



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:02 AM GMT

PDB ID : 2P6R
Title : Crystal structure of superfamily 2 helicase Hel308 in complex with unwound DNA
Authors : Buettner, K.; Nehring, S.; Hopfner, K.P.
Deposited on : 2007-03-19
Resolution : 3.00 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
<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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

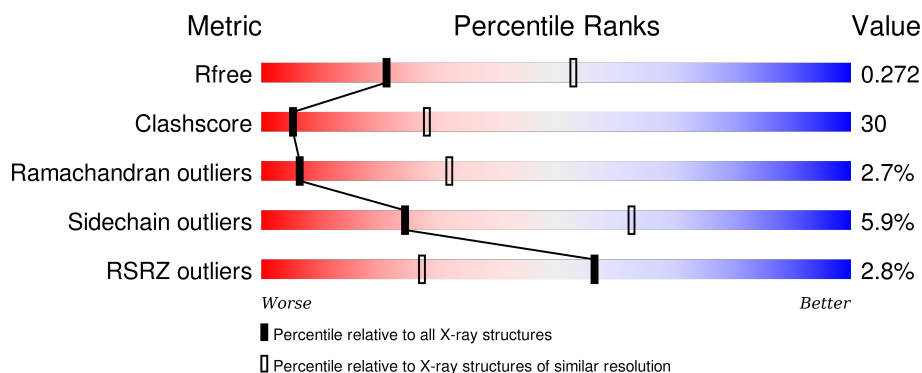
1 Overall quality at a glance

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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (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 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	X	25	<div> <div>36%</div> <div>28%72%</div> </div>
2	Y	15	<div> <div>47%</div> <div>60%33%7%</div> </div>
3	A	702	<div> <div>%</div> <div>48%45%. .</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6219 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 25-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	25	Total	C	N	O	P	0	0	0
			508	244	92	148	24			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Y	15	Total	C	N	O	P	0	0	0
			305	147	57	87	14			

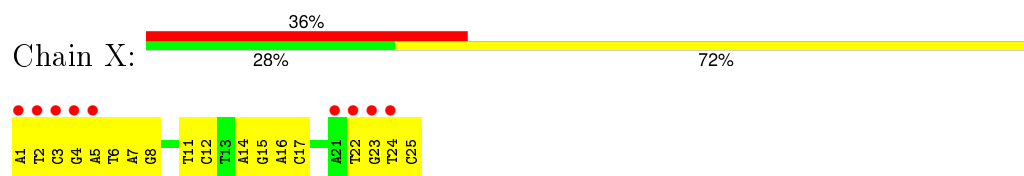
- Molecule 3 is a protein called afUHEL308 HELICASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	683	Total	C	N	O	S	57	0	0
			5406	3417	953	1014	22			

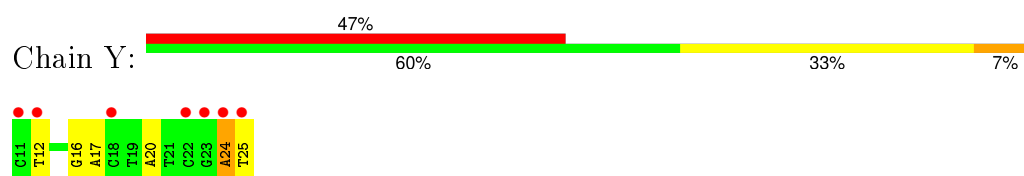
3 Residue-property plots

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.

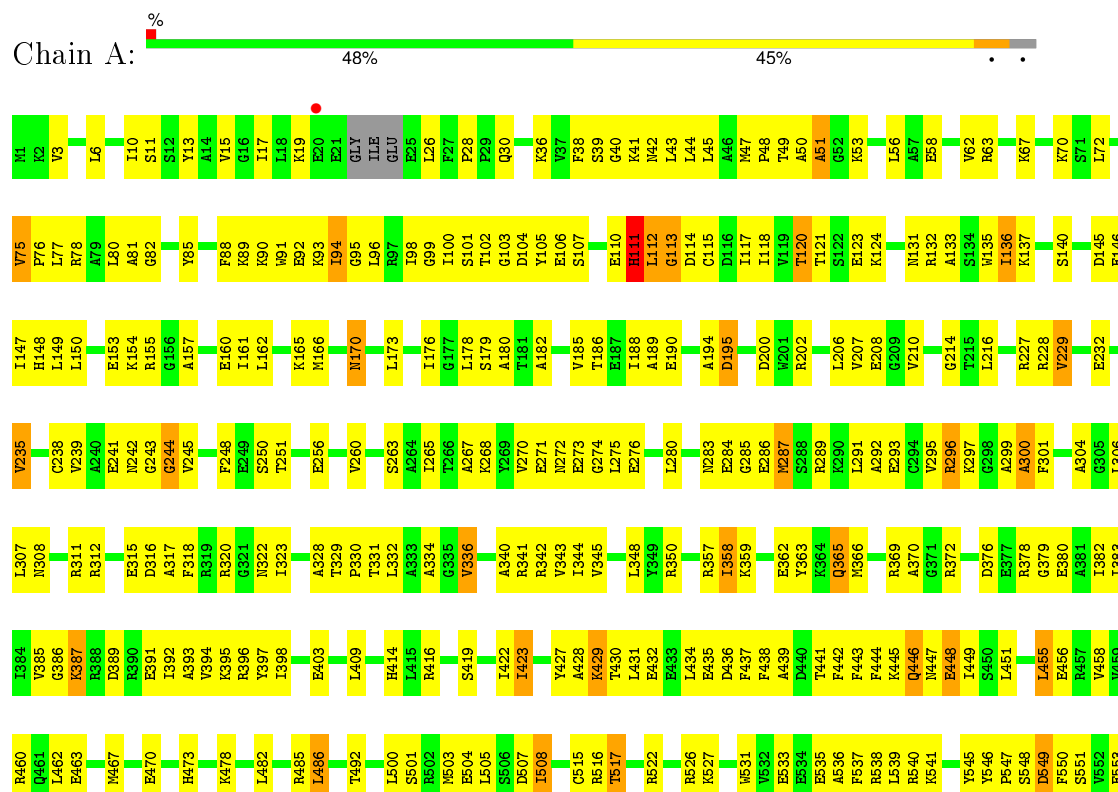
• Molecule 1: 25-MER



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



• Molecule 3: afUHEL308 HELICASE



Y554	D555	W556	F557	L558	S559	E560	Y561	C566	L567	K568	D569	E573	K574	D575	E578	I579	Y583	A586	P587	S588	D589	L590	I593	V594	M604	M605	A608	V611	V616	E621	R622	G626	V627	E630	L631	L632	R636	I637	R638	H639	T640	G641	R642	V643	R644
A645	R646	Y649	I653	R654	M655	V659	V660	R661	R662	R663	E664	K665	G673	I674	A675	E676	R677	V678	V679	I682	S683	V684	K685	S686	LEU	ASN	PRO	GLU	SER	ALA	ALA	ALA	LEU	GLU	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	138.38 Å 138.38 Å 252.97 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.00 29.96 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.9 (20.00-3.00) 100.0 (29.96-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 3.00 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.275 0.234 , 0.272	Depositor DCC
R_{free} test set	2939 reflections (10.04%)	DCC
Wilson B-factor (Å ²)	65.8	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 29357 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6219	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	X	0.62	0/569	0.75	0/876
2	Y	0.71	0/342	0.79	0/526
3	A	0.43	0/5491	0.65	0/7396
All	All	0.47	0/6402	0.67	0/8798

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	Y	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	Y	24	DA	Sidechain

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	X	508	0	284	33	0
2	Y	305	0	171	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	5406	0	5498	337	0
All	All	6219	0	5953	359	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 30.

The worst 5 of 359 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:X:25:DC:H5''	3:A:643:VAL:HB	1.28	1.10
3:A:49:THR:HG22	3:A:51:ALA:H	1.16	1.05
3:A:363:TYR:CE1	3:A:383:ILE:HD11	1.95	1.00
3:A:675:ALA:O	3:A:679:VAL:HG23	1.67	0.95
3:A:77:LEU:HB2	3:A:80:LEU:HD13	1.47	0.94

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	679/702 (97%)	595 (88%)	66 (10%)	18 (3%)	6 32

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	51	ALA
3	A	286	GLU
3	A	3	VAL
3	A	30	GLN
3	A	94	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	576 / 591 (98%)	542 (94%)	34 (6%)	24 63

5 of 34 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	358	ILE
3	A	446	GLN
3	A	611	VAL
3	A	387	LYS
3	A	170	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
3	A	322	ASN
3	A	414	HIS
3	A	461	GLN
3	A	302	HIS
3	A	447	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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	X	25/25 (100%)	1.46	9 (36%) 0 0	40, 78, 110, 113	0
2	Y	15/15 (100%)	2.97	7 (46%) 0 0	82, 102, 112, 113	0
3	A	681/702 (97%)	-0.34	4 (0%) 90 73	19, 53, 88, 109	13 (1%)
All	All	721/742 (97%)	-0.21	20 (2%) 56 27	19, 54, 94, 113	13 (1%)

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	Y	24	DA	9.0
2	Y	23	DG	7.8
2	Y	25	DT	7.6
1	X	1	DA	6.3
1	X	23	DG	5.8

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.