



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 7, 2025 – 10:07 AM JST

PDB ID : 9JCN / pdb_00009jcn
Title : Crystal structure of Zea mays 3-phosphoglycerate dehydrogenase S282L mutant
Authors : Li, R.; Wang, C.
Deposited on : 2024-08-30
Resolution : 3.48 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 : 4-5-2 with Phenix2.0rc1
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

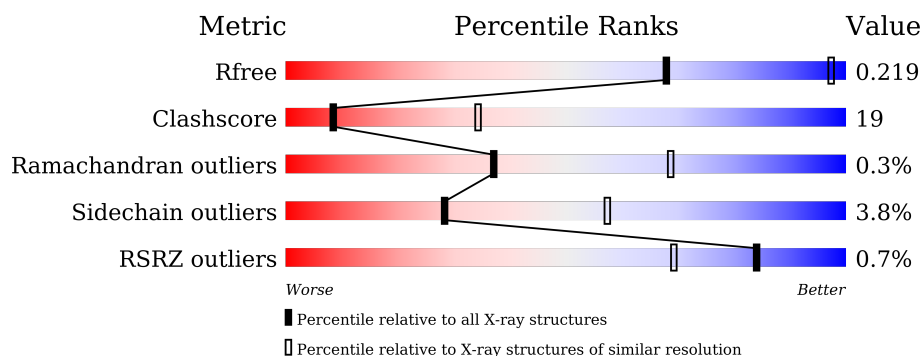
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.48 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}	164625	1099 (3.54-3.42)
Clashscore	180529	1048 (3.52-3.44)
Ramachandran outliers	177936	1033 (3.52-3.44)
Sidechain outliers	177891	1034 (3.52-3.44)
RSRZ outliers	164620	1098 (3.54-3.42)

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	387	
1	B	387	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4432 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 D-3-phosphoglycerate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2216	1390	399	417	10			
1	B	305	Total	C	N	O	S	0	0	0
			2216	1390	399	417	10			

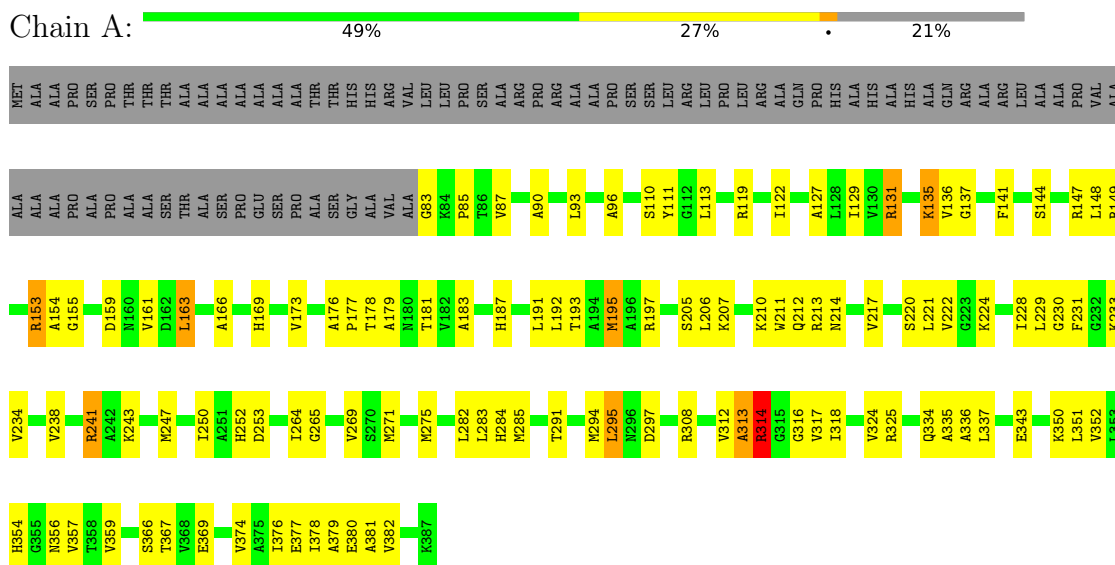
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	282	LEU	SER	engineered mutation	UNP A0A1D6DW07
B	282	LEU	SER	engineered mutation	UNP A0A1D6DW07

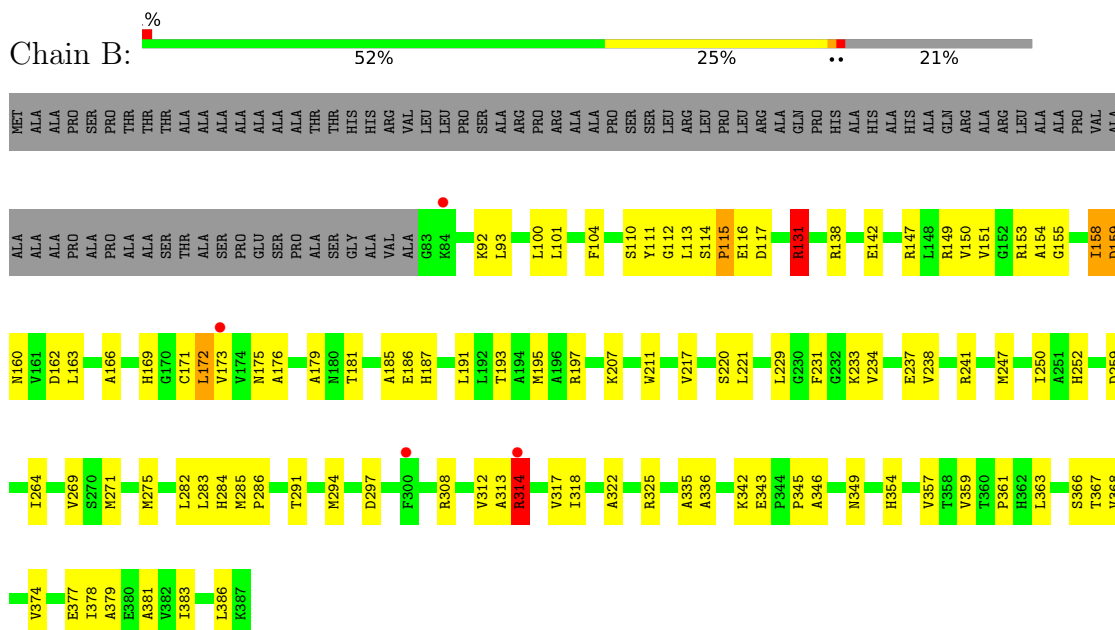
3 Residue-property plots

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: D-3-phosphoglycerate dehydrogenase



• Molecule 1: D-3-phosphoglycerate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	96.34Å 96.34Å 503.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	22.97 – 3.48 22.97 – 3.48	Depositor EDS
% Data completeness (in resolution range)	96.1 (22.97-3.48) 98.9 (22.97-3.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.51 (at 3.96Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: 000)	Depositor
R, R_{free}	0.208 , 0.228 0.194 , 0.219	Depositor DCC
R_{free} test set	566 reflections (4.51%)	wwPDB-VP
Wilson B-factor (Å ²)	114.3	Xtriage
Anisotropy	0.650	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 75.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4432	wwPDB-VP
Average B, all atoms (Å ²)	112.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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

5.1 Standard geometry

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	0.55	0/2242	0.89	4/3037 (0.1%)
1	B	0.57	0/2242	0.91	8/3037 (0.3%)
All	All	0.56	0/4484	0.90	12/6074 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	116	GLU	N-CA-C	8.50	120.55	111.28
1	B	160	ASN	N-CA-C	8.09	121.20	111.82
1	B	163	LEU	N-CA-C	-8.03	103.61	113.41
1	B	162	ASP	N-CA-C	-7.18	99.26	109.96
1	A	163	LEU	N-CA-C	-6.30	104.41	111.28
1	B	259	ASP	N-CA-C	6.15	117.65	111.07
1	A	137	GLY	N-CA-C	-5.77	103.75	112.41
1	A	313	ALA	N-CA-C	5.52	119.59	111.04
1	B	233	LYS	N-CA-C	5.27	117.70	111.33
1	B	112	GLY	N-CA-C	5.24	125.59	113.18
1	B	115	PRO	CB-CA-C	5.09	119.96	111.56
1	A	313	ALA	CB-CA-C	-5.02	103.41	111.39

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	ARG	Sidechain
1	A	314	ARG	Sidechain
1	B	131	ARG	Sidechain
1	B	314	ARG	Sidechain

5.2 Too-close contacts

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	2216	0	2295	92	0
1	B	2216	0	2295	89	0
All	All	4432	0	4590	169	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 19.

All (169) 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:221:LEU:HB3	1:A:247:MET:HE2	1.50	0.89
1:B:234:VAL:HB	1:B:284:HIS:CE1	2.10	0.85
1:B:155:GLY:H	1:B:176:ALA:HB3	1.40	0.85
1:A:87:VAL:HG22	1:A:127:ALA:HB3	1.58	0.83
1:B:110:SER:HB3	1:B:113:LEU:HD21	1.61	0.80
1:B:346:ALA:HB3	1:B:349:ASN:HB2	1.66	0.77
1:A:241:ARG:HH11	1:A:241:ARG:HG2	1.50	0.77
1:B:241:ARG:HH11	1:B:241:ARG:HG2	1.50	0.75
1:B:176:ALA:HB2	1:B:378:ILE:HG23	1.69	0.74
1:B:284:HIS:HA	1:B:313:ALA:HB2	1.71	0.73
1:B:191:LEU:HD22	1:B:336:ALA:HB1	1.68	0.73
1:A:176:ALA:HA	1:A:377:GLU:HG2	1.72	0.71
1:A:93:LEU:HD11	1:A:378:ILE:HD11	1.73	0.70
1:B:297:ASP:OD2	1:B:325:ARG:HD2	1.90	0.70
1:A:187:HIS:HB2	1:B:197:ARG:NH1	2.08	0.69
1:A:178:THR:HG22	1:A:233:LYS:NZ	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:PHE:HZ	1:A:264:ILE:HD11	1.58	0.66
1:A:284:HIS:HA	1:A:313:ALA:HB2	1.77	0.66
1:B:238:VAL:HG11	1:B:282:LEU:CD2	2.27	0.65
1:A:155:GLY:H	1:A:176:ALA:HB3	1.61	0.65
1:B:231:PHE:HZ	1:B:264:ILE:HD11	1.62	0.64
1:B:238:VAL:HG11	1:B:282:LEU:HD21	1.79	0.64
1:A:250:ILE:HD12	1:A:269:VAL:HG21	1.80	0.63
1:A:122:ILE:HD11	1:A:148:LEU:HD22	1.79	0.63
1:B:342:LYS:HG3	1:B:345:PRO:HA	1.82	0.61
1:B:131:ARG:O	1:B:153:ARG:NE	2.35	0.60
1:A:357:VAL:O	1:B:207:LYS:NZ	2.22	0.60
1:A:285:MET:HE3	1:A:317:VAL:HG11	1.83	0.59
1:A:297:ASP:OD2	1:A:325:ARG:HD2	2.02	0.59
1:B:154:ALA:HB2	1:B:378:ILE:HD13	1.86	0.58
1:A:222:VAL:HG23	1:B:186:GLU:OE2	2.03	0.58
1:B:113:LEU:HB3	1:B:117:ASP:HB2	1.85	0.58
1:A:110:SER:HB3	1:A:113:LEU:HD21	1.86	0.57
1:B:291:THR:HA	1:B:294:MET:HE3	1.87	0.57
1:B:150:VAL:HB	1:B:172:LEU:HD12	1.86	0.57
1:B:113:LEU:HD22	1:B:113:LEU:N	2.20	0.56
1:A:191:LEU:HD22	1:A:336:ALA:HB1	1.88	0.56
1:B:176:ALA:HA	1:B:377:GLU:HG2	1.88	0.55
1:A:295:LEU:HB2	1:A:318:ILE:HA	1.87	0.55
1:B:241:ARG:HG2	1:B:241:ARG:NH1	2.22	0.55
1:A:234:VAL:HB	1:A:284:HIS:CE1	2.41	0.55
1:A:228:ILE:HG23	1:A:282:LEU:HD11	1.89	0.54
1:B:252:HIS:CE1	1:B:271:MET:HB2	2.42	0.54
1:A:334:GLN:HB2	1:A:356:ASN:O	2.07	0.54
1:A:316:GLY:C	1:A:318:ILE:N	2.64	0.53
1:A:178:THR:HG22	1:A:233:LYS:HZ3	1.73	0.53
1:A:90:ALA:O	1:A:111:TYR:HA	2.08	0.53
1:A:343:GLU:HG2	1:B:211:TRP:HH2	1.73	0.53
1:B:195:MET:HE3	1:B:336:ALA:HB2	1.90	0.53
1:A:243:LYS:NZ	1:A:265:GLY:O	2.40	0.52
1:A:83:GLY:O	1:A:85:PRO:HD3	2.09	0.52
1:A:187:HIS:HB2	1:B:197:ARG:HH12	1.75	0.52
1:A:159:ASP:N	1:A:159:ASP:OD1	2.41	0.52
1:B:138:ARG:HG2	1:B:142:GLU:OE2	2.09	0.52
1:B:322:ALA:HA	1:B:325:ARG:HE	1.74	0.52
1:A:176:ALA:HB2	1:A:378:ILE:HG23	1.91	0.52
1:A:369:GLU:N	1:A:369:GLU:OE1	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:LEU:HA	1:B:383:ILE:HD11	1.93	0.51
1:A:241:ARG:HG2	1:A:241:ARG:NH1	2.21	0.51
1:A:366:SER:O	1:B:217:VAL:HG13	2.11	0.51
1:A:141:PHE:HB3	1:A:169:HIS:HD2	1.76	0.51
1:A:141:PHE:HB3	1:A:169:HIS:CD2	2.46	0.51
1:B:93:LEU:HD11	1:B:378:ILE:HD11	1.93	0.50
1:B:101:LEU:HD23	1:B:383:ILE:HD11	1.93	0.50
1:A:141:PHE:CZ	1:A:166:ALA:HB2	2.46	0.49
1:B:149:ARG:O	1:B:171:CYS:HA	2.12	0.49
1:B:114:SER:O	1:B:115:PRO:C	2.55	0.49
1:B:179:ALA:HB3	1:B:374:VAL:HG23	1.95	0.49
1:B:153:ARG:O	1:B:175:ASN:HA	2.12	0.49
1:A:177:PRO:HG2	1:A:178:THR:HG23	1.94	0.48
1:B:285:MET:HE3	1:B:317:VAL:HG11	1.96	0.48
1:A:135:LYS:HD3	1:A:135:LYS:HA	1.67	0.48
1:A:220:SER:O	1:A:224:LYS:HE2	2.12	0.48
1:B:237:GLU:O	1:B:241:ARG:HD2	2.13	0.48
1:A:231:PHE:H	1:A:253:ASP:HB2	1.78	0.48
1:A:178:THR:HG22	1:A:233:LYS:HZ1	1.77	0.47
1:A:211:TRP:HH2	1:B:343:GLU:HG2	1.79	0.47
1:B:308:ARG:HG3	1:B:308:ARG:HH11	1.78	0.47
1:B:195:MET:SD	1:B:308:ARG:HB3	2.53	0.47
1:B:252:HIS:CD2	1:B:252:HIS:C	2.92	0.47
1:A:192:LEU:HD21	1:A:247:MET:HE1	1.96	0.47
1:B:335:ALA:O	1:B:357:VAL:HA	2.14	0.47
1:A:131:ARG:HG2	1:A:154:ALA:HB3	1.97	0.47
1:A:176:ALA:CB	1:A:374:VAL:HG13	2.44	0.47
1:A:335:ALA:O	1:A:357:VAL:HA	2.15	0.47
1:A:352:VAL:HA	1:A:359:VAL:HG11	1.96	0.47
1:B:149:ARG:NH2	1:B:386:LEU:O	2.47	0.47
1:A:231:PHE:CZ	1:A:264:ILE:HD11	2.46	0.47
1:B:114:SER:OG	1:B:117:ASP:OD1	2.33	0.46
1:A:197:ARG:NH2	1:B:187:HIS:HB2	2.31	0.46
1:B:322:ALA:O	1:B:325:ARG:HG2	2.15	0.46
1:B:92:LYS:HA	1:B:111:TYR:CE1	2.51	0.46
1:B:172:LEU:H	1:B:172:LEU:HG	1.52	0.46
1:B:142:GLU:HA	1:B:169:HIS:HE1	1.80	0.46
1:B:229:LEU:HD11	1:B:275:MET:HE1	1.98	0.46
1:A:183:ALA:HB1	1:A:367:THR:HG23	1.98	0.46
1:B:101:LEU:HD23	1:B:383:ILE:CD1	2.46	0.46
1:B:220:SER:OG	1:B:221:LEU:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:VAL:HB	1:B:173:VAL:HG22	1.97	0.46
1:B:158:ILE:O	1:B:159:ASP:C	2.59	0.46
1:A:136:VAL:HB	1:A:161:VAL:HA	1.98	0.45
1:A:193:THR:HG22	1:A:221:LEU:HD11	1.97	0.45
1:A:163:LEU:HD12	1:A:173:VAL:HG21	1.97	0.45
1:A:206:LEU:HD12	1:A:206:LEU:HA	1.78	0.45
1:B:113:LEU:N	1:B:113:LEU:CD2	2.80	0.45
1:B:367:THR:O	1:B:368:VAL:C	2.60	0.45
1:A:229:LEU:HD11	1:A:275:MET:HE1	1.99	0.45
1:B:357:VAL:HG12	1:B:359:VAL:HG13	1.99	0.45
1:A:378:ILE:HA	1:A:381:ALA:HB3	1.97	0.45
1:B:187:HIS:O	1:B:187:HIS:ND1	2.38	0.45
1:B:378:ILE:HA	1:B:381:ALA:HB3	1.98	0.45
1:B:221:LEU:HB3	1:B:247:MET:HE2	1.98	0.44
1:B:378:ILE:HG13	1:B:379:ALA:H	1.83	0.44
1:A:238:VAL:HG11	1:A:282:LEU:CD2	2.48	0.44
1:A:195:MET:HG3	1:A:334:GLN:HG3	1.99	0.44
1:B:147:ARG:O	1:B:149:ARG:HG3	2.17	0.44
1:A:314:ARG:HA	1:A:314:ARG:HD2	1.38	0.44
1:A:350:LYS:O	1:A:354:HIS:HB2	2.17	0.44
1:B:181:THR:HG23	1:B:234:VAL:HG22	2.00	0.44
1:A:129:ILE:HG12	1:A:382:VAL:HG21	1.99	0.44
1:A:154:ALA:HB2	1:A:378:ILE:HD13	1.99	0.44
1:B:342:LYS:CE	1:B:349:ASN:HD22	2.31	0.43
1:A:316:GLY:C	1:A:318:ILE:H	2.26	0.43
1:A:324:VAL:HG23	1:A:351:LEU:HG	2.00	0.43
1:B:104:PHE:CD2	1:B:104:PHE:O	2.71	0.43
1:A:147:ARG:O	1:A:149:ARG:HG3	2.17	0.43
1:B:378:ILE:HG13	1:B:379:ALA:N	2.33	0.43
1:A:376:ILE:O	1:A:380:GLU:HG3	2.18	0.43
1:B:131:ARG:O	1:B:153:ARG:CZ	2.67	0.43
1:B:313:ALA:HB1	1:B:314:ARG:HH21	1.83	0.43
1:B:193:THR:HG22	1:B:221:LEU:HD11	2.00	0.43
1:B:386:LEU:HD23	1:B:386:LEU:HA	1.77	0.43
1:A:210:LYS:HB3	1:A:212:GLN:HG3	2.01	0.43
1:A:210:LYS:HD3	1:A:212:GLN:NE2	2.34	0.43
1:A:378:ILE:HG13	1:A:379:ALA:N	2.33	0.43
1:B:221:LEU:HD22	1:B:247:MET:CE	2.49	0.43
1:A:119:ARG:HH11	1:A:119:ARG:HG3	1.83	0.42
1:A:230:GLY:HA3	1:A:284:HIS:O	2.19	0.42
1:A:343:GLU:HG2	1:B:211:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ILE:O	1:A:144:SER:HB2	2.19	0.42
1:A:285:MET:HE3	1:A:285:MET:HB2	1.83	0.42
1:B:166:ALA:O	1:B:171:CYS:N	2.47	0.42
1:B:113:LEU:HD12	1:B:117:ASP:HB3	2.01	0.42
1:B:361:PRO:HD2	1:B:363:LEU:HG	2.02	0.42
1:A:93:LEU:HD11	1:A:378:ILE:CD1	2.45	0.42
1:B:185:ALA:HB1	1:B:238:VAL:HG23	2.02	0.42
1:A:187:HIS:HB2	1:B:197:ARG:CZ	2.50	0.41
1:A:187:HIS:HD1	1:A:187:HIS:C	2.28	0.41
1:A:96:ALA:HB1	1:A:376:ILE:HD11	2.02	0.41
1:A:179:ALA:HB3	1:A:374:VAL:HG23	2.03	0.41
1:B:104:PHE:O	1:B:104:PHE:CG	2.73	0.41
1:B:286:PRO:HA	1:B:314:ARG:CZ	2.50	0.41
1:B:354:HIS:HB3	1:B:357:VAL:CG2	2.51	0.41
1:A:211:TRP:CH2	1:B:343:GLU:HG2	2.56	0.41
1:A:217:VAL:HG13	1:B:366:SER:O	2.20	0.41
1:B:185:ALA:CB	1:B:238:VAL:HG23	2.50	0.41
1:A:212:GLN:O	1:A:214:ASN:N	2.54	0.41
1:A:252:HIS:CE1	1:A:271:MET:HB2	2.56	0.41
1:A:283:LEU:HD12	1:A:318:ILE:HD11	2.01	0.41
1:A:291:THR:HA	1:A:294:MET:HE3	2.02	0.41
1:B:250:ILE:HD12	1:B:269:VAL:HG21	2.02	0.41
1:A:337:LEU:O	1:A:359:VAL:HA	2.21	0.41
1:A:228:ILE:HG12	1:A:282:LEU:HD21	2.03	0.40
1:B:354:HIS:HB3	1:B:357:VAL:HG23	2.02	0.40
1:A:221:LEU:CB	1:A:247:MET:HE2	2.36	0.40
1:A:187:HIS:CE1	1:A:191:LEU:HG	2.57	0.40
1:A:195:MET:SD	1:A:308:ARG:HB3	2.61	0.40
1:A:282:LEU:HD12	1:A:282:LEU:C	2.46	0.40
1:B:283:LEU:HD12	1:B:318:ILE:HD11	2.03	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	303/387 (78%)	277 (91%)	24 (8%)	2 (1%)	19	53
1	B	303/387 (78%)	272 (90%)	31 (10%)	0	100	100
All	All	606/774 (78%)	549 (91%)	55 (9%)	2 (0%)	37	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	213	ARG
1	A	181	THR

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	224/275 (82%)	214 (96%)	10 (4%)	23	53
1	B	224/275 (82%)	217 (97%)	7 (3%)	35	63
All	All	448/550 (82%)	431 (96%)	17 (4%)	28	57

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	131	ARG
1	A	135	LYS
1	A	153	ARG
1	A	195	MET
1	A	205	SER
1	A	207	LYS
1	A	241	ARG
1	A	295	LEU
1	A	312	VAL
1	A	314	ARG
1	B	100	LEU

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Mol	Chain	Res	Type
1	B	131	ARG
1	B	158	ILE
1	B	159	ASP
1	B	172	LEU
1	B	312	VAL
1	B	314	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	ASN
1	A	169	HIS
1	A	212	GLN
1	B	106	ASN
1	B	212	GLN
1	B	252	HIS
1	B	334	GLN
1	B	371	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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	305/387 (78%)	-0.25	0	100 100	77, 106, 131, 168	0
1	B	305/387 (78%)	-0.05	4 (1%)	74 55	30, 111, 163, 190	0
All	All	610/774 (78%)	-0.15	4 (0%)	84 69	30, 108, 156, 190	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	2.5
1	B	314	ARG	2.5
1	B	84	LYS	2.4
1	B	173	VAL	2.1

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 oligosaccharides 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.