



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:52 AM UTC

PDB ID : 2I5G / pdb_00002i5g
Title : Crystal structure of amidohydrolase from *Pseudomonas aeruginosa*
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Deposited on : 2006-08-24
Resolution : 2.60 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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

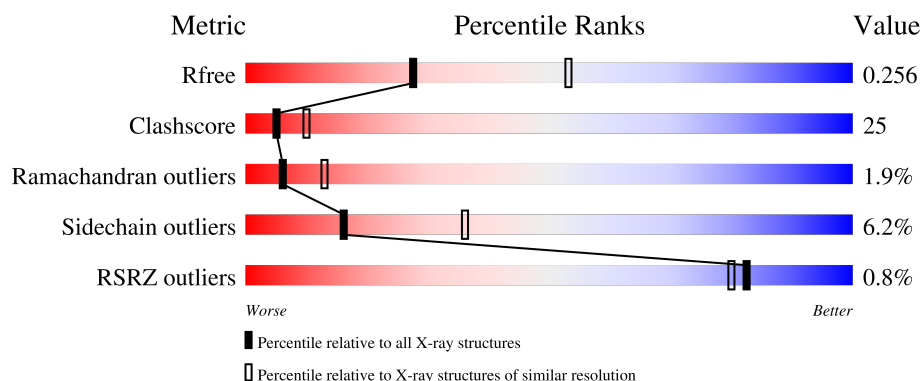
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	325	% 63% 30% 7%
1	B	325	% 56% 37% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5620 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 amidohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	Se	0	0	0
			2538	1603	442	479	6	8			
1	B	325	Total	C	N	O	S	Se	0	0	0
			2538	1603	442	479	6	8			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	LEU	-	cloning artifact	UNP Q9HTG8
A	30	MSE	MET	modified residue	UNP Q9HTG8
A	56	THR	ALA	engineered mutation	UNP Q9HTG8
A	119	MSE	MET	modified residue	UNP Q9HTG8
A	149	MSE	MET	modified residue	UNP Q9HTG8
A	155	MSE	MET	modified residue	UNP Q9HTG8
A	213	MSE	MET	modified residue	UNP Q9HTG8
A	237	MSE	MET	modified residue	UNP Q9HTG8
A	300	LYS	GLU	engineered mutation	UNP Q9HTG8
A	303	MSE	MET	modified residue	UNP Q9HTG8
A	312	MSE	MET	modified residue	UNP Q9HTG8
B	2	LEU	-	cloning artifact	UNP Q9HTG8
B	30	MSE	MET	modified residue	UNP Q9HTG8
B	56	THR	ALA	engineered mutation	UNP Q9HTG8
B	119	MSE	MET	modified residue	UNP Q9HTG8
B	149	MSE	MET	modified residue	UNP Q9HTG8
B	155	MSE	MET	modified residue	UNP Q9HTG8
B	213	MSE	MET	modified residue	UNP Q9HTG8
B	237	MSE	MET	modified residue	UNP Q9HTG8
B	300	LYS	GLU	engineered mutation	UNP Q9HTG8
B	303	MSE	MET	modified residue	UNP Q9HTG8
B	312	MSE	MET	modified residue	UNP Q9HTG8

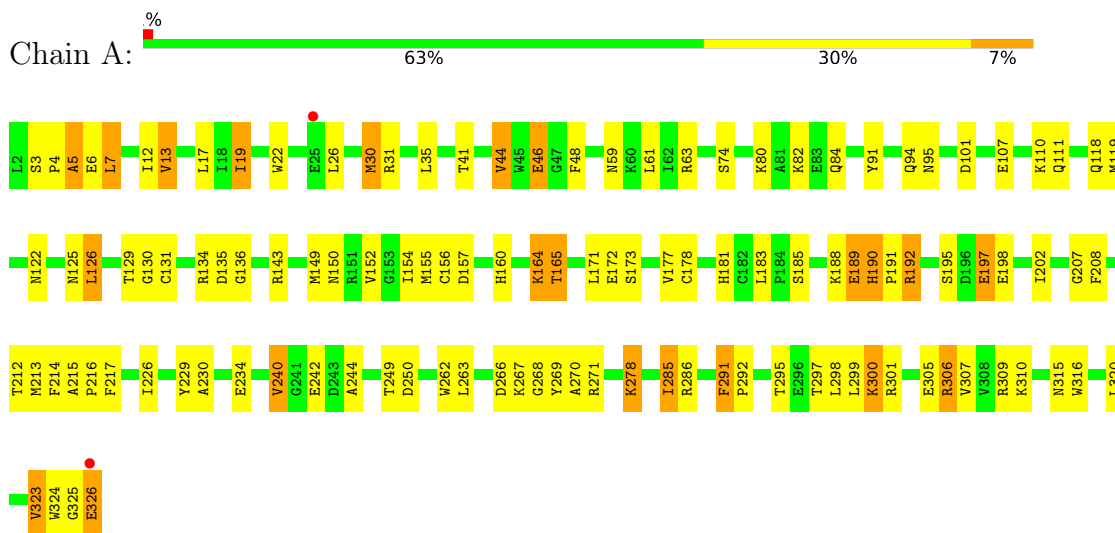
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	257	Total 257	O 257	0	0
2	B	287	Total 287	O 287	0	0

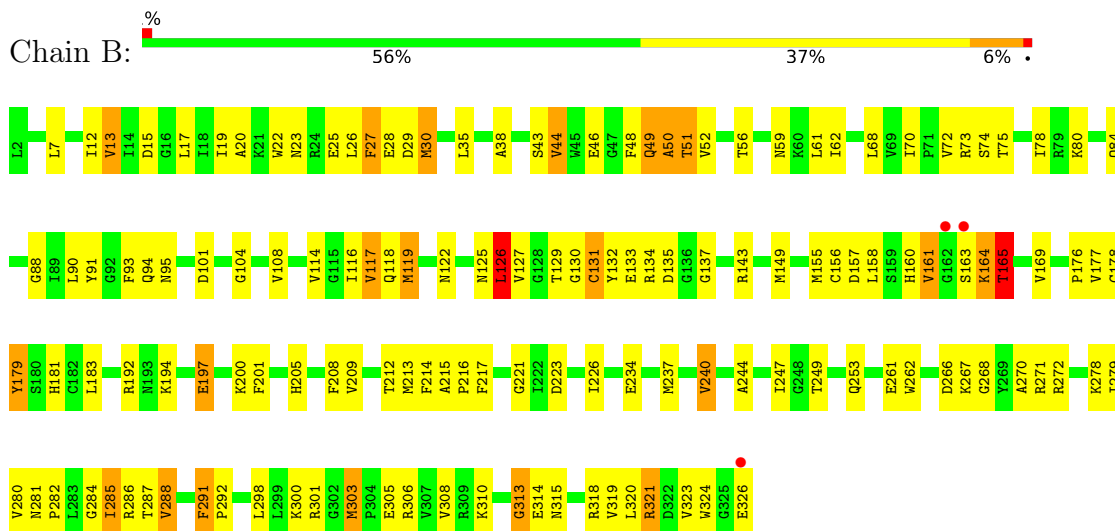
3 Residue-property plots [i](#)

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: amidohydrolase



• Molecule 1: amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	81.64Å 150.62Å 68.83Å 90.00° 97.94° 90.00°	Depositor
Resolution (Å)	19.94 – 2.60 19.94 – 2.60	Depositor EDS
% Data completeness (in resolution range)	86.9 (19.94-2.60) 96.3 (19.94-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.48 (at 2.53Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.181 , 0.250 0.194 , 0.256	Depositor DCC
R_{free} test set	2073 reflections (3.99%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.792	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5620	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.20% 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	0.41	0/2580	0.95	6/3469 (0.2%)
1	B	0.54	5/2580 (0.2%)	0.98	14/3469 (0.4%)
All	All	0.48	5/5160 (0.1%)	0.97	20/6938 (0.3%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	303	MSE	SE-CE	-7.97	1.71	1.95
1	B	149	MSE	SE-CE	-6.80	1.75	1.95
1	B	237	MSE	SE-CE	-6.01	1.77	1.95
1	B	30	MSE	SE-CE	-5.87	1.77	1.95
1	B	213	MSE	CG-SE	-5.65	1.78	1.95

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	PHE	N-CA-C	-11.62	96.12	110.41
1	B	126	LEU	N-CA-C	-8.66	102.84	113.50
1	A	126	LEU	N-CA-C	-8.06	103.59	113.50
1	B	51	THR	N-CA-C	-7.82	102.84	111.36
1	B	165	THR	N-CA-C	-7.32	103.00	112.23
1	A	323	VAL	N-CA-C	6.58	117.33	110.62
1	A	46	GLU	N-CA-C	6.54	119.24	110.35
1	B	29	ASP	N-CA-C	-6.43	103.78	111.69
1	B	119	MSE	N-CA-C	6.20	118.12	111.36
1	B	179	TYR	N-CA-C	-6.11	96.11	107.71
1	B	240	VAL	N-CA-C	5.89	118.42	111.05
1	A	189	GLU	N-CA-C	5.83	117.33	110.97
1	B	117	VAL	N-CA-C	5.73	116.37	108.12
1	B	131	CYS	N-CA-C	-5.32	103.47	110.33
1	B	161	VAL	N-CA-C	5.16	119.63	112.35
1	B	261	GLU	N-CA-C	-5.10	105.91	111.82
1	A	240	VAL	N-CA-C	5.05	115.82	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	49	GLN	N-CA-C	-5.05	102.39	109.96
1	B	209	VAL	N-CA-C	5.04	115.90	107.18
1	A	30	MSE	N-CA-C	-5.01	105.90	111.36

There are no chirality outliers.

There are no planarity outliers.

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	2538	0	2503	125	0
1	B	2538	0	2503	130	0
2	A	257	0	0	6	0
2	B	287	0	0	17	0
All	All	5620	0	5006	250	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 25.

All (250) 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:B:303:MSE:HE3	1:B:308:VAL:HA	1.37	1.04
1:A:101:ASP:HB2	1:B:267:LYS:HD2	1.41	1.02
1:A:135:ASP:HB3	1:A:164:LYS:HE3	1.45	0.93
1:A:129:THR:HG21	1:A:136:GLY:H	1.36	0.90
1:A:324:TRP:C	1:A:326:GLU:H	1.80	0.85
1:A:192:ARG:HG2	1:A:192:ARG:HH11	1.41	0.83
1:B:17:LEU:HD13	1:B:181:HIS:CE1	2.15	0.82
1:A:12:ILE:HD13	1:A:82:LYS:HG3	1.64	0.80
1:B:240:VAL:HG13	1:B:244:ALA:HB3	1.64	0.79
1:A:213:MSE:HE2	1:A:229:TYR:CE2	2.18	0.79
1:A:215:ALA:HB3	1:A:216:PRO:HD3	1.65	0.78
1:B:131:CYS:O	1:B:132:TYR:HB2	1.82	0.78
1:A:135:ASP:OD2	1:A:165:THR:HG22	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:MSE:CE	1:B:308:VAL:HA	2.14	0.75
1:B:135:ASP:OD1	1:B:164:LYS:HB3	1.87	0.73
1:B:253:GLN:HE22	1:B:288:VAL:H	1.37	0.73
1:B:303:MSE:HE3	1:B:308:VAL:CA	2.18	0.72
1:A:195:SER:OG	1:A:197:GLU:HG2	1.90	0.72
1:A:129:THR:HG21	1:A:136:GLY:N	2.04	0.72
1:B:20:ALA:HB1	1:B:253:GLN:O	1.88	0.72
1:A:213:MSE:HE1	1:A:285:ILE:HD13	1.72	0.72
1:B:306:ARG:HH12	1:B:310:LYS:NZ	1.88	0.71
1:B:326:GLU:HG3	2:B:542:HOH:O	1.90	0.71
1:A:300:LYS:HE3	2:A:362:HOH:O	1.90	0.71
1:A:74:SER:HB2	1:A:326:GLU:HG2	1.73	0.71
1:A:195:SER:HB2	1:A:197:GLU:OE2	1.92	0.70
1:A:192:ARG:HG2	1:A:192:ARG:NH1	2.06	0.69
1:A:19:ILE:N	1:A:19:ILE:HD13	2.07	0.69
1:B:126:LEU:HB2	2:B:560:HOH:O	1.92	0.68
1:B:208:PHE:HE1	1:B:315:ASN:HB3	1.56	0.68
1:A:324:TRP:C	1:A:326:GLU:N	2.50	0.68
1:B:215:ALA:HB3	1:B:216:PRO:HD3	1.76	0.67
1:B:129:THR:H	1:B:165:THR:HG21	1.59	0.67
1:B:310:LYS:HA	1:B:314:GLU:HB2	1.77	0.67
1:B:303:MSE:HE2	1:B:308:VAL:HG22	1.77	0.67
1:B:156:CYS:HB2	1:B:177:VAL:HG12	1.77	0.66
1:A:214:PHE:CD2	1:A:216:PRO:HD2	2.30	0.66
1:A:74:SER:HB2	1:A:326:GLU:CG	2.25	0.66
1:B:321:ARG:HG2	1:B:321:ARG:HH11	1.60	0.65
1:A:74:SER:CB	1:A:326:GLU:HG2	2.27	0.65
1:B:27:PHE:O	1:B:28:GLU:CB	2.45	0.65
1:B:26:LEU:HD11	1:B:30:MSE:HE3	1.79	0.64
1:B:321:ARG:HA	1:B:326:GLU:O	1.97	0.64
1:B:12:ILE:O	1:B:313:GLY:HA3	1.98	0.64
1:B:183:LEU:HD12	1:B:217:PHE:HB2	1.79	0.64
1:A:12:ILE:CD1	1:A:82:LYS:HG3	2.28	0.63
1:A:213:MSE:HE3	1:A:249:THR:HG22	1.79	0.63
1:B:116:ILE:HD12	1:B:155:MSE:O	1.99	0.63
1:A:22:TRP:HB3	1:A:61:LEU:HD13	1.80	0.62
1:A:213:MSE:HG3	2:A:359:HOH:O	1.98	0.62
1:B:13:VAL:CG2	1:B:35:LEU:HA	2.29	0.62
1:A:240:VAL:CG1	1:A:244:ALA:HB3	2.30	0.62
1:A:13:VAL:HG22	1:A:35:LEU:HA	1.82	0.62
1:A:185:SER:HA	1:A:188:LYS:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:H	1:A:165:THR:HG21	1.65	0.62
1:B:17:LEU:HD13	1:B:181:HIS:ND1	2.14	0.62
1:A:156:CYS:HB2	1:A:177:VAL:HG12	1.81	0.61
1:B:70:ILE:HG22	2:B:340:HOH:O	2.00	0.61
1:A:234:GLU:OE1	1:A:301:ARG:NH1	2.34	0.61
1:A:44:VAL:HG13	1:A:262:TRP:CH2	2.36	0.61
1:A:101:ASP:CB	1:B:267:LYS:HD2	2.24	0.61
1:B:59:ASN:ND2	1:B:91:TYR:OH	2.33	0.61
1:A:240:VAL:HG13	1:A:244:ALA:HB3	1.84	0.60
1:A:129:THR:HG21	1:A:135:ASP:HA	1.83	0.60
1:B:104:GLY:O	1:B:108:VAL:HG23	2.02	0.59
1:B:119:MSE:HE1	1:B:156:CYS:CB	2.32	0.59
1:B:119:MSE:HE1	1:B:156:CYS:HB3	1.84	0.59
1:B:306:ARG:HD3	2:B:580:HOH:O	2.02	0.59
1:B:27:PHE:O	1:B:28:GLU:HB2	2.03	0.59
1:B:13:VAL:HG23	1:B:35:LEU:HD12	1.85	0.58
1:B:49:GLN:O	1:B:50:ALA:CB	2.51	0.58
1:A:4:PRO:HB3	1:A:299:LEU:HD23	1.83	0.58
1:A:295:THR:O	1:A:299:LEU:HB2	2.03	0.58
1:A:118:GLN:HA	1:A:157:ASP:HB3	1.85	0.58
1:B:165:THR:O	1:B:169:VAL:HG23	2.04	0.58
1:B:130:GLY:HA2	1:B:161:VAL:O	2.04	0.58
1:B:298:LEU:HD13	1:B:303:MSE:HE1	1.86	0.58
1:B:44:VAL:HG13	1:B:262:TRP:CH2	2.39	0.57
1:A:150:ASN:HD21	1:A:173:SER:HA	1.70	0.57
1:A:213:MSE:HE1	1:A:285:ILE:CD1	2.35	0.57
1:B:214:PHE:CD2	1:B:216:PRO:HD2	2.40	0.57
1:B:17:LEU:O	1:B:249:THR:O	2.23	0.57
1:B:74:SER:HB2	2:B:334:HOH:O	2.05	0.57
1:A:26:LEU:O	1:A:30:MSE:HG3	2.05	0.57
1:A:278:LYS:HZ3	1:A:278:LYS:HB2	1.70	0.56
1:A:183:LEU:HD21	1:A:212:THR:O	2.05	0.56
1:A:189:GLU:O	1:A:190:HIS:O	2.23	0.56
1:A:306:ARG:O	1:A:310:LYS:HG2	2.05	0.56
1:A:249:THR:O	1:A:250:ASP:HB3	2.06	0.56
1:A:291:PHE:N	1:A:292:PRO:HD2	2.19	0.56
1:B:12:ILE:O	1:B:12:ILE:HG22	2.05	0.56
1:B:285:ILE:O	1:B:285:ILE:HG13	2.06	0.56
1:B:155:MSE:HB2	1:B:323:VAL:HG11	1.88	0.56
1:A:122:ASN:C	1:A:130:GLY:HA3	2.31	0.56
1:B:131:CYS:O	1:B:132:TYR:CB	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:PRO:HG2	1:B:323:VAL:CG2	2.37	0.55
1:B:319:VAL:O	1:B:323:VAL:HG23	2.06	0.55
1:A:213:MSE:HE3	1:A:249:THR:CG2	2.36	0.55
1:B:326:GLU:C	2:B:334:HOH:O	2.49	0.55
1:A:5:ALA:N	2:A:528:HOH:O	2.39	0.55
1:A:183:LEU:HD12	1:A:217:PHE:HB2	1.90	0.54
1:B:163:SER:O	1:B:164:LYS:CB	2.56	0.54
1:B:43:SER:HB2	1:B:51:THR:HG23	1.89	0.54
1:A:305:GLU:O	1:A:309:ARG:HG2	2.08	0.54
1:A:44:VAL:CG1	1:A:262:TRP:CH2	2.90	0.54
1:B:137:GLY:HA2	1:B:164:LYS:HG3	1.90	0.54
1:B:240:VAL:CG1	1:B:244:ALA:HB3	2.36	0.53
1:B:253:GLN:NE2	1:B:288:VAL:H	2.06	0.53
1:A:17:LEU:HD22	1:A:181:HIS:CE1	2.43	0.53
1:A:324:TRP:O	1:A:326:GLU:N	2.41	0.53
1:A:155:MSE:HE2	1:A:178:CYS:HB3	1.90	0.53
1:B:134:ARG:HD2	2:B:476:HOH:O	2.09	0.52
1:B:44:VAL:HG13	1:B:262:TRP:HH2	1.74	0.52
1:B:80:LYS:HE2	1:B:84:GLN:HE22	1.74	0.52
1:A:135:ASP:CG	1:A:165:THR:HG22	2.34	0.52
1:B:135:ASP:OD2	1:B:165:THR:HG22	2.08	0.52
1:A:242:GLU:OE1	1:A:306:ARG:NH1	2.43	0.52
1:B:13:VAL:HG22	1:B:35:LEU:HA	1.92	0.52
1:B:19:ILE:HG23	1:B:19:ILE:O	2.09	0.52
1:A:160:HIS:ND1	1:A:192:ARG:HD2	2.24	0.52
1:A:230:ALA:O	1:A:234:GLU:HG3	2.09	0.52
1:A:195:SER:HG	1:A:197:GLU:HG2	1.76	0.51
1:A:202:ILE:HG12	1:A:207:GLY:HA3	1.92	0.51
1:B:73:ARG:HD3	2:B:378:HOH:O	2.10	0.51
1:A:136:GLY:HA3	2:B:532:HOH:O	2.09	0.51
1:A:135:ASP:CG	1:A:164:LYS:HB2	2.35	0.51
1:A:80:LYS:CG	1:A:84:GLN:NE2	2.74	0.51
1:B:253:GLN:HE22	1:B:288:VAL:N	2.08	0.51
1:A:285:ILE:HG23	1:A:285:ILE:O	2.10	0.51
1:B:52:VAL:O	1:B:56:THR:HG23	2.10	0.51
1:A:31:ARG:HA	1:A:31:ARG:NE	2.25	0.51
1:B:163:SER:O	1:B:164:LYS:HB2	2.10	0.50
1:A:118:GLN:OE1	1:A:160:HIS:HD2	1.94	0.50
1:A:13:VAL:CG2	1:A:35:LEU:HD12	2.42	0.50
1:A:95:ASN:HB3	2:A:372:HOH:O	2.11	0.50
1:B:95:ASN:HB3	2:B:374:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:ASN:O	1:B:131:CYS:O	2.30	0.50
1:B:201:PHE:O	1:B:205:HIS:HD2	1.95	0.50
1:A:19:ILE:HG22	1:A:41:THR:HB	1.93	0.50
1:A:131:CYS:SG	1:A:192:ARG:HG3	2.52	0.49
1:A:306:ARG:HA	2:A:458:HOH:O	2.12	0.49
1:A:119:MSE:CE	1:A:156:CYS:HB3	2.41	0.49
1:B:118:GLN:HA	1:B:157:ASP:HB3	1.95	0.49
1:B:197:GLU:CD	1:B:197:GLU:N	2.70	0.49
1:A:7:LEU:HD21	1:A:309:ARG:HD2	1.94	0.49
1:B:135:ASP:CG	1:B:164:LYS:HB3	2.36	0.49
1:A:320:LEU:HA	1:A:323:VAL:HG22	1.95	0.49
1:B:116:ILE:HD13	1:B:155:MSE:HG2	1.95	0.49
1:B:176:PRO:HG2	1:B:323:VAL:HG22	1.94	0.49
1:A:3:SER:OG	1:A:6:GLU:HG3	2.12	0.49
1:A:26:LEU:HD11	1:A:30:MSE:HE3	1.94	0.49
1:A:143:ARG:NH2	1:A:172:GLU:OE1	2.46	0.49
1:A:266:ASP:OD2	1:A:267:LYS:HD3	2.12	0.49
1:A:129:THR:CG2	1:A:135:ASP:HA	2.42	0.49
1:B:291:PHE:N	1:B:292:PRO:HD2	2.28	0.49
1:B:272:ARG:NH1	2:B:330:HOH:O	2.45	0.49
1:B:292:PRO:HG2	2:B:412:HOH:O	2.12	0.49
1:B:300:LYS:HE2	2:B:504:HOH:O	2.13	0.49
1:B:279:ILE:HG22	2:B:358:HOH:O	2.12	0.49
1:A:17:LEU:HD22	1:A:181:HIS:CD2	2.48	0.48
1:B:23:ASN:OD1	1:B:25:GLU:HG2	2.13	0.48
1:A:19:ILE:HD11	1:A:250:ASP:OD1	2.13	0.48
1:B:135:ASP:OD2	1:B:165:THR:CG2	2.61	0.48
1:A:208:PHE:HE1	1:A:315:ASN:HB3	1.78	0.48
1:B:26:LEU:C	1:B:27:PHE:O	2.52	0.48
1:B:80:LYS:CE	1:B:84:GLN:HE22	2.27	0.48
1:A:262:TRP:O	1:A:268:GLY:HA2	2.13	0.48
1:A:249:THR:O	1:A:250:ASP:CB	2.60	0.47
1:B:93:PHE:HB2	1:B:117:VAL:HG22	1.96	0.47
1:B:320:LEU:HD22	1:B:324:TRP:CZ2	2.50	0.47
1:A:271:ARG:HH11	1:A:271:ARG:HG2	1.78	0.47
1:A:266:ASP:O	1:A:267:LYS:HB2	2.13	0.47
1:B:155:MSE:HE2	1:B:178:CYS:HB3	1.96	0.47
1:A:19:ILE:CG2	1:A:41:THR:HB	2.45	0.47
1:A:269:TYR:CD1	1:A:269:TYR:C	2.92	0.47
1:B:158:LEU:HD23	1:B:161:VAL:HG21	1.96	0.47
1:A:118:GLN:OE1	1:A:160:HIS:CD2	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:226:ILE:HD12	1:B:284:GLY:O	2.14	0.46
1:A:183:LEU:HG	1:A:212:THR:HG22	1.97	0.46
1:B:181:HIS:HB2	1:B:212:THR:HB	1.96	0.46
1:A:307:VAL:O	1:A:310:LYS:HB2	2.15	0.46
1:A:267:LYS:CG	1:B:101:ASP:HB2	2.46	0.46
1:B:75:THR:O	1:B:78:ILE:HB	2.16	0.46
1:B:93:PHE:CZ	1:B:114:VAL:HG21	2.50	0.46
1:A:59:ASN:ND2	1:A:91:TYR:OH	2.49	0.46
1:A:190:HIS:HA	1:A:191:PRO:HD3	1.87	0.46
1:A:80:LYS:HG3	1:A:84:GLN:NE2	2.31	0.45
1:B:129:THR:HB	1:B:135:ASP:OD2	2.16	0.45
1:B:266:ASP:OD2	1:B:271:ARG:NH2	2.49	0.45
1:B:35:LEU:HD21	1:B:38:ALA:HB2	1.98	0.45
1:A:214:PHE:CE2	1:A:216:PRO:HG2	2.51	0.45
1:B:266:ASP:C	1:B:268:GLY:N	2.75	0.45
1:A:183:LEU:HG	1:A:212:THR:CG2	2.47	0.44
1:A:271:ARG:HG2	1:A:271:ARG:NH1	2.32	0.44
1:A:126:LEU:HG	2:A:426:HOH:O	2.16	0.44
1:B:94:GLN:HA	1:B:118:GLN:HB3	1.98	0.44
1:B:303:MSE:CE	1:B:308:VAL:HG22	2.46	0.44
1:A:119:MSE:HE1	1:A:156:CYS:HA	1.99	0.44
1:B:22:TRP:HB3	1:B:61:LEU:CD1	2.48	0.44
1:A:26:LEU:HD13	1:A:26:LEU:C	2.43	0.43
1:B:281:ASN:HB2	1:B:286:ARG:HG2	2.00	0.43
1:B:28:GLU:HG3	2:B:389:HOH:O	2.18	0.43
1:B:234:GLU:OE2	1:B:301:ARG:HD3	2.17	0.43
1:B:266:ASP:C	1:B:268:GLY:H	2.26	0.43
1:B:306:ARG:HH12	1:B:310:LYS:CE	2.30	0.43
1:B:282:PRO:HG2	1:B:285:ILE:CG2	2.47	0.43
1:B:125:ASN:HB3	1:B:127:VAL:H	1.84	0.43
1:A:213:MSE:HE3	1:A:249:THR:CB	2.48	0.43
1:B:179:TYR:CE1	1:B:194:LYS:HE3	2.54	0.43
1:A:13:VAL:CG2	1:A:35:LEU:HA	2.48	0.43
1:A:80:LYS:HG3	1:A:84:GLN:HE21	1.83	0.43
1:A:267:LYS:HG3	1:B:101:ASP:HB2	2.00	0.43
1:A:119:MSE:HE3	1:A:156:CYS:HB3	2.00	0.43
1:A:17:LEU:HD23	1:A:250:ASP:OD1	2.17	0.43
1:B:133:GLU:O	1:B:134:ARG:C	2.61	0.43
1:B:160:HIS:CE1	1:B:192:ARG:HE	2.37	0.43
1:B:298:LEU:HB3	1:B:303:MSE:HE2	2.01	0.43
1:B:17:LEU:C	1:B:17:LEU:HD23	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:GLN:HG3	2:B:359:HOH:O	2.19	0.42
1:B:278:LYS:HG3	1:B:280:VAL:HG23	2.00	0.42
1:B:49:GLN:O	1:B:50:ALA:HB2	2.19	0.42
1:A:110:LYS:HG2	1:A:152:VAL:CG1	2.49	0.42
1:A:195:SER:OG	1:A:198:GLU:HG3	2.20	0.42
1:B:72:VAL:CG2	1:B:88:GLY:HA3	2.49	0.42
1:A:129:THR:CG2	1:A:135:ASP:OD2	2.68	0.42
1:B:253:GLN:NE2	1:B:287:THR:HA	2.34	0.42
1:A:299:LEU:C	1:A:301:ARG:H	2.27	0.42
1:B:221:GLY:C	1:B:223:ASP:H	2.27	0.42
1:A:286:ARG:HG3	1:A:286:ARG:HH11	1.85	0.42
1:B:26:LEU:C	1:B:26:LEU:HD13	2.44	0.42
1:B:247:ILE:CG2	1:B:249:THR:HG23	2.50	0.42
1:A:149:MSE:HE3	1:A:154:ILE:HG22	2.02	0.42
1:A:213:MSE:HE3	1:A:249:THR:HB	2.02	0.41
1:A:119:MSE:HE1	1:A:156:CYS:CB	2.49	0.41
1:B:306:ARG:HH12	1:B:310:LYS:HZ1	1.66	0.41
1:A:94:GLN:C	1:A:118:GLN:HB3	2.45	0.41
1:B:94:GLN:C	1:B:118:GLN:HB3	2.45	0.41
1:A:134:ARG:NH2	1:B:143:ARG:HH22	2.18	0.41
1:B:118:GLN:HE22	1:B:160:HIS:CD2	2.39	0.41
1:B:321:ARG:HH11	1:B:321:ARG:CG	2.30	0.41
1:A:107:GLU:O	1:A:111:GLN:HG3	2.21	0.41
1:B:15:ASP:OD2	1:B:249:THR:OG1	2.37	0.41
1:B:281:ASN:HB2	1:B:286:ARG:HA	2.03	0.41
1:B:318:ARG:HD3	2:B:539:HOH:O	2.20	0.41
1:B:119:MSE:O	1:B:125:ASN:ND2	2.54	0.41
1:A:4:PRO:O	1:A:5:ALA:HB2	2.20	0.40
1:A:119:MSE:HE1	1:A:156:CYS:CA	2.51	0.40
1:A:208:PHE:HE2	1:A:316:TRP:CE2	2.40	0.40
1:A:129:THR:HG23	1:A:135:ASP:OD2	2.22	0.40
1:A:297:THR:O	1:A:298:LEU:C	2.65	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	323/325 (99%)	296 (92%)	21 (6%)	6 (2%)	6	13
1	B	323/325 (99%)	294 (91%)	23 (7%)	6 (2%)	6	13
All	All	646/650 (99%)	590 (91%)	44 (7%)	12 (2%)	6	13

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	ALA
1	A	190	HIS
1	B	48	PHE
1	B	164	LYS
1	B	270	ALA
1	B	285	ILE
1	B	50	ALA
1	A	285	ILE
1	A	48	PHE
1	A	270	ALA
1	B	313	GLY
1	A	325	GLY

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	272/264 (103%)	253 (93%)	19 (7%)	14	31
1	B	272/264 (103%)	257 (94%)	15 (6%)	19	41
All	All	544/528 (103%)	510 (94%)	34 (6%)	16	36

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	13	VAL
1	A	19	ILE
1	A	44	VAL
1	A	46	GLU
1	A	63	ARG
1	A	125	ASN
1	A	164	LYS
1	A	165	THR
1	A	171	LEU
1	A	192	ARG
1	A	197	GLU
1	A	226	ILE
1	A	263	LEU
1	A	278	LYS
1	A	291	PHE
1	A	300	LYS
1	A	306	ARG
1	A	326	GLU
1	B	7	LEU
1	B	13	VAL
1	B	44	VAL
1	B	46	GLU
1	B	62	ILE
1	B	68	LEU
1	B	90	LEU
1	B	126	LEU
1	B	165	THR
1	B	197	GLU
1	B	200	LYS
1	B	288	VAL
1	B	291	PHE
1	B	305	GLU
1	B	321	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	59	ASN
1	A	84	GLN
1	A	125	ASN
1	A	150	ASN

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Mol	Chain	Res	Type
1	A	293	ASN
1	B	53	ASN
1	B	59	ASN
1	B	84	GLN
1	B	111	GLN
1	B	160	HIS
1	B	205	HIS
1	B	253	GLN

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 oligosaccharides 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 [i](#)

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	317/325 (97%)	-0.25	2 (0%) 85 83	3, 15, 32, 52	0
1	B	317/325 (97%)	-0.31	3 (0%) 81 78	2, 15, 33, 66	0
All	All	634/650 (97%)	-0.28	5 (0%) 82 80	2, 15, 33, 66	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	163	SER	2.8
1	B	326	GLU	2.6
1	A	25	GLU	2.4
1	B	162	GLY	2.1
1	A	326	GLU	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.