



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 5, 2019 – 12:23 PM EST

PDB ID : 1WP9
Title : Crystal structure of Pyrococcus furiosus Hef helicase domain
Authors : Nishino, T.; Komori, K.; Tsuchiya, D.; Ishino, Y.; Morikawa, K.
Deposited on : 2004-08-31
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix) : 1.13
EDS : 2.4
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.4

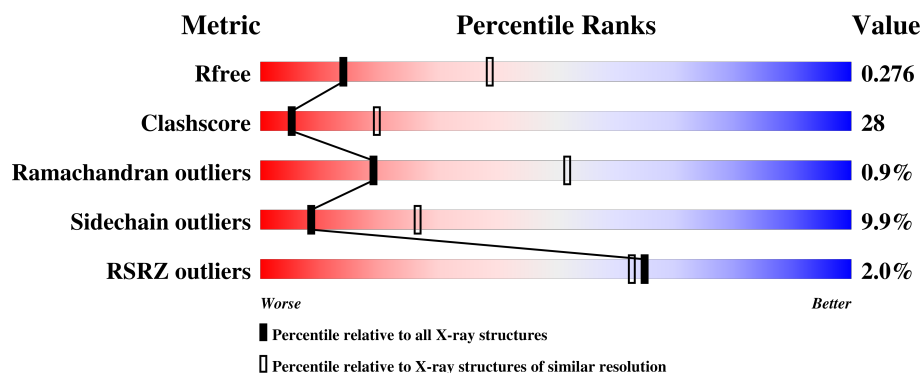
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	111664	1716 (2.90-2.90)
Clashscore	122126	1924 (2.90-2.90)
Ramachandran outliers	120053	1884 (2.90-2.90)
Sidechain outliers	120020	1886 (2.90-2.90)
RSRZ outliers	108989	1669 (2.90-2.90)

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	494	<div> <div>4%</div> <div>51% 40% 6% .</div> </div>
1	B	494	<div> <div>4%</div> <div>51% 38% 7% .</div> </div>
1	C	494	<div> <div>4%</div> <div>47% 43% 6% .</div> </div>
1	D	494	<div> <div>4%</div> <div>50% 40% 5% 5%</div> </div>
1	E	494	<div> <div>3%</div> <div>49% 41% 6% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	494	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	495	-	-	X	-
2	PO4	C	495	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22913 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.

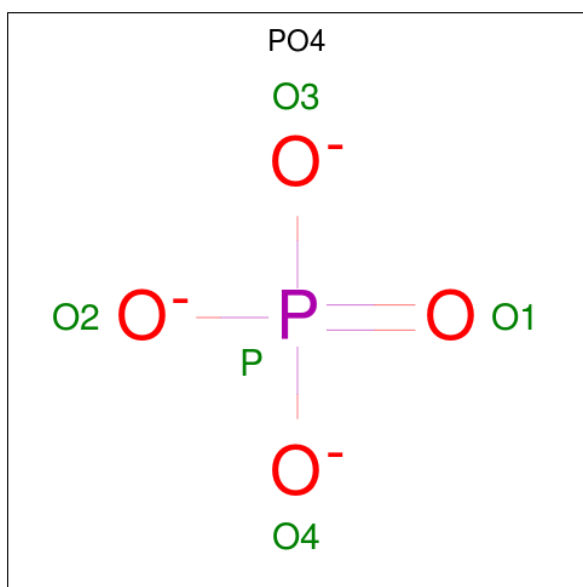
- Molecule 1 is a protein called ATP-dependent RNA helicase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	0	0
			3831	2445	683	689	14			
1	B	477	Total	C	N	O	S	0	0	0
			3817	2436	680	687	14			
1	C	475	Total	C	N	O	S	0	0	0
			3807	2431	678	684	14			
1	D	469	Total	C	N	O	S	0	0	0
			3765	2407	670	674	14			
1	E	475	Total	C	N	O	S	0	0	0
			3807	2431	678	684	14			
1	F	475	Total	C	N	O	S	0	0	0
			3807	2431	678	684	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	LEU	ENGINEERED	UNP Q8TZH8
B	2	VAL	LEU	ENGINEERED	UNP Q8TZH8
C	2	VAL	LEU	ENGINEERED	UNP Q8TZH8
D	2	VAL	LEU	ENGINEERED	UNP Q8TZH8
E	2	VAL	LEU	ENGINEERED	UNP Q8TZH8
F	2	VAL	LEU	ENGINEERED	UNP Q8TZH8

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



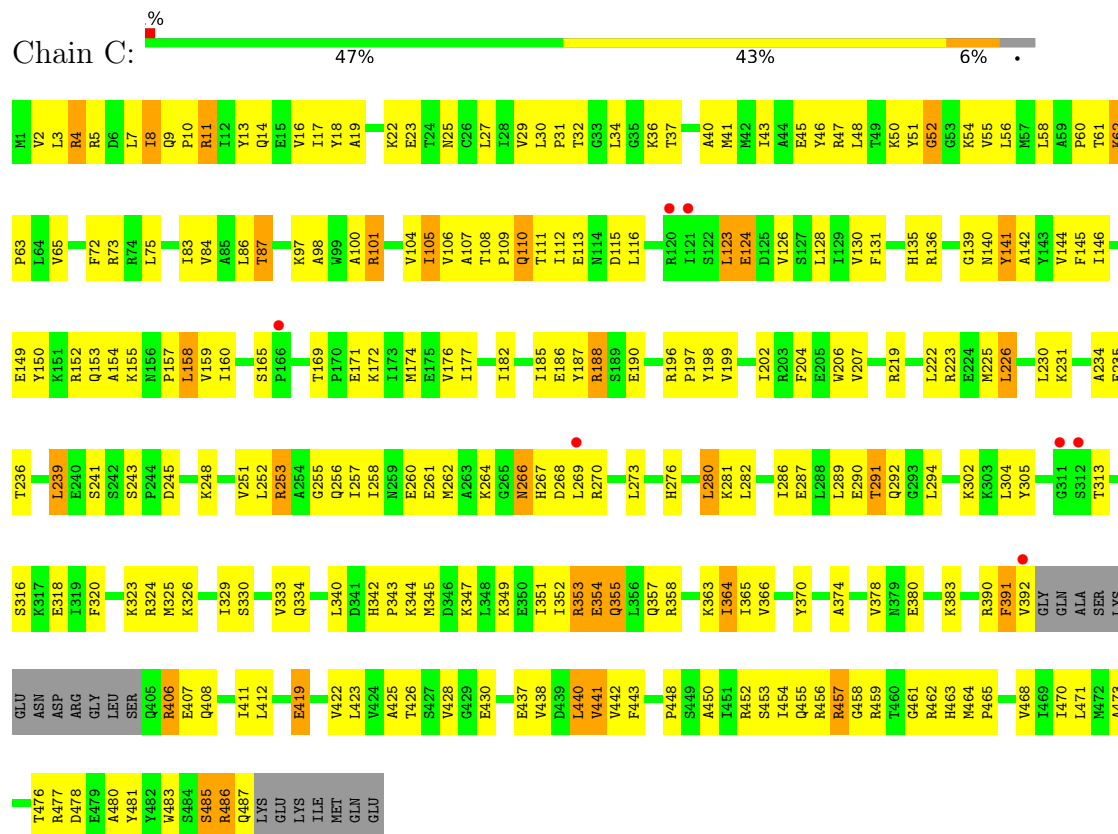
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

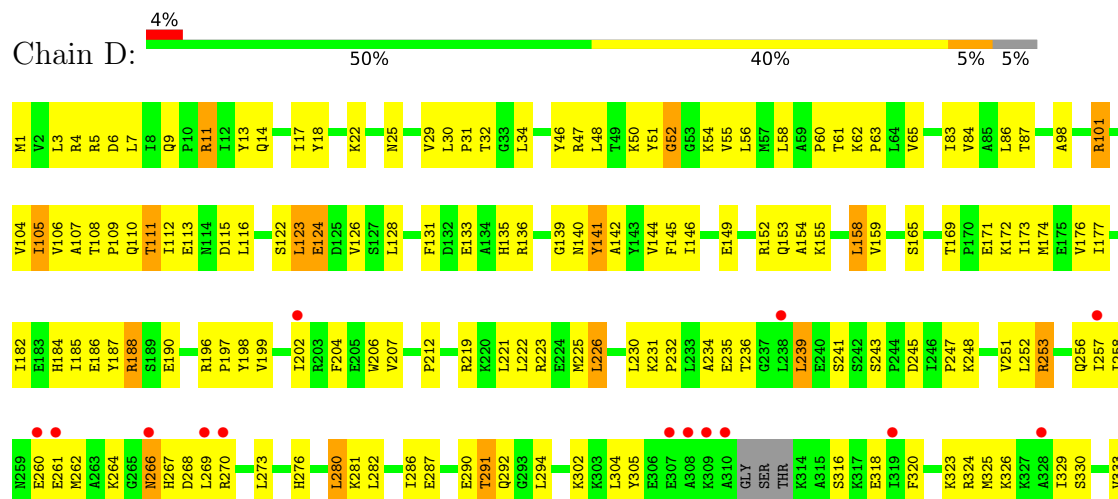
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	B	13	Total	O	0	0
			13	13		
3	C	8	Total	O	0	0
			8	8		
3	D	7	Total	O	0	0
			7	7		
3	E	13	Total	O	0	0
			13	13		
3	F	4	Total	O	0	0
			4	4		

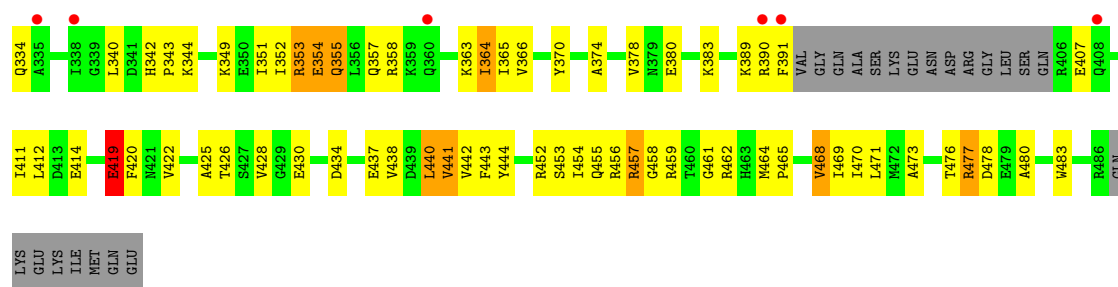


- Molecule 1: ATP-dependent RNA helicase, putative

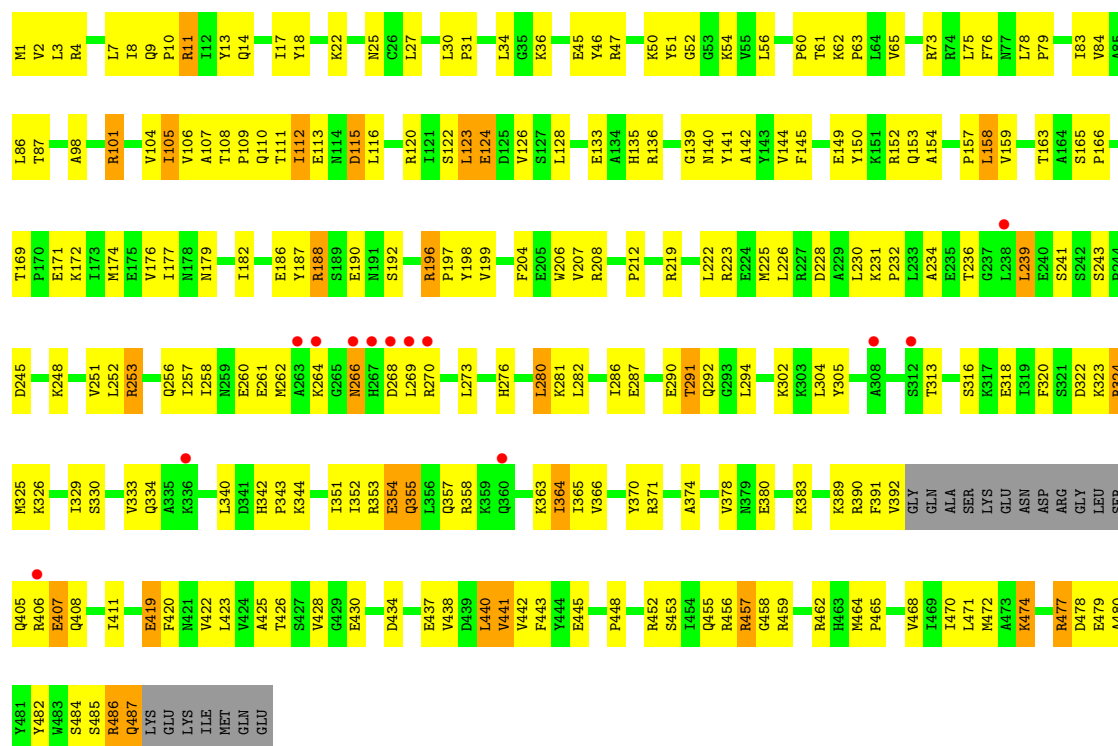


- Molecule 1: ATP-dependent RNA helicase, putative

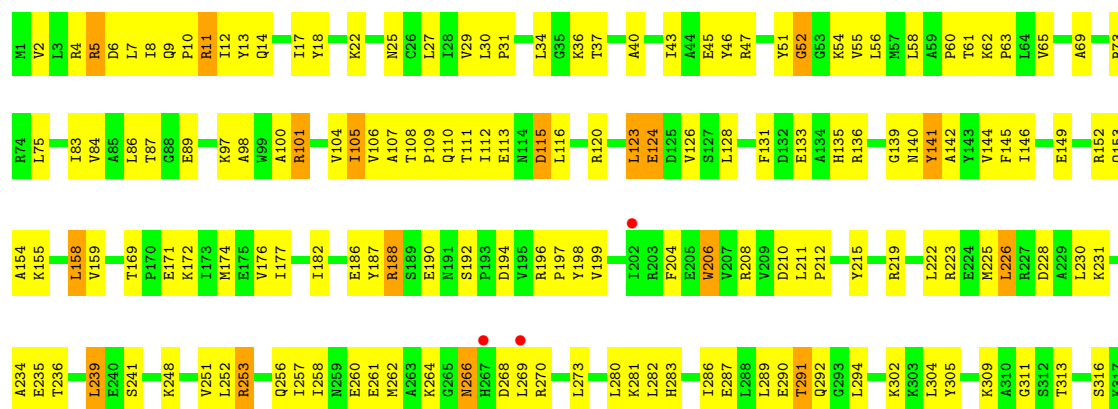


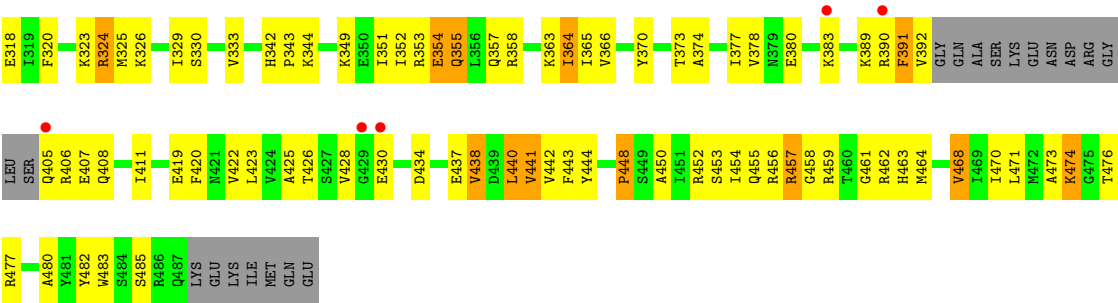


- Molecule 1: ATP-dependent RNA helicase, putative



- Molecule 1: ATP-dependent RNA helicase, putative





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	118.50Å 165.94Å 119.16Å 90.00° 114.85° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 29.85 – 2.78	Depositor EDS
% Data completeness (in resolution range)	97.1 (15.00-2.90) 95.4 (29.85-2.78)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.76Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.257 , 0.286 0.249 , 0.276	Depositor DCC
R_{free} test set	9383 reflections (9.05%)	wwPDB-VP
Wilson B-factor (Å ²)	64.9	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	22913	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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: PO4

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	0.44	0/3894	0.66	0/5241
1	B	0.44	0/3880	0.70	3/5223 (0.1%)
1	C	0.43	0/3870	0.66	0/5210
1	D	0.42	0/3827	0.64	0/5150
1	E	0.44	0/3870	0.70	3/5210 (0.1%)
1	F	0.43	0/3870	0.65	0/5210
All	All	0.43	0/23211	0.67	6/31244 (0.0%)

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	196	ARG	NE-CZ-NH1	-11.98	114.31	120.30
1	B	120	ARG	NE-CZ-NH2	11.49	126.04	120.30
1	B	120	ARG	NE-CZ-NH1	-11.06	114.77	120.30
1	E	196	ARG	NE-CZ-NH2	10.89	125.75	120.30
1	B	120	ARG	CD-NE-CZ	5.90	131.85	123.60

There are no chirality outliers.

There are no planarity outliers.

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	3831	0	4023	234	0
1	B	3817	0	4005	216	1
1	C	3807	0	3997	244	0
1	D	3765	0	3956	228	0
1	E	3807	0	3997	227	1
1	F	3807	0	3997	246	0
2	A	5	0	0	5	0
2	B	5	0	0	1	0
2	C	5	0	0	2	0
2	D	5	0	0	0	0
2	E	5	0	0	1	0
2	F	5	0	0	1	0
3	A	4	0	0	0	0
3	B	13	0	0	0	0
3	C	8	0	0	0	0
3	D	7	0	0	2	0
3	E	13	0	0	0	0
3	F	4	0	0	0	0
All	All	22913	0	23975	1321	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 28.

The worst 5 of 1321 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:F:211:LEU:H	1:F:474:LYS:HE2	1.19	1.06
1:E:405:GLN:HB2	1:E:408:GLN:HG3	1.38	1.00
1:D:123:LEU:HD23	1:D:123:LEU:H	1.25	1.00
1:E:426:THR:HG23	1:E:428:VAL:HG12	1.47	0.97
1:C:476:THR:HG23	1:C:478:ASP:H	1.27	0.97

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:B:268:ASP:OD2	1:E:371:ARG:NH1[1_455]	2.08	0.12

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	475/494 (96%)	442 (93%)	28 (6%)	5 (1%)	16	46
1	B	473/494 (96%)	443 (94%)	26 (6%)	4 (1%)	21	54
1	C	471/494 (95%)	435 (92%)	29 (6%)	7 (2%)	11	37
1	D	463/494 (94%)	433 (94%)	26 (6%)	4 (1%)	19	52
1	E	471/494 (95%)	441 (94%)	27 (6%)	3 (1%)	27	61
1	F	471/494 (95%)	439 (93%)	29 (6%)	3 (1%)	27	61
All	All	2824/2964 (95%)	2633 (93%)	165 (6%)	26 (1%)	19	52

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	52	GLY
1	C	52	GLY
1	F	52	GLY
1	A	52	GLY
1	A	484	SER

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	413/427 (97%)	371 (90%)	42 (10%)	8	24
1	B	412/427 (96%)	370 (90%)	42 (10%)	8	24
1	C	411/427 (96%)	370 (90%)	41 (10%)	8	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	406/427 (95%)	369 (91%)	37 (9%)	10	31
1	E	411/427 (96%)	369 (90%)	42 (10%)	8	24
1	F	411/427 (96%)	371 (90%)	40 (10%)	9	27
All	All	2464/2562 (96%)	2220 (90%)	244 (10%)	8	26

5 of 244 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	355	GLN
1	D	158	LEU
1	F	318	GLU
1	C	370	TYR
1	C	485	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 101 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	360	GLN
1	D	179	ASN
1	F	292	GLN
1	C	408	GLN
1	D	25	ASN

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](#)

6 ligands 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	$\# Z > 2$
2	PO4	A	495	-	4,4,4	1.83	3 (75%)	6,6,6	0.40	0
2	PO4	B	495	-	4,4,4	1.65	0	6,6,6	0.42	0
2	PO4	C	495	-	4,4,4	1.67	1 (25%)	6,6,6	0.46	0
2	PO4	D	495	-	4,4,4	1.60	0	6,6,6	0.41	0
2	PO4	E	495	-	4,4,4	1.64	0	6,6,6	0.42	0
2	PO4	F	495	-	4,4,4	1.60	0	6,6,6	0.42	0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	495	PO4	P-O4	-2.16	1.48	1.54
2	C	495	PO4	P-O3	-2.14	1.48	1.54
2	A	495	PO4	P-O3	-2.05	1.48	1.54
2	A	495	PO4	P-O2	-2.03	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	A	479/494 (96%)	-0.17	4 (0%) 86 85	24, 55, 97, 109	0
1	B	477/494 (96%)	-0.14	6 (1%) 77 76	24, 57, 95, 117	0
1	C	475/494 (96%)	-0.01	7 (1%) 73 72	30, 62, 99, 119	0
1	D	469/494 (94%)	0.04	20 (4%) 35 30	30, 67, 100, 116	0
1	E	475/494 (96%)	-0.06	13 (2%) 54 50	25, 58, 99, 118	0
1	F	475/494 (96%)	-0.06	8 (1%) 70 69	35, 65, 102, 126	0
All	All	2850/2964 (96%)	-0.07	58 (2%) 65 63	24, 61, 99, 126	0

The worst 5 of 58 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	269	LEU	8.2
1	B	269	LEU	5.7
1	D	269	LEU	5.1
1	E	266	ASN	4.6
1	D	308	ALA	4.6

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](#)

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
2	PO4	D	495	5/5	0.73	0.34	124,125,125,126	0
2	PO4	F	495	5/5	0.88	0.18	118,118,119,119	0
2	PO4	C	495	5/5	0.88	0.25	84,86,88,88	0
2	PO4	E	495	5/5	0.90	0.27	111,112,112,113	0
2	PO4	A	495	5/5	0.92	0.23	97,98,98,99	0
2	PO4	B	495	5/5	0.92	0.24	105,105,106,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.