



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:44 AM UTC

PDB ID : 9ISA / pdb_00009isa
Title : Dimeric amylosucrase from *Deinococcus geothermalis*
Authors : Kim, D.S.; Park, J.H.; Seo, D.
Deposited on : 2024-07-17
Resolution : 2.69 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
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

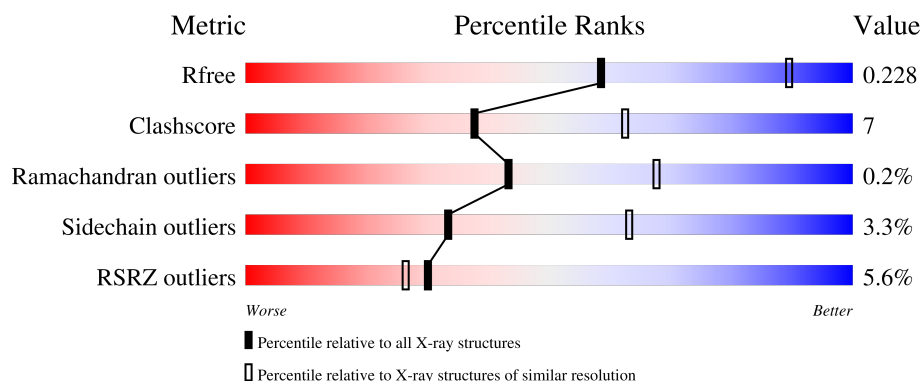
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	645	5% 82% 14% ..
1	B	645	6% 85% 14% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10170 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 Amylosucrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	628	Total	C	N	O	S	0	1	0
			5030	3185	927	906	12			
1	B	639	Total	C	N	O	S	0	0	0
			5100	3234	934	920	12			

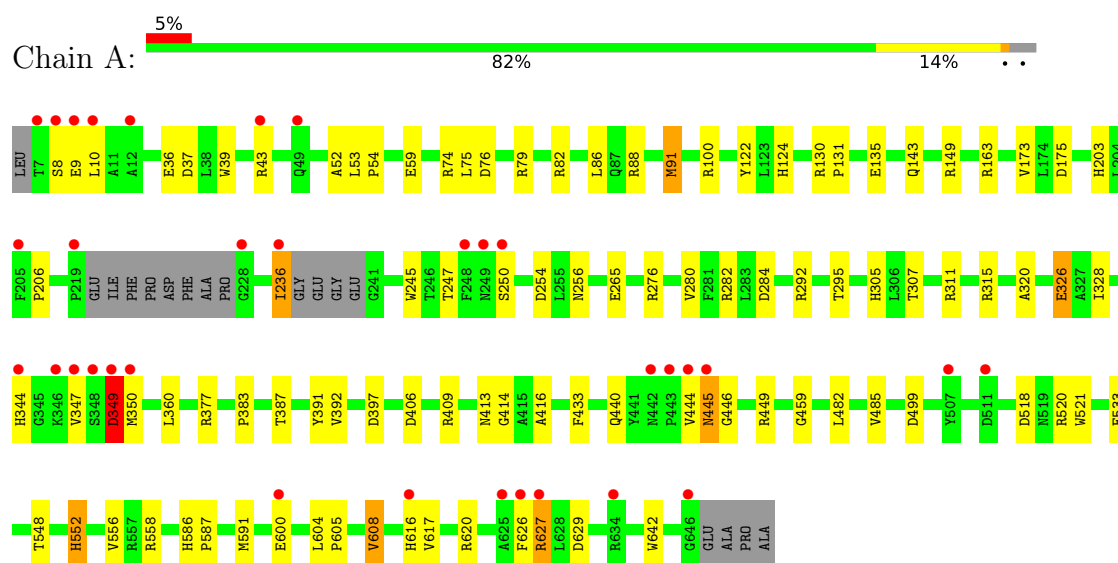
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	24	Total	O	0	0
			24	24		
2	B	16	Total	O	0	0
			16	16		

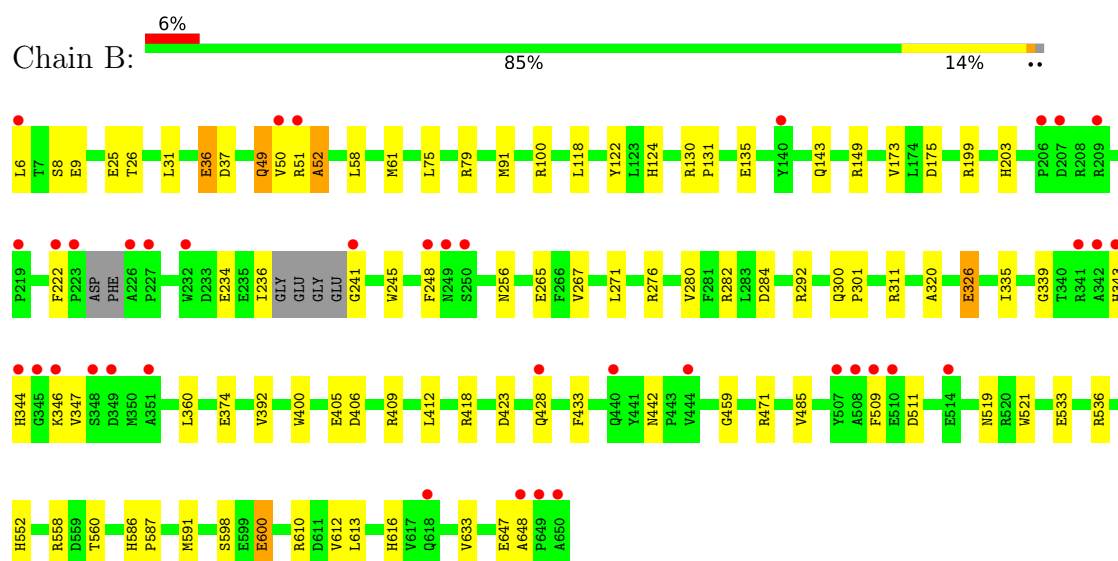
3 Residue-property plots [i](#)

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: Amylosucrase



• Molecule 1: Amylosucrase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	149.92Å 149.92Å 188.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.22 – 2.69 46.22 – 2.69	Depositor EDS
% Data completeness (in resolution range)	86.9 (46.22-2.69) 86.9 (46.22-2.69)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.176 , 0.220 0.192 , 0.228	Depositor DCC
R_{free} test set	2544 reflections (4.25%)	wwPDB-VP
Wilson B-factor (Å ²)	51.3	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10170	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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	0.67	2/5159 (0.0%)	1.15	11/7012 (0.2%)
1	B	0.64	1/5233 (0.0%)	1.15	11/7117 (0.2%)
All	All	0.66	3/10392 (0.0%)	1.15	22/14129 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	552	HIS	CE1-NE2	6.19	1.38	1.32
1	A	305	HIS	CE1-NE2	5.65	1.38	1.32
1	A	552	HIS	CE1-NE2	5.20	1.37	1.32

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	349	ASP	N-CA-C	-7.32	103.50	113.30
1	B	442	ASN	CB-CA-C	6.77	117.68	110.17
1	B	392	VAL	N-CA-C	-6.22	106.46	111.81
1	A	52	ALA	CA-C-N	6.21	127.47	119.78
1	A	52	ALA	C-N-CA	6.21	127.47	119.78
1	A	446	GLY	N-CA-C	-6.07	106.87	115.43
1	A	250	SER	N-CA-C	-6.03	107.75	114.62
1	B	26	THR	CA-CB-OG1	-5.98	100.64	109.60
1	B	406	ASP	CA-CB-CG	5.81	118.41	112.60
1	A	265	GLU	CB-CG-CD	5.66	122.22	112.60
1	A	406	ASP	CA-CB-CG	5.66	118.26	112.60
1	A	76	ASP	CA-CB-CG	5.64	118.24	112.60
1	A	392	VAL	N-CA-C	-5.60	107.00	111.81
1	B	471	ARG	CB-CG-CD	5.54	124.05	111.30
1	B	276	ARG	CB-CG-CD	5.49	123.94	111.30
1	B	265	GLU	CB-CG-CD	5.49	121.92	112.60
1	A	499	ASP	CA-CB-CG	5.47	118.07	112.60
1	B	49	GLN	CB-CA-C	-5.39	110.34	116.54
1	A	276	ARG	CB-CG-CD	5.12	123.08	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	GLU	CB-CG-CD	5.08	121.23	112.60
1	B	52	ALA	CA-C-N	5.07	126.06	119.78
1	B	52	ALA	C-N-CA	5.07	126.06	119.78

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	5030	0	4903	75	0
1	B	5100	0	4970	65	0
2	A	24	0	0	0	0
2	B	16	0	0	0	0
All	All	10170	0	9873	139	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 7.

All (139) 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:339:GLY:HA2	1:B:343:HIS:CD2	1.44	1.50
1:B:339:GLY:CA	1:B:343:HIS:CD2	2.24	1.20
1:A:39:TRP:HB3	1:A:43:ARG:HH12	1.02	1.15
1:A:36:GLU:HG2	1:A:37:ASP:H	1.11	1.14
1:A:344:HIS:CE1	1:A:383:PRO:HB3	1.86	1.09
1:A:79:ARG:HG3	1:A:320:ALA:HA	1.34	1.07
1:B:79:ARG:HG3	1:B:320:ALA:HA	1.35	1.07
1:A:39:TRP:HB3	1:A:43:ARG:NH1	1.73	1.01
1:B:49:GLN:HE21	1:B:51:ARG:HA	1.27	0.98
1:A:344:HIS:CE1	1:A:383:PRO:CB	2.46	0.97
1:A:91:MET:HE1	1:A:122:TYR:HB2	1.48	0.94
1:B:339:GLY:HA2	1:B:343:HIS:HD2	1.12	0.92
1:A:397:ASP:OD2	1:A:449:ARG:HD3	1.71	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:MET:HE3	1:A:642:TRP:HB3	1.50	0.90
1:B:339:GLY:CA	1:B:343:HIS:HD2	1.72	0.90
1:B:339:GLY:HA2	1:B:343:HIS:NE2	1.87	0.89
1:A:36:GLU:HG2	1:A:37:ASP:N	1.89	0.86
1:B:647:GLU:HG3	1:B:648:ALA:H	1.40	0.86
1:B:49:GLN:HG3	1:B:51:ARG:H	1.41	0.85
1:A:444:VAL:HG13	1:A:445:ASN:H	1.44	0.82
1:A:39:TRP:CB	1:A:43:ARG:HH12	1.90	0.82
1:B:37:ASP:OD1	1:B:346:LYS:HD3	1.79	0.82
1:B:49:GLN:HE21	1:B:51:ARG:CA	1.93	0.81
1:B:49:GLN:NE2	1:B:51:ARG:HA	1.96	0.80
1:A:82:ARG:HD3	1:A:88:ARG:HD2	1.70	0.74
1:A:344:HIS:HE1	1:A:383:PRO:HA	1.50	0.74
1:B:558:ARG:HG2	1:B:558:ARG:O	1.88	0.72
1:B:50:VAL:HG13	1:B:52:ALA:H	1.53	0.72
1:B:292:ARG:HH11	1:B:300:GLN:HE21	1.38	0.71
1:A:130:ARG:HD2	1:A:143:GLN:HE21	1.54	0.71
1:A:444:VAL:HG13	1:A:445:ASN:N	2.06	0.70
1:A:9:GLU:HG2	1:A:10:LEU:N	2.07	0.69
1:A:344:HIS:HE1	1:A:383:PRO:CA	2.05	0.69
1:A:591:MET:CE	1:A:642:TRP:HB3	2.23	0.68
1:A:36:GLU:CG	1:A:37:ASP:H	1.99	0.67
1:B:91:MET:HE1	1:B:122:TYR:HB2	1.77	0.66
1:B:130:ARG:HD2	1:B:143:GLN:HE21	1.60	0.66
1:A:344:HIS:CE1	1:A:383:PRO:HA	2.31	0.66
1:A:344:HIS:CE1	1:A:383:PRO:CA	2.78	0.65
1:A:377:ARG:HH11	1:A:377:ARG:HG2	1.62	0.64
1:B:647:GLU:HG3	1:B:648:ALA:N	2.11	0.64
1:A:616:HIS:CE1	1:A:627:ARG:HD3	2.32	0.64
1:A:620:ARG:HH11	1:A:620:ARG:HG2	1.63	0.64
1:B:423:ASP:OD1	1:B:428:GLN:HB2	2.01	0.61
1:A:558:ARG:HG2	1:A:558:ARG:O	1.99	0.61
1:A:344:HIS:CE1	1:A:383:PRO:CG	2.84	0.60
1:A:344:HIS:ND1	1:A:383:PRO:HG3	2.17	0.59
1:A:620:ARG:HG2	1:A:620:ARG:NH1	2.17	0.59
1:A:344:HIS:CE1	1:A:383:PRO:HG3	2.37	0.59
1:A:79:ARG:HG3	1:A:320:ALA:CA	2.22	0.58
1:B:292:ARG:HH11	1:B:300:GLN:NE2	2.01	0.58
1:B:339:GLY:HA3	1:B:343:HIS:CD2	2.32	0.58
1:B:222:PHE:HE2	1:B:248:PHE:HA	1.68	0.58
1:B:173:VAL:HG22	1:B:280:VAL:HB	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:ARG:HG2	1:A:377:ARG:NH1	2.19	0.57
1:B:292:ARG:NH1	1:B:300:GLN:HE21	2.03	0.56
1:A:39:TRP:CB	1:A:43:ARG:NH1	2.58	0.56
1:A:131:PRO:HD2	1:A:149:ARG:HG3	1.87	0.56
1:A:173:VAL:HG22	1:A:280:VAL:HB	1.86	0.56
1:B:6:LEU:O	1:B:6:LEU:HD23	2.07	0.54
1:B:75:LEU:O	1:B:79:ARG:HG2	2.07	0.54
1:B:339:GLY:HA3	1:B:343:HIS:HD2	1.67	0.53
1:B:49:GLN:HG3	1:B:51:ARG:N	2.16	0.53
1:A:627:ARG:HE	1:A:629:ASP:HB2	1.75	0.52
1:B:130:ARG:HD2	1:B:143:GLN:NE2	2.25	0.51
1:A:75:LEU:O	1:A:79:ARG:HG2	2.09	0.50
1:A:86:LEU:HD13	1:A:350:MET:SD	2.52	0.50
1:A:203:HIS:HB2	1:A:245:TRP:HB2	1.93	0.50
1:A:307:THR:HG21	1:A:347:VAL:O	2.12	0.49
1:B:433:PHE:O	1:B:459:GLY:HA2	2.12	0.49
1:A:36:GLU:CG	1:A:37:ASP:N	2.63	0.49
1:B:79:ARG:HG3	1:B:320:ALA:CA	2.25	0.48
1:B:199:ARG:HH12	1:B:236:ILE:HG23	1.78	0.48
1:B:533:GLU:HG3	1:B:536:ARG:HH21	1.78	0.48
1:B:360:LEU:HD21	1:B:485:VAL:HG11	1.95	0.47
1:A:349:ASP:O	1:A:350:MET:HG3	2.14	0.47
1:B:131:PRO:HD2	1:B:149:ARG:HG3	1.97	0.47
1:A:444:VAL:CG1	1:A:445:ASN:N	2.77	0.47
1:B:58:LEU:HD12	1:B:61:MET:HE2	1.97	0.47
1:B:267:VAL:O	1:B:271:LEU:HG	2.15	0.46
1:A:350:MET:HA	1:A:387:THR:O	2.16	0.46
1:A:433:PHE:O	1:A:459:GLY:HA2	2.14	0.46
1:A:413:ASN:O	1:A:414:GLY:C	2.59	0.46
1:A:124:HIS:HA	1:A:173:VAL:HB	1.99	0.45
1:A:9:GLU:HG2	1:A:10:LEU:H	1.81	0.45
1:B:234:GLU:O	1:B:241:GLY:HA2	2.16	0.45
1:A:349:ASP:C	1:A:350:MET:HG3	2.41	0.45
1:B:647:GLU:CG	1:B:648:ALA:H	2.21	0.45
1:B:560:THR:HG21	1:B:591:MET:HE1	1.99	0.45
1:B:598:SER:OG	1:B:600:GLU:HG3	2.16	0.44
1:A:444:VAL:CG1	1:A:445:ASN:H	2.22	0.44
1:A:586:HIS:CG	1:A:587:PRO:HD2	2.53	0.44
1:A:175:ASP:HB3	1:A:284:ASP:OD2	2.18	0.44
1:B:100:ARG:HD2	1:B:521:TRP:HZ3	1.81	0.44
1:A:617:VAL:HG23	1:A:626:PHE:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:HIS:HB2	1:B:245:TRP:HB2	1.99	0.44
1:A:36:GLU:O	1:A:37:ASP:C	2.61	0.44
1:A:131:PRO:CD	1:A:149:ARG:HG3	2.47	0.44
1:B:335:ILE:HD12	1:B:335:ILE:HA	1.85	0.43
1:A:236:ILE:HD12	1:A:236:ILE:HA	1.71	0.43
1:B:131:PRO:CD	1:B:149:ARG:HG3	2.48	0.43
1:B:124:HIS:HA	1:B:173:VAL:HB	2.01	0.43
1:A:91:MET:HE1	1:A:122:TYR:CB	2.34	0.43
1:A:282:ARG:NH2	1:A:326:GLU:HG3	2.34	0.43
1:A:292:ARG:HG2	1:A:295:THR:HG23	2.00	0.43
1:B:51:ARG:O	1:B:51:ARG:HG3	2.19	0.43
1:B:647:GLU:OE2	1:B:647:GLU:HA	2.19	0.43
1:B:222:PHE:CE2	1:B:248:PHE:HA	2.52	0.42
1:B:58:LEU:HD12	1:B:61:MET:CE	2.49	0.42
1:B:423:ASP:OD1	1:B:428:GLN:CB	2.67	0.42
1:B:613:LEU:HD23	1:B:613:LEU:HA	1.90	0.42
1:A:413:ASN:HB3	1:A:416:ALA:HB3	2.02	0.42
1:B:36:GLU:H	1:B:36:GLU:HG2	1.56	0.42
1:A:311:ARG:NH2	1:A:315:ARG:CZ	2.83	0.42
1:A:391:TYR:N	1:A:391:TYR:CD1	2.88	0.41
1:B:400:TRP:O	1:B:418:ARG:HD2	2.20	0.41
1:B:586:HIS:CG	1:B:587:PRO:HD2	2.55	0.41
1:A:100:ARG:HD2	1:A:521:TRP:HZ3	1.85	0.41
1:B:282:ARG:NH2	1:B:326:GLU:HG3	2.35	0.41
1:A:247:THR:HG23	1:A:254:ASP:OD2	2.20	0.41
1:B:175:ASP:HB3	1:B:284:ASP:OD2	2.19	0.41
1:A:360:LEU:HB3	1:A:482:LEU:HD22	2.02	0.41
1:B:616:HIS:CE1	1:B:648:ALA:HB2	2.56	0.41
1:A:552:HIS:O	1:A:556:VAL:HG23	2.21	0.41
1:A:605:PRO:O	1:A:608:VAL:HG13	2.20	0.41
1:A:616:HIS:HE1	1:A:627:ARG:HD3	1.82	0.41
1:B:509:PHE:CE2	1:B:519:ASN:HA	2.56	0.41
1:A:53:LEU:HB3	1:A:54:PRO:HD3	2.03	0.41
1:A:617:VAL:CG2	1:A:626:PHE:HB2	2.51	0.41
1:A:360:LEU:HD21	1:A:485:VAL:HG11	2.02	0.41
1:A:440:GLN:HB2	1:A:449:ARG:HG3	2.03	0.41
1:A:518:ASP:OD2	1:A:520:ARG:NH2	2.54	0.41
1:B:49:GLN:HE21	1:B:51:ARG:N	2.18	0.41
1:B:118:LEU:HD23	1:B:118:LEU:HA	1.91	0.41
1:B:405:GLU:OE1	1:B:405:GLU:N	2.53	0.41
1:B:31:LEU:HD21	1:B:58:LEU:HD11	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:LEU:H	1:A:10:LEU:HG	1.72	0.40
1:A:74:ARG:HD3	1:B:25:GLU:HG3	2.03	0.40
1:B:50:VAL:HG13	1:B:52:ALA:HB3	2.04	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	623/645 (97%)	597 (96%)	25 (4%)	1 (0%)	43	68
1	B	633/645 (98%)	609 (96%)	23 (4%)	1 (0%)	43	68
All	All	1256/1290 (97%)	1206 (96%)	48 (4%)	2 (0%)	43	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	347	VAL
1	A	206	PRO

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	513/524 (98%)	495 (96%)	18 (4%)	32	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	520/524 (99%)	504 (97%)	16 (3%)	35	65
All	All	1033/1048 (99%)	999 (97%)	34 (3%)	33	63

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

Mol	Chain	Res	Type
1	A	8	SER
1	A	59	GLU
1	A	91	MET
1	A	135	GLU
1	A	163	ARG
1	A	236	ILE
1	A	256	ASN
1	A	326	GLU
1	A	328	ILE
1	A	349	ASP
1	A	409	ARG
1	A	445	ASN
1	A	533	GLU
1	A	548	THR
1	A	600	GLU
1	A	604	LEU
1	A	608	VAL
1	A	627	ARG
1	B	8	SER
1	B	9	GLU
1	B	36	GLU
1	B	135	GLU
1	B	256	ASN
1	B	301	PRO
1	B	311	ARG
1	B	326	GLU
1	B	344	HIS
1	B	409	ARG
1	B	412	LEU
1	B	511	ASP
1	B	600	GLU
1	B	610	ARG
1	B	612	VAL
1	B	633	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	124	HIS
1	A	143	GLN
1	A	179	ASN
1	A	249	ASN
1	A	299	ASN
1	A	305	HIS
1	A	445	ASN
1	A	547	ASN
1	A	616	HIS
1	B	49	GLN
1	B	143	GLN
1	B	179	ASN
1	B	299	ASN
1	B	300	GLN
1	B	442	ASN
1	B	445	ASN
1	B	547	ASN
1	B	570	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 ⓘ

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	628/645 (97%)	-0.17	33 (5%) 32 28	11, 43, 84, 128	1 (0%)
1	B	639/645 (99%)	-0.17	38 (5%) 28 25	26, 47, 92, 147	0
All	All	1267/1290 (98%)	-0.17	71 (5%) 30 26	11, 45, 87, 147	1 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	347	VAL	9.2
1	A	348	SER	7.5
1	B	343	HIS	7.4
1	B	223	PRO	6.4
1	B	344	HIS	5.6
1	B	351	ALA	5.0
1	B	345	GLY	4.9
1	B	650	ALA	4.6
1	B	649	PRO	4.6
1	A	346	LYS	4.5
1	B	50	VAL	4.4
1	B	226	ALA	4.4
1	A	442	ASN	4.2
1	A	219	PRO	4.0
1	B	248	PHE	4.0
1	A	646	GLY	3.9
1	A	8	SER	3.8
1	B	209	ARG	3.8
1	A	7	THR	3.8
1	A	12	ALA	3.7
1	B	342	ALA	3.6
1	A	9	GLU	3.5
1	A	616	HIS	3.5
1	B	618	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	227	PRO	3.3
1	A	43	ARG	3.3
1	A	443	PRO	3.3
1	A	444	VAL	3.2
1	A	344	HIS	3.2
1	B	509	PHE	3.1
1	A	600	GLU	3.1
1	A	10	LEU	3.1
1	B	444	VAL	3.0
1	A	250	SER	3.0
1	B	349	ASP	3.0
1	A	248	PHE	3.0
1	B	6	LEU	2.9
1	B	207	ASP	2.9
1	B	348	SER	2.8
1	B	648	ALA	2.8
1	A	634	ARG	2.8
1	B	440	GLN	2.8
1	A	445	ASN	2.8
1	B	232	TRP	2.8
1	A	249	ASN	2.8
1	B	507	TYR	2.8
1	A	236	ILE	2.7
1	B	241	GLY	2.7
1	A	507	TYR	2.6
1	B	510	GLU	2.6
1	B	206	PRO	2.6
1	A	49	GLN	2.5
1	A	350	MET	2.5
1	B	51	ARG	2.4
1	B	514	GLU	2.4
1	B	250	SER	2.4
1	A	625	ALA	2.4
1	A	626	PHE	2.4
1	B	140	TYR	2.3
1	B	428	GLN	2.3
1	A	627	ARG	2.3
1	B	222	PHE	2.3
1	B	219	PRO	2.2
1	B	346	LYS	2.2
1	B	341	ARG	2.1
1	B	249	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	508	ALA	2.1
1	A	205	PHE	2.1
1	A	511	ASP	2.0
1	A	349	ASP	2.0
1	A	228	GLY	2.0

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.