



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 04:32 pm BST

PDB ID : 2PQA
Title : Crystal Structure of Full-length Human RPA 14/32 Heterodimer
Authors : Deng, X.; Borgstahl, G.E.
Deposited on : 2007-05-01
Resolution : 2.50 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	:	2.11
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)	:	2.11

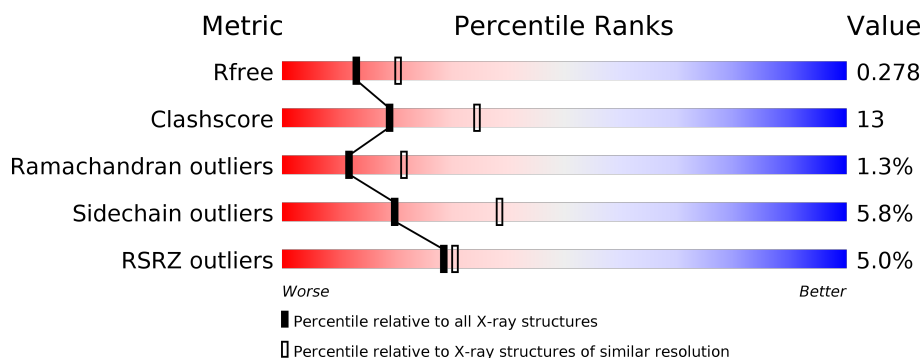
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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 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	131	<div> <div>9%</div> <div>64%</div> <div>31%</div> <div>• •</div> </div>
1	C	131	<div> <div>7%</div> <div>61%</div> <div>33%</div> <div>• •</div> </div>
2	B	142	<div> <div>2%</div> <div>63%</div> <div>14%</div> <div>• •</div> <div>20%</div> </div>
2	D	142	<div> <div>65%</div> <div>13%</div> <div>•</div> <div>20%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3785 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 Replication protein A 32 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			1003	638	171	188	6			
1	C	127	Total	C	N	O	S	0	0	0
			996	632	167	191	6			

- Molecule 2 is a protein called Replication protein A 14 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	114	Total	C	N	O	S	0	0	0
			893	574	145	167	7			
2	D	114	Total	C	N	O	S	0	0	0
			893	576	145	165	7			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MET	-	EXPRESSION TAG	UNP P35244
B	-19	GLY	-	EXPRESSION TAG	UNP P35244
B	-18	HIS	-	EXPRESSION TAG	UNP P35244
B	-17	HIS	-	EXPRESSION TAG	UNP P35244
B	-16	HIS	-	EXPRESSION TAG	UNP P35244
B	-15	HIS	-	EXPRESSION TAG	UNP P35244
B	-14	HIS	-	EXPRESSION TAG	UNP P35244
B	-13	HIS	-	EXPRESSION TAG	UNP P35244
B	-12	HIS	-	EXPRESSION TAG	UNP P35244
B	-11	HIS	-	EXPRESSION TAG	UNP P35244
B	-10	HIS	-	EXPRESSION TAG	UNP P35244
B	-9	HIS	-	EXPRESSION TAG	UNP P35244
B	-8	SER	-	EXPRESSION TAG	UNP P35244
B	-7	SER	-	EXPRESSION TAG	UNP P35244
B	-6	GLY	-	EXPRESSION TAG	UNP P35244
B	-5	HIS	-	EXPRESSION TAG	UNP P35244

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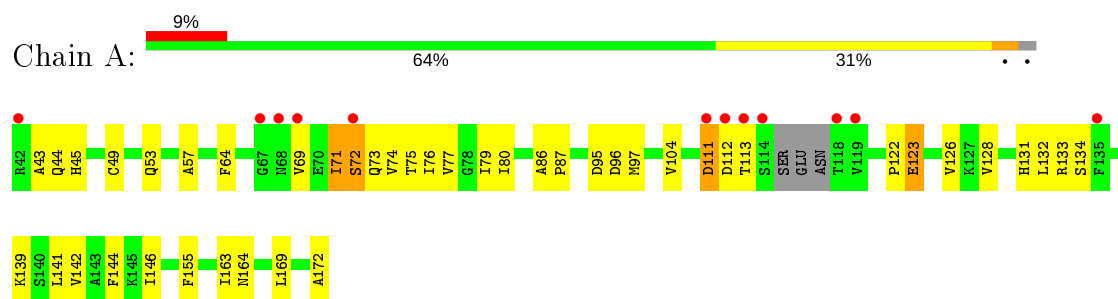
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	ILE	-	EXPRESSION TAG	UNP P35244
B	-3	GLU	-	EXPRESSION TAG	UNP P35244
B	-2	GLY	-	EXPRESSION TAG	UNP P35244
B	-1	ARG	-	EXPRESSION TAG	UNP P35244
B	0	HIS	-	EXPRESSION TAG	UNP P35244
D	-20	MET	-	EXPRESSION TAG	UNP P35244
D	-19	GLY	-	EXPRESSION TAG	UNP P35244
D	-18	HIS	-	EXPRESSION TAG	UNP P35244
D	-17	HIS	-	EXPRESSION TAG	UNP P35244
D	-16	HIS	-	EXPRESSION TAG	UNP P35244
D	-15	HIS	-	EXPRESSION TAG	UNP P35244
D	-14	HIS	-	EXPRESSION TAG	UNP P35244
D	-13	HIS	-	EXPRESSION TAG	UNP P35244
D	-12	HIS	-	EXPRESSION TAG	UNP P35244
D	-11	HIS	-	EXPRESSION TAG	UNP P35244
D	-10	HIS	-	EXPRESSION TAG	UNP P35244
D	-9	HIS	-	EXPRESSION TAG	UNP P35244
D	-8	SER	-	EXPRESSION TAG	UNP P35244
D	-7	SER	-	EXPRESSION TAG	UNP P35244
D	-6	GLY	-	EXPRESSION TAG	UNP P35244
D	-5	HIS	-	EXPRESSION TAG	UNP P35244
D	-4	ILE	-	EXPRESSION TAG	UNP P35244
D	-3	GLU	-	EXPRESSION TAG	UNP P35244
D	-2	GLY	-	EXPRESSION TAG	UNP P35244
D	-1	ARG	-	EXPRESSION TAG	UNP P35244
D	0	HIS	-	EXPRESSION TAG	UNP P35244

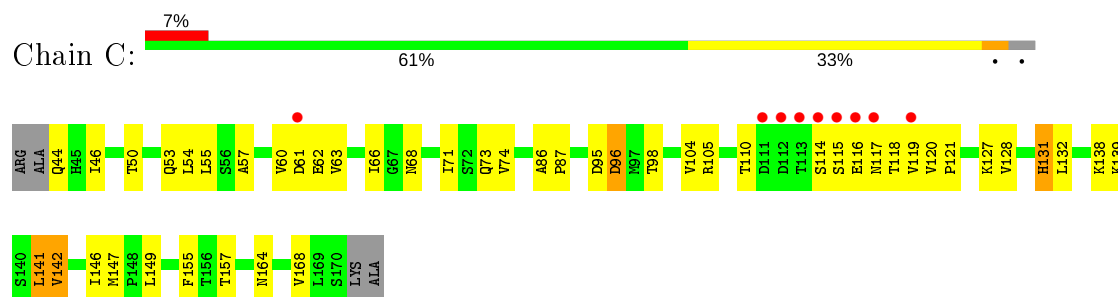
3 Residue-property plots [i](#)

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

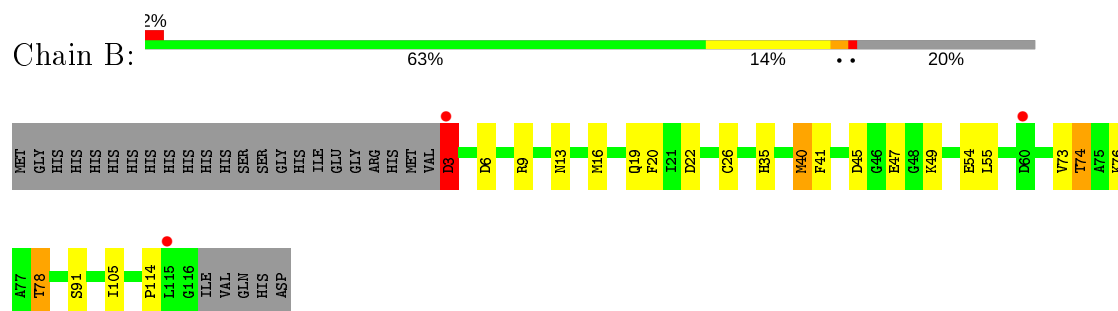
- Molecule 1: Replication protein A 32 kDa subunit



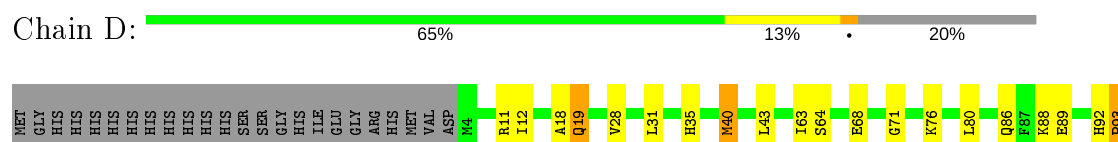
- Molecule 1: Replication protein A 32 kDa subunit



- Molecule 2: Replication protein A 14 kDa subunit



- Molecule 2: Replication protein A 14 kDa subunit



117	VAL
	GLN
	HIS
	ASP

4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	63.38Å 63.38Å 272.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.50 31.76 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.0 (25.00-2.50) 96.2 (31.76-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.228 , 0.280 0.231 , 0.278	Depositor DCC
R_{free} test set	1195 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 29.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.086 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3785	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% 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	0.56	0/1020	0.76	1/1387 (0.1%)
1	C	0.60	0/1014	0.71	1/1382 (0.1%)
2	B	0.61	0/912	0.69	1/1231 (0.1%)
2	D	0.55	0/912	0.75	1/1231 (0.1%)
All	All	0.58	0/3858	0.73	4/5231 (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
1	A	1	1
2	B	1	1
All	All	2	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	6	ASP	CB-CG-OD1	5.93	123.64	118.30
2	D	92	HIS	N-CA-C	5.63	126.21	111.00
1	A	95	ASP	CB-CG-OD1	5.28	123.06	118.30
1	C	96	ASP	N-CA-C	-5.02	97.45	111.00

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	112	ASP	CA
2	B	3	ASP	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	71	ILE	Peptide
2	B	3	ASP	Peptide

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	1003	0	1024	45	0
1	C	996	0	1006	32	0
2	B	893	0	892	19	0
2	D	893	0	899	11	0
All	All	3785	0	3821	99	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 13.

All (99) 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:B:20:PHE:O	2:B:73:VAL:HG21	1.33	1.28
1:A:111:ASP:O	1:A:112:ASP:OD1	1.68	1.12
1:A:111:ASP:O	1:A:112:ASP:CG	1.93	1.07
1:C:66:ILE:HD12	1:C:71:ILE:HD13	1.43	0.98
1:A:164:ASN:HD21	1:C:46:ILE:H	1.07	0.98
1:C:66:ILE:HD12	1:C:71:ILE:CD1	1.94	0.97
2:B:20:PHE:O	2:B:73:VAL:CG2	2.16	0.93
1:A:49:CYS:O	1:A:76:ILE:HD13	1.69	0.93
1:A:76:ILE:HD12	1:A:77:VAL:H	1.35	0.91
1:C:61:ASP:N	1:C:62:GLU:HA	1.88	0.88
2:B:22:ASP:N	2:B:73:VAL:HG23	1.95	0.81
1:A:133:ARG:HG3	1:A:142:VAL:CG2	2.15	0.76
2:B:22:ASP:H	2:B:73:VAL:HG23	1.52	0.73
1:C:54:LEU:HD21	1:C:74:VAL:HG11	1.70	0.73
1:A:126:VAL:CG1	1:A:146:ILE:HG23	2.22	0.69
1:A:133:ARG:HG3	1:A:142:VAL:HG23	1.74	0.68
2:D:89:GLU:OE2	2:D:93:PRO:HB2	1.93	0.68
1:A:164:ASN:ND2	1:C:46:ILE:H	1.88	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ILE:CD1	1:A:77:VAL:H	2.06	0.66
2:B:3:ASP:N	2:B:3:ASP:OD1	2.30	0.65
1:A:126:VAL:HG11	1:A:146:ILE:HG23	1.78	0.65
1:C:61:ASP:HB2	1:C:63:VAL:HG12	1.78	0.64
1:A:45:HIS:HB3	1:C:164:ASN:CG	2.17	0.64
2:B:20:PHE:C	2:B:73:VAL:HG21	2.15	0.64
2:D:35:HIS:HB2	2:D:40:MET:HG3	1.79	0.64
1:A:122:PRO:O	1:A:123:GLU:HB3	1.98	0.64
2:B:54:GLU:OE2	2:B:78:THR:HG21	1.99	0.63
1:C:57:ALA:O	1:C:139:LYS:NZ	2.29	0.63
2:D:11:ARG:NH2	2:D:68:GLU:OE2	2.24	0.63
1:A:72:SER:OG	1:A:73:GLN:N	2.32	0.61
1:A:133:ARG:HG3	1:A:142:VAL:HG21	1.81	0.61
1:A:122:PRO:O	1:A:123:GLU:CB	2.51	0.58
2:B:74:THR:HG22	2:B:76:LYS:H	1.68	0.58
2:D:31:LEU:HA	2:D:43:LEU:HD23	1.86	0.57
1:C:53:GLN:HB3	1:C:66:ILE:HD13	1.85	0.57
2:D:64:SER:H	2:D:86:GLN:NE2	2.03	0.57
1:C:66:ILE:HD12	1:C:71:ILE:HD12	1.83	0.56
1:A:163:ILE:HD13	2:B:105:ILE:HG22	1.88	0.56
1:A:71:ILE:HD12	1:A:74:VAL:CG1	2.36	0.56
1:A:75:THR:HA	1:A:128:VAL:O	2.05	0.56
1:C:73:GLN:HG2	1:C:131:HIS:CE1	2.41	0.55
1:C:60:VAL:C	1:C:62:GLU:HA	2.27	0.55
1:A:44:GLN:O	1:A:44:GLN:HG3	2.07	0.54
1:C:95:ASP:OD1	1:C:96:ASP:O	2.26	0.54
2:B:16:MET:HA	2:B:19:GLN:OE1	2.08	0.54
1:A:97:MET:HE2	2:B:26:CYS:HB2	1.91	0.53
1:C:61:ASP:H	1:C:63:VAL:H	1.55	0.53
1:C:131:HIS:HB2	1:C:142:VAL:HG12	1.93	0.51
1:A:111:ASP:O	1:A:112:ASP:CB	2.58	0.51
1:A:111:ASP:C	1:A:112:ASP:OD1	2.45	0.51
1:A:76:ILE:HD12	1:A:77:VAL:N	2.16	0.50
2:D:28:VAL:HG22	2:D:68:GLU:HG3	1.93	0.50
1:A:126:VAL:HG11	1:A:146:ILE:CG2	2.43	0.49
1:C:61:ASP:N	1:C:62:GLU:CA	2.70	0.49
1:A:132:LEU:HA	1:A:141:LEU:HD23	1.95	0.48
2:B:47:GLU:HG3	2:B:49:LYS:HG3	1.94	0.48
1:A:64:PHE:HB2	1:A:71:ILE:O	2.13	0.48
1:A:131:HIS:CD2	1:A:144:PHE:CE1	3.02	0.47
2:B:13:ASN:OD1	2:B:16:MET:HG2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:THR:HG22	1:C:119:VAL:H	1.79	0.47
2:D:31:LEU:HD22	2:D:63:ILE:O	2.15	0.47
1:C:73:GLN:C	1:C:74:VAL:HG23	2.34	0.47
1:A:97:MET:CE	2:B:9:ARG:HD2	2.45	0.47
1:A:53:GLN:HG3	2:B:114:PRO:HG2	1.95	0.47
2:D:71:GLY:HA3	2:D:80:LEU:O	2.15	0.46
2:B:35:HIS:HB2	2:B:40:MET:HG3	1.97	0.46
1:A:69:VAL:HG11	1:A:169:LEU:HD22	1.97	0.46
1:C:127:LYS:HB2	1:C:149:LEU:HD21	1.98	0.45
1:C:71:ILE:HG23	1:C:74:VAL:HG22	1.99	0.45
1:A:80:ILE:HG13	1:A:126:VAL:HG21	1.98	0.44
1:A:64:PHE:HZ	1:A:134:SER:HB2	1.83	0.44
1:A:86:ALA:HB1	1:A:87:PRO:HD2	1.99	0.44
1:A:104:VAL:HG22	1:A:141:LEU:HB2	2.00	0.44
1:C:68:ASN:HA	1:C:68:ASN:HD22	1.63	0.44
1:C:66:ILE:CD1	1:C:71:ILE:CD1	2.83	0.43
1:A:86:ALA:HB1	1:A:87:PRO:CD	2.49	0.43
1:A:57:ALA:O	1:A:139:LYS:NZ	2.51	0.43
1:A:96:ASP:C	1:A:96:ASP:OD1	2.58	0.43
1:A:131:HIS:CD2	1:A:144:PHE:HE1	2.37	0.42
1:C:54:LEU:HD22	1:C:132:LEU:HD13	2.00	0.42
1:A:79:ILE:HD11	1:A:123:GLU:HA	2.01	0.42
1:A:133:ARG:CG	1:A:142:VAL:HG23	2.47	0.42
1:C:66:ILE:CG1	1:C:71:ILE:HD12	2.49	0.42
2:B:13:ASN:HB2	2:B:45:ASP:OD2	2.20	0.42
1:C:50:THR:HB	1:C:98:THR:HG21	2.02	0.41
2:B:47:GLU:HG3	2:B:49:LYS:CG	2.50	0.41
1:C:120:VAL:HA	1:C:121:PRO:HD3	1.94	0.41
2:D:88:LYS:HE2	2:D:88:LYS:HB2	1.88	0.41
1:C:66:ILE:CD1	1:C:71:ILE:HD12	2.49	0.41
1:C:104:VAL:HG22	1:C:141:LEU:HB2	2.02	0.41
1:A:126:VAL:CG1	1:A:146:ILE:CG2	2.94	0.41
2:B:41:PHE:CE2	2:B:55:LEU:HD11	2.55	0.41
1:A:172:ALA:HB2	1:C:168:VAL:HG13	2.02	0.41
2:D:18:ALA:HB3	2:D:19:GLN:HE21	1.86	0.41
2:D:19:GLN:H	2:D:19:GLN:NE2	2.19	0.41
1:C:86:ALA:HB1	1:C:87:PRO:HD2	2.02	0.41
1:A:43:ALA:HA	1:A:44:GLN:HA	1.91	0.40
1:A:76:ILE:CG1	1:A:77:VAL:N	2.84	0.40
1:C:128:VAL:HG22	1:C:146:ILE:HG12	2.02	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	
1	A	124/131 (95%)	118 (95%)	4 (3%)	2 (2%)	9	17
1	C	125/131 (95%)	112 (90%)	10 (8%)	3 (2%)	6	9
2	B	112/142 (79%)	108 (96%)	4 (4%)	0	100	100
2	D	112/142 (79%)	107 (96%)	4 (4%)	1 (1%)	17	31
All	All	473/546 (87%)	445 (94%)	22 (5%)	6 (1%)	12	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	SER
1	C	117	ASN
1	C	114	SER
1	C	115	SER
1	A	123	GLU
2	D	93	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	115/118 (98%)	112 (97%)	3 (3%)	46	72
1	C	116/118 (98%)	104 (90%)	12 (10%)	7	14
2	B	99/124 (80%)	94 (95%)	5 (5%)	24	45
2	D	99/124 (80%)	94 (95%)	5 (5%)	24	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	429/484 (89%)	404 (94%)	25 (6%)	20	38

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	ASP
1	A	113	THR
1	A	155	PHE
2	B	3	ASP
2	B	40	MET
2	B	74	THR
2	B	78	THR
2	B	91	SER
1	C	44	GLN
1	C	55	LEU
1	C	105	ARG
1	C	110	THR
1	C	116	GLU
1	C	131	HIS
1	C	138	LYS
1	C	141	LEU
1	C	142	VAL
1	C	147	MET
1	C	155	PHE
1	C	157	THR
2	D	12	ILE
2	D	19	GLN
2	D	40	MET
2	D	76	LYS
2	D	117	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	131	HIS
1	A	164	ASN
1	C	68	ASN
1	C	73	GLN
1	C	131	HIS
1	C	164	ASN

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Mol	Chain	Res	Type
2	D	19	GLN
2	D	86	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 ⓘ

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	128/131 (97%)	0.33	12 (9%) 8 8	22, 37, 59, 67	0
1	C	127/131 (96%)	0.33	9 (7%) 16 16	27, 37, 65, 73	0
2	B	114/142 (80%)	-0.19	3 (2%) 56 59	20, 33, 48, 52	0
2	D	114/142 (80%)	-0.25	0 100 100	26, 34, 51, 54	0
All	All	483/546 (88%)	0.07	24 (4%) 28 30	20, 35, 55, 73	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	SER	5.5
1	C	114	SER	5.4
1	C	61	ASP	5.0
1	C	113	THR	4.9
1	A	112	ASP	4.6
1	C	111	ASP	4.0
1	C	115	SER	4.0
1	C	116	GLU	3.9
1	A	118	THR	3.6
1	A	119	VAL	3.5
1	C	119	VAL	3.4
1	A	135	PHE	3.3
1	A	113	THR	3.1
1	C	117	ASN	3.1
2	B	115	LEU	3.0
1	A	111	ASP	2.9
2	B	60	ASP	2.9
1	C	112	ASP	2.8
1	A	42	ARG	2.8
2	B	3	ASP	2.8
1	A	68	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	72	SER	2.5
1	A	67	GLY	2.4
1	A	69	VAL	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.