



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 07:12 PM UTC

PDB ID : 9BJ5 / pdb\_00009bj5  
Title : Crystal structure of the HEPN family member AbiV (P212121 space group)  
Authors : Zhu, X.; Moineau, S.; Shi, R.  
Deposited on : 2024-04-25  
Resolution : 2.55 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](mailto: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>.

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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

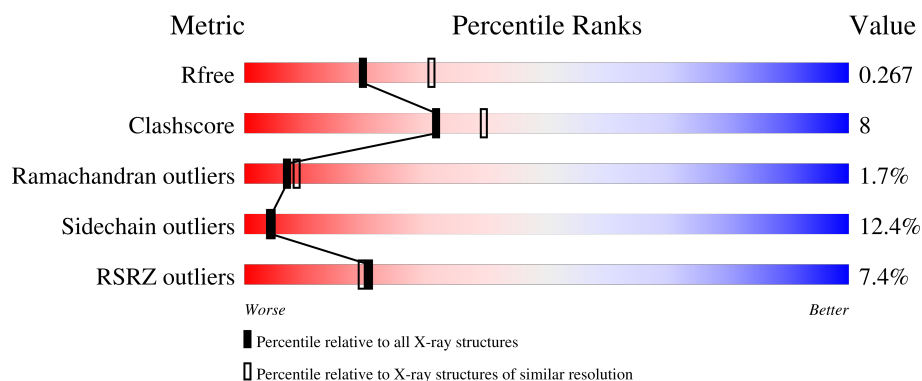
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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  $\geq 3$ , 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	A	207	<div> <div>3%</div> <div>51%</div> <div>21%</div> <div>•</div> <div>25%</div> </div>
1	D	207	<div> <div>9%</div> <div>48%</div> <div>21%</div> <div>•</div> <div>29%</div> </div>
2	B	207	<div> <div>7%</div> <div>53%</div> <div>14%</div> <div>5%</div> <div>28%</div> </div>
3	C	207	<div> <div>2%</div> <div>60%</div> <div>12%</div> <div>•</div> <div>25%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4823 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 AbiV family abortive infection protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1219	778	186	251	4			
1	D	147	Total	C	N	O	S	0	0	0
			1156	735	177	239	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	202	HIS	-	expression tag	UNP Q9AGY3
A	203	HIS	-	expression tag	UNP Q9AGY3
A	204	HIS	-	expression tag	UNP Q9AGY3
A	205	HIS	-	expression tag	UNP Q9AGY3
A	206	HIS	-	expression tag	UNP Q9AGY3
A	207	HIS	-	expression tag	UNP Q9AGY3
D	202	HIS	-	expression tag	UNP Q9AGY3
D	203	HIS	-	expression tag	UNP Q9AGY3
D	204	HIS	-	expression tag	UNP Q9AGY3
D	205	HIS	-	expression tag	UNP Q9AGY3
D	206	HIS	-	expression tag	UNP Q9AGY3
D	207	HIS	-	expression tag	UNP Q9AGY3

- Molecule 2 is a protein called AbiV family abortive infection protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	150	Total	C	N	O	S	0	0	0
			1180	754	181	241	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	202	HIS	-	expression tag	UNP Q9AGY3
B	203	HIS	-	expression tag	UNP Q9AGY3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	204	HIS	-	expression tag	UNP Q9AGY3
B	205	HIS	-	expression tag	UNP Q9AGY3
B	206	HIS	-	expression tag	UNP Q9AGY3
B	207	HIS	-	expression tag	UNP Q9AGY3

- Molecule 3 is a protein called AbiV family abortive infection protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	156	Total	C	N	O	S	0	0	0
			1227	784	187	251	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	202	HIS	-	expression tag	UNP Q9AGY3
C	203	HIS	-	expression tag	UNP Q9AGY3
C	204	HIS	-	expression tag	UNP Q9AGY3
C	205	HIS	-	expression tag	UNP Q9AGY3
C	206	HIS	-	expression tag	UNP Q9AGY3
C	207	HIS	-	expression tag	UNP Q9AGY3

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

- Molecule 5 is BROMIDE ION (CCD ID: BR) (formula: Br) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Br	0	0
			1	1		
5	B	1	Total	Br	0	0
			1	1		
5	D	1	Total	Br	0	0
			1	1		

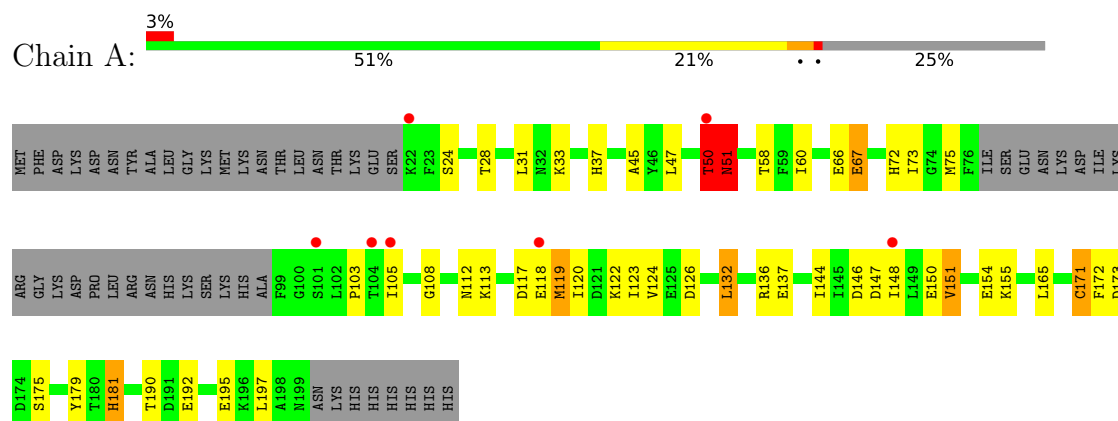
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	12	Total 12	O 12	0	0
6	B	4	Total 4	O 4	0	0
6	C	15	Total 15	O 15	0	0
6	D	6	Total 6	O 6	0	0

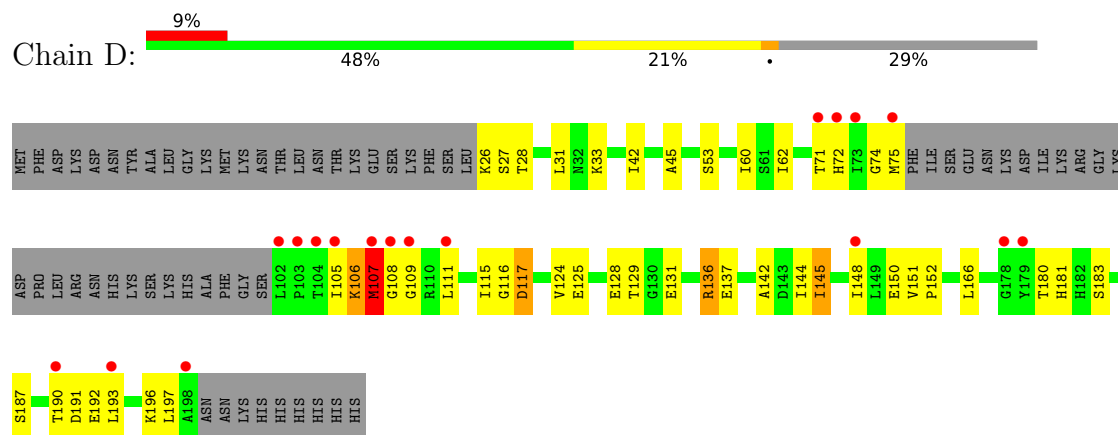
### 3 Residue-property plots [i](#)

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.

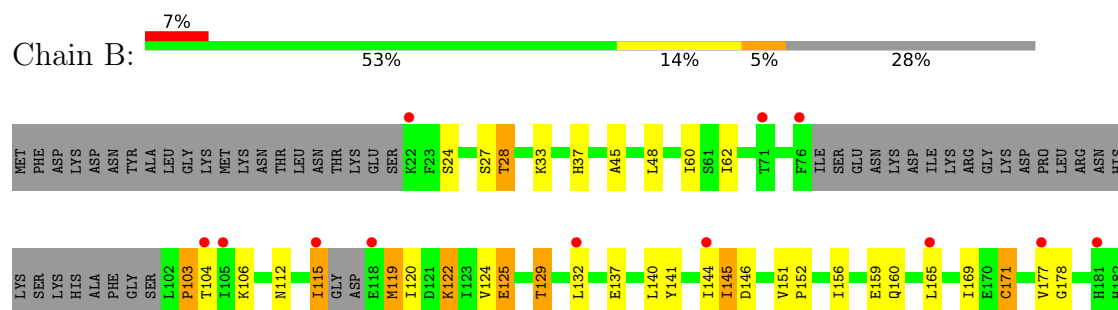
- Molecule 1: AbiV family abortive infection protein

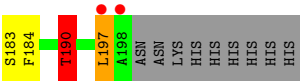


- Molecule 1: AbiV family abortive infection protein

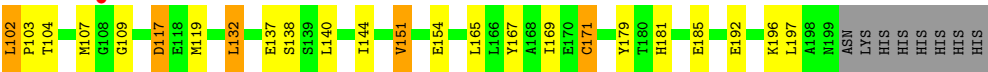


- Molecule 2: AbiV family abortive infection protein





● Molecule 3: AbiV family abortive infection protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.69Å 79.17Å 136.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.31 – 2.55 42.31 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.31-2.55) 99.7 (42.31-2.55)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.40 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.219 , 0.267 0.222 , 0.267	Depositor DCC
$R_{free}$ test set	1307 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.0	Xtriage
Anisotropy	0.702	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4823	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.75 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.9341e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NA, MLY, BR

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	1.22	1/1192 (0.1%)	1.64	14/1616 (0.9%)
1	D	1.23	0/1126	1.71	1/1525 (0.1%)
2	B	1.20	1/1151 (0.1%)	1.70	2/1557 (0.1%)
3	C	1.27	1/1211 (0.1%)	1.59	6/1638 (0.4%)
All	All	1.23	3/4680 (0.1%)	1.66	23/6336 (0.4%)

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	D	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	138	SER	C-O	-5.91	1.16	1.24
2	B	37	HIS	CE1-NE2	5.86	1.38	1.32
1	A	181	HIS	CE1-NE2	5.44	1.38	1.32

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	154	GLU	CB-CA-C	7.69	124.71	110.10
1	A	50	THR	CA-CB-OG1	-7.58	98.22	109.60
3	C	171	CYS	CB-CA-C	-6.97	99.66	110.81
1	A	51	ASN	CB-CA-C	6.54	123.44	110.42
2	B	190	THR	CA-CB-OG1	-6.45	99.93	109.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	116	GLY	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	1219	0	1187	21	0
1	D	1156	0	1146	23	0
2	B	1180	0	1160	22	0
3	C	1227	0	1211	15	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	D	1	0	0	1	0
6	A	12	0	0	1	0
6	B	4	0	0	0	0
6	C	15	0	0	2	0
6	D	6	0	0	2	0
All	All	4823	0	4704	77	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 8.

The worst 5 of 77 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:124:VAL:O	1:D:128:GLU:HG2	1.86	0.76
1:D:145:ILE:HG13	1:D:150:GLU:OE1	1.86	0.75
3:C:192:GLU:OE2	3:C:196:MLY:HH13	1.87	0.75
1:A:47:LEU:HD23	2:B:48:LEU:HD23	1.69	0.74
1:D:192:GLU:O	1:D:196:MLY:HG2	1.89	0.73

There are no symmetry-related clashes.

## 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	148/207 (72%)	137 (93%)	8 (5%)	3 (2%)	6	6
1	D	139/207 (67%)	129 (93%)	7 (5%)	3 (2%)	5	5
2	B	140/207 (68%)	132 (94%)	6 (4%)	2 (1%)	9	11
3	C	149/207 (72%)	143 (96%)	4 (3%)	2 (1%)	9	12
All	All	576/828 (70%)	541 (94%)	25 (4%)	10 (2%)	7	8

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	119	MET
1	D	108	GLY
1	D	117	ASP
3	C	101	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	132/183 (72%)	115 (87%)	17 (13%)	4	4
1	D	127/183 (69%)	110 (87%)	17 (13%)	4	3
2	B	128/183 (70%)	111 (87%)	17 (13%)	4	4
3	C	136/184 (74%)	122 (90%)	14 (10%)	7	8
All	All	523/733 (71%)	458 (88%)	65 (12%)	4	4

5 of 65 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	129	THR
1	D	136	ARG
2	B	122	LYS
2	B	119	MET
1	D	137	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
1	A	112	ASN
3	C	72	HIS
3	C	182	HIS
1	D	72	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

15 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	$\# Z  > 2$
3	MLY	C	196	3	9,10,11	0.51	0	6,11,13	0.46	0
1	MLY	D	196	1	9,10,11	0.41	0	6,11,13	0.23	0
2	MLY	B	196	2	9,10,11	0.43	0	6,11,13	0.15	0
1	MLY	A	155	1	9,10,11	0.54	0	6,11,13	0.68	0
2	MLY	B	33	2	9,10,11	0.51	0	6,11,13	0.47	0
1	MLY	A	43	1	9,10,11	0.48	0	6,11,13	0.49	0
1	MLY	A	158	1	9,10,11	0.55	0	6,11,13	0.40	0
1	MLY	D	158	1	9,10,11	0.49	0	6,11,13	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MLY	B	155	2	9,10,11	0.47	0	6,11,13	0.22	0
3	MLY	C	155	3	9,10,11	0.42	0	6,11,13	0.43	0
1	MLY	A	196	1	9,10,11	0.51	0	6,11,13	0.62	0
3	MLY	C	158	3	9,10,11	0.57	0	6,11,13	0.34	0
2	MLY	B	158	2	9,10,11	0.48	0	6,11,13	0.61	0
1	MLY	D	155	1	9,10,11	0.51	0	6,11,13	0.34	0
1	MLY	D	43	1	9,10,11	0.58	0	6,11,13	0.22	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
3	MLY	C	196	3	-	3/8/9/11	-
1	MLY	D	196	1	-	4/8/9/11	-
2	MLY	B	196	2	-	3/8/9/11	-
1	MLY	A	155	1	-	1/8/9/11	-
2	MLY	B	33	2	-	3/8/9/11	-
1	MLY	A	43	1	-	4/8/9/11	-
1	MLY	A	158	1	-	4/8/9/11	-
1	MLY	D	158	1	-	1/8/9/11	-
2	MLY	B	155	2	-	0/8/9/11	-
3	MLY	C	155	3	-	3/8/9/11	-
1	MLY	A	196	1	-	5/8/9/11	-
3	MLY	C	158	3	-	3/8/9/11	-
2	MLY	B	158	2	-	1/8/9/11	-
1	MLY	D	155	1	-	3/8/9/11	-
1	MLY	D	43	1	-	4/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 42 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	43	MLY	C-CA-CB-CG
2	B	33	MLY	C-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
2	B	196	MLY	O-C-CA-CB
1	D	43	MLY	CD-CE-NZ-CH1
1	A	196	MLY	CD-CE-NZ-CH1

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	196	MLY	2	0
1	D	196	MLY	2	0
1	A	155	MLY	1	0
2	B	33	MLY	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	152/207 (73%)	0.05	7 (4%) 37 38	34, 50, 84, 122	0
1	D	143/207 (69%)	0.69	18 (12%) 8 7	46, 75, 117, 167	0
2	B	146/207 (70%)	0.65	14 (9%) 13 13	49, 76, 108, 136	0
3	C	153/207 (73%)	0.13	5 (3%) 49 50	35, 50, 95, 123	0
All	All	594/828 (71%)	0.37	44 (7%) 20 19	34, 63, 108, 167	0

The worst 5 of 44 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	103	PRO	4.8
1	D	102	LEU	4.8
1	D	108	GLY	4.5
1	D	198	ALA	4.3
2	B	198	ALA	4.1

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	MLY	B	196	11/12	0.77	0.20	95,100,117,117	0
2	MLY	B	155	11/12	0.85	0.19	75,80,111,113	0
1	MLY	A	196	11/12	0.85	0.15	53,59,87,87	0
2	MLY	B	158	11/12	0.86	0.15	68,79,99,105	0
1	MLY	D	155	11/12	0.87	0.19	70,79,119,121	0
1	MLY	D	196	11/12	0.88	0.16	102,108,119,120	0
1	MLY	D	158	11/12	0.90	0.15	69,76,88,89	0

*Continued on next page...*

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	MLY	D	43	11/12	0.90	0.17	52,60,77,79	0
2	MLY	B	33	11/12	0.92	0.14	55,59,79,86	0
3	MLY	C	158	11/12	0.93	0.14	44,52,70,72	0
3	MLY	C	196	11/12	0.93	0.10	58,66,78,80	0
1	MLY	A	158	11/12	0.93	0.13	42,51,77,78	0
1	MLY	A	43	11/12	0.95	0.11	45,55,86,89	0
1	MLY	A	155	11/12	0.95	0.09	43,50,57,62	0
3	MLY	C	155	11/12	0.95	0.09	46,56,63,65	0

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides 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, 95<sup>th</sup> 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.

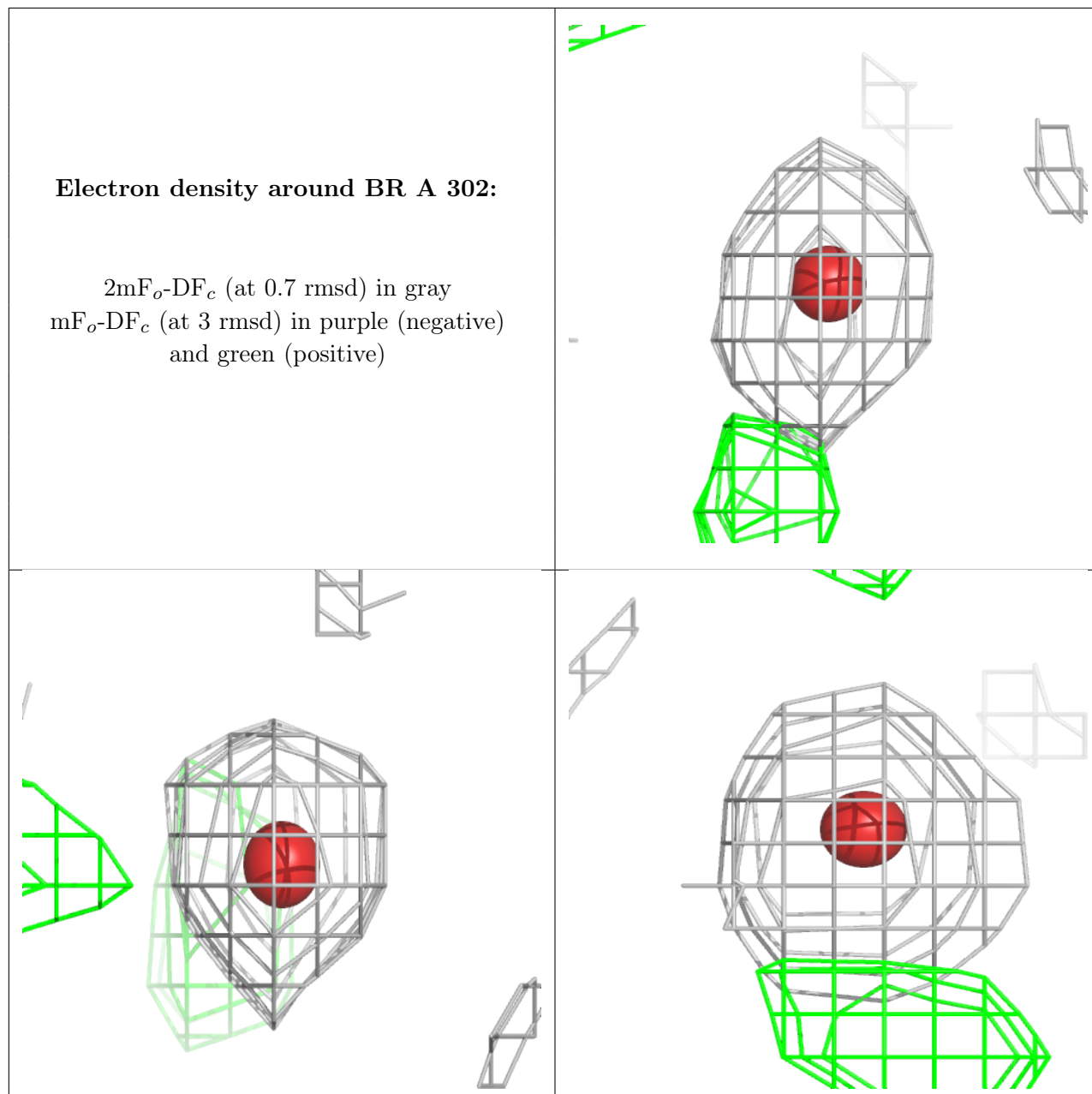
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	NA	A	301	1/1	0.92	0.16	43,43,43,43	0
5	BR	A	302	1/1	0.93	0.15	43,43,43,43	1
5	BR	D	301	1/1	0.94	0.06	52,52,52,52	1
5	BR	B	301	1/1	0.98	0.08	50,50,50,50	1

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



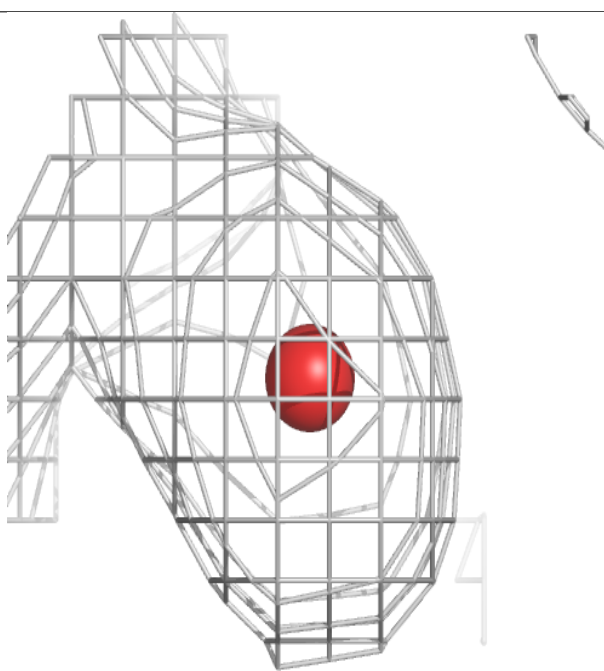
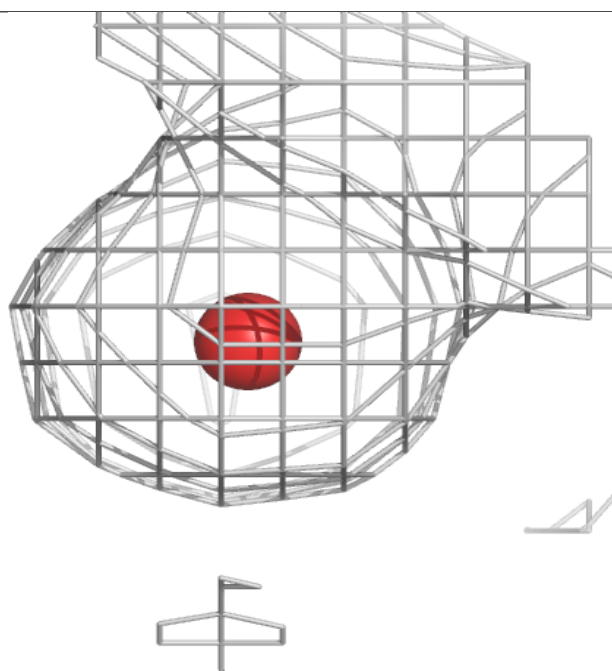
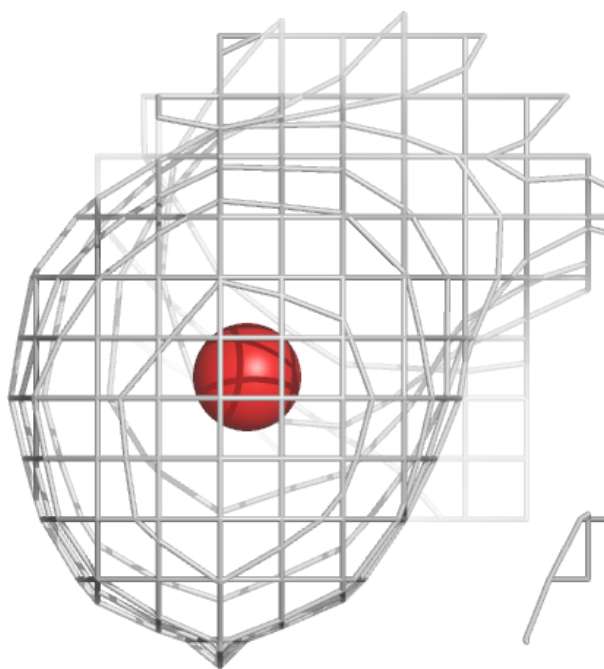
**Electron density around BR A 302:**

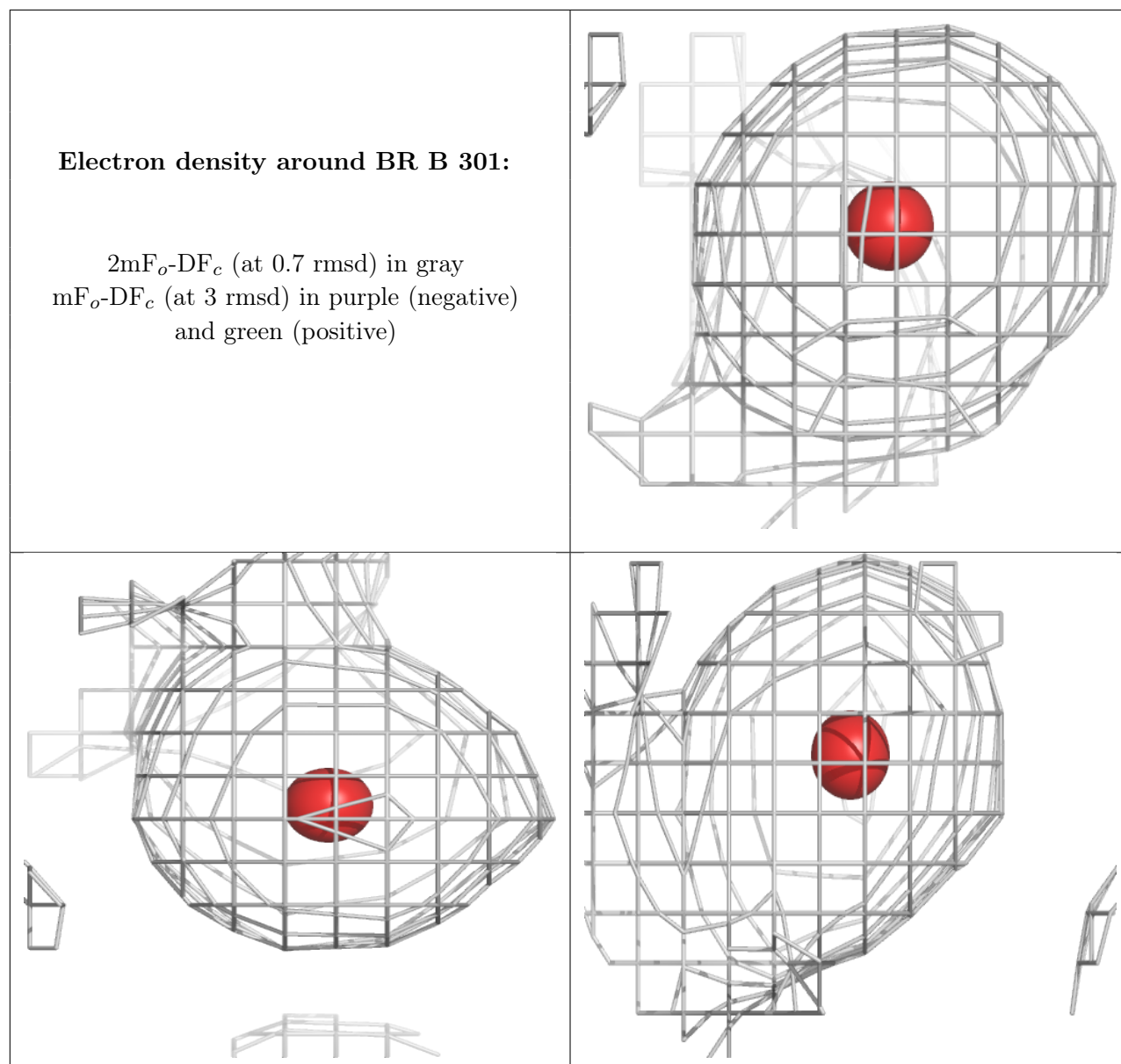
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around BR D 301:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers ⓘ

There are no such residues in this entry.