Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:02 PM GMT

PDB ID	:	4BBN
Title	:	NEDD4 HECT-Ub:Ub complex
Authors	:	Maspero, E.; Valentini, E.; Mari, S.; Cecatiello, V.; Polo, S.; Pasqualato, S.
Deposited on	:	2012-09-27
Resolution	:	2.51 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:

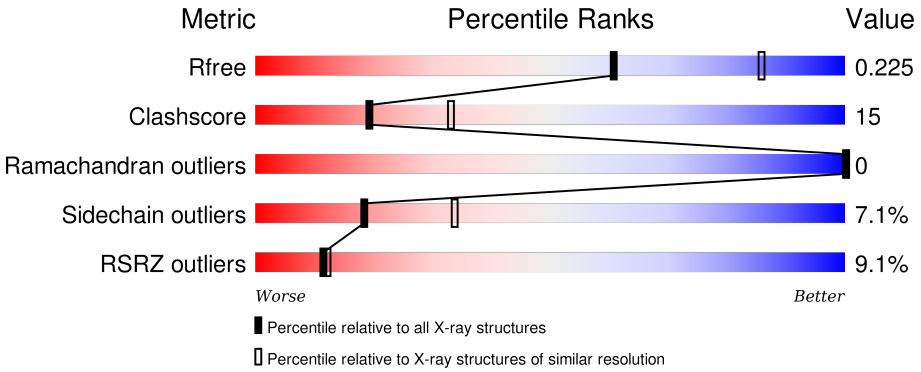
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	385	9% 73% 21%
2	C	76	14% 58% 37% 5%
3	F	76	% 87% 13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4492 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 E3 UBIQUITIN-PROTEIN LIGASE NEDD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	0	0
			3176	2054	538	570	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	515	GLY	-	EXPRESSION TAG	UNP P46934
A	516	PRO	-	EXPRESSION TAG	UNP P46934
A	517	LEU	-	EXPRESSION TAG	UNP P46934
A	518	GLY	-	EXPRESSION TAG	UNP P46934
A	626	ILE	LEU	ENGINEERED MUTATION	UNP P46934
A	.	-	CYS	DELETION	UNP P46934
A	628	PRO	GLU	ENGINEERED MUTATION	UNP P46934
A	777	SER	CYS	ENGINEERED MUTATION	UNP P46934

- Molecule 2 is a protein called POLYUBIQUITIN-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	76	Total	C	N	O	S	0	0	0
			602	378	105	118	1			

- Molecule 3 is a protein called POLYUBIQUITIN-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	76	Total	C	N	O	S	0	0	0
			604	379	105	118	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	76	CYS	GLY	ENGINEERED MUTATION	UNP P46934

- Molecule 4 is water.

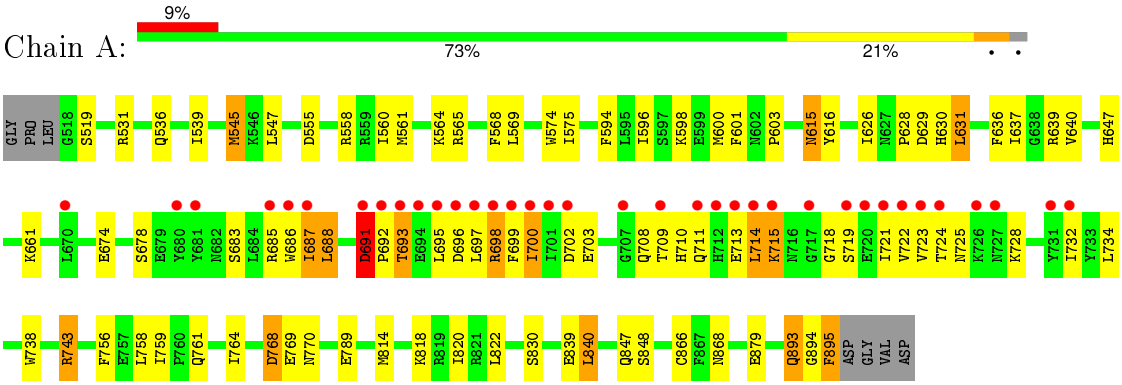
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	92	Total O 92 92	0	0
4	C	5	Total O 5 5	0	0
4	F	13	Total O 13 13	0	0

3 Residue-property plots (i)

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.

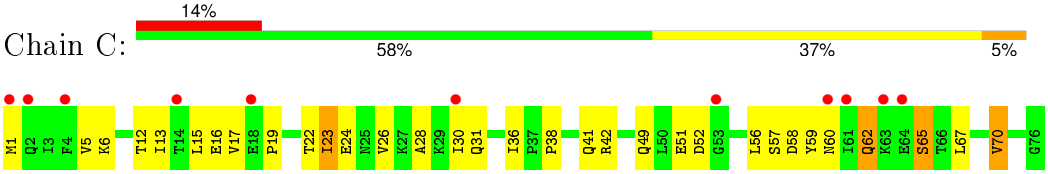
- Molecule 1: E3 UBIQUITIN-PROTEIN LIGASE NEDD4

Chain A:



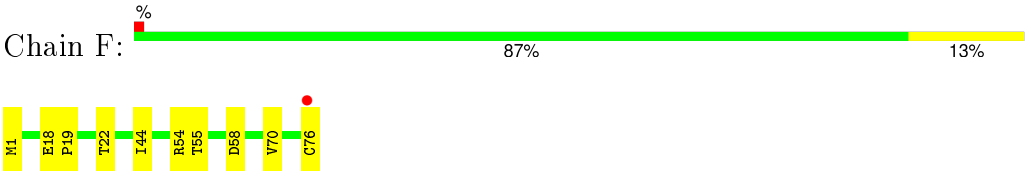
- Molecule 2: POLYUBIQUITIN-B

Chain C:



- Molecule 3: POLYUBIQUITIN-B

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	196.54Å 196.54Å 98.77Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	98.27 – 2.51 56.74 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.9 (98.27-2.51) 100.0 (56.74-2.51)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.25 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.176 , 0.224 0.177 , 0.225	Depositor DCC
R_{free} test set	1244 reflections (5.23%)	DCC
Wilson B-factor (Å^2)	47.3	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 41.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 25122 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4492	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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, 0.375 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	0.59	2/3263 (0.1%)	0.62	1/4397 (0.0%)
2	C	0.40	0/608	0.58	0/816
3	F	0.48	0/610	0.63	0/819
All	All	0.55	2/4481 (0.0%)	0.61	1/6032 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	574	TRP	CD2-CE2	5.18	1.47	1.41
1	A	738	TRP	CD2-CE2	5.12	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	691	ASP	C-N-CD	6.08	141.17	128.40

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	3176	0	3096	96	0
2	C	602	0	629	30	0
3	F	604	0	630	8	0
4	A	92	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	5	0	0	1	0
4	F	13	0	0	0	0
All	All	4492	0	4355	132	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 15.

All (132) 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)
2:C:23:ILE:CD1	2:C:52:ASP:HA	1.83	1.08
1:A:700:ILE:HD12	1:A:713:GLU:HG2	1.34	1.05
2:C:23:ILE:HD11	2:C:52:ASP:HA	1.06	1.02
2:C:23:ILE:HD11	2:C:52:ASP:CA	1.90	1.01
1:A:698:ARG:HG2	1:A:719:SER:O	1.67	0.95
2:C:1:MET:N	2:C:17:VAL:O	2.02	0.92
1:A:561:MET:HE1	1:A:639:ARG:HD2	1.54	0.89
1:A:866:CYS:HB2	3:F:76:CYS:OXT	1.73	0.89
2:C:19:PRO:HA	2:C:56:LEU:HD11	1.55	0.88
1:A:713:GLU:OE1	1:A:718:GLY:HA3	1.72	0.88
1:A:698:ARG:HH11	1:A:698:ARG:HG2	1.37	0.88
1:A:560:ILE:HG22	1:A:561:MET:HE3	1.57	0.85
1:A:561:MET:CE	1:A:639:ARG:HD2	2.07	0.85
1:A:699:PHE:CD2	1:A:715:LYS:HD3	2.10	0.85
1:A:661:LYS:HG2	1:A:769:GLU:HG2	1.59	0.84
2:C:59:TYR:O	2:C:60:ASN:ND2	2.11	0.82
1:A:696:ASP:O	1:A:698:ARG:HD2	1.80	0.82
1:A:560:ILE:HG22	1:A:561:MET:CE	2.10	0.80
1:A:894:GLY:O	1:A:895:PHE:CG	2.39	0.76
1:A:820:ILE:HG23	1:A:830:SER:HB2	1.67	0.76
1:A:895:PHE:C	1:A:895:PHE:HD1	1.91	0.75
1:A:895:PHE:HD1	1:A:895:PHE:O	1.70	0.74
1:A:713:GLU:OE1	1:A:718:GLY:CA	2.35	0.74
1:A:700:ILE:CD1	1:A:713:GLU:HG2	2.14	0.72
1:A:895:PHE:CD1	1:A:895:PHE:C	2.64	0.71
2:C:62:GLN:O	2:C:65:SER:HB2	1.91	0.70
1:A:688:LEU:HG	1:A:732:ILE:HD12	1.74	0.69
1:A:600:MET:HE1	1:A:640:VAL:HB	1.75	0.68
1:A:868:ASN:OD1	4:A:2073:HOH:O	2.10	0.68
2:C:23:ILE:HD11	2:C:51:GLU:O	1.94	0.68
1:A:895:PHE:CD1	1:A:895:PHE:O	2.47	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:721:ILE:H	1:A:721:ILE:HD12	1.59	0.67
1:A:714:LEU:HD23	1:A:734:LEU:HB2	1.76	0.67
1:A:714:LEU:HD12	1:A:714:LEU:N	2.09	0.67
1:A:536:GLN:OE1	1:A:565:ARG:NH2	2.28	0.66
1:A:714:LEU:CD1	1:A:714:LEU:N	2.60	0.65
1:A:691:ASP:CG	1:A:725:ASN:HD21	2.00	0.64
1:A:743:ARG:HG2	4:A:2059:HOH:O	1.96	0.64
1:A:698:ARG:NH1	1:A:698:ARG:HG2	2.10	0.64
3:F:18:GLU:CG	3:F:19:PRO:HD2	2.28	0.64
1:A:629:ASP:HB2	4:A:2041:HOH:O	1.97	0.64
3:F:54:ARG:NH2	3:F:58:ASP:O	2.30	0.64
1:A:893:GLN:HG3	1:A:895:PHE:CE1	2.32	0.63
2:C:23:ILE:CD1	2:C:51:GLU:O	2.47	0.63
1:A:691:ASP:OD1	1:A:725:ASN:ND2	2.29	0.63
2:C:15:LEU:HD21	2:C:30:ILE:HG13	1.82	0.61
2:C:42:ARG:HB2	2:C:70:VAL:HG13	1.83	0.60
1:A:555:ASP:OD1	1:A:558:ARG:NH2	2.35	0.60
1:A:661:LYS:CG	1:A:769:GLU:HG2	2.29	0.60
1:A:600:MET:HE1	1:A:637:ILE:HA	1.84	0.59
1:A:596:ILE:O	1:A:600:MET:HG3	2.02	0.58
1:A:697:LEU:O	1:A:722:VAL:HA	2.04	0.57
1:A:699:PHE:HD2	1:A:715:LYS:HD3	1.69	0.57
2:C:5:VAL:HG13	2:C:13:ILE:HG13	1.85	0.57
1:A:594:PHE:CZ	1:A:598:LYS:HE3	2.40	0.56
1:A:893:GLN:CG	1:A:895:PHE:CE1	2.89	0.56
1:A:814:MET:HE1	1:A:822:LEU:HB2	1.88	0.56
2:C:28:ALA:HA	2:C:31:GLN:HB3	1.89	0.55
3:F:18:GLU:HG3	3:F:19:PRO:HD2	1.87	0.55
1:A:600:MET:HE3	1:A:640:VAL:HG21	1.88	0.55
1:A:686:TRP:HE3	1:A:687:ILE:HG23	1.72	0.55
2:C:49:GLN:OE1	4:C:2004:HOH:O	2.18	0.54
1:A:600:MET:CE	1:A:637:ILE:HA	2.38	0.54
2:C:23:ILE:CG1	2:C:52:ASP:HA	2.38	0.53
1:A:713:GLU:CD	1:A:718:GLY:HA3	2.29	0.53
1:A:545:MET:HG3	1:A:575:ILE:HG12	1.89	0.53
1:A:703:GLU:OE2	1:A:710:HIS:ND1	2.34	0.53
1:A:691:ASP:CG	1:A:725:ASN:ND2	2.61	0.53
1:A:693:THR:HG23	1:A:724:THR:HG22	1.91	0.52
1:A:839:GLU:O	1:A:839:GLU:HG3	2.11	0.51
1:A:687:ILE:HG22	1:A:692:PRO:HB3	1.92	0.51
2:C:6:LYS:HE2	2:C:12:THR:HG22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:702:ASP:HB3	1:A:711:GLN:HG2	1.92	0.51
1:A:714:LEU:CD1	1:A:714:LEU:H	2.25	0.50
2:C:5:VAL:HA	2:C:67:LEU:O	2.12	0.50
1:A:561:MET:HE2	1:A:639:ARG:HD2	1.88	0.50
2:C:23:ILE:HD11	2:C:51:GLU:C	2.32	0.49
1:A:601:PHE:O	1:A:603:PRO:HD3	2.12	0.49
1:A:814:MET:HE2	1:A:818:LYS:C	2.33	0.49
2:C:19:PRO:HA	2:C:56:LEU:CD1	2.34	0.49
1:A:820:ILE:CG2	1:A:830:SER:HB2	2.39	0.49
1:A:893:GLN:CD	1:A:893:GLN:O	2.51	0.48
3:F:18:GLU:HG2	3:F:19:PRO:HD2	1.94	0.48
1:A:698:ARG:CG	1:A:698:ARG:NH1	2.73	0.47
1:A:626:ILE:O	1:A:628:PRO:HD3	2.13	0.47
1:A:687:ILE:CG2	1:A:692:PRO:HB3	2.44	0.47
1:A:698:ARG:HG3	1:A:721:ILE:O	2.14	0.47
1:A:539:ILE:HD12	1:A:568:PHE:HD1	1.80	0.47
1:A:686:TRP:CE3	1:A:687:ILE:HG23	2.50	0.46
1:A:866:CYS:CB	3:F:76:CYS:OXT	2.54	0.46
1:A:545:MET:HE2	1:A:547:LEU:HD21	1.98	0.46
1:A:702:ASP:N	1:A:702:ASP:OD1	2.44	0.46
2:C:23:ILE:CD1	2:C:51:GLU:C	2.84	0.46
3:F:22:THR:HA	3:F:55:THR:HA	1.98	0.46
2:C:24:GLU:HB2	2:C:52:ASP:OD2	2.16	0.46
1:A:713:GLU:HB3	1:A:718:GLY:HA2	1.98	0.46
1:A:569:LEU:O	1:A:647:HIS:HE1	1.99	0.46
2:C:16:GLU:OE1	2:C:16:GLU:N	2.37	0.46
1:A:840:LEU:HB3	1:A:847:GLN:O	2.16	0.45
1:A:839:GLU:HG2	4:A:2077:HOH:O	2.16	0.45
1:A:615:ASN:ND2	1:A:616:TYR:H	2.15	0.45
2:C:6:LYS:HE3	2:C:6:LYS:HB2	1.82	0.44
2:C:22:THR:O	2:C:26:VAL:HG23	2.18	0.44
1:A:768:ASP:OD1	1:A:770:ASN:N	2.47	0.44
1:A:769:GLU:CD	1:A:769:GLU:H	2.20	0.44
1:A:708:GLN:HG3	1:A:709:THR:H	1.81	0.44
1:A:761:GLN:O	1:A:764:ILE:HG22	2.18	0.44
2:C:38:PRO:HA	2:C:41:GLN:NE2	2.33	0.43
1:A:685:ARG:HA	1:A:685:ARG:HD2	1.90	0.43
1:A:692:PRO:HG2	1:A:723:VAL:HG12	2.00	0.43
1:A:696:ASP:O	1:A:698:ARG:CD	2.60	0.43
1:A:698:ARG:CZ	1:A:722:VAL:HG23	2.47	0.43
1:A:725:ASN:HA	1:A:728:LYS:HD3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:687:ILE:H	1:A:687:ILE:HG12	1.65	0.42
1:A:697:LEU:O	1:A:723:VAL:N	2.41	0.42
1:A:596:ILE:HA	1:A:596:ILE:HD12	1.91	0.42
1:A:698:ARG:HA	1:A:721:ILE:O	2.18	0.42
1:A:894:GLY:O	1:A:895:PHE:CB	2.67	0.42
2:C:23:ILE:HG13	2:C:52:ASP:OD1	2.20	0.42
1:A:691:ASP:OD2	1:A:725:ASN:ND2	2.54	0.41
2:C:36:ILE:O	2:C:41:GLN:NE2	2.52	0.41
3:F:44:ILE:HD12	3:F:70:VAL:CG2	2.51	0.41
2:C:23:ILE:CD1	2:C:52:ASP:CA	2.70	0.41
1:A:683:SER:O	1:A:687:ILE:HG12	2.20	0.41
1:A:761:GLN:NE2	4:A:2053:HOH:O	2.21	0.41
1:A:561:MET:HE1	1:A:636:PHE:HA	2.03	0.41
1:A:631:LEU:HD12	1:A:631:LEU:HA	1.82	0.41
2:C:23:ILE:HD11	2:C:52:ASP:N	2.33	0.41
1:A:697:LEU:HD23	1:A:697:LEU:HA	1.84	0.41
1:A:756:PHE:HA	1:A:759:ILE:O	2.19	0.41
2:C:24:GLU:HB2	2:C:52:ASP:CG	2.41	0.40
1:A:698:ARG:NH1	1:A:719:SER:O	2.52	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	376/385 (98%)	369 (98%)	7 (2%)	0	100	100
2	C	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
3	F	74/76 (97%)	72 (97%)	2 (3%)	0	100	100
All	All	524/537 (98%)	514 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	339/344 (98%)	312 (92%)	27 (8%)	15 28
2	C	68/68 (100%)	62 (91%)	6 (9%)	12 23
3	F	69/69 (100%)	68 (99%)	1 (1%)	74 91
All	All	476/481 (99%)	442 (93%)	34 (7%)	18 34

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	519	SER
1	A	531	ARG
1	A	545	MET
1	A	564	LYS
1	A	615	ASN
1	A	630	HIS
1	A	631	LEU
1	A	674	GLU
1	A	678	SER
1	A	687	ILE
1	A	688	LEU
1	A	691	ASP
1	A	693	THR
1	A	695	LEU
1	A	698	ARG
1	A	700	ILE
1	A	714	LEU
1	A	715	LYS
1	A	743	ARG
1	A	758	LEU
1	A	768	ASP
1	A	789	GLU
1	A	840	LEU
1	A	848	SER
1	A	879	GLU
1	A	893	GLN

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Mol	Chain	Res	Type
1	A	895	PHE
2	C	23	ILE
2	C	57	SER
2	C	58	ASP
2	C	62	GLN
2	C	65	SER
2	C	70	VAL
3	F	1	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	615	ASN
1	A	690	ASN
1	A	864	HIS
1	A	868	ASN
1	A	893	GLN
2	C	2	GLN
2	C	60	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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, 95^{th} 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	378/385 (98%)	0.33	36 (9%) 10 11	26, 44, 104, 147	0
2	C	76/76 (100%)	0.70	11 (14%) 3 3	41, 84, 115, 150	0
3	F	76/76 (100%)	-0.10	1 (1%) 79 82	29, 49, 74, 86	0
All	All	530/537 (98%)	0.33	48 (9%) 11 12	26, 47, 105, 150	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	695	LEU	12.0
1	A	723	VAL	7.9
1	A	699	PHE	6.0
1	A	722	VAL	5.2
1	A	702	ASP	5.2
1	A	697	LEU	5.1
1	A	700	ILE	4.7
1	A	692	PRO	4.6
1	A	686	TRP	4.4
2	C	4	PHE	4.3
1	A	698	ARG	4.1
1	A	694	GLU	3.9
1	A	720	GLU	3.9
2	C	64	GLU	3.8
1	A	696	ASP	3.8
1	A	717	GLY	3.6
1	A	724	THR	3.4
1	A	714	LEU	3.4
1	A	701	ILE	3.2
3	F	76	CYS	3.2
1	A	726	LYS	3.1
1	A	719	SER	3.0
1	A	707	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	731	TYR	3.0
1	A	713	GLU	3.0
2	C	2	GLN	2.9
1	A	727	ASN	2.9
1	A	732	ILE	2.7
1	A	687	ILE	2.7
2	C	1	MET	2.7
2	C	61	ILE	2.6
1	A	715	LYS	2.6
2	C	53	GLY	2.6
1	A	680	TYR	2.6
2	C	18	GLU	2.6
1	A	711	GLN	2.5
1	A	721	ILE	2.5
1	A	709	THR	2.4
1	A	670	LEU	2.4
1	A	693	THR	2.4
1	A	681	TYR	2.4
1	A	712	HIS	2.4
1	A	691	ASP	2.2
2	C	14	THR	2.2
2	C	30	ILE	2.1
2	C	60	ASN	2.1
1	A	685	ARG	2.1
2	C	63	LYS	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.