



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:54 am GMT

PDB ID : 4FQ3
Title : Crystal structure of transportin/FUS-NLS
Authors : Gong, W.; Niu, C.; Jia, M.; Gao, F.
Deposited on : 2012-06-25
Resolution : 3.00 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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

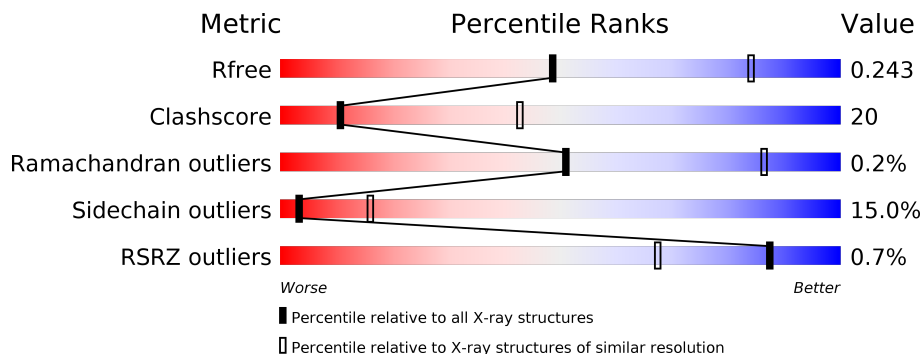
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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	890	
2	B	37	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6805 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	834	Total	C	N	O	S	0	3	0
			6631	4259	1106	1215	51			

- Molecule 2 is a protein called Fusion (Involved in t(12;16) in malignant liposarcoma).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	19	Total	C	N	O	S	0	1	0
			174	101	41	31	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	490	GLY	-	EXPRESSION TAG	UNP Q8TBR3
B	491	PRO	-	EXPRESSION TAG	UNP Q8TBR3
B	492	LEU	-	EXPRESSION TAG	UNP Q8TBR3
B	494	SER	PHE	CONFLICT	UNP Q8TBR3

Chain B:



GLY	PRO	LEU	GLY	SER	ARG	GLY	GLY	ARG	GLY	GLY	GLY	ASP	ARG	ARG	GLY	GLY	PHE	GLY	P508	E516	D520	R521	R522	Y526
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	128.81Å 158.29Å 68.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.94 – 3.00 48.01 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.94-3.00) 99.1 (48.01-2.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.198 , 0.251 0.189 , 0.243	Depositor DCC
R_{free} test set	1456 reflections (5.10%)	DCC
Wilson B-factor (Å ²)	55.8	Xtriage
Anisotropy	0.739	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6805	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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 [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	0.32	0/6784	0.51	0/9218
2	B	0.32	0/180	0.48	0/235
All	All	0.32	0/6964	0.51	0/9453

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	7

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	GLY	Peptide
1	A	168	LEU	Peptide
1	A	746	TYR	Peptide
1	A	8	ASP	Peptide
1	A	841	SER	Peptide
1	A	867	GLU	Peptide
1	A	868	ASN	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6631	0	6687	279	0
2	B	174	0	169	1	0
All	All	6805	0	6856	279	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 20.

All (279) 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:6:LYS:HG2	1:A:7:PRO:HD2	1.31	1.11
1:A:659:ILE:HD11	1:A:697:ALA:HB1	1.48	0.96
1:A:794:ILE:HA	1:A:797:TRP:HB3	1.50	0.91
1:A:204:SER:HB2	1:A:206:THR:HG22	1.58	0.85
1:A:535:HIS:HA	1:A:538:LEU:HB2	1.60	0.83
1:A:787:ALA:HB2	1:A:820:MET:HG2	1.58	0.83
1:A:732:ILE:HD13	1:A:750:VAL:HG11	1.60	0.82
1:A:66:SER:O	1:A:70:LEU:HD22	1.82	0.80
1:A:789:MET:SD	1:A:792:GLN:NE2	2.55	0.80
1:A:790:LEU:O	1:A:791:GLN:HG2	1.82	0.80
1:A:708:ASP:N	1:A:708:ASP:OD1	2.15	0.79
1:A:737:ILE:HD11	1:A:777:ARG:HD2	1.65	0.78
1:A:494:LYS:HE2	1:A:534:GLN:HG3	1.65	0.78
1:A:810:LYS:HA	1:A:813:ALA:HB3	1.68	0.76
1:A:521:TYR:HD1	1:A:521:TYR:H	1.34	0.75
1:A:460:TRP:CZ2	1:A:464:ARG:HD2	2.21	0.75
1:A:815:ARG:NH2	1:A:848:ASP:OD1	2.20	0.75
1:A:848:ASP:OD2	1:A:848:ASP:N	2.18	0.74
1:A:422:GLY:HA3	1:A:460:TRP:CZ3	2.22	0.73
1:A:785:GLU:OE1	1:A:786:VAL:N	2.22	0.72
1:A:6:LYS:CG	1:A:7:PRO:HD2	2.15	0.71
1:A:523:LEU:O	1:A:527:VAL:HG13	1.91	0.70
1:A:885:ALA:HA	1:A:889:GLY:HA2	1.73	0.70
1:A:85:ASN:O	1:A:88:THR:N	2.26	0.69
1:A:821:ILE:HD11	1:A:828:VAL:HG21	1.74	0.69
1:A:673:TYR:O	1:A:677:GLN:NE2	2.26	0.69
1:A:578:ASP:OD2	1:A:611:ARG:NH1	2.24	0.68
1:A:703:LYS:O	1:A:706:ILE:HG13	1.94	0.68
1:A:195:VAL:HG22	1:A:234:ASN:HB3	1.76	0.68
1:A:245:VAL:HG12	1:A:246:ARG:HG2	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:ILE:HD11	1:A:394:LEU:HD11	1.76	0.68
1:A:47:ASN:HD21	1:A:75:ASN:HD22	1.43	0.67
1:A:30:GLN:HA	1:A:33:VAL:HG22	1.77	0.67
1:A:204:SER:HB2	1:A:206:THR:CG2	2.25	0.66
1:A:512:CYS:O	1:A:513:THR:OG1	2.11	0.66
1:A:794:ILE:HA	1:A:797:TRP:CB	2.23	0.66
1:A:661:GLN:O	1:A:665:ARG:HD3	1.95	0.66
1:A:749:MET:O	1:A:753:GLN:NE2	2.28	0.66
1:A:759:ASN:ND2	1:A:792:GLN:O	2.28	0.65
1:A:304:LEU:O	1:A:308:MET:HG3	1.96	0.65
1:A:762:ASN:N	1:A:762:ASN:OD1	2.29	0.65
1:A:760:ARG:HG2	1:A:761:PRO:O	1.96	0.64
1:A:55:THR:OG1	1:A:94:GLU:HG3	1.96	0.64
1:A:770:ASN:O	1:A:774:THR:HG23	1.98	0.64
1:A:706:ILE:HD12	1:A:707:ALA:N	2.13	0.64
1:A:70:LEU:HD13	1:A:109:THR:HG23	1.80	0.64
1:A:31:ARG:O	1:A:35:GLN:NE2	2.31	0.64
1:A:455:ARG:HD2	1:A:488:ARG:HD3	1.79	0.64
1:A:747:ILE:N	1:A:748:PRO:CD	2.60	0.64
1:A:789:MET:HA	1:A:789:MET:CE	2.29	0.63
1:A:108:ALA:O	1:A:112:ILE:HG13	1.98	0.63
1:A:558:LYS:O	1:A:562:ILE:HG12	1.97	0.63
1:A:736:SER:HA	1:A:743:MET:HG3	1.81	0.63
1:A:470:VAL:HG11	1:A:511:ALA:HB2	1.80	0.62
1:A:21:GLU:HB2	1:A:33:VAL:HG11	1.81	0.62
1:A:513:THR:HA	1:A:516:VAL:HG23	1.82	0.62
1:A:606:GLU:O	1:A:610:GLN:HG2	2.00	0.61
1:A:729:THR:O	1:A:774:THR:HG21	1.99	0.61
1:A:502:SER:HB3	2:B:522:ARG:HG2	1.82	0.61
1:A:251:LEU:HA	1:A:254:MET:HG2	1.82	0.61
1:A:395:LEU:CD1	1:A:431:GLY:HA3	2.31	0.60
1:A:746:TYR:HA	1:A:748:PRO:HD2	1.82	0.60
1:A:11:GLY:O	1:A:15:ILE:HG12	2.00	0.60
1:A:534:GLN:O	1:A:535:HIS:HB3	2.00	0.59
1:A:416:SER:O	1:A:420:VAL:HG23	2.03	0.59
1:A:742:GLU:O	1:A:745:PRO:HD2	2.02	0.59
1:A:787:ALA:HB2	1:A:820:MET:CG	2.33	0.59
1:A:80:PHE:HA	1:A:83:PHE:CE1	2.36	0.59
1:A:62:GLU:HG3	1:A:106:ILE:HD13	1.84	0.58
1:A:381:ALA:O	1:A:385:VAL:HG23	2.02	0.58
1:A:787:ALA:C	1:A:789:MET:H	2.07	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:HIS:CD2	1:A:413:VAL:HG11	2.39	0.58
1:A:395:LEU:HD11	1:A:431:GLY:HA3	1.85	0.58
1:A:130:LEU:HB2	1:A:131:PRO:HD3	1.86	0.57
1:A:203:ILE:HA	1:A:245:VAL:HG21	1.86	0.57
1:A:684:ARG:HD2	1:A:716:ASN:OD1	2.04	0.57
1:A:747:ILE:HG22	1:A:751:LEU:HB2	1.87	0.57
1:A:121:GLY:O	1:A:122:GLU:HB2	2.05	0.57
1:A:602:LEU:HD11	1:A:658:ASN:HD22	1.70	0.57
1:A:227:GLU:O	1:A:232:ARG:NH1	2.31	0.57
1:A:723:SER:O	1:A:727:ASN:ND2	2.38	0.56
1:A:790:LEU:C	1:A:792:GLN:H	2.07	0.56
1:A:100:GLY:HA3	1:A:144:THR:HG22	1.87	0.56
1:A:141:ASP:OD2	1:A:144:THR:HG23	2.05	0.56
1:A:787:ALA:O	1:A:789:MET:N	2.38	0.56
1:A:311:SER:O	1:A:315:ILE:HG13	2.05	0.56
1:A:565:LEU:O	1:A:568:PRO:HD2	2.06	0.55
1:A:176:ILE:HB	1:A:177:PRO:HD3	1.88	0.55
1:A:482:MET:HG2	1:A:504:PHE:HE1	1.71	0.55
1:A:732:ILE:HD13	1:A:750:VAL:CG1	2.32	0.55
1:A:264:ARG:NH1	1:A:267:ASP:OD2	2.40	0.55
1:A:619:THR:HG23	1:A:636:PRO:HB2	1.89	0.55
1:A:443:LEU:HB2	1:A:462:LEU:HD21	1.90	0.54
1:A:795:ARG:N	1:A:796:PRO:HD2	2.22	0.54
1:A:48:ASN:ND2	1:A:87:VAL:HG13	2.23	0.54
1:A:786:VAL:HG12	1:A:820:MET:SD	2.48	0.53
1:A:6:LYS:HD2	1:A:48:ASN:OD1	2.08	0.53
1:A:79:HIS:HA	1:A:81:GLN:OE1	2.08	0.53
1:A:659:ILE:HD12	1:A:659:ILE:H	1.73	0.53
1:A:167:VAL:HG23	1:A:168:LEU:HD23	1.91	0.53
1:A:657:GLY:H	1:A:659:ILE:HD11	1.74	0.53
1:A:159:SER:HB3	1:A:162:ILE:HD13	1.90	0.52
1:A:479:LYS:HB3	1:A:480:PRO:HD3	1.90	0.52
1:A:877:PRO:HD2	1:A:880:LEU:HD23	1.90	0.52
1:A:718:ASN:ND2	1:A:720:GLU:OE1	2.43	0.52
1:A:821:ILE:HG21	1:A:856:ILE:HD13	1.92	0.52
1:A:584:PHE:HB2	1:A:585:PRO:HD3	1.91	0.52
1:A:793:PHE:O	1:A:794:ILE:HG22	2.09	0.52
1:A:809:GLU:O	1:A:810:LYS:HG2	2.08	0.52
1:A:516:VAL:HB	1:A:517:PRO:HD3	1.92	0.52
1:A:604:TYR:O	1:A:608:VAL:HG13	2.09	0.52
1:A:470:VAL:CG1	1:A:511:ALA:HB2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:TYR:CD1	1:A:521:TYR:N	2.78	0.52
1:A:629:GLN:HB2	1:A:633:TYR:HD1	1.75	0.51
1:A:649:SER:OG	1:A:693:ASP:OD2	2.27	0.51
1:A:713:LEU:HD13	1:A:732:ILE:HG13	1.92	0.51
1:A:744:GLN:N	1:A:745:PRO:HD2	2.26	0.51
1:A:535:HIS:HA	1:A:538:LEU:CB	2.36	0.51
1:A:848:ASP:O	1:A:852:MET:HG3	2.10	0.51
1:A:126:TRP:CE3	1:A:129:LEU:HD22	2.46	0.51
1:A:725:CYS:O	1:A:729:THR:HG23	2.11	0.51
1:A:167:VAL:CG2	1:A:168:LEU:HD23	2.41	0.51
1:A:747:ILE:N	1:A:748:PRO:HD3	2.25	0.50
1:A:297:LEU:N	1:A:298:PRO:CD	2.74	0.50
1:A:659:ILE:N	1:A:659:ILE:HD12	2.26	0.50
1:A:247:MET:CE	1:A:288:ILE:HD11	2.41	0.50
1:A:99:ILE:HG23	1:A:100:GLY:N	2.27	0.50
1:A:508:GLU:HA	1:A:515:LEU:HD11	1.94	0.50
1:A:494:LYS:CE	1:A:534:GLN:HG3	2.40	0.49
1:A:795:ARG:HB2	1:A:796:PRO:HD3	1.94	0.49
1:A:801:LEU:HD22	1:A:804:ILE:HD12	1.93	0.49
1:A:582:ASP:O	1:A:585:PRO:HD2	2.13	0.49
1:A:581:LYS:HE3	1:A:637:ASP:OD1	2.11	0.49
1:A:47:ASN:ND2	1:A:75:ASN:HD22	2.08	0.49
1:A:731:ALA:O	1:A:735:ILE:HG12	2.12	0.49
1:A:797:TRP:CZ2	1:A:801:LEU:HD12	2.48	0.49
1:A:452:ALA:HB1	1:A:493:ASN:HD22	1.78	0.49
1:A:787:ALA:C	1:A:789:MET:N	2.65	0.49
1:A:203:ILE:HD11	1:A:241:MET:HG2	1.95	0.48
1:A:76:VAL:O	1:A:80:PHE:HB2	2.13	0.48
1:A:527:VAL:O	1:A:530:PHE:HB2	2.13	0.48
1:A:467:HIS:ND1	1:A:467:HIS:C	2.66	0.48
1:A:866:ASP:OD1	1:A:867:GLU:N	2.46	0.48
1:A:195:VAL:CG2	1:A:234:ASN:HB3	2.43	0.48
1:A:533:TYR:HB2	1:A:538:LEU:HD13	1.96	0.48
1:A:460:TRP:CH2	1:A:464:ARG:HD2	2.48	0.48
1:A:254:MET:HE2	1:A:292:VAL:HG11	1.96	0.48
1:A:294:VAL:O	1:A:295:ARG:HD2	2.13	0.48
1:A:78:ALA:O	1:A:79:HIS:CG	2.67	0.48
1:A:860:PHE:CZ	1:A:864:VAL:HG11	2.49	0.48
1:A:80:PHE:CD1	1:A:83:PHE:CZ	3.02	0.48
1:A:808:GLU:O	1:A:811:ASP:HB3	2.14	0.48
1:A:493:ASN:OD1	1:A:494:LYS:N	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ILE:H	1:A:29:ILE:HG12	1.30	0.47
1:A:32:THR:HG22	1:A:36:LYS:HE3	1.95	0.47
1:A:672:MET:O	1:A:676:MET:HG3	2.14	0.47
1:A:312:ASP:O	1:A:316:ILE:HG13	2.14	0.47
1:A:83:PHE:HB2	1:A:84:PRO:HD2	1.95	0.47
1:A:587:LEU:HD22	1:A:647:LEU:HD12	1.97	0.47
1:A:19:LEU:HA	1:A:22:SER:OG	2.15	0.47
1:A:213:ILE:HG23	1:A:214:ASP:N	2.31	0.46
1:A:21:GLU:CB	1:A:33:VAL:HG11	2.44	0.46
1:A:242:LEU:HD12	1:A:242:LEU:HA	1.68	0.46
1:A:488:ARG:NE	1:A:488:ARG:HA	2.31	0.46
1:A:706:ILE:HD12	1:A:707:ALA:H	1.81	0.46
1:A:685:GLN:HG3	1:A:727:ASN:ND2	2.31	0.46
1:A:795:ARG:HB2	1:A:796:PRO:CD	2.45	0.46
1:A:659:ILE:CD1	1:A:659:ILE:H	2.29	0.46
1:A:747:ILE:H	1:A:748:PRO:HD3	1.80	0.46
1:A:246:ARG:NH2	1:A:248:ASP:OD2	2.31	0.46
1:A:233:LYS:HG3	1:A:275:GLU:HG3	1.96	0.46
1:A:314:ASP:O	1:A:318:LEU:HG	2.16	0.46
1:A:625:LEU:HD22	1:A:633:TYR:CD1	2.50	0.46
1:A:622:GLN:HB3	1:A:636:PRO:HB3	1.98	0.45
1:A:710:MET:HB2	1:A:711:PRO:HD3	1.97	0.45
1:A:440:ILE:HB	1:A:441:PRO:HD3	1.98	0.45
1:A:54:LEU:HD22	1:A:69:GLY:HA3	1.98	0.45
1:A:248:ASP:OD1	1:A:248:ASP:N	2.40	0.45
1:A:28:THR:O	1:A:31:ARG:HG2	2.16	0.45
1:A:462:LEU:HD13	1:A:481:LEU:CD1	2.46	0.45
1:A:810:LYS:HE2	1:A:842:TRP:CZ2	2.52	0.45
1:A:96:LEU:HA	1:A:96:LEU:HD12	1.80	0.45
1:A:109:THR:HA	1:A:112:ILE:HD12	1.99	0.45
1:A:299:LYS:HD3	1:A:299:LYS:HA	1.66	0.45
1:A:61:ASP:HB2	1:A:63:PRO:HD2	1.98	0.45
1:A:587:LEU:HB2	1:A:643:VAL:CG1	2.46	0.45
1:A:700:GLN:HG2	1:A:701:HIS:CD2	2.51	0.45
1:A:784:GLN:O	1:A:788:PRO:HD2	2.16	0.45
1:A:113:LEU:HA	1:A:113:LEU:HD12	1.79	0.45
1:A:192:SER:OG	1:A:230:GLU:HB3	2.17	0.45
1:A:699:PHE:CZ	1:A:739:MET:HG2	2.52	0.45
1:A:174:ILE:HD13	1:A:174:ILE:HA	1.73	0.45
1:A:283:LEU:HD22	1:A:283:LEU:O	2.16	0.45
1:A:399:LEU:HD13	1:A:399:LEU:HA	1.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:869:TRP:O	1:A:872:PHE:HB3	2.17	0.45
1:A:168:LEU:HB3	1:A:171:PRO:HG3	1.99	0.44
1:A:374:ASN:OD1	1:A:377:LYS:HG3	2.17	0.44
1:A:867:GLU:O	1:A:868:ASN:HB3	2.17	0.44
1:A:254:MET:CE	1:A:292:VAL:HG11	2.48	0.44
1:A:373:TRP:CZ2	1:A:378:CYS:HA	2.52	0.44
1:A:751:LEU:HD23	1:A:751:LEU:HA	1.68	0.44
1:A:567:PRO:HB2	1:A:568:PRO:HD3	1.99	0.44
1:A:6:LYS:HG2	1:A:7:PRO:CD	2.23	0.44
1:A:889:GLY:N	1:A:890:VAL:HA	2.32	0.44
1:A:793:PHE:CD2	1:A:794:ILE:N	2.85	0.44
1:A:535:HIS:NE2	1:A:582:ASP:OD2	2.49	0.44
1:A:810:LYS:O	1:A:814:PHE:HB2	2.17	0.44
1:A:789:MET:SD	1:A:792:GLN:HB2	2.58	0.44
1:A:99:ILE:HG23	1:A:100:GLY:H	1.83	0.43
1:A:717:LEU:O	1:A:718:ASN:C	2.56	0.43
1:A:41:ASN:ND2	1:A:75:ASN:OD1	2.51	0.43
1:A:533:TYR:CB	1:A:538:LEU:HD13	2.48	0.43
1:A:783:PRO:O	1:A:787:ALA:HB3	2.18	0.43
1:A:790:LEU:C	1:A:792:GLN:N	2.71	0.43
1:A:810:LYS:HE2	1:A:842:TRP:CE2	2.53	0.43
1:A:23:GLN:H	1:A:23:GLN:HG2	1.68	0.43
1:A:78:ALA:O	1:A:79:HIS:CD2	2.71	0.43
1:A:293:LEU:HA	1:A:293:LEU:HD23	1.86	0.43
1:A:274:LEU:HD13	1:A:375:LEU:HA	2.00	0.43
1:A:185:HIS:CD2	1:A:190:ILE:HG21	2.54	0.43
1:A:205:ARG:HG2	1:A:245:VAL:CG1	2.49	0.43
1:A:431:GLY:O	1:A:434:PRO:HD2	2.19	0.43
1:A:651:LEU:O	1:A:655:LEU:CB	2.67	0.43
1:A:83:PHE:HA	1:A:84:PRO:HD3	1.89	0.43
1:A:203:ILE:HA	1:A:245:VAL:CG2	2.47	0.43
1:A:422:GLY:HA3	1:A:460:TRP:HZ3	1.78	0.43
1:A:204:SER:CB	1:A:206:THR:HG22	2.39	0.42
1:A:769:GLU:HB3	1:A:804:ILE:HD11	2.01	0.42
1:A:309:LYS:HG2	1:A:309:LYS:H	1.70	0.42
1:A:318:LEU:O	1:A:319:LYS:HB3	2.18	0.42
1:A:779:GLY:HA3	1:A:820:MET:SD	2.58	0.42
1:A:824:ASN:N	1:A:825:PRO:HD3	2.34	0.42
1:A:100:GLY:HA3	1:A:144:THR:CG2	2.49	0.42
1:A:395:LEU:HD21	1:A:432:MET:HG2	2.00	0.42
1:A:181:GLN:O	1:A:184:LYS:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ARG:O	1:A:267:ASP:HB2	2.20	0.42
1:A:85:ASN:HB3	1:A:86:GLY:H	1.67	0.42
1:A:123:LEU:HD13	1:A:123:LEU:HA	1.88	0.42
1:A:652:ALA:CB	1:A:693:ASP:HB3	2.49	0.42
1:A:651:LEU:O	1:A:655:LEU:HB3	2.19	0.42
1:A:700:GLN:H	1:A:700:GLN:CD	2.21	0.42
1:A:779:GLY:CA	1:A:786:VAL:HG11	2.50	0.42
1:A:635:ALA:HB1	1:A:636:PRO:HD2	2.02	0.42
1:A:583:LEU:HA	1:A:583:LEU:HD23	1.82	0.41
1:A:887:PHE:HD2	1:A:887:PHE:HA	1.78	0.41
1:A:301:ILE:HD11	1:A:394:LEU:CD1	2.48	0.41
1:A:493:ASN:O	1:A:497:GLN:HG3	2.20	0.41
1:A:849:LEU:HD23	1:A:849:LEU:HA	1.73	0.41
1:A:859:GLY:O	1:A:863:GLN:HB2	2.20	0.41
1:A:878:LEU:N	1:A:879:PRO:HD2	2.35	0.41
1:A:876:PHE:HB3	1:A:881:LYS:HB2	2.02	0.41
1:A:573:TRP:HZ3	1:A:587:LEU:CD2	2.33	0.41
1:A:589:CYS:O	1:A:593:VAL:HG23	2.19	0.41
1:A:187:SER:HA	1:A:188:PRO:HD2	1.92	0.41
1:A:220:LEU:HD21	1:A:238:ALA:HB1	2.02	0.41
1:A:512:CYS:C	1:A:514:GLU:H	2.22	0.41
1:A:566:MET:HG2	1:A:604:TYR:CZ	2.55	0.41
1:A:666:SER:HB3	1:A:668:ILE:HG22	2.02	0.41
1:A:298:PRO:O	1:A:302:PRO:HD2	2.20	0.41
1:A:482:MET:HB3	1:A:482:MET:HE3	1.80	0.41
1:A:452:ALA:CB	1:A:493:ASN:HD22	2.33	0.41
1:A:560:GLU:CD	1:A:560:GLU:H	2.24	0.41
1:A:6:LYS:HE2	1:A:7:PRO:CD	2.51	0.41
1:A:162:ILE:O	1:A:165:SER:OG	2.36	0.41
1:A:179:PHE:CE2	1:A:197:CYS:HB3	2.56	0.41
1:A:247:MET:HE2	1:A:288:ILE:HD11	2.02	0.41
1:A:739:MET:CB	1:A:743:MET:HG2	2.50	0.41
1:A:866:ASP:O	1:A:868:ASN:ND2	2.54	0.41
1:A:194:ALA:O	1:A:198:VAL:HG23	2.20	0.41
1:A:402:LEU:HD23	1:A:402:LEU:HA	1.87	0.41
1:A:406:LEU:HA	1:A:406:LEU:HD23	1.82	0.41
1:A:421:LEU:HB3	1:A:461:THR:HG21	2.03	0.41
1:A:482:MET:HE1	1:A:518:TYR:CB	2.51	0.41
1:A:852:MET:O	1:A:856:ILE:HG13	2.20	0.41
1:A:116:THR:O	1:A:120:LYS:HB2	2.21	0.40
1:A:318:LEU:HG	1:A:318:LEU:H	1.65	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASP:HA	1:A:169:ASP:O	2.22	0.40
1:A:606:GLU:N	1:A:607:PRO:HD2	2.36	0.40
1:A:62:GLU:N	1:A:63:PRO:CD	2.85	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	833/890 (94%)	795 (95%)	36 (4%)	2 (0%)	51	86
2	B	18/37 (49%)	17 (94%)	1 (6%)	0	100	100
All	All	851/927 (92%)	812 (95%)	37 (4%)	2 (0%)	51	86

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	794	ILE
1	A	786	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	745/802 (93%)	634 (85%)	111 (15%)	3	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	18/25 (72%)	15 (83%)	3 (17%)	2	13
All	All	763/827 (92%)	649 (85%)	114 (15%)	3	16

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	16	LEU
1	A	20	LYS
1	A	29	ILE
1	A	48	ASN
1	A	50	LEU
1	A	70	LEU
1	A	81	GLN
1	A	88	THR
1	A	94	GLU
1	A	96	LEU
1	A	113	LEU
1	A	115	THR
1	A	119	SER
1	A	123	LEU
1	A	124	GLN
1	A	128	ASP
1	A	132	LYS
1	A	136	LEU
1	A	137	LEU
1	A	146	GLU
1	A	149	PHE
1	A	159	SER
1	A	161	GLU
1	A	166	ASP
1	A	167	VAL
1	A	168	LEU
1	A	174	ILE
1	A	186	SER
1	A	192	SER
1	A	195	VAL
1	A	204	SER
1	A	206	THR
1	A	220	LEU
1	A	242	LEU
1	A	247	MET

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Mol	Chain	Res	Type
1	A	248	ASP
1	A	258	VAL
1	A	262	LEU
1	A	268	GLN
1	A	270	GLU
1	A	283	LEU
1	A	288	ILE
1	A	295	ARG
1	A	297	LEU
1	A	299	LYS
1	A	300	LEU
1	A	309	LYS
1	A	318	LEU
1	A	372	ASP
1	A	399	LEU
1	A	401	LEU
1	A	421	LEU
1	A	424	ILE
1	A	456	SER
1	A	463	SER
1	A	476	THR
1	A	483	THR
1	A	488	ARG
1	A	490	LEU
1	A	501	CYS
1	A	512	CYS
1	A	519	LEU
1	A	521	TYR
1	A	524	ASP
1	A	536	LYS
1	A	539	LEU
1	A	560	GLU
1	A	563	GLN
1	A	564	MET
1	A	566	MET
1	A	570	ILE
1	A	586	LEU
1	A	595	THR
1	A	606	GLU
1	A	608	VAL
1	A	615	LEU
1	A	625	LEU

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Mol	Chain	Res	Type
1	A	634	GLU
1	A	638	LYS
1	A	646	ASP
1	A	659	ILE
1	A	687	SER
1	A	708	ASP
1	A	718	ASN
1	A	746	TYR
1	A	749	MET
1	A	750	VAL
1	A	751	LEU
1	A	762	ASN
1	A	767	LEU
1	A	774	THR
1	A	777	ARG
1	A	785	GLU
1	A	789	MET
1	A	792	GLN
1	A	794	ILE
1	A	800	SER
1	A	817	ILE
1	A	821	ILE
1	A	831	ASP
1	A	833	ILE
1	A	841	SER
1	A	847	ASP
1	A	848	ASP
1	A	849	LEU
1	A	850	ARG
1	A	869	TRP
1	A	875	GLN
1	A	887	PHE
1	A	890	VAL
2	B	516	GLU
2	B	520	ASP
2	B	522	ARG

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	23	GLN
1	A	41	ASN

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Mol	Chain	Res	Type
1	A	47	ASN
1	A	79	HIS
1	A	81	GLN
1	A	658	ASN
1	A	718	ASN
1	A	792	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 [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	834/890 (93%)	-0.31	6 (0%) 87 67	24, 60, 99, 147	0
2	B	19/37 (51%)	-0.45	0 100 100	35, 55, 83, 98	0
All	All	853/927 (92%)	-0.31	6 (0%) 87 67	24, 60, 98, 147	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	791	GLN	2.9
1	A	371	SER	2.9
1	A	792	GLN	2.7
1	A	167	VAL	2.5
1	A	752	HIS	2.4
1	A	168	LEU	2.1

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.