Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2023 – 07:10 AM EDT

PDB ID	:	2I3R
Title	:	Engineered catalytic domain of protein tyrosine phosphatase HPTPbeta
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Deposited on	:	2006-08-20
Resolution	:	1.85 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:

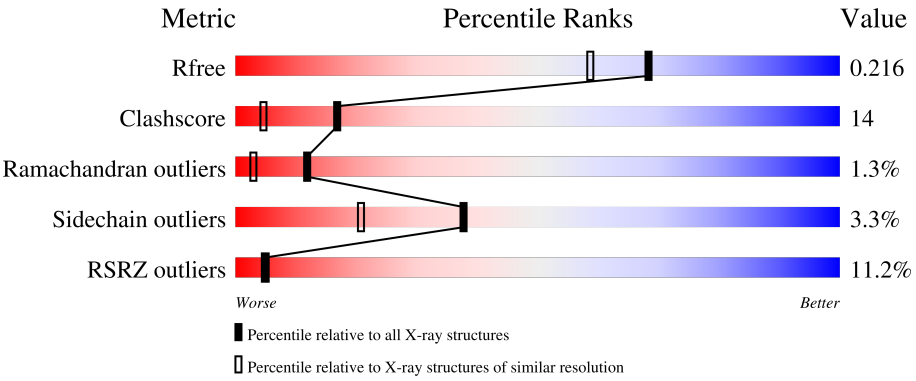
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 <=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	313	8% 72% 17% • 9%
1	B	313	12% 67% 19% • 11%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4870 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 Receptor-type tyrosine-protein phosphatase beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total 2376	C 1502	N 426	O 432	S 16	0	9	0
1	B	278	Total 2302	C 1459	N 409	O 419	S 15	0	7	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1661	SER	-	cloning artifact	UNP P23467
B	1661	SER	-	cloning artifact	UNP P23467

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is water.

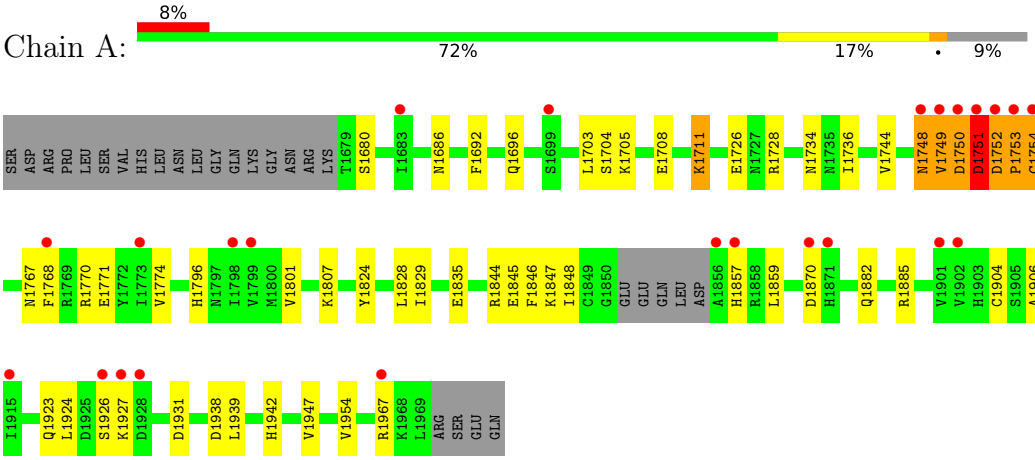
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total 123	O 123	0	0
3	B	67	Total 67	O 67	0	0

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.

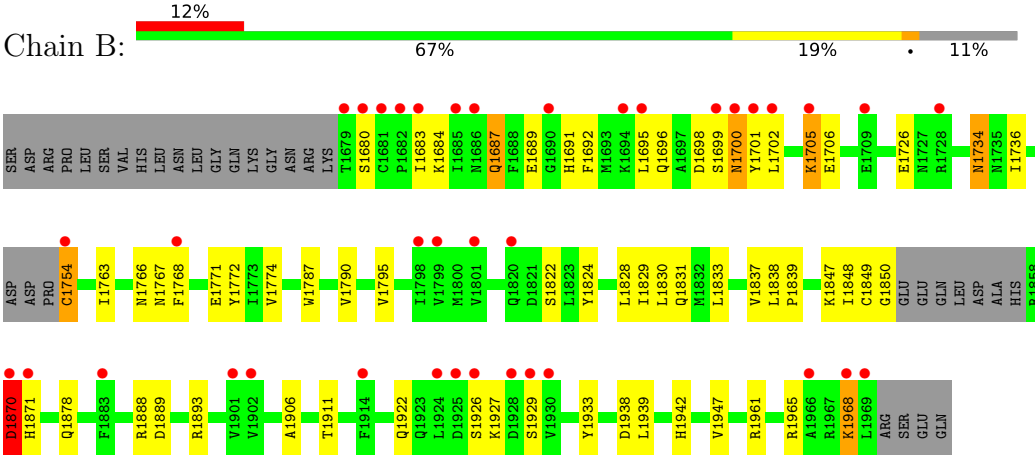
• Molecule 1: Receptor-type tyrosine-protein phosphatase beta

Chain A:



• Molecule 1: Receptor-type tyrosine-protein phosphatase beta

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	61.86Å 71.64Å 70.53Å 90.00° 93.58° 90.00°	Depositor
Resolution (Å)	31.90 – 1.85 31.93 – 1.85	Depositor EDS
% Data completeness (in resolution range)	98.6 (31.90-1.85) 98.6 (31.93-1.85)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.88 (at 1.85Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.217 0.182 , 0.216	Depositor DCC
R_{free} test set	2648 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	26.8	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 54.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4870	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.75% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.73	0/2459	0.76	0/3336
1	B	0.51	0/2377	0.61	0/3221
All	All	0.63	0/4836	0.69	0/6557

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	1
1	B	0	3
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1751	ASP	Peptide
1	B	1699	SER	Peptide
1	B	1869	PRO	Peptide
1	B	1870	ASP	Peptide

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	2376	0	2345	67	3
1	B	2302	0	2268	64	3
2	A	2	0	0	2	0
3	A	123	0	0	5	0
3	B	67	0	0	3	0
All	All	4870	0	4613	131	3

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 14.

All (131) 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:1726:GLU:OE2	1:A:1754:CYS:HB2	1.23	1.30
1:B:1726:GLU:OE2	1:B:1754:CYS:CB	1.84	1.24
1:B:1768[B]:PHE:CE1	1:B:1771:GLU:HG2	1.72	1.24
1:B:1726:GLU:OE2	1:B:1754:CYS:HB2	0.91	1.07
1:B:1684:LYS:HB2	1:B:1687:GLN:HG3	1.34	1.05
1:B:1696:GLN:HA	1:B:1700:ASN:HB3	1.40	1.04
1:B:1828:LEU:HD13	1:B:1848:ILE:HD11	1.46	0.98
1:A:1726:GLU:OE2	1:A:1754:CYS:CB	2.13	0.96
1:B:1726:GLU:CD	1:B:1754:CYS:HB2	1.85	0.95
1:B:1768[B]:PHE:CZ	1:B:1771:GLU:HG2	2.02	0.94
1:B:1768[B]:PHE:CE1	1:B:1771:GLU:CG	2.59	0.83
1:A:1807:LYS:HG3	1:A:1870:ASP:OD2	1.80	0.81
1:A:1748:ASN:ND2	1:A:1752:ASP:CB	2.43	0.81
1:B:1726:GLU:HG2	3:B:21:HOH:O	1.78	0.81
1:A:1923:GLN:O	1:A:1927:LYS:HG3	1.86	0.76
1:B:1696:GLN:HA	1:B:1700:ASN:CB	2.16	0.75
1:A:1726:GLU:CD	1:A:1754:CYS:HB2	2.05	0.75
1:A:1748:ASN:ND2	1:A:1752:ASP:HB2	2.00	0.74
1:B:1684:LYS:HB2	1:B:1687:GLN:CG	2.17	0.74
1:A:1752:ASP:H	1:A:1753:PRO:HD3	1.51	0.74
1:A:1752:ASP:HA	1:A:1755:SER:OG	1.88	0.73
1:B:1726:GLU:CD	1:B:1754:CYS:CB	2.51	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1683:ILE:HG13	1:B:1691:HIS:CD2	2.23	0.73
1:B:1692:PHE:O	1:B:1696:GLN:HG3	1.88	0.72
1:A:1680:SER:HB2	1:A:1931:ASP:OD2	1.91	0.71
1:B:1869:PRO:HA	1:B:1870:ASP:HB2	1.74	0.70
1:B:1705:LYS:HG2	1:B:1706:GLU:N	2.07	0.67
1:A:1750:ASP:O	1:A:1751:ASP:HB2	1.98	0.63
1:B:1838:LEU:HB3	1:B:1839:PRO:CD	2.28	0.63
1:A:1924:LEU:O	1:A:1967:ARG:NH2	2.29	0.63
1:A:1726:GLU:OE1	1:A:1754:CYS:N	2.22	0.63
1:B:1768[B]:PHE:CD1	1:B:1771:GLU:CG	2.82	0.63
1:B:1766:ASN:HD22	1:B:1922:GLN:HB3	1.65	0.62
1:B:1768[B]:PHE:CD1	1:B:1771:GLU:HG2	2.31	0.61
1:A:1726:GLU:CD	1:A:1754:CYS:H	2.03	0.61
1:B:1838:LEU:HB3	1:B:1839:PRO:HD2	1.83	0.61
1:A:1752:ASP:H	1:A:1753:PRO:CD	2.13	0.61
1:B:1870:ASP:O	1:B:1871:HIS:HD2	1.84	0.61
1:B:1961:ARG:HG2	1:B:1965:ARG:HD2	1.83	0.61
1:A:1749:VAL:HG12	1:A:1750:ASP:N	2.16	0.61
1:A:1847:LYS:HE3	1:A:1857:HIS:CG	2.36	0.60
1:B:1870:ASP:O	1:B:1871:HIS:CD2	2.55	0.60
1:A:1726:GLU:CD	1:A:1754:CYS:CB	2.68	0.59
1:A:1750:ASP:O	1:A:1751:ASP:CB	2.49	0.59
1:A:1845:GLU:OE2	1:A:1859[B]:LEU:HD21	2.03	0.58
1:A:1748:ASN:ND2	1:A:1752:ASP:HB3	2.18	0.58
1:A:1767[B]:ASN:HB2	1:A:1771:GLU:OE1	2.05	0.57
1:B:1766:ASN:ND2	1:B:1922:GLN:CB	2.68	0.57
1:A:1752:ASP:N	1:A:1753:PRO:CD	2.67	0.57
1:A:1762:TYR:CD1	1:A:1770[A]:ARG:HG2	2.40	0.57
1:B:1830:LEU:HD12	1:B:1848:ILE:HB	1.87	0.57
1:B:1698:ASP:HB3	1:B:1701:TYR:HB3	1.88	0.56
1:A:1748:ASN:HD22	1:A:1752:ASP:HB3	1.71	0.56
1:A:1923:GLN:NE2	3:A:85:HOH:O	2.20	0.55
1:B:1683:ILE:O	1:B:1929:SER:HB2	2.06	0.55
1:B:1889:ASP:O	1:B:1893:ARG:HG3	2.06	0.55
1:B:1869:PRO:CA	1:B:1870:ASP:HB2	2.37	0.55
1:B:1702:LEU:HD23	1:B:1705:LYS:HD3	1.87	0.55
1:A:1828:LEU:HD13	1:A:1848:ILE:HD11	1.87	0.54
1:B:1766:ASN:ND2	1:B:1922:GLN:HB2	2.22	0.54
1:A:1845:GLU:HB3	1:A:1859[B]:LEU:HD11	1.89	0.54
1:A:1752:ASP:O	1:A:1753:PRO:O	2.26	0.54
1:B:1768[B]:PHE:CZ	1:B:1771:GLU:HA	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1744:VAL:O	2:A:301:CL:CL	2.64	0.53
1:A:1726:GLU:OE1	1:A:1753:PRO:HB2	2.09	0.52
1:B:1766:ASN:OD1	3:B:104:HOH:O	2.19	0.52
1:A:1938:ASP:O	1:A:1942:HIS:HD2	1.92	0.51
1:A:1770[B]:ARG:NE	3:A:128:HOH:O	2.41	0.50
1:A:1711:LYS:HD3	3:A:4:HOH:O	2.11	0.50
1:A:1750:ASP:N	1:A:1750:ASP:OD1	2.34	0.49
1:B:1774:VAL:HG22	1:B:1911:THR:CG2	2.42	0.49
1:B:1965:ARG:O	1:B:1968:LYS:HG3	2.12	0.49
1:B:1766:ASN:ND2	1:B:1922:GLN:HB3	2.26	0.49
1:B:1695:LEU:HA	1:B:1702:LEU:HB2	1.95	0.49
1:A:1763:ILE:HG21	1:A:1939:LEU:HD22	1.94	0.49
1:A:1763:ILE:CG2	1:A:1939:LEU:HD22	2.42	0.48
1:A:1708:GLU:O	1:A:1711:LYS:HB2	2.14	0.48
1:B:1695:LEU:O	1:B:1700:ASN:HA	2.14	0.48
1:A:1846:PHE:O	1:A:1859[B]:LEU:HD12	2.12	0.48
1:B:1830:LEU:CD1	1:B:1848:ILE:HB	2.44	0.48
1:B:1774:VAL:HG22	1:B:1911:THR:HG21	1.96	0.47
1:A:1703:LEU:CD1	1:A:1954:VAL:HG22	2.44	0.47
1:A:1774[A]:VAL:HG22	1:A:1911:THR:CG2	2.44	0.47
1:A:1801:VAL:HG23	1:A:1904:CYS:HB3	1.97	0.47
1:B:1833:LEU:HD12	1:B:1859:LEU:HD21	1.95	0.47
1:B:1754:CYS:HA	3:B:96:HOH:O	2.14	0.47
1:B:1926:SER:C	1:B:1927:LYS:HD3	2.36	0.47
1:A:1726:GLU:CD	1:A:1754:CYS:N	2.68	0.46
1:A:1748:ASN:HD22	1:A:1752:ASP:CB	2.23	0.46
1:B:1938:ASP:O	1:B:1942:HIS:HD2	1.98	0.46
1:A:1807:LYS:CG	1:A:1870:ASP:OD2	2.59	0.46
1:B:1689:GLU:OE1	1:B:1961:ARG:HG3	2.16	0.46
1:A:1824:TYR:CE1	1:A:1829:ILE:HG12	2.51	0.46
1:A:1910:ARG:HG3	3:A:65:HOH:O	2.16	0.46
1:A:1686:ASN:OD1	1:A:1686:ASN:C	2.54	0.45
1:B:1795:VAL:O	1:B:1860:ILE:CD1	2.64	0.45
1:B:1736:ILE:HG21	1:B:1906:ALA:HB1	1.97	0.45
1:A:1736:ILE:HG21	1:A:1906:ALA:HB1	1.98	0.45
1:B:1831[B]:GLN:HG3	1:B:1847:LYS:HB3	1.98	0.45
1:B:1822[B]:SER:OG	1:B:1831[B]:GLN:HB3	2.17	0.45
1:B:1828:LEU:CD2	1:B:1850:GLY:HA3	2.47	0.45
1:A:1923:GLN:NE2	2:A:302:CL:CL	2.76	0.44
1:A:1848:ILE:O	1:A:1857:HIS:HA	2.18	0.44
1:A:1923:GLN:O	1:A:1927:LYS:CG	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1774[A]:VAL:HG22	1:A:1911:THR:HG21	2.00	0.43
1:A:1926:SER:OG	1:A:1927:LYS:HG2	2.18	0.43
1:B:1706:GLU:HG2	1:B:1933:TYR:CD1	2.53	0.43
1:B:1837:VAL:O	1:B:1838:LEU:HD23	2.18	0.43
1:A:1752:ASP:N	1:A:1753:PRO:HD3	2.24	0.43
1:A:1824:TYR:CD1	1:A:1829:ILE:HG12	2.54	0.43
1:A:1692:PHE:O	1:A:1696:GLN:HG3	2.19	0.42
1:A:1726:GLU:OE1	1:A:1753:PRO:CB	2.67	0.42
1:B:1768[B]:PHE:CZ	1:B:1771:GLU:CG	2.90	0.42
1:B:1824:TYR:CE1	1:B:1829:ILE:HG12	2.55	0.42
1:B:1837:VAL:C	1:B:1838:LEU:HD23	2.40	0.42
1:B:1702:LEU:CD2	1:B:1705:LYS:HD3	2.48	0.42
1:A:1752:ASP:HA	1:A:1755:SER:HG	1.84	0.42
1:A:1835:GLU:HG3	1:A:1844:ARG:HG2	2.02	0.42
1:B:1734:ASN:HD22	1:B:1734:ASN:HA	1.66	0.42
1:A:1704:SER:O	1:A:1708:GLU:HG2	2.20	0.41
1:A:1923:GLN:HA	1:A:1927:LYS:HG3	2.03	0.41
1:B:1766:ASN:HD22	1:B:1922:GLN:CB	2.27	0.41
1:A:1882:GLN:HE22	1:A:1885:ARG:HH11	1.68	0.41
1:B:1763:ILE:HB	1:B:1772:TYR:HB2	2.03	0.41
1:B:1888:ARG:O	1:B:1888:ARG:HD3	2.20	0.41
1:A:1734:ASN:HD22	1:A:1734:ASN:HA	1.73	0.41
1:B:1787:TRP:CH2	1:B:1830:LEU:HD21	2.56	0.40
1:A:1696:GLN:NE2	3:A:131:HOH:O	2.53	0.40
1:A:1751:ASP:N	1:A:1752:ASP:OD1	2.50	0.40
1:B:1790:VAL:HG13	1:B:1795:VAL:HB	2.04	0.40

All (3) 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:1768:PHE:CE1	1:B:1768[B]:PHE:CE1[2_646]	2.07	0.13
1:A:1768:PHE:CD1	1:B:1768[B]:PHE:CZ[2_646]	2.10	0.10
1:A:1768:PHE:CZ	1:B:1768[B]:PHE:CE1[2_646]	2.12	0.08

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	291/313 (93%)	278 (96%)	8 (3%)	5 (2%)	9 2
1	B	279/313 (89%)	263 (94%)	14 (5%)	2 (1%)	22 9
All	All	570/626 (91%)	541 (95%)	22 (4%)	7 (1%)	12 3

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1753	PRO
1	B	1700	ASN
1	A	1751	ASP
1	A	1752	ASP
1	A	1749	VAL
1	A	1947	VAL
1	B	1947	VAL

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	268/284 (94%)	262 (98%)	6 (2%)	52 36
1	B	259/284 (91%)	248 (96%)	11 (4%)	30 13
All	All	527/568 (93%)	510 (97%)	17 (3%)	38 22

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1705	LYS
1	A	1711	LYS
1	A	1728	ARG
1	A	1748	ASN
1	A	1750	ASP
1	A	1754	CYS
1	B	1680	SER
1	B	1687	GLN
1	B	1705	LYS
1	B	1734	ASN
1	B	1754	CYS
1	B	1767	ASN
1	B	1849	CYS
1	B	1870	ASP
1	B	1878	GLN
1	B	1939	LEU
1	B	1968	LYS

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	1696	GLN
1	A	1734	ASN
1	A	1748	ASN
1	A	1882	GLN
1	A	1942	HIS
1	B	1691	HIS
1	B	1734	ASN
1	B	1871	HIS
1	B	1878	GLN
1	B	1882	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 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, 95^{th} 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	286/313 (91%)	0.52	24 (8%)	11	10	13, 23, 40, 54	6 (2%)
1	B	278/313 (88%)	0.71	39 (14%)	2	3	22, 32, 48, 60	0
All	All	564/626 (90%)	0.62	63 (11%)	5	5	13, 28, 46, 60	6 (1%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1753	PRO	13.1
1	A	1749	VAL	12.7
1	A	1750	ASP	10.7
1	A	1752	ASP	8.7
1	B	1754	CYS	7.7
1	B	1701	TYR	6.1
1	A	1751	ASP	6.0
1	A	1754	CYS	5.2
1	B	1702	LEU	5.1
1	A	1856	ALA	4.7
1	A	1857	HIS	4.7
1	B	1695	LEU	4.6
1	B	1681	CYS	4.3
1	B	1705	LYS	4.2
1	B	1925	ASP	4.2
1	A	1927	LYS	3.9
1	A	1748	ASN	3.9
1	A	1902	VAL	3.8
1	B	1968	LYS	3.6
1	B	1871	HIS	3.4
1	B	1799	VAL	3.3
1	A	1798	ILE	3.3
1	B	1928	ASP	3.3
1	A	1901	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	1966	ALA	3.1
1	B	1902	VAL	3.0
1	B	1685	ILE	3.0
1	B	1683	ILE	2.9
1	B	1798	ILE	2.9
1	B	1680	SER	2.9
1	B	1820	GLN	2.9
1	A	1870	ASP	2.7
1	B	1914	PHE	2.7
1	B	1686	ASN	2.7
1	A	1799	VAL	2.7
1	A	1928	ASP	2.7
1	A	1768	PHE	2.6
1	B	1901	VAL	2.6
1	B	1969	LEU	2.6
1	B	1768[A]	PHE	2.5
1	A	1915	ILE	2.5
1	B	1700	ASN	2.5
1	B	1926	SER	2.5
1	A	1967	ARG	2.5
1	A	1699	SER	2.5
1	A	1773	ILE	2.4
1	B	1694	LYS	2.4
1	A	1683	ILE	2.4
1	B	1801	VAL	2.3
1	B	1924	LEU	2.3
1	A	1871	HIS	2.3
1	A	1926	SER	2.3
1	B	1870	ASP	2.3
1	B	1690	GLY	2.2
1	B	1930	VAL	2.2
1	B	1699	SER	2.2
1	B	1679	THR	2.2
1	B	1728	ARG	2.1
1	B	1929	SER	2.1
1	B	1682	PRO	2.1
1	B	1859	LEU	2.1
1	B	1709	GLU	2.0
1	B	1883	PHE	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 monosaccharides 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, 95^{th} 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
2	CL	A	302	1/1	0.97	0.27	43,43,43,43	0
2	CL	A	301	1/1	0.98	0.18	30,30,30,30	0

6.5 Other polymers (i)

There are no such residues in this entry.