



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:26 PM UTC

PDB ID : 4N2C / pdb_00004n2c
Title : Crystal structure of Protein Arginine Deiminase 2 (F221/222A, 10 mM Ca²⁺)
Authors : Slade, D.J.; Zhang, X.; Fang, P.; Dreyton, C.J.; Zhang, Y.; Gross, M.L.; Guo, M.; Coonrod, S.A.; Thompson, P.R.
Deposited on : 2013-10-04
Resolution : 3.02 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

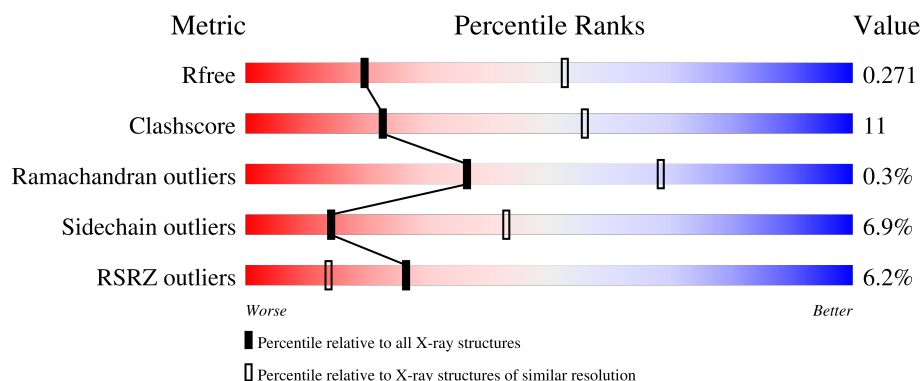
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	690	6% 66% 24% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5039 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 Protein-arginine deiminase type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	645	4997	3204	826	939	28	17	2	0

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q9Y2J8
A	-19	GLY	-	expression tag	UNP Q9Y2J8
A	-18	HIS	-	expression tag	UNP Q9Y2J8
A	-17	HIS	-	expression tag	UNP Q9Y2J8
A	-16	HIS	-	expression tag	UNP Q9Y2J8
A	-15	HIS	-	expression tag	UNP Q9Y2J8
A	-14	HIS	-	expression tag	UNP Q9Y2J8
A	-13	HIS	-	expression tag	UNP Q9Y2J8
A	-12	HIS	-	expression tag	UNP Q9Y2J8
A	-11	HIS	-	expression tag	UNP Q9Y2J8
A	-10	HIS	-	expression tag	UNP Q9Y2J8
A	-9	HIS	-	expression tag	UNP Q9Y2J8
A	-8	SER	-	expression tag	UNP Q9Y2J8
A	-7	SER	-	expression tag	UNP Q9Y2J8
A	-6	GLY	-	expression tag	UNP Q9Y2J8
A	-5	HIS	-	expression tag	UNP Q9Y2J8
A	-4	ILE	-	expression tag	UNP Q9Y2J8
A	-3	GLU	-	expression tag	UNP Q9Y2J8
A	-2	GLY	-	expression tag	UNP Q9Y2J8
A	-1	ARG	-	expression tag	UNP Q9Y2J8
A	0	HIS	-	expression tag	UNP Q9Y2J8
A	221	ALA	PHE	engineered mutation	UNP Q9Y2J8
A	222	ALA	PHE	engineered mutation	UNP Q9Y2J8
A	666	SER	-	expression tag	UNP Q9Y2J8
A	667	ARG	-	expression tag	UNP Q9Y2J8
A	668	ARG	-	expression tag	UNP Q9Y2J8
A	669	SER	-	expression tag	UNP Q9Y2J8

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total 6	Ca 6	0	0

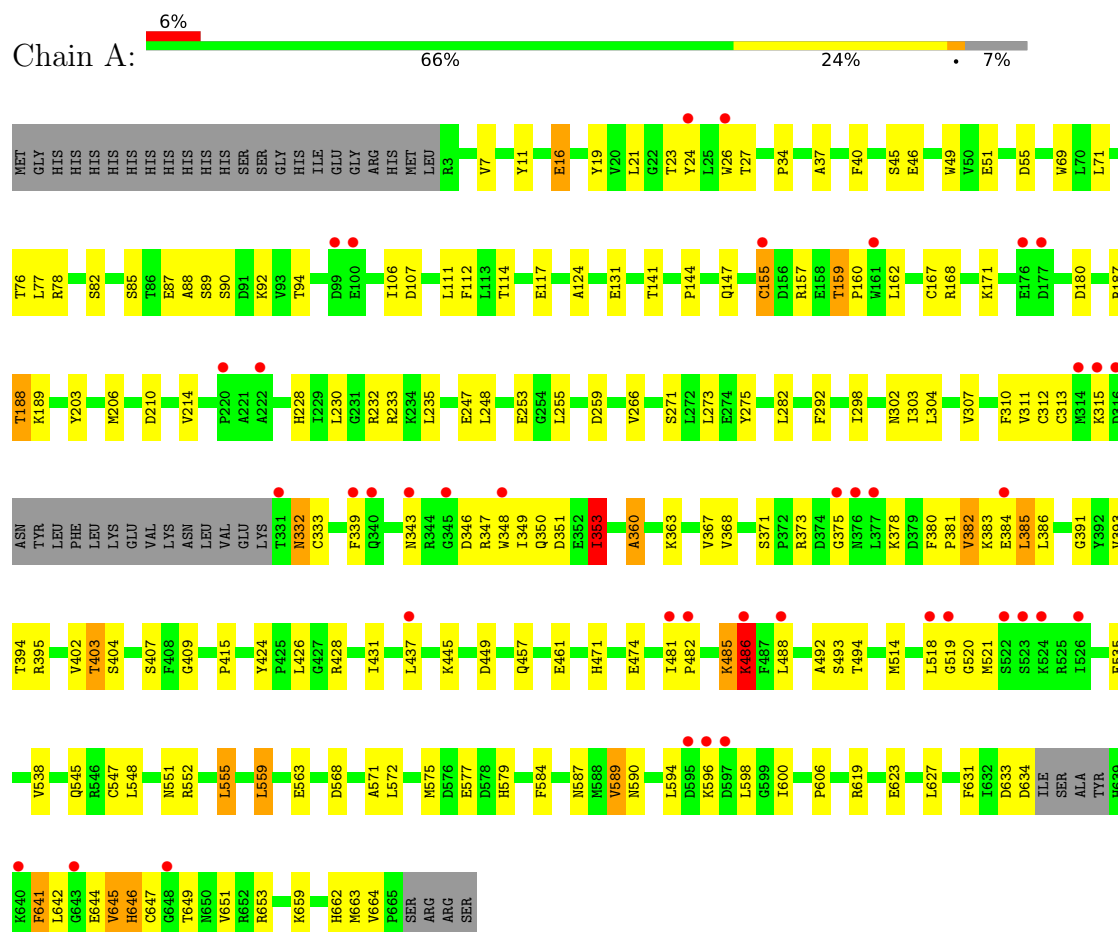
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	36	Total 36	O 36	0	0

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: Protein-arginine deiminase type-2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	202.93Å 52.18Å 75.82Å 90.00° 105.21° 90.00°	Depositor
Resolution (Å)	48.95 – 3.02 48.95 – 3.02	Depositor EDS
% Data completeness (in resolution range)	88.3 (48.95-3.02) 88.3 (48.95-3.02)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.76 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.7.2 _869	Depositor
R, R_{free}	0.245 , 0.273 0.242 , 0.271	Depositor DCC
R_{free} test set	659 reflections (4.27%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	5039	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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: CA

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	0.36	2/5123 (0.0%)	0.84	15/6974 (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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	471	HIS	CG-ND1	-6.18	1.31	1.38
1	A	471	HIS	CD2-NE2	-5.91	1.31	1.37

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	385	LEU	N-CA-C	-7.62	104.00	113.38
1	A	641	PHE	N-CA-C	7.19	120.53	111.69
1	A	518	LEU	CB-CA-C	-6.58	109.00	116.63
1	A	159	THR	CA-C-N	6.09	125.77	119.56
1	A	159	THR	C-N-CA	6.09	125.77	119.56
1	A	424	TYR	CA-C-N	5.99	125.54	119.19
1	A	424	TYR	C-N-CA	5.99	125.54	119.19
1	A	371	SER	CA-C-N	5.85	125.56	119.82
1	A	371	SER	C-N-CA	5.85	125.56	119.82
1	A	360	ALA	CA-C-N	-5.58	113.81	119.83
1	A	360	ALA	C-N-CA	-5.58	113.81	119.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	409	GLY	N-CA-C	-5.42	107.61	114.16
1	A	353	ILE	N-CA-C	5.39	116.42	108.45
1	A	313	CYS	N-CA-C	5.20	117.90	108.69
1	A	646	HIS	N-CA-C	-5.00	105.54	111.69

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	339	PHE	Peptide
1	A	486	LYS	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	4997	0	4808	103	0
2	A	6	0	0	1	0
3	A	36	0	0	2	0
All	All	5039	0	4808	103	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 11.

All (103) 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:253:GLU:OE1	3:A:834:HOH:O	1.74	1.02
1:A:381:PRO:O	1:A:385:LEU:N	2.08	0.86
1:A:380:PHE:CD2	1:A:380:PHE:O	2.30	0.85
1:A:485:LYS:HE2	1:A:486:LYS:H	1.51	0.75
1:A:380:PHE:CD2	1:A:380:PHE:C	2.70	0.69
1:A:555:LEU:HD12	1:A:559:LEU:HD22	1.76	0.67
1:A:259:ASP:OD2	1:A:395:ARG:NH1	2.32	0.63
1:A:144:PRO:HA	1:A:662:HIS:HD2	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:GLU:HG3	1:A:112:PHE:HB2	1.80	0.62
1:A:375:GLY:O	1:A:378:LYS:HG2	2.01	0.61
1:A:383:LYS:N	1:A:384:GLU:HA	2.14	0.61
1:A:303:ILE:HD13	1:A:415:PRO:HB3	1.83	0.61
1:A:619:ARG:NH1	1:A:623:GLU:OE2	2.33	0.60
1:A:481:ILE:HD13	1:A:488:LEU:HG	1.83	0.60
1:A:180:ASP:OD1	2:A:701:CA:CA	1.79	0.59
1:A:157:ARG:NH2	1:A:386:LEU:O	2.36	0.58
1:A:481:ILE:HG13	1:A:481:ILE:O	2.03	0.58
1:A:304:LEU:HB2	1:A:653:ARG:HG2	1.87	0.57
1:A:373:ARG:HA	1:A:402:VAL:HG11	1.87	0.57
1:A:634:ASP:OD1	1:A:644:GLU:HG3	2.05	0.56
1:A:37:ALA:HB2	1:A:106:ILE:HD11	1.86	0.55
1:A:382:VAL:C	1:A:384:GLU:HA	2.32	0.55
1:A:311:VAL:HG22	1:A:312:CYS:O	2.07	0.55
1:A:89:SER:O	1:A:90:SER:HB3	2.07	0.54
1:A:351:ASP:HB3	1:A:407:SER:HB3	1.89	0.54
1:A:403:THR:OG1	1:A:404:SER:N	2.37	0.54
1:A:547:CYS:O	1:A:551:ASN:ND2	2.40	0.54
1:A:19:TYR:OH	1:A:23:THR:O	2.21	0.54
1:A:27:THR:HB	1:A:77:LEU:HB2	1.90	0.54
1:A:302:ASN:OD1	1:A:653:ARG:NH1	2.41	0.53
1:A:124:ALA:N	1:A:131:GLU:OE1	2.42	0.53
1:A:232:ARG:CZ	1:A:255:LEU:HD13	2.39	0.53
1:A:106:ILE:HG13	1:A:107:ASP:H	1.74	0.53
1:A:492:ALA:HB2	1:A:589:VAL:HG21	1.90	0.52
1:A:600:ILE:O	1:A:631:PHE:HA	2.09	0.52
1:A:428:ARG:HA	1:A:457:GLN:O	2.09	0.52
1:A:514:MET:HG2	1:A:606:PRO:HD3	1.92	0.52
1:A:431:ILE:HG13	1:A:461:GLU:HG2	1.91	0.52
1:A:348:TRP:CE3	1:A:647:CYS:HB3	2.45	0.52
1:A:598:LEU:HD21	1:A:627:LEU:HD12	1.91	0.51
1:A:310:PHE:HB3	1:A:349:ILE:HD13	1.92	0.51
1:A:445:LYS:NZ	1:A:449:ASP:OD2	2.40	0.50
1:A:76:THR:OG1	1:A:77:LEU:N	2.44	0.50
1:A:214:VAL:HB	1:A:230:LEU:HB2	1.94	0.50
1:A:228:HIS:NE2	1:A:230:LEU:O	2.44	0.50
1:A:360:ALA:HB3	1:A:363:LYS:HB2	1.92	0.49
1:A:167:CYS:HB2	1:A:255:LEU:HA	1.94	0.49
1:A:273:LEU:HB3	1:A:282:LEU:HB3	1.94	0.49
1:A:141:THR:OG1	1:A:147:GLN:N	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ARG:HG2	1:A:247:GLU:HG2	1.93	0.48
1:A:7:VAL:O	1:A:27:THR:HA	2.14	0.48
1:A:21:LEU:HD12	1:A:117:GLU:HB2	1.94	0.48
1:A:275:TYR:HA	1:A:282:LEU:HD12	1.95	0.48
1:A:645:VAL:O	1:A:649:THR:OG1	2.30	0.48
1:A:545:GLN:HA	1:A:548:LEU:HD12	1.95	0.48
1:A:49:TRP:CZ2	1:A:82:SER:HB3	2.49	0.47
1:A:572:LEU:HB3	1:A:584:PHE:HB3	1.95	0.47
1:A:552:ARG:NH1	1:A:568:ASP:OD1	2.38	0.47
1:A:353:ILE:HD12	1:A:368:VAL:HG22	1.96	0.47
1:A:659:LYS:HB2	1:A:662:HIS:ND1	2.29	0.47
1:A:520:GLY:N	1:A:521:MET:HA	2.30	0.46
1:A:485:LYS:HE2	1:A:486:LYS:N	2.27	0.46
1:A:51:GLU:HB2	1:A:78:ARG:HB3	1.98	0.46
1:A:117:GLU:HB3	1:A:189:LYS:HB2	1.98	0.45
1:A:203:TYR:HE2	1:A:271:SER:HB2	1.81	0.45
1:A:494:THR:HA	1:A:571:ALA:O	2.17	0.45
1:A:117:GLU:O	1:A:188:THR:HA	2.16	0.45
1:A:486:LYS:O	1:A:488:LEU:N	2.40	0.45
1:A:493:SER:HB2	1:A:568:ASP:HB3	1.99	0.45
1:A:11:TYR:HE1	1:A:34:PRO:HA	1.80	0.44
1:A:45:SER:OG	1:A:46:GLU:N	2.50	0.44
1:A:88:ALA:O	1:A:114:THR:OG1	2.31	0.44
1:A:282:LEU:HD22	1:A:282:LEU:H	1.83	0.43
1:A:332:ASN:O	1:A:332:ASN:ND2	2.47	0.43
1:A:644:GLU:O	1:A:646:HIS:N	2.43	0.43
1:A:26:TRP:HE1	1:A:76:THR:HG21	1.83	0.43
1:A:481:ILE:HG22	1:A:627:LEU:HD21	1.99	0.43
1:A:307:VAL:O	1:A:333:CYS:HB2	2.18	0.43
1:A:381:PRO:O	1:A:384:GLU:HA	2.19	0.43
1:A:210:ASP:O	1:A:255:LEU:HB2	2.18	0.42
1:A:395:ARG:NH2	3:A:812:HOH:O	2.52	0.42
1:A:641:PHE:N	1:A:642:LEU:HA	2.34	0.42
1:A:519:GLY:HA2	1:A:520:GLY:C	2.44	0.42
1:A:552:ARG:NH2	1:A:563:GLU:OE2	2.44	0.42
1:A:663:MET:HE2	1:A:663:MET:HB3	1.96	0.42
1:A:85:SER:HB3	1:A:87:GLU:O	2.19	0.42
1:A:485:LYS:HZ1	1:A:486:LYS:HD2	1.83	0.42
1:A:367:VAL:HG22	1:A:391:GLY:HA3	2.02	0.41
1:A:350:GLN:HB2	1:A:646:HIS:O	2.21	0.41
1:A:40:PHE:CE2	1:A:69:TRP:HB2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:TRP:HE3	1:A:647:CYS:HB3	1.83	0.41
1:A:600:ILE:H	1:A:600:ILE:HG13	1.59	0.41
1:A:266:VAL:HG23	1:A:292:PHE:HB2	2.02	0.41
1:A:535:GLU:O	1:A:538:VAL:HG12	2.21	0.41
1:A:590:ASN:O	1:A:646:HIS:HB2	2.20	0.41
1:A:633:ASP:OD1	1:A:634:ASP:N	2.54	0.41
1:A:298:ILE:HD13	1:A:426:LEU:HD13	2.02	0.40
1:A:474:GLU:O	1:A:587:ASN:ND2	2.54	0.40
1:A:488:LEU:HD23	1:A:488:LEU:HA	1.80	0.40
1:A:159:THR:HA	1:A:160:PRO:HD3	1.86	0.40
1:A:575:MET:HG2	1:A:579:HIS:HA	2.03	0.40
1:A:590:ASN:HB3	1:A:644:GLU:O	2.21	0.40
1:A:155:CYS:HB2	1:A:391:GLY:CA	2.52	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	641/690 (93%)	600 (94%)	39 (6%)	2 (0%)	36 68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	645	VAL
1	A	482	PRO

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	536/610 (88%)	499 (93%)	37 (7%)	14	43

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	24	TYR
1	A	55	ASP
1	A	71	LEU
1	A	92	LYS
1	A	94	THR
1	A	111	LEU
1	A	155	CYS
1	A	162	LEU
1	A	168	ARG
1	A	171	LYS
1	A	188	THR
1	A	206	MET
1	A	233	ARG
1	A	235	LEU
1	A	248	LEU
1	A	315	LYS
1	A	332	ASN
1	A	343	ASN
1	A	346	ASP
1	A	347	ARG
1	A	353	ILE
1	A	382	VAL
1	A	393	VAL
1	A	394	THR
1	A	403	THR
1	A	437	LEU
1	A	485	LYS
1	A	486	LYS
1	A	555	LEU
1	A	559	LEU
1	A	577	GLU
1	A	589	VAL
1	A	594	LEU

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Mol	Chain	Res	Type
1	A	596	LYS
1	A	651	VAL
1	A	664	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	579	HIS
1	A	607	GLN
1	A	662	HIS

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 6 ligands modelled in this entry, 6 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	644/690 (93%)	0.58	40 (6%) 26 14	13, 34, 66, 91	3 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	ASP	8.3
1	A	331	THR	5.0
1	A	595	ASP	4.0
1	A	314	MET	3.8
1	A	596	LYS	3.5
1	A	222	ALA	3.2
1	A	26	TRP	3.0
1	A	648	GLY	3.0
1	A	643	GLY	3.0
1	A	482	PRO	2.9
1	A	340	GLN	2.8
1	A	488	LEU	2.7
1	A	518	LEU	2.7
1	A	220	PRO	2.7
1	A	597	ASP	2.7
1	A	375	GLY	2.6
1	A	376	ASN	2.6
1	A	161	TRP	2.6
1	A	524	LYS	2.6
1	A	377	LEU	2.6
1	A	315	LYS	2.5
1	A	486	LYS	2.5
1	A	522	SER	2.5
1	A	519	GLY	2.5
1	A	345	GLY	2.4
1	A	437	LEU	2.4
1	A	177	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	481	ILE	2.3
1	A	339	PHE	2.3
1	A	176	GLU	2.3
1	A	348	TRP	2.2
1	A	24	TYR	2.2
1	A	523	SER	2.2
1	A	640	LYS	2.1
1	A	384	GLU	2.1
1	A	526	ILE	2.1
1	A	155	CYS	2.1
1	A	343	ASN	2.0
1	A	99	ASP	2.0
1	A	100	GLU	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](#)

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	CA	A	701	1/1	0.76	0.13	46,46,46,46	0
2	CA	A	702	1/1	0.86	0.21	43,43,43,43	0
2	CA	A	706	1/1	0.91	0.07	36,36,36,36	0
2	CA	A	704	1/1	0.93	0.06	22,22,22,22	0
2	CA	A	703	1/1	0.94	0.14	44,44,44,44	0
2	CA	A	705	1/1	0.96	0.05	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.