



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:37 AM GMT

PDB ID : 3FEY  
Title : Crystal structure of the CBC-importin alpha complex.  
Authors : Dias, S.M.G.; Ambrosio, A.L.B.; Cerione, R.A.  
Deposited on : 2008-12-01  
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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

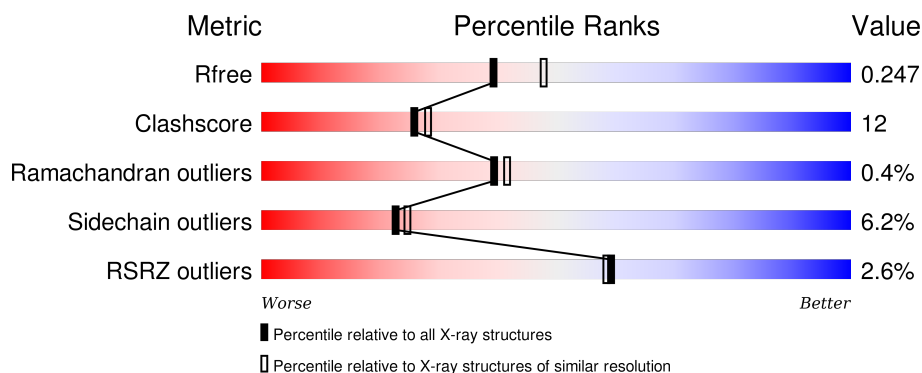
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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  $\geq 3$ , 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	790	<div> <div>2%</div> <div>74%</div> <div>19%</div> <div>• •</div> </div>
2	B	156	<div> <div>5%</div> <div>47%</div> <div>12%</div> <div>•</div> <div>38%</div> </div>
3	C	467	<div> <div>2%</div> <div>68%</div> <div>22%</div> <div>•</div> <div>8%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10746 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 Nuclear cap-binding protein subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	761	Total	C	N	O	S	0	3	0
			6242	4007	1067	1130	38			

- Molecule 2 is a protein called Nuclear cap-binding protein subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	97	Total	C	N	O	S	0	0	0
			792	501	129	157	5			

- Molecule 3 is a protein called Importin subunit alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	428	Total	C	N	O	S	0	1	0
			3258	2074	547	626	11			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	69	MET	-	EXPRESSION TAG	UNP P52292
C	530	HIS	-	EXPRESSION TAG	UNP P52292
C	531	HIS	-	EXPRESSION TAG	UNP P52292
C	532	HIS	-	EXPRESSION TAG	UNP P52292
C	533	HIS	-	EXPRESSION TAG	UNP P52292
C	534	HIS	-	EXPRESSION TAG	UNP P52292
C	535	HIS	-	EXPRESSION TAG	UNP P52292

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	282	Total	O	0	0
			282	282		

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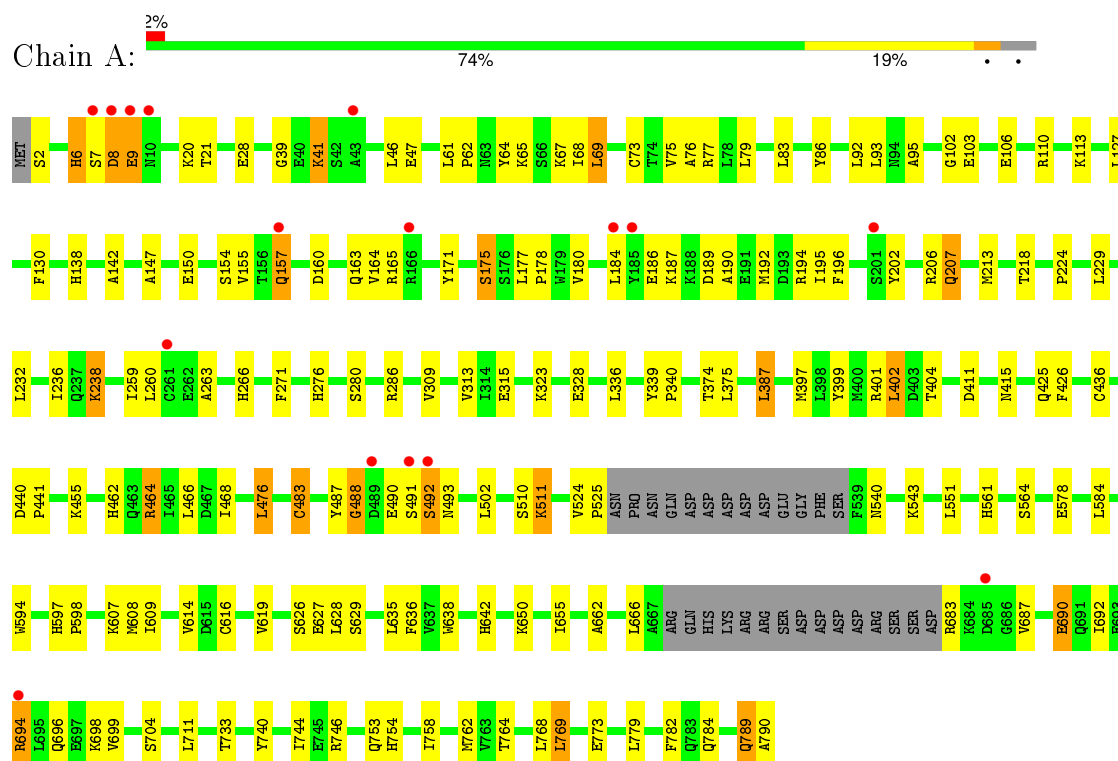
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	43	Total 43	O 43	0	0
4	C	129	Total 129	O 129	0	0

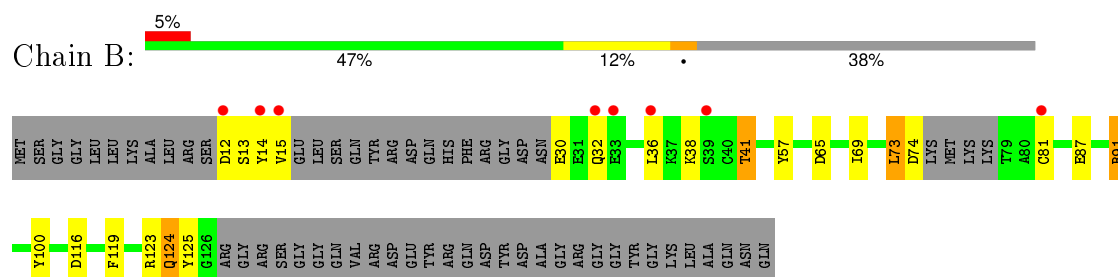
### 3 Residue-property plots [i](#)

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: Nuclear cap-binding protein subunit 1



- Molecule 2: Nuclear cap-binding protein subunit 2



- Molecule 3: Importin subunit alpha-2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.84Å 104.50Å 108.56Å 90.00° 109.04° 90.00°	Depositor
Resolution (Å)	15.00 – 2.20 14.98 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.2 (15.00-2.20) 97.2 (14.98-2.20)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.97 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, $R_{free}$	0.192 , 0.238 0.203 , 0.247	Depositor DCC
$R_{free}$ test set	4376 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.5	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 51.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 87017 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10746	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [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	0.52	2/6407 (0.0%)	0.61	0/8685
2	B	0.60	1/804 (0.1%)	0.66	0/1075
3	C	0.53	0/3317	0.61	0/4525
All	All	0.53	3/10528 (0.0%)	0.62	0/14285

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	483	CYS	CB-SG	-6.32	1.71	1.82
1	A	616	CYS	CB-SG	5.18	1.91	1.82
2	B	81	CYS	CB-SG	-5.07	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	8	ASP	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	6242	0	6221	136	0
2	B	792	0	756	20	0
3	C	3258	0	3323	108	0
4	A	282	0	0	12	0
4	B	43	0	0	4	0
4	C	129	0	0	13	0
All	All	10746	0	10300	255	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 12.

All (255) 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)
3:C:246:ILE:H	3:C:246:ILE:CD1	1.55	1.15
1:A:6[A]:HIS:NE2	3:C:354:GLU:OE2	1.83	1.11
3:C:219:SER:HB2	3:C:220:SER:CA	1.80	1.10
3:C:246:ILE:HD13	3:C:246:ILE:N	1.67	1.07
3:C:219:SER:CB	3:C:220:SER:HA	1.73	1.05
3:C:246:ILE:H	3:C:246:ILE:HD13	0.92	1.04
2:B:14:TYR:HE1	4:B:428:HOH:O	1.42	1.02
3:C:261:HIS:HE1	4:C:594:HOH:O	1.45	1.00
3:C:132:ASP:OD2	3:C:133:CYS:N	1.95	0.99
3:C:219:SER:HB2	3:C:220:SER:HA	0.99	0.96
1:A:662:ALA:CB	1:A:692:ILE:CD1	2.50	0.90
1:A:180:VAL:HG22	1:A:184:LEU:HD13	1.53	0.90
2:B:119:PHE:CE1	2:B:124:GLN:HG3	2.07	0.89
1:A:374:THR:HG21	2:B:100:TYR:O	1.75	0.87
3:C:435:LYS:HB2	4:C:583:HOH:O	1.77	0.84
3:C:207[B]:ASP:HB3	3:C:208:PRO:HD3	1.60	0.84
3:C:261:HIS:CE1	4:C:594:HOH:O	2.26	0.83
1:A:662:ALA:HB1	1:A:692:ILE:CD1	2.08	0.83
1:A:694:ARG:HH21	1:A:694:ARG:HA	1.43	0.82
1:A:490:GLU:O	1:A:492:SER:N	2.13	0.82
1:A:6[A]:HIS:CE1	3:C:354:GLU:OE2	2.33	0.82
3:C:292:THR:CG2	3:C:294:VAL:HG13	2.09	0.81
1:A:662:ALA:CB	1:A:692:ILE:HD11	2.08	0.81
1:A:662:ALA:HB1	1:A:692:ILE:HD11	1.63	0.80
1:A:142:ALA:HB3	1:A:187:LYS:HD2	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:411:GLU:H	3:C:411:GLU:CD	1.85	0.79
1:A:6[A]:HIS:HE2	3:C:354:GLU:CD	1.85	0.79
3:C:292:THR:HG22	3:C:294:VAL:H	1.48	0.79
1:A:69:LEU:HD21	1:A:103:GLU:HG2	1.64	0.79
1:A:77:ARG:HB3	1:A:127:LEU:HD21	1.64	0.79
1:A:690:GLU:O	1:A:694:ARG:HG2	1.81	0.79
3:C:290:VAL:HG12	3:C:291:LYS:HE2	1.65	0.79
3:C:287:GLY:O	3:C:291:LYS:HD2	1.83	0.78
3:C:89:ASN:HB2	3:C:92:ASN:OD1	1.84	0.78
2:B:14:TYR:CE1	4:B:428:HOH:O	2.23	0.77
2:B:119:PHE:CZ	2:B:124:GLN:HG3	2.18	0.77
1:A:627:GLU:O	4:A:985:HOH:O	2.02	0.77
1:A:662:ALA:HB3	1:A:692:ILE:CD1	2.14	0.76
1:A:399:TYR:HA	1:A:402:LEU:HD22	1.68	0.76
2:B:57:TYR:CD2	2:B:69:ILE:HD12	2.21	0.76
1:A:75:VAL:HG13	1:A:86:TYR:CD2	2.20	0.76
1:A:75:VAL:HG22	1:A:79:LEU:HD12	1.68	0.75
1:A:224:PRO:HD2	4:A:959:HOH:O	1.87	0.75
3:C:458:GLU:O	3:C:462:ILE:HG12	1.85	0.75
3:C:207[A]:ASP:HB2	3:C:208:PRO:HD3	1.66	0.75
3:C:134:SER:HB3	3:C:135:PRO:HD3	1.69	0.74
1:A:259:ILE:N	1:A:259:ILE:HD12	2.03	0.73
3:C:165:PRO:HG3	3:C:203:TYR:CE2	2.23	0.73
3:C:121:ILE:HB	3:C:122:PRO:HD3	1.70	0.73
3:C:395:LYS:HG2	3:C:436:ILE:HG13	1.70	0.72
1:A:483:CYS:HB3	1:A:594:TRP:CH2	2.24	0.72
3:C:292:THR:HG21	3:C:294:VAL:HG13	1.72	0.71
3:C:435:LYS:CB	4:C:583:HOH:O	2.37	0.71
3:C:291:LYS:HE3	4:C:620:HOH:O	1.91	0.70
3:C:352:GLN:O	3:C:356:THR:HG22	1.91	0.70
1:A:28:GLU:HB3	1:A:67:LYS:HE2	1.73	0.70
2:B:30:GLU:OE1	2:B:30:GLU:HA	1.92	0.70
1:A:387:LEU:O	1:A:387:LEU:HD13	1.91	0.70
1:A:510:SER:O	1:A:511:LYS:HB2	1.94	0.68
1:A:655:ILE:HD11	1:A:698:LYS:HE2	1.76	0.68
2:B:119:PHE:HE1	2:B:124:GLN:HG3	1.56	0.68
3:C:473:ILE:HG22	3:C:492:ILE:HD11	1.77	0.67
2:B:73:LEU:O	2:B:125:TYR:HA	1.94	0.67
1:A:315:GLU:OE1	4:A:972:HOH:O	2.13	0.66
1:A:608:MET:HE3	1:A:614:VAL:HG13	1.76	0.66
1:A:483:CYS:HB3	1:A:594:TRP:CZ3	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:753:GLN:HG2	4:A:980:HOH:O	1.96	0.66
1:A:154:SER:O	1:A:157:GLN:HB2	1.96	0.65
3:C:260:LEU:HD13	3:C:297:GLN:HG3	1.78	0.65
3:C:246:ILE:N	3:C:246:ILE:CD1	2.34	0.65
2:B:36:LEU:O	2:B:91:ARG:NH1	2.29	0.65
3:C:352:GLN:O	3:C:356:THR:CG2	2.45	0.65
3:C:101:ARG:HD3	3:C:139:GLU:OE1	1.98	0.64
1:A:171:TYR:O	1:A:175:SER:HB2	1.97	0.64
3:C:250:GLU:HA	3:C:250:GLU:OE1	1.97	0.63
1:A:662:ALA:CB	1:A:692:ILE:HD13	2.26	0.63
1:A:411:ASP:CG	1:A:635:LEU:HD11	2.19	0.62
2:B:116:ASP:OD1	2:B:123:ARG:NH2	2.31	0.62
3:C:257:VAL:HG22	3:C:294:VAL:CG1	2.30	0.61
3:C:499:GLU:C	4:C:597:HOH:O	2.38	0.61
1:A:399:TYR:HA	1:A:402:LEU:CD2	2.29	0.61
3:C:263:ASP:OD2	3:C:263:ASP:N	2.34	0.61
1:A:609:ILE:HD11	1:A:619:VAL:HG21	1.82	0.61
1:A:75:VAL:CG1	1:A:86:TYR:CD2	2.84	0.61
3:C:426:MET:HA	3:C:426:MET:HE2	1.83	0.60
1:A:259:ILE:H	1:A:259:ILE:HD12	1.64	0.60
2:B:119:PHE:CZ	2:B:124:GLN:CG	2.85	0.60
3:C:101:ARG:CD	3:C:139:GLU:OE1	2.49	0.60
1:A:157:GLN:HA	1:A:157:GLN:OE1	2.03	0.58
3:C:426:MET:HE1	3:C:444:ILE:HD11	1.85	0.58
1:A:662:ALA:HB1	1:A:692:ILE:HD13	1.81	0.58
3:C:129:GLY:O	3:C:131:THR:N	2.36	0.57
3:C:164:ILE:HB	3:C:165:PRO:HD3	1.85	0.57
1:A:180:VAL:CG2	1:A:184:LEU:HD13	2.32	0.57
3:C:257:VAL:HG22	3:C:294:VAL:HG13	1.87	0.57
1:A:286:ARG:NE	4:A:1035:HOH:O	2.17	0.57
2:B:57:TYR:CE2	2:B:69:ILE:HD12	2.40	0.57
1:A:7:SER:O	1:A:8:ASP:HB2	2.03	0.56
3:C:260:LEU:HD13	3:C:297:GLN:CG	2.36	0.55
1:A:61:LEU:N	1:A:62:PRO:HD2	2.21	0.55
1:A:41:LYS:HB2	1:A:41:LYS:NZ	2.20	0.55
3:C:477:GLN:HB2	4:C:543:HOH:O	2.05	0.55
1:A:662:ALA:C	1:A:692:ILE:HD11	2.27	0.55
1:A:138:HIS:CD2	4:A:1029:HOH:O	2.60	0.55
1:A:764:THR:HG23	1:A:768:LEU:HD12	1.87	0.54
3:C:411:GLU:HG2	4:C:550:HOH:O	2.07	0.54
3:C:307:LEU:N	3:C:308:PRO:CD	2.70	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:HIS:HD2	4:A:1029:HOH:O	1.89	0.54
3:C:401:VAL:HG21	3:C:421:ILE:HD11	1.90	0.53
1:A:157:GLN:OE1	1:A:157:GLN:CA	2.56	0.53
3:C:219:SER:CB	3:C:220:SER:CA	2.57	0.53
2:B:30:GLU:OE1	2:B:30:GLU:CA	2.57	0.53
3:C:101:ARG:O	3:C:101:ARG:HG2	2.07	0.53
1:A:426:PHE:HB3	1:A:468:ILE:HD12	1.90	0.53
1:A:476:LEU:HD23	1:A:476:LEU:N	2.23	0.53
3:C:79:ASP:N	3:C:79:ASP:OD1	2.41	0.52
1:A:6[A]:HIS:CD2	3:C:315:ARG:CZ	2.93	0.52
1:A:73:CYS:O	1:A:76:ALA:HB3	2.09	0.52
1:A:733:THR:HG21	4:A:1067:HOH:O	2.08	0.52
1:A:650:LYS:HE3	4:B:202:HOH:O	2.08	0.52
1:A:608:MET:CE	1:A:614:VAL:HG13	2.40	0.52
3:C:366:ARG:HB2	3:C:368:ASP:OD1	2.09	0.52
3:C:142:TRP:HD1	3:C:142:TRP:O	1.92	0.52
3:C:219:SER:HB2	3:C:220:SER:CB	2.39	0.52
3:C:175:HIS:HB3	3:C:177:HIS:CE1	2.44	0.52
1:A:102:GLY:N	1:A:266:HIS:CE1	2.79	0.51
1:A:655:ILE:CD1	1:A:698:LYS:HE2	2.40	0.51
1:A:177:LEU:N	1:A:178:PRO:CD	2.72	0.51
3:C:311:THR:HB	3:C:312:PRO:HD3	1.91	0.51
1:A:238:LYS:HD3	1:A:238:LYS:O	2.10	0.51
1:A:462:HIS:NE2	1:A:466:LEU:HD21	2.25	0.51
1:A:415:ASN:OD1	1:A:455:LYS:NZ	2.28	0.51
1:A:694:ARG:HH21	1:A:694:ARG:CA	2.20	0.51
1:A:259:ILE:H	1:A:259:ILE:CD1	2.23	0.51
1:A:259:ILE:N	1:A:259:ILE:CD1	2.71	0.51
3:C:311:THR:HB	3:C:312:PRO:CD	2.41	0.51
1:A:397:MET:O	1:A:401:ARG:HG3	2.12	0.50
1:A:655:ILE:HG22	1:A:699:VAL:HG22	1.93	0.50
1:A:789:GLN:O	1:A:789:GLN:HG3	2.12	0.50
3:C:477:GLN:HA	3:C:477:GLN:NE2	2.27	0.49
3:C:213:LEU:O	3:C:258:ARG:NH2	2.43	0.49
3:C:250:GLU:HG2	3:C:288:MET:CE	2.43	0.49
1:A:75:VAL:HG11	1:A:86:TYR:CG	2.48	0.49
1:A:189:ASP:OD1	1:A:190:ALA:N	2.45	0.49
1:A:147:ALA:O	1:A:150:GLU:HB2	2.12	0.49
2:B:119:PHE:HZ	2:B:124:GLN:CG	2.25	0.48
3:C:292:THR:HG22	3:C:294:VAL:N	2.23	0.48
3:C:94:LEU:HD12	3:C:94:LEU:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6[A]:HIS:CD2	3:C:315:ARG:NE	2.81	0.48
1:A:309:VAL:O	1:A:313:VAL:HG23	2.13	0.48
3:C:85:ILE:HD12	3:C:120:LEU:HD22	1.94	0.48
1:A:75:VAL:CG1	1:A:86:TYR:CG	2.96	0.48
1:A:102:GLY:CA	1:A:266:HIS:CE1	2.97	0.48
1:A:655:ILE:HG22	1:A:699:VAL:CG2	2.44	0.48
3:C:411:GLU:N	3:C:411:GLU:CD	2.58	0.48
3:C:426:MET:CE	3:C:444:ILE:HD11	2.44	0.48
1:A:510:SER:O	1:A:511:LYS:CB	2.60	0.47
3:C:250:GLU:OE1	3:C:250:GLU:CA	2.62	0.47
3:C:499:GLU:O	4:C:597:HOH:O	2.20	0.47
1:A:387:LEU:C	1:A:387:LEU:HD13	2.35	0.47
3:C:346:ASN:HA	4:C:618:HOH:O	2.15	0.47
1:A:106:GLU:OE1	1:A:110:ARG:NH1	2.48	0.47
1:A:155:VAL:C	1:A:157:GLN:H	2.18	0.47
3:C:273:TRP:CD2	3:C:312:PRO:HB3	2.50	0.47
1:A:607:LYS:NZ	4:A:1049:HOH:O	2.43	0.47
1:A:540:ASN:ND2	1:A:543:LYS:HE3	2.29	0.47
3:C:340:PHE:N	3:C:341:PRO:CD	2.78	0.47
1:A:462:HIS:CE1	1:A:466:LEU:CD2	2.98	0.46
3:C:307:LEU:N	3:C:308:PRO:HD2	2.30	0.46
1:A:769:LEU:N	1:A:769:LEU:CD2	2.78	0.46
1:A:740:TYR:O	1:A:744:ILE:HG22	2.15	0.46
3:C:101:ARG:HD2	3:C:139:GLU:OE1	2.16	0.46
1:A:64:TYR:O	1:A:68:ILE:HG13	2.15	0.46
1:A:164:VAL:HG13	1:A:790:ALA:OXT	2.16	0.46
1:A:694:ARG:NH2	1:A:694:ARG:HA	2.23	0.46
1:A:524:VAL:HG13	1:A:525:PRO:HD2	1.98	0.46
1:A:683:ARG:O	1:A:687:VAL:HG23	2.15	0.46
1:A:232:LEU:O	1:A:236:ILE:HG12	2.16	0.46
1:A:218:THR:OG1	1:A:404:THR:HB	2.15	0.46
3:C:207[A]:ASP:OD2	3:C:251:GLN:NE2	2.39	0.45
1:A:75:VAL:HG13	1:A:86:TYR:CE2	2.51	0.45
1:A:6[A]:HIS:HD2	3:C:315:ARG:NE	2.13	0.45
3:C:448:PHE:O	3:C:452:GLU:HG3	2.17	0.45
1:A:638:TRP:O	1:A:642:HIS:HD2	1.99	0.45
1:A:83:LEU:HD11	1:A:130:PHE:HA	1.98	0.45
1:A:165:ARG:HD2	1:A:276:HIS:HB2	1.99	0.45
1:A:95:ALA:HA	1:A:263:ALA:HB2	1.97	0.45
3:C:199:LEU:O	3:C:203:TYR:HD1	2.00	0.44
3:C:85:ILE:CD1	3:C:120:LEU:HD22	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:282:PRO:HB2	3:C:284:GLU:OE2	2.16	0.44
1:A:476:LEU:HD22	4:A:1063:HOH:O	2.17	0.44
3:C:435:LYS:O	3:C:439:VAL:HG23	2.17	0.44
1:A:61:LEU:HD22	1:A:65:LYS:HG3	2.00	0.44
1:A:106:GLU:CG	1:A:110:ARG:NH1	2.81	0.44
1:A:762:MET:HE1	1:A:782:PHE:CZ	2.52	0.44
1:A:411:ASP:CG	1:A:635:LEU:CD1	2.85	0.43
1:A:608:MET:CE	1:A:614:VAL:CG1	2.95	0.43
1:A:192:MET:HE3	1:A:196:PHE:CE1	2.53	0.43
1:A:150:GLU:HG3	1:A:195:ILE:HD11	1.99	0.43
1:A:75:VAL:HG11	1:A:86:TYR:CD1	2.52	0.43
1:A:487:TYR:O	1:A:488:GLY:O	2.37	0.43
3:C:289:VAL:O	3:C:292:THR:HB	2.19	0.43
3:C:426:MET:HE1	3:C:444:ILE:CD1	2.49	0.43
3:C:168:ILE:HD11	3:C:205:ALA:HB2	2.00	0.43
1:A:61:LEU:N	1:A:62:PRO:CD	2.81	0.43
3:C:115:ILE:HG23	3:C:120:LEU:HD12	2.01	0.43
1:A:754:HIS:O	1:A:758:ILE:HG13	2.19	0.43
1:A:323:LYS:NZ	4:A:862:HOH:O	2.37	0.43
1:A:425:GLN:O	1:A:426:PHE:HB2	2.17	0.43
1:A:762:MET:CE	1:A:782:PHE:HZ	2.31	0.43
3:C:306:GLU:HG3	4:C:572:HOH:O	2.19	0.43
1:A:339:TYR:HA	1:A:340:PRO:HD3	1.82	0.42
3:C:165:PRO:HG3	3:C:203:TYR:CD2	2.53	0.42
2:B:38:LYS:CG	4:B:203:HOH:O	2.67	0.42
1:A:155:VAL:C	1:A:157:GLN:N	2.72	0.42
3:C:337:LEU:HA	3:C:340:PHE:HD1	1.84	0.42
1:A:202:TYR:CE2	1:A:206:ARG:HD2	2.55	0.42
3:C:221:LEU:HD23	3:C:221:LEU:N	2.34	0.42
1:A:551:LEU:HA	1:A:551:LEU:HD12	1.74	0.42
2:B:116:ASP:CG	2:B:123:ARG:HH22	2.21	0.42
3:C:360:SER:HA	3:C:400:ALA:HA	2.01	0.42
1:A:150:GLU:CG	1:A:195:ILE:HD11	2.48	0.42
1:A:229:LEU:HA	1:A:229:LEU:HD23	1.85	0.42
3:C:91:GLU:O	3:C:95:GLN:HG3	2.19	0.42
3:C:367:GLN:HG3	4:C:571:HOH:O	2.20	0.42
3:C:223:CYS:O	3:C:224:GLY:C	2.57	0.42
1:A:598:PRO:HB3	1:A:636:PHE:CG	2.55	0.42
1:A:113:LYS:HE3	1:A:271:PHE:CE2	2.55	0.42
3:C:485:TYR:OH	4:C:545:HOH:O	2.05	0.42
1:A:440:ASP:OD1	1:A:441:PRO:HD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:VAL:HG22	1:A:180:VAL:O	2.20	0.41
2:B:12:ASP:OD2	2:B:15:VAL:HB	2.20	0.41
3:C:473:ILE:CG2	3:C:492:ILE:HD11	2.47	0.41
1:A:77:ARG:HB3	1:A:127:LEU:CD2	2.44	0.41
3:C:170:LEU:HD13	3:C:178:ILE:HG23	2.03	0.41
1:A:328[A]:GLU:OE2	2:B:13:SER:HB3	2.20	0.41
3:C:373:VAL:HG13	3:C:378:LEU:HB2	2.02	0.41
3:C:250:GLU:HG2	3:C:288:MET:HE2	2.03	0.41
1:A:387:LEU:C	1:A:387:LEU:CD1	2.89	0.41
1:A:20:LYS:O	3:C:105:SER:HB3	2.20	0.41
3:C:110:PRO:HA	3:C:111:PRO:HD3	1.88	0.41
1:A:39:GLY:HA2	1:A:46:LEU:HD13	2.02	0.41
1:A:163:GLN:OE1	1:A:207:GLN:HG2	2.21	0.41
3:C:385:VAL:CG1	3:C:393:THR:HG22	2.51	0.40
3:C:130:ARG:CZ	3:C:130:ARG:HB3	2.51	0.40
1:A:464:ARG:NH2	4:A:827:HOH:O	2.53	0.40
3:C:134:SER:CB	3:C:135:PRO:HD3	2.46	0.40
2:B:41:THR:HG22	2:B:87:GLU:HG3	2.03	0.40
3:C:207[B]:ASP:HB3	3:C:208:PRO:CD	2.41	0.40
1:A:411:ASP:OD1	1:A:635:LEU:CD1	2.69	0.40
1:A:213:MET:HB2	1:A:784:GLN:HB3	2.03	0.40
1:A:426:PHE:HB3	1:A:468:ILE:CD1	2.51	0.40
3:C:346:ASN:OD1	3:C:347:PRO:HD2	2.22	0.40
3:C:296:PRO:HA	3:C:334:ALA:HB1	2.03	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	758/790 (96%)	740 (98%)	14 (2%)	4 (0%)	34 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	91/156 (58%)	91 (100%)	0	0	100	100
3	C	425/467 (91%)	416 (98%)	8 (2%)	1 (0%)	52	59
All	All	1274/1413 (90%)	1247 (98%)	22 (2%)	5 (0%)	39	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	491	SER
3	C	130	ARG
1	A	488	GLY
1	A	9	GLU
1	A	511	LYS

### 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	699/724 (96%)	652 (93%)	47 (7%)	20	21
2	B	84/130 (65%)	77 (92%)	7 (8%)	14	13
3	C	362/396 (91%)	344 (95%)	18 (5%)	30	35
All	All	1145/1250 (92%)	1073 (94%)	72 (6%)	23	24

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	6[A]	HIS
1	A	6[B]	HIS
1	A	9	GLU
1	A	21	THR
1	A	41	LYS
1	A	47	GLU
1	A	69	LEU
1	A	92	LEU

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Mol	Chain	Res	Type
1	A	93	LEU
1	A	157	GLN
1	A	160	ASP
1	A	175	SER
1	A	186	GLU
1	A	194	ARG
1	A	207	GLN
1	A	238	LYS
1	A	260	LEU
1	A	280	SER
1	A	336	LEU
1	A	375	LEU
1	A	387	LEU
1	A	402	LEU
1	A	436	CYS
1	A	464	ARG
1	A	476	LEU
1	A	492	SER
1	A	493	ASN
1	A	502	LEU
1	A	561	HIS
1	A	564	SER
1	A	578	GLU
1	A	584	LEU
1	A	597	HIS
1	A	626	SER
1	A	628	LEU
1	A	629	SER
1	A	666	LEU
1	A	690	GLU
1	A	694	ARG
1	A	696	GLN
1	A	704	SER
1	A	711	LEU
1	A	769	LEU
1	A	773	GLU
1	A	779	LEU
1	A	789	GLN
2	B	32	GLN
2	B	41	THR
2	B	65	ASP
2	B	73	LEU

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Mol	Chain	Res	Type
2	B	74	ASP
2	B	91	ARG
2	B	124	GLN
3	C	79	ASP
3	C	101	ARG
3	C	105	SER
3	C	221	LEU
3	C	225	TYR
3	C	246	ILE
3	C	247	ASP
3	C	256	LEU
3	C	260	LEU
3	C	263	ASP
3	C	291	LYS
3	C	292	THR
3	C	339	VAL
3	C	356	THR
3	C	392	LYS
3	C	434	THR
3	C	453	LYS
3	C	457	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	138	HIS
1	A	207	GLN
1	A	367	HIS
3	C	137	GLN
3	C	261	HIS

### 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 [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

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	761/790 (96%)	-0.16	16 (2%) 67 65	9, 23, 30, 44	41 (5%)
2	B	97/156 (62%)	0.34	8 (8%) 14 14	11, 17, 93, 111	5 (5%)
3	C	428/467 (91%)	-0.16	10 (2%) 64 63	10, 20, 35, 42	25 (5%)
All	All	1286/1413 (91%)	-0.12	34 (2%) 59 58	9, 22, 33, 111	71 (5%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	15	VAL	8.0
1	A	491	SER	6.8
2	B	14	TYR	6.5
2	B	39	SER	6.0
1	A	8	ASP	5.9
2	B	12	ASP	4.5
3	C	498	VAL	3.9
2	B	32	GLN	3.9
3	C	89	ASN	3.6
2	B	36	LEU	3.6
3	C	85	ILE	3.2
1	A	685	ASP	3.2
3	C	220	SER	3.2
2	B	81	CYS	3.1
3	C	129	GLY	2.9
1	A	489	ASP	2.8
3	C	499	GLU	2.8
2	B	33	GLU	2.6
1	A	7	SER	2.4
1	A	43	ALA	2.4
3	C	88	SER	2.3
1	A	694	ARG	2.3
1	A	166	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	157	GLN	2.3
3	C	72	GLY	2.3
3	C	95	GLN	2.3
3	C	90	VAL	2.2
1	A	261	CYS	2.2
1	A	201	SER	2.2
1	A	9	GLU	2.2
1	A	184	LEU	2.2
1	A	10	ASN	2.1
1	A	185	TYR	2.1
1	A	492	SER	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.