Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 07:31 AM EDT

PDB ID	:	5M5P
Title	:	S. cerevisiae spliceosomal helicase Brr2 (271-end) in complex with the Jab/MPN domain of S. cerevisiae Prp8
Authors	:	Becke, C.; Absmeier, E.; Wollenhaupt, J.; Santos, K.F.; Wahl, M.C.
Deposited on	:	2016-10-22
Resolution	:	4.20 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 (i) 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 (1)) were used in the production of this report:

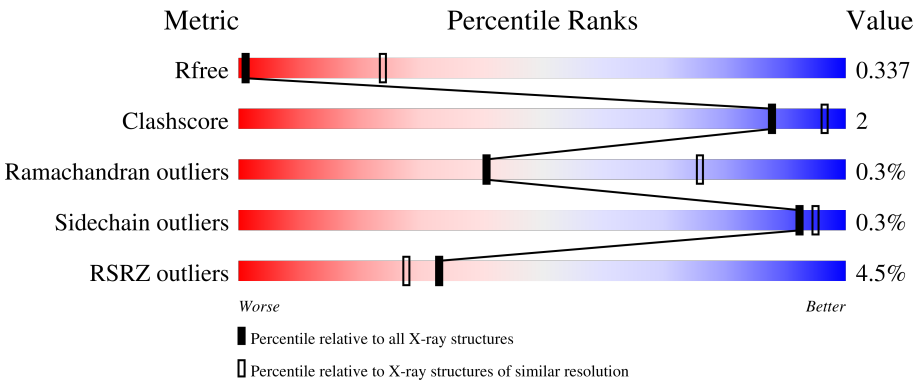
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 4.20 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	1005 (4.62-3.78)
Clashscore	141614	1044 (4.60-3.80)
Ramachandran outliers	138981	1000 (4.60-3.80)
Sidechain outliers	138945	1007 (4.62-3.78)
RSRZ outliers	127900	1063 (4.70-3.70)

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	1897	3% 92% 5% •
1	C	1897	5% 92% 5% •
2	B	270	8% 89% • 8%
2	D	270	8% 85% 7% 8%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 67309 atoms, of which 33658 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 Pre-mRNA-splicing helicase BRR2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	1850	29711	9505	14876	2465	2808	57	0	0	0
1	C	1850	29716	9505	14878	2467	2809	57	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	267	GLY	-	expression tag	UNP P32639
A	268	ALA	-	expression tag	UNP P32639
A	269	GLU	-	expression tag	UNP P32639
A	270	PHE	-	expression tag	UNP P32639
C	267	GLY	-	expression tag	UNP P32639
C	268	ALA	-	expression tag	UNP P32639
C	269	GLU	-	expression tag	UNP P32639
C	270	PHE	-	expression tag	UNP P32639

- Molecule 2 is a protein called Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	248	3941	1284	1952	322	377	6	0	0	0
2	D	248	3941	1284	1952	322	377	6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	2144	GLY	-	expression tag	UNP P33334
B	2145	ALA	-	expression tag	UNP P33334
B	2146	MET	-	expression tag	UNP P33334
D	2144	GLY	-	expression tag	UNP P33334

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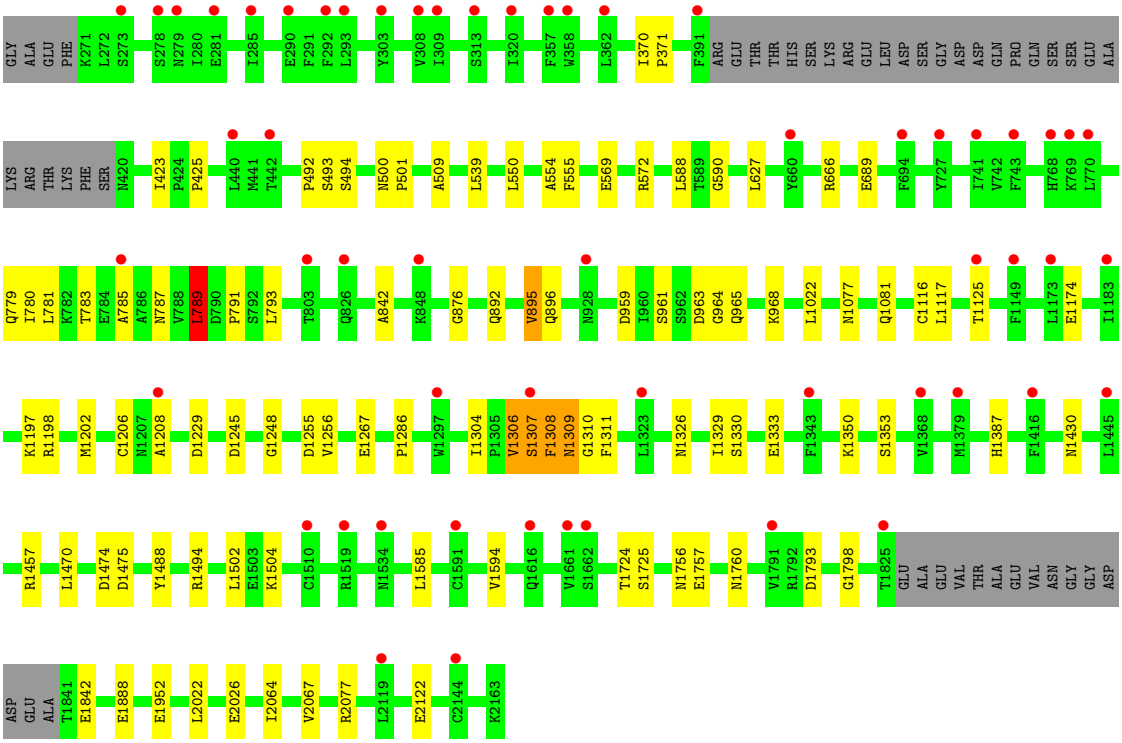
Chain	Residue	Modelled	Actual	Comment	Reference
D	2145	ALA	-	expression tag	UNP P33334
D	2146	MET	-	expression tag	UNP P33334

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.

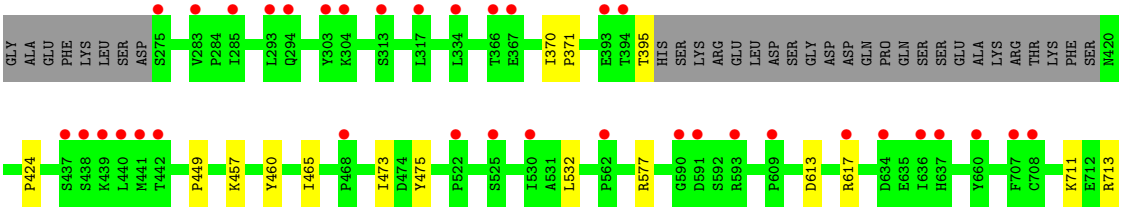
• Molecule 1: Pre-mRNA-splicing helicase BRR2

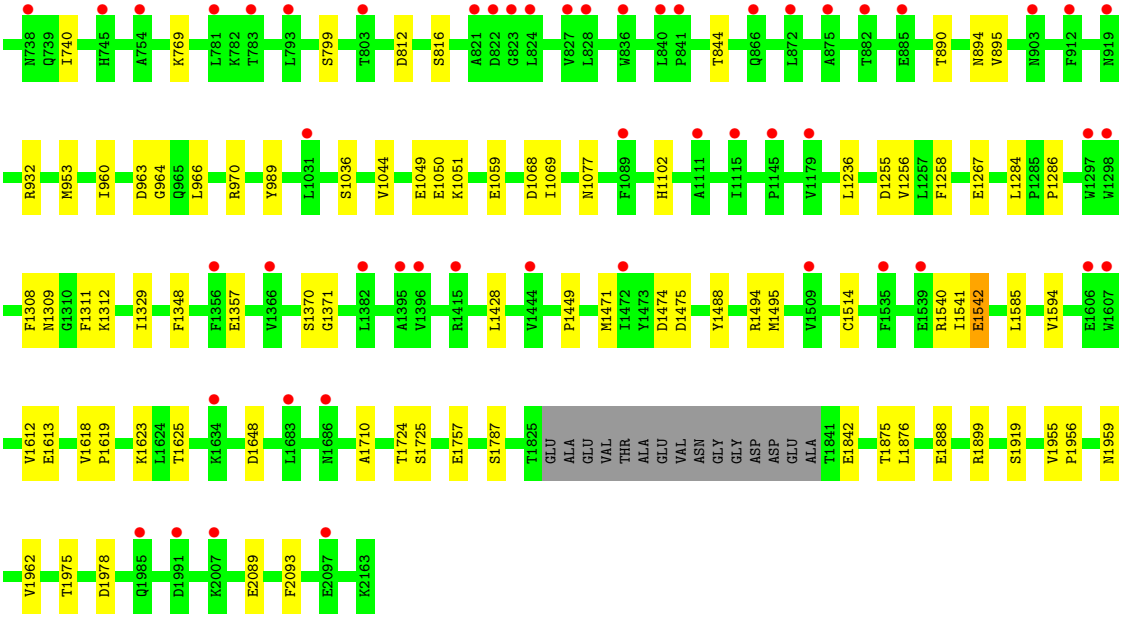
Chain A: 3% 92% 5%



• Molecule 1: Pre-mRNA-splicing helicase BRR2

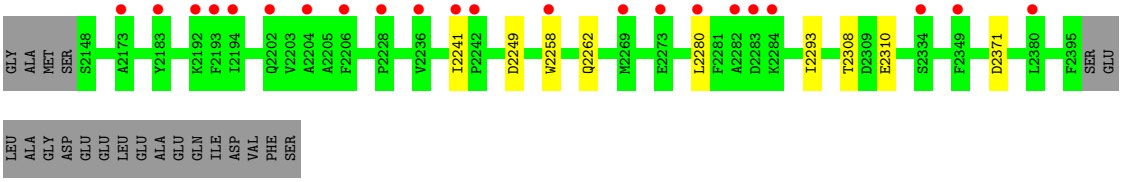
Chain C: 5% 92% 5%





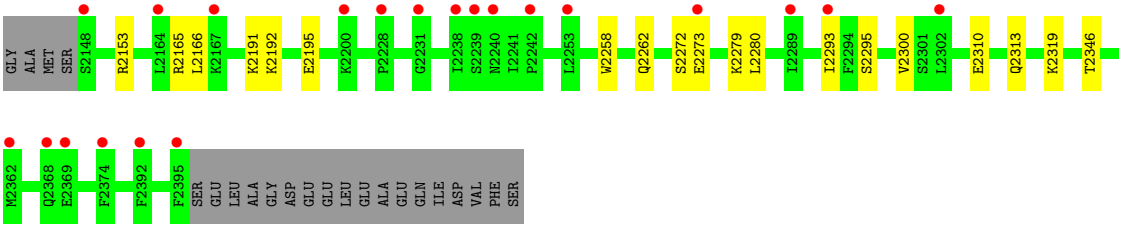
• Molecule 2: Pre-mRNA-splicing factor 8

Chain B:



• Molecule 2: Pre-mRNA-splicing factor 8

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	185.97Å 196.87Å 191.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	95.50 – 4.20 95.52 – 4.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (95.50-4.20) 99.9 (95.52-4.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.32	Depositor
$< I/\sigma(I) >$ [1]	1.61 (at 4.15Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.307 , 0.335 0.311 , 0.337	Depositor DCC
R_{free} test set	2667 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å^2)	147.0	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 94.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	0.038 for l,-k,h	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	67309	wwPDB-VP
Average B, all atoms (Å^2)	157.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% 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 (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	0.23	0/15149	0.43	1/20529 (0.0%)
1	C	0.24	0/15152	0.42	0/20534
2	B	0.22	0/2038	0.40	0/2764
2	D	0.23	0/2038	0.41	0/2764
All	All	0.23	0/34377	0.42	1/46591 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	789	LEU	CA-CB-CG	5.17	127.20	115.30

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	14835	14876	14874	65	0
1	C	14838	14878	14875	52	1
2	B	1989	1952	1951	6	0
2	D	1989	1952	1951	13	0
All	All	33651	33658	33651	131	1

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 2.

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:1208:ALA:CB	1:A:1307:SER:OG	1.67	1.38
1:A:1208:ALA:HB3	1:A:1307:SER:OG	1.02	1.17
1:A:1286:PRO:O	1:A:1308:PHE:CD1	2.06	0.99
1:A:1309:ASN:OD1	2:B:2249:ASP:OD1	1.81	0.96
1:A:1208:ALA:HB2	1:A:1307:SER:OG	1.72	0.88
1:A:1307:SER:HB3	1:A:1311:PHE:CE2	2.13	0.82
1:C:1540:ARG:NH2	1:C:1710:ALA:O	2.13	0.82
1:A:963:ASP:O	1:A:965:GLN:N	2.11	0.82
1:A:1206:CYS:SG	1:A:1306:VAL:HG13	2.23	0.79
1:A:1286:PRO:O	1:A:1308:PHE:CG	2.36	0.78
1:A:1308:PHE:O	1:A:1310:GLY:N	2.16	0.78
1:A:787:ASN:OD1	1:A:789:LEU:N	2.20	0.74
1:A:1307:SER:HB3	1:A:1311:PHE:HE2	1.54	0.72
1:A:1077:ASN:OD1	1:A:1081:GLN:NE2	2.25	0.69
2:D:2165:ARG:NH1	2:D:2295:SER:O	2.25	0.69
1:C:1059:GLU:O	2:D:2153:ARG:NH2	2.26	0.68
1:C:769:LYS:NZ	1:C:799:SER:O	2.27	0.67
1:A:1174:GLU:OE1	1:A:1197:LYS:NZ	2.31	0.64
1:C:963:ASP:O	1:C:966:LEU:N	2.27	0.64
1:A:1307:SER:CB	1:A:1311:PHE:CE2	2.80	0.63
1:A:959:ASP:OD1	1:A:961:SER:OG	2.16	0.63
1:C:1625:THR:OG1	1:C:1648:ASP:OD1	2.15	0.63
1:C:890:THR:OG1	1:C:894:ASN:OD1	2.17	0.61
1:C:1050:GLU:N	1:C:1050:GLU:OE1	2.33	0.61
1:A:1307:SER:CB	1:A:1311:PHE:HE2	2.15	0.60
2:D:2273:GLU:N	2:D:2273:GLU:OE1	2.35	0.59
1:A:779:GLN:O	1:A:783:THR:N	2.35	0.58
1:A:1494:ARG:NH1	1:A:1757:GLU:OE2	2.37	0.58
1:C:395:THR:O	1:C:395:THR:HG22	2.04	0.57
1:A:492:PRO:O	1:A:494:SER:N	2.37	0.57
1:C:2089:GLU:N	1:C:2089:GLU:OE1	2.37	0.56
1:A:1888:GLU:OE1	1:A:1888:GLU:N	2.38	0.56
1:A:1330:SER:OG	1:A:1333:GLU:OE1	2.19	0.55
1:A:1326:ASN:ND2	1:A:1350:LYS:O	2.39	0.55
1:A:1286:PRO:O	1:A:1308:PHE:HA	2.07	0.55
1:C:1888:GLU:N	1:C:1888:GLU:OE1	2.38	0.55
1:A:1198:ARG:NH1	1:A:1229:ASP:OD2	2.40	0.54
1:C:960:ILE:O	1:C:964:GLY:N	2.40	0.54
1:A:1309:ASN:OD1	2:B:2249:ASP:CG	2.46	0.54
1:C:1267:GLU:N	1:C:1267:GLU:OE1	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1036:SER:OG	1:C:1077:ASN:OD1	2.27	0.53
1:A:370:ILE:HB	1:A:371:PRO:HD3	1.92	0.52
2:D:2272:SER:OG	2:D:2273:GLU:OE1	2.25	0.52
1:A:781:LEU:HA	1:A:785:ALA:HB3	1.90	0.52
2:D:2310:GLU:OE1	2:D:2313:GLN:NE2	2.43	0.52
1:A:689:GLU:OE1	1:A:689:GLU:N	2.44	0.51
1:A:1842:GLU:OE1	1:A:1842:GLU:N	2.42	0.51
1:C:1312:LYS:N	1:C:1787:SER:OG	2.36	0.51
2:B:2371:ASP:OD2	2:D:2319:LYS:NZ	2.42	0.51
1:C:960:ILE:HA	1:C:963:ASP:HB3	1.93	0.51
1:C:1284:LEU:HD11	1:C:1308:PHE:HB3	1.94	0.50
1:A:539:LEU:HD22	1:A:550:LEU:HD21	1.94	0.50
1:A:892:GLN:HA	1:A:895:VAL:HG23	1.94	0.49
1:C:966:LEU:O	1:C:970:ARG:NE	2.39	0.49
1:A:1308:PHE:C	1:A:1310:GLY:H	2.12	0.49
1:A:1255:ASP:OD1	1:A:1256:VAL:N	2.44	0.49
1:C:1449:PRO:HG3	1:C:1488:TYR:HD2	1.78	0.48
1:C:370:ILE:HB	1:C:371:PRO:HD3	1.96	0.47
1:C:1068:ASP:OD1	1:C:1069:ILE:N	2.47	0.47
2:D:2165:ARG:HB3	2:D:2300:VAL:HG13	1.97	0.47
1:C:1474:ASP:OD1	1:C:1475:ASP:N	2.47	0.47
1:C:740:ILE:HG22	1:C:844:THR:HB	1.96	0.46
1:A:780:ILE:O	1:A:785:ALA:N	2.46	0.46
1:A:1267:GLU:OE1	1:A:1267:GLU:N	2.44	0.46
1:C:532:LEU:HD22	1:C:577:ARG:HD3	1.98	0.46
1:A:1022:LEU:HD21	1:A:1116:CYS:SG	2.56	0.45
1:C:1724:THR:OG1	1:C:1725:SER:N	2.49	0.45
1:A:1474:ASP:OD1	1:A:1475:ASP:N	2.49	0.45
1:A:1724:THR:OG1	1:A:1725:SER:N	2.48	0.45
1:C:1370:SER:O	1:C:1514:CYS:N	2.49	0.45
1:A:1502:LEU:O	1:A:1504:LYS:N	2.43	0.45
1:C:1255:ASP:OD1	1:C:1256:VAL:N	2.47	0.45
1:C:1875:THR:OG1	1:C:1876:LEU:N	2.48	0.45
2:D:2262:GLN:N	2:D:2293:ILE:O	2.50	0.45
1:C:1612:VAL:HG22	1:C:1613:GLU:N	2.31	0.44
1:C:473:ILE:HG12	1:C:475:TYR:H	1.82	0.44
1:C:932:ARG:NH2	1:C:989:TYR:OH	2.50	0.44
1:A:842:ALA:O	1:A:876:GLY:N	2.49	0.44
1:A:509:ALA:O	1:A:666:ARG:NH2	2.50	0.44
1:A:1202:MET:SD	1:A:1202:MET:N	2.90	0.44
1:C:812:ASP:O	1:C:816:SER:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1585:LEU:HD21	1:C:1594:VAL:HG21	1.99	0.44
1:C:1623:LYS:O	1:C:1625:THR:N	2.48	0.44
1:A:588:LEU:O	1:A:590:GLY:N	2.51	0.44
1:A:787:ASN:ND2	1:A:789:LEU:HG	2.32	0.44
1:C:1049:GLU:OE1	1:C:1049:GLU:N	2.45	0.44
1:A:1245:ASP:OD1	1:A:1248:GLY:N	2.47	0.43
1:C:1311:PHE:HD1	1:C:1787:SER:HB3	1.83	0.43
1:C:1541:ILE:HG23	1:C:1542:GLU:HG2	2.01	0.43
1:C:1975:THR:O	1:C:1978:ASP:N	2.52	0.43
2:B:2308:THR:OG1	2:B:2310:GLU:OE1	2.23	0.43
1:A:539:LEU:HD23	1:A:555:PHE:CE2	2.54	0.43
2:D:2279:LYS:HE3	2:D:2319:LYS:HA	2.01	0.43
1:C:1494:ARG:NH1	1:C:1757:GLU:OE2	2.51	0.42
1:C:1842:GLU:OE1	1:C:1842:GLU:N	2.43	0.42
1:C:1899:ARG:NH2	1:C:1919:SER:OG	2.47	0.42
1:A:423:ILE:O	1:A:425:PRO:HD3	2.19	0.42
1:A:569:GLU:OE1	1:A:572:ARG:NH2	2.47	0.42
1:C:711:LYS:O	1:C:713:ARG:NH1	2.52	0.42
1:A:500:ASN:HB2	1:A:501:PRO:HD2	2.02	0.42
1:A:1326:ASN:ND2	1:A:1353:SER:OG	2.52	0.42
2:D:2192:LYS:HA	2:D:2195:GLU:HG2	2.01	0.42
1:A:789:LEU:HD11	1:A:793:LEU:HD12	2.01	0.42
1:A:1793:ASP:HB2	1:A:1798:GLY:HA3	2.02	0.42
1:C:457:LYS:HD2	1:C:460:TYR:CE1	2.55	0.42
2:B:2241:ILE:HG13	2:B:2280:LEU:HD21	2.02	0.42
1:C:449:PRO:HG2	1:C:465:ILE:HG23	2.00	0.42
1:A:1457:ARG:O	1:A:1760:ASN:ND2	2.42	0.41
1:C:1370:SER:OG	1:C:1371:GLY:N	2.53	0.41
1:A:1387:HIS:HB2	1:A:1470:LEU:HD22	2.01	0.41
1:A:1585:LEU:HD21	1:A:1594:VAL:HG21	2.01	0.41
1:A:1117:LEU:HA	1:A:1125:THR:HG21	2.03	0.41
1:C:1286:PRO:HD3	2:D:2346:THR:HB	2.02	0.41
1:A:554:ALA:HA	1:A:627:LEU:HD22	2.03	0.41
1:A:895:VAL:HG12	1:A:896:GLN:N	2.36	0.41
2:B:2262:GLN:N	2:B:2293:ILE:O	2.53	0.41
1:C:613:ASP:OD2	1:C:617:ARG:NH2	2.54	0.41
1:C:1236:LEU:HD11	1:C:1258:PHE:HB3	2.02	0.41
1:C:1428:LEU:HD21	1:C:2093:PHE:HB3	2.02	0.41
1:C:1471:MET:SD	1:C:1495:MET:HG3	2.61	0.41
1:C:1618:VAL:HB	1:C:1619:PRO:HD3	2.03	0.41
1:C:1959:ASN:HA	1:C:1962:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:2295:SER:HA	2:D:2300:VAL:HG12	2.03	0.41
1:A:2077:ARG:N	1:A:2122:GLU:O	2.46	0.41
1:A:963:ASP:OD2	1:A:968:LYS:N	2.54	0.40
1:A:2022:LEU:HG	1:A:2026:GLU:HB2	2.02	0.40
2:D:2166:LEU:HD13	2:D:2191:LYS:HA	2.02	0.40
1:C:1955:VAL:HB	1:C:1956:PRO:HD3	2.03	0.40
1:A:1494:ARG:NH2	1:A:1756:ASN:OD1	2.54	0.40
1:A:2064:ILE:HB	1:A:2067:VAL:HB	2.03	0.40
1:A:1430:ASN:ND2	1:A:1952:GLU:O	2.54	0.40

All (1) 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:C:1051:LYS:NZ	1:C:1357:GLU:OE1[3_444]	2.17	0.03

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	1844/1897 (97%)	1704 (92%)	132 (7%)	8 (0%)	34	72
1	C	1844/1897 (97%)	1698 (92%)	141 (8%)	5 (0%)	41	76
2	B	246/270 (91%)	229 (93%)	17 (7%)	0	100	100
2	D	246/270 (91%)	225 (92%)	20 (8%)	1 (0%)	34	72
All	All	4180/4334 (96%)	3856 (92%)	310 (7%)	14 (0%)	41	76

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	493	SER

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Mol	Chain	Res	Type
1	A	895	VAL
1	A	964	GLY
1	A	1307	SER
1	A	1308	PHE
1	A	1309	ASN
1	C	895	VAL
1	C	1309	ASN
1	C	1329	ILE
1	A	1329	ILE
2	D	2280	LEU
1	C	424	PRO
1	A	791	PRO
1	C	1044	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	1671/1709 (98%)	1667 (100%)	4 (0%)	93 96
1	C	1671/1709 (98%)	1667 (100%)	4 (0%)	93 96
2	B	220/237 (93%)	219 (100%)	1 (0%)	88 93
2	D	220/237 (93%)	219 (100%)	1 (0%)	88 93
All	All	3782/3892 (97%)	3772 (100%)	10 (0%)	92 95

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	789	LEU
1	A	1304	ILE
1	A	1306	VAL
1	A	1488	TYR
2	B	2258	TRP
1	C	953	MET
1	C	1102	HIS
1	C	1348	PHE

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Mol	Chain	Res	Type
1	C	1542	GLU
2	D	2258	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

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	1850/1897 (97%)	0.40	56 (3%)	50	39	141, 156, 170, 178	0
1	C	1850/1897 (97%)	0.44	88 (4%)	30	26	143, 158, 169, 180	0
2	B	248/270 (91%)	0.64	22 (8%)	9	8	148, 159, 168, 175	0
2	D	248/270 (91%)	0.62	21 (8%)	10	10	148, 160, 169, 177	0
All	All	4196/4334 (96%)	0.45	187 (4%)	33	27	141, 157, 169, 180	0

All (187) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	440	LEU	8.1
1	C	438	SER	5.7
1	C	437	SER	4.9
1	A	1825	THR	4.8
2	D	2239	SER	4.6
1	A	1662	SER	4.2
1	A	303	TYR	4.2
1	A	442	THR	4.1
1	A	1616	GLN	4.1
1	A	769	LYS	4.1
1	C	394	THR	3.9
1	C	525	SER	3.9
1	A	768	HIS	3.9
1	C	590	GLY	3.9
1	C	821	ALA	3.9
1	C	439	LYS	3.8
1	A	362	LEU	3.7
1	C	875	ALA	3.7
1	A	440	LEU	3.7
1	C	442	THR	3.7
1	C	1145	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	1179	VAL	3.6
1	C	823	GLY	3.6
1	C	1535	PHE	3.6
1	C	1356	PHE	3.5
1	A	660	TYR	3.4
2	B	2258	TRP	3.3
1	C	841	PRO	3.3
1	C	1382	LEU	3.3
1	A	285	ILE	3.2
2	B	2334	SER	3.2
1	C	840	LEU	3.2
1	C	636	ILE	3.2
2	B	2280	LEU	3.2
1	A	1173	LEU	3.2
1	C	317	LEU	3.1
1	C	1509	VAL	3.1
2	D	2200	LYS	3.1
1	A	278	SER	3.1
2	D	2395	PHE	3.1
2	D	2362	MET	3.0
2	B	2284	LYS	3.0
1	C	637	HIS	3.0
1	A	273	SER	3.0
2	D	2369	GLU	3.0
1	C	609	PRO	2.9
1	A	826	GLN	2.9
1	A	308	VAL	2.9
1	C	1395	ALA	2.9
1	C	2097	GLU	2.9
2	D	2368	GLN	2.9
1	C	1607	TRP	2.9
2	D	2164	LEU	2.9
1	C	303	TYR	2.9
1	A	391	PHE	2.9
1	C	275	SER	2.9
1	C	1991	ASP	2.9
1	C	522	PRO	2.9
2	B	2242	PRO	2.9
2	B	2204	ALA	2.9
2	B	2192	LYS	2.8
1	C	1366	VAL	2.8
1	A	1661	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	281	GLU	2.8
1	A	785	ALA	2.8
1	C	745	HIS	2.8
1	C	468	PRO	2.7
1	A	1307	SER	2.7
2	D	2240	ASN	2.7
2	B	2269	MET	2.7
1	C	1115	ILE	2.7
2	D	2242	PRO	2.7
1	C	783	THR	2.7
1	A	1125	THR	2.7
2	B	2236	VAL	2.7
1	C	660	TYR	2.6
1	C	1683	LEU	2.6
1	A	290	GLU	2.6
2	B	2282	ALA	2.6
2	D	2148	SER	2.6
1	C	285	ILE	2.6
1	C	530	ILE	2.6
1	C	919	ASN	2.6
1	C	803	THR	2.6
1	A	694	PHE	2.6
2	D	2167	LYS	2.6
1	C	1297	TRP	2.6
1	A	727	TYR	2.5
1	C	707	PHE	2.5
1	A	357	PHE	2.5
1	C	754	ALA	2.5
1	A	1510	CYS	2.5
1	C	822	ASP	2.5
1	C	1539	GLU	2.5
2	B	2273	GLU	2.5
1	A	279	ASN	2.5
1	C	294	GLN	2.5
1	A	1591	CYS	2.5
1	C	824	LEU	2.5
1	C	1985	GLN	2.5
1	C	313	SER	2.5
1	C	781	LEU	2.5
1	A	848	LYS	2.5
1	C	885	GLU	2.5
1	A	770	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	1472	ILE	2.4
1	A	1519	ARG	2.4
1	C	866	GLN	2.4
1	C	283	VAL	2.4
1	C	708	CYS	2.4
1	C	827	VAL	2.4
1	C	836	TRP	2.4
1	A	1297	TRP	2.4
1	A	803	THR	2.4
1	A	309	ILE	2.3
1	A	1208	ALA	2.3
1	A	292	PHE	2.3
1	A	1534	ASN	2.3
2	D	2238	ILE	2.3
1	A	1379	MET	2.3
1	C	366	THR	2.3
1	C	393	GLU	2.3
2	B	2241	ILE	2.3
2	D	2374	PHE	2.3
1	C	872	LEU	2.3
1	A	743	PHE	2.3
1	A	358	TRP	2.3
2	B	2183	TYR	2.3
2	D	2392	PHE	2.3
2	B	2349	PHE	2.2
1	C	593	ARG	2.2
1	A	2119	LEU	2.2
2	B	2202	GLN	2.2
1	A	1149	PHE	2.2
1	A	928	ASN	2.2
1	C	334	LEU	2.2
2	B	2193	PHE	2.2
1	C	591	ASP	2.2
1	C	912	PHE	2.2
1	C	882	THR	2.2
1	C	1396	VAL	2.2
1	A	1445	LEU	2.2
1	C	1634	LYS	2.2
1	A	313	SER	2.2
1	C	903	ASN	2.2
1	C	304	LYS	2.2
1	A	1323	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	1416	PHE	2.2
1	C	441	MET	2.2
1	A	741	ILE	2.2
1	C	1111	ALA	2.1
1	C	1298	TRP	2.1
2	D	2231	GLY	2.1
1	C	562	PRO	2.1
1	C	634	ASP	2.1
1	C	293	LEU	2.1
1	C	367	GLU	2.1
2	B	2194	ILE	2.1
2	D	2293	ILE	2.1
1	C	1031	LEU	2.1
1	A	1368	VAL	2.1
1	C	1089	PHE	2.1
2	B	2173	ALA	2.1
1	C	1415	ARG	2.1
1	C	1606	GLU	2.1
2	B	2380	LEU	2.1
2	D	2228	PRO	2.1
2	B	2206	PHE	2.1
1	A	1791	VAL	2.1
1	A	1183	ILE	2.1
1	C	617	ARG	2.1
2	B	2228	PRO	2.1
1	A	2144	CYS	2.0
1	C	1686	ASN	2.0
1	C	2007	LYS	2.0
2	B	2283	ASP	2.0
2	D	2253	LEU	2.0
1	A	293	LEU	2.0
2	D	2302	LEU	2.0
2	D	2273	GLU	2.0
1	A	1343	PHE	2.0
1	C	793	LEU	2.0
1	C	738	ASN	2.0
1	C	1444	VAL	2.0
1	A	320	ILE	2.0
2	D	2289	ILE	2.0
1	C	828	LEU	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.