1:B:443:ASN:OD1	1:B:444:GLY:N	2.51	0.43
1:B:715:SER:O	1:B:719:VAL:HG23	2.19	0.43
1:A:205:GLU:CA	1:A:386:LEU:HD21	2.48	0.43
1:A:715:SER:O	1:A:719:VAL:HG23	2.19	0.43
1:B:346:VAL:HA	1:B:349:VAL:HG22	2.00	0.43
1:B:425:LYS:O	1:B:430:THR:HG23	2.18	0.43
1:A:133:LEU:HB3	1:A:134:PRO:HD3	2.01	0.43
1:A:124:GLU:O	1:A:128:GLN:HG3	2.18	0.43
1:A:159:ASP:OD2	1:A:202:ARG:NH2	2.52	0.43
1:A:457:GLU:HB2	1:A:468:ILE:HD11	2.01	0.43
1:A:714:SER:O	1:A:718:VAL:HG13	2.18	0.43
1:B:123:TYR:HE1	1:B:126:PHE:HB2	1.83	0.43
1:B:439:LEU:HD23	1:B:544:ILE:HG23	2.01	0.43
1:B:700:VAL:HG12	1:B:700:VAL:O	2.19	0.43
1:A:329:PRO:HD2	1:A:332:GLN:CB	2.48	0.42
1:B:378:VAL:HG23	1:B:707:ALA:HB1	2.01	0.42
1:A:190:THR:C	1:A:192:VAL:H	2.21	0.42
1:A:446:GLU:N	1:A:446:GLU:OE1	2.52	0.42
1:B:216:ILE:HG12	1:B:322:GLU:HG2	2.01	0.42
1:B:245:VAL:HG23	1:B:308:VAL:CG1	2.49	0.42
1:A:284:LYS:HE3	1:A:290:VAL:HG12	2.02	0.42
1:B:263:VAL:HG21	1:B:301:LEU:HD22	2.00	0.42
1:A:439:LEU:HD13	1:A:446:GLU:HB3	2.01	0.42
1:B:415:ALA:HA	1:B:418:VAL:HG23	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:LEU:HD11	1:B:160:VAL:HG13	2.02	0.42
1:B:204:LEU:HD23	1:B:204:LEU:HA	1.83	0.42
1:B:236:ILE:HG12	1:B:238:VAL:HG12	2.02	0.42
1:B:434:PRO:HA	1:B:548:ASP:CG	2.40	0.42
1:B:499:LEU:HD12	1:B:511:VAL:HG21	2.01	0.42
1:B:335:ALA:HA	1:B:338:VAL:HG22	2.02	0.42
1:A:236:ILE:HG12	1:A:238:VAL:HG12	2.02	0.42
1:A:308:VAL:H	1:A:311:GLU:HB2	1.85	0.42
1:A:471:LYS:HD2	1:A:475:HIS:ND1	2.34	0.42
1:A:699:GLY:O	1:A:701:VAL:N	2.53	0.42
1:B:199:LEU:HD23	1:B:199:LEU:HA	1.76	0.41
1:B:513:VAL:O	1:B:513:VAL:HG13	2.20	0.41
1:A:513:VAL:O	1:A:513:VAL:HG13	2.20	0.41
1:A:576:TRP:HD1	1:A:594:GLU:OE2	2.03	0.41
1:A:715:SER:O	1:A:718:VAL:HG22	2.20	0.41
1:B:343:ILE:HA	1:B:346:VAL:HG22	2.02	0.41
1:B:687:ILE:N	1:B:688:PRO:HD2	2.35	0.41
1:A:158:MET:HB3	1:A:202:ARG:NH1	2.34	0.41
1:B:441:PRO:HB2	1:B:443:ASN:OD1	2.20	0.41
1:A:190:THR:O	1:A:192:VAL:N	2.52	0.41
1:A:451:ARG:O	1:A:455:ILE:HG12	2.21	0.41
1:B:367:PHE:O	1:B:371:THR:HG23	2.20	0.41
1:B:384:PHE:HE2	1:B:680:LEU:HG	1.85	0.41
1:B:428:THR:CG2	1:B:617:GLY:HA2	2.51	0.41
1:A:120:SER:O	1:A:121:LEU:HD23	2.20	0.41
1:B:381:PRO:HA	1:B:711:MET:CE	2.50	0.41
1:B:652:ARG:NH1	1:B:657:ASP:OD2	2.48	0.41
1:A:98:MET:HB3	1:A:203:THR:HG22	2.03	0.41
1:A:204:LEU:HA	1:A:204:LEU:HD23	1.75	0.41
1:A:468:ILE:HD12	1:A:469:VAL:H	1.86	0.41
1:B:127:VAL:O	1:B:131:ILE:HG13	2.20	0.41
1:B:529:LYS:HG2	1:B:547:SER:HA	2.03	0.41
1:A:471:LYS:HD2	1:A:475:HIS:CE1	2.55	0.41
1:B:531:ALA:HA	1:B:544:ILE:O	2.21	0.41
1:A:199:LEU:HA	1:A:199:LEU:HD23	1.84	0.41
1:A:375:VAL:CG2	1:A:691:ALA:HB2	2.51	0.41
1:A:156:LEU:HB3	1:A:343:ILE:HG21	2.03	0.41
1:A:169:ALA:HB1	1:A:187:PHE:HB3	2.03	0.41
1:A:384:PHE:HE2	1:A:680:LEU:HG	1.86	0.41
1:B:375:VAL:HG22	1:B:690:ALA:HB3	2.02	0.41
1:B:656:ARG:HG2	1:B:732:ILE:CG1	2.50	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:689:ALA:HA	1:B:694:LEU:CD2	2.51	0.41
1:B:365:LEU:HD12	1:B:365:LEU:O	2.21	0.41
1:B:470:LYS:NZ	1:B:474:GLU:OE1	2.35	0.41
1:A:260:ASP:OD1	1:A:292:GLY:N	2.32	0.40
1:A:443:ASN:OD1	1:A:444:GLY:N	2.54	0.40
1:B:381:PRO:C	1:B:383:ALA:H	2.23	0.40
1:A:700:VAL:O	1:A:700:VAL:HG12	2.21	0.40
1:B:211:ARG:HA	1:B:214:GLU:CD	2.42	0.40
1:B:692:GLY:O	1:B:694:LEU:HG	2.21	0.40
1:A:357:TRP:HZ2	1:A:692:GLY:HA3	1.87	0.40
1:B:619:GLY:HA2	1:B:626:LEU:CD1	2.52	0.40
1:A:414:VAL:O	1:A:418:VAL:HG23	2.21	0.40
1:B:126:PHE:O	1:B:129:LEU:HG	2.22	0.40
1:B:184:GLU:O	1:B:186:SER:N	2.53	0.40
1:B:701:VAL:O	1:B:701:VAL:HG12	2.21	0.40
1:A:238:VAL:HG23	1:A:239:PRO:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	652/658 (99%)	618 (95%)	32 (5%)	2 (0%)	41 75
1	B	650/658 (99%)	610 (94%)	35 (5%)	5 (1%)	19 58
All	All	1302/1316 (99%)	1228 (94%)	67 (5%)	7 (0%)	29 67

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	693	LEU
1	A	700	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	434	PRO
1	B	702	PHE
1	A	434	PRO
1	B	224	ALA
1	B	701	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	521/523 (100%)	506 (97%)	15 (3%)	42 65
1	B	519/523 (99%)	506 (98%)	13 (2%)	47 69
All	All	1040/1046 (99%)	1012 (97%)	28 (3%)	44 67

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	TYR
1	A	148	PHE
1	A	211	ARG
1	A	225	LYS
1	A	333	ARG
1	A	410	ASP
1	A	462	HIS
1	A	471	LYS
1	A	486	GLU
1	A	504	ARG
1	A	610	LYS
1	A	682	TYR
1	A	703	ARG
1	A	706	PHE
1	A	711	MET
1	B	123	TYR
1	B	148	PHE
1	B	232	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	328	LYS
1	B	410	ASP
1	B	471	LYS
1	B	486	GLU
1	B	504	ARG
1	B	610	LYS
1	B	637	SER
1	B	682	TYR
1	B	694	LEU
1	B	711	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	683	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	654/658 (99%)	-0.22	9 (1%) 75 65	134, 192, 275, 295	0
1	B	652/658 (99%)	-0.35	7 (1%) 80 72	139, 196, 273, 295	0
All	All	1306/1316 (99%)	-0.28	16 (1%) 79 70	134, 195, 274, 295	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	87	ALA	8.7
1	B	86	ALA	6.0
1	A	544	ILE	5.7
1	A	531	ALA	4.3
1	A	439	LEU	4.1
1	A	545	ALA	4.1
1	A	488	ILE	3.2
1	A	543	ILE	2.9
1	B	483	GLU	2.8
1	B	459	ARG	2.6
1	B	233	GLY	2.5
1	A	534	VAL	2.5
1	B	85	GLN	2.4
1	A	538	GLY	2.3
1	B	235	GLU	2.2
1	A	434	PRO	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.