



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2023 – 06:24 AM EDT

PDB ID : 2FL3
Title : Binary Complex of Restriction Endonuclease HinP1I with Cognate DNA
Authors : Horton, J.R.
Deposited on : 2006-01-05
Resolution : 2.39 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 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 : 4.02b-467
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

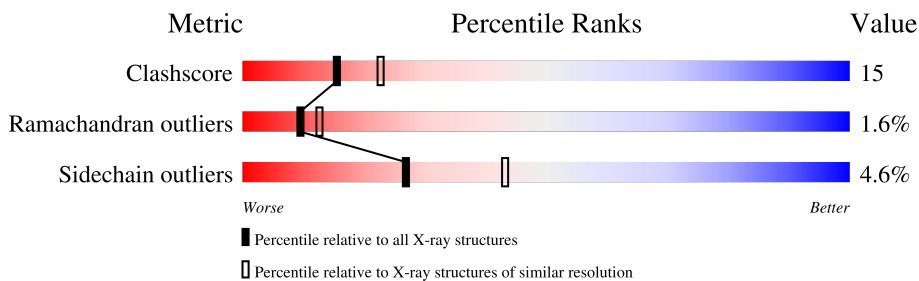
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	10	
1	D	10	
2	A	247	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2352 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 5'-D(*CP*CP*AP*GP*CP*GP*CP*TP*GP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	10	202	96	39	58	9	0	0	0
1	D	10	201	96	39	57	9	0	0	0

- Molecule 2 is a protein called R.HinP1I Restriction Endonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	247	1939	1224	351	359	5	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total 1 O 1	0	0
3	A	9	Total 9 O 9	0	0

4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.86Å 100.86Å 139.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.17 – 2.39	Depositor
% Data completeness (in resolution range)	(Not available) (34.17-2.39)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.246 , 0.284	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	2352	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

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	C	0.41	0/226	0.95	0/347
1	D	0.40	0/225	0.81	0/346
2	A	0.40	0/1979	0.64	0/2681
All	All	0.40	0/2430	0.70	0/3374

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	C	0	6
1	D	0	3
All	All	0	9

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1	DC	Sidechain
1	C	2	DC	Sidechain
1	C	3	DA	Sidechain
1	C	4	DG	Sidechain
1	C	5	DC	Sidechain
1	C	6	DG	Sidechain
1	D	14	DG	Sidechain
1	D	15	DC	Sidechain
1	D	16	DG	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	C	202	0	113	2	0
1	D	201	0	110	9	0
2	A	1939	0	1820	60	0
3	A	9	0	0	0	0
3	D	1	0	0	0	0
All	All	2352	0	2043	64	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 15.

All (64) 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:D:13:DA:H5 [?]	2:A:210:ARG:HD2	1.56	0.86
2:A:127:HIS:HD2	2:A:129:ASN:H	1.35	0.74
2:A:127:HIS:CD2	2:A:129:ASN:H	2.07	0.73
2:A:35:GLN:O	2:A:39:VAL:HG23	1.90	0.70
1:D:15:DC:H2 [?]	2:A:225:GLY:HA2	1.76	0.67
2:A:154:GLU:HA	2:A:205:ILE:HD12	1.75	0.66
1:D:17:DC:H5 ^{??}	2:A:4:VAL:HG13	1.79	0.64
1:D:15:DC:H2 [?]	2:A:225:GLY:CA	2.28	0.63
1:D:13:DA:H5 ^{??}	2:A:210:ARG:HB2	1.80	0.63
2:A:28:TRP:O	2:A:30:GLU:N	2.27	0.62
2:A:105:MET:CE	2:A:163:ASP:HB3	2.29	0.62
2:A:111:ASN:O	2:A:115:ILE:HG13	2.02	0.59
2:A:28:TRP:C	2:A:30:GLU:H	2.05	0.59
2:A:191:ARG:HD2	2:A:247:ILE:HD11	1.84	0.58
2:A:200:TYR:CZ	2:A:239:ILE:HD12	2.39	0.57
2:A:177:LEU:HB2	2:A:190:LEU:HD13	1.87	0.57
2:A:208:SER:HB2	2:A:209:PRO:HD2	1.87	0.57
2:A:117:ARG:HG2	2:A:122:GLU:CB	2.34	0.56
2:A:157:ARG:NH2	2:A:203:GLY:O	2.34	0.56
2:A:44:ASN:HD21	2:A:46:ASP:HB2	1.69	0.56
2:A:17:ASN:O	2:A:21:ILE:HD13	2.08	0.54
1:D:13:DA:C5 [?]	2:A:210:ARG:HD2	2.32	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:63:ILE:HG22	2:A:65:VAL:HG23	1.90	0.52
2:A:52:LYS:HE2	2:A:54:VAL:HG22	1.91	0.52
2:A:112:LEU:HD12	2:A:112:LEU:O	2.10	0.52
1:D:16:DG:H2''	1:D:17:DC:OP2	2.10	0.51
2:A:80:ILE:HG12	2:A:175:TRP:HB2	1.93	0.51
2:A:117:ARG:HG2	2:A:122:GLU:HB2	1.93	0.51
2:A:28:TRP:C	2:A:30:GLU:N	2.66	0.50
2:A:154:GLU:HA	2:A:205:ILE:CD1	2.42	0.50
2:A:157:ARG:NH1	2:A:205:ILE:HG13	2.27	0.50
2:A:189:ILE:HD11	2:A:191:ARG:CZ	2.43	0.49
2:A:135:ARG:HG2	2:A:222:ARG:CD	2.43	0.48
2:A:231:THR:O	2:A:234:MET:HE2	2.13	0.48
2:A:135:ARG:C	2:A:135:ARG:HD3	2.34	0.48
2:A:1:MET:O	2:A:2:ASN:CG	2.52	0.48
2:A:208:SER:HB2	2:A:209:PRO:CD	2.43	0.48
2:A:157:ARG:HG3	2:A:205:ILE:HD11	1.95	0.47
1:D:16:DG:H5''	2:A:3:LEU:HB3	1.97	0.47
2:A:135:ARG:HG2	2:A:222:ARG:HD3	1.97	0.46
2:A:170:ASP:O	2:A:171:PHE:HB2	2.16	0.46
2:A:38:LEU:O	2:A:43:HIS:HB2	2.15	0.46
2:A:94:ILE:HD12	2:A:165:LEU:HD13	1.97	0.46
2:A:127:HIS:CD2	2:A:129:ASN:HB2	2.50	0.45
1:C:1:DC:H2''	1:C:2:DC:O5''	2.15	0.45
2:A:23:ASP:O	2:A:25:PHE:N	2.50	0.44
2:A:120:THR:O	2:A:222:ARG:NE	2.51	0.44
2:A:183:SER:OG	2:A:184:ASN:N	2.50	0.44
2:A:217:ARG:HH21	2:A:243:GLU:CD	2.22	0.43
2:A:44:ASN:HD22	2:A:47:GLU:HB2	1.83	0.43
2:A:59:TYR:CD1	2:A:171:PHE:HB2	2.54	0.42
2:A:127:HIS:HD2	2:A:129:ASN:N	2.09	0.42
1:C:5:DC:H2''	1:C:6:DG:C5''	2.49	0.42
2:A:213:ILE:HB	2:A:220:ILE:HB	2.00	0.42
2:A:32:SER:O	2:A:36:ASP:OD2	2.38	0.41
2:A:91:PHE:CE1	2:A:238:LYS:HE2	2.55	0.41
2:A:96:LYS:HD2	2:A:234:MET:HE3	2.02	0.41
2:A:214:ASN:HD22	2:A:214:ASN:HA	1.67	0.41
2:A:189:ILE:HB	2:A:245:PHE:CD1	2.56	0.41
2:A:241:PRO:O	2:A:244:LEU:HB2	2.20	0.41
2:A:96:LYS:CD	2:A:234:MET:HE3	2.52	0.40
1:D:11:DC:H2''	1:D:12:DC:O5''	2.22	0.40
2:A:99:LEU:HB3	2:A:113:LEU:CD1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:125:PRO:HB3	2:A:130:THR:HG21	2.02	0.40

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
2	A	245/247 (99%)	220 (90%)	21 (9%)	4 (2%)	9 13

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	73	ASP
2	A	24	ARG
2	A	28	TRP
2	A	70	PHE

5.3.2 Protein sidechains [i](#)

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
2	A	195/221 (88%)	186 (95%)	9 (5%)	27 43

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	95	ASP
2	A	99	LEU
2	A	109	ASP
2	A	135	ARG
2	A	159	LEU
2	A	165	LEU
2	A	227	ASN
2	A	238	LYS
2	A	244	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	31	ASN
2	A	43	HIS
2	A	44	ASN
2	A	64	ASN
2	A	79	ASN
2	A	87	ASN
2	A	127	HIS
2	A	214	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 monosaccharides 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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.