



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:08 PM UTC

PDB ID : 2FA1 / pdb\_00002fa1  
Title : Crystal structure of PhnF C-terminal domain  
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Deposited on : 2005-12-06  
Resolution : 1.70 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](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

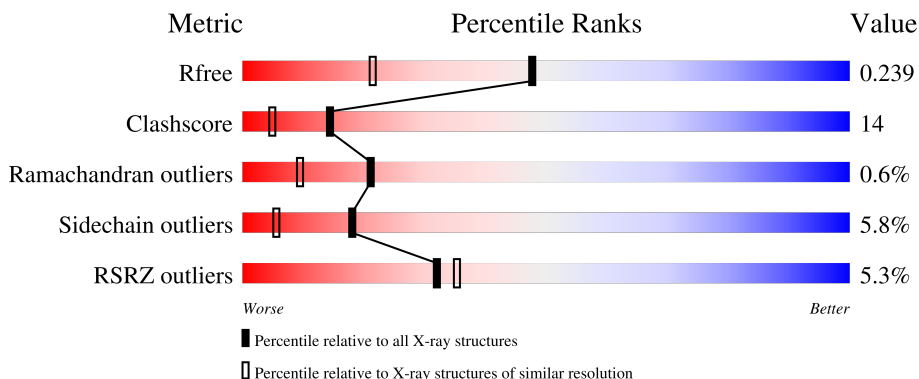
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	160	
1	B	160	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2960 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.

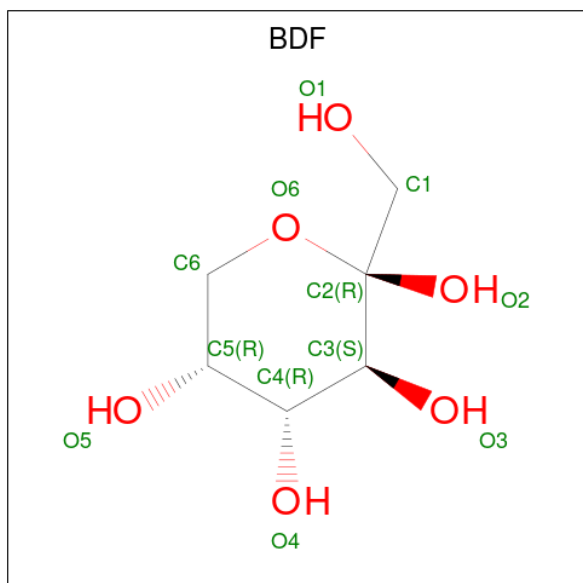
- Molecule 1 is a protein called Probable transcriptional regulator phnF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	159	1333	818	254	253	8	0	11	0
1	B	159	1288	791	249	241	7	0	4	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	82	GLY	-	cloning artifact	UNP P16684
A	83	HIS	-	cloning artifact	UNP P16684
A	84	MET	-	cloning artifact	UNP P16684
B	82	GLY	-	cloning artifact	UNP P16684
B	83	HIS	-	cloning artifact	UNP P16684
B	84	MET	-	cloning artifact	UNP P16684

- Molecule 2 is beta-D-fructopyranose (CCD ID: BDF) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			12	6	6		


- Molecule 3 is water.

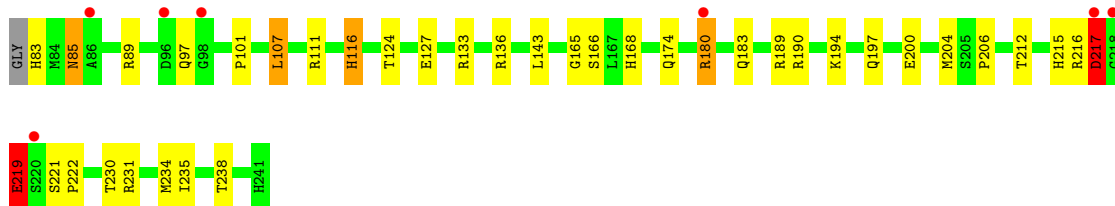
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	163	Total	O	0	0
			163	163		
3	B	164	Total	O	0	0
			164	164		

### 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.

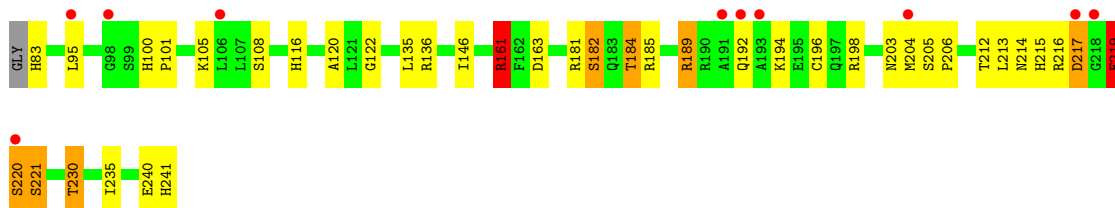
- Molecule 1: Probable transcriptional regulator phnF

Chain A: 



- Molecule 1: Probable transcriptional regulator phnF

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.77Å 51.57Å 63.96Å 90.00° 103.64° 90.00°	Depositor
Resolution (Å)	50.00 – 1.70 50.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-1.70) 99.9 (50.00-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 1.60Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.187 , 0.241 0.187 , 0.239	Depositor DCC
$R_{free}$ test set	1938 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.4	Xtrriage
Anisotropy	0.195	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 47.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2960	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.38 % 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 4.7117e-03. 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	1.42	4/1368 (0.3%)	1.35	5/1846 (0.3%)
1	B	1.40	4/1315 (0.3%)	1.34	7/1776 (0.4%)
All	All	1.41	8/2683 (0.3%)	1.35	12/3622 (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	B	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	146	ILE	CA-CB	-5.73	1.47	1.54
1	B	184	THR	N-CA	5.43	1.52	1.46
1	A	200	GLU	C-O	5.36	1.30	1.23
1	B	135	LEU	N-CA	5.19	1.52	1.46
1	B	230	THR	N-CA	5.09	1.52	1.46
1	A	212	THR	C-O	-5.07	1.17	1.23
1	A	136	ARG	C-O	5.05	1.30	1.24
1	A	166	SER	N-CA	-5.03	1.40	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	SER	CA-C-N	-10.33	109.81	120.03
1	A	221	SER	C-N-CA	-10.33	109.81	120.03
1	B	219	GLU	N-CA-C	8.82	122.78	108.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	161	ARG	NE-CZ-NH1	8.27	129.77	121.50
1	B	161	ARG	CD-NE-CZ	6.93	134.11	124.40
1	B	221	SER	N-CA-C	-6.39	94.92	108.73
1	A	111	ARG	NE-CZ-NH2	6.22	124.80	119.20
1	A	111	ARG	NE-CZ-NH1	-6.11	115.39	121.50
1	B	205[A]	SER	N-CA-C	-5.13	103.31	109.83
1	B	205[B]	SER	N-CA-C	-5.13	103.31	109.83
1	A	219	GLU	N-CA-C	5.06	117.38	110.55
1	B	161	ARG	NE-CZ-NH2	-5.04	114.67	119.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	217	ASP	Peptide
1	B	219	GLU	Peptide

## 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	A	1333	0	1305	39	4
1	B	1288	0	1278	36	0
2	B	12	0	12	0	0
3	A	163	0	0	7	2
3	B	164	0	0	18	2
All	All	2960	0	2595	74	4

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 14.

All (74) 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:B:189:ARG:HG3	1:B:189:ARG:HH11	1.15	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:HIS:CD2	3:A:385:HOH:O	2.19	0.95
1:A:206[B]:PRO:HG3	3:B:326:HOH:O	1.81	0.80
1:B:189:ARG:HG3	1:B:189:ARG:NH1	1.91	0.78
1:A:216:ARG:HB2	1:A:219:GLU:OE1	1.85	0.77
1:A:180[A]:ARG:HH11	1:A:180[A]:ARG:CG	1.98	0.77
1:A:85[A]:ASN:HD22	1:A:85[A]:ASN:H	1.36	0.73
1:B:219:GLU:HG2	1:B:220:SER:CB	2.19	0.72
1:A:174:GLN:OE1	3:A:339:HOH:O	2.09	0.70
1:A:206[B]:PRO:CG	3:B:326:HOH:O	2.37	0.69
1:B:216:ARG:HD3	3:B:373:HOH:O	1.93	0.69
1:B:120:ALA:HB1	1:B:213:LEU:HD21	1.74	0.68
1:A:107[B]:LEU:HD11	1:A:133:ARG:NH1	2.09	0.68
1:B:198:ARG:NH1	3:B:399:HOH:O	2.22	0.67
1:B:219:GLU:HG2	1:B:220:SER:OG	1.94	0.67
1:B:203:ASN:ND2	1:B:204:MET:HG2	2.11	0.65
1:B:83:HIS:N	3:B:380:HOH:O	2.31	0.64
1:B:182:SER:OG	1:B:215:HIS:CE1	2.51	0.64
1:A:124:THR:HB	1:A:127[B]:GLU:OE1	1.99	0.63
1:A:116:HIS:HD2	3:A:385:HOH:O	1.71	0.63
1:A:85[A]:ASN:HD22	1:A:85[A]:ASN:N	2.00	0.59
1:A:124:THR:CB	1:A:127[B]:GLU:OE1	2.50	0.59
1:A:180[A]:ARG:HG2	1:A:180[A]:ARG:NH1	2.17	0.59
1:A:143:LEU:CD2	1:A:234[B]:MET:SD	2.91	0.58
1:A:89:ARG:HB2	3:A:301:HOH:O	2.04	0.58
1:B:83:HIS:HB3	3:B:350:HOH:O	2.03	0.57
1:B:182:SER:HB3	1:B:215:HIS:CE1	2.40	0.57
1:A:143:LEU:HD22	1:A:234[B]:MET:SD	2.46	0.56
1:A:97:GLN:HB3	1:A:168:HIS:ND1	2.21	0.56
1:A:180[A]:ARG:CG	1:A:180[A]:ARG:NH1	2.61	0.55
1:B:182:SER:HB3	1:B:215:HIS:ND1	2.21	0.55
1:B:241:HIS:OXT	3:B:388:HOH:O	2.18	0.55
1:B:206:PRO:CG	3:B:318:HOH:O	2.55	0.54
1:A:180[A]:ARG:HH11	1:A:180[A]:ARG:HG2	1.70	0.53
1:B:182:SER:CB	1:B:215:HIS:CE1	2.91	0.53
1:A:190:ARG:HD2	1:A:204:MET:HE3	1.91	0.52
1:A:216:ARG:O	1:A:217[A]:ASP:C	2.53	0.51
1:A:107[B]:LEU:CD1	1:A:133:ARG:NH1	2.74	0.51
1:A:180[A]:ARG:HH11	1:A:180[A]:ARG:HG3	1.71	0.49
1:B:240:GLU:CD	3:B:363:HOH:O	2.55	0.49
1:A:215:HIS:CD2	1:A:222:PRO:HA	2.46	0.49
1:A:83:HIS:HA	3:B:379:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180[B]:ARG:HD2	1:A:215:HIS:HB2	1.96	0.48
1:B:181:ARG:HA	1:B:214:ASN:HD22	1.78	0.47
1:B:161:ARG:NH2	3:B:286:HOH:O	2.45	0.47
1:B:206:PRO:HG2	3:B:318:HOH:O	2.12	0.47
1:A:238:THR:HG22	1:B:185:ARG:HG2	1.97	0.47
1:A:183:GLN:HG2	3:B:363:HOH:O	2.15	0.47
1:B:219:GLU:CG	1:B:220:SER:OG	2.63	0.46
1:B:163:ASP:HB2	3:B:330:HOH:O	2.14	0.46
1:B:230:THR:HG23	1:B:235:ILE:HD12	1.97	0.46
1:A:234[A]:MET:HE3	1:A:234[A]:MET:HB3	1.69	0.46
1:B:196:CYS:SG	1:B:203:ASN:HA	2.56	0.45
1:B:182:SER:OG	1:B:215:HIS:HE1	1.95	0.45
1:A:230[B]:THR:HG23	1:A:235:ILE:HD12	1.98	0.45
1:A:231:ARG:CB	1:A:234[B]:MET:HG3	2.47	0.44
1:A:234[B]:MET:HE2	1:A:234[B]:MET:HB3	1.78	0.44
1:B:136[B]:ARG:CG	3:B:332:HOH:O	2.66	0.44
1:B:181:ARG:HD3	1:B:184:THR:OG1	2.18	0.43
1:B:100:HIS:ND1	1:B:101:PRO:HD2	2.32	0.43
1:A:85[A]:ASN:H	1:A:85[A]:ASN:ND2	2.12	0.43
1:A:116:HIS:NE2	3:A:385:HOH:O	2.36	0.43
1:B:122:GLY:HA3	3:B:396:HOH:O	2.19	0.43
1:A:197[A]:GLN:CD	3:A:319:HOH:O	2.62	0.43
1:A:189:ARG:NH2	3:A:362:HOH:O	2.52	0.42
1:B:189:ARG:NH1	1:B:189:ARG:CG	2.70	0.41
1:A:206[B]:PRO:HG2	3:B:326:HOH:O	2.13	0.41
1:A:216:ARG:O	1:A:217[B]:ASP:C	2.58	0.41
1:B:105:LYS:HD2	1:B:108:SER:HB3	2.03	0.41
1:B:215:HIS:CD2	1:B:220:SER:HA	2.56	0.41
1:B:181:ARG:HG3	1:B:212[B]:THR:CG2	2.50	0.41
1:A:101:PRO:O	1:A:165:GLY:HA2	2.20	0.41
1:B:136[B]:ARG:NE	3:B:332:HOH:O	2.52	0.40
1:B:194:LYS:HA	1:B:194:LYS:HD2	1.92	0.40

All (4) 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:A:219:GLU:CD	3:B:367:HOH:O[2_646]	1.97	0.23
1:A:127[B]:GLU:OE1	3:A:320:HOH:O[2_646]	1.98	0.22
1:A:219:GLU:OE1	3:B:367:HOH:O[2_646]	2.01	0.19
1:A:127[B]:GLU:OE2	3:A:320:HOH:O[2_646]	2.13	0.07

## 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	
1	A	168/160 (105%)	164 (98%)	2 (1%)	2 (1%)	10	2
1	B	161/160 (101%)	154 (96%)	6 (4%)	1 (1%)	21	9
All	All	329/320 (103%)	318 (97%)	8 (2%)	3 (1%)	21	4

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217[A]	ASP
1	A	217[B]	ASP
1	B	217	ASP

### 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	
1	A	148/138 (107%)	137 (93%)	11 (7%)	13	3
1	B	142/138 (103%)	133 (94%)	9 (6%)	16	5
All	All	290/276 (105%)	270 (93%)	20 (7%)	18	4

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85[A]	ASN
1	A	85[B]	ASN
1	A	107[A]	LEU

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Mol	Chain	Res	Type
1	A	107[B]	LEU
1	A	116	HIS
1	A	180[A]	ARG
1	A	180[B]	ARG
1	A	194	LYS
1	A	217[A]	ASP
1	A	217[B]	ASP
1	A	219	GLU
1	B	95	LEU
1	B	116	HIS
1	B	161	ARG
1	B	182	SER
1	B	189	ARG
1	B	192	GLN
1	B	219	GLU
1	B	220	SER
1	B	221	SER

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

Mol	Chain	Res	Type
1	A	87	GLN
1	A	160	GLN
1	A	183	GLN
1	A	215	HIS
1	B	174	GLN
1	B	203	ASN
1	B	214	ASN
1	B	215	HIS

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	BDF	B	1	-	12,12,12	1.38	2 (16%)	18,18,18	2.92	9 (50%)

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
2	BDF	B	1	-	-	0/3/23/23	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	BDF	O5-C5	2.44	1.48	1.43
2	B	1	BDF	C2-C3	2.43	1.56	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	BDF	O6-C6-C5	-6.01	101.48	111.06
2	B	1	BDF	C6-C5-C4	5.54	117.71	109.64
2	B	1	BDF	O6-C2-C1	4.58	111.76	105.41
2	B	1	BDF	O4-C4-C3	-3.97	101.49	109.77
2	B	1	BDF	O2-C2-C1	-3.47	105.30	111.35
2	B	1	BDF	O2-C2-C3	3.18	113.61	107.93
2	B	1	BDF	C5-C4-C3	-2.50	105.91	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	BDF	O1-C1-C2	-2.31	106.55	111.67
2	B	1	BDF	C1-C2-C3	-2.23	106.88	111.63

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	159/160 (99%)	0.16	7 (4%) 39 42	6, 19, 37, 43	10 (6%)
1	B	159/160 (99%)	0.45	10 (6%) 26 28	8, 24, 42, 60	4 (2%)
All	All	318/320 (99%)	0.30	17 (5%) 32 35	6, 22, 41, 60	14 (4%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	191	ALA	3.1
1	A	98	GLY	2.9
1	B	98	GLY	2.8
1	B	204	MET	2.7
1	B	217	ASP	2.7
1	A	180[A]	ARG	2.6
1	B	95	LEU	2.6
1	B	192	GLN	2.6
1	A	86	ALA	2.6
1	A	218	GLY	2.4
1	B	193	ALA	2.4
1	B	220	SER	2.3
1	B	218	GLY	2.2
1	A	220	SER	2.2
1	A	217[A]	ASP	2.2
1	B	106	LEU	2.2
1	A	96	ASP	2.1

### 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 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(Å <sup>2</sup> )	Q<0.9
2	BDF	B	1	12/12	0.95	0.06	16,17,20,20	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.