



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 02:29 PM EDT

PDB ID : 2I5G
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

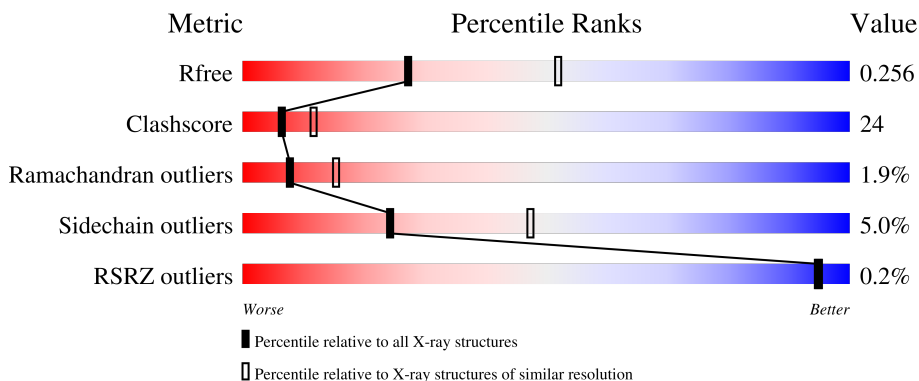
1 Overall quality at a glance

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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (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	
1	B	325	

2 Entry composition [i](#)

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
			Total	C	N	O	S	Se			
1	A	325	2538	1603	442	479	6	8	0	0	0
1	B	325	2538	1603	442	479	6	8	0	0	0

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

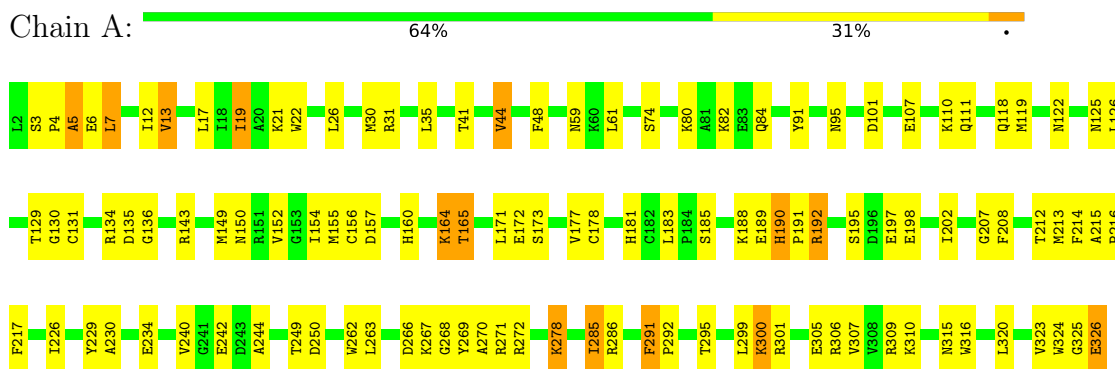
- Molecule 2 is water.

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

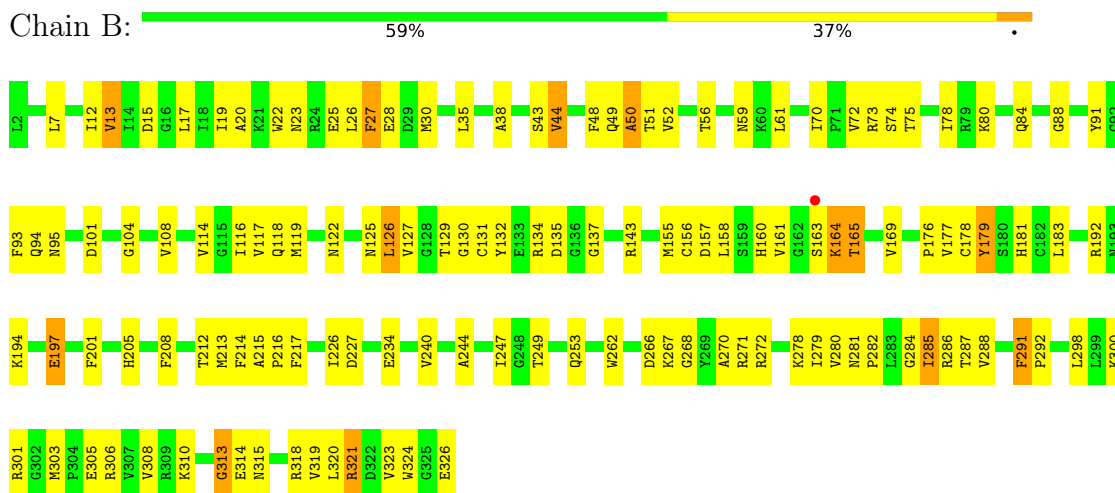
3 Residue-property plots [i](#)

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 20.09 – 2.52	Depositor EDS
% Data completeness (in resolution range)	86.9 (19.94-2.60) 96.3 (20.09-2.52)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.48 (at 2.53Å)	Xtrriage
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	Xtrriage
Anisotropy	0.792	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5620	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtrriage'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 [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	0.34	0/2580	0.60	0/3469
1	B	0.42	0/2580	0.62	2/3469 (0.1%)
All	All	0.38	0/5160	0.61	2/6938 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	TYR	N-CA-C	-5.52	96.11	111.00
1	B	27	PHE	N-CA-C	-5.51	96.12	111.00

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	2538	0	2503	122	0
1	B	2538	0	2503	124	0
2	A	257	0	0	7	0
2	B	287	0	0	17	0
All	All	5620	0	5006	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 24.

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: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: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:324:TRP:C	1:A:326:GLU:H	1.80	0.81
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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:324:TRP:C	1:A:326:GLU:N	2.50	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:B:131:CYS:O	1:B:132:TYR:CB	2.54	0.55
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: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: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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:122:ASN:C	1:A:130:GLY:HA3	2.31	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:B:326:GLU:C	2:B:334:HOH:O	2.49	0.50
1:A:13:VAL:CG2	1:A:35:LEU:HD12	2.42	0.50
1:A:195:SER:HG	1:A:197:GLU:HG2	1.77	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
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:A:7:LEU:HD21	1:A:309:ARG:HD2	1.94	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: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:135:ASP:CG	1:A:165:THR:HG22	2.34	0.47
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
1:A:135:ASP:CG	1:A:164:LYS:HB2	2.35	0.46
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:A:21:LYS:NZ	2:A:572:HOH:O	2.50	0.45
1:B:35:LEU:HD21	1:B:38:ALA:HB2	1.98	0.45
1:B:197:GLU:CD	1:B:197:GLU:N	2.70	0.45
1:A:214:PHE:CE2	1:A:216:PRO:HG2	2.51	0.45
1:B:135:ASP:CG	1:B:164:LYS:HB3	2.36	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:B:281:ASN:HB2	1:B:286:ARG:HG2	2.00	0.43
1:B:26:LEU:C	1:B:27:PHE:O	2.52	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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:A:278:LYS:HZ3	1:A:278:LYS:HB2	1.83	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:A:269:TYR:CD1	1:A:269:TYR:C	2.92	0.42
1:B:253:GLN:NE2	1:B:287:THR:HA	2.34	0.42
1:A:286:ARG:HG3	1:A:286:ARG:HH11	1.85	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:B:15:ASP:OD2	1:B:249:THR:OG1	2.37	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: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: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:B:266:ASP:C	1:B:268:GLY:N	2.75	0.40
1:A:129:THR:HG23	1:A:135:ASP:OD2	2.22	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	323/325 (99%)	296 (92%)	21 (6%)	6 (2%)	8	15
1	B	323/325 (99%)	294 (91%)	23 (7%)	6 (2%)	8	15
All	All	646/650 (99%)	590 (91%)	44 (7%)	12 (2%)	8	15

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 [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	272/264 (103%)	256 (94%)	16 (6%)	19	39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	272/264 (103%)	261 (96%)	11 (4%)	31	57
All	All	544/528 (103%)	517 (95%)	27 (5%)	24	47

All (27) 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	125	ASN
1	A	164	LYS
1	A	165	THR
1	A	171	LEU
1	A	192	ARG
1	A	226	ILE
1	A	263	LEU
1	A	272	ARG
1	A	278	LYS
1	A	291	PHE
1	A	300	LYS
1	A	326	GLU
1	B	7	LEU
1	B	13	VAL
1	B	44	VAL
1	B	126	LEU
1	B	165	THR
1	B	197	GLU
1	B	213	MSE
1	B	227	ASP
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

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Mol	Chain	Res	Type
1	A	150	ASN
1	A	293	ASN
1	B	53	ASN
1	B	59	ASN
1	B	84	GLN
1	B	160	HIS
1	B	205	HIS
1	B	238	ASN
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 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 [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%)	-0.51	0 100 100	3, 15, 32, 52	0
1	B	317/325 (97%)	-0.56	1 (0%) 94 93	2, 15, 33, 66	0
All	All	634/650 (97%)	-0.54	1 (0%) 95 95	2, 15, 33, 66	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	163	SER	2.4

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.