



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:48 AM UTC

PDB ID : 8HPP / pdb_00008hpp
Title : Crystal structure of human INTS3 with SAGE1
Authors : Deng, W.; Wu, J.; Lei, M.
Deposited on : 2022-12-12
Resolution : 3.00 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

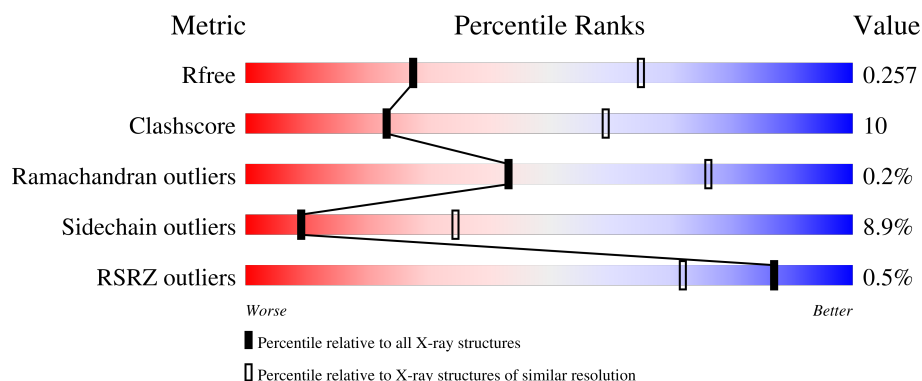
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	471	
1	B	471	
2	C	90	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7102 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 Integrator complex subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			3195	2026	528	615	26			
1	B	394	Total	C	N	O	S	0	0	0
			3189	2023	527	613	26			

- Molecule 2 is a protein called Sarcoma antigen 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	86	Total	C	N	O	S	0	0	0
			718	461	131	122	4			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	815	PRO	-	expression tag	UNP Q9NXZ1
C	816	GLY	-	expression tag	UNP Q9NXZ1
C	817	SER	-	expression tag	UNP Q9NXZ1

- Molecule 1: Integrator complex subunit 3



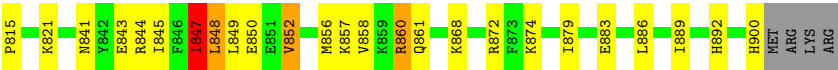
● Molecule 2: Sarcoma antigen 1

Chain C:

68%

23%

...



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	260.12Å 46.24Å 103.10Å 90.00° 113.28° 90.00°	Depositor
Resolution (Å)	47.40 – 3.00 47.40 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.4 (47.40-3.00) 97.4 (47.40-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.215 , 0.257 0.218 , 0.257	Depositor DCC
R_{free} test set	1102 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	65.5	Xtriage
Anisotropy	0.782	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7102	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% 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	0.43	0/3251	1.10	8/4392 (0.2%)
1	B	0.43	0/3245	1.07	4/4384 (0.1%)
2	C	0.41	0/728	1.21	3/965 (0.3%)
All	All	0.43	0/7224	1.10	15/9741 (0.2%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	900	SER	N-CA-C	19.80	140.56	112.04
1	B	628	ALA	CB-CA-C	13.38	133.43	110.68
1	B	629	HIS	N-CA-CB	-13.05	90.81	110.26
2	C	847	ILE	CB-CA-C	12.64	131.75	112.16
1	A	901	LEU	N-CA-CB	-9.94	93.60	110.50
1	A	701	ALA	N-CA-C	-9.21	93.29	108.67
2	C	848	LEU	N-CA-CB	-9.03	96.19	110.19
1	A	900	SER	CB-CA-C	-8.71	94.62	110.56
1	B	629	HIS	N-CA-C	8.11	123.72	112.45
1	A	956	ALA	CB-CA-C	5.61	121.62	110.17
2	C	850	GLU	N-CA-C	-5.47	106.11	112.89
1	A	957	SER	N-CA-CB	-5.22	99.95	111.69
1	A	701	ALA	CB-CA-C	5.12	124.20	117.23
1	B	873	THR	CB-CA-C	5.12	120.83	110.38
1	A	866	HIS	CB-CA-C	5.01	116.85	110.13

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	3195	0	3178	70	0
1	B	3189	0	3173	67	0
2	C	718	0	763	13	0
All	All	7102	0	7114	146	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 10.

All (146) 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:832:LYS:H	1:A:836:HIS:HD2	0.99	0.97
1:B:832:LYS:H	1:B:836:HIS:HD2	1.19	0.89
1:B:727:MET:HB3	1:B:762:MET:HE1	1.53	0.88
1:A:832:LYS:H	1:A:836:HIS:CD2	1.90	0.86
1:B:777:CYS:O	1:B:781:MET:HG2	1.76	0.85
1:A:825:ILE:N	1:A:826:PRO:HD2	1.94	0.82
1:B:825:ILE:N	1:B:826:PRO:HD2	1.95	0.82
1:B:821:LEU:HD23	1:B:847:LEU:HD23	1.60	0.81
1:A:821:LEU:HD21	1:A:847:LEU:HD23	1.60	0.81
1:B:695:ARG:HH11	1:B:744:SER:HB2	1.50	0.77
1:A:832:LYS:N	1:A:836:HIS:HD2	1.80	0.77
1:A:898:ASN:HB3	1:A:920:LEU:HD12	1.69	0.74
1:B:821:LEU:HD21	1:B:846:GLN:HB3	1.71	0.72
1:B:645:GLU:HA	1:B:648:VAL:HG22	1.74	0.70
1:A:831:LEU:O	1:A:863:ARG:NH2	2.25	0.69
1:B:832:LYS:H	1:B:836:HIS:CD2	2.09	0.68
1:A:821:LEU:CD2	1:A:847:LEU:HD23	2.23	0.68
1:A:866:HIS:HD2	1:A:867:PRO:HD2	1.59	0.68
2:C:843:GLU:O	2:C:847:ILE:HG23	1.93	0.68
1:A:803:GLU:HG2	2:C:874:LYS:O	1.95	0.66
1:A:615:LEU:HA	1:A:618:LEU:HB3	1.77	0.65
1:A:820:PRO:HB2	1:A:822:GLU:HG2	1.76	0.65
1:B:634:VAL:HG21	1:B:737:LEU:HD22	1.78	0.65
1:B:727:MET:HA	1:B:727:MET:HE2	1.79	0.64
1:B:766:VAL:O	2:C:872:ARG:NH2	2.31	0.64
2:C:849:LEU:O	2:C:852:VAL:HG12	1.97	0.64
1:A:937:THR:HG22	1:A:938:LYS:H	1.62	0.64
1:A:743:PRO:HA	1:A:784:LEU:HD13	1.78	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:824:ILE:C	1:B:826:PRO:HD2	2.23	0.62
1:B:847:LEU:HD22	1:B:856:MET:HE3	1.80	0.62
1:A:825:ILE:N	1:A:826:PRO:CD	2.63	0.62
1:B:825:ILE:N	1:B:826:PRO:CD	2.62	0.62
1:B:828:LEU:HG	1:B:863:ARG:HD2	1.82	0.61
1:A:626:PHE:C	1:A:628:ALA:H	2.09	0.60
1:B:635:LEU:HD23	1:B:636:PRO:HD2	1.85	0.59
1:B:950:ALA:O	1:B:954:VAL:HG23	2.03	0.59
1:B:899:ASN:HA	1:B:920:LEU:HD11	1.87	0.57
1:A:952:GLN:HE22	1:A:976:TYR:HD1	1.51	0.56
1:A:941:PHE:O	1:A:944:GLN:HG2	2.06	0.56
1:B:694:LEU:HD21	1:B:706:LEU:HD23	1.86	0.56
1:B:815:LEU:HD11	1:B:846:GLN:HG3	1.86	0.56
1:A:606:LEU:O	1:A:607:GLU:C	2.49	0.56
1:B:899:ASN:HD22	1:B:953:HIS:CD2	2.24	0.55
1:A:853:SER:O	1:A:857:VAL:HG23	2.05	0.55
1:B:659:LEU:HA	1:B:662:MET:HE2	1.87	0.55
2:C:856:MET:HE3	2:C:889:ILE:HG23	1.88	0.55
1:B:862:SER:C	1:B:922:GLN:HE22	2.15	0.55
1:B:838:GLU:H	1:B:838:GLU:CD	2.14	0.54
1:A:892:LYS:HG3	1:A:950:ALA:HB2	1.90	0.54
1:B:641:GLU:HA	1:B:644:LEU:HD12	1.90	0.53
1:B:861:LEU:O	1:B:922:GLN:NE2	2.41	0.53
1:A:877:ARG:HD2	1:A:933:ASN:OD1	2.08	0.53
1:A:826:PRO:HA	1:A:859:MET:HE2	1.88	0.53
1:A:620:SER:O	1:A:624:GLU:HG2	2.09	0.53
1:A:939:GLN:HE21	1:A:939:GLN:N	2.07	0.53
1:A:748:GLU:C	1:A:750:PRO:HD3	2.35	0.52
1:B:803:GLU:HB2	1:B:806:GLU:CD	2.33	0.52
1:A:680:TYR:CD1	1:A:710:PHE:HZ	2.28	0.52
1:A:870:GLN:NE2	1:A:870:GLN:HA	2.24	0.51
2:C:860:ARG:HB2	2:C:889:ILE:HG21	1.91	0.51
1:B:826:PRO:HA	1:B:859:MET:HE2	1.93	0.51
1:B:835:GLU:C	1:B:837:PRO:HD3	2.36	0.51
1:B:913:SER:HA	1:B:916:ALA:HB3	1.93	0.51
1:A:666:ASN:HB3	1:A:669:PHE:HB2	1.92	0.51
1:A:799:SER:HA	1:A:802:TRP:CD2	2.46	0.50
1:B:795:ILE:O	1:B:798:GLN:HB2	2.12	0.50
2:C:841:ASN:OD1	2:C:843:GLU:HG2	2.10	0.50
1:B:731:GLN:CB	1:B:738:LEU:HD22	2.42	0.50
2:C:857:LYS:O	2:C:861:GLN:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:788:ARG:HD3	1:B:790:ASP:OD1	2.10	0.50
1:B:727:MET:CB	1:B:762:MET:HE1	2.34	0.50
1:A:711:ALA:C	1:A:713:ALA:H	2.19	0.50
1:B:824:ILE:O	1:B:827:ILE:HG22	2.11	0.50
1:B:933:ASN:O	1:B:936:ASN:ND2	2.42	0.49
1:B:731:GLN:HB2	1:B:738:LEU:HD22	1.95	0.48
1:A:578:LEU:HD21	1:A:604:GLN:HB3	1.95	0.48
2:C:845:ILE:O	2:C:848:LEU:HB3	2.13	0.48
2:C:815:PRO:N	2:C:821:LYS:HZ2	2.12	0.48
1:A:892:LYS:O	1:A:893:SER:C	2.56	0.48
1:A:626:PHE:C	1:A:628:ALA:N	2.70	0.48
1:A:865:CYS:SG	1:A:925:GLU:HB3	2.53	0.47
1:B:764:VAL:O	1:B:806:GLU:HG2	2.15	0.47
1:B:891:ILE:O	1:B:895:LEU:HB2	2.15	0.47
1:A:641:GLU:HA	1:A:644:LEU:HD12	1.96	0.47
1:B:937:THR:HB	1:B:939:GLN:HG3	1.96	0.46
1:A:585:LEU:HD22	1:A:625:LEU:HD21	1.96	0.46
1:A:663:GLN:HG3	1:A:666:ASN:HB2	1.98	0.46
1:B:626:PHE:O	1:B:629:HIS:HB3	2.16	0.46
1:A:824:ILE:C	1:A:826:PRO:HD2	2.41	0.46
1:A:694:LEU:HD21	1:A:706:LEU:HD23	1.98	0.46
1:B:954:VAL:HG12	1:B:954:VAL:O	2.15	0.46
2:C:879:ILE:O	2:C:883:GLU:HG2	2.16	0.46
1:A:741:LEU:O	1:A:742:THR:C	2.59	0.45
1:A:589:SER:HA	1:A:594:GLN:NE2	2.32	0.45
1:A:691:LEU:HD12	1:A:691:LEU:HA	1.79	0.45
1:B:802:TRP:HB3	1:B:806:GLU:OE1	2.17	0.45
1:B:891:ILE:HB	1:B:947:ILE:HD11	1.99	0.44
1:A:598:MET:HG3	1:A:651:PRO:HB3	2.00	0.44
1:A:818:ASN:ND2	1:B:848:ARG:HH12	2.15	0.44
1:A:825:ILE:HG23	1:A:859:MET:HG3	1.99	0.44
1:A:768:ASP:OD1	1:A:771:GLN:HG3	2.18	0.43
1:A:574:LEU:HB3	1:A:579:ARG:HD2	1.99	0.43
1:A:752:GLU:HA	1:A:755:ARG:NH2	2.33	0.43
1:A:850:GLU:OE1	1:A:850:GLU:HA	2.19	0.43
1:A:948:LEU:HD22	1:A:973:ALA:HA	2.00	0.43
1:A:839:ALA:O	1:A:843:LEU:HB2	2.18	0.43
1:A:939:GLN:HE21	1:A:939:GLN:H	1.67	0.43
1:B:720:HIS:O	1:B:724:MET:HG2	2.18	0.43
1:B:767:ILE:HA	1:B:771:GLN:OE1	2.18	0.43
1:B:594:GLN:O	1:B:598:MET:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:703:LYS:HB3	1:B:705:ASN:OD1	2.18	0.43
1:B:808:TYR:O	1:B:812:GLN:HG2	2.19	0.43
1:A:833:TYR:CD2	1:A:866:HIS:CE1	3.06	0.42
1:A:870:GLN:O	1:A:874:SER:HB2	2.18	0.42
1:B:860:VAL:HG13	1:B:872:THR:HG23	2.00	0.42
1:A:775:LEU:HD12	1:A:775:LEU:HA	1.82	0.42
2:C:856:MET:HE2	2:C:892:HIS:CD2	2.54	0.42
1:B:760:LEU:HD23	1:B:802:TRP:HH2	1.84	0.42
1:B:787:PHE:HD2	1:B:817:HIS:ND1	2.18	0.42
1:B:821:LEU:CD2	1:B:847:LEU:HD23	2.43	0.42
1:A:611:ASP:HB3	1:A:612:SER:H	1.73	0.42
1:B:761:ASN:HB2	1:B:802:TRP:CZ2	2.55	0.42
1:A:755:ARG:H	1:A:755:ARG:HG3	1.76	0.41
1:A:773:GLN:NE2	1:B:840:LEU:HB3	2.34	0.41
1:A:585:LEU:HA	1:A:594:GLN:HG2	2.02	0.41
2:C:886:LEU:HD12	2:C:886:LEU:HA	1.87	0.41
1:A:728:LYS:HE3	1:A:731:GLN:OE1	2.19	0.41
1:B:877:ARG:NH1	1:B:936:ASN:HD21	2.18	0.41
1:A:731:GLN:HG3	1:A:732:GLU:N	2.33	0.41
1:A:931:ARG:HH11	1:A:931:ARG:HG2	1.85	0.41
1:B:860:VAL:CG1	1:B:872:THR:HG23	2.50	0.41
1:A:799:SER:C	1:A:801:ASP:N	2.78	0.41
1:B:672:LEU:HD12	1:B:672:LEU:HA	1.85	0.41
1:B:764:VAL:HG23	1:B:806:GLU:HB3	2.02	0.41
1:B:803:GLU:O	1:B:804:THR:C	2.63	0.41
1:A:779:VAL:HG23	1:A:784:LEU:HD23	2.03	0.41
1:B:707:TYR:OH	1:B:723:LEU:HA	2.20	0.41
1:B:578:LEU:HD23	1:B:581:LYS:HG3	2.03	0.40
1:B:871:PHE:CD2	1:B:871:PHE:C	2.99	0.40
1:A:728:LYS:HB2	1:A:762:MET:HE2	2.03	0.40
1:A:955:GLN:O	1:A:963:LYS:HE2	2.21	0.40
1:A:876:LEU:O	1:A:877:ARG:C	2.65	0.40
1:A:575:ASP:HB3	1:A:578:LEU:HD12	2.04	0.40
1:A:796:LEU:HD12	1:A:796:LEU:HA	1.93	0.40
1:B:763:ILE:C	1:B:765:ALA:H	2.30	0.40
1:B:959:ASP:H	1:B:962:HIS:CE1	2.40	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	391/471 (83%)	355 (91%)	35 (9%)	1 (0%)	36	70
1	B	390/471 (83%)	364 (93%)	25 (6%)	1 (0%)	36	70
2	C	84/90 (93%)	84 (100%)	0	0	100	100
All	All	865/1032 (84%)	803 (93%)	60 (7%)	2 (0%)	43	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	591	THR
1	B	588	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	366/433 (84%)	330 (90%)	36 (10%)	7	30
1	B	365/433 (84%)	336 (92%)	29 (8%)	11	39
2	C	78/82 (95%)	71 (91%)	7 (9%)	9	34
All	All	809/948 (85%)	737 (91%)	72 (9%)	9	34

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	574	LEU

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Mol	Chain	Res	Type
1	A	604	GLN
1	A	609	ASP
1	A	610	PHE
1	A	613	GLU
1	A	615	LEU
1	A	627	LYS
1	A	678	GLU
1	A	691	LEU
1	A	716	LEU
1	A	731	GLN
1	A	735	VAL
1	A	755	ARG
1	A	763	ILE
1	A	769	SER
1	A	774	GLU
1	A	775	LEU
1	A	787	PHE
1	A	789	LYS
1	A	796	LEU
1	A	815	LEU
1	A	818	ASN
1	A	825	ILE
1	A	828	LEU
1	A	834	LYS
1	A	843	LEU
1	A	873	THR
1	A	874	SER
1	A	880	CYS
1	A	881	MET
1	A	901	LEU
1	A	919	THR
1	A	937	THR
1	A	939	GLN
1	A	949	GLN
1	A	954	VAL
1	B	608	GLU
1	B	627	LYS
1	B	634	VAL
1	B	637	GLU
1	B	674	ASP
1	B	676	LEU
1	B	681	GLN

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Mol	Chain	Res	Type
1	B	691	LEU
1	B	695	ARG
1	B	712	GLN
1	B	716	LEU
1	B	731	GLN
1	B	735	VAL
1	B	775	LEU
1	B	789	LYS
1	B	797	ILE
1	B	804	THR
1	B	815	LEU
1	B	817	HIS
1	B	818	ASN
1	B	833	TYR
1	B	843	LEU
1	B	847	LEU
1	B	873	THR
1	B	937	THR
1	B	939	GLN
1	B	947	ILE
1	B	959	ASP
1	B	962	HIS
2	C	844	ARG
2	C	847	ILE
2	C	852	VAL
2	C	858	VAL
2	C	860	ARG
2	C	868	LYS
2	C	900	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	594	GLN
1	A	599	GLN
1	A	623	GLN
1	A	663	GLN
1	A	681	GLN
1	A	683	GLN
1	A	689	HIS
1	A	705	ASN
1	A	818	ASN

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Mol	Chain	Res	Type
1	A	836	HIS
1	A	866	HIS
1	A	870	GLN
1	A	922	GLN
1	A	939	GLN
1	A	940	ASN
1	A	944	GLN
1	A	949	GLN
1	A	952	GLN
1	A	953	HIS
1	B	573	GLN
1	B	604	GLN
1	B	623	GLN
1	B	683	GLN
1	B	689	HIS
1	B	731	GLN
1	B	761	ASN
1	B	836	HIS
1	B	870	GLN
1	B	890	HIS
1	B	899	ASN
1	B	922	GLN
1	B	940	ASN
1	B	949	GLN
2	C	861	GLN
2	C	892	HIS

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

There are no ligands in this entry.

5.7 Other polymers

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 [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	395/471 (83%)	-0.38	1 (0%) 90 79	44, 102, 200, 240	0
1	B	394/471 (83%)	-0.32	3 (0%) 82 64	41, 108, 203, 239	0
2	C	86/90 (95%)	-0.80	0 100 100	43, 67, 105, 152	0
All	All	875/1032 (84%)	-0.39	4 (0%) 87 72	41, 98, 201, 240	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	915	LEU	3.2
1	A	915	LEU	2.3
1	B	918	LEU	2.2
1	B	580	ASP	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.