



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:49 AM GMT

PDB ID : 3JV5  
Title : Crystal structure of the dimerization domains p52 homodimer  
Authors : Vu, D.; Huang, D.B.; Ghosh, G.  
Deposited on : 2009-09-15  
Resolution : 2.65 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

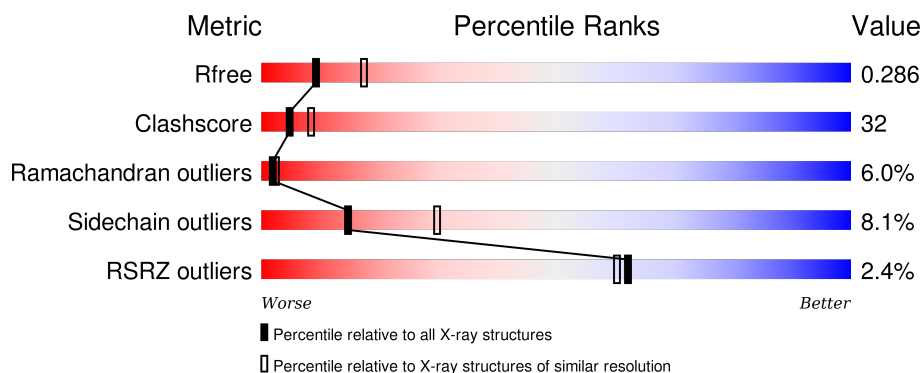
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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	104	<div> <div>2%</div> <div>51% 40% 7%</div> </div>
1	B	104	<div> <div>3%</div> <div>45% 44% 10%</div> </div>
1	C	104	<div> <div>%</div> <div>47% 38% 13%</div> </div>
1	D	104	<div> <div>4%</div> <div>41% 49% 8%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3562 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 factor NF-kappa-B p100 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	104	Total	C	N	O	S	0	0	0
			852	542	147	160	3			
1	B	103	Total	C	N	O	S	0	0	0
			847	539	146	159	3			
1	C	102	Total	C	N	O	S	0	0	0
			841	536	145	157	3			
1	D	102	Total	C	N	O	S	0	0	0
			839	533	145	158	3			

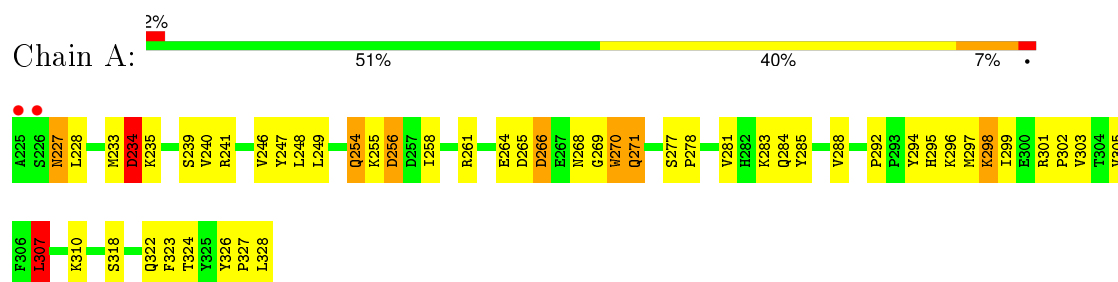
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	54	Total	O	0	0
			54	54		
2	B	39	Total	O	0	0
			39	39		
2	C	44	Total	O	0	0
			44	44		
2	D	46	Total	O	0	0
			46	46		

### 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.

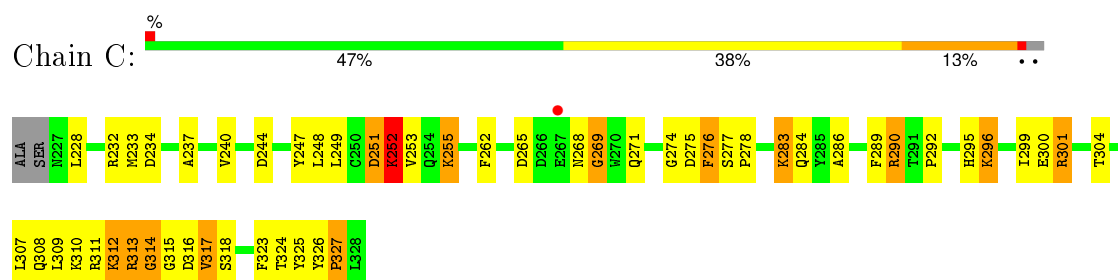
- Molecule 1: Nuclear factor NF-kappa-B p100 subunit



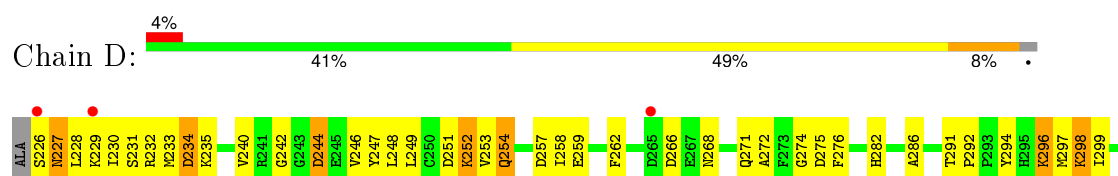
- Molecule 1: Nuclear factor NF-kappa-B p100 subunit

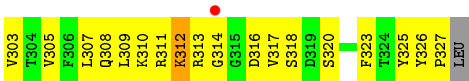


- Molecule 1: Nuclear factor NF-kappa-B p100 subunit



- Molecule 1: Nuclear factor NF-kappa-B p100 subunit





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.43 Å   56.86 Å   102.19 Å 90.00°   97.77°   90.00°	Depositor
Resolution (Å)	19.71 – 2.65 19.71 – 2.65	Depositor EDS
% Data completeness (in resolution range)	75.9 (19.71-2.65) 86.9 (19.71-2.65)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.99	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.13 (at 2.63 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.209   ,   0.268 0.231   ,   0.286	Depositor DCC
$R_{free}$ test set	943 reflections (6.92%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.2	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 62.3	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	1 of 14651 reflections (0.007%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3562	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% 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	0.45	0/872	0.73	1/1174 (0.1%)
1	B	0.43	0/867	0.72	0/1167
1	C	0.43	0/861	0.71	0/1159
1	D	0.36	0/859	0.61	0/1156
All	All	0.42	0/3459	0.70	1/4656 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	307	LEU	CA-CB-CG	5.13	127.09	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	852	0	833	50	0
1	B	847	0	828	59	0
1	C	841	0	823	52	0
1	D	839	0	817	58	0
2	A	54	0	0	3	0
2	B	39	0	0	1	0
2	C	44	0	0	1	0
2	D	46	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3562	0	3301	211	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 32.

All (211) 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:313:ARG:H	1:B:313:ARG:HD2	1.36	0.91
1:C:313:ARG:H	1:C:313:ARG:HD2	1.38	0.89
1:D:229:LYS:HE3	1:D:251:ASP:OD2	1.72	0.88
1:C:307:LEU:HD23	1:C:323:PHE:HB2	1.55	0.88
1:D:299:ILE:HG13	1:D:327:PRO:HG3	1.55	0.85
1:A:301:ARG:HB3	1:A:302:PRO:HD2	1.64	0.78
1:D:240:VAL:HG13	1:D:294:TYR:HB3	1.66	0.78
1:B:264:GLU:HG3	1:B:270:TRP:HE3	1.50	0.77
1:D:228:LEU:HB3	1:D:318:SER:HB3	1.66	0.76
1:A:254:GLN:HB3	2:A:450:HOH:O	1.83	0.76
1:B:234:ASP:HB2	1:B:247:TYR:H	1.51	0.74
1:B:234:ASP:CB	1:B:247:TYR:H	2.00	0.74
1:D:308:GLN:NE2	1:D:317:VAL:HB	2.03	0.72
1:A:277:SER:HB2	1:A:278:PRO:HD2	1.73	0.71
1:B:295:HIS:CD2	1:B:296:LYS:HG2	2.27	0.69
1:B:307:LEU:HD13	1:B:307:LEU:O	1.94	0.68
1:A:234:ASP:HB2	1:A:247:TYR:H	1.59	0.67
1:B:311:ARG:HB2	1:B:314:GLY:HA3	1.76	0.66
1:B:264:GLU:HG3	1:B:270:TRP:CE3	2.30	0.66
1:D:240:VAL:HG23	1:D:326:TYR:O	1.96	0.65
1:C:304:THR:HG22	1:C:324:THR:OG1	1.96	0.65
1:C:313:ARG:H	1:C:313:ARG:CD	2.03	0.65
1:D:240:VAL:HG22	1:D:325:TYR:HB3	1.79	0.64
1:D:251:ASP:O	1:D:253:VAL:HG13	1.98	0.64
1:A:239:SER:OG	1:A:241:ARG:HG2	1.98	0.63
1:B:313:ARG:HH11	1:B:313:ARG:HG3	1.63	0.63
1:A:270:TRP:CD1	1:A:295:HIS:HB3	2.33	0.63
1:C:251:ASP:O	1:C:253:VAL:HG13	1.98	0.63
1:A:298:LYS:HE3	2:A:504:HOH:O	1.97	0.63
1:D:297:MET:O	1:D:299:ILE:N	2.31	0.62
1:B:240:VAL:HG12	1:B:326:TYR:O	1.98	0.62
1:B:324:THR:HG1	1:B:326:TYR:HE1	1.44	0.62
1:B:313:ARG:N	1:B:313:ARG:HD2	2.12	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:VAL:HG13	1:D:294:TYR:CB	2.31	0.61
1:A:264:GLU:HG2	1:A:305:VAL:HG12	1.82	0.60
1:D:311:ARG:HH21	1:D:316:ASP:CB	2.15	0.60
1:C:311:ARG:HD3	1:C:316:ASP:HB3	1.83	0.60
1:C:275:ASP:O	1:C:276:PHE:HB3	2.02	0.60
1:D:240:VAL:CG1	1:D:294:TYR:HB3	2.32	0.59
1:A:281:VAL:O	1:A:281:VAL:HG12	2.00	0.59
1:C:255:LYS:HZ2	1:C:278:PRO:HB2	1.67	0.59
1:C:234:ASP:HB2	1:C:247:TYR:HB2	1.85	0.59
1:B:310:LYS:CB	1:B:317:VAL:HG12	2.33	0.59
1:C:299:ILE:HG13	1:C:327:PRO:HG3	1.84	0.59
1:D:233:MET:HE2	1:D:248:LEU:CD1	2.32	0.59
1:B:235:LYS:CG	1:B:246:VAL:HG22	2.33	0.58
1:B:246:VAL:HB	1:B:289:PHE:CE1	2.38	0.58
1:D:257:ASP:OD2	1:D:311:ARG:HB3	2.03	0.58
1:A:268:ASN:O	1:A:270:TRP:N	2.36	0.58
1:C:255:LYS:NZ	1:C:278:PRO:HB2	2.20	0.57
1:B:315:GLY:O	1:C:317:VAL:HG21	2.04	0.57
1:B:248:LEU:HD23	1:B:248:LEU:C	2.25	0.57
1:B:277:SER:HB2	1:B:278:PRO:CD	2.35	0.57
1:C:295:HIS:HE1	1:C:296:LYS:HE2	1.70	0.57
1:B:275:ASP:O	1:B:276:PHE:HB3	2.05	0.56
1:C:309:LEU:HB2	1:C:318:SER:HB3	1.87	0.56
1:D:274:GLY:HA2	1:D:292:PRO:HD3	1.87	0.56
1:A:322:GLN:HG3	2:A:432:HOH:O	2.04	0.56
1:A:248:LEU:HD23	1:A:248:LEU:C	2.25	0.56
1:A:299:ILE:HD12	1:A:303:VAL:HG23	1.86	0.56
1:C:290:ARG:HH11	1:C:290:ARG:HG2	1.71	0.56
1:D:234:ASP:HB2	1:D:247:TYR:HB2	1.87	0.56
1:C:249:LEU:HD13	1:D:282:HIS:HB2	1.88	0.55
1:C:296:LYS:HB3	1:C:296:LYS:NZ	2.21	0.55
1:D:272:ALA:HB1	1:D:292:PRO:HB2	1.89	0.55
1:B:240:VAL:HG23	1:B:294:TYR:HB3	1.89	0.55
1:B:272:ALA:HB1	1:B:292:PRO:HB2	1.89	0.55
1:C:313:ARG:HD2	1:C:313:ARG:N	2.13	0.55
1:A:228:LEU:HB3	1:A:318:SER:HB3	1.88	0.55
1:A:307:LEU:O	1:A:307:LEU:HD13	2.07	0.54
1:D:305:VAL:CG2	1:D:323:PHE:HB3	2.38	0.54
1:A:285:TYR:CD1	1:B:283:LYS:HG3	2.42	0.54
1:C:311:ARG:O	1:C:312:LYS:O	2.25	0.54
1:D:258:ILE:HD12	1:D:310:LYS:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234:ASP:HB2	1:B:247:TYR:HB2	1.90	0.54
1:B:304:THR:OG1	1:B:324:THR:HG22	2.08	0.54
1:D:229:LYS:HG2	1:D:251:ASP:HB2	1.91	0.53
1:A:307:LEU:O	1:A:307:LEU:HD22	2.08	0.53
1:A:299:ILE:HD12	1:A:301:ARG:O	2.09	0.53
1:D:233:MET:HE2	1:D:248:LEU:HD11	1.91	0.52
1:B:251:ASP:O	1:B:253:VAL:HG13	2.08	0.52
1:A:283:LYS:O	1:A:284:GLN:HB2	2.08	0.52
1:C:295:HIS:CE1	1:C:296:LYS:HE2	2.45	0.52
1:A:301:ARG:HB3	1:A:302:PRO:CD	2.39	0.51
1:C:255:LYS:HZ2	1:C:255:LYS:HB2	1.76	0.51
1:D:282:HIS:HB3	1:D:286:ALA:HB3	1.92	0.51
1:A:254:GLN:NE2	1:A:254:GLN:N	2.58	0.51
1:A:240:VAL:HG12	1:A:328:LEU:HD21	1.91	0.51
1:D:307:LEU:O	1:D:320:SER:HA	2.11	0.51
1:B:240:VAL:HG21	1:B:297:MET:HA	1.93	0.51
1:B:235:LYS:HG2	1:B:246:VAL:HG22	1.92	0.50
1:A:298:LYS:HD3	1:A:298:LYS:N	2.26	0.50
1:D:235:LYS:CG	1:D:246:VAL:HG22	2.41	0.49
1:D:258:ILE:HG13	1:D:259:GLU:N	2.26	0.49
1:D:262:PHE:CZ	1:D:292:PRO:HD2	2.47	0.49
1:C:274:GLY:HA2	1:C:292:PRO:HD3	1.93	0.49
1:D:235:LYS:HG2	1:D:246:VAL:HG22	1.95	0.49
1:D:258:ILE:HD11	1:D:309:LEU:HD13	1.94	0.49
1:D:228:LEU:CB	1:D:318:SER:HB3	2.40	0.49
1:A:261:ARG:O	1:A:307:LEU:HA	2.13	0.49
1:C:275:ASP:O	1:C:289:PHE:HB2	2.12	0.49
1:A:255:LYS:HG3	1:A:256:ASP:H	1.78	0.49
1:B:277:SER:HB2	1:B:278:PRO:HD2	1.94	0.49
1:D:308:GLN:CD	1:D:317:VAL:HB	2.33	0.49
1:C:234:ASP:CB	1:C:247:TYR:H	2.26	0.48
1:D:232:ARG:H	1:D:249:LEU:HB2	1.77	0.48
1:D:291:THR:HG21	1:D:323:PHE:CZ	2.49	0.48
1:A:265:ASP:O	1:A:266:ASP:O	2.31	0.48
1:B:261:ARG:HB2	1:B:273:PHE:HE1	1.78	0.48
1:C:301:ARG:HG2	1:C:301:ARG:HH21	1.78	0.48
1:B:237:ALA:HA	1:B:324:THR:O	2.13	0.48
1:B:313:ARG:NH1	1:B:313:ARG:HG3	2.29	0.48
1:D:307:LEU:HD23	1:D:307:LEU:N	2.27	0.48
1:B:262:PHE:O	1:B:271:GLN:HA	2.13	0.48
1:D:313:ARG:HG3	1:D:313:ARG:NH1	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:LYS:NZ	1:C:255:LYS:HB2	2.29	0.48
1:A:288:VAL:HG11	1:B:232:ARG:NH2	2.29	0.48
1:C:237:ALA:HA	1:C:324:THR:O	2.14	0.47
1:A:228:LEU:CB	1:A:318:SER:HB3	2.43	0.47
1:B:310:LYS:HB3	1:B:317:VAL:HG12	1.97	0.47
1:C:312:LYS:C	1:C:314:GLY:N	2.68	0.47
1:D:232:ARG:O	1:D:248:LEU:HD12	2.15	0.47
1:B:227:ASN:HB3	1:B:316:ASP:OD1	2.14	0.47
1:B:234:ASP:HB2	1:B:247:TYR:N	2.22	0.47
1:D:244:ASP:O	1:D:246:VAL:HG23	2.14	0.47
1:B:311:ARG:O	1:B:312:LYS:O	2.32	0.47
1:B:307:LEU:C	1:B:307:LEU:HD22	2.35	0.47
1:B:301:ARG:HH11	1:B:301:ARG:HG3	1.80	0.47
1:C:317:VAL:HG23	1:C:318:SER:O	2.16	0.46
1:A:240:VAL:HG13	1:A:297:MET:HE1	1.97	0.46
1:A:233:MET:HE1	1:A:246:VAL:HG11	1.95	0.46
1:A:249:LEU:HB3	1:B:282:HIS:CD2	2.50	0.46
1:D:251:ASP:O	1:D:252:LYS:C	2.53	0.46
1:D:299:ILE:HD11	1:D:327:PRO:N	2.30	0.46
1:A:258:ILE:HD12	1:A:310:LYS:O	2.15	0.46
1:C:315:GLY:O	2:C:404:HOH:O	2.21	0.46
1:B:228:LEU:HD12	1:B:317:VAL:HA	1.98	0.46
1:A:233:MET:CE	1:A:246:VAL:HG11	2.46	0.46
1:B:301:ARG:HD3	2:B:607:HOH:O	2.16	0.45
1:B:301:ARG:HG3	1:B:301:ARG:NH1	2.31	0.45
1:C:252:LYS:HZ1	1:C:284:GLN:CD	2.20	0.45
1:A:255:LYS:HA	1:A:281:VAL:HG11	1.99	0.45
1:B:248:LEU:O	1:B:286:ALA:HA	2.17	0.45
1:C:296:LYS:HB3	1:C:296:LYS:HZ3	1.81	0.45
1:A:261:ARG:HE	1:A:271:GLN:CG	2.30	0.45
1:D:258:ILE:HD11	1:D:309:LEU:HB3	1.98	0.45
1:B:231:SER:N	1:B:249:LEU:O	2.48	0.45
1:D:266:ASP:HB2	1:D:268:ASN:OD1	2.17	0.45
1:D:305:VAL:HG23	1:D:323:PHE:HB3	1.98	0.45
1:C:240:VAL:HA	1:C:325:TYR:CD1	2.52	0.44
1:C:233:MET:CE	1:C:307:LEU:HD21	2.48	0.44
1:A:234:ASP:CB	1:A:247:TYR:H	2.29	0.44
1:B:310:LYS:HB2	1:B:317:VAL:HG12	1.99	0.44
1:C:283:LYS:O	1:C:284:GLN:HB2	2.17	0.44
1:D:294:TYR:OH	1:D:303:VAL:HG11	2.18	0.44
1:B:313:ARG:O	1:B:314:GLY:O	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ASP:HB2	1:A:247:TYR:N	2.29	0.44
1:A:234:ASP:HB3	1:A:235:LYS:H	1.31	0.44
1:C:277:SER:HB2	1:C:278:PRO:HD2	2.00	0.44
1:B:299:ILE:O	1:B:299:ILE:HG13	2.19	0.43
1:C:228:LEU:HD23	1:C:228:LEU:HA	1.77	0.43
1:D:228:LEU:HD12	1:D:317:VAL:C	2.38	0.43
1:D:311:ARG:O	1:D:312:LYS:C	2.56	0.43
1:C:249:LEU:HD13	1:D:282:HIS:CB	2.48	0.43
1:B:301:ARG:HD2	1:B:301:ARG:H	1.82	0.43
1:C:262:PHE:O	1:C:271:GLN:HA	2.19	0.43
1:C:234:ASP:HB2	1:C:247:TYR:H	1.84	0.43
1:D:262:PHE:CE1	1:D:292:PRO:HD2	2.54	0.43
1:C:252:LYS:NZ	1:C:284:GLN:CD	2.72	0.43
1:B:315:GLY:HA3	1:C:308:GLN:HE22	1.84	0.43
1:C:244:ASP:O	1:C:290:ARG:HA	2.19	0.43
1:B:309:LEU:HB2	1:B:318:SER:HB3	2.01	0.43
1:A:248:LEU:HD23	1:A:248:LEU:O	2.20	0.42
1:B:317:VAL:HG21	1:C:315:GLY:O	2.19	0.42
1:C:232:ARG:HD3	1:C:249:LEU:HD12	2.01	0.42
1:A:323:PHE:CD2	1:A:323:PHE:C	2.92	0.42
1:D:296:LYS:HE3	1:D:298:LYS:HB2	2.01	0.42
1:D:275:ASP:O	1:D:276:PHE:HB3	2.18	0.42
1:B:310:LYS:NZ	1:B:312:LYS:O	2.52	0.42
1:C:251:ASP:O	1:C:252:LYS:C	2.58	0.42
1:A:255:LYS:HG3	1:A:256:ASP:N	2.33	0.42
1:B:283:LYS:O	1:B:284:GLN:HB2	2.20	0.42
1:B:299:ILE:HG13	1:B:327:PRO:HG3	2.01	0.42
1:A:296:LYS:HE2	1:A:296:LYS:HB2	1.85	0.42
1:A:254:GLN:NE2	1:A:254:GLN:H	2.17	0.42
1:D:262:PHE:HB2	1:D:272:ALA:HB3	2.02	0.41
1:D:313:ARG:HH11	1:D:313:ARG:HG3	1.84	0.41
1:C:311:ARG:HB2	1:C:314:GLY:HA3	2.02	0.41
1:D:230:ILE:CG2	1:D:248:LEU:HD11	2.50	0.41
1:A:227:ASN:HA	1:A:227:ASN:HD22	1.63	0.41
1:B:282:HIS:HB3	1:B:286:ALA:HB3	2.02	0.41
1:B:261:ARG:HB2	1:B:273:PHE:CE1	2.55	0.41
1:C:326:TYR:HA	1:C:327:PRO:HD3	1.82	0.41
1:C:317:VAL:HG23	1:C:318:SER:N	2.36	0.41
1:A:323:PHE:CD2	1:A:324:THR:N	2.89	0.41
1:D:226:SER:O	1:D:227:ASN:C	2.59	0.41
1:D:233:MET:HE2	1:D:248:LEU:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:ALA:HB1	1:B:326:TYR:CE1	2.56	0.41
1:C:290:ARG:NH1	1:C:290:ARG:HG2	2.36	0.41
1:A:261:ARG:NE	1:A:271:GLN:OE1	2.54	0.41
1:A:326:TYR:HA	1:A:327:PRO:HD3	1.73	0.41
1:D:251:ASP:O	1:D:252:LYS:O	2.39	0.40
1:A:294:TYR:CZ	1:A:295:HIS:CE1	3.09	0.40
1:A:264:GLU:OE2	1:A:295:HIS:HE1	2.04	0.40
1:C:268:ASN:O	1:C:269:GLY:C	2.59	0.40
1:D:254:GLN:O	1:D:258:ILE:HG22	2.21	0.40
1:D:240:VAL:C	1:D:242:GLY:H	2.24	0.40
1:A:295:HIS:CD2	1:A:296:LYS:HE2	2.57	0.40
1:C:248:LEU:O	1:C:286:ALA:HA	2.21	0.40
1:B:251:ASP:O	1:B:252:LYS:C	2.58	0.40
1:D:244:ASP:O	1:D:291:THR:HG23	2.22	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	102/104 (98%)	82 (80%)	15 (15%)	5 (5%)	3	4
1	B	101/104 (97%)	90 (89%)	6 (6%)	5 (5%)	3	4
1	C	100/104 (96%)	89 (89%)	4 (4%)	7 (7%)	1	1
1	D	100/104 (96%)	82 (82%)	11 (11%)	7 (7%)	1	1
All	All	403/416 (97%)	343 (85%)	36 (9%)	24 (6%)	2	2

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	234	ASP

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Mol	Chain	Res	Type
1	A	266	ASP
1	A	269	GLY
1	B	312	LYS
1	B	314	GLY
1	B	327	PRO
1	C	269	GLY
1	C	312	LYS
1	C	327	PRO
1	D	234	ASP
1	D	252	LYS
1	D	298	LYS
1	D	312	LYS
1	C	283	LYS
1	C	314	GLY
1	D	227	ASN
1	D	231	SER
1	D	314	GLY
1	B	283	LYS
1	C	252	LYS
1	A	270	TRP
1	B	276	PHE
1	C	276	PHE
1	A	292	PRO

### 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	93/93 (100%)	86 (92%)	7 (8%)	17	35
1	B	93/93 (100%)	85 (91%)	8 (9%)	13	26
1	C	92/93 (99%)	81 (88%)	11 (12%)	6	13
1	D	92/93 (99%)	88 (96%)	4 (4%)	35	63
All	All	370/372 (100%)	340 (92%)	30 (8%)	15	30

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	234	ASP
1	A	254	GLN
1	A	256	ASP
1	A	271	GLN
1	A	298	LYS
1	A	307	LEU
1	B	240	VAL
1	B	251	ASP
1	B	267	GLU
1	B	268	ASN
1	B	290	ARG
1	B	301	ARG
1	B	307	LEU
1	B	313	ARG
1	C	251	ASP
1	C	252	LYS
1	C	255	LYS
1	C	265	ASP
1	C	290	ARG
1	C	296	LYS
1	C	300	GLU
1	C	301	ARG
1	C	310	LYS
1	C	313	ARG
1	C	317	VAL
1	D	244	ASP
1	D	254	GLN
1	D	271	GLN
1	D	296	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	284	GLN
1	A	295	HIS
1	B	227	ASN
1	B	295	HIS
1	B	308	GLN
1	B	322	GLN
1	C	227	ASN
1	C	295	HIS

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Mol	Chain	Res	Type
1	D	271	GLN
1	D	308	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, 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	104/104 (100%)	-0.16	2 (1%) 70 69	16, 32, 72, 116	0
1	B	103/104 (99%)	-0.17	3 (2%) 55 53	13, 31, 63, 135	0
1	C	102/104 (98%)	-0.19	1 (0%) 84 84	16, 41, 75, 87	0
1	D	102/104 (98%)	0.32	4 (3%) 43 41	18, 57, 112, 129	0
All	All	411/416 (98%)	-0.05	10 (2%) 62 60	13, 40, 86, 135	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	225	ALA	4.4
1	A	226	SER	3.8
1	D	226	SER	3.0
1	B	227	ASN	2.9
1	B	315	GLY	2.7
1	D	314	GLY	2.3
1	D	229	LYS	2.1
1	B	301	ARG	2.1
1	C	267	GLU	2.1
1	D	265	ASP	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.