



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:26 AM UTC

PDB ID : 5SDS / pdb_00005sds
Title : PanDDA analysis group deposition of ground-state model of Porphyromonas gingivalis DPP11
Authors : Tham, C.T.; Coker, J.A.; Krojer, T.; Foster, W.R.; Koekemoer, L.; Douangamath, A.; Talon, R.; Fearon, D.; von Delft, F.; Yue, W.W.; Bountra, C.; Bezerra, G.A.
Deposited on : 2022-01-20
Resolution : 1.91 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

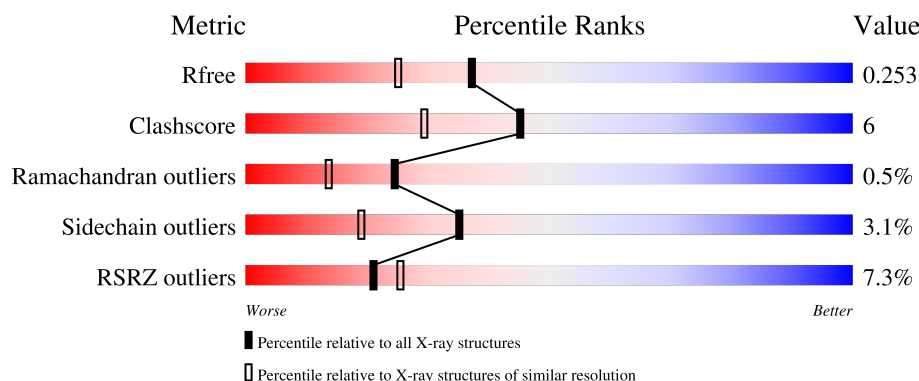
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

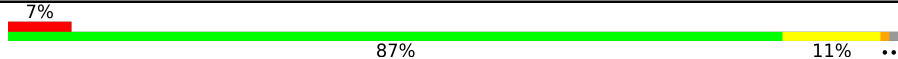
The reported resolution of this entry is 1.91 Å.

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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	707	
1	B	707	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11220 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 Asp/Glu-specific dipeptidyl-peptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	697	Total	C	N	O	S	0	0	0
			5493	3488	948	1030	27			
1	B	695	Total	C	N	O	S	0	0	0
			5458	3465	940	1026	27			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	initiating methionine	UNP B2RID1
A	721	ALA	-	expression tag	UNP B2RID1
A	722	HIS	-	expression tag	UNP B2RID1
A	723	HIS	-	expression tag	UNP B2RID1
A	724	HIS	-	expression tag	UNP B2RID1
A	725	HIS	-	expression tag	UNP B2RID1
A	726	HIS	-	expression tag	UNP B2RID1
A	727	HIS	-	expression tag	UNP B2RID1
B	21	MET	-	initiating methionine	UNP B2RID1
B	721	ALA	-	expression tag	UNP B2RID1
B	722	HIS	-	expression tag	UNP B2RID1
B	723	HIS	-	expression tag	UNP B2RID1
B	724	HIS	-	expression tag	UNP B2RID1
B	725	HIS	-	expression tag	UNP B2RID1
B	726	HIS	-	expression tag	UNP B2RID1
B	727	HIS	-	expression tag	UNP B2RID1

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cl	0	0
			2	2		
2	B	1	Total	Cl	0	0
			1	1		

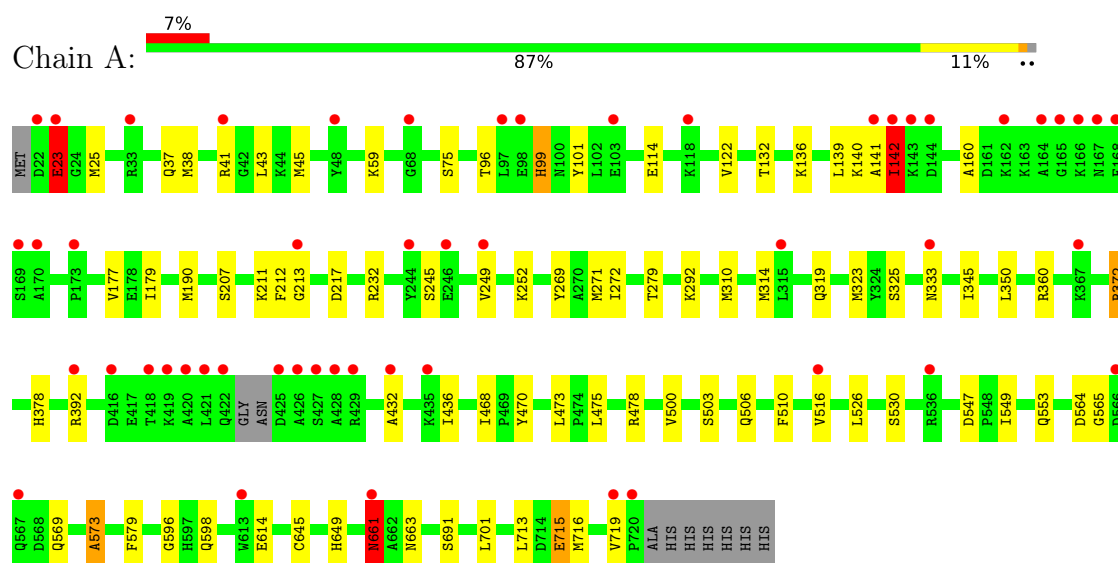
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	140	Total 140	O 140	0	0
3	B	126	Total 126	O 126	0	0

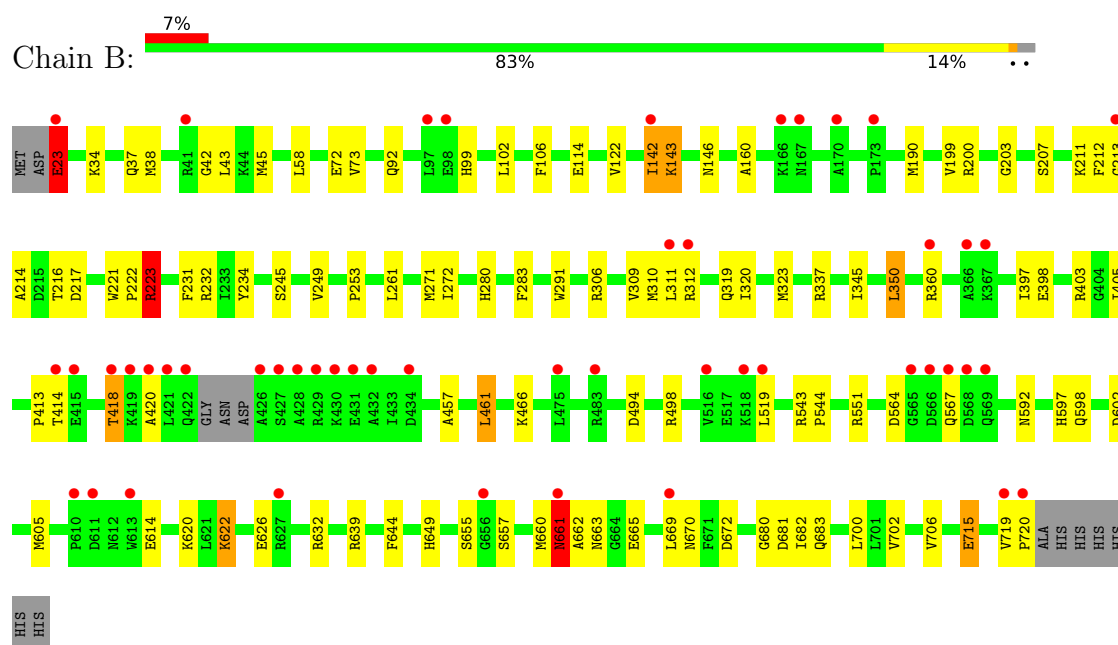
3 Residue-property plots [i](#)

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: Asp/Glu-specific dipeptidyl-peptidase



- Molecule 1: Asp/Glu-specific dipeptidyl-peptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.77Å 117.33Å 147.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.83 – 1.91 91.83 – 1.91	Depositor EDS
% Data completeness (in resolution range)	73.4 (91.83-1.91) 73.8 (91.83-1.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.213 , 0.247 0.220 , 0.253	Depositor DCC
R_{free} test set	6987 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 22.3	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	0.95	EDS
Total number of atoms	11220	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CL

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	1.03	1/5617 (0.0%)	1.38	3/7607 (0.0%)
1	B	1.03	1/5581 (0.0%)	1.38	8/7561 (0.1%)
All	All	1.03	2/11198 (0.0%)	1.38	11/15168 (0.1%)

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	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	223	ARG	NE-CZ	-5.21	1.27	1.33
1	A	468	ILE	N-CA	5.03	1.50	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	223	ARG	CB-CG-CD	8.98	131.96	111.30
1	A	661	ASN	CB-CA-C	-6.71	97.06	110.42
1	B	661	ASN	CA-CB-CG	-5.90	106.70	112.60
1	B	23	GLU	CB-CG-CD	5.26	121.54	112.60
1	B	681	ASP	CA-CB-CG	5.26	117.86	112.60
1	A	565	GLY	CA-C-O	-5.25	118.04	122.29
1	B	661	ASN	CB-CA-C	-5.23	100.01	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	418	THR	CB-CA-C	5.21	119.72	110.85
1	A	573	ALA	CA-C-O	-5.02	116.02	121.44
1	B	306	ARG	CA-C-N	5.01	127.00	120.28
1	B	306	ARG	C-N-CA	5.01	127.00	120.28

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ILE	Peptide
1	A	23	GLU	Peptide
1	B	660	MET	Peptide

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	5493	0	5282	61	0
1	B	5458	0	5231	67	0
2	A	2	0	0	1	0
2	B	1	0	0	0	0
3	A	140	0	0	6	1
3	B	126	0	0	5	1
All	All	11220	0	10513	127	1

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 6.

All (127) 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:269:TYR:OH	1:A:271:MET:HE3	1.32	1.23
1:A:271:MET:SD	3:A:1029:HOH:O	2.03	1.15
1:B:142:ILE:HA	3:B:1013:HOH:O	1.61	0.97
1:A:25:MET:HE1	1:A:573:ALA:HB2	1.46	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:GLN:HG3	1:B:323:MET:HE2	1.49	0.95
1:B:23:GLU:O	1:B:649:HIS:ND1	2.02	0.93
1:B:337:ARG:NH1	3:B:901:HOH:O	2.08	0.86
1:B:223:ARG:NH2	1:B:672:ASP:OD2	2.11	0.84
1:A:160:ALA:HB1	1:A:177:VAL:HG12	1.64	0.78
1:B:99:HIS:CB	3:B:1014:HOH:O	2.32	0.78
1:A:25:MET:CE	1:A:573:ALA:HB2	2.15	0.75
1:A:269:TYR:OH	1:A:271:MET:CE	2.26	0.75
1:A:319:GLN:HG3	1:A:323:MET:HE2	1.68	0.73
1:A:23:GLU:O	1:A:649:HIS:ND1	2.18	0.73
1:B:320:ILE:HD13	1:B:323:MET:HE3	1.71	0.72
1:B:655:SER:O	1:B:670:ASN:O	2.08	0.72
1:B:280:HIS:HB2	1:B:682:ILE:CD1	2.23	0.68
1:B:37:GLN:HG3	3:B:964:HOH:O	1.91	0.68
1:A:114:GLU:OE2	1:A:232:ARG:NH1	2.27	0.68
1:B:626:GLU:O	3:B:902:HOH:O	2.12	0.68
1:B:160:ALA:HA	1:B:190:MET:CE	2.25	0.67
1:B:494:ASP:OD2	1:B:498:ARG:HD3	1.96	0.66
1:A:212:PHE:O	1:A:614:GLU:O	2.13	0.66
1:B:114:GLU:OE2	1:B:232:ARG:NH1	2.28	0.65
1:A:564:ASP:HB2	1:A:569:GLN:NE2	2.13	0.63
1:A:392:ARG:NH1	3:A:906:HOH:O	2.32	0.63
1:A:232:ARG:NH2	1:A:715:GLU:OE2	2.32	0.62
1:A:661:ASN:HB3	1:A:663:ASN:H	1.65	0.62
1:B:661:ASN:HB3	1:B:663:ASN:H	1.64	0.62
1:B:43:LEU:HD11	1:B:45:MET:HG2	1.81	0.61
1:A:213:GLY:HA3	1:A:217:ASP:HB2	1.82	0.61
1:A:160:ALA:HA	1:A:190:MET:CE	2.31	0.61
1:B:605:MET:HE1	1:B:622:LYS:HG2	1.82	0.61
1:A:38:MET:CE	1:A:271:MET:CE	2.80	0.60
1:B:34:LYS:NZ	1:B:567:GLN:O	2.27	0.60
1:B:212:PHE:O	1:B:614:GLU:O	2.20	0.60
1:A:269:TYR:HH	1:A:271:MET:HE3	1.58	0.60
1:A:319:GLN:HG3	1:A:323:MET:CE	2.33	0.59
1:B:232:ARG:NH2	1:B:715:GLU:OE1	2.35	0.59
1:B:72:GLU:OE2	1:B:669:LEU:HD13	2.02	0.58
1:A:38:MET:CE	1:A:271:MET:HE1	2.34	0.57
1:A:37:GLN:HB3	3:A:988:HOH:O	2.04	0.56
1:B:232:ARG:NH2	1:B:715:GLU:OE2	2.36	0.56
1:B:160:ALA:HA	1:B:190:MET:HE1	1.87	0.56
1:B:216:THR:OG1	1:B:614:GLU:OE2	2.24	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:GLY:HA3	1:B:217:ASP:HB2	1.89	0.54
1:B:320:ILE:HD13	1:B:323:MET:CE	2.37	0.54
1:A:232:ARG:NH2	1:A:715:GLU:CD	2.67	0.53
1:A:526:LEU:O	1:A:530:SER:HB2	2.08	0.53
1:A:314:MET:HE1	1:A:325:SER:HA	1.91	0.52
1:A:232:ARG:NH2	1:A:715:GLU:OE1	2.42	0.52
1:B:232:ARG:NH2	1:B:715:GLU:CD	2.68	0.52
1:A:139:LEU:O	1:A:142:ILE:HG13	2.11	0.51
1:A:475:LEU:HA	1:A:478:ARG:HD2	1.92	0.51
1:B:661:ASN:ND2	1:B:665:GLU:OE1	2.44	0.51
1:A:310:MET:O	1:A:314:MET:HG3	2.12	0.50
1:B:312:ARG:HG3	1:B:312:ARG:HH11	1.77	0.50
1:B:661:ASN:HD22	1:B:665:GLU:HB2	1.77	0.50
1:A:547:ASP:CG	3:A:965:HOH:O	2.55	0.50
1:B:200:ARG:HD2	1:B:234:TYR:CE1	2.46	0.49
1:A:564:ASP:HB2	1:A:569:GLN:HE22	1.75	0.49
1:A:503:SER:OG	1:A:506:GLN:HG3	2.13	0.49
1:B:261:LEU:HD13	1:B:632:ARG:HG3	1.95	0.48
1:B:309:VAL:HA	1:B:312:ARG:HD2	1.94	0.48
1:A:292:LYS:HB2	1:A:345:ILE:HD13	1.96	0.48
1:B:211:LYS:HG2	1:B:214:ALA:HA	1.96	0.48
1:B:413:PRO:CG	1:B:519:LEU:HD13	2.44	0.48
1:B:73:VAL:HG11	1:B:253:PRO:HG3	1.95	0.47
1:B:661:ASN:ND2	1:B:665:GLU:HB2	2.29	0.47
1:B:114:GLU:CD	1:B:232:ARG:HH11	2.22	0.47
1:A:38:MET:HE2	1:A:271:MET:HE1	1.96	0.47
1:A:269:TYR:CZ	1:A:271:MET:HE3	2.40	0.47
1:A:140:LYS:C	1:A:142:ILE:H	2.23	0.47
1:B:142:ILE:O	1:B:143:LYS:CB	2.63	0.46
1:B:38:MET:CE	1:B:271:MET:HE1	2.46	0.46
1:B:146:ASN:HB2	1:B:551:ARG:CZ	2.44	0.46
1:B:92:GLN:HG3	1:B:102:LEU:HD22	1.98	0.46
1:B:398:GLU:O	1:B:403:ARG:HG2	2.16	0.46
1:A:470:TYR:HA	1:A:473:LEU:HD12	1.97	0.46
1:A:43:LEU:HD11	1:A:45:MET:HG2	1.97	0.46
1:B:122:VAL:HG13	1:B:199:VAL:HG21	1.98	0.46
1:B:414:THR:O	1:B:418:THR:HG22	2.17	0.45
1:B:291:TRP:CZ3	1:B:682:ILE:HG13	2.52	0.45
1:A:59:LYS:HD2	1:A:59:LYS:C	2.42	0.45
1:A:713:LEU:HA	1:A:716:MET:HE2	1.99	0.45
1:B:261:LEU:HD21	1:B:700:LEU:HD12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ARG:NE	1:A:41:ARG:HA	2.32	0.44
1:B:719:VAL:O	1:B:720:PRO:C	2.61	0.44
1:A:25:MET:HE1	1:A:279:THR:CG2	2.48	0.44
1:B:221:TRP:CD2	1:B:222:PRO:HA	2.53	0.44
1:B:405:ILE:HD11	1:B:461:LEU:HD13	2.00	0.43
1:B:350:LEU:HD22	1:B:683:GLN:HB3	2.00	0.43
1:A:160:ALA:HB1	1:A:177:VAL:CG1	2.43	0.43
1:A:99:HIS:HD2	1:A:101:TYR:OH	2.01	0.42
1:A:432:ALA:O	1:A:436:ILE:HG13	2.20	0.42
1:A:25:MET:HE1	1:A:279:THR:HG23	2.01	0.42
1:A:549:ILE:O	1:A:553:GLN:HG3	2.20	0.42
1:B:602:ASP:HB2	1:B:639:ARG:HH21	1.84	0.42
1:A:179:ILE:HD11	1:A:190:MET:HE3	2.01	0.42
1:A:378:HIS:CE1	3:A:967:HOH:O	2.72	0.42
1:A:500:VAL:HG22	1:A:510:PHE:HB2	2.03	0.41
1:B:280:HIS:HB2	1:B:682:ILE:HD12	1.99	0.41
1:B:543:ARG:N	1:B:544:PRO:HD2	2.35	0.41
1:A:372:ARG:HG2	2:A:801:CL:CL	2.57	0.41
1:B:543:ARG:N	1:B:544:PRO:CD	2.83	0.41
1:B:680:GLY:HA2	1:B:683:GLN:O	2.21	0.41
1:A:177:VAL:CG1	1:A:190:MET:HE2	2.50	0.41
1:A:378:HIS:HE1	3:A:967:HOH:O	2.02	0.41
1:A:645:CYS:HA	1:A:691:SER:O	2.21	0.41
1:B:272:ILE:HG23	1:B:657:SER:HB3	2.03	0.41
1:B:291:TRP:CE3	1:B:345:ILE:HD11	2.56	0.41
1:B:106:PHE:O	1:B:203:GLY:HA2	2.21	0.41
1:B:42:GLY:O	1:B:662:ALA:HB1	2.21	0.41
1:B:360:ARG:HH21	1:B:564:ASP:CG	2.28	0.41
1:B:702:VAL:O	1:B:706:VAL:HB	2.20	0.41
1:A:25:MET:HE1	1:A:573:ALA:CB	2.33	0.41
1:A:25:MET:CE	1:A:279:THR:HG21	2.50	0.41
1:A:96:THR:OG1	1:A:99:HIS:ND1	2.52	0.41
1:A:132:THR:CG2	1:A:136:LYS:HE3	2.51	0.41
1:A:160:ALA:HA	1:A:190:MET:HE2	2.02	0.41
1:A:596:GLY:CA	1:B:592:ASN:OD1	2.69	0.41
1:B:200:ARG:O	1:B:231:PHE:HA	2.20	0.41
1:B:597:HIS:O	1:B:644:PHE:HA	2.21	0.41
1:A:75:SER:HB2	1:A:715:GLU:HG3	2.03	0.41
1:A:272:ILE:O	1:A:579:PHE:HA	2.21	0.41
1:B:320:ILE:HA	1:B:323:MET:HE3	2.02	0.40
1:B:457:ALA:O	1:B:461:LEU:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1035:HOH:O	3:B:1024:HOH:O[4_555]	1.79	0.41

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	693/707 (98%)	665 (96%)	25 (4%)	3 (0%)	30	19
1	B	691/707 (98%)	665 (96%)	22 (3%)	4 (1%)	21	10
All	All	1384/1414 (98%)	1330 (96%)	47 (3%)	7 (0%)	24	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	661	ASN
1	A	661	ASN
1	B	143	LYS
1	B	420	ALA
1	B	142	ILE
1	A	141	ALA
1	A	142	ILE

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	560/597 (94%)	543 (97%)	17 (3%)	36	20
1	B	554/597 (93%)	537 (97%)	17 (3%)	35	19
All	All	1114/1194 (93%)	1080 (97%)	34 (3%)	35	19

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

Mol	Chain	Res	Type
1	A	23	GLU
1	A	99	HIS
1	A	122	VAL
1	A	207	SER
1	A	211	LYS
1	A	245	SER
1	A	249	VAL
1	A	252	LYS
1	A	333	ASN
1	A	350	LEU
1	A	360	ARG
1	A	372	ARG
1	A	516	VAL
1	A	598	GLN
1	A	701	LEU
1	A	715	GLU
1	A	719	VAL
1	B	23	GLU
1	B	58	LEU
1	B	207	SER
1	B	223	ARG
1	B	245	SER
1	B	249	VAL
1	B	283	PHE
1	B	310	MET
1	B	311	LEU
1	B	350	LEU
1	B	397	ILE
1	B	461	LEU
1	B	466	LYS
1	B	598	GLN
1	B	620	LYS
1	B	622	LYS
1	B	715	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	157	GLN
1	A	472	ASN
1	A	506	GLN
1	A	569	GLN
1	B	146	ASN
1	B	307	GLN
1	B	355	GLN
1	B	583	GLN
1	B	598	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 ⓘ

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	697/707 (98%)	0.61	52 (7%) 20 24	26, 40, 64, 92	0
1	B	695/707 (98%)	0.46	49 (7%) 22 26	25, 38, 62, 85	0
All	All	1392/1414 (98%)	0.54	101 (7%) 21 25	25, 39, 63, 92	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	428	ALA	6.8
1	A	142	ILE	6.7
1	A	22	ASP	6.0
1	B	420	ALA	5.9
1	A	720	PRO	5.6
1	A	425	ASP	5.2
1	A	167	ASN	5.2
1	B	566	ASP	5.2
1	A	23	GLU	5.2
1	B	720	PRO	5.1
1	A	426	ALA	4.9
1	A	48	TYR	4.5
1	B	613	TRP	4.5
1	A	97	LEU	4.4
1	B	426	ALA	4.3
1	B	415	GLU	4.2
1	A	422	GLN	4.0
1	B	567	GLN	4.0
1	A	170	ALA	3.8
1	B	568	ASP	3.8
1	B	427	SER	3.7
1	A	169	SER	3.6
1	A	427	SER	3.6
1	A	98	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	418	THR	3.6
1	B	167	ASN	3.6
1	B	422	GLN	3.6
1	B	418	THR	3.6
1	A	613	TRP	3.5
1	A	719	VAL	3.5
1	B	166	LYS	3.5
1	A	420	ALA	3.3
1	B	419	LYS	3.2
1	A	566	ASP	3.2
1	B	97	LEU	3.2
1	B	429	ARG	3.2
1	A	213	GLY	3.1
1	A	103	GLU	3.1
1	B	421	LEU	3.0
1	B	431	GLU	3.0
1	B	432	ALA	2.9
1	B	669	LEU	2.9
1	A	249	VAL	2.8
1	A	143	LYS	2.8
1	A	118	LYS	2.8
1	A	516	VAL	2.8
1	B	719	VAL	2.7
1	A	33	ARG	2.7
1	B	516	VAL	2.7
1	B	366	ALA	2.7
1	B	430	LYS	2.7
1	B	41	ARG	2.6
1	A	367	LYS	2.6
1	B	519	LEU	2.6
1	B	656	GLY	2.6
1	A	428	ALA	2.6
1	B	611	ASP	2.6
1	A	166	LYS	2.6
1	B	475	LEU	2.6
1	A	168	PHE	2.6
1	B	170	ALA	2.6
1	A	333	ASN	2.6
1	B	661	ASN	2.6
1	A	567	GLN	2.5
1	A	244	TYR	2.5
1	B	23	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	429	ARG	2.5
1	A	536	ARG	2.5
1	B	142	ILE	2.5
1	B	312	ARG	2.4
1	B	627	ARG	2.4
1	A	164	ALA	2.4
1	B	213	GLY	2.4
1	A	41	ARG	2.3
1	B	414	THR	2.3
1	A	141	ALA	2.3
1	B	367	LYS	2.3
1	A	173	PRO	2.3
1	A	432	ALA	2.3
1	A	246	GLU	2.3
1	A	144	ASP	2.2
1	B	569	GLN	2.2
1	B	434	ASP	2.2
1	A	315	LEU	2.2
1	A	421	LEU	2.2
1	A	416	ASP	2.1
1	B	518	LYS	2.1
1	B	98	GLU	2.1
1	A	661	ASN	2.1
1	A	419	LYS	2.1
1	B	173	PRO	2.1
1	A	435	LYS	2.1
1	A	68	GLY	2.1
1	A	165	GLY	2.1
1	A	392	ARG	2.1
1	B	311	LEU	2.0
1	B	483	ARG	2.0
1	A	162	LYS	2.0
1	B	565	GLY	2.0
1	B	610	PRO	2.0
1	B	360	ARG	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	A	801	1/1	0.85	0.23	71,71,71,71	0
2	CL	B	801	1/1	0.89	0.20	60,60,60,60	0
2	CL	A	802	1/1	0.97	0.08	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.