



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 10:18 AM EDT

PDB ID : 7WCY / pdb_00007wcy
Title : Crystal Structure of H-2Kb with Cryptosporidium parvum gp40/15 epitope
Authors : Wang, Y.L.; Gao, M.H.; Zhang, L.X.; Fan, S.H.
Deposited on : 2021-12-20
Resolution : 2.36 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

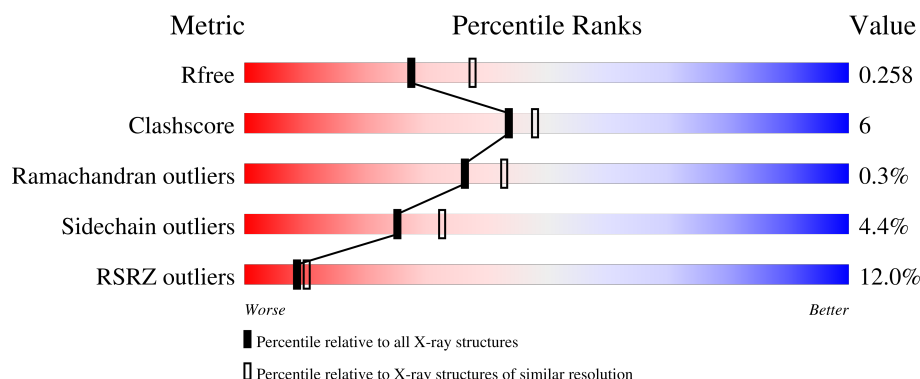
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	275	14% 84% 15%
1	D	275	12% 85% 14%
2	B	99	9% 87% 13%
2	E	99	8% 86% 14%
3	C	9	22% 56% 44%

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Mol	Chain	Length	Quality of chain
3	F	9	22% 33% 33% 33%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6329 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 H-2 class I histocompatibility antigen, K-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2240	1413	394	424	9			
1	D	275	Total	C	N	O	S	0	0	0
			2240	1413	394	424	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			818	523	138	150	7			
2	E	99	Total	C	N	O	S	0	0	0
			818	523	138	150	7			

- Molecule 3 is a protein called SER-VAL-PHE-ALA-ILE-PHE-ALA-ALA-LEU.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			66	47	9	10			
3	F	9	Total	C	N	O	0	2	0
			79	56	11	12			

- Molecule 4 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ni	0	0
			1	1		
4	B	2	Total	Ni	0	0
			2	2		
4	D	3	Total	Ni	0	0
			3	3		
4	E	2	Total	Ni	0	0
			2	2		

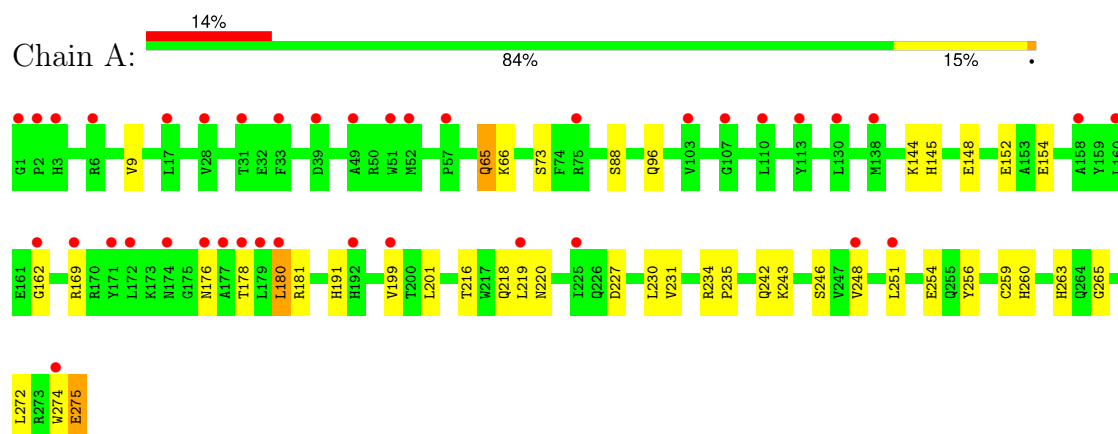
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total 21	O 21	0	0
5	B	8	Total 8	O 8	0	0
5	C	2	Total 2	O 2	0	0
5	D	20	Total 20	O 20	0	0
5	E	9	Total 9	O 9	0	0

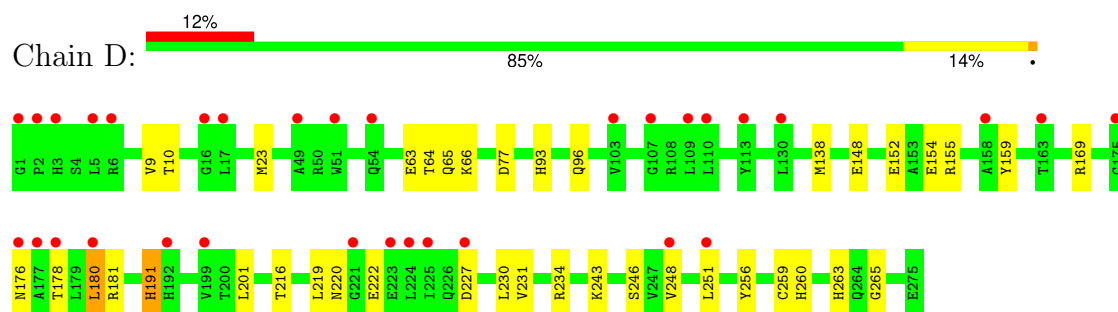
3 Residue-property plots [i](#)

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.

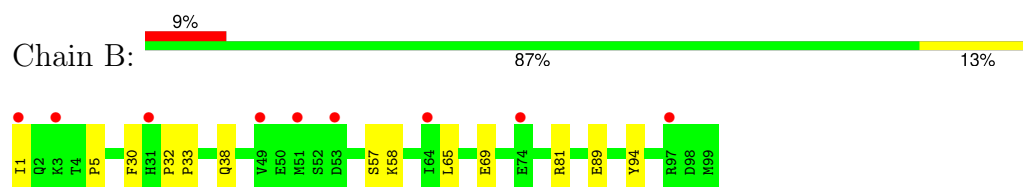
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



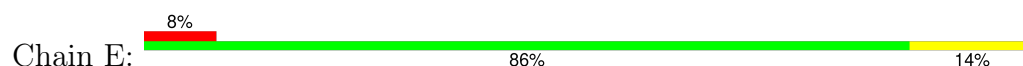
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain

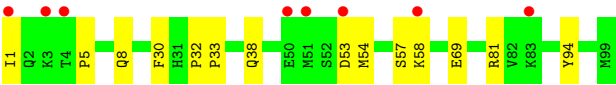


- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin

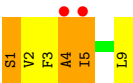




● Molecule 3: SER-VAL-PHE-ALA-ILE-PHE-ALA-ALA-LEU



● Molecule 3: SER-VAL-PHE-ALA-ILE-PHE-ALA-ALA-LEU



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.92Å 84.77Å 95.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.47 – 2.36 49.47 – 2.36	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.47-2.36) 99.6 (49.47-2.36)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.226 , 0.258 0.226 , 0.258	Depositor DCC
R_{free} test set	2079 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	66.5	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.346 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6329	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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: NI

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	0.99	0/2301	1.35	3/3125 (0.1%)
1	D	1.01	0/2301	1.36	4/3125 (0.1%)
2	B	0.99	0/844	1.26	0/1144
2	E	1.01	0/844	1.28	0/1144
3	C	0.92	0/67	1.55	0/90
3	F	0.99	0/80	1.20	0/108
All	All	1.00	0/6437	1.34	7/8736 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	275	GLU	CA-C-O	6.37	131.62	120.80
1	D	191	HIS	CA-CB-CG	5.58	119.38	113.80
1	D	64	THR	CA-CB-OG1	-5.42	101.47	109.60
1	D	93	HIS	CA-C-N	5.06	130.13	122.99
1	D	93	HIS	C-N-CA	5.06	130.13	122.99
1	A	162	GLY	CA-C-N	5.01	127.25	120.38
1	A	162	GLY	C-N-CA	5.01	127.25	120.38

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	2240	0	2129	31	1
1	D	2240	0	2129	28	1
2	B	818	0	797	5	1
2	E	818	0	797	8	1
3	C	66	0	71	8	0
3	F	79	0	86	10	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	D	3	0	0	0	0
4	E	2	0	0	0	0
5	A	21	0	0	9	0
5	B	8	0	0	0	0
5	C	2	0	0	4	0
5	D	20	0	0	3	0
5	E	9	0	0	0	0
All	All	6329	0	6009	75	2

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 (75) 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:145:HIS:NE2	5:A:401:HOH:O	1.85	1.07
3:F:4[A]:ALA:O	3:F:5[A]:ILE:HG13	1.63	0.98
1:A:191:HIS:NE2	1:A:275:GLU:O	1.97	0.98
1:A:145:HIS:CE1	5:A:401:HOH:O	2.14	0.96
1:A:88:SER:O	5:A:402:HOH:O	1.86	0.93
1:D:63:GLU:OE2	3:F:1:SER:OG	1.93	0.87
3:C:5:ILE:CG2	5:C:701:HOH:O	2.26	0.83
1:A:144:LYS:NZ	5:A:404:HOH:O	2.16	0.78
3:C:5:ILE:HG21	5:C:701:HOH:O	1.85	0.76
1:D:152:GLU:HG3	5:D:712:HOH:O	1.88	0.72
1:D:66:LYS:NZ	3:F:2:VAL:O	2.27	0.67
1:A:191:HIS:HE2	1:A:275:GLU:C	2.01	0.63
3:C:5:ILE:HG22	5:C:700:HOH:O	1.98	0.63
3:C:5:ILE:HG22	5:C:701:HOH:O	1.93	0.62
3:F:4[A]:ALA:O	3:F:5[A]:ILE:CG1	2.42	0.61
1:A:66:LYS:NZ	3:C:2:VAL:O	2.35	0.60
1:D:230:LEU:HD23	1:D:231:VAL:O	2.04	0.58
1:D:230:LEU:HD21	1:D:243:LYS:HE3	1.86	0.58
1:D:176:ASN:OD1	1:D:180:LEU:HB2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:23:MET:HE1	2:E:54:MET:CE	2.34	0.58
5:A:401:HOH:O	1:D:260:HIS:HE1	1.86	0.57
5:A:401:HOH:O	1:D:260:HIS:CE1	2.58	0.57
1:A:230:LEU:HD23	1:A:231:VAL:O	2.05	0.57
1:A:176:ASN:OD1	1:A:180:LEU:HB2	2.05	0.56
1:A:230:LEU:HD21	1:A:243:LYS:HE3	1.88	0.55
1:A:66:LYS:HG2	3:C:4:ALA:CB	2.38	0.53
1:A:263:HIS:CD2	1:A:265:GLY:H	2.27	0.53
1:D:23:MET:HE1	2:E:54:MET:HE2	1.90	0.52
3:F:4[A]:ALA:C	3:F:5[A]:ILE:CG1	2.83	0.51
1:D:263:HIS:CD2	1:D:265:GLY:H	2.28	0.51
1:A:230:LEU:HD23	1:A:230:LEU:C	2.36	0.51
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.93	0.50
2:E:5:PRO:HB3	2:E:30:PHE:HB3	1.93	0.50
1:A:230:LEU:HD21	1:A:243:LYS:CE	2.41	0.50
1:D:230:LEU:HD21	1:D:243:LYS:CE	2.42	0.49
3:F:4[A]:ALA:C	3:F:5[A]:ILE:HG13	2.35	0.49
1:A:234:ARG:HE	1:A:242:GLN:HE21	1.61	0.48
2:E:38:GLN:OE1	2:E:81:ARG:NH1	2.46	0.48
1:A:216:THR:OG1	1:A:260:HIS:HB2	2.14	0.48
1:A:235:PRO:HG2	2:B:65:LEU:HD22	1.96	0.47
1:D:230:LEU:HD23	1:D:230:LEU:C	2.40	0.47
2:B:38:GLN:OE1	2:B:81:ARG:NH1	2.47	0.47
1:A:148:GLU:CD	5:A:404:HOH:O	2.57	0.47
1:A:73:SER:OG	5:A:403:HOH:O	2.11	0.46
1:D:178:THR:O	1:D:181:ARG:NH2	2.43	0.46
1:A:178:THR:O	1:A:181:ARG:NH2	2.43	0.46
1:D:159:TYR:CG	3:F:3:PHE:HD1	2.34	0.45
1:A:274:TRP:O	1:A:275:GLU:HB3	2.17	0.45
1:A:219:LEU:HD11	1:A:256:TYR:HB3	1.99	0.45
1:A:66:LYS:HG2	3:C:4:ALA:HB1	1.99	0.45
1:D:148:GLU:OE2	5:D:701:HOH:O	2.20	0.44
2:E:32:PRO:HB2	2:E:33:PRO:HD2	2.00	0.44
1:D:23:MET:HE1	2:E:54:MET:HE3	1.99	0.44
1:D:263:HIS:HD2	1:D:265:GLY:H	1.64	0.44
1:D:77:ASP:HB3	3:F:9:LEU:HD12	1.99	0.43
1:D:155:ARG:NH1	3:F:5[B]:ILE:HG23	2.33	0.43
1:A:263:HIS:HD2	1:A:265:GLY:H	1.64	0.43
1:D:216:THR:OG1	1:D:260:HIS:HB2	2.18	0.43
2:E:1:ILE:O	2:E:1:ILE:HG23	2.19	0.43
1:D:216:THR:O	1:D:259:CYS:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LEU:O	1:A:246:SER:HA	2.19	0.43
1:A:9:VAL:O	1:A:96:GLN:HA	2.19	0.42
1:A:216:THR:O	1:A:259:CYS:HA	2.19	0.42
1:D:148:GLU:CD	5:D:701:HOH:O	2.61	0.42
2:B:32:PRO:HB2	2:B:33:PRO:HD2	2.02	0.42
2:B:1:ILE:HG23	2:B:1:ILE:O	2.19	0.42
1:A:218:GLN:NE2	5:A:408:HOH:O	2.51	0.42
1:D:201:LEU:O	1:D:246:SER:HA	2.19	0.42
1:A:152:GLU:HG2	3:C:7:ALA:CB	2.50	0.42
1:D:9:VAL:O	1:D:96:GLN:HA	2.20	0.41
1:D:219:LEU:HD11	1:D:256:TYR:HB3	2.02	0.41
1:A:230:LEU:HD21	1:A:243:LYS:NZ	2.35	0.41
1:D:155:ARG:HH12	3:F:5[B]:ILE:HG23	1.86	0.41
1:A:199:VAL:HG21	1:A:254:GLU:CG	2.52	0.40
1:D:234:ARG:HH11	2:E:8:GLN:NE2	2.19	0.40

All (2) 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)
1:A:65:GLN:OE1	2:B:94:TYR:OH[2_646]	1.95	0.25
1:D:65:GLN:OE1	2:E:94:TYR:OH[2_755]	2.04	0.16

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	273/275 (99%)	257 (94%)	16 (6%)	0	100	100
1	D	273/275 (99%)	256 (94%)	17 (6%)	0	100	100
2	B	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
2	E	97/99 (98%)	94 (97%)	2 (2%)	1 (1%)	12	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	9/9 (100%)	7 (78%)	0	2 (22%)	0	0
All	All	756/766 (99%)	714 (94%)	39 (5%)	3 (0%)	36	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	4[A]	ALA
3	F	4[B]	ALA
2	E	53	ASP

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	233/233 (100%)	224 (96%)	9 (4%)	28	39
1	D	233/233 (100%)	222 (95%)	11 (5%)	23	30
2	B	93/93 (100%)	89 (96%)	4 (4%)	26	34
2	E	93/93 (100%)	90 (97%)	3 (3%)	34	46
3	C	6/6 (100%)	6 (100%)	0	100	100
3	F	7/6 (117%)	4 (57%)	3 (43%)	0	0
All	All	665/664 (100%)	635 (96%)	30 (4%)	25	32

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	154	GLU
1	A	169	ARG
1	A	180	LEU
1	A	220	ASN
1	A	227	ASP
1	A	248	VAL

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Mol	Chain	Res	Type
1	A	251	LEU
1	A	272	LEU
2	B	57	SER
2	B	58	LYS
2	B	69	GLU
2	B	89	GLU
1	D	10	THR
1	D	138	MET
1	D	154	GLU
1	D	169	ARG
1	D	180	LEU
1	D	191	HIS
1	D	220	ASN
1	D	222	GLU
1	D	227	ASP
1	D	248	VAL
1	D	251	LEU
2	E	57	SER
2	E	58	LYS
2	E	69	GLU
3	F	1	SER
3	F	5[A]	ILE
3	F	5[B]	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	86	ASN
1	A	127	ASN
1	A	145	HIS
1	A	149	GLN
1	A	218	GLN
1	A	242	GLN
1	A	260	HIS
1	A	263	HIS
1	D	54	GLN
1	D	86	ASN
1	D	127	ASN
1	D	263	HIS
2	E	8	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](#)

Of 8 ligands modelled in this entry, 8 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	275/275 (100%)	0.92	39 (14%) 6 7	45, 71, 117, 153	0
1	D	275/275 (100%)	0.81	32 (11%) 9 11	42, 67, 112, 151	0
2	B	99/99 (100%)	0.96	9 (9%) 15 17	53, 71, 110, 128	0
2	E	99/99 (100%)	0.70	8 (8%) 18 20	46, 61, 93, 127	0
3	C	9/9 (100%)	1.04	2 (22%) 2 2	53, 59, 72, 82	0
3	F	9/9 (100%)	1.27	2 (22%) 2 2	29, 49, 60, 66	2 (22%)
All	All	766/766 (100%)	0.87	92 (12%) 9 10	29, 68, 112, 153	2 (0%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	ILE	7.8
2	E	1	ILE	7.1
1	D	1	GLY	5.6
2	E	3	LYS	5.2
1	A	1	GLY	5.1
2	B	3	LYS	4.8
1	D	223	GLU	4.4
1	D	2	PRO	4.4
1	D	107	GLY	4.2
1	A	171	TYR	4.2
1	A	2	PRO	4.1
1	D	109	LEU	4.0
1	A	51	TRP	3.8
1	A	49	ALA	3.6
3	F	5[A]	ILE	3.5
1	D	221	GLY	3.5
2	B	53	ASP	3.5
2	B	64	ILE	3.4
1	A	274	TRP	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	251	LEU	3.3
1	A	225	ILE	3.3
1	D	180	LEU	3.3
2	E	51	MET	3.2
1	D	17	LEU	3.2
1	A	103	VAL	3.1
1	A	180	LEU	3.1
1	D	178	THR	3.0
1	A	199	VAL	3.0
1	D	175	GLY	3.0
1	A	52	MET	3.0
1	D	225	ILE	3.0
1	A	179	LEU	2.9
1	A	248	VAL	2.9
1	A	17	LEU	2.9
1	D	192	HIS	2.8
2	B	31	HIS	2.8
1	A	33	PHE	2.8
1	A	176	ASN	2.8
1	A	107	GLY	2.8
1	A	6	ARG	2.7
1	A	192	HIS	2.7
1	D	103	VAL	2.7
2	E	53	ASP	2.6
1	D	110	LEU	2.6
2	B	51	MET	2.6
1	D	251	LEU	2.6
1	A	158	ALA	2.6
1	D	177	ALA	2.6
1	D	248	VAL	2.6
1	A	113	TYR	2.6
3	C	5	ILE	2.5
1	D	113	TYR	2.5
2	B	74	GLU	2.4
1	A	219	LEU	2.4
1	A	75	ARG	2.4
2	B	49	VAL	2.4
1	D	6	ARG	2.4
1	D	49	ALA	2.4
1	A	3	HIS	2.4
1	A	174	ASN	2.4
1	A	31	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	57	PRO	2.3
1	A	130	LEU	2.3
2	E	4	THR	2.3
3	C	1	SER	2.3
3	F	4[A]	ALA	2.3
1	A	172	LEU	2.3
1	A	138	MET	2.3
1	A	110	LEU	2.2
1	D	227	ASP	2.2
2	E	50	GLU	2.2
1	D	51	TRP	2.2
1	D	199	VAL	2.2
2	E	83	LYS	2.2
1	D	16	GLY	2.2
1	A	178	THR	2.2
1	D	158	ALA	2.2
1	A	169	ARG	2.1
1	D	224	LEU	2.1
2	E	58	LYS	2.1
1	D	54	GLN	2.1
1	D	163	THR	2.1
1	A	162	GLY	2.1
1	A	177	ALA	2.1
1	D	5	LEU	2.1
1	D	176	ASN	2.1
1	A	39	ASP	2.0
1	A	160	LEU	2.0
1	A	28	VAL	2.0
1	D	3	HIS	2.0
1	D	130	LEU	2.0
2	B	97	ARG	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(Å ²)	Q<0.9
4	NI	A	301	1/1	0.92	0.07	105,105,105,105	0
4	NI	E	601	1/1	0.92	0.08	92,92,92,92	0
4	NI	D	603	1/1	0.96	0.07	88,88,88,88	0
4	NI	B	601	1/1	0.98	0.04	80,80,80,80	0
4	NI	B	602	1/1	0.98	0.04	85,85,85,85	0
4	NI	E	602	1/1	0.98	0.04	89,89,89,89	0
4	NI	D	602	1/1	0.99	0.06	74,74,74,74	0
4	NI	D	601	1/1	0.99	0.07	79,79,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.