



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:32 PM GMT

PDB ID : 1U4Q
Title : Crystal Structure of Repeats 15, 16 and 17 of Chicken Brain Alpha Spectrin
Authors : Kusunoki, H.; Minasov, G.; MacDonald, R.I.; Mondragon, A.
Deposited on : 2004-07-26
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

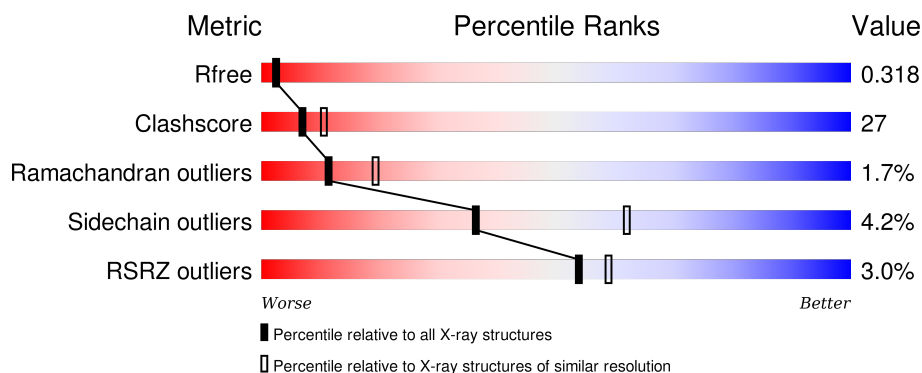
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	322	<div> <div>3%</div> <div>58%</div> <div>37%</div> <div>..</div> </div>
1	B	322	<div> <div>2%</div> <div>52%</div> <div>44%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5636 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 Spectrin alpha chain, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	5	0
			2604	1611	477	510	6			
1	B	317	Total	C	N	O	S	0	4	0
			2587	1599	473	509	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1661	MET	-	INITIATING METHIONINE	UNP P07751
B	1661	MET	-	INITIATING METHIONINE	UNP P07751

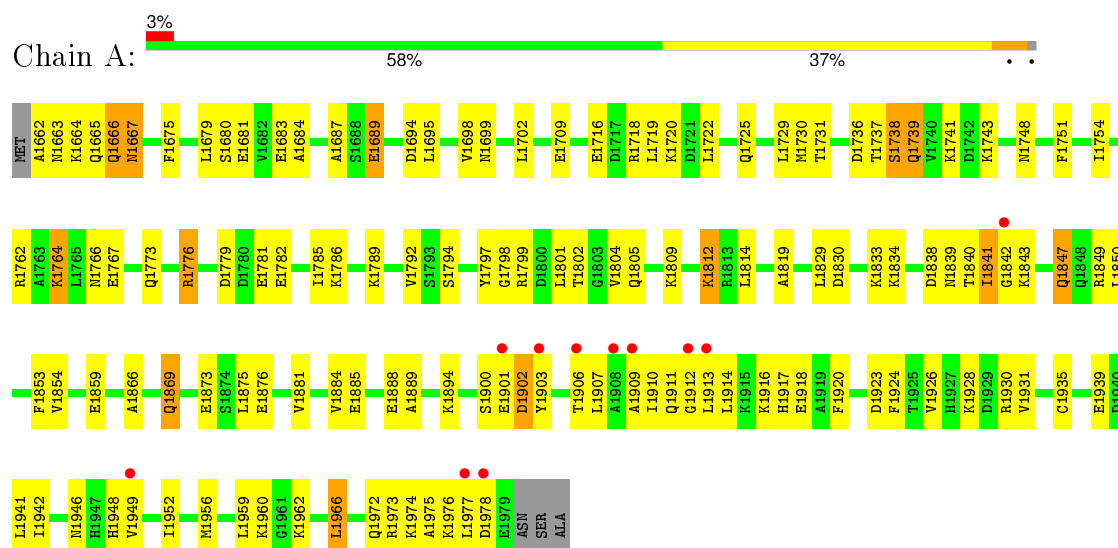
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	230	Total	O	0	0
			230	230		
2	B	215	Total	O	0	0
			215	215		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Spectrin alpha chain, brain



- Molecule 1: Spectrin alpha chain, brain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	150.03 Å 74.86 Å 66.61 Å 90.00° 97.70° 90.00°	Depositor
Resolution (Å)	14.94 – 2.50 26.70 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (14.94-2.50) 98.8 (26.70-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 2.20 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.313 0.229 , 0.318	Depositor DCC
R_{free} test set	2497 reflections (9.96%)	DCC
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.630	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 36806 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5636	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0/2638	0.53	0/3532
1	B	0.36	0/2621	0.55	0/3511
All	All	0.36	0/5259	0.54	0/7043

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 ⓘ

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	2604	0	2581	137	0
1	B	2587	0	2556	142	0
2	A	230	0	0	25	0
2	B	215	0	0	31	0
All	All	5636	0	5137	275	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 27.

All (275) 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:1764:LYS:HB3	1:A:1841:ILE:HD11	1.35	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1739:GLN:HA	1:A:1739:GLN:HE21	1.21	1.01
1:A:1847:GLN:HE21	1:A:1847:GLN:HA	1.25	0.99
1:A:1719:LEU:HD11	2:A:76:HOH:O	1.61	0.97
1:B:1939:GLU:HA	1:B:1942:ILE:HD12	1.52	0.89
1:A:1924:PHE:O	1:A:1928:LYS:HG2	1.74	0.87
1:A:1948:HIS:O	1:A:1952:ILE:HG13	1.77	0.84
1:A:1812:LYS:HA	1:A:1812:LYS:HE3	1.59	0.84
1:B:1925:THR:HG21	2:B:225:HOH:O	1.79	0.82
1:B:1801:LEU:O	1:B:1805:GLN:HG3	1.80	0.81
1:A:1789:LYS:HD3	1:A:1859:GLU:OE1	1.82	0.80
1:A:1776:ARG:HD2	2:A:183:HOH:O	1.82	0.78
1:A:1889:ALA:HA	2:A:122:HOH:O	1.84	0.78
1:B:1830:ASP:HB2	2:B:150:HOH:O	1.84	0.78
1:A:1912:GLY:O	1:A:1916:LYS:HG2	1.84	0.77
1:A:1702:LEU:HD11	2:A:353:HOH:O	1.83	0.77
1:B:1739:GLN:NE2	1:B:1740:VAL:HG23	2.00	0.76
1:B:1938:GLY:O	1:B:1942:ILE:HG13	1.86	0.76
1:A:1809:LYS:HD2	2:A:238:HOH:O	1.85	0.76
1:B:1962:LYS:HB3	1:B:1962:LYS:NZ	2.00	0.76
1:A:1663:ASN:O	1:A:1666:GLN:HG2	1.85	0.76
1:B:1895:MET:CE	1:B:1966:LEU:HA	2.16	0.76
1:B:1946:ASN:O	1:B:1949:VAL:HG23	1.87	0.75
1:A:1812:LYS:HE2	2:A:359:HOH:O	1.85	0.75
1:B:1866:ALA:O	1:B:1869:GLN:HG2	1.88	0.74
1:B:1949:VAL:HA	1:B:1952:ILE:HG22	1.71	0.73
1:B:1895:MET:HE1	1:B:1966:LEU:HA	1.70	0.72
1:A:1946:ASN:O	1:A:1949:VAL:HG23	1.89	0.72
1:A:1911:GLN:OE1	1:A:1914:LEU:HD23	1.88	0.72
1:B:1936:ALA:HA	1:B:1939:GLU:HG2	1.72	0.71
1:A:1942:ILE:HD12	1:A:1949:VAL:HG13	1.73	0.71
1:B:1923:ASP:O	1:B:1926:VAL:HG22	1.91	0.70
1:B:1942:ILE:HD11	1:B:1956:MET:SD	2.30	0.70
1:A:1739:GLN:HA	1:A:1739:GLN:NE2	2.03	0.69
1:A:1866:ALA:O	1:A:1869:GLN:HG3	1.94	0.68
1:A:1743:LYS:HG3	2:A:421:HOH:O	1.91	0.68
1:A:1935:CYS:SG	1:A:1960:LYS:HE3	2.34	0.68
1:B:1926:VAL:O	1:B:1929:ASP:HB2	1.94	0.67
1:A:1773:GLN:HB2	2:A:391:HOH:O	1.95	0.67
1:A:1812:LYS:CE	1:A:1812:LYS:HA	2.25	0.67
1:B:1679:LEU:O	1:B:1683:GLU:HG3	1.95	0.67
1:B:1754:ILE:HG21	2:B:222:HOH:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1824:ALA:O	1:B:1828:VAL:HG23	1.95	0.66
1:B:1943:LYS:HG3	1:B:1944:LYS:N	2.09	0.66
1:B:1756:SER:HB3	2:B:295:HOH:O	1.95	0.66
1:A:1935:CYS:O	1:A:1939:GLU:HG2	1.96	0.65
1:B:1689:GLU:HG2	2:B:54:HOH:O	1.96	0.65
1:B:1808:ARG:O	1:B:1812:LYS:HG2	1.96	0.65
1:A:1776:ARG:HA	2:A:183:HOH:O	1.97	0.65
1:A:1833[B]:LYS:HG3	1:A:1834:LYS:N	2.11	0.65
1:B:1781:GLU:HG3	1:B:1828:VAL:HG21	1.79	0.65
1:B:1907:LEU:O	1:B:1910:ILE:HG22	1.97	0.64
1:A:1679:LEU:O	1:A:1683:GLU:HG3	1.98	0.64
1:B:1739:GLN:HE22	1:B:1740:VAL:HG23	1.62	0.63
1:B:1980:ASN:HB3	2:B:336:HOH:O	1.98	0.63
1:B:1850:LEU:O	1:B:1854:VAL:HG23	1.99	0.63
1:B:1788:LYS:HE3	1:B:1817:GLU:OE2	1.98	0.63
1:A:1914:LEU:O	1:A:1918:GLU:HG2	1.98	0.62
1:B:1914:LEU:HD23	1:B:1977:LEU:HD13	1.80	0.62
1:B:1822:GLU:N	1:B:1823:PRO:CD	2.63	0.61
1:B:1699:ASN:O	1:B:1703[A]:LYS:HG2	2.00	0.61
1:A:1843:LYS:O	1:A:1847:GLN:HG2	2.01	0.61
1:A:1942:ILE:HD11	1:A:1952:ILE:HB	1.82	0.61
1:A:1766:ASN:HA	2:A:353:HOH:O	2.00	0.60
1:B:1903:TYR:HB3	2:B:336:HOH:O	2.00	0.60
1:B:1835:LEU:HD22	1:B:1840:THR:OG1	2.01	0.60
1:B:1943:LYS:HG3	1:B:1944:LYS:H	1.66	0.60
1:B:1773:GLN:O	1:B:1776:ARG:HB3	2.00	0.60
1:A:1748:ASN:HA	2:A:76:HOH:O	2.02	0.60
1:B:1859:GLU:O	1:B:1863:LEU:HG	2.02	0.59
1:B:1667:ASN:HD22	1:B:1667:ASN:N	1.99	0.59
1:B:1819:ALA:O	1:B:1822:GLU:HG2	2.03	0.59
1:A:1914:LEU:HD13	1:A:1977:LEU:HB3	1.84	0.59
1:A:1909:ALA:O	1:A:1913:LEU:HG	2.02	0.59
1:A:1762[B]:ARG:HD2	2:A:289:HOH:O	2.03	0.59
1:A:1797:TYR:HE1	2:A:221:HOH:O	1.86	0.59
1:B:1948:HIS:O	1:B:1952:ILE:HB	2.03	0.58
1:A:1773:GLN:O	1:A:1776:ARG:HB3	2.03	0.58
1:A:1664:LYS:HB2	1:A:1729:LEU:HD21	1.85	0.58
1:B:1879:GLN:HB2	2:B:198:HOH:O	2.03	0.58
1:B:1781:GLU:OE1	1:B:1821:HIS:HE1	1.86	0.58
1:A:1924:PHE:CZ	1:A:1966:LEU:HD13	2.39	0.57
1:A:1830:ASP:O	1:A:1833[B]:LYS:HG2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1907:LEU:O	1:A:1907:LEU:HD23	2.04	0.57
1:B:1867:ARG:HA	2:B:191:HOH:O	2.02	0.57
1:B:1761:ARG:HA	2:B:206:HOH:O	2.05	0.57
1:A:1751:PHE:O	1:A:1754:ILE:HG12	2.04	0.57
1:A:1689:GLU:OE1	1:A:1689:GLU:N	2.33	0.57
1:A:1731:THR:HG22	1:A:1731:THR:O	2.05	0.57
1:B:1959:LEU:O	1:B:1963:VAL:HG23	2.04	0.57
1:A:1876:GLU:HA	2:A:305:HOH:O	2.05	0.56
1:B:1953:THR:HA	1:B:1956:MET:CE	2.35	0.56
1:B:1942:ILE:HG23	2:B:291:HOH:O	2.03	0.56
1:A:1737:THR:O	1:A:1739:GLN:N	2.39	0.56
1:B:1838:ASP:O	1:B:1839:ASN:C	2.43	0.56
1:B:1939:GLU:O	1:B:1942:ILE:HB	2.06	0.56
1:B:1970:ALA:HB1	2:B:215:HOH:O	2.04	0.56
1:B:1925:THR:O	1:B:1928:LYS:HG2	2.05	0.56
1:A:1942:ILE:HG23	1:A:1949:VAL:HG13	1.86	0.56
1:B:1872:GLU:HG2	2:B:185:HOH:O	2.05	0.56
1:B:1942:ILE:HA	2:B:428:HOH:O	2.05	0.56
1:B:1764:LYS:HD2	2:B:403:HOH:O	2.05	0.56
1:B:1945:ASN:OD1	1:B:1945:ASN:O	2.24	0.56
1:A:1814:LEU:HD23	1:A:1814:LEU:O	2.06	0.55
1:B:1770:ARG:HD3	1:B:1835:LEU:HD21	1.88	0.55
1:B:1881:VAL:HG22	1:B:1955:LYS:HE3	1.88	0.55
1:B:1940:ASP:O	1:B:1943:LYS:HG2	2.05	0.55
1:A:1847:GLN:HE21	1:A:1847:GLN:CA	2.03	0.55
1:A:1664:LYS:HD3	1:A:1729:LEU:HD23	1.89	0.55
1:A:1801:LEU:O	1:A:1805:GLN:HG3	2.07	0.55
1:A:1767:GLU:HG2	1:A:1840:THR:HG23	1.90	0.54
1:B:1903:TYR:O	1:B:1910:ILE:HD13	2.08	0.54
1:A:1849[A]:ARG:HB3	1:A:1849[A]:ARG:NH1	2.22	0.54
1:A:1952:ILE:HG22	1:A:1956:MET:CE	2.38	0.54
1:A:1869:GLN:O	1:A:1873:GLU:HG3	2.08	0.54
1:B:1976:LYS:HE3	2:B:376:HOH:O	2.07	0.54
1:B:1718[A]:ARG:NH1	2:B:126:HOH:O	2.40	0.54
1:A:1952:ILE:HG22	1:A:1956:MET:HE1	1.90	0.54
1:B:1962:LYS:HB3	1:B:1962:LYS:HZ3	1.70	0.54
1:A:1829:LEU:N	1:A:1829:LEU:HD23	2.23	0.54
1:B:1870:ARG:HH21	1:B:1947:HIS:CG	2.26	0.53
1:A:1767:GLU:OE2	1:A:1840:THR:HA	2.07	0.53
1:B:1925:THR:HA	1:B:1928:LYS:HD3	1.90	0.53
1:A:1702:LEU:HD21	2:A:353:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1910:ILE:HG22	1:A:1911:GLN:N	2.22	0.53
1:A:1939:GLU:OE2	1:A:1956:MET:SD	2.67	0.53
1:B:1926:VAL:CG2	1:B:1927:HIS:N	2.72	0.53
1:A:1792:VAL:O	1:A:1792:VAL:HG12	2.08	0.53
1:B:1821:HIS:C	1:B:1823:PRO:HD2	2.30	0.52
1:B:1889:ALA:HB3	2:B:313:HOH:O	2.09	0.52
1:B:1799:ARG:HB3	2:B:400:HOH:O	2.09	0.52
1:B:1771:LEU:O	1:B:1774:PHE:HB3	2.09	0.52
1:A:1680:SER:HA	1:A:1683:GLU:OE1	2.10	0.52
1:B:1808:ARG:NH1	2:B:390:HOH:O	2.43	0.52
1:A:1917:HIS:CE1	1:A:1974:LYS:HB2	2.44	0.52
1:A:1794:SER:HB2	1:B:1790:LEU:HD11	1.91	0.51
1:A:1847:GLN:NE2	1:A:1847:GLN:HA	2.08	0.51
1:A:1903:TYR:HA	1:A:1913:LEU:HD11	1.91	0.51
1:B:1962:LYS:HB3	1:B:1962:LYS:HZ2	1.74	0.51
1:A:1694:ASP:O	1:A:1698:VAL:HG23	2.11	0.51
1:A:1911:GLN:HA	1:A:1914:LEU:HB3	1.91	0.51
1:B:1926:VAL:HG23	1:B:1927:HIS:N	2.25	0.51
1:A:1907:LEU:C	1:A:1907:LEU:HD23	2.32	0.51
1:A:1972:GLN:O	1:A:1975:ALA:HB3	2.10	0.51
1:B:1760:ALA:O	1:B:1763:ALA:HB3	2.10	0.50
1:A:1840:THR:O	1:A:1842:GLY:N	2.44	0.50
1:A:1730:MET:HG2	1:A:1737:THR:CG2	2.40	0.50
1:A:1709:GLU:OE2	1:A:1762[A]:ARG:HD2	2.12	0.50
1:A:1804:VAL:HG11	1:A:1875:LEU:CA	2.41	0.50
1:B:1767:GLU:HG2	2:B:197:HOH:O	2.10	0.50
1:A:1894:LYS:NZ	2:A:342:HOH:O	2.43	0.50
1:B:1910:ILE:O	1:B:1914:LEU:HG	2.11	0.50
1:B:1961:GLY:O	1:B:1964:SER:N	2.45	0.50
1:B:1768:SER:HB2	1:B:1841:ILE:HD11	1.94	0.50
1:B:1784:TRP:CH2	1:B:1818:LEU:HD23	2.47	0.50
1:A:1901:GLU:O	1:A:1902:ASP:C	2.49	0.50
1:B:1953:THR:HA	1:B:1956:MET:HE3	1.93	0.49
1:B:1750:ARG:O	1:B:1754:ILE:HG23	2.12	0.49
1:B:1781:GLU:O	1:B:1785:ILE:HG13	2.12	0.49
1:B:1767:GLU:HB3	1:B:1840:THR:HG23	1.94	0.49
1:B:1739:GLN:O	1:B:1743:LYS:HG3	2.12	0.49
1:B:1976:LYS:O	1:B:1980:ASN:HB2	2.12	0.49
1:A:1931:VAL:HG13	1:A:1959:LEU:HD21	1.94	0.49
1:A:1829:LEU:HD21	1:A:1853:PHE:CD2	2.46	0.49
1:B:1910:ILE:CD1	1:B:1977:LEU:HD22	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:159:HOH:O	1:B:1916:LYS:HG2	2.13	0.49
1:B:1891:ILE:HG23	1:B:1966:LEU:HD22	1.95	0.48
1:A:1850:LEU:O	1:A:1854:VAL:HG23	2.14	0.48
1:B:1869:GLN:O	1:B:1873:GLU:HG3	2.13	0.48
1:A:1739:GLN:HE21	1:A:1739:GLN:CA	2.06	0.48
1:B:1895:MET:HA	1:B:1895:MET:HE2	1.96	0.48
1:B:1731:THR:HG21	2:B:395:HOH:O	2.13	0.47
1:A:1764:LYS:CB	1:A:1841:ILE:HD11	2.26	0.47
1:A:1906:THR:O	1:A:1910:ILE:HD13	2.14	0.47
1:A:1802:THR:HA	1:A:1805:GLN:HE21	1.79	0.47
1:A:1974:LYS:NZ	1:A:1978:ASP:HB3	2.30	0.47
1:A:1894:LYS:HB3	1:A:1920:PHE:HE1	1.78	0.47
1:A:1926:VAL:O	1:A:1930:ARG:HG3	2.15	0.47
1:B:1921:GLU:O	1:B:1925:THR:HG23	2.15	0.47
1:B:1941:LEU:HB2	1:B:1952:ILE:HD11	1.96	0.47
1:A:1767:GLU:OE2	1:A:1841:ILE:N	2.43	0.47
1:B:1933:ASP:O	1:B:1936:ALA:HB3	2.15	0.47
1:A:1973:ARG:HH11	1:A:1973:ARG:HG2	1.78	0.47
1:A:1730:MET:HG2	1:A:1737:THR:HG23	1.96	0.47
1:A:1675:PHE:CZ	1:A:1719:LEU:HD13	2.49	0.47
1:A:1702:LEU:HD22	1:A:1762[B]:ARG:NH1	2.29	0.47
1:B:1849[B]:ARG:HG3	1:B:1849[B]:ARG:HH11	1.80	0.46
1:A:1776:ARG:O	1:A:1779:ASP:HB2	2.16	0.46
1:A:1730:MET:HE2	1:A:1741:LYS:HA	1.97	0.46
1:A:1840:THR:O	1:A:1841:ILE:C	2.54	0.46
1:B:1914:LEU:HD23	1:B:1977:LEU:CD1	2.44	0.46
1:A:1881:VAL:O	1:A:1885:GLU:HG3	2.15	0.46
1:B:1809:LYS:HE2	2:B:44:HOH:O	2.16	0.46
1:A:1665:GLN:NE2	2:A:392:HOH:O	2.47	0.46
1:A:1804:VAL:HG11	1:A:1875:LEU:HA	1.98	0.46
1:A:1781:GLU:O	1:A:1785:ILE:HG13	2.14	0.46
1:B:1943:LYS:HG3	1:B:1944:LYS:HG3	1.97	0.46
1:A:1924:PHE:CE1	1:A:1966:LEU:HD13	2.51	0.46
1:B:1869:GLN:HB2	2:B:420:HOH:O	2.15	0.45
1:B:1903:TYR:CD1	1:B:1903:TYR:N	2.83	0.45
1:A:1814:LEU:C	1:A:1814:LEU:HD23	2.37	0.45
1:A:1884:VAL:O	1:A:1888:GLU:HG3	2.17	0.45
1:B:1869:GLN:HG3	1:B:1870:ARG:N	2.30	0.45
1:B:1766:ASN:O	1:B:1770:ARG:HG2	2.16	0.45
1:A:1684:ALA:O	1:A:1687:ALA:HB3	2.17	0.45
1:B:1917:HIS:O	1:B:1921:GLU:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1764:LYS:HE3	2:B:206:HOH:O	2.17	0.45
1:A:1974:LYS:HZ2	1:A:1978:ASP:HB3	1.82	0.45
1:B:1898:VAL:HG12	1:B:1898:VAL:O	2.17	0.45
1:B:1820:ALA:O	1:B:1823:PRO:HD2	2.17	0.45
1:B:1883:ASN:HB3	1:B:1934:VAL:HG22	1.99	0.45
1:A:1900:SER:HA	2:A:160:HOH:O	2.16	0.45
1:A:1819:ALA:O	2:A:92:HOH:O	2.20	0.44
1:B:1909:ALA:O	1:B:1913:LEU:HG	2.17	0.44
1:B:1753:ARG:HG2	1:B:1757:MET:CE	2.47	0.44
1:B:1903:TYR:CE2	1:B:1976:LYS:HD3	2.53	0.44
1:B:1822:GLU:HG2	1:B:1823:PRO:HD3	2.00	0.44
1:B:1953:THR:HA	1:B:1956:MET:HE2	1.99	0.44
1:B:1754:ILE:HD13	2:B:222:HOH:O	2.18	0.44
1:B:1821:HIS:C	1:B:1823:PRO:CD	2.87	0.44
1:B:1903:TYR:CB	2:B:336:HOH:O	2.64	0.44
1:B:1700:ASN:O	1:B:1703[A]:LYS:HB2	2.18	0.44
1:A:1894:LYS:HB3	1:A:1920:PHE:CE1	2.53	0.44
1:B:1768:SER:CB	1:B:1841:ILE:HD11	2.48	0.43
1:A:1716:GLU:O	1:A:1720:LYS:HG3	2.17	0.43
1:B:1797:TYR:O	1:B:1804:VAL:HG12	2.18	0.43
1:A:1903:TYR:OH	1:A:1976:LYS:CD	2.67	0.43
1:A:1681:GLU:OE2	1:B:1916:LYS:HE3	2.18	0.43
1:B:1800:ASP:OD1	1:B:1803:GLY:N	2.47	0.43
1:A:1662:ALA:N	1:A:1666:GLN:HG3	2.34	0.43
1:B:1665:GLN:N	2:B:385:HOH:O	2.51	0.43
1:A:1663:ASN:O	1:A:1667:ASN:HB2	2.19	0.43
1:A:1962:LYS:HE3	2:A:429:HOH:O	2.18	0.43
1:A:1782:GLU:O	1:A:1786:LYS:HG3	2.19	0.43
1:A:1664:LYS:HD3	1:A:1729:LEU:CD2	2.49	0.42
1:A:1695:LEU:O	1:A:1699:ASN:ND2	2.51	0.42
1:A:1776:ARG:HG2	2:A:171:HOH:O	2.18	0.42
1:B:1911:GLN:HA	1:B:1911:GLN:OE1	2.20	0.42
1:B:1665:GLN:O	1:B:1669:ASN:ND2	2.52	0.42
1:B:1814:LEU:HD23	1:B:1814:LEU:C	2.40	0.42
1:B:1924:PHE:CD1	1:B:1928:LYS:HD2	2.55	0.42
1:B:1963:VAL:O	1:B:1967:GLU:HG3	2.20	0.42
1:A:1928:LYS:NZ	2:A:74:HOH:O	2.48	0.42
1:B:1695:LEU:O	1:B:1698:VAL:HG12	2.19	0.42
1:A:1736:ASP:OD1	1:A:1738:SER:HB3	2.18	0.42
1:A:1838:ASP:O	1:A:1840:THR:N	2.53	0.42
1:A:1719:LEU:HA	1:A:1719:LEU:HD12	1.73	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1924:PHE:O	1:B:1928:LYS:HG2	2.20	0.41
1:B:1670:THR:O	1:B:1673:LYS:HB3	2.20	0.41
1:A:1941:LEU:HB2	1:A:1952:ILE:HD13	2.02	0.41
1:B:1949:VAL:HA	1:B:1952:ILE:CG2	2.45	0.41
1:A:1722:LEU:HD23	1:A:1725:GLN:NE2	2.34	0.41
1:A:1847:GLN:NE2	1:A:1847:GLN:CA	2.76	0.41
1:A:1829:LEU:HD21	1:A:1853:PHE:HD2	1.85	0.41
1:A:1718[B]:ARG:HH12	1:A:1725:GLN:HE22	1.68	0.41
1:B:1914:LEU:HD21	1:B:1977:LEU:HB3	2.03	0.41
1:B:1881:VAL:O	1:B:1885:GLU:HG3	2.20	0.41
1:B:1840:THR:O	1:B:1841:ILE:C	2.59	0.41
1:A:1794:SER:HB2	1:B:1790:LEU:CD1	2.50	0.41
1:B:1734:ALA:HB1	2:B:273:HOH:O	2.19	0.41
1:A:1833[A]:LYS:HE3	2:A:106:HOH:O	2.20	0.41
1:B:1954:ALA:O	1:B:1957:LYS:HB3	2.19	0.41
1:A:1942:ILE:HD11	1:A:1952:ILE:CB	2.48	0.41
1:A:1903:TYR:HA	1:A:1913:LEU:CD1	2.50	0.41
1:A:1781:GLU:OE1	1:A:1781:GLU:HA	2.21	0.41
1:B:1945:ASN:HB3	2:B:178:HOH:O	2.21	0.41
1:B:1755:LYS:HE3	2:B:161:HOH:O	2.19	0.41
1:B:1880:PHE:O	1:B:1883:ASN:HB2	2.20	0.40
1:A:1794:SER:CB	1:B:1790:LEU:HD11	2.51	0.40
1:B:1727:ASP:OD1	1:B:1744:ARG:NH1	2.54	0.40
1:A:1773:GLN:HG2	2:A:144:HOH:O	2.22	0.40
1:A:1702:LEU:HD22	1:A:1762[B]:ARG:HH11	1.86	0.40
1:A:1804:VAL:HG11	1:A:1875:LEU:N	2.36	0.40
1:B:1928:LYS:O	1:B:1932:ASN:ND2	2.54	0.40
1:A:1829:LEU:CD2	1:A:1853:PHE:CD2	3.04	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	321/322 (100%)	296 (92%)	20 (6%)	5 (2%)	12	21
1	B	319/322 (99%)	295 (92%)	18 (6%)	6 (2%)	10	16
All	All	640/644 (99%)	591 (92%)	38 (6%)	11 (2%)	11	19

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1738	SER
1	A	1839	ASN
1	A	1841	ILE
1	B	1944	LYS
1	A	1902	ASP
1	B	1952	ILE
1	B	1957	LYS
1	B	1839	ASN
1	A	1798	GLY
1	B	1842	GLY
1	B	1841	ILE

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	277/275 (101%)	265 (96%)	12 (4%)	35	61
1	B	276/275 (100%)	264 (96%)	12 (4%)	35	61
All	All	553/550 (100%)	529 (96%)	24 (4%)	36	61

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1666	GLN
1	A	1667	ASN
1	A	1689	GLU
1	A	1739	GLN
1	A	1764	LYS

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Mol	Chain	Res	Type
1	A	1776	ARG
1	A	1799	ARG
1	A	1812	LYS
1	A	1847	GLN
1	A	1869	GLN
1	A	1923	ASP
1	A	1966	LEU
1	B	1667	ASN
1	B	1676	ASP
1	B	1707	LEU
1	B	1779	ASP
1	B	1848	GLN
1	B	1849[A]	ARG
1	B	1849[B]	ARG
1	B	1896	THR
1	B	1905	ASP
1	B	1924	PHE
1	B	1972	GLN
1	B	1973	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1665	GLN
1	A	1666	GLN
1	A	1725	GLN
1	A	1739	GLN
1	A	1805	GLN
1	A	1806	ASN
1	A	1839	ASN
1	A	1869	GLN
1	A	1879	GLN
1	A	1883	ASN
1	A	1892	ASN
1	A	1937	ASN
1	A	1945	ASN
1	A	1947	HIS
1	A	1948	HIS
1	A	1951	ASN
1	B	1667	ASN
1	B	1739	GLN
1	B	1805	GLN

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Mol	Chain	Res	Type
1	B	1806	ASN
1	B	1821	HIS
1	B	1852	GLN
1	B	1869	GLN
1	B	1883	ASN
1	B	1892	ASN
1	B	1932	ASN
1	B	1945	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 carbohydrates 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 [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	318/322 (98%)	0.06	11 (3%) 48 53	28, 53, 99, 122	0
1	B	317/322 (98%)	0.10	8 (2%) 61 65	30, 60, 93, 113	0
All	All	635/644 (98%)	0.08	19 (2%) 54 59	28, 55, 95, 122	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1977	LEU	9.6
1	A	1912	GLY	6.5
1	A	1909	ALA	4.3
1	B	1952	ILE	4.2
1	B	1949	VAL	4.2
1	A	1903	TYR	4.1
1	A	1949	VAL	3.7
1	B	1840	THR	3.5
1	A	1913	LEU	2.8
1	A	1978	ASP	2.7
1	A	1908	ALA	2.6
1	B	1820	ALA	2.6
1	A	1906	THR	2.3
1	A	1901	GLU	2.3
1	B	1951	ASN	2.3
1	B	1953	THR	2.2
1	A	1842	GLY	2.1
1	B	1950	GLU	2.1
1	B	1945	ASN	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 carbohydrates 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.