



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:59 PM UTC

PDB ID : 5YVG / pdb_00005yvg
Title : Crystal structure of Karyopherin beta2 in complex with FUS(full length)
Authors : Yoshizawa, T.; Fung, H.Y.J.; Chook, Y.M.
Deposited on : 2017-11-25
Resolution : 4.05 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

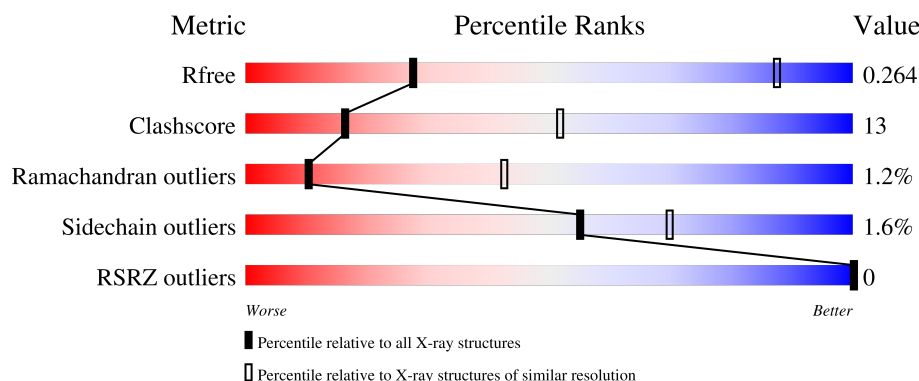
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

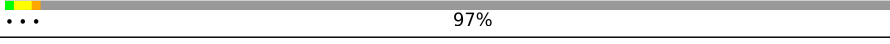
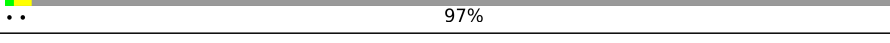
The reported resolution of this entry is 4.05 Å.

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	1183 (4.32-3.80)
Clashscore	190562	1231 (4.32-3.80)
Ramachandran outliers	187476	1154 (4.32-3.80)
Sidechain outliers	187428	1144 (4.32-3.80)
RSRZ outliers	180081	1182 (4.32-3.80)

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	868	 71% 24% .
1	B	868	 63% 23% . 12%
2	X	528	 ... 97%
2	Y	528	 .. 97%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12966 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 Transportin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	830	Total	C	N	O	S	0	0	0
			6620	4248	1100	1220	52			
1	B	764	Total	C	N	O	S	0	0	0
			6057	3904	1001	1104	48			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q92973
A	0	SER	-	expression tag	UNP Q92973
A	361	GLY	-	linker	UNP Q92973
A	362	GLY	-	linker	UNP Q92973
A	363	SER	-	linker	UNP Q92973
A	364	GLY	-	linker	UNP Q92973
A	365	GLY	-	linker	UNP Q92973
A	366	SER	-	linker	UNP Q92973
A	367	GLY	-	linker	UNP Q92973
B	-1	GLY	-	expression tag	UNP Q92973
B	0	SER	-	expression tag	UNP Q92973
B	361	GLY	-	linker	UNP Q92973
B	362	GLY	-	linker	UNP Q92973
B	363	SER	-	linker	UNP Q92973
B	364	GLY	-	linker	UNP Q92973
B	365	GLY	-	linker	UNP Q92973
B	366	SER	-	linker	UNP Q92973
B	367	GLY	-	linker	UNP Q92973

- Molecule 2 is a protein called RNA-binding protein FUS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	18	Total	C	N	O	S	0	0	0
			159	91	37	30	1			

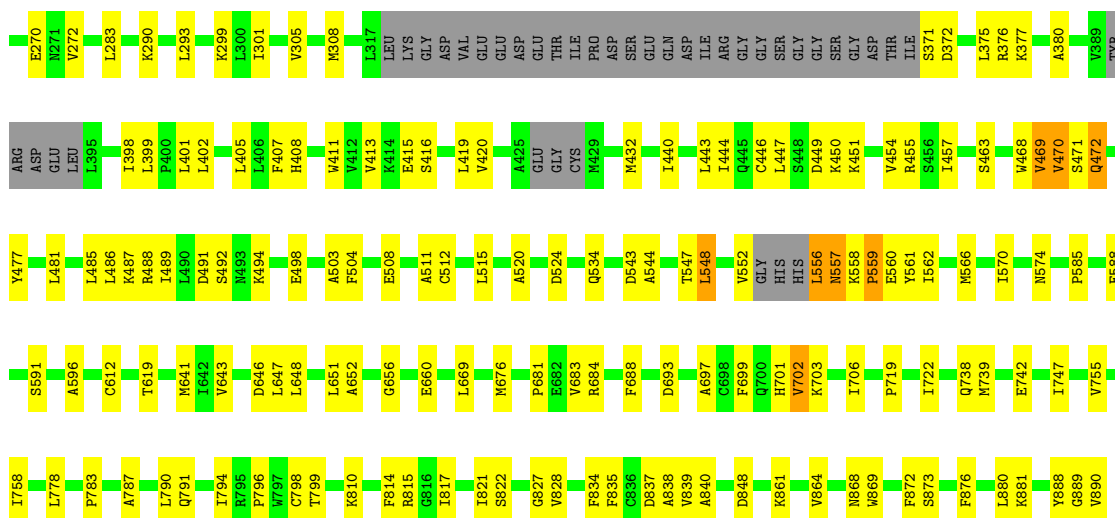
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	Y	14	Total	C	N	O	0	0	0
			130	74	32	24			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	-1	GLY	-	expression tag	UNP P35637
X	0	SER	-	expression tag	UNP P35637
Y	-1	GLY	-	expression tag	UNP P35637
Y	0	SER	-	expression tag	UNP P35637

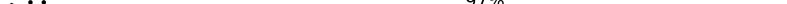


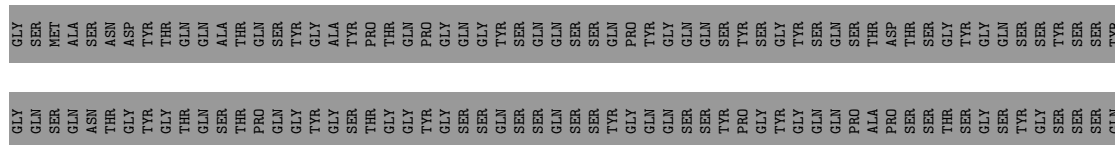
- Molecule 2: RNA-binding protein FUS

Chain X: ... 97%



- Molecule 2: RNA-binding protein FUS

Chain Y:  97%





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	108.22Å 145.01Å 108.68Å 90.00° 94.55° 90.00°	Depositor
Resolution (Å)	41.33 – 4.05 41.33 – 4.05	Depositor EDS
% Data completeness (in resolution range)	87.4 (41.33-4.05) 86.5 (41.33-4.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 4.00Å)	Xtriage
Refinement program	phenix.refine 1.13_2998, PHENIX 1.13_2998	Depositor
R, R_{free}	0.251 , 0.260 0.262 , 0.264	Depositor DCC
R_{free} test set	1189 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	38.6	Xtriage
Anisotropy	0.593	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.278 for l,-k,h	Xtriage
Reported twinning fraction	0.360 for l,-k,h	Depositor
Outliers	0 of 23952 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12966	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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

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	0.18	0/6760	0.38	0/9176
1	B	0.19	0/6174	0.41	1/8370 (0.0%)
2	X	0.23	0/161	0.55	0/209
2	Y	0.17	0/132	0.40	0/172
All	All	0.19	0/13227	0.40	1/17927 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	561	TYR	N-CA-C	-7.34	103.36	111.36

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	6620	0	6675	149	0
1	B	6057	0	6164	180	0
2	X	159	0	148	35	0
2	Y	130	0	119	7	0
All	All	12966	0	13106	340	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 13.

All (340) 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:B:103:SER:HB2	1:B:104:PRO:HD2	1.19	1.16
1:A:854:CYS:O	1:A:858:HIS:HB2	1.50	1.11
1:B:508:GLU:HG3	1:B:548:LEU:HD12	1.25	1.10
2:Y:516:GLU:O	2:Y:520:ASP:HB2	1.57	1.05
1:B:512:CYS:O	1:B:552:VAL:HG22	1.56	1.04
1:B:508:GLU:OE2	1:B:548:LEU:HD11	1.61	1.01
1:B:103:SER:HB2	1:B:104:PRO:CD	1.90	1.01
1:B:508:GLU:HG3	1:B:548:LEU:CD1	2.00	0.91
1:B:99:ILE:CD1	1:B:114:ILE:HD11	2.01	0.90
1:B:556:LEU:HD21	1:B:596:ALA:HB1	1.55	0.89
1:A:460:TRP:CH2	2:X:526:TYR:HB3	2.06	0.89
2:X:513:SER:O	2:X:517:HIS:HB2	1.73	0.89
1:B:556:LEU:HD11	1:B:596:ALA:CA	2.08	0.83
1:B:556:LEU:HD11	1:B:596:ALA:HA	1.60	0.82
1:B:511:ALA:CB	1:B:515:LEU:HD21	2.14	0.78
1:A:853:PHE:O	1:A:857:LEU:HB3	1.84	0.78
1:A:685:GLN:HE21	2:X:510:LYS:HE2	1.49	0.77
1:A:543:ASP:OD1	2:X:518:ARG:HB2	1.82	0.77
1:B:508:GLU:CG	1:B:548:LEU:HD12	2.11	0.77
1:B:415:GLU:HG2	1:B:457:ILE:HG21	1.64	0.77
1:B:99:ILE:CD1	1:B:114:ILE:CD1	2.63	0.76
1:B:105:LEU:HD12	1:B:105:LEU:O	1.86	0.74
1:A:464:ARG:NH2	2:X:526:TYR:CE1	2.55	0.74
1:B:511:ALA:HB1	1:B:515:LEU:HD21	1.69	0.73
1:B:99:ILE:HD12	1:B:114:ILE:HD11	1.70	0.73
1:A:145:CYS:O	1:A:149:PHE:HB2	1.88	0.73
1:B:377:LYS:HG3	2:Y:525:PRO:O	1.87	0.72
1:B:66:SER:OG	1:B:105:LEU:HD11	1.89	0.71
1:B:65:ARG:HE	1:B:106:ILE:HD11	1.55	0.71
1:A:458:THR:O	1:A:462:LEU:HB2	1.90	0.71
1:B:652:ALA:HB3	1:B:693:ASP:HB3	1.74	0.70
1:B:556:LEU:O	1:B:557:ASN:HB2	1.90	0.70
1:A:869:TRP:O	1:A:873:SER:HB2	1.92	0.69
1:B:376:ARG:NH2	1:B:415:GLU:OE2	2.24	0.69
1:B:641:MET:HE3	1:B:683:VAL:HG21	1.75	0.68
1:A:460:TRP:CE2	2:X:524:ARG:HB2	2.28	0.68
2:X:510:LYS:O	2:X:512:ASP:N	2.27	0.68
1:A:495:ARG:HA	1:A:498:GLU:HG2	1.76	0.67
1:B:556:LEU:HD13	1:B:596:ALA:O	1.95	0.67
1:A:423:ALA:HA	2:X:526:TYR:CD2	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:SER:CB	1:B:104:PRO:CD	2.67	0.66
1:B:270:GLU:HG3	1:B:375:LEU:HD12	1.75	0.66
1:B:790:LEU:HD21	1:B:821:ILE:HG22	1.78	0.66
1:A:760:ARG:NH2	1:A:762:ASN:O	2.28	0.66
1:B:556:LEU:CD1	1:B:596:ALA:O	2.44	0.66
1:A:554:HIS:ND1	1:A:598:GLN:OE1	2.27	0.66
1:B:701:HIS:O	1:B:703:LYS:N	2.23	0.66
1:A:18:LEU:HD23	1:A:37:LEU:HD13	1.79	0.65
1:A:543:ASP:CG	2:X:518:ARG:HB2	2.22	0.65
1:A:685:GLN:NE2	2:X:510:LYS:HE2	2.12	0.65
1:B:512:CYS:O	1:B:552:VAL:CG2	2.41	0.65
1:B:556:LEU:C	1:B:556:LEU:HD22	2.22	0.65
1:A:869:TRP:O	1:A:873:SER:CB	2.45	0.64
1:A:703:LYS:NZ	1:A:739:MET:SD	2.71	0.64
1:B:796:PRO:HA	1:B:799:THR:HB	1.80	0.64
1:A:213:ILE:HA	1:A:216:PHE:HB3	1.80	0.64
1:B:214:ASP:O	1:B:218:GLU:CB	2.46	0.64
1:B:224:ALA:O	1:B:264:ARG:NH2	2.30	0.64
1:B:398:ILE:HG13	1:B:401:LEU:HD13	1.80	0.63
1:B:559:PRO:O	1:B:560:GLU:OE2	2.17	0.63
1:A:509:GLU:CD	2:X:522:ARG:HH12	2.05	0.63
1:A:853:PHE:O	1:A:857:LEU:CB	2.46	0.63
1:B:126:TRP:HB3	1:B:127:PRO:HD3	1.81	0.63
1:B:504:PHE:O	1:B:508:GLU:HG2	1.99	0.63
1:B:515:LEU:HD22	1:B:515:LEU:N	2.14	0.62
1:B:238:ALA:O	1:B:242:LEU:HB2	1.98	0.62
1:B:233:LYS:HB2	1:B:272:VAL:HG22	1.81	0.62
1:A:185:HIS:O	1:A:191:ARG:NH1	2.33	0.62
1:A:460:TRP:NE1	2:X:524:ARG:CB	2.63	0.61
1:A:460:TRP:NE1	2:X:524:ARG:HB2	2.15	0.61
1:A:293:LEU:HG	1:A:297:LEU:HD21	1.83	0.61
1:A:464:ARG:NH2	2:X:526:TYR:HE1	1.98	0.60
1:A:652:ALA:HB3	1:A:693:ASP:HB3	1.82	0.60
1:A:62:GLU:HB3	1:A:105:LEU:HD23	1.83	0.60
1:B:17:GLN:NE2	1:B:21:GLU:OE1	2.34	0.60
1:B:99:ILE:HG23	1:B:99:ILE:O	2.00	0.60
1:A:588:GLU:O	1:A:591:SER:HB3	2.01	0.60
1:A:660:GLU:HG3	1:A:701:HIS:HE1	1.65	0.60
1:A:263:GLN:OE1	1:A:264:ARG:NH1	2.34	0.60
1:B:481:LEU:O	1:B:485:LEU:HB2	2.02	0.59
1:B:840:ALA:HB1	1:B:880:LEU:HD22	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:794:ILE:HG12	1:B:828:VAL:HG12	1.84	0.59
1:B:790:LEU:HD11	1:B:794:ILE:HD13	1.84	0.59
1:A:850:ARG:HH21	1:A:887:PHE:HE1	1.47	0.59
1:B:511:ALA:HB3	1:B:515:LEU:HD21	1.85	0.59
1:B:214:ASP:O	1:B:218:GLU:HB2	2.03	0.58
1:B:508:GLU:CD	1:B:548:LEU:HD11	2.28	0.58
1:B:755:VAL:HA	1:B:758:ILE:HG22	1.84	0.58
1:A:486:LEU:HD23	1:A:489:ILE:HD12	1.84	0.58
1:A:546:GLY:C	2:X:518:ARG:HH12	2.11	0.58
1:A:163:LEU:HB3	1:A:172:LEU:HD21	1.86	0.58
1:B:588:GLU:O	1:B:591:SER:HB3	2.04	0.57
1:A:21:GLU:HG3	1:A:33:VAL:HG21	1.87	0.57
1:B:103:SER:CB	1:B:104:PRO:HD2	2.11	0.57
1:B:508:GLU:CG	1:B:548:LEU:CD1	2.76	0.57
1:B:179:PHE:HA	1:B:182:PHE:CZ	2.38	0.57
1:B:556:LEU:O	1:B:557:ASN:CB	2.52	0.57
1:B:869:TRP:O	1:B:873:SER:HB3	2.05	0.57
1:A:390:TYR:HB2	1:A:394:LEU:HD23	1.88	0.56
1:A:681:PRO:HB3	1:A:721:PHE:HD2	1.70	0.56
1:A:55:THR:HG21	1:A:94:GLU:HB3	1.85	0.56
1:A:183:PHE:HZ	1:A:220:LEU:HA	1.70	0.56
1:A:376:ARG:NH2	1:A:415:GLU:OE2	2.38	0.56
1:A:739:MET:O	1:A:742:GLU:HB2	2.05	0.56
2:Y:516:GLU:O	2:Y:520:ASP:CB	2.42	0.56
1:A:436:LEU:HD23	1:A:439:LEU:HD12	1.87	0.56
1:A:610:GLN:O	1:A:614:ASN:ND2	2.27	0.56
1:B:864:VAL:HG13	1:B:868:ASN:HB2	1.87	0.56
1:B:508:GLU:OE1	1:B:515:LEU:HG	2.05	0.56
1:B:648:LEU:HD23	1:B:651:LEU:HD12	1.87	0.56
1:A:588:GLU:OE1	2:X:514:ARG:HB2	2.06	0.56
1:B:411:TRP:HB2	1:B:454:VAL:HG21	1.87	0.55
1:A:444:ILE:HD13	1:A:462:LEU:HD11	1.89	0.55
1:A:464:ARG:CZ	2:X:526:TYR:CE1	2.89	0.55
1:A:192:SER:O	1:A:234:ASN:ND2	2.39	0.55
1:A:273:ALA:HB3	1:A:375:LEU:HD11	1.89	0.55
1:B:95:CYS:O	1:B:98:ASN:OD1	2.24	0.55
1:A:76:VAL:HG11	1:A:117:ILE:HD13	1.87	0.55
1:B:399:LEU:HD13	1:B:432:MET:HE1	1.89	0.55
1:B:449:ASP:OD1	1:B:450:LYS:N	2.40	0.55
1:B:570:ILE:O	1:B:574:ASN:ND2	2.31	0.55
1:B:556:LEU:HD11	1:B:596:ALA:CB	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:X:512:ASP:N	2:X:512:ASP:OD1	2.41	0.54
1:A:460:TRP:HH2	2:X:526:TYR:HB3	1.70	0.54
1:A:796:PRO:HA	1:A:799:THR:HB	1.88	0.54
1:A:491:ASP:OD1	1:A:493:ASN:N	2.34	0.54
1:A:546:GLY:C	2:X:518:ARG:NH1	2.66	0.54
1:B:57:LEU:O	1:B:65:ARG:NH2	2.40	0.53
1:A:588:GLU:HG2	1:A:643:VAL:HG23	1.91	0.53
1:B:407:PHE:O	1:B:408:HIS:ND1	2.42	0.53
1:A:449:ASP:OD1	1:A:450:LYS:N	2.42	0.53
1:B:798:CYS:HB3	1:B:838:ALA:HB2	1.91	0.53
1:B:498:GLU:HG2	2:Y:523:GLU:HG2	1.90	0.53
1:A:267:ASP:OD1	1:A:268:GLN:N	2.41	0.53
1:B:95:CYS:HA	1:B:98:ASN:HD21	1.73	0.52
1:B:557:ASN:O	1:B:557:ASN:ND2	2.43	0.52
1:A:4:GLU:OE2	1:A:56:LYS:NZ	2.43	0.52
1:A:850:ARG:HD2	1:A:887:PHE:CE1	2.44	0.52
1:B:869:TRP:O	1:B:873:SER:CB	2.58	0.52
1:A:727:ASN:HD21	2:X:510:LYS:HA	1.75	0.52
1:B:612:CYS:SG	1:B:647:LEU:HD23	2.50	0.52
1:B:643:VAL:O	1:B:646:ASP:HB2	2.09	0.52
1:B:676:MET:HE3	1:B:688:PHE:HE1	1.75	0.52
1:B:556:LEU:CD2	1:B:596:ALA:HB1	2.33	0.52
2:X:524:ARG:HD3	2:X:524:ARG:N	2.25	0.51
1:A:423:ALA:HA	2:X:526:TYR:HD2	1.71	0.51
1:B:876:PHE:HE1	1:B:881:LYS:HA	1.74	0.51
1:B:214:ASP:O	1:B:218:GLU:HB3	2.09	0.51
1:B:810:LYS:H	1:B:810:LYS:HD2	1.74	0.51
1:B:402:LEU:HD21	1:B:420:VAL:HG11	1.93	0.51
1:B:861:LYS:NZ	1:B:890:VAL:O	2.43	0.51
1:A:51:ILE:HG21	1:A:91:ILE:HD13	1.93	0.51
1:B:888:TYR:O	1:B:890:VAL:N	2.43	0.51
1:B:308:MET:HE3	1:B:380:ALA:HA	1.93	0.50
1:A:301:ILE:HG21	1:A:398:ILE:HG12	1.94	0.50
1:B:491:ASP:OD1	1:B:492:SER:N	2.44	0.50
1:A:426:GLU:HB3	2:X:526:TYR:OH	2.11	0.50
1:B:301:ILE:HG21	1:B:398:ILE:HG21	1.93	0.50
1:B:791:GLN:HG3	1:B:827:GLY:HA2	1.93	0.50
2:X:513:SER:O	2:X:517:HIS:CB	2.55	0.50
1:B:216:PHE:O	1:B:220:LEU:HB2	2.12	0.50
1:A:835:PHE:O	1:A:839:VAL:HG23	2.10	0.50
1:B:817:ILE:O	1:B:821:ILE:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:LEU:HB2	1:A:171:PRO:HG3	1.93	0.49
1:B:669:LEU:HD21	1:B:702:VAL:HA	1.94	0.49
1:B:99:ILE:HD13	1:B:114:ILE:HD11	1.91	0.49
1:B:262:LEU:HG	1:B:299:LYS:HE3	1.95	0.49
1:B:443:LEU:HA	1:B:446:CYS:HB2	1.94	0.49
1:B:556:LEU:HD21	1:B:596:ALA:CB	2.37	0.49
1:A:460:TRP:NE1	2:X:524:ARG:HB3	2.26	0.49
1:A:669:LEU:HA	1:A:672:MET:HE2	1.94	0.49
1:A:126:TRP:CH2	1:A:129:LEU:HD13	2.48	0.49
1:B:699:PHE:CD2	1:B:738:GLN:HB3	2.47	0.49
1:A:612:CYS:SG	1:A:647:LEU:HD23	2.53	0.49
1:A:815:ARG:NH2	1:A:848:ASP:OD2	2.45	0.49
1:A:814:PHE:HA	1:A:817:ILE:HG22	1.94	0.48
1:A:26:ASP:OD1	1:A:27:THR:N	2.46	0.48
1:A:237:ARG:O	1:A:241:MET:HG2	2.13	0.48
1:A:23:GLN:NE2	1:A:63:PRO:O	2.46	0.48
1:B:556:LEU:C	1:B:556:LEU:CD2	2.86	0.48
1:B:257:ILE:O	1:B:261:MET:HB2	2.12	0.48
1:B:413:VAL:O	1:B:416:SER:OG	2.26	0.48
2:X:510:LYS:HD3	2:X:512:ASP:HB3	1.95	0.48
1:A:232:ARG:NH1	1:A:267:ASP:OD2	2.47	0.48
1:A:460:TRP:CH2	2:X:526:TYR:CB	2.89	0.48
1:A:573:TRP:CZ2	1:A:611:ARG:HD3	2.49	0.48
1:A:233:LYS:HG3	1:A:275:GLU:HG3	1.96	0.47
1:B:515:LEU:HB2	1:B:552:VAL:HG21	1.95	0.47
1:A:180:LEU:HA	1:A:183:PHE:HB2	1.97	0.47
1:A:297:LEU:O	1:A:300:LEU:N	2.47	0.47
1:B:681:PRO:HA	1:B:684:ARG:HE	1.79	0.47
1:B:703:LYS:O	1:B:706:ILE:HB	2.15	0.47
1:A:681:PRO:HB3	1:A:721:PHE:CD2	2.49	0.47
1:A:850:ARG:NH2	1:A:887:PHE:CD1	2.83	0.47
1:B:51:ILE:O	1:B:55:THR:HG22	2.15	0.47
1:B:487:LYS:HA	1:B:487:LYS:HD3	1.71	0.47
1:B:834:PHE:O	1:B:837:ASP:HB3	2.15	0.47
1:A:183:PHE:CZ	1:A:220:LEU:HA	2.48	0.47
1:A:558:LYS:HB2	1:A:561:TYR:CD2	2.50	0.47
1:A:590:LEU:O	1:A:594:ALA:N	2.40	0.47
1:A:850:ARG:HD2	1:A:887:PHE:HE1	1.79	0.47
1:A:403:LYS:HE3	1:A:403:LYS:HB2	1.77	0.46
1:B:66:SER:CB	1:B:105:LEU:HD11	2.45	0.46
1:B:258:VAL:O	1:B:262:LEU:CB	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:719:PRO:O	1:B:722:ILE:HG13	2.15	0.46
1:A:491:ASP:OD1	1:A:492:SER:N	2.48	0.46
1:A:646:ASP:O	1:A:649:SER:HB3	2.15	0.46
1:A:710:MET:HE2	1:A:743:MET:HE1	1.98	0.46
1:B:544:ALA:O	1:B:548:LEU:HB2	2.16	0.46
1:B:216:PHE:CZ	1:B:242:LEU:HD21	2.51	0.46
1:A:192:SER:OG	1:A:230:GLU:HB3	2.15	0.46
1:A:221:PHE:CE1	1:A:257:ILE:HD11	2.50	0.46
1:B:868:ASN:O	1:B:872:PHE:HB2	2.16	0.46
1:A:798:CYS:HB2	1:A:838:ALA:HB2	1.98	0.46
1:A:100:GLY:HA3	1:A:144:THR:HG22	1.98	0.46
1:B:835:PHE:O	1:B:839:VAL:HG23	2.16	0.46
1:B:494:LYS:NZ	1:B:534:GLN:OE1	2.49	0.46
1:B:211:LEU:HD13	1:B:211:LEU:HA	1.84	0.46
1:B:469:VAL:C	1:B:471:SER:H	2.24	0.46
1:B:647:LEU:O	1:B:651:LEU:N	2.39	0.46
1:B:660:GLU:HG2	1:B:701:HIS:NE2	2.31	0.46
1:A:408:HIS:CE1	1:A:410:GLU:HB3	2.51	0.45
1:A:703:LYS:HB3	1:A:704:PRO:HD3	1.98	0.45
1:B:411:TRP:CE2	1:B:451:LYS:HG3	2.51	0.45
1:B:739:MET:O	1:B:742:GLU:HB2	2.16	0.45
1:B:65:ARG:NE	1:B:106:ILE:HD11	2.26	0.45
1:B:58:LYS:HA	1:B:65:ARG:NH2	2.31	0.45
1:B:562:ILE:HG23	1:B:566:MET:HG3	1.98	0.45
1:B:504:PHE:CZ	1:B:508:GLU:OE2	2.69	0.45
1:B:747:ILE:HD12	1:B:778:LEU:HD22	1.99	0.45
2:Y:522:ARG:HA	2:Y:522:ARG:HD3	1.35	0.45
1:B:556:LEU:HD11	1:B:596:ALA:HB1	1.99	0.45
1:A:370:ILE:HG23	1:A:371:SER:O	2.16	0.45
1:A:463:SER:HB3	1:A:503:ALA:HB1	1.99	0.45
1:B:65:ARG:HE	1:B:106:ILE:CD1	2.28	0.45
1:B:98:ASN:OD1	1:B:99:ILE:N	2.37	0.45
1:A:440:ILE:O	1:A:444:ILE:HG12	2.17	0.45
1:B:99:ILE:HD13	1:B:114:ILE:CD1	2.47	0.45
2:Y:517:HIS:O	2:Y:521:ARG:N	2.37	0.45
1:A:416:SER:O	1:A:420:VAL:HG23	2.17	0.45
1:B:30:GLN:O	1:B:34:GLN:HB2	2.17	0.45
1:A:96:LEU:HD21	1:A:132:LYS:HG2	1.99	0.44
1:A:192:SER:HB2	1:A:231:VAL:HG22	1.99	0.44
1:B:32:THR:O	1:B:36:LYS:HG2	2.16	0.44
1:B:873:SER:HA	1:B:876:PHE:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:LEU:HD12	1:B:105:LEU:C	2.39	0.44
1:B:371:SER:OG	1:B:372:ASP:N	2.45	0.44
1:B:463:SER:HB3	1:B:503:ALA:HB1	1.99	0.44
1:B:481:LEU:O	1:B:485:LEU:CB	2.65	0.44
1:A:401:LEU:O	1:A:405:LEU:HG	2.17	0.44
1:B:814:PHE:CZ	1:B:838:ALA:HB1	2.52	0.44
1:A:1:MET:H	1:A:1:MET:HG2	1.72	0.44
1:A:509:GLU:OE1	2:X:522:ARG:NH1	2.50	0.44
1:B:648:LEU:HA	1:B:651:LEU:HB2	2.00	0.44
1:A:794:ILE:HG12	1:A:828:VAL:HG12	1.99	0.44
1:B:656:GLY:HA2	1:B:697:ALA:HA	2.00	0.44
1:A:304:LEU:O	1:A:308:MET:HG3	2.18	0.43
1:B:27:THR:HA	1:B:30:GLN:HG2	2.01	0.43
1:A:53:VAL:HA	1:A:57:LEU:HB2	2.01	0.43
1:A:727:ASN:ND2	2:X:509:GLY:O	2.51	0.43
1:A:660:GLU:HG3	1:A:701:HIS:CE1	2.50	0.43
1:A:779:GLY:HA3	1:A:820:MET:HE2	1.98	0.43
1:B:127:PRO:HB2	1:B:128:ASP:H	1.59	0.43
1:A:171:PRO:O	1:A:174:ILE:HG22	2.18	0.43
1:B:176:ILE:HA	1:B:179:PHE:HD2	1.83	0.43
1:B:471:SER:HB3	1:B:477:TYR:HE2	1.84	0.43
1:B:520:ALA:O	1:B:524:ASP:HB2	2.19	0.43
1:B:243:LEU:HD21	1:B:283:LEU:HB2	2.01	0.43
1:B:301:ILE:O	1:B:305:VAL:HG23	2.17	0.43
1:A:460:TRP:HH2	2:X:526:TYR:CB	2.31	0.43
1:A:509:GLU:CD	2:X:522:ARG:NH1	2.75	0.43
1:A:662:LEU:HD23	1:A:662:LEU:HA	1.83	0.43
1:A:790:LEU:O	1:A:794:ILE:HB	2.18	0.43
1:A:834:PHE:O	1:A:837:ASP:HB3	2.18	0.43
1:B:504:PHE:CE1	1:B:508:GLU:OE2	2.72	0.43
1:A:18:LEU:O	1:A:22:SER:HB2	2.19	0.43
1:A:76:VAL:O	1:A:80:PHE:HB2	2.19	0.43
1:B:376:ARG:NE	1:B:419:LEU:HD22	2.33	0.43
2:Y:521:ARG:HA	2:Y:524:ARG:HH21	1.83	0.43
1:A:77:LYS:HA	1:A:120:LYS:HG2	2.00	0.43
1:A:723:SER:OG	2:X:511:MET:HE3	2.19	0.43
1:A:883:ARG:O	1:A:887:PHE:CB	2.67	0.42
1:A:887:PHE:CD1	1:A:887:PHE:O	2.70	0.42
1:B:211:LEU:HB3	1:B:212:HIS:H	1.54	0.42
1:A:498:GLU:HB3	1:A:537:ASN:OD1	2.18	0.42
1:B:99:ILE:HD11	1:B:114:ILE:CD1	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:LYS:HE3	1:B:290:LYS:HB3	1.81	0.42
1:B:449:ASP:O	1:B:455:ARG:NE	2.38	0.42
1:B:783:PRO:O	1:B:787:ALA:HB2	2.19	0.42
1:A:273:ALA:O	1:A:277:CYS:HB2	2.19	0.42
1:B:472:GLN:N	1:B:472:GLN:OE1	2.52	0.42
1:A:154:LYS:HD2	1:A:154:LYS:HA	1.92	0.42
1:B:440:ILE:O	1:B:444:ILE:HG13	2.20	0.42
1:A:70:LEU:O	1:A:73:LYS:HB3	2.20	0.42
1:B:283:LEU:HD22	1:B:293:LEU:HD11	2.02	0.42
1:B:868:ASN:O	1:B:872:PHE:CB	2.67	0.42
1:A:371:SER:OG	1:A:372:ASP:N	2.52	0.42
1:A:783:PRO:O	1:A:787:ALA:HB2	2.19	0.42
1:B:651:LEU:HD23	1:B:651:LEU:HA	1.91	0.42
1:A:19:LEU:HA	1:A:22:SER:HB3	2.01	0.42
1:A:53:VAL:HG22	1:A:57:LEU:HD12	2.02	0.42
1:B:447:LEU:HD13	1:B:488:ARG:NH1	2.34	0.42
1:A:306:ASN:O	1:A:309:LYS:NZ	2.48	0.41
1:A:383:LEU:HD23	1:A:423:ALA:HB3	2.02	0.41
1:A:840:ALA:HB1	1:A:880:LEU:HD22	2.01	0.41
1:A:387:ALA:HB1	1:A:427:GLY:HA3	2.03	0.41
1:A:848:ASP:N	1:A:848:ASP:OD1	2.53	0.41
1:B:486:LEU:HD23	1:B:489:ILE:HD12	2.02	0.41
1:B:815:ARG:NH1	1:B:848:ASP:OD2	2.53	0.41
1:A:547:THR:HG21	2:X:522:ARG:NH2	2.34	0.41
1:B:258:VAL:O	1:B:262:LEU:HB2	2.20	0.41
1:A:5:TRP:CH2	1:A:48:ASN:HB3	2.56	0.41
1:B:51:ILE:HD11	1:B:98:ASN:ND2	2.35	0.41
1:B:192:SER:HB2	1:B:231:VAL:HG22	2.01	0.41
1:B:405:LEU:HA	1:B:408:HIS:CD2	2.55	0.41
1:A:105:LEU:O	1:A:109:THR:HG22	2.21	0.41
1:B:107:ARG:HH12	1:B:146:GLU:HB2	1.84	0.41
1:B:468:TRP:C	1:B:470:VAL:H	2.28	0.41
1:B:494:LYS:HG2	1:B:534:GLN:HE22	1.85	0.41
1:B:848:ASP:OD1	1:B:848:ASP:N	2.53	0.41
1:B:508:GLU:CD	1:B:548:LEU:CD1	2.93	0.41
1:A:582:ASP:O	1:A:585:PRO:HD2	2.20	0.41
1:B:556:LEU:HD11	1:B:596:ALA:C	2.46	0.41
1:A:462:LEU:HD22	1:A:481:LEU:HD13	2.02	0.41
1:B:19:LEU:HB3	1:B:64:THR:HG23	2.03	0.41
1:B:543:ASP:O	1:B:547:THR:OG1	2.22	0.41
1:A:385:VAL:O	1:A:389:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:719:PRO:O	1:A:722:ILE:HG12	2.21	0.41
1:A:107:ARG:NH1	1:A:143:ASN:OD1	2.54	0.40
1:A:243:LEU:HD22	1:A:279:PHE:CE1	2.56	0.40
1:A:502:SER:CB	2:X:523:GLU:H	2.34	0.40
1:B:18:LEU:HD13	1:B:37:LEU:HD22	2.01	0.40
1:B:512:CYS:O	1:B:515:LEU:HD23	2.20	0.40
1:B:34:GLN:O	1:B:38:GLU:HG3	2.22	0.40
1:B:515:LEU:N	1:B:515:LEU:CD2	2.83	0.40
1:B:68:SER:O	1:B:71:ILE:HG22	2.21	0.40
1:B:585:PRO:O	1:B:588:GLU:HB2	2.21	0.40
1:A:387:ALA:HA	1:A:394:LEU:HD21	2.03	0.40
1:B:301:ILE:HG21	1:B:398:ILE:HG12	2.03	0.40
1:B:619:THR:HG21	1:B:641:MET:HB2	2.04	0.40

There are no symmetry-related clashes.

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	820/868 (94%)	753 (92%)	64 (8%)	3 (0%)	30	65
1	B	738/868 (85%)	675 (92%)	50 (7%)	13 (2%)	6	35
2	X	16/528 (3%)	10 (62%)	4 (25%)	2 (12%)	0	4
2	Y	12/528 (2%)	9 (75%)	2 (17%)	1 (8%)	0	10
All	All	1586/2792 (57%)	1447 (91%)	120 (8%)	19 (1%)	10	42

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	470	VAL
1	B	557	ASN

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Mol	Chain	Res	Type
2	X	511	MET
1	B	127	PRO
1	B	162	ILE
1	B	469	VAL
1	B	889	GLY
1	A	10	GLN
1	B	12	LEU
1	B	228	GLU
2	X	525	PRO
1	A	394	LEU
1	B	103	SER
2	Y	522	ARG
1	B	559	PRO
1	B	472	GLN
1	B	702	VAL
1	A	474	PRO
1	B	213	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	746/776 (96%)	737 (99%)	9 (1%)	63	73
1	B	683/776 (88%)	674 (99%)	9 (1%)	61	72
2	X	16/362 (4%)	13 (81%)	3 (19%)	1	10
2	Y	13/362 (4%)	10 (77%)	3 (23%)	1	6
All	All	1458/2276 (64%)	1434 (98%)	24 (2%)	55	70

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	428	CYS
1	A	513	THR
1	A	706	ILE

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Mol	Chain	Res	Type
1	A	822	SER
1	A	858	HIS
1	A	861	LYS
1	A	863	GLN
1	A	871	ARG
1	B	101	ASP
1	B	156	CYS
1	B	211	LEU
1	B	213	ILE
1	B	220	LEU
1	B	548	LEU
1	B	556	LEU
1	B	558	LYS
1	B	822	SER
2	X	517	HIS
2	X	524	ARG
2	X	526	TYR
2	Y	514	ARG
2	Y	518	ARG
2	Y	522	ARG

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	97	ASN
1	A	124	GLN
1	A	153	GLN
1	A	271	ASN
1	A	685	GLN
1	A	701	HIS
1	A	727	ASN
1	B	296	HIS
1	B	430	GLN
1	B	493	ASN
1	B	557	ASN
1	B	629	GLN
1	B	762	ASN
1	B	824	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 [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	830/868 (95%)	-1.14	0 100 100	4, 33, 70, 97	0
1	B	764/868 (88%)	-1.05	0 100 100	6, 59, 109, 137	0
2	X	18/528 (3%)	-1.05	0 100 100	14, 24, 32, 32	0
2	Y	14/528 (2%)	-1.13	0 100 100	40, 59, 77, 78	0
All	All	1626/2792 (58%)	-1.10	0 100 100	4, 38, 101, 137	0

There are no RSRZ outliers to report.

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.