



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:11 am GMT

PDB ID : 4D2G
Title : Crystal structure of human PCNA in complex with p15 peptide
Authors : DeBiasio, A.; Ibanez, A.; Mortuza, G.; Molina, R.; Cordeiro, T.N.; Castillo, F.; Villate, M.; Merino, N.; Lelli, M.; Diercks, T.; Luque, I.; Bernardo, P.; Montoya, G.; Blanco, F.J.
Deposited on : 2014-05-09
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

A user guide is available at

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

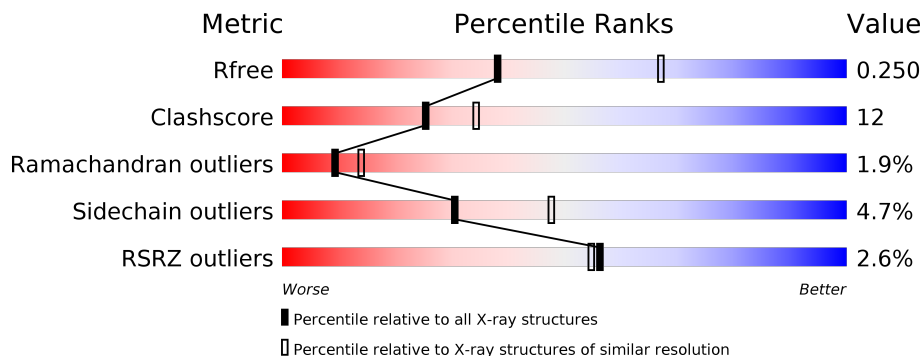
1 Overall quality at a glance

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}	100719	3491 (2.70-2.62)
Clashscore	112137	1026 (2.68-2.64)
Ramachandran outliers	110173	1010 (2.68-2.64)
Sidechain outliers	110143	1010 (2.68-2.64)
RSRZ outliers	101464	3511 (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 ≥ 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	264	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>26%</div> <div>• •</div> </div> </div>
1	B	264	<div> <div>2%</div> <div> <div></div> <div>66%</div> <div>28%</div> <div>• •</div> </div> </div>
1	C	264	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• •</div> </div> </div>
2	D	21	<div> <div></div> <div> <div>95%</div> <div>5%</div> </div> </div>
2	E	21	<div> <div>10%</div> <div> <div></div> <div>71%</div> <div>19%</div> <div>10%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6461 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 PROLIFERATING CELL NUCLEAR ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1984	1246	327	395	16			
1	B	256	Total	C	N	O	S	0	0	0
			1973	1239	325	393	16			
1	C	258	Total	C	N	O	S	0	0	0
			1984	1246	327	395	16			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P12004
A	-1	PRO	-	EXPRESSION TAG	UNP P12004
A	0	HIS	-	EXPRESSION TAG	UNP P12004
B	-2	GLY	-	EXPRESSION TAG	UNP P12004
B	-1	PRO	-	EXPRESSION TAG	UNP P12004
B	0	HIS	-	EXPRESSION TAG	UNP P12004
C	-2	GLY	-	EXPRESSION TAG	UNP P12004
C	-1	PRO	-	EXPRESSION TAG	UNP P12004
C	0	HIS	-	EXPRESSION TAG	UNP P12004

- Molecule 2 is a protein called P15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	21	Total	C	N	O	S	0	0	0
			162	108	28	25	1			
2	E	21	Total	C	N	O	S	0	0	0
			162	108	28	25	1			

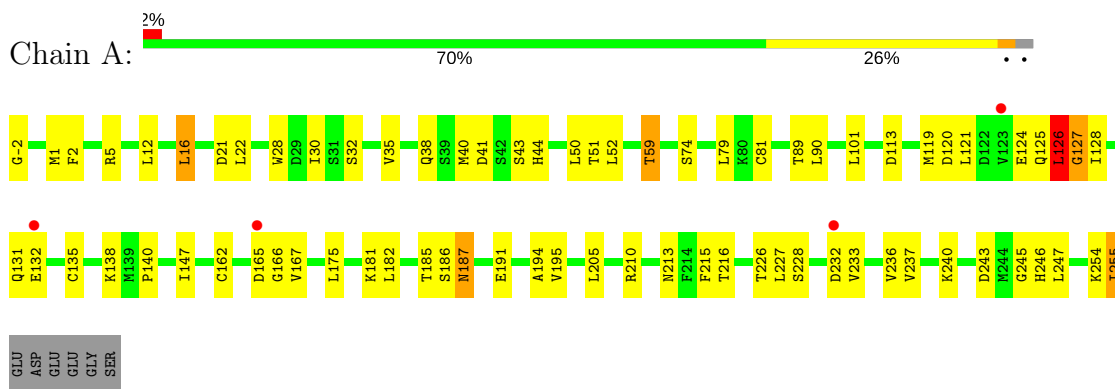
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	75	Total 75	O 75	0	0
3	B	58	Total 58	O 58	0	0
3	C	60	Total 60	O 60	0	0
3	D	2	Total 2	O 2	0	0
3	E	1	Total 1	O 1	0	0

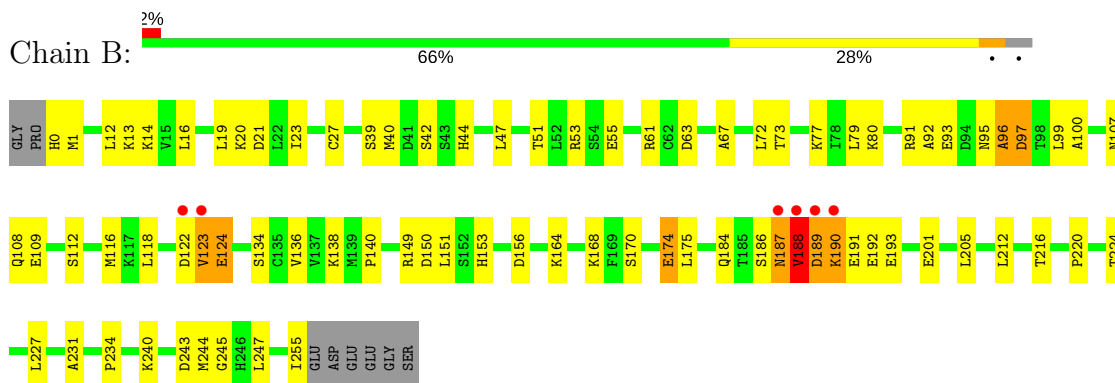
3 Residue-property plots [i](#)

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

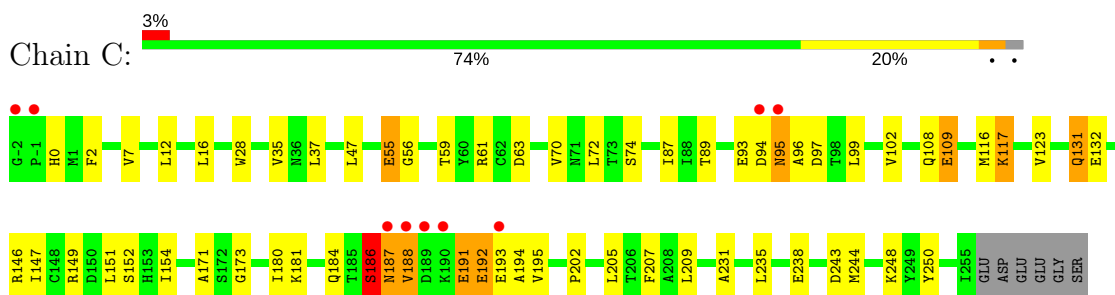
• Molecule 1: PROLIFERATING CELL NUCLEAR ANTIGEN



• Molecule 1: PROLIFERATING CELL NUCLEAR ANTIGEN



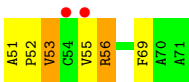
• Molecule 1: PROLIFERATING CELL NUCLEAR ANTIGEN



• Molecule 2: P15



● Molecule 2: P15



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.05Å 41.23Å 141.12Å 90.00° 105.39° 90.00°	Depositor
Resolution (Å)	44.35 – 2.65 44.35 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.5 (44.35-2.65) 98.6 (44.35-2.65)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.65Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.170 , 0.250 0.173 , 0.250	Depositor DCC
R_{free} test set	1240 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.401	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6461	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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 [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	0.73	2/2012 (0.1%)	0.86	6/2719 (0.2%)
1	B	0.77	0/2000	0.83	2/2702 (0.1%)
1	C	0.76	0/2012	0.84	0/2719
2	D	0.88	0/168	0.82	0/228
2	E	0.80	0/168	0.91	0/228
All	All	0.76	2/6360 (0.0%)	0.85	8/8596 (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	0	1
1	B	0	3
1	C	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	127	GLY	C-O	6.24	1.33	1.23
1	A	135	CYS	CB-SG	-5.16	1.73	1.81

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	189	ASP	CB-CG-OD1	7.56	125.10	118.30
1	A	22	LEU	CA-CB-CG	6.38	129.98	115.30
1	B	16	LEU	CA-CB-CG	-5.93	101.67	115.30
1	A	120	ASP	CB-CG-OD1	5.69	123.42	118.30
1	A	16	LEU	CA-CB-CG	-5.46	102.75	115.30
1	A	126	LEU	CA-CB-CG	5.42	127.77	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	LEU	CB-CG-CD2	-5.18	102.20	111.00
1	A	205	LEU	CA-CB-CG	5.08	126.97	115.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	232	ASP	Peptide
1	B	187	ASN	Peptide
1	B	188	VAL	Peptide
1	B	190	LYS	Peptide
1	C	186	SER	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	1984	0	1990	46	1
1	B	1973	0	1978	62	0
1	C	1984	0	1990	47	1
2	D	162	0	163	1	0
2	E	162	0	163	5	0
3	A	75	0	0	6	0
3	B	58	0	0	6	0
3	C	60	0	0	10	0
3	D	2	0	0	0	0
3	E	1	0	0	0	0
All	All	6461	0	6284	151	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 12.

All (151) 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:C:202:PRO:CB	3:C:2044:HOH:O	1.81	1.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:PRO:CG	3:C:2044:HOH:O	1.99	1.04
1:C:202:PRO:HG3	3:C:2044:HOH:O	1.58	0.99
1:B:107:ASN:ND2	3:B:2022:HOH:O	2.00	0.93
1:C:202:PRO:HB3	3:C:2044:HOH:O	1.51	0.89
1:B:97:ASP:HB2	1:B:118:LEU:HB2	1.56	0.88
1:A:138:LYS:NZ	1:A:191:GLU:OE2	2.14	0.79
1:C:93:GLU:HG2	1:C:95:ASN:HB2	1.63	0.79
1:A:126:LEU:O	1:A:128:ILE:N	2.19	0.76
1:B:190:LYS:HG2	1:B:193:GLU:HB2	1.67	0.76
1:B:190:LYS:HE2	1:B:193:GLU:HG2	1.67	0.75
1:A:228:SER:HB2	1:A:236:VAL:HG22	1.67	0.75
1:A:185:THR:HG22	1:C:109:GLU:HG2	1.69	0.74
1:B:190:LYS:HE2	1:B:193:GLU:CG	2.17	0.74
1:C:202:PRO:CA	3:C:2044:HOH:O	2.22	0.71
1:A:38:GLN:NE2	1:A:126:LEU:HB2	2.07	0.70
1:A:243:ASP:OD1	3:A:2072:HOH:O	2.09	0.70
1:B:184:GLN:OE1	3:B:2039:HOH:O	2.07	0.70
1:A:191:GLU:OE2	3:A:2043:HOH:O	2.10	0.69
1:A:5:ARG:HB3	1:A:59:THR:HB	1.73	0.69
1:A:119:MET:HE3	1:A:121:LEU:HD21	1.76	0.67
1:C:0:HIS:HA	1:C:63:ASP:OD2	1.96	0.66
1:B:0:HIS:HA	1:B:63:ASP:OD2	1.96	0.66
1:B:97:ASP:N	1:B:97:ASP:OD1	2.29	0.66
1:C:131:GLN:HG2	3:C:2031:HOH:O	1.94	0.66
1:B:138:LYS:NZ	1:B:240:LYS:NZ	2.43	0.65
1:B:187:ASN:C	1:B:189:ASP:H	1.99	0.65
1:B:164:LYS:O	3:B:2039:HOH:O	2.13	0.65
1:A:-2:GLY:N	3:A:2001:HOH:O	2.21	0.64
1:A:254:LYS:HD2	1:A:255:ILE:H	1.64	0.63
1:B:1:MET:HB3	1:B:61:ARG:HH12	1.61	0.63
1:A:40:MET:HG2	1:A:44:HIS:HA	1.80	0.63
1:B:140:PRO:HG3	1:B:193:GLU:HA	1.80	0.63
1:B:138:LYS:HZ2	1:B:240:LYS:NZ	1.96	0.63
1:A:131:GLN:NE2	1:A:132:GLU:O	2.32	0.62
1:C:188:VAL:HG11	1:C:193:GLU:HB2	1.82	0.61
1:B:187:ASN:HA	1:B:189:ASP:OD1	2.00	0.60
1:A:125:GLN:HG3	1:A:126:LEU:N	2.16	0.60
1:A:186:SER:OG	1:A:187:ASN:N	2.35	0.60
1:B:138:LYS:NZ	1:B:240:LYS:HZ2	2.01	0.58
1:C:186:SER:OG	1:C:187:ASN:N	2.30	0.58
1:B:134:SER:HB3	1:B:201:GLU:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ASN:OD1	1:B:109:GLU:N	2.20	0.57
1:C:238:GLU:OE2	1:C:248:LYS:HE3	2.05	0.57
1:A:125:GLN:O	1:A:126:LEU:HB3	2.04	0.56
1:C:202:PRO:HA	3:C:2044:HOH:O	1.98	0.56
1:B:107:ASN:OD1	1:B:108:GLN:N	2.40	0.55
1:B:138:LYS:HZ2	1:B:240:LYS:HZ2	1.53	0.55
1:B:39:SER:O	1:B:47:LEU:HD12	2.06	0.55
1:C:191:GLU:O	1:C:193:GLU:N	2.40	0.54
1:B:168:LYS:NZ	3:B:2042:HOH:O	2.40	0.54
1:C:47:LEU:HB2	2:D:65:ILE:HG21	1.90	0.54
1:B:91:ARG:HB3	1:B:100:ALA:HB3	1.89	0.53
1:A:185:THR:HG23	1:A:195:VAL:HB	1.90	0.53
1:B:156:ASP:HB3	2:E:56:ARG:HD2	1.90	0.53
1:A:124:GLU:N	1:A:124:GLU:OE1	2.42	0.53
1:C:151:LEU:HD22	1:C:171:ALA:HB3	1.91	0.53
1:B:73:THR:O	1:B:77:LYS:HG3	2.08	0.53
1:A:89:THR:O	1:A:90:LEU:HD23	2.09	0.53
1:B:187:ASN:C	1:B:189:ASP:N	2.60	0.53
1:B:19:LEU:HD21	1:B:247:LEU:HD11	1.91	0.53
1:A:21:ASP:HB2	3:A:2007:HOH:O	2.08	0.52
1:B:190:LYS:CD	1:B:191:GLU:H	2.22	0.52
1:A:81:CYS:SG	1:B:150:ASP:HB3	2.50	0.52
1:C:108:GLN:NE2	3:C:2027:HOH:O	2.31	0.52
1:C:184:GLN:HG3	1:C:195:VAL:O	2.10	0.52
1:A:50:LEU:HD13	1:A:247:LEU:HD13	1.91	0.51
1:B:53:ARG:HB3	1:B:55:GLU:OE1	2.10	0.51
1:C:7:VAL:HG22	1:C:87:ILE:HD12	1.92	0.51
1:C:205:LEU:HD21	1:C:231:ALA:HA	1.93	0.51
1:A:140:PRO:HD3	1:A:194:ALA:O	2.11	0.50
1:C:97:ASP:OD2	1:C:117:LYS:HE3	2.12	0.50
1:B:1:MET:HB3	1:B:61:ARG:NH1	2.27	0.49
1:A:35:VAL:HB	1:A:52:LEU:HB2	1.95	0.49
1:C:188:VAL:HG12	1:C:191:GLU:H	1.78	0.49
1:A:227:LEU:HD23	1:A:237:VAL:HG22	1.94	0.49
1:B:93:GLU:O	1:B:95:ASN:N	2.46	0.48
1:C:154:ILE:O	1:C:173:GLY:HA3	2.12	0.48
1:A:246:HIS:CD2	3:A:2014:HOH:O	2.65	0.48
1:A:51:THR:O	1:A:245:GLY:HA3	2.13	0.48
1:B:190:LYS:HE3	1:B:192:GLU:N	2.28	0.48
1:B:190:LYS:NZ	1:B:192:GLU:HG2	2.28	0.48
1:C:235:LEU:O	1:C:250:TYR:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:PHE:CD1	1:A:30:ILE:HG21	2.48	0.48
1:B:174:GLU:HG2	3:B:2034:HOH:O	2.13	0.48
1:C:146:ARG:HG3	1:C:146:ARG:HH11	1.79	0.48
1:B:136:VAL:HA	1:B:227:LEU:O	2.13	0.47
1:B:23:ILE:HB	1:B:72:LEU:HD12	1.96	0.47
1:B:190:LYS:HE3	1:B:193:GLU:N	2.29	0.47
1:C:188:VAL:HG12	1:C:191:GLU:N	2.29	0.47
1:B:27:CYS:SG	1:B:67:ALA:HB1	2.55	0.47
1:C:70:VAL:HG12	1:C:72:LEU:HD23	1.97	0.47
1:B:187:ASN:OD1	1:B:188:VAL:N	2.48	0.46
1:C:56:GLY:HA3	1:C:244:MET:HG3	1.97	0.46
1:B:153:HIS:NE2	2:E:51:ALA:HB1	2.30	0.46
1:A:165:ASP:OD1	1:A:166:GLY:N	2.48	0.46
1:B:123:VAL:H	1:B:123:VAL:HG22	1.35	0.46
1:C:108:GLN:HA	3:C:2028:HOH:O	2.16	0.46
1:A:185:THR:HG22	1:C:109:GLU:CG	2.44	0.46
1:B:234:PRO:HG2	2:E:69:PHE:CZ	2.50	0.46
1:B:122:ASP:O	1:B:124:GLU:HG3	2.16	0.46
1:C:99:LEU:HD23	1:C:116:MET:CE	2.45	0.46
1:B:20:LYS:HG3	1:B:21:ASP:N	2.31	0.45
1:C:152:SER:HA	1:C:209:LEU:HD13	1.97	0.45
1:A:113:ASP:OD1	1:A:113:ASP:N	2.49	0.45
1:A:79:LEU:HD23	1:A:79:LEU:HA	1.76	0.45
1:B:205:LEU:HD21	1:B:231:ALA:HA	1.98	0.45
1:B:99:LEU:HD23	1:B:116:MET:CE	2.46	0.45
1:C:55:GLU:HG2	1:C:56:GLY:N	2.31	0.45
1:B:12:LEU:HD12	1:B:12:LEU:HA	1.67	0.45
1:B:14:LYS:HD3	1:B:220:PRO:HB2	1.99	0.45
1:B:138:LYS:HZ1	1:B:240:LYS:NZ	2.15	0.44
1:A:226:THR:HG22	3:A:2069:HOH:O	2.16	0.44
1:A:28:TRP:CD1	1:A:28:TRP:N	2.85	0.44
1:B:149:ARG:HG3	2:E:53:VAL:CG2	2.48	0.44
1:B:212:LEU:O	1:B:216:THR:HG23	2.18	0.44
1:B:51:THR:O	1:B:245:GLY:HA3	2.18	0.44
1:B:80:LYS:HE2	3:B:2004:HOH:O	2.17	0.43
1:C:47:LEU:HD23	1:C:250:TYR:CD1	2.53	0.43
1:C:146:ARG:CG	1:C:146:ARG:HH11	2.31	0.43
1:A:12:LEU:HA	1:A:12:LEU:HD12	1.79	0.43
1:C:2:PHE:HA	1:C:61:ARG:O	2.18	0.43
1:A:162:CYS:HB3	1:A:167:VAL:HB	2.01	0.43
1:A:213:ASN:O	1:A:216:THR:OG1	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:GLU:O	1:C:194:ALA:N	2.47	0.42
2:E:52:PRO:O	2:E:53:VAL:HB	2.19	0.42
1:C:89:THR:HB	1:C:102:VAL:HB	2.00	0.42
1:C:28:TRP:CE2	1:C:37:LEU:HD12	2.54	0.42
1:A:147:ILE:HD11	1:A:182:LEU:HD11	2.01	0.42
1:C:192:GLU:N	1:C:192:GLU:OE1	2.39	0.42
1:C:191:GLU:C	1:C:193:GLU:H	2.21	0.42
1:A:38:GLN:CD	1:A:126:LEU:HB2	2.40	0.42
1:B:13:LYS:HE3	1:B:79:LEU:O	2.20	0.42
1:A:175:LEU:HD22	1:C:74:SER:HA	2.02	0.42
1:A:254:LYS:O	1:A:255:ILE:HG13	2.19	0.42
1:C:0:HIS:HB2	1:C:94:ASP:HB3	2.01	0.41
1:A:41:ASP:OD2	1:A:43:SER:OG	2.33	0.41
1:B:1:MET:HG3	1:B:92:ALA:O	2.19	0.41
1:C:28:TRP:HE3	1:C:35:VAL:CG1	2.33	0.41
1:B:190:LYS:CG	1:B:191:GLU:N	2.83	0.41
1:A:90:LEU:HD22	1:A:101:LEU:HG	2.02	0.41
1:A:74:SER:HB3	1:B:175:LEU:HD22	2.03	0.41
1:B:95:ASN:C	1:B:96:ALA:O	2.59	0.41
1:C:149:ARG:NH1	3:C:2040:HOH:O	2.52	0.41
1:C:207:PHE:CE1	1:C:235:LEU:HB2	2.56	0.41
1:A:240:LYS:HD3	1:A:240:LYS:HA	1.90	0.40
1:B:40:MET:HE1	1:B:44:HIS:HB3	2.04	0.40
1:C:147:ILE:HG23	1:C:180:ILE:HD12	2.02	0.40
1:B:151:LEU:HA	1:B:151:LEU:HD23	1.82	0.40
1:B:190:LYS:HG3	1:B:191:GLU:N	2.36	0.40
1:A:215:PHE:N	1:A:215:PHE:CD1	2.87	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:44:HIS:ND1	1:C:109:GLU:OE2[1_565]	2.07	0.13

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	256/264 (97%)	243 (95%)	10 (4%)	3 (1%)	15	24
1	B	254/264 (96%)	242 (95%)	9 (4%)	3 (1%)	15	24
1	C	256/264 (97%)	239 (93%)	10 (4%)	7 (3%)	6	8
2	D	19/21 (90%)	18 (95%)	1 (5%)	0	100	100
2	E	19/21 (90%)	16 (84%)	1 (5%)	2 (10%)	0	0
All	All	804/834 (96%)	758 (94%)	31 (4%)	15 (2%)	9	14

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	LEU
1	A	127	GLY
1	A	187	ASN
1	B	96	ALA
1	B	188	VAL
1	C	95	ASN
1	C	96	ALA
1	C	186	SER
2	E	56	ARG
1	C	187	ASN
2	E	53	VAL
1	B	42	SER
1	C	192	GLU
1	C	243	ASP
1	C	191	GLU

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	225/230 (98%)	216 (96%)	9 (4%)	36	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	224/230 (97%)	213 (95%)	11 (5%)	29	46
1	C	225/230 (98%)	213 (95%)	12 (5%)	26	42
2	D	16/16 (100%)	16 (100%)	0	100	100
2	E	16/16 (100%)	15 (94%)	1 (6%)	21	34
All	All	706/722 (98%)	673 (95%)	33 (5%)	30	48

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	16	LEU
1	A	32	SER
1	A	59	THR
1	A	126	LEU
1	A	181	LYS
1	A	210	ARG
1	A	233	VAL
1	A	255	ILE
1	B	97	ASP
1	B	112	SER
1	B	123	VAL
1	B	124	GLU
1	B	170	SER
1	B	174	GLU
1	B	186	SER
1	B	224	THR
1	B	243	ASP
1	B	244	MET
1	B	255	ILE
1	C	12	LEU
1	C	16	LEU
1	C	55	GLU
1	C	59	THR
1	C	109	GLU
1	C	117	LYS
1	C	123	VAL
1	C	131	GLN
1	C	132	GLU
1	C	181	LYS
1	C	186	SER
1	C	188	VAL

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Mol	Chain	Res	Type
2	E	55	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	184	GLN
1	C	153	HIS

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	258/264 (97%)	-0.13	4 (1%) 72 72	27, 48, 99, 132	0
1	B	256/264 (96%)	-0.10	6 (2%) 61 59	25, 45, 99, 169	0
1	C	258/264 (97%)	-0.05	9 (3%) 44 43	22, 46, 99, 155	0
2	D	21/21 (100%)	-0.44	0 100 100	32, 50, 68, 78	0
2	E	21/21 (100%)	0.51	2 (9%) 9 7	45, 63, 89, 121	0
All	All	814/834 (97%)	-0.09	21 (2%) 56 55	22, 47, 99, 169	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	190	LYS	6.0
1	C	188	VAL	5.2
2	E	55	VAL	4.4
1	C	190	LYS	4.0
1	C	95	ASN	4.0
1	A	165	ASP	3.3
1	C	94	ASP	3.1
1	C	193	GLU	2.7
1	A	232	ASP	2.7
1	B	187	ASN	2.6
1	A	123	VAL	2.5
1	C	-1	PRO	2.5
1	B	188	VAL	2.4
1	A	132	GLU	2.3
1	B	189	ASP	2.3
1	C	-2	GLY	2.2
2	E	54	CYS	2.2
1	B	123	VAL	2.2
1	C	189	ASP	2.1
1	C	187	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	122	ASP	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.