



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:39 PM GMT

PDB ID : 1A3Q  
Title : HUMAN NF-KAPPA-B P52 BOUND TO DNA  
Authors : Cramer, P.; Larson, C.J.; Verdine, G.L.; Muller, C.W.  
Deposited on : 1998-01-23  
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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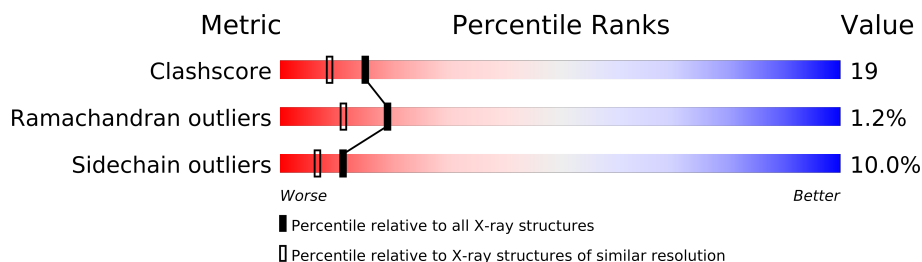
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.10 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))
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	11	
2	D	11	
3	A	285	
3	B	285	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5748 atoms, of which 0 are hydrogen and 0 are deuterium.

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 DNA chain called DNA (5'-D(\*GP\*GP\*GP\*GP\*AP\*AP\*TP\*CP\*CP\*CP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	11	Total	C	N	O	P	0	0	0
			223	106	44	63	10			

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*GP\*GP\*GP\*AP\*TP\*TP\*CP\*CP\*CP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	11	Total	C	N	O	P	0	0	0
			222	106	41	65	10			

- Molecule 3 is a protein called PROTEIN (NUCLEAR FACTOR KAPPA-B P52).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	285	Total	C	N	O	S	0	0	0
			2259	1423	409	415	12			
3	B	285	Total	C	N	O	S	0	0	0
			2259	1423	409	415	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	213	ILE	THR	CONFLICT	UNP Q00653
B	213	ILE	THR	CONFLICT	UNP Q00653

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	377	Total	O	0	0
			377	377		
4	B	335	Total	O	0	0
			335	335		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	32	Total 32	O 32	0	0
4	D	41	Total 41	O 41	0	0

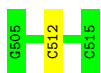
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

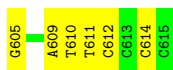
- Molecule 1: DNA (5'-D(\*GP\*GP\*GP\*GP\*AP\*AP\*TP\*CP\*CP\*CP\*C)-3')

Chain C: 



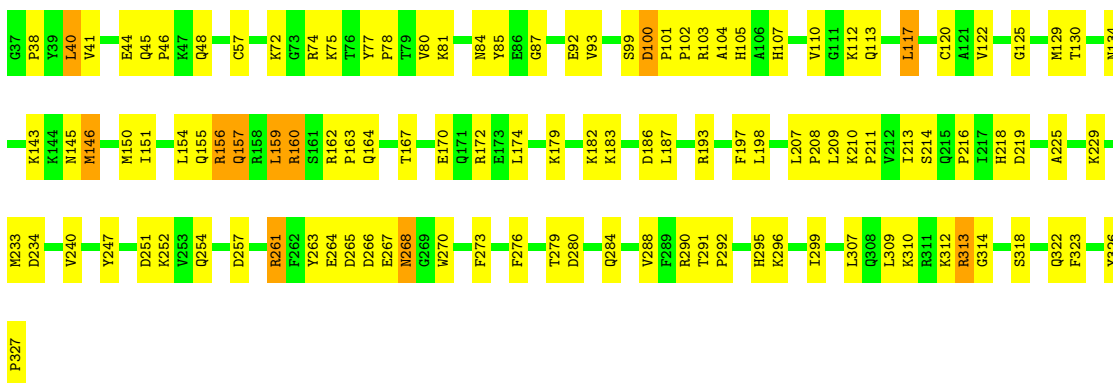
- Molecule 2: DNA (5'-D(\*GP\*GP\*GP\*GP\*AP\*TP\*TP\*CP\*CP\*CP\*C)-3')

Chain D: 



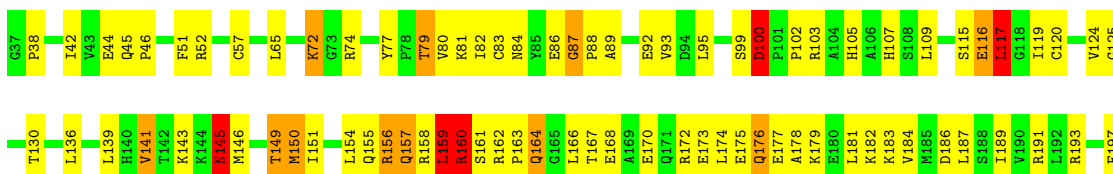
- Molecule 3: PROTEIN (NUCLEAR FACTOR KAPPA-B P52)

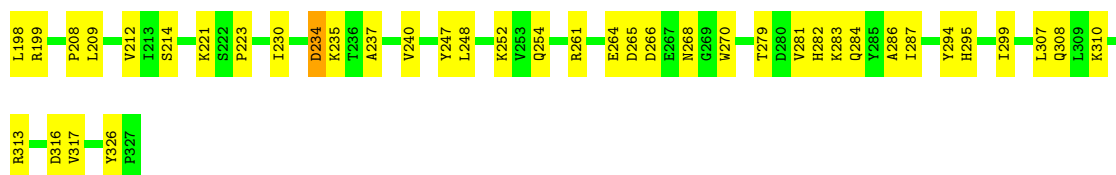
Chain A: 



- Molecule 3: PROTEIN (NUCLEAR FACTOR KAPPA-B P52)

Chain B: 





## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.20Å 121.00Å 134.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.10	Depositor
% Data completeness (in resolution range)	91.8 (10.00-2.10)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	6.10	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.219 , 0.320	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5748	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

## 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	C	0.86	0/250	1.04	0/384
2	D	0.89	0/248	1.13	2/381 (0.5%)
3	A	0.60	0/2307	0.89	2/3104 (0.1%)
3	B	0.63	0/2307	0.91	6/3104 (0.2%)
All	All	0.65	0/5112	0.92	10/6973 (0.1%)

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
2	D	0	1
3	A	0	1
All	All	0	2

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	164	GLN	N-CA-C	7.45	131.12	111.00
3	B	117	LEU	CB-CG-CD1	-7.38	98.46	111.00
3	B	164	GLN	N-CA-C	6.82	129.42	111.00
3	A	117	LEU	CB-CG-CD1	-6.44	100.05	111.00
2	D	605	DG	O5'-C5'-C4'	5.79	125.47	111.00
3	B	117	LEU	CB-CG-CD2	5.55	120.43	111.00
3	B	100	ASP	N-CA-C	5.50	125.85	111.00
2	D	609	DA	OP2-P-O3'	5.24	116.73	105.20
3	B	159	LEU	CA-CB-CG	5.15	127.14	115.30
3	B	117	LEU	CA-CB-CG	5.15	127.14	115.30

There are no chirality outliers.



All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	263	TYR	Sidechain
2	D	614	DC	Sidechain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	223	0	124	2	0
2	D	222	0	125	3	0
3	A	2259	0	2279	87	0
3	B	2259	0	2279	95	0
4	A	377	0	0	12	0
4	B	335	0	0	8	0
4	C	32	0	0	0	0
4	D	41	0	0	2	0
All	All	5748	0	4807	185	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (185) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:B:117:LEU:HD11	3:B:156:ARG:HD2	1.14	1.14
3:B:117:LEU:CD1	3:B:156:ARG:HD2	2.04	0.84
3:A:102:PRO:HG3	3:A:213:ILE:HD11	1.62	0.82
3:A:218:HIS:HB3	3:A:225:ALA:HB1	1.60	0.82
3:A:162:ARG:HB3	3:A:163:PRO:HD2	1.63	0.80
3:B:156:ARG:HD3	3:B:160:ARG:HH12	1.47	0.79
3:A:210:LYS:HG3	3:A:211:PRO:HD2	1.67	0.76
3:B:265:ASP:HB3	4:B:2644:HOH:O	1.87	0.75
3:A:45:GLN:OE1	3:A:214:SER:HB2	1.88	0.74
3:B:310:LYS:HG3	3:B:317:VAL:HG12	1.70	0.73
3:A:167:THR:HG22	3:A:170:GLU:HG3	1.71	0.73
3:B:167:THR:HG22	3:B:170:GLU:HG3	1.71	0.73
2:D:612:DC:H2'	3:B:57:CYS:SG	2.29	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:117:LEU:HD11	3:A:156:ARG:HD2	1.72	0.72
3:A:48:GLN:HG3	3:A:216:PRO:O	1.90	0.72
3:B:105:HIS:HD2	3:B:107:HIS:H	1.35	0.71
3:B:51:PHE:CD2	3:B:65:LEU:HD12	2.26	0.71
3:B:159:LEU:HD12	3:B:164:GLN:O	1.90	0.71
3:A:313:ARG:O	3:A:313:ARG:HG2	1.92	0.70
3:B:117:LEU:HG	3:B:156:ARG:NH1	2.07	0.69
3:B:230:ILE:HG23	3:B:248:LEU:CD1	2.23	0.69
3:A:103:ARG:HD3	4:A:2493:HOH:O	1.92	0.69
3:B:158:ARG:HH22	3:B:173:GLU:HG3	1.59	0.67
3:B:240:VAL:HG13	3:B:294:TYR:HB3	1.77	0.66
3:A:229:LYS:HE2	3:A:251:ASP:OD2	1.97	0.65
3:B:237:ALA:HB1	3:B:326:TYR:HE2	1.59	0.65
2:D:611:DT:H72	4:D:2637:HOH:O	1.97	0.65
3:A:117:LEU:HD11	3:A:156:ARG:NH1	2.11	0.64
3:A:113:GLN:NE2	3:A:122:VAL:HG12	2.12	0.64
3:A:167:THR:HG22	3:A:170:GLU:CG	2.29	0.63
3:B:313:ARG:HG3	4:B:2640:HOH:O	1.97	0.63
3:A:117:LEU:HD11	3:A:156:ARG:CZ	2.29	0.63
3:A:117:LEU:CD1	3:A:156:ARG:HD2	2.29	0.62
3:A:105:HIS:HD2	3:A:107:HIS:H	1.47	0.62
3:A:146:MET:SD	3:A:182:LYS:HG3	2.40	0.62
3:B:145:ASN:O	3:B:149:THR:HG23	1.99	0.62
3:A:264:GLU:OE2	3:A:295:HIS:HE1	1.82	0.62
3:B:197:PHE:CE2	3:B:208:PRO:HB3	2.34	0.62
3:B:157:GLN:O	3:B:160:ARG:HB2	1.99	0.62
3:A:151:ILE:O	3:A:155:GLN:HG3	2.00	0.61
3:B:237:ALA:HB1	3:B:326:TYR:CE2	2.35	0.61
3:A:117:LEU:HD11	3:A:156:ARG:CD	2.30	0.61
3:B:86:GLU:O	3:B:198:LEU:HD22	2.01	0.60
3:B:158:ARG:HH11	3:B:166:LEU:HD22	1.66	0.60
3:B:44:GLU:HB3	3:B:79:THR:HG23	1.84	0.59
3:A:266:ASP:HB2	3:A:268:ASN:OD1	2.02	0.59
3:A:75:LYS:HE2	3:A:77:TYR:CE1	2.38	0.59
3:A:103:ARG:HE	3:A:104:ALA:H	1.49	0.59
3:A:155:GLN:O	3:A:159:LEU:HD22	2.03	0.58
3:B:77:TYR:O	3:B:79:THR:HG22	2.03	0.58
3:A:240:VAL:HG21	3:A:299:ILE:HG12	1.85	0.58
3:B:158:ARG:C	3:B:160:ARG:H	2.06	0.58
3:A:160:ARG:HA	3:A:160:ARG:HE	1.69	0.58
3:B:42:ILE:HA	3:B:80:VAL:HG12	1.86	0.57
3:B:230:ILE:HG23	3:B:248:LEU:HD11	1.87	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:92:GLU:HA	3:A:120:CYS:O	2.03	0.57
3:B:240:VAL:HG23	3:B:326:TYR:O	2.05	0.57
3:A:107:HIS:CE1	3:A:187:LEU:HD22	2.40	0.56
3:A:75:LYS:HE2	3:A:77:TYR:HE1	1.71	0.56
3:B:141:VAL:HG22	3:B:145:ASN:HB3	1.88	0.55
3:B:156:ARG:O	3:B:160:ARG:NE	2.40	0.55
3:B:156:ARG:HD3	3:B:160:ARG:NH1	2.18	0.55
3:A:219:ASP:O	3:A:225:ALA:HB3	2.07	0.55
3:A:143:LYS:HE3	4:A:2019:HOH:O	2.07	0.55
3:B:92:GLU:HA	3:B:120:CYS:O	2.07	0.54
3:B:105:HIS:CD2	3:B:107:HIS:H	2.23	0.54
3:A:167:THR:HG23	3:A:170:GLU:H	1.73	0.54
3:B:197:PHE:HE2	3:B:208:PRO:HB3	1.71	0.54
3:A:45:GLN:OE1	3:A:46:PRO:HD2	2.08	0.53
3:B:282:HIS:HB3	3:B:286:ALA:HB3	1.88	0.53
3:A:198:LEU:HD21	3:A:209:LEU:HD11	1.89	0.53
1:C:512:DC:H2'	3:A:57:CYS:SG	2.49	0.53
3:B:155:GLN:O	3:B:159:LEU:HD22	2.09	0.52
3:A:261:ARG:HG3	3:A:273:PHE:HE1	1.74	0.52
3:A:261:ARG:HG3	3:A:273:PHE:CE1	2.44	0.52
3:A:162:ARG:CB	3:A:163:PRO:HD2	2.35	0.52
3:B:235:LYS:NZ	4:B:2417:HOH:O	2.43	0.52
3:B:52:ARG:HG3	3:B:221:LYS:HB2	1.91	0.52
3:A:322:GLN:HG3	4:A:2588:HOH:O	2.09	0.52
3:A:234:ASP:HB2	3:A:247:TYR:H	1.74	0.51
2:D:610:DT:H72	4:B:2066:HOH:O	2.10	0.51
3:B:151:ILE:CD1	3:B:175:GLU:HG2	2.41	0.51
3:A:167:THR:CG2	3:A:170:GLU:HG3	2.40	0.51
3:B:264:GLU:HG2	3:B:270:TRP:HE3	1.75	0.51
3:B:93:VAL:HA	3:B:193:ARG:O	2.10	0.51
3:B:155:GLN:HG2	3:B:166:LEU:HD11	1.92	0.51
3:A:117:LEU:CD1	3:A:156:ARG:NH1	2.74	0.50
3:A:81:LYS:HB2	3:A:130:THR:HG22	1.94	0.50
3:A:156:ARG:HA	3:A:159:LEU:HD21	1.92	0.50
3:B:81:LYS:HG3	3:B:130:THR:HG22	1.92	0.50
3:A:87:GLY:HA3	4:A:2554:HOH:O	2.11	0.50
3:B:223:PRO:HB2	3:B:254:GLN:NE2	2.27	0.50
3:B:117:LEU:CD1	3:B:156:ARG:HH11	2.24	0.50
3:A:112:LYS:HE3	4:A:2152:HOH:O	2.12	0.49
3:B:184:VAL:O	3:B:184:VAL:HG12	2.11	0.49
3:A:198:LEU:O	3:A:207:LEU:HB2	2.11	0.49
3:A:233:MET:HE1	3:A:307:LEU:HD21	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:264:GLU:HG2	3:A:270:TRP:HE3	1.76	0.49
3:A:257:ASP:HA	3:A:312:LYS:HE2	1.94	0.49
3:B:261:ARG:HD3	3:B:308:GLN:OE1	2.12	0.49
3:A:156:ARG:HA	3:A:159:LEU:CD2	2.43	0.49
3:B:117:LEU:CG	3:B:156:ARG:NH1	2.74	0.48
3:B:81:LYS:NZ	4:B:2639:HOH:O	2.46	0.48
3:A:103:ARG:NE	3:A:104:ALA:H	2.09	0.48
3:B:162:ARG:HB3	3:B:163:PRO:HD2	1.96	0.48
3:A:198:LEU:HD11	3:A:209:LEU:HD21	1.96	0.47
3:A:85:TYR:HB2	3:A:129:MET:CE	2.44	0.47
3:B:38:PRO:O	3:B:209:LEU:HD22	2.14	0.47
3:B:151:ILE:HD11	3:B:175:GLU:HG2	1.96	0.47
3:B:117:LEU:CG	3:B:156:ARG:HH11	2.27	0.47
3:A:213:ILE:HD12	4:A:2521:HOH:O	2.14	0.47
3:A:146:MET:CE	3:A:182:LYS:HG3	2.45	0.47
3:B:283:LYS:HE2	4:B:2453:HOH:O	2.14	0.47
3:B:87:GLY:O	3:B:88:PRO:C	2.53	0.47
3:B:157:GLN:O	3:B:160:ARG:CB	2.63	0.47
3:B:100:ASP:O	3:B:102:PRO:HD3	2.15	0.47
3:B:156:ARG:O	3:B:160:ARG:HD2	2.15	0.46
3:A:210:LYS:HD3	4:A:2479:HOH:O	2.15	0.46
3:A:252:LYS:HA	3:A:284:GLN:O	2.15	0.46
3:B:179:LYS:O	3:B:183:LYS:HG2	2.14	0.46
3:B:234:ASP:HB2	3:B:247:TYR:H	1.80	0.46
3:B:230:ILE:HG23	3:B:248:LEU:HD13	1.95	0.45
3:B:160:ARG:HE	3:B:160:ARG:HA	1.82	0.45
3:A:102:PRO:HB3	3:A:213:ILE:HD13	1.97	0.45
3:B:95:LEU:HG	3:B:109:LEU:HG	1.99	0.45
3:A:290:ARG:HG2	3:A:290:ARG:HH11	1.82	0.45
3:B:83:CYS:O	3:B:84:ASN:HB2	2.17	0.45
3:B:99:SER:O	3:B:191:ARG:NH2	2.49	0.45
3:B:264:GLU:OE2	3:B:295:HIS:HE1	2.00	0.44
3:B:45:GLN:OE1	3:B:46:PRO:HD2	2.17	0.44
3:A:276:PHE:HB2	3:A:280:ASP:HB2	1.99	0.44
3:B:156:ARG:HA	3:B:159:LEU:CD2	2.48	0.44
3:B:186:ASP:OD2	3:B:189:ILE:HD12	2.18	0.44
3:B:281:VAL:HG13	3:B:287:ILE:HG12	2.00	0.44
3:A:40:LEU:HG	3:A:41:VAL:N	2.32	0.44
3:A:159:LEU:HD13	4:A:2318:HOH:O	2.18	0.43
3:B:143:LYS:HD2	4:B:2502:HOH:O	2.18	0.43
3:B:176:GLN:C	3:B:178:ALA:H	2.21	0.43
3:B:42:ILE:HG13	3:B:212:VAL:HG11	2.01	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:309:LEU:HB2	3:A:318:SER:OG	2.18	0.43
3:A:146:MET:O	3:A:150:MET:HB2	2.18	0.43
3:B:45:GLN:OE1	3:B:214:SER:HB2	2.19	0.43
3:A:326:TYR:HA	3:A:327:PRO:HD3	1.79	0.43
3:A:179:LYS:O	3:A:183:LYS:HG2	2.19	0.43
3:A:100:ASP:HB3	3:A:101:PRO:HD3	2.00	0.43
3:B:146:MET:HE3	3:B:182:LYS:HG3	2.00	0.43
3:B:115:SER:O	3:B:117:LEU:N	2.52	0.42
3:B:146:MET:CE	3:B:182:LYS:HG3	2.49	0.42
3:B:158:ARG:C	3:B:160:ARG:N	2.72	0.42
3:B:234:ASP:CG	3:B:247:TYR:HB2	2.40	0.42
3:A:40:LEU:HD11	3:A:80:VAL:HB	2.00	0.42
3:B:156:ARG:O	3:B:160:ARG:CD	2.67	0.42
3:A:117:LEU:HD11	3:A:156:ARG:NE	2.33	0.42
3:B:310:LYS:HE3	4:B:2385:HOH:O	2.19	0.42
3:A:38:PRO:HG3	3:A:84:ASN:O	2.18	0.42
3:B:177:GLU:O	3:B:181:LEU:HB2	2.19	0.42
3:B:150:MET:HG3	3:B:181:LEU:HD23	2.00	0.42
3:A:291:THR:HG21	3:A:323:PHE:CZ	2.55	0.42
3:A:113:GLN:NE2	4:A:2152:HOH:O	2.50	0.42
1:C:512:DC:H5'	4:D:2010:HOH:O	2.20	0.42
3:B:89:ALA:HB3	3:B:124:VAL:HG23	2.02	0.41
3:A:197:PHE:CD1	3:A:208:PRO:HA	2.54	0.41
3:A:157:GLN:HB3	3:A:157:GLN:HE21	1.65	0.41
3:B:240:VAL:CG2	3:B:299:ILE:HD11	2.50	0.41
3:A:280:ASP:HB3	3:A:288:VAL:O	2.21	0.41
3:A:40:LEU:HD11	3:A:80:VAL:CG2	2.49	0.41
3:B:116:GLU:O	3:B:116:GLU:HG2	2.21	0.41
3:A:156:ARG:HG3	4:A:2470:HOH:O	2.20	0.41
3:A:38:PRO:O	3:A:209:LEU:HD22	2.20	0.41
3:B:160:ARG:HB3	3:B:161:SER:H	1.57	0.41
3:B:92:GLU:HG3	3:B:119:ILE:CG2	2.50	0.41
3:A:296:LYS:HE2	4:A:2171:HOH:O	2.19	0.41
3:B:115:SER:C	3:B:117:LEU:H	2.24	0.41
3:A:44:GLU:O	3:A:78:PRO:HA	2.21	0.41
3:B:72:LYS:O	3:B:72:LYS:HD2	2.21	0.41
3:A:273:PHE:O	3:A:292:PRO:HB3	2.21	0.41
3:A:261:ARG:HD2	3:A:310:LYS:HD2	2.02	0.41
3:B:146:MET:HE2	3:B:187:LEU:HD11	2.03	0.40
3:A:117:LEU:HD13	4:A:2423:HOH:O	2.21	0.40
3:A:93:VAL:HA	3:A:193:ARG:O	2.21	0.40
3:A:105:HIS:HE1	3:A:186:ASP:O	2.05	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:B:181:LEU:HA	3:B:181:LEU:HD12	1.87	0.40
3:B:82:ILE:HD12	3:B:124:VAL:HG21	2.03	0.40
3:B:252:LYS:HA	3:B:284:GLN:O	2.21	0.40
3:B:136:LEU:HA	3:B:136:LEU:HD23	1.89	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	
3	A	281/285 (99%)	250 (89%)	29 (10%)	2 (1%)	30	23
3	B	281/285 (99%)	252 (90%)	24 (8%)	5 (2%)	13	5
All	All	562/570 (99%)	502 (89%)	53 (9%)	7 (1%)	19	11

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	87	GLY
3	B	125	GLY
3	A	125	GLY
3	B	116	GLU
3	B	145	ASN
3	B	160	ARG
3	A	314	GLY

### 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	
3	A	249/249 (100%)	226 (91%)	23 (9%)	13	8
3	B	249/249 (100%)	222 (89%)	27 (11%)	9	5
All	All	498/498 (100%)	448 (90%)	50 (10%)	11	7

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	40	LEU
3	A	72	LYS
3	A	74	ARG
3	A	99	SER
3	A	100	ASP
3	A	110	VAL
3	A	134	ASN
3	A	145	ASN
3	A	146	MET
3	A	154	LEU
3	A	156	ARG
3	A	157	GLN
3	A	159	LEU
3	A	160	ARG
3	A	172	ARG
3	A	174	LEU
3	A	254	GLN
3	A	261	ARG
3	A	265	ASP
3	A	267	GLU
3	A	268	ASN
3	A	279	THR
3	A	313	ARG
3	B	72	LYS
3	B	74	ARG
3	B	79	THR
3	B	100	ASP
3	B	103	ARG
3	B	117	LEU
3	B	139	LEU
3	B	141	VAL
3	B	145	ASN
3	B	149	THR
3	B	150	MET
3	B	154	LEU

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Mol	Chain	Res	Type
3	B	156	ARG
3	B	157	GLN
3	B	159	LEU
3	B	160	ARG
3	B	168	GLU
3	B	172	ARG
3	B	174	LEU
3	B	176	GLN
3	B	199	ARG
3	B	234	ASP
3	B	266	ASP
3	B	268	ASN
3	B	279	THR
3	B	307	LEU
3	B	316	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	105	HIS
3	A	113	GLN
3	A	134	ASN
3	A	145	ASN
3	A	157	GLN
3	A	164	GLN
3	A	295	HIS
3	A	322	GLN
3	B	84	ASN
3	B	98	HIS
3	B	105	HIS
3	B	145	ASN
3	B	157	GLN
3	B	176	GLN
3	B	254	GLN
3	B	322	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.