



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 12:21 PM UTC

PDB ID : 2P5O / pdb_00002p5o
Title : Crystal structure of RB69 GP43 in complex with DNA containing an abasic site analog
Authors : Hogg, M.; Wallace, S.S.; Doublié, S.
Deposited on : 2007-03-15
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

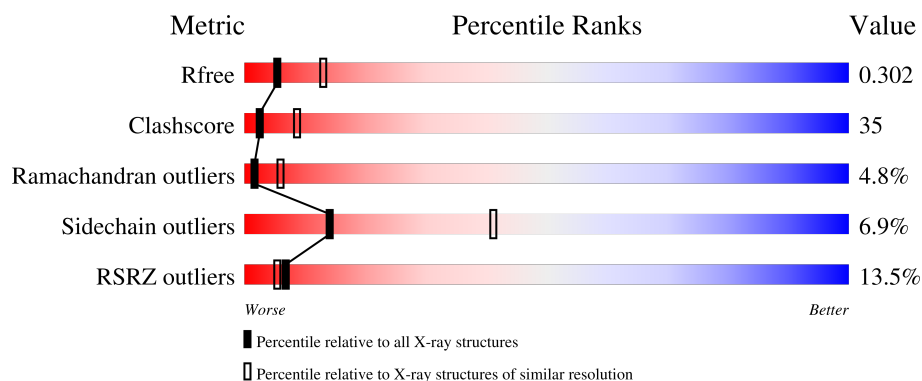
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	E	18	
1	G	18	
1	I	18	
1	K	18	
2	F	15	

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Mol	Chain	Length	Quality of chain
2	H	15	47% 40% 60%
2	J	15	20% 33% 67%
2	L	15	33% 7% 47% 47%
3	A	903	3% 53% 36% 5% 5%
3	B	903	12% 37% 41% 7% 15%
3	C	903	3% 48% 40% 7% 6%
3	D	903	26% 28% 38% 8% 26%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 27752 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 DNA chain called Template DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	18	Total	C	N	O	P	0	0	0
			355	169	64	105	17			
1	G	13	Total	C	N	O	P	0	0	0
			264	126	51	75	12			
1	I	18	Total	C	N	O	P	0	0	0
			355	169	64	105	17			
1	K	9	Total	C	N	O	P	0	0	0
			181	86	37	50	8			

- Molecule 2 is a DNA chain called Primer DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	15	Total	C	N	O	P	0	0	0
			308	147	60	87	14			
2	H	15	Total	C	N	O	P	0	0	0
			308	147	60	87	14			
2	J	15	Total	C	N	O	P	0	0	0
			308	147	60	87	14			
2	L	8	Total	C	N	O	P	0	0	0
			163	78	30	48	7			

- Molecule 3 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	A	856	Total	C	N	O	S	Se	0	0	0
			6941	4460	1154	1295	8	24			
3	B	771	Total	C	N	O	S	Se	0	0	0
			6245	4013	1034	1167	6	25			
3	C	853	Total	C	N	O	S	Se	0	0	0
			6895	4430	1143	1290	8	24			
3	D	671	Total	C	N	O	S	Se	0	0	0
			4934	3144	818	948	6	18			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	modified residue	UNP Q38087
A	65	MSE	MET	modified residue	UNP Q38087
A	75	MSE	MET	modified residue	UNP Q38087
A	85	MSE	MET	modified residue	UNP Q38087
A	189	MSE	MET	modified residue	UNP Q38087
A	199	MSE	MET	modified residue	UNP Q38087
A	222	ALA	ASP	engineered mutation	UNP Q38087
A	256	MSE	MET	modified residue	UNP Q38087
A	327	ALA	ASP	engineered mutation	UNP Q38087
A	347	MSE	MET	modified residue	UNP Q38087
A	408	MSE	MET	modified residue	UNP Q38087
A	461	MSE	MET	modified residue	UNP Q38087
A	462	MSE	MET	modified residue	UNP Q38087
A	489	MSE	MET	modified residue	UNP Q38087
A	541	MSE	MET	modified residue	UNP Q38087
A	553	MSE	MET	modified residue	UNP Q38087
A	592	MSE	MET	modified residue	UNP Q38087
A	659	MSE	MET	modified residue	UNP Q38087
A	670	MSE	MET	modified residue	UNP Q38087
A	674	MSE	MET	modified residue	UNP Q38087
A	681	MSE	MET	modified residue	UNP Q38087
A	683	MSE	MET	modified residue	UNP Q38087
A	715	MSE	MET	modified residue	UNP Q38087
A	728	MSE	MET	modified residue	UNP Q38087
A	752	MSE	MET	modified residue	UNP Q38087
A	866	MSE	MET	modified residue	UNP Q38087
A	900	MSE	MET	modified residue	UNP Q38087
B	1	MSE	-	modified residue	UNP Q38087
B	65	MSE	MET	modified residue	UNP Q38087
B	75	MSE	MET	modified residue	UNP Q38087
B	85	MSE	MET	modified residue	UNP Q38087
B	189	MSE	MET	modified residue	UNP Q38087
B	199	MSE	MET	modified residue	UNP Q38087
B	222	ALA	ASP	engineered mutation	UNP Q38087
B	256	MSE	MET	modified residue	UNP Q38087
B	327	ALA	ASP	engineered mutation	UNP Q38087
B	347	MSE	MET	modified residue	UNP Q38087
B	408	MSE	MET	modified residue	UNP Q38087
B	461	MSE	MET	modified residue	UNP Q38087
B	462	MSE	MET	modified residue	UNP Q38087
B	489	MSE	MET	modified residue	UNP Q38087
B	541	MSE	MET	modified residue	UNP Q38087

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Chain	Residue	Modelled	Actual	Comment	Reference
B	553	MSE	MET	modified residue	UNP Q38087
B	592	MSE	MET	modified residue	UNP Q38087
B	659	MSE	MET	modified residue	UNP Q38087
B	670	MSE	MET	modified residue	UNP Q38087
B	674	MSE	MET	modified residue	UNP Q38087
B	681	MSE	MET	modified residue	UNP Q38087
B	683	MSE	MET	modified residue	UNP Q38087
B	715	MSE	MET	modified residue	UNP Q38087
B	728	MSE	MET	modified residue	UNP Q38087
B	752	MSE	MET	modified residue	UNP Q38087
B	866	MSE	MET	modified residue	UNP Q38087
B	900	MSE	MET	modified residue	UNP Q38087
C	1	MSE	-	modified residue	UNP Q38087
C	65	MSE	MET	modified residue	UNP Q38087
C	75	MSE	MET	modified residue	UNP Q38087
C	85	MSE	MET	modified residue	UNP Q38087
C	189	MSE	MET	modified residue	UNP Q38087
C	199	MSE	MET	modified residue	UNP Q38087
C	222	ALA	ASP	engineered mutation	UNP Q38087
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C	327	ALA	ASP	engineered mutation	UNP Q38087
C	347	MSE	MET	modified residue	UNP Q38087
C	408	MSE	MET	modified residue	UNP Q38087
C	461	MSE	MET	modified residue	UNP Q38087
C	462	MSE	MET	modified residue	UNP Q38087
C	489	MSE	MET	modified residue	UNP Q38087
C	541	MSE	MET	modified residue	UNP Q38087
C	553	MSE	MET	modified residue	UNP Q38087
C	592	MSE	MET	modified residue	UNP Q38087
C	659	MSE	MET	modified residue	UNP Q38087
C	670	MSE	MET	modified residue	UNP Q38087
C	674	MSE	MET	modified residue	UNP Q38087
C	681	MSE	MET	modified residue	UNP Q38087
C	683	MSE	MET	modified residue	UNP Q38087
C	715	MSE	MET	modified residue	UNP Q38087
C	728	MSE	MET	modified residue	UNP Q38087
C	752	MSE	MET	modified residue	UNP Q38087
C	866	MSE	MET	modified residue	UNP Q38087
C	900	MSE	MET	modified residue	UNP Q38087
D	1	MSE	-	modified residue	UNP Q38087
D	65	MSE	MET	modified residue	UNP Q38087
D	75	MSE	MET	modified residue	UNP Q38087

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Chain	Residue	Modelled	Actual	Comment	Reference
D	85	MSE	MET	modified residue	UNP Q38087
D	189	MSE	MET	modified residue	UNP Q38087
D	199	MSE	MET	modified residue	UNP Q38087
D	222	ALA	ASP	engineered mutation	UNP Q38087
D	256	MSE	MET	modified residue	UNP Q38087
D	327	ALA	ASP	engineered mutation	UNP Q38087
D	347	MSE	MET	modified residue	UNP Q38087
D	408	MSE	MET	modified residue	UNP Q38087
D	461	MSE	MET	modified residue	UNP Q38087
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D	489	MSE	MET	modified residue	UNP Q38087
D	541	MSE	MET	modified residue	UNP Q38087
D	553	MSE	MET	modified residue	UNP Q38087
D	592	MSE	MET	modified residue	UNP Q38087
D	659	MSE	MET	modified residue	UNP Q38087
D	670	MSE	MET	modified residue	UNP Q38087
D	674	MSE	MET	modified residue	UNP Q38087
D	681	MSE	MET	modified residue	UNP Q38087
D	683	MSE	MET	modified residue	UNP Q38087
D	715	MSE	MET	modified residue	UNP Q38087
D	728	MSE	MET	modified residue	UNP Q38087
D	752	MSE	MET	modified residue	UNP Q38087
D	866	MSE	MET	modified residue	UNP Q38087
D	900	MSE	MET	modified residue	UNP Q38087

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	3	Total O 3 3	0	0
4	F	5	Total O 5 5	0	0
4	G	7	Total O 7 7	0	0
4	H	2	Total O 2 2	0	0
4	I	17	Total O 17 17	0	0
4	J	9	Total O 9 9	0	0
4	K	8	Total O 8 8	0	0

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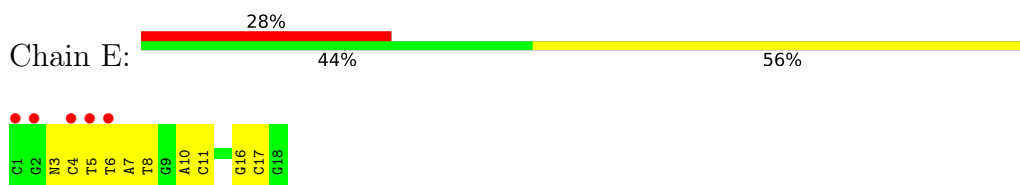
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	6	Total 6	O 6	0	0
4	A	139	Total 139	O 139	0	0
4	B	106	Total 106	O 106	0	0
4	C	147	Total 147	O 147	0	0
4	D	46	Total 46	O 46	0	0

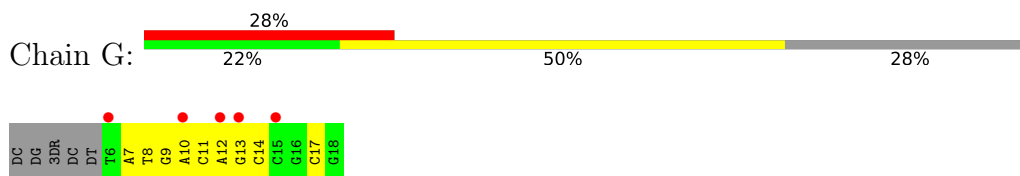
3 Residue-property plots [i](#)

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.

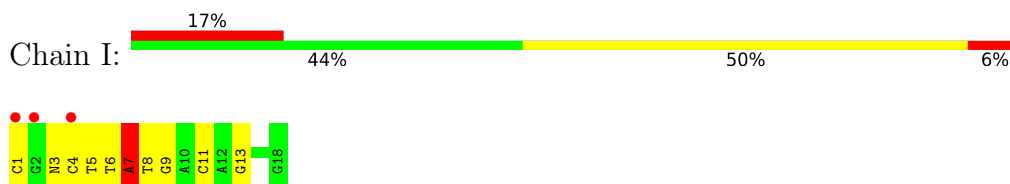
- Molecule 1: Template DNA



- Molecule 1: Template DNA



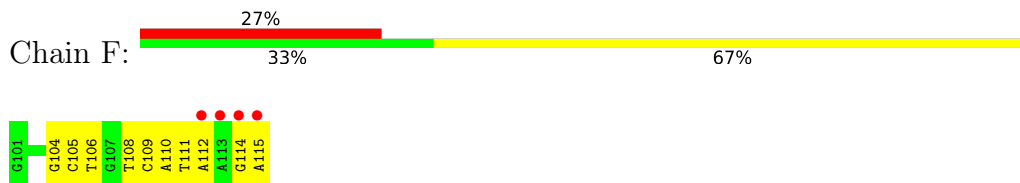
- Molecule 1: Template DNA



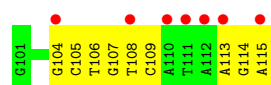
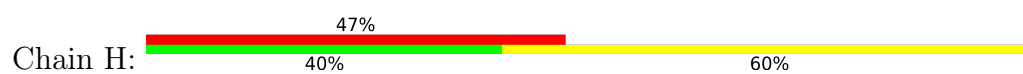
- Molecule 1: Template DNA



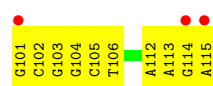
- Molecule 2: Primer DNA



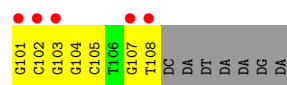
- Molecule 2: Primer DNA



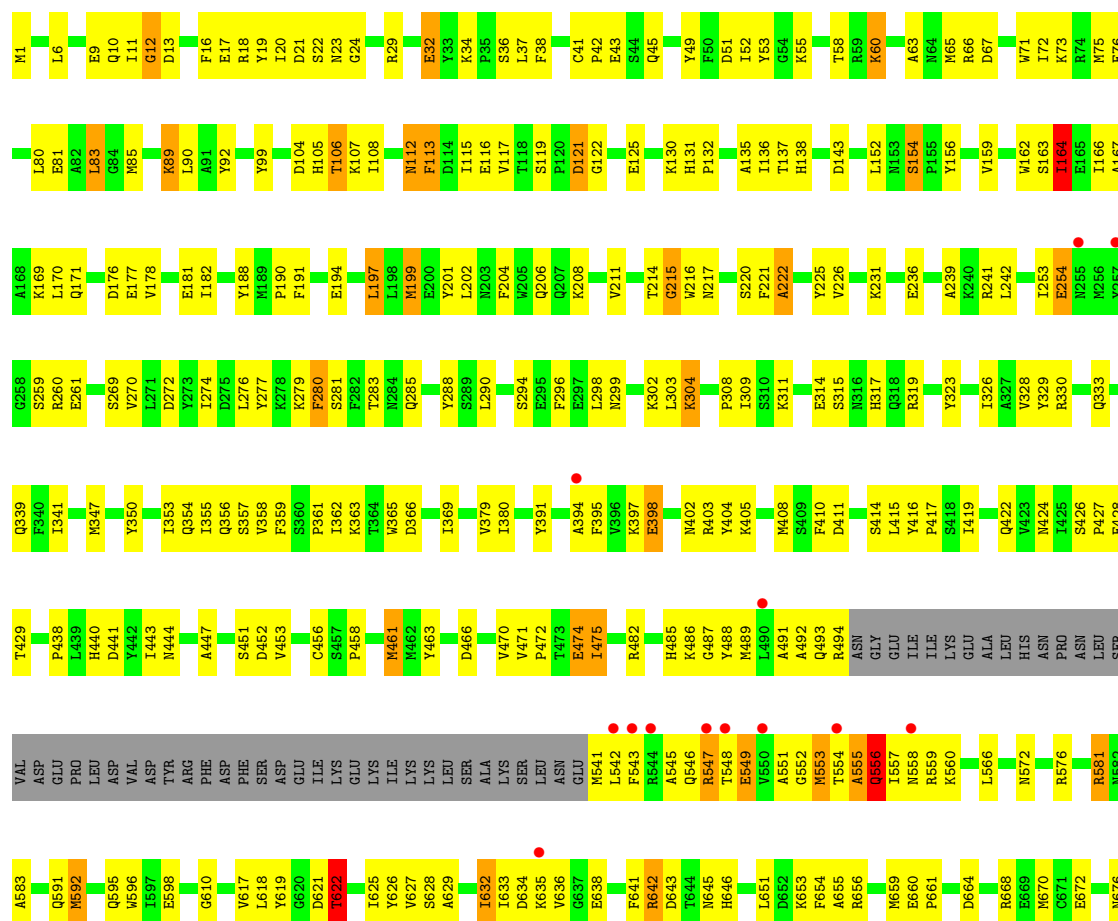
• Molecule 2: Primer DNA

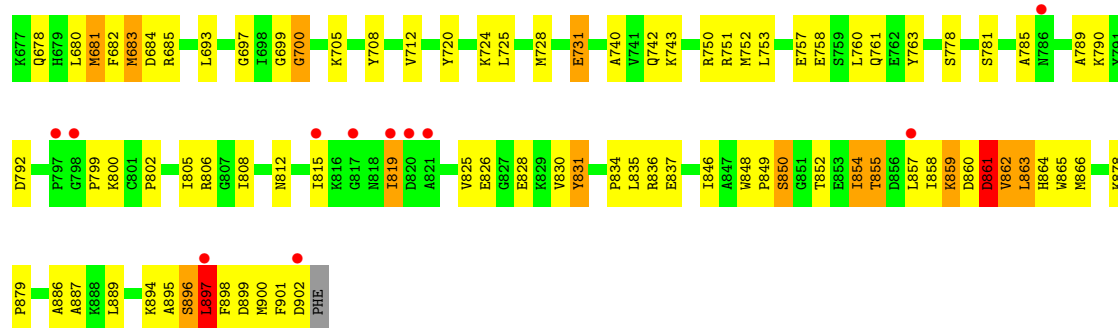


• Molecule 2: Primer DNA

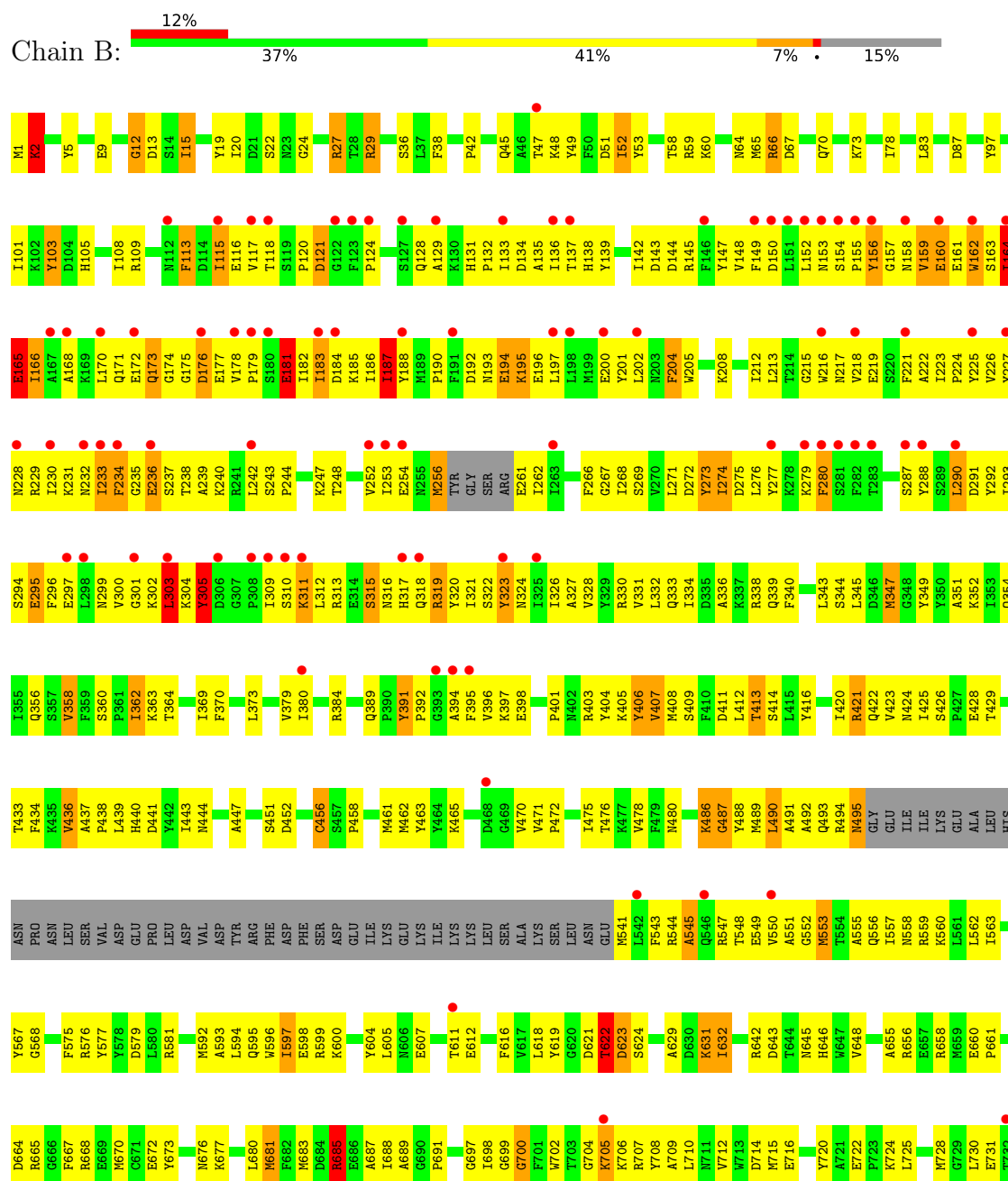


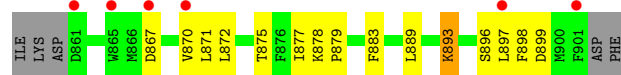
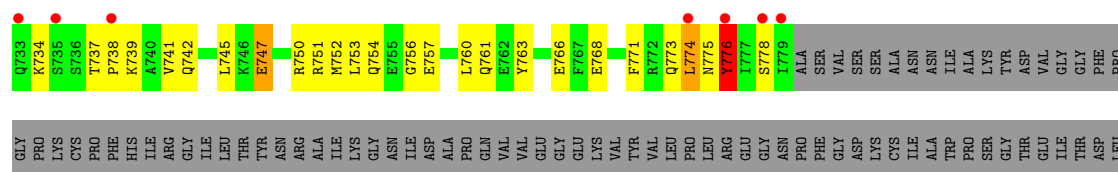
• Molecule 3: DNA polymerase



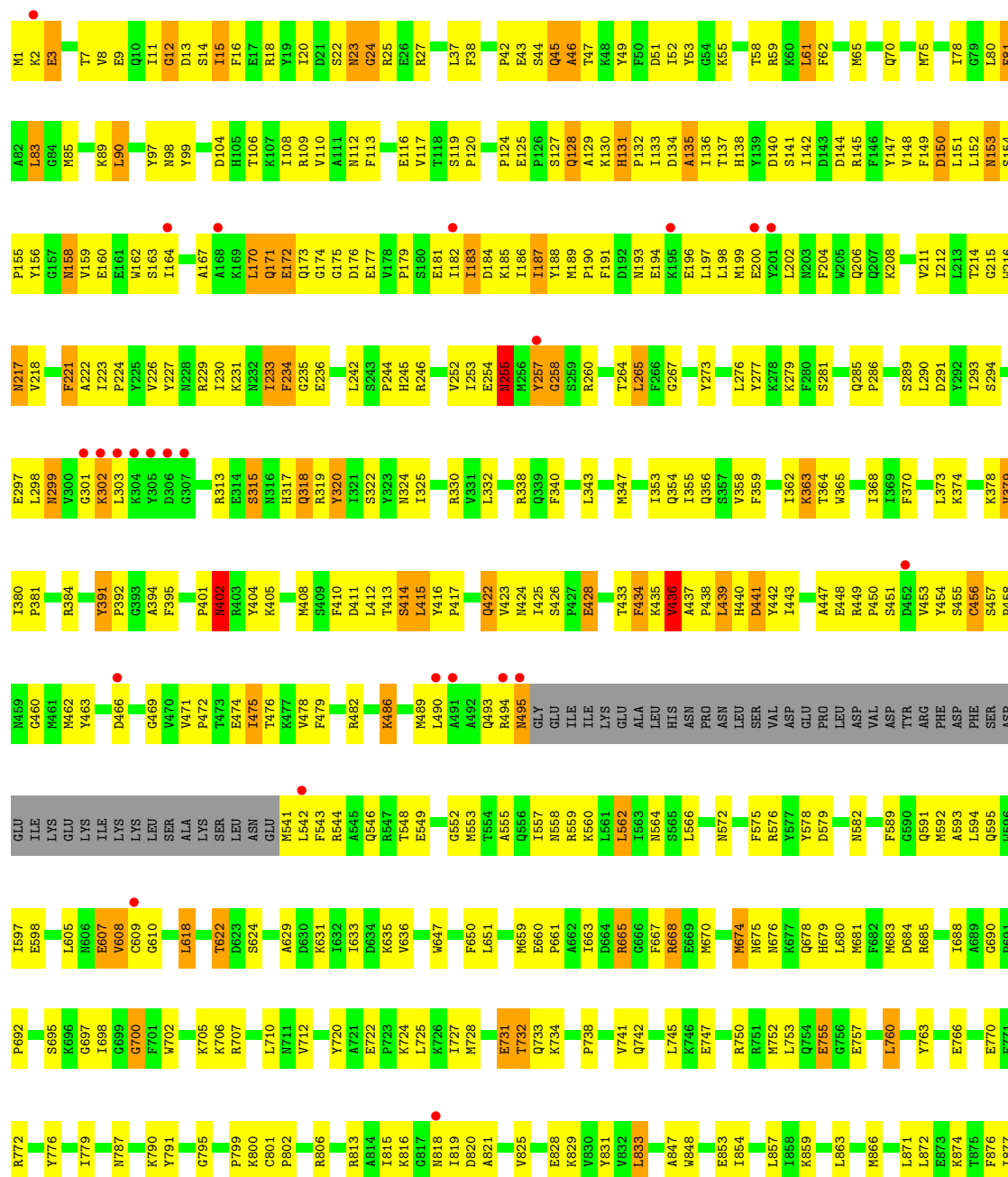


• Molecule 3: DNA polymerase



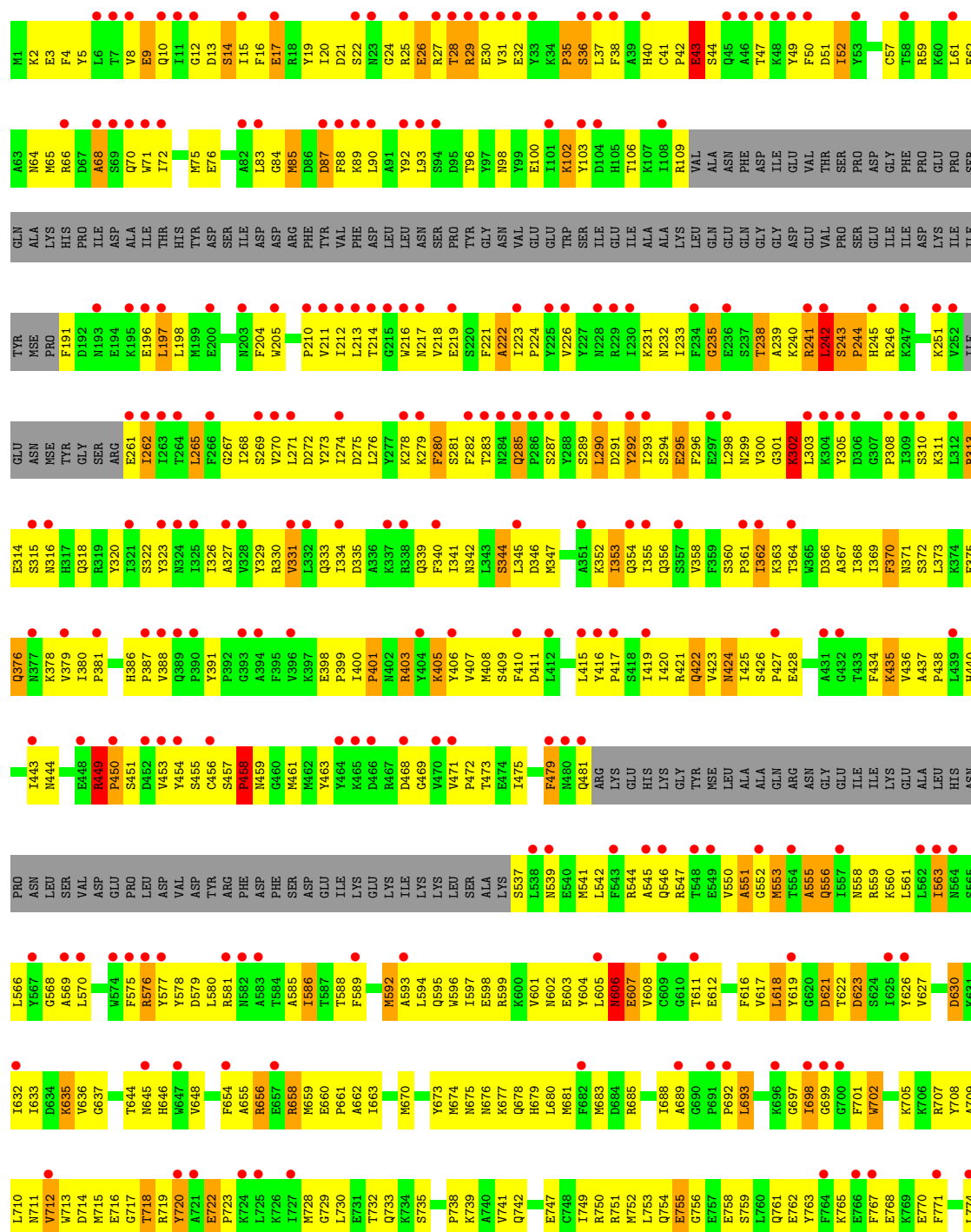
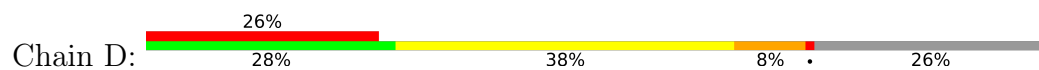


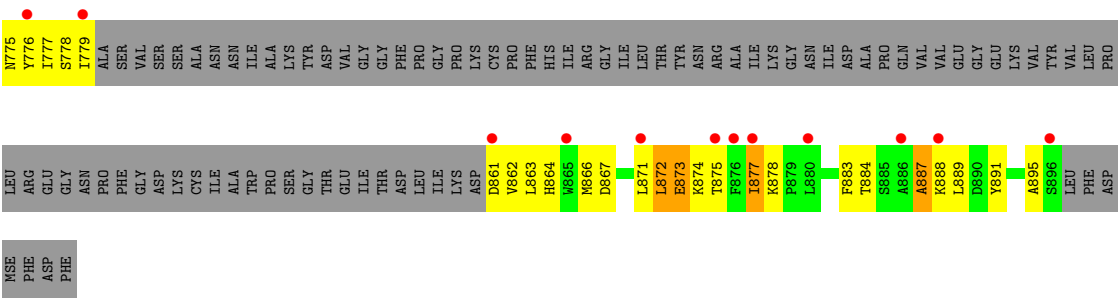
• Molecule 3: DNA polymerase





● Molecule 3: DNA polymerase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	133.02Å 123.21Å 165.62Å 90.00° 95.82° 90.00°	Depositor
Resolution (Å)	44.00 – 2.80 44.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.9 (44.00-2.80) 92.1 (44.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.245 , 0.290 0.263 , 0.302	Depositor DCC
R_{free} test set	24426 reflections (9.67%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	27752	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.47% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 3DR

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	E	0.30	0/384	0.61	0/588
1	G	0.31	0/296	0.72	0/455
1	I	0.39	0/384	0.83	2/588 (0.3%)
1	K	0.30	0/203	0.55	0/311
2	F	0.28	0/346	0.58	0/533
2	H	0.27	0/346	0.59	0/533
2	J	0.33	0/346	0.67	0/533
2	L	0.26	0/182	0.54	0/280
3	A	0.57	0/7090	1.05	32/9555 (0.3%)
3	B	0.50	0/6376	0.97	23/8593 (0.3%)
3	C	0.54	0/7045	1.01	28/9502 (0.3%)
3	D	0.39	0/5028	0.92	19/6831 (0.3%)
All	All	0.50	0/28026	0.96	104/38302 (0.3%)

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	G	0	1
1	I	0	1
3	B	0	1
All	All	0	3

There are no bond length outliers.

The worst 5 of 104 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	819	ILE	N-CA-C	-9.92	103.38	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	87	ASP	N-CA-C	-8.81	101.79	112.54
3	B	486	LYS	N-CA-C	-8.64	102.66	113.55
3	A	456	CYS	N-CA-C	8.26	122.55	109.50
3	A	211	VAL	N-CA-C	-8.12	103.96	111.67

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	273	TYR	Sidechain
1	G	17	DC	Sidechain
1	I	7	DA	Sidechain

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	E	355	0	200	18	0
1	G	264	0	147	12	0
1	I	355	0	200	13	0
1	K	181	0	101	15	0
2	F	308	0	170	20	0
2	H	308	0	170	18	0
2	J	308	0	170	13	0
2	L	163	0	92	24	0
3	A	6941	0	6787	328	0
3	B	6245	0	6024	515	0
3	C	6895	0	6710	406	0
3	D	4934	0	4297	454	0
4	A	139	0	0	17	0
4	B	106	0	0	24	0
4	C	147	0	0	16	0
4	D	46	0	0	9	0
4	E	3	0	0	1	0
4	F	5	0	0	0	0
4	G	7	0	0	0	0
4	H	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	17	0	0	0	0
4	J	9	0	0	0	0
4	K	8	0	0	1	0
4	L	6	0	0	2	0
All	All	27752	0	25068	1810	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 35.

The worst 5 of 1810 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:L:107:DG:C2'	2:L:108:DT:H71	1.77	1.14
1:G:12:DA:H2''	1:G:13:DG:H5'	1.30	1.14
3:D:14:SER:HA	3:D:32:GLU:HA	1.32	1.11
3:B:164:ILE:HD12	3:B:164:ILE:H	1.10	1.10
1:E:16:DG:H2''	1:E:17:DC:H5'	1.27	1.09

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	
3	A	852/903 (94%)	760 (89%)	76 (9%)	16 (2%)	6	22
3	B	763/903 (84%)	621 (81%)	110 (14%)	32 (4%)	2	7
3	C	849/903 (94%)	745 (88%)	74 (9%)	30 (4%)	3	10
3	D	661/903 (73%)	433 (66%)	155 (23%)	73 (11%)	0	1
All	All	3125/3612 (86%)	2559 (82%)	415 (13%)	151 (5%)	2	6

5 of 151 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	12	GLY
3	A	542	LEU
3	A	859	LYS
3	A	894	LYS
3	B	164	ILE

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	745/775 (96%)	700 (94%)	45 (6%)	17	47
3	B	663/775 (86%)	608 (92%)	55 (8%)	10	32
3	C	737/775 (95%)	692 (94%)	45 (6%)	17	46
3	D	453/775 (58%)	419 (92%)	34 (8%)	12	37
All	All	2598/3100 (84%)	2419 (93%)	179 (7%)	14	41

5 of 179 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	255	ASN
3	C	829	LYS
3	C	299	ASN
3	C	441	ASP
3	D	85	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 81 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	354	GLN
3	D	70	GLN
3	C	377	ASN
3	C	546	GLN
3	D	606	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	3DR	E	3	1	8,11,12	0.33	0	7,14,17	0.83	0
1	3DR	I	3	1	8,11,12	0.31	0	7,14,17	0.87	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	3DR	E	3	1	-	0/3/15/16	0/1/1/1
1	3DR	I	3	1	-	1/3/15/16	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	I	3	3DR	O4'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	3	3DR	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	I	3	3DR	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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, 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	E	17/18 (94%)	1.10	5 (29%) 1 1	57, 95, 121, 134	0
1	G	13/18 (72%)	1.71	5 (38%) 1 0	48, 103, 143, 144	0
1	I	17/18 (94%)	0.14	3 (17%) 4 3	27, 43, 112, 123	0
1	K	9/18 (50%)	1.72	5 (55%) 0 0	38, 140, 155, 159	0
2	F	15/15 (100%)	1.41	4 (26%) 1 1	77, 104, 136, 140	0
2	H	15/15 (100%)	1.80	7 (46%) 0 0	75, 118, 131, 139	0
2	J	15/15 (100%)	0.59	3 (20%) 3 2	28, 57, 97, 114	0
2	L	8/15 (53%)	2.11	5 (62%) 0 0	144, 148, 149, 149	0
3	A	831/903 (92%)	0.15	24 (2%) 53 43	16, 40, 83, 133	0
3	B	746/903 (82%)	0.80	104 (13%) 6 5	19, 57, 112, 140	0
3	C	829/903 (91%)	0.29	25 (3%) 52 42	18, 46, 86, 119	0
3	D	650/903 (71%)	1.77	238 (36%) 1 0	78, 102, 134, 142	0
All	All	3165/3744 (84%)	0.71	428 (13%) 7 5	16, 56, 123, 159	0

The worst 5 of 428 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	896	SER	6.9
3	B	779	ILE	6.2
3	D	90	LEU	5.7
3	D	216	TRP	5.5
3	D	569	ALA	5.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	3DR	E	3	11/12	0.79	0.24	117,125,130,133	0
1	3DR	I	3	11/12	0.89	0.18	101,102,103,103	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.