



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 17, 2022 – 08:07 PM JST

PDB ID : 7W7L
Title : Structure of NF- κ B p52 homodimer bound to 13-mer A/T-centric P-Selectin
kB DNA fragment
Authors : Meshcheryakov, V.A.; Wang, V.Y.-F.
Deposited on : 2021-12-05
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

<https://www.wwpdb.org/validation/2017/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)	:	1.13
EDS	:	2.25
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.25

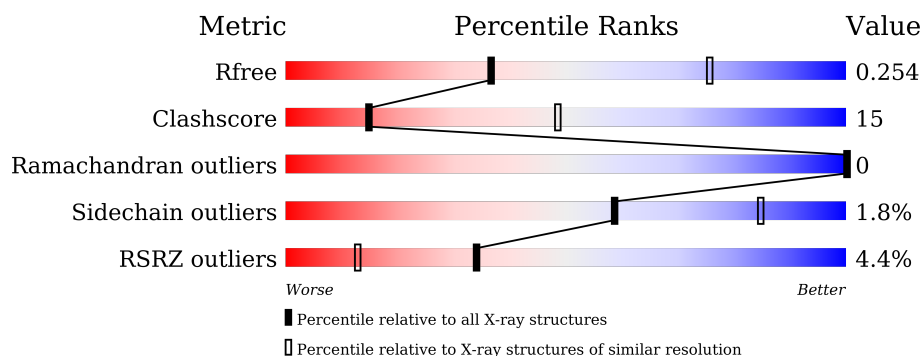
1 Overall quality at a glance ⓘ

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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	327	<div> <div>%</div> <div> <div></div> <div>68%</div> <div>21%</div> <div>10%</div> </div> </div>
1	B	327	<div> <div>7%</div> <div> <div></div> <div>66%</div> <div>23%</div> <div>10%</div> </div> </div>
2	C	13	<div> <div>38%</div> <div>46%</div> <div>15%</div> </div>
3	D	13	<div> <div>46%</div> <div>31%</div> <div>23%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5191 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 p52 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	0	0
			2321	1458	418	433	12			
1	B	294	Total	C	N	O	S	0	0	0
			2319	1458	418	431	12			

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*GP*GP*GP*GP*TP*AP*AP*CP*CP*CP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	P	0	0	0
			265	126	51	76	12			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	13	Total	C	N	O	P	0	0	0
			262	125	49	76	12			

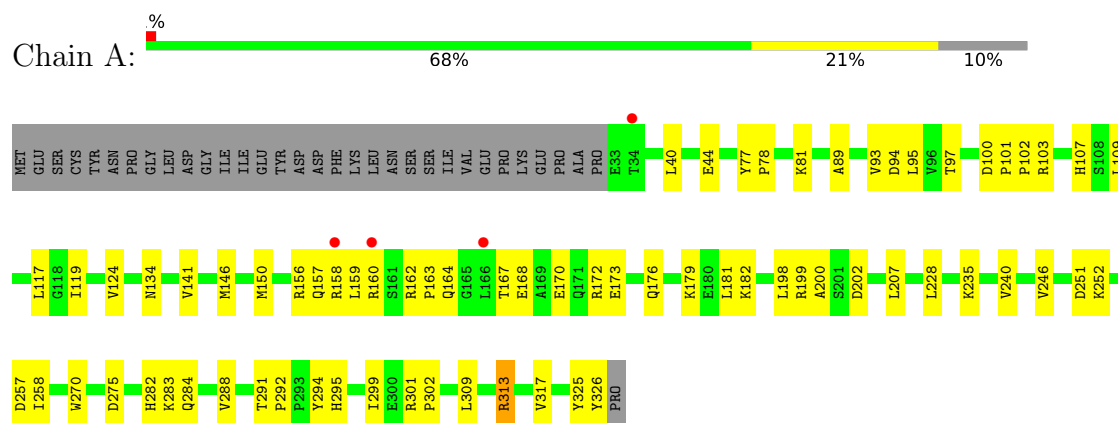
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	17	Total	O	0	0
			17	17		
4	C	1	Total	O	0	0
			1	1		
4	B	6	Total	O	0	0
			6	6		

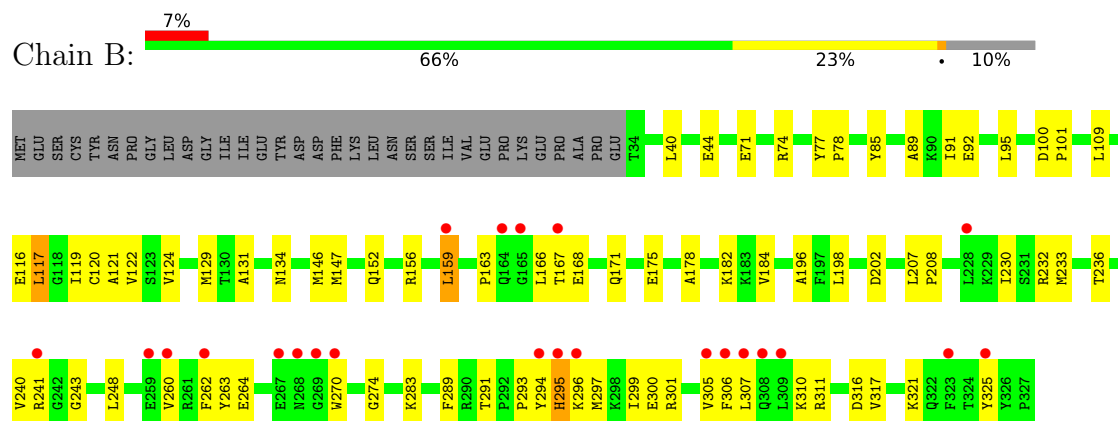
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

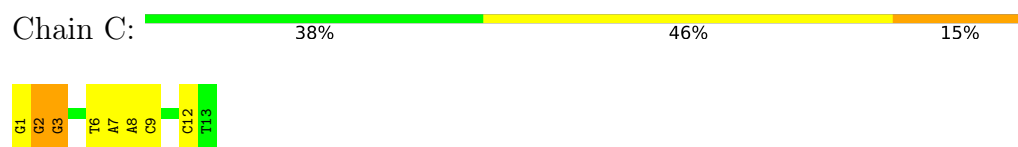
- Molecule 1: Nuclear factor NF-kappa-B p52 subunit



- Molecule 1: Nuclear factor NF-kappa-B p52 subunit

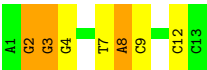


- Molecule 2: DNA (5'-D(*GP*GP*GP*GP*GP*TP*AP*AP*CP*CP*CP*CP*T)-3')



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





4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	225.28Å 225.28Å 96.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.77 – 3.00 48.77 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.77-3.00) 100.0 (48.77-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.238 , 0.258 0.240 , 0.254	Depositor DCC
R_{free} test set	1529 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å ²)	98.9	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 87.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5191	wwPDB-VP
Average B, all atoms (Å ²)	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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

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	0.67	0/2370	0.91	0/3190
1	B	0.67	0/2369	0.90	0/3190
2	C	0.95	0/297	1.64	6/457 (1.3%)
3	D	0.94	0/293	1.57	7/450 (1.6%)
All	All	0.70	0/5329	1.02	13/7287 (0.2%)

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

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	DG	P-O3'-C3'	-8.69	109.28	119.70
3	D	3	DG	P-O3'-C3'	-7.43	110.78	119.70
3	D	12	DC	P-O3'-C3'	-6.92	111.40	119.70
2	C	3	DG	P-O3'-C3'	-6.91	111.41	119.70
3	D	7	DT	P-O3'-C3'	-6.65	111.72	119.70
2	C	2	DG	C1'-O4'-C4'	-6.12	103.98	110.10
2	C	7	DA	P-O3'-C3'	-5.67	112.90	119.70
2	C	12	DC	P-O3'-C3'	-5.57	113.02	119.70
3	D	4	DG	OP2-P-O3'	5.42	117.13	105.20
3	D	2	DG	P-O3'-C3'	-5.32	113.32	119.70
3	D	2	DG	C1'-O4'-C4'	-5.32	104.78	110.10
3	D	8	DA	C1'-O4'-C4'	-5.07	105.03	110.10
2	C	6	DT	OP2-P-O3'	5.06	116.34	105.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	179	LYS	Mainchain

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	2321	0	2324	63	1
1	B	2319	0	2325	91	0
2	C	265	0	147	2	0
3	D	262	0	147	2	0
4	A	17	0	0	0	0
4	B	6	0	0	0	0
4	C	1	0	0	0	0
All	All	5191	0	4943	154	1

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 (154) 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:200:ALA:HA	1:A:207:LEU:HD23	1.41	1.01
1:A:198:LEU:HB2	1:A:207:LEU:HD11	1.50	0.94
1:B:171:GLN:O	1:B:175:GLU:HG3	1.70	0.91
1:A:167:THR:OG1	1:A:170:GLU:HG2	1.71	0.89
1:B:260:VAL:CG2	1:B:274:GLY:HA3	2.02	0.89
1:A:168:GLU:OE2	1:A:172:ARG:HG3	1.73	0.88
1:A:207:LEU:HD12	1:A:207:LEU:O	1.73	0.87
1:B:296:LYS:HB3	1:B:299:ILE:HD11	1.60	0.84
1:A:240:VAL:HG13	1:A:294:TYR:HB3	1.64	0.78
1:A:173:GLU:O	1:A:176:GLN:HB2	1.84	0.78
1:A:240:VAL:HG22	1:A:325:TYR:HB3	1.67	0.76
1:B:122:VAL:CG2	1:B:131:ALA:HB1	2.16	0.76
1:B:262:PHE:CE1	1:B:307:LEU:HD13	2.23	0.74
1:A:288:VAL:HG11	1:B:232:ARG:NH1	2.03	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:VAL:HG21	1:B:131:ALA:HB1	1.73	0.71
1:B:116:GLU:OE2	1:B:116:GLU:N	2.22	0.70
1:A:93:VAL:O	1:A:119:ILE:HG13	1.92	0.70
2:C:2:DG:H2''	2:C:3:DG:H5''	1.74	0.70
3:D:2:DG:H2''	3:D:3:DG:H5''	1.72	0.70
1:B:163:PRO:O	1:B:167:THR:HG23	1.92	0.70
1:B:241:ARG:HA	1:B:293:PRO:HB3	1.75	0.69
1:B:230:ILE:CG2	1:B:248:LEU:HD11	2.22	0.69
3:D:8:DA:H2''	3:D:9:DC:O5'	1.93	0.68
1:B:74:ARG:O	1:B:74:ARG:HG2	1.92	0.68
1:A:100:ASP:O	1:A:102:PRO:HD3	1.96	0.66
1:A:158:ARG:NH1	1:A:170:GLU:OE1	2.28	0.66
1:B:117:LEU:HD12	1:B:119:ILE:H	1.61	0.66
1:B:230:ILE:HG23	1:B:248:LEU:HD11	1.78	0.66
1:B:92:GLU:HA	1:B:120:CYS:O	1.96	0.65
1:B:147:MET:HE3	1:B:178:ALA:C	2.16	0.65
2:C:8:DA:H2''	2:C:9:DC:O5'	1.96	0.65
1:B:260:VAL:HG23	1:B:274:GLY:H	1.61	0.65
1:A:200:ALA:CA	1:A:207:LEU:HD23	2.23	0.65
1:B:240:VAL:HB	1:B:325:TYR:HB3	1.79	0.64
1:A:146:MET:HE2	1:A:182:LYS:HA	1.80	0.64
1:B:152:GLN:O	1:B:156:ARG:HG3	1.98	0.64
1:B:184:VAL:HG12	1:B:184:VAL:O	1.96	0.63
1:B:168:GLU:OE2	1:B:168:GLU:N	2.31	0.63
1:B:260:VAL:HG23	1:B:274:GLY:HA3	1.81	0.63
1:B:240:VAL:HG23	1:B:294:TYR:HB2	1.80	0.62
1:B:260:VAL:HG21	1:B:274:GLY:HA3	1.81	0.62
1:B:117:LEU:CD1	1:B:119:ILE:HB	2.30	0.62
1:A:299:ILE:HD11	1:A:326:TYR:HB2	1.80	0.62
1:A:117:LEU:O	1:A:117:LEU:HD23	1.99	0.62
1:A:198:LEU:O	1:A:207:LEU:HG	2.00	0.61
1:B:230:ILE:HD13	1:B:248:LEU:HD11	1.83	0.61
1:A:258:ILE:HD11	1:A:309:LEU:HD22	1.81	0.61
1:A:159:LEU:CD1	1:A:164:GLN:HA	2.30	0.61
1:B:240:VAL:HG23	1:B:294:TYR:CB	2.31	0.61
1:A:162:ARG:HB3	1:A:163:PRO:HD2	1.84	0.59
1:B:240:VAL:HA	1:B:325:TYR:CD1	2.38	0.58
1:B:262:PHE:CZ	1:B:307:LEU:HD13	2.38	0.58
1:B:310:LYS:CB	1:B:317:VAL:HG12	2.34	0.58
1:B:167:THR:OG1	1:B:168:GLU:OE2	2.20	0.57
1:B:166:LEU:C	1:B:166:LEU:HD23	2.24	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:VAL:HG23	1:B:274:GLY:N	2.19	0.56
1:B:117:LEU:HA	1:B:156:ARG:HH22	1.71	0.56
1:A:282:HIS:CE1	1:A:283:LYS:HG2	2.41	0.56
1:A:240:VAL:HG21	1:A:299:ILE:HD13	1.87	0.55
1:A:167:THR:O	1:A:170:GLU:HB2	2.06	0.55
1:B:260:VAL:CG2	1:B:274:GLY:CA	2.82	0.55
1:B:260:VAL:HG23	1:B:274:GLY:CA	2.36	0.55
1:B:159:LEU:HD21	1:B:167:THR:HG22	1.88	0.54
1:A:117:LEU:HD22	1:A:156:ARG:HH12	1.72	0.54
1:A:313:ARG:HG3	1:A:313:ARG:HH11	1.72	0.54
1:B:311:ARG:HB2	1:B:316:ASP:OD1	2.06	0.54
1:B:91:ILE:O	1:B:121:ALA:HA	2.08	0.54
1:B:230:ILE:HG21	1:B:248:LEU:HD11	1.89	0.54
1:A:251:ASP:OD1	1:B:283:LYS:HE2	2.08	0.54
1:B:243:GLY:HA2	1:B:291:THR:O	2.07	0.54
1:A:159:LEU:HD13	1:A:164:GLN:HA	1.91	0.53
1:B:310:LYS:HA	1:B:317:VAL:HG12	1.90	0.53
1:A:301:ARG:HG3	1:A:302:PRO:HD2	1.91	0.53
1:A:301:ARG:CG	1:A:302:PRO:HD2	2.39	0.52
1:A:159:LEU:HG	1:A:159:LEU:O	2.09	0.52
1:B:77:TYR:CD1	1:B:134:ASN:HA	2.44	0.52
1:A:77:TYR:CD1	1:A:134:ASN:HA	2.44	0.51
1:B:270:TRP:CD1	1:B:295:HIS:HB3	2.44	0.51
1:B:95:LEU:HD21	1:B:109:LEU:HD21	1.93	0.50
1:B:305:VAL:HG12	1:B:306:PHE:N	2.27	0.50
1:A:95:LEU:HG	1:A:109:LEU:HG	1.92	0.50
1:A:158:ARG:HH12	1:A:170:GLU:HB3	1.76	0.50
1:B:264:GLU:HG3	1:B:270:TRP:HE3	1.76	0.50
1:A:95:LEU:HD21	1:A:109:LEU:HD21	1.94	0.49
1:A:235:LYS:HB2	1:A:246:VAL:HG22	1.93	0.49
1:A:291:THR:HG23	1:A:292:PRO:HD2	1.93	0.49
1:A:200:ALA:HA	1:A:207:LEU:CD2	2.28	0.49
1:B:124:VAL:HG22	1:B:129:MET:HA	1.94	0.49
1:A:240:VAL:CG2	1:A:299:ILE:HD13	2.42	0.49
1:A:107:HIS:ND1	1:A:141:VAL:HG12	2.28	0.48
1:A:288:VAL:HG11	1:B:232:ARG:HH11	1.77	0.48
1:B:40:LEU:HD11	1:B:196:ALA:HB2	1.94	0.48
1:B:233:MET:HE3	1:B:321:LYS:HD3	1.96	0.48
1:B:241:ARG:HG2	1:B:297:MET:SD	2.53	0.48
1:B:240:VAL:CG2	1:B:294:TYR:CB	2.91	0.48
1:A:182:LYS:O	1:A:182:LYS:HG3	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:GLY:CA	1:B:291:THR:O	2.62	0.47
1:A:89:ALA:HB3	1:A:124:VAL:CG2	2.44	0.47
1:B:92:GLU:HB3	1:B:119:ILE:CG2	2.45	0.46
1:B:310:LYS:HB2	1:B:317:VAL:HG12	1.98	0.46
1:B:236:THR:O	1:B:236:THR:HG22	2.15	0.46
1:B:310:LYS:CA	1:B:317:VAL:HG12	2.45	0.46
1:B:40:LEU:HD11	1:B:196:ALA:CB	2.46	0.46
1:B:159:LEU:HD13	1:B:163:PRO:HB3	1.97	0.46
1:A:100:ASP:N	1:A:101:PRO:HD2	2.31	0.46
1:A:252:LYS:HA	1:A:284:GLN:O	2.16	0.46
1:B:77:TYR:CG	1:B:134:ASN:HA	2.51	0.46
1:A:240:VAL:HG21	1:A:299:ILE:CD1	2.46	0.45
1:B:270:TRP:HZ3	1:B:305:VAL:HG11	1.81	0.45
1:B:184:VAL:O	1:B:184:VAL:CG1	2.62	0.45
1:B:296:LYS:O	1:B:299:ILE:HD11	2.16	0.45
1:B:95:LEU:HG	1:B:109:LEU:HG	1.98	0.45
1:A:258:ILE:HD11	1:A:309:LEU:CD2	2.47	0.45
1:A:40:LEU:HD12	1:A:81:LYS:O	2.17	0.45
1:A:288:VAL:CG1	1:B:232:ARG:NH1	2.75	0.45
1:A:240:VAL:HG13	1:A:294:TYR:CB	2.42	0.44
1:B:89:ALA:HB3	1:B:124:VAL:CG1	2.47	0.44
1:B:146:MET:CE	1:B:182:LYS:HA	2.47	0.44
1:A:94:ASP:HB3	1:A:119:ILE:HG12	1.99	0.44
1:B:89:ALA:HB3	1:B:124:VAL:HG12	2.00	0.44
1:B:124:VAL:CG2	1:B:129:MET:HA	2.47	0.43
1:B:262:PHE:CD1	1:B:307:LEU:HD13	2.53	0.43
1:B:100:ASP:N	1:B:101:PRO:CD	2.81	0.43
1:B:270:TRP:HZ3	1:B:305:VAL:CG1	2.30	0.43
1:A:100:ASP:N	1:A:101:PRO:CD	2.82	0.43
1:A:157:GLN:O	1:A:160:ARG:HB3	2.18	0.43
1:B:166:LEU:C	1:B:166:LEU:CD2	2.87	0.43
1:B:240:VAL:CG2	1:B:294:TYR:HB3	2.48	0.43
1:A:109:LEU:HB2	1:A:119:ILE:HD11	2.00	0.43
1:B:270:TRP:CZ3	1:B:305:VAL:HG11	2.54	0.43
1:B:85:TYR:CE2	1:B:198:LEU:HD13	2.54	0.43
1:B:248:LEU:HB2	1:B:289:PHE:HE1	1.84	0.43
1:A:270:TRP:CD1	1:A:295:HIS:HB3	2.53	0.42
1:B:116:GLU:H	1:B:116:GLU:CD	2.19	0.42
1:B:207:LEU:HD12	1:B:208:PRO:CD	2.49	0.42
1:A:181:LEU:HD12	1:A:181:LEU:HA	1.92	0.42
1:B:117:LEU:HD12	1:B:119:ILE:N	2.30	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:GLU:CG	1:B:270:TRP:HE3	2.32	0.42
1:A:44:GLU:O	1:A:78:PRO:HA	2.20	0.42
1:A:257:ASP:OD1	1:A:313:ARG:HG2	2.20	0.41
1:B:116:GLU:N	1:B:116:GLU:CD	2.74	0.41
1:B:198:LEU:O	1:B:207:LEU:HB3	2.19	0.41
1:B:44:GLU:O	1:B:78:PRO:HA	2.19	0.41
1:B:71:GLU:O	1:B:74:ARG:O	2.38	0.41
1:B:146:MET:HE1	1:B:182:LYS:HA	2.01	0.41
1:A:100:ASP:O	1:A:102:PRO:CD	2.68	0.41
1:B:117:LEU:H	1:B:117:LEU:HG	1.63	0.41
1:A:77:TYR:CG	1:A:134:ASN:HA	2.56	0.41
1:A:97:THR:HG23	1:A:103:ARG:O	2.21	0.41
1:B:147:MET:CE	1:B:178:ALA:HB1	2.51	0.41
1:A:199:ARG:O	1:A:207:LEU:HD21	2.21	0.40
1:A:157:GLN:HG2	1:A:160:ARG:HE	1.87	0.40
1:A:228:LEU:HD12	1:A:317:VAL:HA	2.03	0.40
1:B:300:GLU:HG2	1:B:301:ARG:HG2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:TYR:O	1:A:326:TYR:O[10_444]	1.84	0.36

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	292/327 (89%)	272 (93%)	20 (7%)	0	100	100
1	B	292/327 (89%)	272 (93%)	20 (7%)	0	100	100
All	All	584/654 (89%)	544 (93%)	40 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	255/285 (90%)	251 (98%)	4 (2%)	62	86
1	B	255/285 (90%)	250 (98%)	5 (2%)	55	83
All	All	510/570 (90%)	501 (98%)	9 (2%)	59	85

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	MET
1	A	202	ASP
1	A	275	ASP
1	A	313	ARG
1	B	117	LEU
1	B	159	LEU
1	B	202	ASP
1	B	263	TYR
1	B	295	HIS

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

Mol	Chain	Res	Type
1	A	113	GLN
1	A	152	GLN
1	A	282	HIS
1	A	308	GLN
1	B	152	GLN
1	B	227	ASN

5.3.3 RNA ⓘ

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 monosaccharides 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, 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	294/327 (89%)	0.10	4 (1%) 75 49	60, 117, 196, 242	0
1	B	294/327 (89%)	0.34	23 (7%) 13 4	73, 130, 198, 238	0
2	C	13/13 (100%)	0.08	0 100 100	74, 80, 87, 88	0
3	D	13/13 (100%)	0.14	0 100 100	74, 78, 84, 90	0
All	All	614/680 (90%)	0.22	27 (4%) 34 13	60, 120, 197, 242	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	165	GLY	5.7
1	B	260	VAL	4.6
1	B	305	VAL	4.0
1	B	307	LEU	3.4
1	B	308	GLN	3.1
1	A	166	LEU	2.9
1	B	325	TYR	2.9
1	B	167	THR	2.9
1	B	294	TYR	2.9
1	B	159	LEU	2.8
1	A	160	ARG	2.8
1	B	270	TRP	2.7
1	B	259	GLU	2.6
1	A	34	THR	2.6
1	B	309	LEU	2.6
1	B	164	GLN	2.5
1	A	158	ARG	2.5
1	B	269	GLY	2.5
1	B	268	ASN	2.5
1	B	228	LEU	2.3
1	B	262	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	295	HIS	2.1
1	B	296	LYS	2.1
1	B	267	GLU	2.1
1	B	306	PHE	2.0
1	B	241	ARG	2.0
1	B	323	PHE	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 monosaccharides 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.