



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 07:42 AM EDT

PDB ID : 7BA2
Title : D319A mutant of the PilB minor pilin from *Streptococcus sanguinis*
Authors : Pelicic, V.; Sheppard, D.
Deposited on : 2020-12-15
Resolution : 3.00 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
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

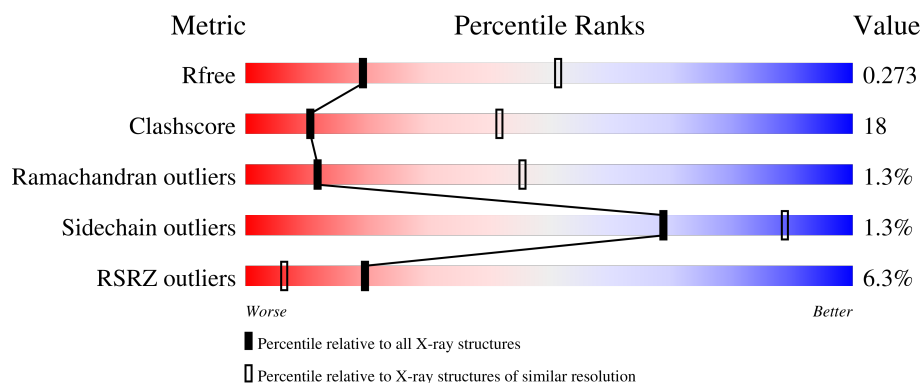
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	426	<div> <div>5%</div> <div>63%</div> <div>34%</div> <div>..</div> </div>
1	B	426	<div> <div>7%</div> <div>69%</div> <div>28%</div> <div>..</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6561 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 Type IV pilus biogenesis protein PilB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	0	0	0
			3276	2061	565	640	10			
1	B	420	Total	C	N	O	S	0	0	0
			3285	2066	566	643	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	319	ALA	ASP	engineered mutation	UNP A0A0B7GP99
B	319	ALA	ASP	engineered mutation	UNP A0A0B7GP99

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	120.95Å 120.95Å 151.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.31 – 3.00 61.31 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (61.31-3.00) 99.5 (61.31-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.17.1	Depositor
R, R_{free}	0.241 , 0.283 0.251 , 0.273	Depositor DCC
R_{free} test set	1237 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	99.3	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 61.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.439 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6561	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% 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	0.30	0/3329	0.66	2/4489 (0.0%)
1	B	0.32	0/3338	0.69	3/4501 (0.1%)
All	All	0.31	0/6667	0.67	5/8990 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	LEU	CA-CB-CG	8.57	135.02	115.30
1	A	213	LEU	CA-CB-CG	6.75	130.83	115.30
1	B	128	LEU	CA-CB-CG	6.29	129.75	115.30
1	B	111	TRP	CA-CB-CG	5.77	124.66	113.70
1	A	460	GLY	C-N-CD	5.18	139.28	128.40

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	3276	0	3276	144	0
1	B	3285	0	3282	97	0
All	All	6561	0	6558	241	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 18.

All (241) 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:149:LYS:NZ	1:A:154:ASN:HA	1.04	1.36
1:A:86:GLU:OE1	1:A:187:ARG:CZ	1.72	1.35
1:A:149:LYS:CE	1:A:153:THR:O	1.75	1.32
1:A:149:LYS:HE3	1:A:153:THR:O	1.09	1.26
1:A:149:LYS:NZ	1:A:154:ASN:CA	2.01	1.23
1:A:149:LYS:HB2	1:A:156:LYS:HG2	1.23	1.17
1:A:149:LYS:HZ2	1:A:154:ASN:CA	1.56	1.15
1:A:86:GLU:OE1	1:A:187:ARG:NH1	1.78	1.14
1:A:86:GLU:OE1	1:A:187:ARG:NE	1.88	1.05
1:A:149:LYS:CB	1:A:156:LYS:HG2	1.91	1.00
1:B:97:LYS:HG2	1:B:124:MET:O	1.66	0.95
1:A:86:GLU:CD	1:A:187:ARG:CZ	2.38	0.91
1:A:149:LYS:HB2	1:A:156:LYS:CG	2.00	0.90
1:A:86:GLU:OE1	1:A:187:ARG:CD	2.22	0.88
1:A:149:LYS:HZ1	1:A:154:ASN:CA	1.77	0.87
1:A:149:LYS:HZ1	1:A:154:ASN:HA	1.04	0.85
1:B:240:ASP:HB3	1:B:448:LYS:HE2	1.57	0.84
1:A:99:VAL:HG12	1:A:116:LEU:HD12	1.61	0.82
1:A:86:GLU:OE1	1:A:187:ARG:HD2	1.79	0.81
1:A:86:GLU:CD	1:A:187:ARG:NH1	2.33	0.81
1:B:71:LEU:HD11	1:B:185:ALA:HB2	1.63	0.79
1:A:187:ARG:NH1	1:A:191:ILE:HA	1.97	0.78
1:B:83:LEU:HD11	1:B:111:TRP:CZ2	2.19	0.77
1:B:305:GLN:NE2	1:B:395:LEU:O	2.17	0.76
1:A:212:ASP:OD2	1:A:216:ARG:HB2	1.87	0.75
1:B:78:GLY:HA2	1:B:93:SER:HB2	1.69	0.75
1:B:135:ALA:HA	1:B:141:LEU:HG	1.69	0.74
1:B:43:ARG:NH1	1:B:122:TYR:HE2	1.85	0.73
1:A:194:GLN:HG3	1:A:459:SER:HB2	1.71	0.72
1:B:104:TRP:HE3	1:B:111:TRP:CE3	2.07	0.71
1:A:68:ILE:HB	1:A:171:VAL:HG23	1.71	0.71
1:A:104:TRP:CD1	1:A:106:LYS:HG2	2.26	0.71
1:B:75:LYS:NZ	1:B:461:PRO:O	2.23	0.70
1:A:233:LYS:HD2	1:A:441:GLN:HB2	1.74	0.69
1:A:93:SER:HB3	1:A:98:LYS:O	1.92	0.69
1:A:201:PHE:HE2	1:A:238:ILE:HD11	1.57	0.69
1:A:334:ALA:O	1:A:336:ASN:N	2.21	0.69
1:A:268:LEU:HB3	1:A:272:THR:HG22	1.75	0.69
1:A:439:ALA:O	1:A:443:THR:HG23	1.94	0.68
1:A:149:LYS:NZ	1:A:153:THR:O	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LYS:HB2	1:A:156:LYS:HE2	1.77	0.67
1:B:156:LYS:HG2	1:B:156:LYS:O	1.96	0.66
1:B:176:ALA:HB3	1:B:179:LYS:HZ3	1.60	0.66
1:A:217:ASN:OD1	1:A:218:VAL:N	2.29	0.65
1:B:104:TRP:HE3	1:B:111:TRP:HE3	1.43	0.65
1:A:92:LEU:HG	1:A:128:LEU:HB3	1.80	0.64
1:A:52:ARG:HE	1:A:159:ILE:HG23	1.64	0.63
1:A:110:ASP:OD1	1:A:111:TRP:N	2.32	0.62
1:B:445:SER:HA	1:B:448:LYS:HD2	1.81	0.62
1:A:199:VAL:HB	1:A:249:VAL:HG12	1.81	0.62
1:A:86:GLU:OE2	1:A:187:ARG:NH1	2.32	0.62
1:B:249:VAL:HG23	1:B:275:ILE:HD13	1.82	0.62
1:B:135:ALA:HB1	1:B:139:ASN:O	1.99	0.62
1:A:66:THR:HA	1:A:168:THR:HG21	1.81	0.62
1:A:187:ARG:NH2	1:A:192:GLN:CD	2.54	0.61
1:A:44:GLU:HG2	1:A:47:ILE:HD11	1.81	0.61
1:B:150:TYR:HE1	1:B:157:LEU:HD13	1.66	0.61
1:A:51:MET:HB2	1:A:159:ILE:HD11	1.82	0.61
1:B:199:VAL:HB	1:B:249:VAL:HG12	1.83	0.60
1:A:149:LYS:HG2	1:A:150:TYR:H	1.67	0.60
1:A:149:LYS:HG2	1:A:150:TYR:N	2.16	0.60
1:A:149:LYS:HZ2	1:A:154:ASN:HA	0.77	0.59
1:B:251:LEU:HG	1:B:261:ILE:HB	1.84	0.59
1:B:227:MET:O	1:B:231:ARG:HG3	2.03	0.59
1:A:93:SER:HB3	1:A:100:MET:HG2	1.84	0.58
1:A:399:ARG:NH1	1:A:428:GLU:OE1	2.37	0.58
1:A:408:VAL:HG12	1:A:411:GLU:H	1.68	0.58
1:B:213:LEU:HD23	1:B:226:ARG:HG2	1.86	0.58
1:A:54:SER:O	1:A:58:VAL:HG23	2.03	0.57
1:A:103:VAL:O	1:A:111:TRP:HA	2.04	0.57
1:B:237:MET:HA	1:B:448:LYS:HE3	1.86	0.57
1:A:304:LEU:HD13	1:A:312:LYS:HB3	1.87	0.57
1:B:102:TYR:HB3	1:B:111:TRP:HE1	1.70	0.57
1:A:86:GLU:HA	1:A:104:TRP:CD1	2.40	0.57
1:A:217:ASN:HB3	1:A:220:LYS:HB2	1.86	0.57
1:B:208:SER:O	1:B:226:ARG:NH1	2.34	0.56
1:B:68:ILE:HD13	1:B:166:LEU:HD13	1.87	0.56
1:A:142:ILE:HG13	1:A:166:LEU:HD21	1.87	0.56
1:B:240:ASP:HB3	1:B:448:LYS:CE	2.32	0.56
1:B:197:ILE:HB	1:B:247:ILE:HG13	1.87	0.56
1:B:80:LYS:HA	1:B:83:LEU:HG	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:TRP:CE3	1:B:187:ARG:HG3	2.40	0.55
1:B:62:VAL:HG13	1:B:166:LEU:HD21	1.88	0.55
1:A:187:ARG:HH22	1:A:192:GLN:CD	2.10	0.55
1:B:314:ILE:HB	1:B:400:VAL:HG22	1.88	0.55
1:A:101:ASN:HB2	1:A:116:LEU:HD21	1.89	0.55
1:A:149:LYS:HB2	1:A:156:LYS:CE	2.37	0.54
1:B:72:ASP:OD2	1:B:179:LYS:HB3	2.08	0.54
1:A:129:GLU:OE1	1:A:131:LYS:HE3	2.07	0.53
1:A:80:LYS:HD3	1:A:111:TRP:HB2	1.90	0.53
1:A:68:ILE:HG13	1:A:166:LEU:HD13	1.90	0.53
1:A:71:LEU:O	1:A:181:GLY:HA2	2.08	0.53
1:A:97:LYS:HG2	1:A:124:MET:O	2.09	0.53
1:B:76:PHE:CZ	1:B:91:GLY:HA3	2.44	0.53
1:B:87:TRP:CD2	1:B:187:ARG:HD2	2.44	0.53
1:A:80:LYS:HA	1:A:83:LEU:HG	1.91	0.52
1:A:149:LYS:HZ1	1:A:154:ASN:CB	2.21	0.52
1:A:187:ARG:NH2	1:A:192:GLN:OE1	2.41	0.52
1:B:337:ARG:HE	1:B:347:VAL:HG11	1.75	0.52
1:B:72:ASP:CG	1:B:179:LYS:HB3	2.30	0.51
1:B:124:MET:HA	1:B:151:PRO:HD3	1.91	0.51
1:B:228:ASP:HA	1:B:231:ARG:HD2	1.91	0.51
1:A:140:ARG:HD3	1:A:171:VAL:CG1	2.41	0.51
1:B:97:LYS:HD3	1:B:123:ASN:C	2.31	0.51
1:B:97:LYS:HG2	1:B:124:MET:C	2.30	0.51
1:B:213:LEU:HD21	1:B:405:PHE:CZ	2.45	0.51
1:A:135:ALA:HA	1:A:141:LEU:HG	1.92	0.51
1:B:175:VAL:HG22	1:B:179:LYS:HB2	1.92	0.51
1:A:224:GLU:HB2	1:A:229:ILE:HD11	1.91	0.50
1:A:231:ARG:NH1	1:A:279:ILE:O	2.44	0.50
1:A:83:LEU:HD13	1:A:89:TYR:CZ	2.46	0.50
1:A:149:LYS:HE3	1:A:153:THR:C	2.13	0.50
1:B:437:GLU:CD	1:B:437:GLU:H	2.15	0.50
1:A:52:ARG:HG3	1:A:159:ILE:HG12	1.93	0.50
1:A:149:LYS:HB2	1:A:156:LYS:CD	2.42	0.50
1:B:205:THR:OG1	1:B:255:SER:OG	2.30	0.50
1:B:419:THR:OG1	1:B:431:TYR:HB2	2.11	0.49
1:A:278:THR:HA	1:A:281:LYS:HE3	1.94	0.49
1:A:199:VAL:O	1:A:249:VAL:HA	2.13	0.49
1:A:256:THR:HA	1:A:290:VAL:HG23	1.94	0.49
1:B:76:PHE:CE2	1:B:91:GLY:HA3	2.48	0.48
1:B:233:LYS:HG3	1:B:440:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:SER:O	1:A:96:GLY:N	2.46	0.48
1:B:370:ASP:OD2	1:B:374:ARG:N	2.46	0.48
1:B:233:LYS:HD2	1:B:441:GLN:HB2	1.96	0.48
1:B:265:PHE:CZ	1:B:304:LEU:HD23	2.49	0.48
1:B:299:TYR:CE1	1:B:360:LEU:HB3	2.49	0.48
1:A:143:SER:O	1:A:143:SER:OG	2.32	0.48
1:A:187:ARG:CD	1:A:189:ASP:O	2.62	0.48
1:B:55:MET:HE2	1:B:146:LEU:H	1.78	0.47
1:A:195:MET:HE3	1:A:310:GLN:HB2	1.95	0.47
1:A:149:LYS:CA	1:A:156:LYS:HG2	2.43	0.47
1:A:76:PHE:CZ	1:A:91:GLY:HA3	2.49	0.47
1:A:93:SER:CB	1:A:100:MET:HG2	2.44	0.47
1:B:408:VAL:HB	1:B:411:GLU:HG3	1.97	0.47
1:A:72:ASP:HB3	1:A:175:VAL:HG12	1.95	0.47
1:B:116:LEU:HD23	1:B:116:LEU:HA	1.60	0.47
1:A:227:MET:O	1:A:231:ARG:HG3	2.14	0.47
1:A:249:VAL:HG13	1:A:268:LEU:HD11	1.95	0.47
1:A:75:LYS:NZ	1:A:461:PRO:HB2	2.30	0.46
1:B:104:TRP:HA	1:B:111:TRP:CG	2.51	0.46
1:A:195:MET:HE1	1:A:311:LEU:HG	1.98	0.46
1:B:278:THR:HA	1:B:281:LYS:HE3	1.97	0.46
1:A:295:ASP:OD1	1:A:298:ARG:NH2	2.36	0.46
1:B:372:TYR:HB2	1:B:374:ARG:HH11	1.80	0.46
1:A:206:SER:O	1:A:287:PRO:HB3	2.15	0.46
1:A:321:ILE:HG12	1:A:414:TYR:CD2	2.50	0.46
1:B:87:TRP:CE2	1:B:187:ARG:HD2	2.51	0.46
1:B:292:ASN:HB2	1:B:324:ALA:HB3	1.98	0.46
1:B:301:MET:HB3	1:B:393:PHE:O	2.16	0.46
1:A:125:LYS:HE3	1:A:151:PRO:HA	1.98	0.45
1:A:213:LEU:O	1:A:407:GLY:N	2.40	0.45
1:B:87:TRP:CZ2	1:B:460:GLY:HA2	2.50	0.45
1:A:149:LYS:CB	1:A:156:LYS:HE2	2.46	0.45
1:A:72:ASP:OD1	1:A:72:ASP:N	2.49	0.45
1:A:294:GLY:HA3	1:A:385:TYR:CD2	2.52	0.45
1:A:55:MET:SD	1:A:144:TYR:HB2	2.57	0.45
1:A:76:PHE:CD1	1:A:89:TYR:HD2	2.35	0.45
1:B:372:TYR:HD2	1:B:374:ARG:HH12	1.64	0.45
1:A:140:ARG:HG3	1:A:169:LYS:HD2	1.98	0.45
1:A:195:MET:CE	1:A:311:LEU:HG	2.46	0.45
1:A:92:LEU:HD11	1:A:127:ASP:HA	1.98	0.45
1:B:204:ASP:HB2	1:B:318:THR:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ILE:HD12	1:A:452:GLN:HB2	1.99	0.45
1:A:135:ALA:HB1	1:A:139:ASN:O	2.17	0.44
1:A:302:ILE:HD12	1:A:302:ILE:HA	1.88	0.44
1:B:104:TRP:CZ2	1:B:106:LYS:HA	2.52	0.44
1:B:176:ALA:CB	1:B:179:LYS:HZ3	2.30	0.44
1:A:195:MET:HE3	1:A:195:MET:HB2	1.73	0.44
1:B:104:TRP:CE2	1:B:106:LYS:HA	2.52	0.44
1:B:324:ALA:HA	1:B:368:GLY:O	2.17	0.44
1:B:43:ARG:NH1	1:B:122:TYR:CE2	2.76	0.44
1:B:104:TRP:HA	1:B:111:TRP:CD2	2.52	0.44
1:A:301:MET:SD	1:A:390:SER:HA	2.57	0.44
1:A:321:ILE:HG12	1:A:414:TYR:HD2	1.83	0.44
1:B:83:LEU:HD22	1:B:89:TYR:CZ	2.53	0.44
1:B:86:GLU:HA	1:B:104:TRP:CD1	2.53	0.44
1:B:129:GLU:HG2	1:B:130:PHE:N	2.33	0.44
1:B:104:TRP:CE3	1:B:111:TRP:CE3	2.97	0.44
1:A:251:LEU:HG	1:A:261:ILE:HB	2.00	0.44
1:A:399:ARG:HA	1:A:428:GLU:O	2.18	0.44
1:B:297:LEU:HD23	1:B:297:LEU:HA	1.88	0.43
1:A:95:ASP:OD2	1:A:98:LYS:HE3	2.18	0.43
1:A:249:VAL:HG22	1:A:275:ILE:HD13	2.00	0.43
1:A:75:LYS:NZ	1:A:461:PRO:C	2.71	0.43
1:A:301:MET:HB3	1:A:393:PHE:O	2.18	0.43
1:A:327:VAL:O	1:A:366:ARG:NH1	2.46	0.43
1:A:237:MET:O	1:A:241:LEU:HB2	2.19	0.43
1:A:85:ARG:O	1:A:87:TRP:HD1	2.01	0.43
1:B:133:GLU:HG3	1:B:163:ILE:HD11	2.01	0.43
1:B:67:SER:O	1:B:186:TYR:HA	2.19	0.43
1:B:72:ASP:O	1:B:76:PHE:N	2.51	0.42
1:A:116:LEU:HA	1:A:116:LEU:HD23	1.52	0.42
1:A:140:ARG:HD2	1:A:169:LYS:HE3	2.01	0.42
1:A:225:SER:O	1:A:229:ILE:HG13	2.20	0.42
1:A:408:VAL:HG11	1:A:411:GLU:HG3	2.01	0.42
1:B:237:MET:O	1:B:241:LEU:HB2	2.19	0.42
1:A:79:SER:C	1:A:81:GLN:H	2.23	0.42
1:B:92:LEU:C	1:B:92:LEU:HD12	2.40	0.42
1:A:99:VAL:CG1	1:A:116:LEU:HD12	2.42	0.42
1:B:62:VAL:CG1	1:B:166:LEU:HD21	2.48	0.42
1:B:150:TYR:HB2	1:B:153:THR:HG23	2.02	0.42
1:A:189:ASP:OD1	1:A:189:ASP:N	2.37	0.42
1:A:233:LYS:HE3	1:A:437:GLU:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:SER:HB2	1:A:288:ASP:O	2.19	0.42
1:B:121:LEU:H	1:B:121:LEU:HG	1.60	0.42
1:B:237:MET:HA	1:B:448:LYS:CE	2.50	0.42
1:A:392:LYS:HD3	1:A:393:PHE:CE2	2.55	0.42
1:B:173:SER:OG	1:B:175:VAL:HG12	2.20	0.42
1:A:83:LEU:HB3	1:A:111:TRP:CH2	2.54	0.42
1:A:84:THR:O	1:A:111:TRP:HZ2	2.02	0.41
1:A:229:ILE:HG13	1:A:229:ILE:H	1.60	0.41
1:A:337:ARG:CZ	1:A:347:VAL:HG11	2.50	0.41
1:B:166:LEU:HA	1:B:166:LEU:HD23	1.61	0.41
1:B:399:ARG:NE	1:B:428:GLU:OE1	2.40	0.41
1:B:201:PHE:HE2	1:B:238:ILE:HD11	1.85	0.41
1:B:233:LYS:CD	1:B:441:GLN:HB2	2.50	0.41
1:B:309:ALA:O	1:B:312:LYS:HE3	2.20	0.41
1:A:93:SER:O	1:A:95:ASP:N	2.53	0.41
1:B:105:ASN:C	1:B:107:GLN:H	2.23	0.41
1:A:255:SER:O	1:A:256:THR:C	2.59	0.41
1:A:304:LEU:HD11	1:A:313:TYR:O	2.20	0.41
1:B:192:GLN:H	1:B:192:GLN:CD	2.24	0.41
1:B:200:SER:HA	1:B:250:ASN:OD1	2.21	0.41
1:A:177:LYS:HG2	1:A:178:GLY:N	2.33	0.41
1:B:223:ASN:N	1:B:223:ASN:OD1	2.54	0.41
1:A:79:SER:N	1:A:94:ALA:HB2	2.36	0.41
1:A:187:ARG:NE	1:A:189:ASP:O	2.54	0.41
1:A:399:ARG:HG2	1:A:400:VAL:N	2.36	0.41
1:A:159:ILE:HD12	1:A:159:ILE:H	1.86	0.41
1:A:258:ALA:HB2	1:A:295:ASP:HB3	2.02	0.41
1:A:78:GLY:HA3	1:A:94:ALA:HB2	2.02	0.40
1:A:299:TYR:O	1:A:302:ILE:HG22	2.21	0.40
1:B:299:TYR:CZ	1:B:360:LEU:HB3	2.56	0.40
1:A:83:LEU:HD13	1:A:89:TYR:CE2	2.56	0.40
1:A:128:LEU:HD21	1:A:130:PHE:CZ	2.56	0.40
1:B:140:ARG:HD3	1:B:171:VAL:HG22	2.03	0.40
1:B:133:GLU:HG3	1:B:143:SER:HB3	2.03	0.40
1:B:213:LEU:HD23	1:B:213:LEU:HA	1.73	0.40
1:A:148:GLY:O	1:A:156:LYS:HB3	2.21	0.40
1:A:309:ALA:O	1:A:312:LYS:HE3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	417/426 (98%)	372 (89%)	37 (9%)	8 (2%)	8	36
1	B	418/426 (98%)	380 (91%)	35 (8%)	3 (1%)	22	60
All	All	835/852 (98%)	752 (90%)	72 (9%)	11 (1%)	12	45

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	ALA
1	A	334	ALA
1	A	335	GLY
1	B	160	ASP
1	B	80	LYS
1	A	177	LYS
1	A	80	LYS
1	A	93	SER
1	A	156	LYS
1	A	160	ASP
1	B	137	GLN

5.3.2 Protein sidechains

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	356/363 (98%)	354 (99%)	2 (1%)	86	95
1	B	357/363 (98%)	350 (98%)	7 (2%)	55	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	713/726 (98%)	704 (99%)	9 (1%)	69 89

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	MET
1	A	250	ASN
1	B	110	ASP
1	B	114	SER
1	B	152	ASP
1	B	196	ASN
1	B	223	ASN
1	B	241	LEU
1	B	399	ARG

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	145	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	419/426 (98%)	0.50	23 (5%)	25 9	47, 97, 142, 173	0
1	B	420/426 (98%)	0.54	30 (7%)	16 5	50, 96, 143, 168	0
All	All	839/852 (98%)	0.52	53 (6%)	20 6	47, 96, 143, 173	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	150	TYR	4.9
1	B	153	THR	3.8
1	B	155	ASN	3.7
1	B	93	SER	3.5
1	B	154	ASN	3.4
1	B	116	LEU	3.4
1	A	124	MET	3.3
1	B	230	LEU	3.3
1	B	90	ILE	3.2
1	A	97	LYS	3.2
1	A	55	MET	3.0
1	B	128	LEU	2.9
1	B	117	GLY	2.9
1	B	146	LEU	2.9
1	A	238	ILE	2.9
1	A	275	ILE	2.9
1	B	272	THR	2.8
1	B	118	THR	2.8
1	A	150	TYR	2.8
1	A	444	PHE	2.8
1	A	213	LEU	2.8
1	A	251	LEU	2.7
1	A	146	LEU	2.7
1	B	201	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	220	LYS	2.7
1	B	47	ILE	2.6
1	B	68	ILE	2.6
1	B	157	LEU	2.6
1	A	316	LEU	2.5
1	B	149	LYS	2.5
1	A	237	MET	2.5
1	B	55	MET	2.5
1	B	251	LEU	2.4
1	A	69	PHE	2.4
1	B	129	GLU	2.4
1	B	316	LEU	2.4
1	B	237	MET	2.4
1	A	317	LEU	2.3
1	A	51	MET	2.3
1	B	405	PHE	2.2
1	A	159	ILE	2.2
1	B	145	ASN	2.2
1	A	230	LEU	2.2
1	B	76	PHE	2.2
1	A	144	TYR	2.2
1	A	285	LEU	2.1
1	B	112	ASP	2.1
1	B	241	LEU	2.1
1	A	52	ARG	2.1
1	B	203	PHE	2.1
1	A	92	LEU	2.0
1	A	126	LEU	2.0
1	A	234	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.