



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:36 PM GMT

PDB ID : 4PD3
Title : Crystal Structure of Rigor-Like Human Nonmuscle Myosin-2B
Authors : Munnich, S.; Pathan-Chhatbar, S.; Manstein, D.J.
Deposited on : 2014-04-17
Resolution : 2.84 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

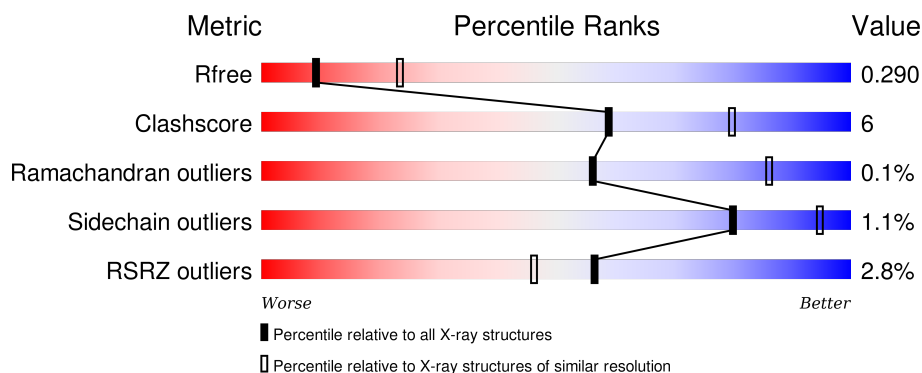
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	3170 (2.88-2.80)
Clashscore	102246	3658 (2.88-2.80)
Ramachandran outliers	100387	3591 (2.88-2.80)
Sidechain outliers	100360	3594 (2.88-2.80)
RSRZ outliers	91569	3184 (2.88-2.80)

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 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	1032	<div> <div>2%</div> <div>76%</div> <div>13%</div> <div>11%</div> </div>
1	B	1032	<div> <div>3%</div> <div>74%</div> <div>16%</div> <div>10%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14828 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 Nonmuscle myosin heavy chain B, Alpha-actinin A Chimera Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	922	Total	C	N	O	S	0	0	0
			7379	4672	1284	1388	35			
1	B	930	Total	C	N	O	S	0	0	0
			7416	4698	1288	1395	35			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	783	ALA	-	linker	UNP P35580
A	784	SER	-	linker	UNP P35580
A	1023	ALA	-	expression tag	UNP P05095
A	1024	LEU	-	expression tag	UNP P05095
A	1025	HIS	-	expression tag	UNP P05095
A	1026	HIS	-	expression tag	UNP P05095
A	1027	HIS	-	expression tag	UNP P05095
A	1028	HIS	-	expression tag	UNP P05095
A	1029	HIS	-	expression tag	UNP P05095
A	1030	HIS	-	expression tag	UNP P05095
A	1031	HIS	-	expression tag	UNP P05095
A	1032	HIS	-	expression tag	UNP P05095
B	783	ALA	-	linker	UNP P35580
B	784	SER	-	linker	UNP P35580
B	1023	ALA	-	expression tag	UNP P05095
B	1024	LEU	-	expression tag	UNP P05095
B	1025	HIS	-	expression tag	UNP P05095
B	1026	HIS	-	expression tag	UNP P05095
B	1027	HIS	-	expression tag	UNP P05095
B	1028	HIS	-	expression tag	UNP P05095
B	1029	HIS	-	expression tag	UNP P05095
B	1030	HIS	-	expression tag	UNP P05095
B	1031	HIS	-	expression tag	UNP P05095
B	1032	HIS	-	expression tag	UNP P05095

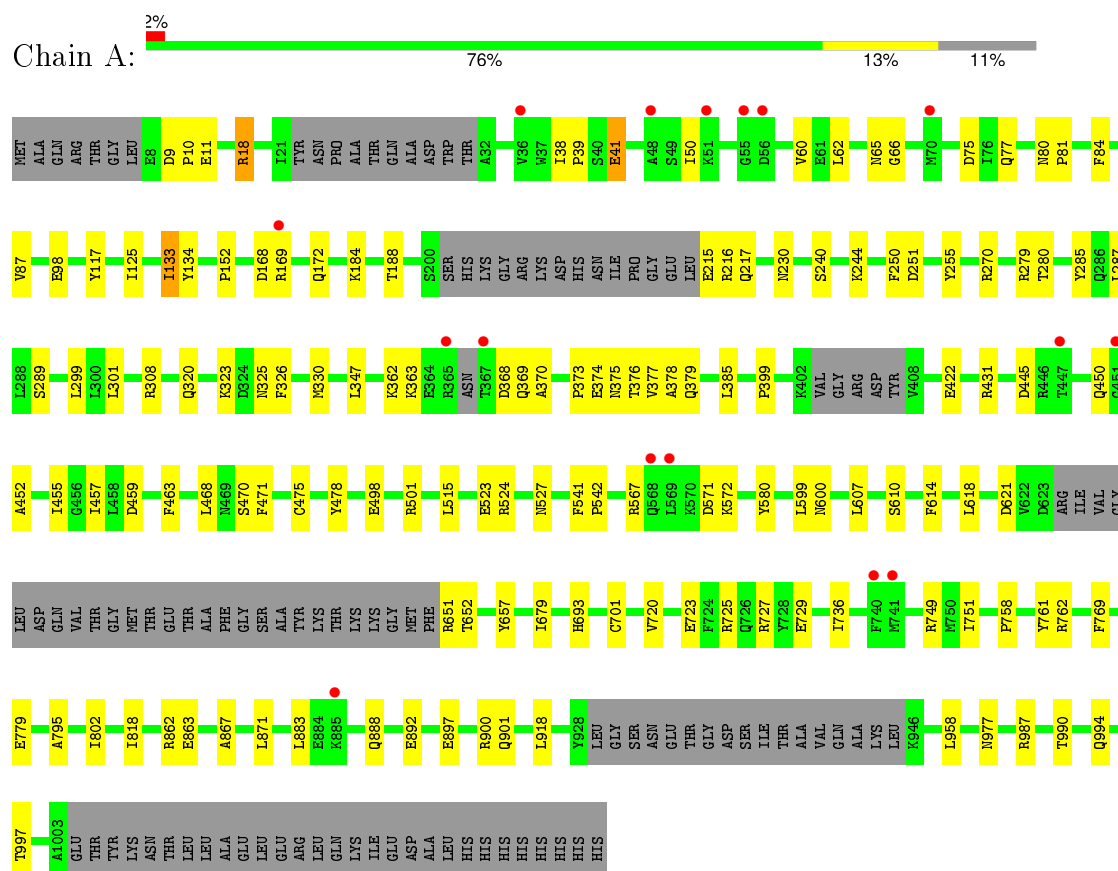
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	18	Total 18	O 18	0	0
2	B	15	Total 15	O 15	0	0

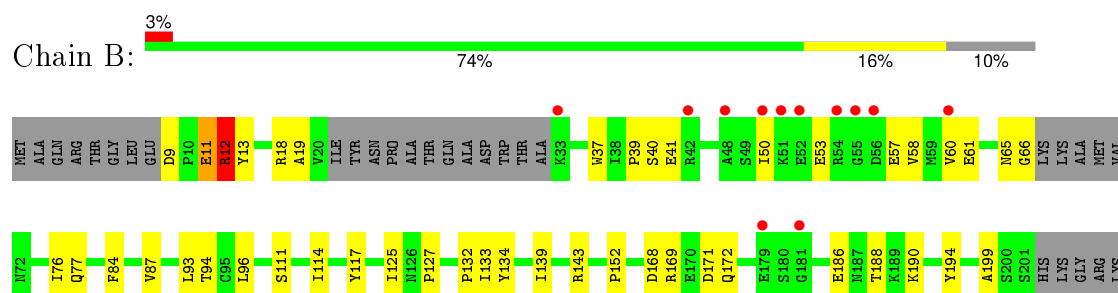
3 Residue-property plots

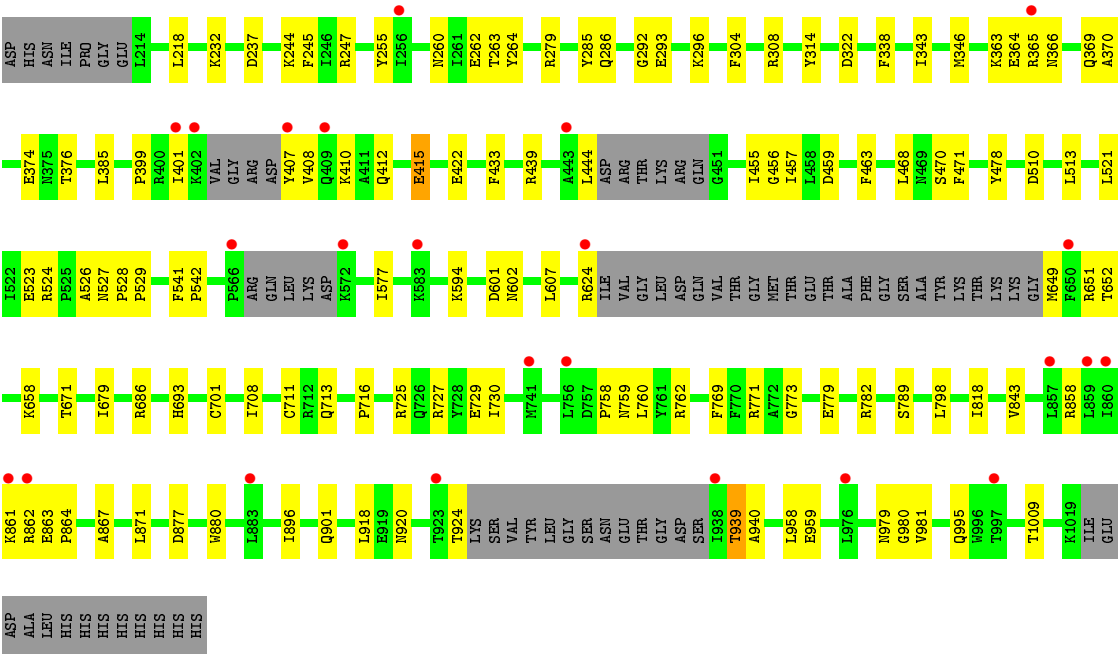
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Nonmuscle myosin heavy chain B, Alpha-actinin A Chimera Protein



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.85Å 157.87Å 143.48Å 90.00° 94.21° 90.00°	Depositor
Resolution (Å)	30.00 – 2.84 30.00 – 2.84	Depositor EDS
% Data completeness (in resolution range)	95.1 (30.00-2.84) 83.2 (30.00-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.85Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.253 , 0.288 0.253 , 0.290	Depositor DCC
R_{free} test set	1769 reflections (3.09%)	DCC
Wilson B-factor (Å ²)	56.6	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 32.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 65526 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14828	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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, 0.375 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	0.21	0/7515	0.39	0/10128
1	B	0.21	0/7551	0.41	2/10180 (0.0%)
All	All	0.21	0/15066	0.40	2/20308 (0.0%)

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	2
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	981	VAL	C-N-CD	-6.83	105.58	120.60
1	B	981	VAL	C-N-CA	5.15	143.62	122.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	41	GLU	Peptide
1	B	61	GLU	Peptide
1	B	858	ARG	Peptide

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	7379	0	7292	71	0
1	B	7416	0	7300	101	0
2	A	18	0	0	0	0
2	B	15	0	0	0	0
All	All	14828	0	14592	171	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 6.

All (171) 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:12:ARG:HD3	1:B:18:ARG:HB3	1.46	0.97
1:B:12:ARG:HB3	1:B:18:ARG:CB	2.01	0.90
1:B:12:ARG:CB	1:B:18:ARG:HB2	2.10	0.82
1:B:11:GLU:O	1:B:13:TYR:N	2.13	0.81
1:B:12:ARG:HB3	1:B:18:ARG:HB2	1.62	0.77
1:A:818:ILE:HG22	1:A:901:GLN:HB3	1.67	0.75
1:B:12:ARG:CB	1:B:18:ARG:CB	2.63	0.75
1:A:62:LEU:HD23	1:A:65:ASN:HD21	1.52	0.74
1:B:76:ILE:HA	1:B:77:GLN:HB2	1.71	0.71
1:B:818:ILE:HG22	1:B:901:GLN:HB3	1.71	0.70
1:B:12:ARG:HB3	1:B:18:ARG:HB3	1.73	0.70
1:B:510:ASP:HB3	1:B:513:LEU:HD13	1.75	0.68
1:B:308:ARG:NH2	1:B:422:GLU:OE2	2.26	0.68
1:A:431:ARG:NH2	1:A:621:ASP:O	2.28	0.67
1:A:320:GLN:HG2	1:A:325:ASN:HD21	1.60	0.66
1:B:39:PRO:HD3	1:B:76:ILE:HD11	1.76	0.66
1:A:729:GLU:HG3	1:A:736:ILE:HD12	1.78	0.65
1:A:215:GLU:HG2	1:A:216:ARG:H	1.62	0.64
1:A:900:ARG:NH2	1:A:977:ASN:O	2.31	0.64
1:B:727:ARG:O	1:B:782:ARG:NH1	2.31	0.64
1:A:308:ARG:NH2	1:A:422:GLU:OE2	2.31	0.63
1:B:363:LYS:HZ3	1:B:369:GLN:H	1.46	0.62
1:A:50:ILE:HA	1:A:60:VAL:HG12	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:527:ASN:O	1:B:529:PRO:HD3	2.00	0.61
1:B:12:ARG:HB2	1:B:18:ARG:HB2	1.82	0.61
1:B:861:LYS:HG3	1:B:862:ARG:H	1.67	0.60
1:B:12:ARG:CB	1:B:18:ARG:HB3	2.29	0.60
1:B:524:ARG:HD3	1:B:528:PRO:HG2	1.83	0.59
1:B:959:GLU:HB2	1:B:995:GLN:HE22	1.66	0.59
1:A:374:GLU:HG3	1:A:376:THR:HG22	1.84	0.58
1:A:217:GLN:NE2	1:A:445:ASP:OD2	2.36	0.58
1:A:720:VAL:HG22	1:A:723:GLU:HG2	1.86	0.58
1:A:897:GLU:OE2	1:A:900:ARG:NH1	2.37	0.58
1:A:862:ARG:HG3	1:A:863:GLU:H	1.68	0.57
1:B:374:GLU:HG3	1:B:376:THR:HG22	1.86	0.57
1:B:41:GLU:O	1:B:693:HIS:NE2	2.30	0.56
1:B:65:ASN:OD1	1:B:66:GLY:N	2.39	0.56
1:A:188:THR:HG23	1:A:457:ILE:HG21	1.88	0.55
1:B:365:ARG:HG2	1:B:366:ASN:H	1.71	0.55
1:A:802:ILE:HG23	1:A:883:LEU:HD22	1.89	0.55
1:B:401:ILE:HD12	1:B:601:ASP:HB2	1.88	0.54
1:B:199:ALA:HB1	1:B:255:TYR:HA	1.90	0.54
1:A:369:GLN:OE1	1:A:370:ALA:N	2.39	0.54
1:B:188:THR:HG23	1:B:457:ILE:HG21	1.90	0.54
1:B:247:ARG:HD3	1:B:260:ASN:HD21	1.73	0.54
1:A:918:LEU:HD22	1:A:958:LEU:HD21	1.90	0.53
1:B:759:ASN:HA	1:B:762:ARG:HH12	1.74	0.53
1:B:244:LYS:HB3	1:B:459:ASP:HB3	1.91	0.52
1:B:50:ILE:HA	1:B:60:VAL:HG11	1.90	0.52
1:B:370:ALA:N	1:B:412:GLN:O	2.43	0.52
1:B:53:GLU:HG2	1:B:58:VAL:HG13	1.92	0.52
1:A:599:LEU:HB2	1:A:651:ARG:HD3	1.92	0.51
1:B:343:ILE:HA	1:B:346:MET:HE2	1.92	0.51
1:A:727:ARG:NH2	1:A:779:GLU:OE2	2.35	0.51
1:B:364:GLU:O	1:B:366:ASN:ND2	2.45	0.50
1:B:478:TYR:OH	1:B:523:GLU:OE1	2.28	0.50
1:A:250:PHE:O	1:A:452:ALA:HB1	2.12	0.50
1:A:399:PRO:HG3	1:A:600:ASN:HD21	1.77	0.49
1:A:363:LYS:NZ	1:A:368:ASP:HA	2.28	0.49
1:B:117:TYR:CZ	1:B:152:PRO:HA	2.47	0.49
1:B:9:ASP:HB2	1:B:132:PRO:HG3	1.94	0.49
1:A:289:SER:O	1:A:323:LYS:NZ	2.45	0.49
1:B:713:GLN:O	1:B:771:ARG:NH1	2.45	0.49
1:A:270:ARG:HG2	1:A:280:THR:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ASN:OD1	1:A:66:GLY:N	2.46	0.49
1:B:877:ASP:HA	1:B:880:TRP:HB3	1.93	0.48
1:B:524:ARG:NH1	1:B:526:ALA:HB3	2.27	0.48
1:A:39:PRO:HG2	1:A:75:ASP:HB3	1.94	0.48
1:B:369:GLN:OE1	1:B:370:ALA:N	2.47	0.48
1:B:94:THR:O	1:B:773:GLY:N	2.46	0.48
1:A:463:PHE:HB3	1:A:701:CYS:HB3	1.95	0.48
1:B:918:LEU:HD22	1:B:958:LEU:HD21	1.96	0.48
1:A:117:TYR:CZ	1:A:152:PRO:HA	2.48	0.48
1:B:725:ARG:O	1:B:729:GLU:HB2	2.14	0.48
1:A:184:LYS:NZ	1:A:459:ASP:OD1	2.34	0.48
1:A:41:GLU:O	1:A:693:HIS:NE2	2.40	0.47
1:A:77:GLN:HE22	1:A:98:GLU:HG2	1.80	0.47
1:A:377:VAL:HG12	1:A:378:ALA:H	1.79	0.47
1:A:230:ASN:ND2	1:A:240:SER:HA	2.30	0.47
1:A:385:LEU:HD13	1:A:607:LEU:HD21	1.97	0.47
1:B:232:LYS:HG3	1:B:237:ASP:HA	1.97	0.47
1:A:38:ILE:HD11	1:A:62:LEU:HD11	1.96	0.47
1:B:415:GLU:CD	1:B:415:GLU:H	2.18	0.47
1:B:727:ARG:NH2	1:B:779:GLU:OE2	2.44	0.47
1:B:521:LEU:HD23	1:B:577:ILE:HG23	1.96	0.47
1:B:407:TYR:CD1	1:B:408:VAL:HG13	2.49	0.46
1:A:11:GLU:HB3	1:A:18:ARG:HG2	1.97	0.46
1:B:263:THR:HG21	1:B:433:PHE:HE2	1.79	0.46
1:B:12:ARG:HB2	1:B:18:ARG:CB	2.41	0.46
1:A:279:ARG:HB3	1:A:285:TYR:CE2	2.50	0.46
1:B:365:ARG:CG	1:B:366:ASN:H	2.26	0.46
1:A:987:ARG:O	1:A:990:THR:OG1	2.32	0.46
1:B:762:ARG:HB2	1:B:769:PHE:HB2	1.97	0.46
1:B:399:PRO:HA	1:B:602:ASN:HD22	1.80	0.45
1:B:114:ILE:HG21	1:B:127:PRO:HB3	1.98	0.45
1:A:867:ALA:HB1	1:A:871:LEU:HB2	1.97	0.45
1:A:377:VAL:C	1:A:379:GLN:H	2.20	0.45
1:A:888:GLN:O	1:A:892:GLU:N	2.38	0.45
1:A:362:LYS:HD3	1:A:373:PRO:HG2	1.97	0.45
1:B:58:VAL:HG12	1:B:60:VAL:HG13	1.98	0.45
1:B:470:SER:OG	1:B:594:LYS:NZ	2.42	0.45
1:B:463:PHE:HB3	1:B:701:CYS:HB3	1.99	0.45
1:A:541:PHE:HA	1:A:542:PRO:HD3	1.80	0.45
1:B:262:GLU:OE1	1:B:264:TYR:OH	2.28	0.45
1:A:125:ILE:HG12	1:A:679:ILE:HB	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:TYR:OH	1:A:523:GLU:OE1	2.31	0.44
1:A:758:PRO:HA	1:A:761:TYR:CE2	2.52	0.44
1:B:168:ASP:O	1:B:169:ARG:HG2	2.16	0.44
1:A:168:ASP:O	1:A:169:ARG:HG2	2.17	0.44
1:B:338:PHE:CE1	1:B:439:ARG:HG3	2.52	0.44
1:B:711:CYS:HB3	1:B:716:PRO:HB3	1.98	0.44
1:B:867:ALA:HB1	1:B:871:LEU:HB2	1.99	0.44
1:B:798:LEU:HD21	1:B:843:VAL:HG22	2.00	0.44
1:B:125:ILE:HG12	1:B:679:ILE:HB	1.99	0.44
1:A:9:ASP:HA	1:A:10:PRO:HA	1.83	0.44
1:A:172:GLN:HB2	1:A:455:ILE:HG12	1.99	0.44
1:B:186:GLU:O	1:B:190:LYS:HG2	2.18	0.44
1:B:279:ARG:HB3	1:B:285:TYR:CE2	2.53	0.43
1:A:498:GLU:OE1	1:A:501:ARG:NH2	2.51	0.43
1:B:686:ARG:HD2	1:B:686:ARG:HA	1.90	0.43
1:B:172:GLN:HB2	1:B:455:ILE:HG12	2.00	0.43
1:A:725:ARG:O	1:A:729:GLU:HB2	2.17	0.43
1:B:293:GLU:HA	1:B:296:LYS:HE2	2.00	0.43
1:B:40:SER:OG	1:B:41:GLU:N	2.51	0.43
1:A:795:ALA:HB1	1:A:871:LEU:HD22	2.00	0.43
1:A:133:ILE:HG22	1:A:134:TYR:H	1.84	0.43
1:B:304:PHE:CG	1:B:314:TYR:HB2	2.54	0.43
1:A:614:PHE:O	1:A:618:LEU:HD13	2.18	0.43
1:B:12:ARG:CD	1:B:18:ARG:HB3	2.34	0.43
1:A:524:ARG:NH1	1:A:527:ASN:HB2	2.34	0.43
1:A:251:ASP:OD1	1:A:255:TYR:N	2.33	0.43
1:B:410:LYS:HE3	1:B:412:GLN:HE22	1.84	0.42
1:B:470:SER:OG	1:B:471:PHE:N	2.50	0.42
1:B:385:LEU:HD13	1:B:607:LEU:HD21	2.00	0.42
1:B:19:ALA:HB1	1:B:111:SER:HB3	2.01	0.42
1:B:524:ARG:HH12	1:B:526:ALA:HB3	1.83	0.42
1:A:751:ILE:HG21	1:A:761:TYR:CZ	2.55	0.42
1:B:541:PHE:HA	1:B:542:PRO:HD3	1.81	0.42
1:A:470:SER:OG	1:A:471:PHE:N	2.50	0.42
1:A:84:PHE:O	1:A:87:VAL:HG13	2.18	0.42
1:B:84:PHE:O	1:B:87:VAL:HG13	2.20	0.42
1:A:80:ASN:HA	1:A:81:PRO:HD3	1.95	0.42
1:B:286:GLN:HA	1:B:322:ASP:HB3	2.02	0.42
1:A:994:GLN:O	1:A:997:THR:OG1	2.33	0.42
1:B:730:ILE:HD11	1:B:789:SER:HB3	2.02	0.42
1:B:758:PRO:C	1:B:760:LEU:H	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:571:ASP:OD1	1:A:572:LYS:N	2.53	0.42
1:A:762:ARG:HB2	1:A:769:PHE:HB2	2.01	0.42
1:B:171:ASP:HB2	1:B:671:THR:HG22	2.02	0.42
1:B:863:GLU:HA	1:B:864:PRO:HD3	1.80	0.42
1:B:979:ASN:HA	1:B:980:GLY:HA2	1.69	0.42
1:B:920:ASN:O	1:B:924:THR:HG22	2.20	0.42
1:B:292:GLY:O	1:B:296:LYS:HG3	2.20	0.42
1:B:939:THR:HB	1:B:940:ALA:H	1.65	0.42
1:B:624:ARG:HD2	1:B:624:ARG:HA	1.87	0.42
1:B:218:LEU:HD13	1:B:444:LEU:HD21	2.01	0.41
1:B:37:TRP:O	1:B:76:ILE:HD12	2.19	0.41
1:A:475:CYS:HB3	1:A:657:TYR:CZ	2.54	0.41
1:B:896:ILE:HD13	1:B:896:ILE:HA	1.92	0.41
1:A:326:PHE:O	1:A:330:MET:HG2	2.20	0.41
1:B:245:PHE:CZ	1:B:456:GLY:HA3	2.55	0.41
1:B:139:ILE:HG23	1:B:194:TYR:CD1	2.56	0.41
1:A:375:ASN:HD21	1:B:143:ARG:HB2	1.86	0.41
1:B:133:ILE:HG22	1:B:134:TYR:H	1.85	0.41
1:B:96:LEU:HB3	1:B:708:ILE:HG23	2.01	0.41
1:B:523:GLU:HG3	1:B:658:LYS:HE3	2.03	0.41
1:A:244:LYS:HB3	1:A:459:ASP:HB3	2.02	0.41
1:A:299:LEU:HB2	1:A:301:LEU:HG	2.03	0.40
1:A:287:LEU:HD13	1:A:347:LEU:HD22	2.03	0.40
1:A:515:LEU:HD11	1:A:580:TYR:HB3	2.04	0.40
1:B:624:ARG:NH2	1:B:651:ARG:HB2	2.37	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	908/1032 (88%)	874 (96%)	34 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	912/1032 (88%)	878 (96%)	33 (4%)	1 (0%)	56 86
All	All	1820/2064 (88%)	1752 (96%)	67 (4%)	1 (0%)	56 86

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	12	ARG

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	792/904 (88%)	784 (99%)	8 (1%)	82 95
1	B	791/904 (88%)	781 (99%)	10 (1%)	76 93
All	All	1583/1808 (88%)	1565 (99%)	18 (1%)	80 95

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	133	ILE
1	A	450	GLN
1	A	468	LEU
1	A	567	ARG
1	A	610	SER
1	A	652	THR
1	A	749	ARG
1	B	11	GLU
1	B	12	ARG
1	B	57	GLU
1	B	93	LEU
1	B	415	GLU
1	B	468	LEU
1	B	649	MET

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Mol	Chain	Res	Type
1	B	652	THR
1	B	939	THR
1	B	1009	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	GLN
1	A	325	ASN
1	A	379	GLN
1	A	568	GLN
1	B	217	GLN
1	B	602	ASN
1	B	826	ASN
1	B	995	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 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	922/1032 (89%)	-0.05	16 (1%) 73 65	41, 77, 120, 166	0
1	B	930/1032 (90%)	0.04	36 (3%) 43 32	37, 82, 128, 161	0
All	All	1852/2064 (89%)	-0.01	52 (2%) 56 46	37, 79, 124, 166	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	407	TYR	12.2
1	A	70	MET	8.1
1	B	861	LYS	7.1
1	B	50	ILE	6.3
1	B	862	ARG	6.3
1	A	56	ASP	5.9
1	B	51	LYS	5.1
1	B	56	ASP	4.4
1	B	860	ILE	4.1
1	B	923	THR	3.8
1	A	367	THR	3.8
1	B	54	ARG	3.5
1	A	740	PHE	3.5
1	A	569	LEU	3.4
1	A	447	THR	3.4
1	B	48	ALA	3.4
1	B	401	ILE	3.3
1	B	976	LEU	3.3
1	B	624	ARG	3.2
1	B	997	THR	3.1
1	B	60	VAL	3.1
1	A	365	ARG	3.1
1	A	451	GLY	3.0
1	B	181	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	859	LEU	2.9
1	B	583	LYS	2.8
1	B	52	GLU	2.6
1	A	885	LYS	2.6
1	B	33	LYS	2.5
1	B	756	LEU	2.5
1	B	402	LYS	2.5
1	B	42	ARG	2.5
1	B	256	ILE	2.4
1	B	741	MET	2.4
1	B	650	PHE	2.4
1	B	938	ILE	2.4
1	A	568	GLN	2.4
1	B	443	ALA	2.4
1	A	741	MET	2.3
1	B	55	GLY	2.3
1	A	51	LYS	2.3
1	B	572	LYS	2.2
1	B	365	ARG	2.2
1	A	55	GLY	2.2
1	B	857	LEU	2.2
1	A	48	ALA	2.1
1	A	36	VAL	2.1
1	B	409	GLN	2.1
1	B	883	LEU	2.1
1	B	179	GLU	2.0
1	B	566	PRO	2.0
1	A	169	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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.