



# Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 11:24 pm BST

PDB ID : 3LYQ  
Title : Crystal structure of IpgB2 from Shigella flexneri  
Authors : Klink, B.U.; Barden, S.; Heidler, T.V.; Borchers, C.; Ladwein, M.; Stradal, T.E.B.; Rottner, K.; Heinz, D.W.  
Deposited on : 2010-02-28  
Resolution : 2.30 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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

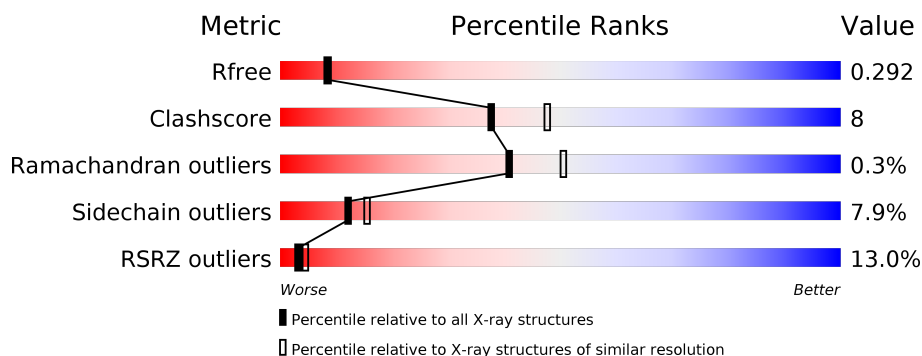
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	192	<div> <div>13%</div> <div> <div></div> <div>78%</div> <div>15%</div> <div>• 5%</div> </div> </div>
1	B	192	<div> <div>13%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>• •</div> </div> </div>

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	FLC	A	301	-	-	X	-
2	FLC	B	302	-	-	X	-
3	FEO	A	401	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3117 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.

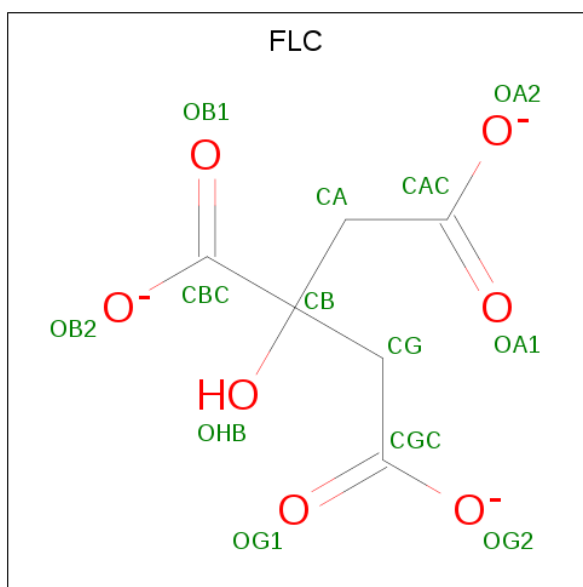
- Molecule 1 is a protein called IpgB2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	1	0
			1477	919	261	289	8			
1	B	186	Total	C	N	O	S	0	0	0
			1489	928	263	290	8			

There are 8 discrepancies between the modelled and reference sequences:

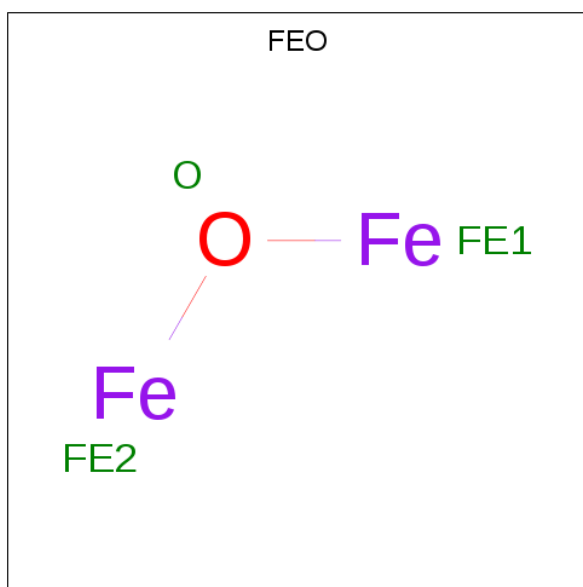
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP Q9AJW7
A	-2	ALA	-	EXPRESSION TAG	UNP Q9AJW7
A	-1	MET	-	EXPRESSION TAG	UNP Q9AJW7
A	0	ASP	-	EXPRESSION TAG	UNP Q9AJW7
B	-3	GLY	-	EXPRESSION TAG	UNP Q9AJW7
B	-2	ALA	-	EXPRESSION TAG	UNP Q9AJW7
B	-1	MET	-	EXPRESSION TAG	UNP Q9AJW7
B	0	ASP	-	EXPRESSION TAG	UNP Q9AJW7

- Molecule 2 is CITRATE ANION (three-letter code: FLC) (formula: C<sub>6</sub>H<sub>5</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is MU-OXO-DIIRON (three-letter code: FEO) (formula:  $\text{Fe}_2\text{O}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	O	0	0
			3	2	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	O	0	0
			3	2	1		

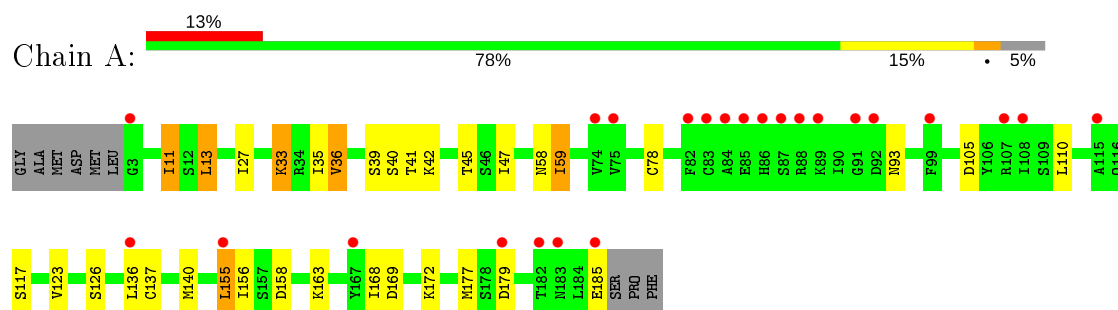
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total	O	0	0
			49	49		
4	B	57	Total	O	0	0
			57	57		

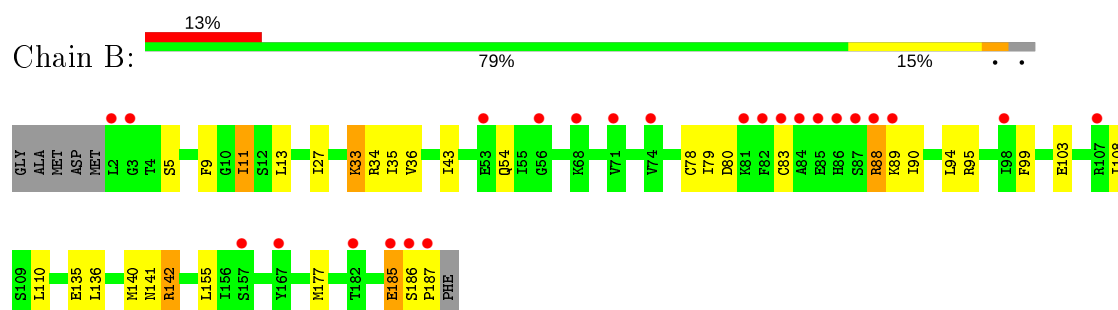
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: IpgB2



#### • Molecule 1: IpgB2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	114.08 Å   114.08 Å   88.77 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	22.88 – 2.30 22.82 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.6 (22.88-2.30) 95.6 (22.82-2.30)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.31 Å)	Xtriage
Refinement program	REFMAC refmac_5.5.0109	Depositor
R, $R_{free}$	0.229   ,   0.292 0.229   ,   0.292	Depositor DCC
$R_{free}$ test set	1267 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.4	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 54.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3117	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: FEO, FLC

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	0.67	0/1495	0.72	1/2000 (0.1%)
1	B	0.70	2/1508 (0.1%)	0.74	0/2019
All	All	0.69	2/3003 (0.1%)	0.73	1/4019 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	135	GLU	CG-CD	6.95	1.62	1.51
1	B	185	GLU	CD-OE2	5.25	1.31	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	LEU	CA-CB-CG	5.15	127.14	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	1477	0	1483	19	0
1	B	1489	0	1501	24	0
2	A	13	0	4	3	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	26	0	9	6	3
3	A	6	0	0	2	2
4	A	49	0	0	0	0
4	B	57	0	0	1	0
All	All	3117	0	2997	46	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 8.

All (46) 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)
2:A:301:FLC:CA	3:A:401:FEO:O	1.67	1.32
1:B:33:LYS:HD3	1:B:33:LYS:N	1.63	1.13
1:B:33:LYS:HD3	1:B:33:LYS:H	1.16	1.05
2:B:302:FLC:OG2	2:B:302:FLC:OHB	1.64	0.98
1:B:88:ARG:HB2	1:B:88:ARG:HH11	1.31	0.95
2:A:301:FLC:OA2	3:A:401:FEO:O	1.85	0.84
1:B:33:LYS:H	1:B:33:LYS:CD	1.92	0.83
1:A:58:ASN:O	1:A:59:ILE:HG22	1.80	0.82
1:B:88:ARG:HH11	1:B:88:ARG:CB	1.99	0.75
1:A:33:LYS:HD3	4:B:198:HOH:O	1.88	0.72
1:B:11:ILE:CD1	1:B:27:ILE:HG23	2.22	0.69
1:A:11:ILE:HD12	1:A:27:ILE:HG23	1.77	0.66
1:B:99:PHE:O	1:B:103:GLU:HG3	2.01	0.60
1:B:88:ARG:HG2	1:B:89:LYS:H	1.67	0.59
1:A:140:MET:CE	1:B:9:PHE:HZ	2.17	0.58
2:A:301:FLC:OA2	2:B:302:FLC:OA2	2.21	0.57
1:A:36:VAL:HG13	1:A:136:LEU:HD13	1.87	0.56
1:B:141:ASN:HB2	2:B:302:FLC:OB2	2.06	0.55
1:A:47:ILE:HD11	1:A:123:VAL:HG22	1.88	0.55
1:A:93:ASN:H	1:A:93:ASN:HD22	1.56	0.52
2:B:302:FLC:CAC	2:B:302:FLC:OB2	2.56	0.52
1:B:83:CYS:SG	1:B:90:ILE:HG12	2.49	0.52
1:B:34:ARG:C	1:B:35:ILE:HD12	2.30	0.52
1:A:137:CYS:O	1:A:140:MET:HG2	2.10	0.51
1:A:140:MET:HE3	1:B:9:PHE:HZ	1.75	0.51
1:B:36:VAL:HG13	1:B:136:LEU:HD13	1.93	0.50
1:B:80:ASP:OD1	1:B:95:ARG:NH1	2.44	0.50
1:B:79:ILE:HD13	1:B:95:ARG:HG2	1.94	0.49
1:A:11:ILE:HD12	1:A:27:ILE:CG2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:THR:O	1:A:45:THR:HG23	2.14	0.48
1:B:11:ILE:HD13	1:B:27:ILE:CG2	2.43	0.48
1:A:47:ILE:HD11	1:A:123:VAL:CG2	2.43	0.47
1:B:11:ILE:HD13	1:B:27:ILE:HG23	1.93	0.47
1:B:88:ARG:HB2	1:B:88:ARG:NH1	2.14	0.46
1:A:169:ASP:OD1	1:A:172:LYS:HB2	2.15	0.46
1:B:141:ASN:OD1	1:B:141:ASN:C	2.53	0.46
1:B:186:SER:HA	1:B:187:PRO:HD3	1.89	0.45
1:A:39:SER:HB3	1:A:42:LYS:HB2	1.98	0.44
1:A:35:ILE:N	1:A:35:ILE:HD12	2.34	0.43
1:B:78:CYS:CB	1:B:177:MET:HG3	2.49	0.43
1:A:78:CYS:SG	1:A:177:MET:HG3	2.59	0.42
1:A:156:ILE:HG12	1:B:27:ILE:HD13	2.01	0.42
1:A:155:LEU:O	1:A:158:ASP:HB2	2.20	0.42
1:A:13:LEU:HD12	1:A:13:LEU:HA	1.71	0.41
2:B:303:FLC:OA1	2:B:303:FLC:OHB	2.38	0.41
1:B:142:ARG:NH1	2:B:302:FLC:CGC	2.85	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)
2:A:301:FLC:OB1	2:B:302:FLC:OG2[8_555]	1.82	0.38
2:A:301:FLC:OA2	3:A:401:FEO:FE1[8_555]	1.83	0.37
3:A:401:FEO:FE1	2:B:302:FLC:OHB[8_555]	1.87	0.33
2:B:302:FLC:OG1	2:B:302:FLC:OG1[8_555]	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	182/192 (95%)	178 (98%)	3 (2%)	1 (0%)	29 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	184/192 (96%)	178 (97%)	6 (3%)	0	100	100
All	All	366/384 (95%)	356 (97%)	9 (2%)	1 (0%)	41	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	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	
1	A	171/177 (97%)	158 (92%)	13 (8%)	13	16
1	B	173/177 (98%)	159 (92%)	14 (8%)	11	15
All	All	344/354 (97%)	317 (92%)	27 (8%)	12	16

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	13	LEU
1	A	33	LYS
1	A	36	VAL
1	A	40	SER
1	A	105	ASP
1	A	110	LEU
1	A	117	SER
1	A	126	SER
1	A	163	LYS
1	A	168	ILE
1	A	179	ASP
1	A	185	GLU
1	B	5	SER
1	B	11	ILE

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Mol	Chain	Res	Type
1	B	13	LEU
1	B	33	LYS
1	B	43	ILE
1	B	54	GLN
1	B	88	ARG
1	B	94	LEU
1	B	108	ILE
1	B	110	LEU
1	B	140	MET
1	B	142	ARG
1	B	155	LEU
1	B	185	GLU

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

Mol	Chain	Res	Type
1	A	93	ASN
1	B	8	ASN
1	B	113	ASN
1	B	120	ASN
1	B	183	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](#)

5 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	FLC	B	302	3	3,12,12	3.66	2 (66%)	3,17,17	8.32	3 (100%)
3	FEO	A	401	2	0,2,2	0.00	-	-		
3	FEO	A	402	-	0,2,2	0.00	-	-		
2	FLC	A	301	3	3,12,12	2.93	2 (66%)	3,17,17	3.99	3 (100%)
2	FLC	B	303	-	3,12,12	0.54	0	3,17,17	0.27	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
2	FLC	B	302	3	-	4/6/16/16	-
2	FLC	A	301	3	-	3/6/16/16	-
2	FLC	B	303	-	-	4/6/16/16	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	302	FLC	CG-CB	-5.05	1.47	1.54
2	A	301	FLC	CA-CB	-4.59	1.48	1.54
2	B	302	FLC	OHB-CB	3.51	1.48	1.43
2	A	301	FLC	OHB-CB	2.14	1.46	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	302	FLC	CB-CA-CAC	-12.03	95.71	114.98
2	B	302	FLC	CG-CB-CA	6.71	127.27	109.33
2	A	301	FLC	CB-CG-CGC	4.74	122.57	114.98
2	B	302	FLC	CB-CG-CGC	-4.23	108.22	114.98
2	A	301	FLC	CB-CA-CAC	-4.04	108.51	114.98
2	A	301	FLC	CG-CB-CA	-3.01	101.28	109.33

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	302	FLC	CAC-CA-CB-CBC
2	B	302	FLC	CAC-CA-CB-OHB
2	B	302	FLC	CA-CB-CG-CGC
2	A	301	FLC	OHB-CB-CG-CGC
2	B	303	FLC	CAC-CA-CB-CBC
2	B	303	FLC	CAC-CA-CB-CG
2	B	303	FLC	CAC-CA-CB-OHB
2	B	302	FLC	OHB-CB-CG-CGC
2	A	301	FLC	CA-CB-CG-CGC
2	A	301	FLC	CBC-CB-CG-CGC
2	B	303	FLC	CA-CB-CG-CGC

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	FLC	5	3
3	A	401	FEO	2	2
2	A	301	FLC	3	2
2	B	303	FLC	1	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	183/192 (95%)	0.74	24 (13%) 3 4	18, 39, 82, 84	0
1	B	186/192 (96%)	0.69	24 (12%) 3 5	15, 37, 79, 102	0
All	All	369/384 (96%)	0.71	48 (13%) 3 4	15, 38, 81, 102	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	187	PRO	8.0
1	A	87	SER	7.6
1	B	186	SER	7.4
1	A	84	ALA	5.4
1	B	2	LEU	4.8
1	A	83	CYS	4.8
1	B	83	CYS	4.7
1	A	107	ARG	4.3
1	B	185	GLU	4.0
1	B	182	THR	3.8
1	A	86	HIS	3.7
1	B	3	GLY	3.4
1	A	3	GLY	3.3
1	A	74	VAL	3.3
1	B	87	SER	3.3
1	A	88	ARG	3.3
1	A	183	ASN	3.2
1	B	82	PHE	3.2
1	B	107	ARG	3.1
1	A	89	LYS	2.8
1	B	84	ALA	2.8
1	A	92	ASP	2.8
1	A	179	ASP	2.8
1	A	91	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	74	VAL	2.7
1	A	85	GLU	2.7
1	B	71	VAL	2.7
1	A	185	GLU	2.7
1	B	56	GLY	2.7
1	A	115	ALA	2.6
1	B	157	SER	2.5
1	B	167	TYR	2.5
1	B	88	ARG	2.5
1	A	167	TYR	2.4
1	B	98	ILE	2.4
1	B	86	HIS	2.4
1	B	89	LYS	2.4
1	A	75	VAL	2.4
1	B	53	GLU	2.3
1	B	81	LYS	2.3
1	A	182	THR	2.3
1	A	99	PHE	2.3
1	A	136	LEU	2.2
1	B	68	LYS	2.2
1	A	82	PHE	2.1
1	A	108	ILE	2.1
1	A	155	LEU	2.0
1	B	85	GLU	2.0

## 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, 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
3	FEO	A	402	3/3	0.45	0.15	118,118,118,119	3
2	FLC	B	303	13/13	0.76	0.19	107,109,112,113	0
2	FLC	B	302	13/13	0.92	0.16	34,48,52,52	0
2	FLC	A	301	13/13	0.97	0.14	48,56,62,65	0
3	FEO	A	401	3/3	0.99	0.11	42,42,43,44	3

## 6.5 Other polymers [i](#)

There are no such residues in this entry.