



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2026 – 11:23 PM JST

PDB ID : 9WP8 / pdb_00009wp8
Title : Crystal structure of the SARS-CoV-2 spike protein HR1 in complex with a designed peptide inhibitor dP-S2FI-0089 at 3.1 angstrom resolution
Authors : Wang, K.; Lyu, Z.; Ye, S.
Deposited on : 2025-09-08
Resolution : 3.10 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

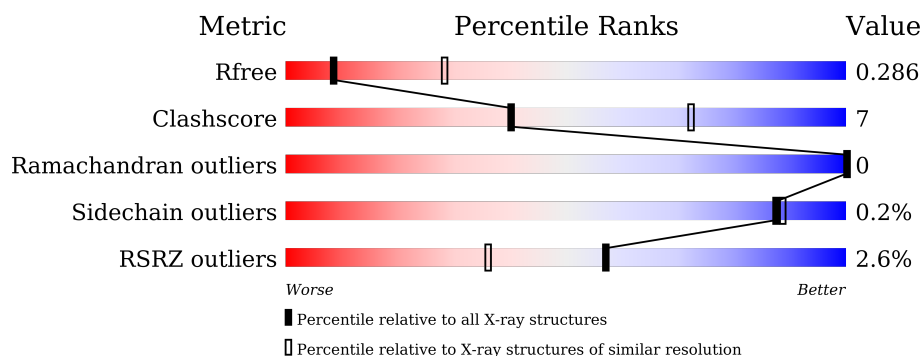
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	120	0% 68% 21% 11%
1	C	120	73% 18% 9%
1	E	120	2% 81% 12% 8%
1	G	120	3% 80% 12% 8%
1	I	120	3% 73% 18% 9%
1	K	120	5% 78% 17% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5001 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 dP-S2FI-0089,Spike protein S2'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	107	Total	C	N	O	S	0	0	0
			806	502	140	163	1			
1	C	109	Total	C	N	O	S	0	1	0
			821	512	143	165	1			
1	E	111	Total	C	N	O	S	0	0	0
			831	517	145	168	1			
1	G	110	Total	C	N	O	S	0	0	0
			827	515	144	167	1			
1	I	109	Total	C	N	O	S	0	0	0
			818	510	142	165	1			
1	K	113	Total	C	N	O	S	0	0	0
			842	523	147	171	1			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	903C	GLY	-	linker	PDB ?
A	903D	SER	-	linker	PDB ?
A	903E	SER	-	linker	PDB ?
A	903F	GLY	-	linker	PDB ?
A	903G	GLY	-	linker	PDB ?
A	903H	SER	-	linker	PDB ?
A	903I	GLY	-	linker	PDB ?
C	904B	GLY	-	linker	PDB ?
C	904C	SER	-	linker	PDB ?
C	904D	SER	-	linker	PDB ?
C	904E	GLY	-	linker	PDB ?
C	904F	GLY	-	linker	PDB ?
C	904G	SER	-	linker	PDB ?
C	904H	GLY	-	linker	PDB ?
E	904B	GLY	-	linker	PDB ?
E	904C	SER	-	linker	PDB ?
E	904D	SER	-	linker	PDB ?

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Chain	Residue	Modelled	Actual	Comment	Reference
E	904E	GLY	-	linker	PDB ?
E	904F	GLY	-	linker	PDB ?
E	904G	SER	-	linker	PDB ?
E	904H	GLY	-	linker	PDB ?
G	903C	GLY	-	linker	PDB ?
G	903D	SER	-	linker	PDB ?
G	903E	SER	-	linker	PDB ?
G	903F	GLY	-	linker	PDB ?
G	903G	GLY	-	linker	PDB ?
G	903H	SER	-	linker	PDB ?
G	903I	GLY	-	linker	PDB ?
I	903C	GLY	-	linker	PDB ?
I	903D	SER	-	linker	PDB ?
I	903E	SER	-	linker	PDB ?
I	903F	GLY	-	linker	PDB ?
I	903G	GLY	-	linker	PDB ?
I	903H	SER	-	linker	PDB ?
I	903I	GLY	-	linker	PDB ?
K	904B	GLY	-	linker	PDB ?
K	904C	SER	-	linker	PDB ?
K	904D	SER	-	linker	PDB ?
K	904E	GLY	-	linker	PDB ?
K	904F	GLY	-	linker	PDB ?
K	904G	SER	-	linker	PDB ?
K	911	GLY	-	linker	PDB ?

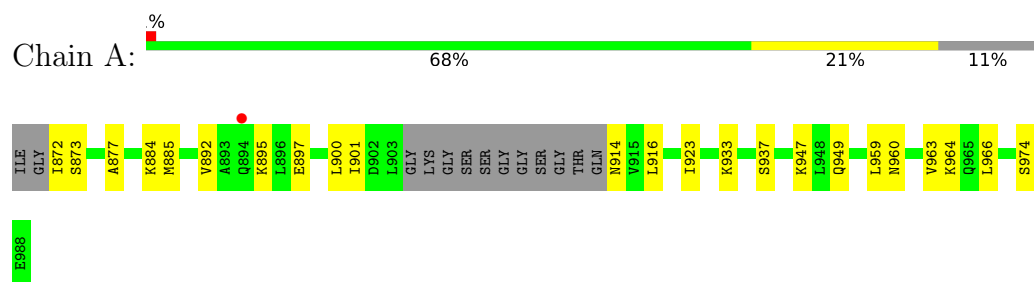
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	11	Total O 11 11	0	0
2	C	8	Total O 8 8	0	0
2	E	9	Total O 9 9	0	0
2	G	6	Total O 6 6	0	0
2	I	5	Total O 5 5	0	0
2	K	17	Total O 17 17	0	0

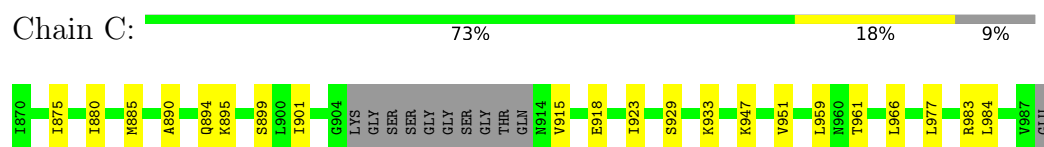
3 Residue-property plots [i](#)

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.

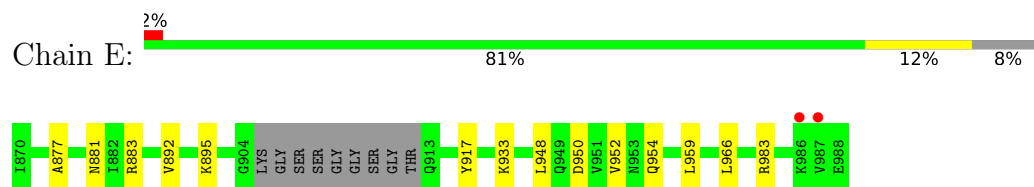
- Molecule 1: dP-S2FI-0089,Spike protein S2'



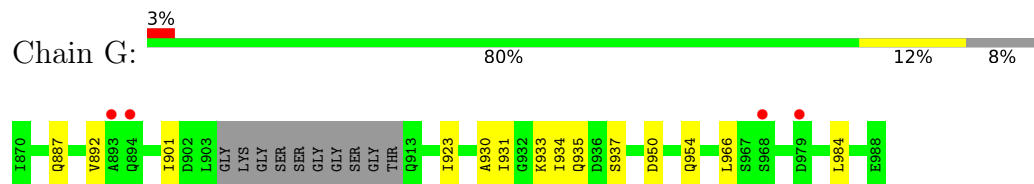
- Molecule 1: dP-S2FI-0089,Spike protein S2'



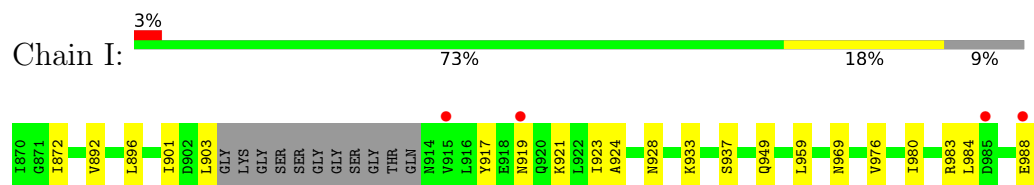
- Molecule 1: dP-S2FI-0089,Spike protein S2'



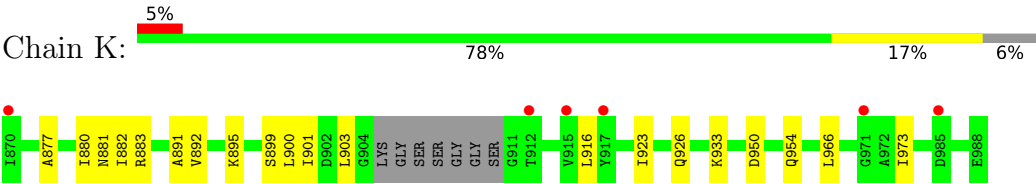
- Molecule 1: dP-S2FI-0089,Spike protein S2'



- Molecule 1: dP-S2FI-0089,Spike protein S2'



● Molecule 1: dP-S2FI-0089,Spike protein S2'



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.64Å 63.79Å 63.86Å 64.46° 64.54° 80.66°	Depositor
Resolution (Å)	30.68 – 3.10 30.68 – 3.10	Depositor EDS
% Data completeness (in resolution range)	95.2 (30.68-3.10) 95.1 (30.68-3.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.49 (at 3.11Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.237 , 0.290 0.239 , 0.286	Depositor DCC
R_{free} test set	754 reflections (4.41%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5001	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.35% 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	0.09	0/807	0.22	0/1086
1	C	0.08	0/822	0.19	0/1107
1	E	0.08	0/832	0.19	0/1119
1	G	0.09	0/828	0.24	0/1114
1	I	0.09	0/819	0.25	0/1102
1	K	0.10	0/843	0.25	0/1134
All	All	0.09	0/4951	0.22	0/6662

There are no bond length outliers.

There are no bond angle outliers.

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	806	0	842	17	0
1	C	821	0	858	15	0
1	E	831	0	867	13	0
1	G	827	0	864	9	0
1	I	818	0	856	15	0
1	K	842	0	877	15	0
2	A	11	0	0	1	0
2	C	8	0	0	0	0
2	E	9	0	0	0	0
2	G	6	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	5	0	0	0	0
2	K	17	0	0	0	0
All	All	5001	0	5164	67	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 (67) 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:K:901:ILE:HG21	1:K:923:ILE:HG22	1.66	0.78
1:C:895:LYS:HE3	1:C:933:LYS:HD3	1.75	0.68
1:A:916:LEU:HD11	1:E:917:TYR:HB2	1.77	0.67
1:A:885:MET:HE1	1:A:947:LYS:HD2	1.75	0.67
1:A:959:LEU:HD23	1:E:959:LEU:HD11	1.77	0.67
1:G:966:LEU:HD21	1:K:966:LEU:HD13	1.80	0.62
1:A:895:LYS:HB3	1:A:933:LYS:NZ	2.13	0.62
1:A:897:GLU:HA	1:A:900:LEU:HD23	1.81	0.62
1:A:884:LYS:NZ	2:A:1003:HOH:O	2.34	0.60
1:G:892:VAL:HG11	1:G:937:SER:HB3	1.83	0.60
1:E:892:VAL:HG23	1:E:933:LYS:HB3	1.85	0.58
1:A:974:SER:HA	1:A:977:LEU:HD12	1.86	0.58
1:G:984:LEU:HD23	1:I:983:ARG:HH21	1.69	0.58
1:G:887:GLN:NE2	2:G:1001:HOH:O	2.37	0.57
1:I:921:LYS:HE2	1:K:903:LEU:HD12	1.87	0.56
1:C:880:ILE:HD11	1:C:947:LYS:HB3	1.88	0.56
1:E:895:LYS:HB3	1:E:933:LYS:HD3	1.88	0.56
1:A:959:LEU:O	1:A:963:VAL:HG23	2.07	0.55
1:I:984:LEU:O	1:I:988:GLU:N	2.40	0.54
1:E:895:LYS:HG2	1:E:933:LYS:HE2	1.89	0.54
1:A:901:ILE:HD13	1:A:923:ILE:HG13	1.90	0.54
1:A:984:LEU:HB3	1:C:983:ARG:HH21	1.73	0.53
1:C:901:ILE:HD13	1:C:923:ILE:HG12	1.91	0.53
1:I:959:LEU:HD23	1:K:877:ALA:HB2	1.91	0.52
1:I:903:LEU:HG	1:I:919:ASN:ND2	2.25	0.52
1:C:875:ILE:HD11	1:C:961:THR:HB	1.92	0.51
1:G:950:ASP:O	1:G:954:GLN:HG3	2.12	0.50
1:C:984:LEU:HD13	1:E:983:ARG:HD2	1.94	0.50
1:A:877:ALA:HB2	1:E:959:LEU:HD23	1.93	0.50
1:C:915:VAL:HA	1:C:918[A]:GLU:HG2	1.94	0.49
1:A:966:LEU:HD21	1:E:966:LEU:HD13	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:924:ALA:O	1:I:928:ASN:ND2	2.46	0.48
1:K:950:ASP:O	1:K:954:GLN:HG3	2.13	0.48
1:A:872:ILE:HG22	1:A:873:SER:H	1.79	0.48
1:K:973:ILE:H	1:K:973:ILE:HD12	1.79	0.48
1:C:885:MET:H	1:C:885:MET:HE3	1.79	0.48
1:C:959:LEU:HD23	1:E:877:ALA:HB2	1.96	0.47
1:A:949:GLN:HG3	1:C:880:ILE:HG23	1.97	0.47
1:A:960:ASN:O	1:A:964:LYS:HD2	2.14	0.47
1:I:892:VAL:HG11	1:I:937:SER:HB3	1.98	0.46
1:I:949:GLN:HG3	1:K:880:ILE:HG23	1.97	0.46
1:K:892:VAL:HG13	1:K:933:LYS:HB3	1.97	0.45
1:I:896:LEU:HD23	1:I:933:LYS:HD3	1.98	0.45
1:C:890:ALA:O	1:C:894:GLN:HG3	2.16	0.45
1:C:966:LEU:HD13	1:E:966:LEU:HD21	1.97	0.45
1:C:899:SER:OG	1:C:929:SER:OG	2.34	0.44
1:E:948:LEU:O	1:E:952:VAL:HG23	2.17	0.44
1:I:976:VAL:O	1:I:980:ILE:HD12	2.16	0.44
1:I:917:TYR:HD1	1:K:916:LEU:HD21	1.82	0.44
1:I:901:ILE:HD13	1:I:923:ILE:HG12	2.00	0.44
1:G:901:ILE:HG12	1:G:923:ILE:HD13	2.00	0.43
1:G:933:LYS:HA	1:G:933:LYS:HD3	1.79	0.43
1:E:950:ASP:O	1:E:954:GLN:HG3	2.19	0.43
1:G:930:ALA:O	1:G:934:ILE:HD12	2.19	0.42
1:I:949:GLN:HB2	1:K:882:ILE:HD11	2.01	0.42
1:K:891:ALA:O	1:K:895:LYS:HG3	2.18	0.42
1:K:899:SER:HA	1:K:926:GLN:OE1	2.20	0.42
1:A:892:VAL:HG11	1:A:937:SER:HB3	2.02	0.41
1:A:914:ASN:N	1:A:914:ASN:OD1	2.52	0.41
1:E:881:ASN:OD1	1:E:883:ARG:HG3	2.20	0.41
1:I:928:ASN:OD1	1:K:900:LEU:HB3	2.20	0.41
1:K:903:LEU:HD23	1:K:903:LEU:HA	1.89	0.41
1:C:880:ILE:HD13	1:C:951:VAL:HG21	2.03	0.41
1:C:984:LEU:HD23	1:C:984:LEU:HA	1.86	0.40
1:G:931:ILE:O	1:G:935:GLN:HG3	2.22	0.40
1:I:872:ILE:HG13	1:I:969:ASN:ND2	2.36	0.40
1:K:881:ASN:HD21	1:K:883:ARG:NH2	2.19	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	103/120 (86%)	100 (97%)	3 (3%)	0	100	100
1	C	106/120 (88%)	105 (99%)	1 (1%)	0	100	100
1	E	107/120 (89%)	106 (99%)	1 (1%)	0	100	100
1	G	106/120 (88%)	105 (99%)	1 (1%)	0	100	100
1	I	105/120 (88%)	104 (99%)	1 (1%)	0	100	100
1	K	109/120 (91%)	108 (99%)	1 (1%)	0	100	100
All	All	636/720 (88%)	628 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	92/99 (93%)	92 (100%)	0	100	100
1	C	93/99 (94%)	92 (99%)	1 (1%)	65	78
1	E	94/99 (95%)	94 (100%)	0	100	100
1	G	94/99 (95%)	94 (100%)	0	100	100
1	I	93/99 (94%)	93 (100%)	0	100	100
1	K	95/99 (96%)	95 (100%)	0	100	100
All	All	561/594 (94%)	560 (100%)	1 (0%)	87	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	977	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	955	ASN
1	A	969	ASN
1	C	957	GLN
1	E	876	ASN
1	E	978	ASN
1	G	955	ASN
1	K	881	ASN
1	K	914	ASN

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 [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	107/120 (89%)	0.25	1 (0%) 81 63	29, 47, 88, 103	0
1	C	109/120 (90%)	0.10	0 100 100	29, 43, 81, 90	1 (0%)
1	E	111/120 (92%)	0.19	2 (1%) 67 47	28, 42, 84, 103	0
1	G	110/120 (91%)	0.35	4 (3%) 46 26	28, 49, 90, 106	0
1	I	109/120 (90%)	0.16	4 (3%) 45 25	28, 43, 91, 126	0
1	K	113/120 (94%)	0.38	6 (5%) 32 17	27, 47, 99, 119	0
All	All	659/720 (91%)	0.24	17 (2%) 57 36	27, 46, 90, 126	1 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	894	GLN	3.4
1	K	985	ASP	3.0
1	G	893	ALA	2.9
1	G	979	ASP	2.7
1	K	870	ILE	2.7
1	I	915	VAL	2.7
1	E	987	VAL	2.6
1	I	988	GLU	2.5
1	G	968	SER	2.5
1	E	986	LYS	2.4
1	A	894	GLN	2.4
1	K	917	TYR	2.3
1	I	985	ASP	2.2
1	K	971	GLY	2.2
1	K	912	THR	2.2
1	I	919	ASN	2.1
1	K	915	VAL	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.