



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 11:25 pm GMT

PDB ID : 3D6N  
Title : Crystal Structure of Aquifex Dihydroorotase Activated by Aspartate Transcarbamoylase  
Authors : Edwards, B.F.P.  
Deposited on : 2008-05-20  
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.

---

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
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) : trunk30967

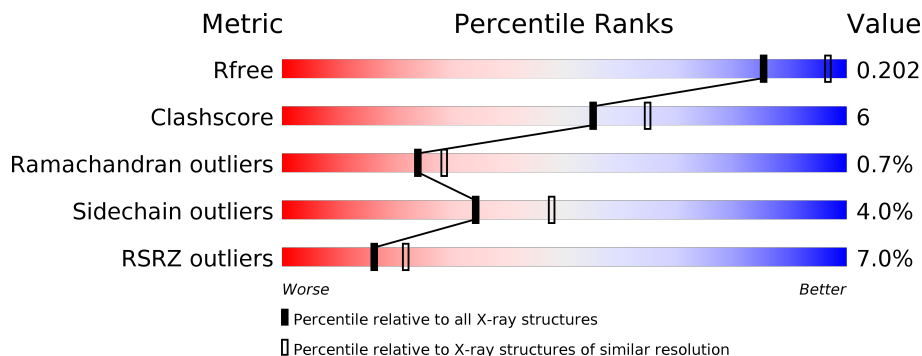
# 1 Overall quality at a glance ⓘ

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))
$R_{free}$	111664	4477 (2.30-2.30)
Clashscore	122126	5072 (2.30-2.30)
Ramachandran outliers	120053	5022 (2.30-2.30)
Sidechain outliers	120020	5021 (2.30-2.30)
RSRZ outliers	108989	4374 (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. 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	422	<div> <div>11%</div> <div> <div></div> <div>86%</div> <div>11%</div> </div> </div>
2	B	291	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>13%</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
4	FLC	A	424	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6081 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 Dihydroorotase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3256	2071	554	615	16			

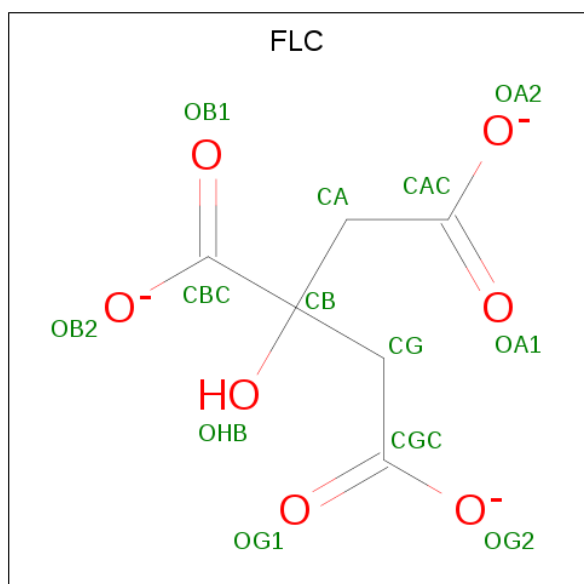
- Molecule 2 is a protein called Aspartate carbamoyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	291	Total	C	N	O	S	0	0	0
			2376	1541	395	436	4			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 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
4	A	1	Total 13	C 6	O 7	0	0
4	B	1	Total 13	C 6	O 7	0	0

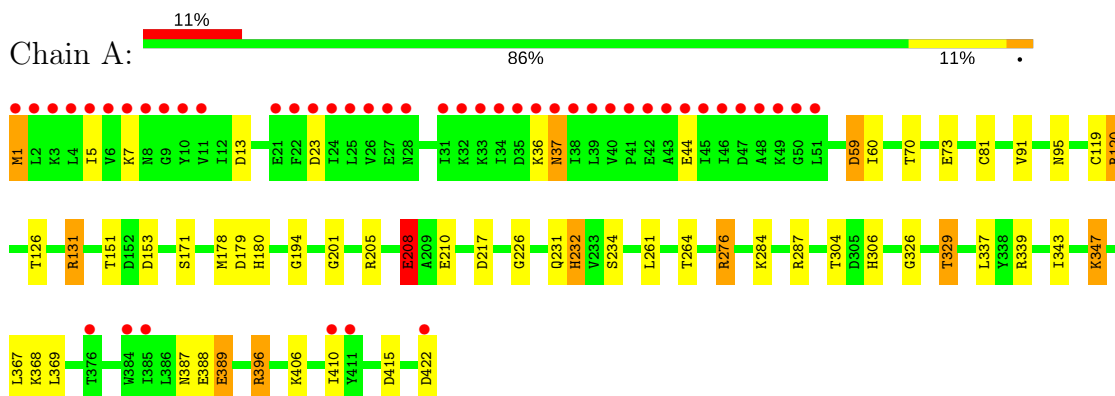
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	209	Total 209	O 209	0	0
5	B	213	Total 213	O 213	0	0

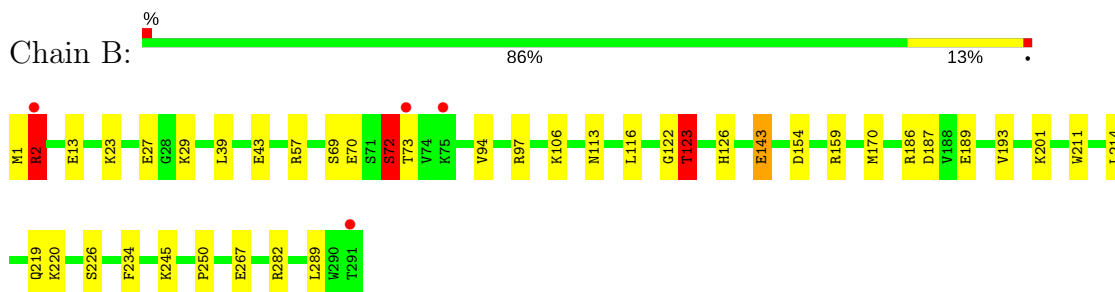
### 3 Residue-property plots [i](#)

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: Dihydroorotase



#### • Molecule 2: Aspartate carbamoyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	158.44Å 158.44Å 233.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	65.00 – 2.30 53.73 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (65.00-2.30) 100.0 (53.73-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.80	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.02 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.164 , 0.204 0.163 , 0.202	Depositor DCC
$R_{free}$ test set	2544 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.3	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 48.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6081	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.85	3/3310 (0.1%)	0.85	6/4472 (0.1%)
2	B	1.10	4/2431 (0.2%)	0.95	8/3274 (0.2%)
All	All	0.97	7/5741 (0.1%)	0.89	14/7746 (0.2%)

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
2	B	1	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	143	GLU	CB-CG	9.72	1.70	1.52
2	B	143	GLU	CG-CD	8.23	1.64	1.51
1	A	119	CYS	CB-SG	-7.00	1.70	1.82
1	A	81	CYS	CB-SG	-6.58	1.71	1.82
2	B	13	GLU	CG-CD	5.79	1.60	1.51
2	B	245	LYS	CE-NZ	5.46	1.62	1.49
1	A	208	GLU	CB-CG	5.06	1.61	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ARG	NE-CZ-NH2	-9.04	115.78	120.30
1	A	205	ARG	NE-CZ-NH2	-8.94	115.83	120.30
2	B	123	THR	OG1-CB-CG2	6.23	124.33	110.00
2	B	123	THR	N-CA-CB	-6.19	98.54	110.30

*Continued on next page...*



Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	205	ARG	NE-CZ-NH1	5.98	123.29	120.30
2	B	57	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	A	231	GLN	C-N-CA	-5.75	107.34	121.70
1	A	59	ASP	CB-CG-OD1	5.62	123.35	118.30
2	B	214	LEU	CA-CB-CG	-5.52	102.59	115.30
2	B	282	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	A	131	ARG	NE-CZ-NH1	5.37	122.99	120.30
2	B	143	GLU	OE1-CD-OE2	-5.26	116.98	123.30
2	B	154	ASP	CB-CG-OD1	5.18	122.96	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	123	THR	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	72	SER	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	3256	0	3362	35	0
2	B	2376	0	2393	32	0
3	A	1	0	0	0	0
4	A	13	0	5	1	0
4	B	13	0	5	0	0
5	A	209	0	0	9	0
5	B	213	0	0	8	1
All	All	6081	0	5765	64	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 6.

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:A:120:ARG:HD3	5:A:579:HOH:O	1.36	1.24
2:B:69:SER:HB3	2:B:72:SER:HA	1.20	1.09
2:B:70:GLU:OE1	5:B:491:HOH:O	1.81	0.97
2:B:1:MET:HE1	5:B:425:HOH:O	1.65	0.95
2:B:2:ARG:HG3	2:B:289:LEU:HD22	1.58	0.85
2:B:69:SER:CB	2:B:72:SER:HA	2.05	0.84
2:B:211:TRP:HZ3	2:B:234:PHE:O	1.64	0.79
1:A:201:GLY:O	2:B:123:THR:CG2	2.33	0.77
2:B:2:ARG:HG3	2:B:289:LEU:CD2	2.17	0.74
1:A:59:ASP:OD2	1:A:304:THR:HG23	1.91	0.71
1:A:232:HIS:NE2	5:A:577:HOH:O	2.15	0.70
2:B:69:SER:HB3	2:B:72:SER:CA	2.12	0.70
1:A:339:ARG:NH2	1:A:406:LYS:O	2.26	0.69
2:B:126:HIS:HD2	5:B:427:HOH:O	1.76	0.69
2:B:2:ARG:HH11	2:B:2:ARG:HG2	1.58	0.67
2:B:70:GLU:HA	5:B:494:HOH:O	1.97	0.65
1:A:179:ASP:OD2	1:A:217:ASP:OD2	2.15	0.65
2:B:2:ARG:CG	2:B:2:ARG:HH11	2.12	0.62
1:A:264:THR:HG21	1:A:284:LYS:HE3	1.83	0.61
2:B:23:LYS:O	2:B:27:GLU:HG3	1.99	0.61
1:A:178:MET:HE3	5:A:468:HOH:O	2.01	0.60
1:A:178:MET:CE	5:A:468:HOH:O	2.51	0.59
1:A:276:ARG:NH2	5:A:547:HOH:O	2.27	0.58
1:A:70:THR:HA	1:A:73:GLU:O	2.05	0.57
2:B:2:ARG:CG	2:B:289:LEU:CD2	2.83	0.56
1:A:201:GLY:O	2:B:123:THR:HG21	2.07	0.54
4:A:424:FLC:OA1	4:A:424:FLC:CGC	2.56	0.53
2:B:2:ARG:NH1	2:B:2:ARG:HG2	2.22	0.53
1:A:387:ASN:OD1	1:A:389:GLU:HG2	2.09	0.52
2:B:219:GLN:HG3	2:B:226:SER:HA	1.92	0.52
1:A:388:GLU:OE1	1:A:396:ARG:NH2	2.43	0.52
1:A:343:ILE:HB	1:A:347:LYS:HG2	1.91	0.51
2:B:211:TRP:CZ3	2:B:234:PHE:O	2.55	0.50
2:B:43:GLU:HB3	2:B:97:ARG:HG2	1.92	0.50
2:B:122:GLY:O	2:B:159:ARG:HD3	2.11	0.50
1:A:120:ARG:NH2	5:A:562:HOH:O	2.45	0.50
1:A:261:LEU:HD13	1:A:337:LEU:HD11	1.95	0.49
1:A:7:LYS:HA	1:A:23:ASP:OD1	2.13	0.49
1:A:208:GLU:OE2	2:B:186:ARG:NH2	2.43	0.48
1:A:329:THR:HG23	5:A:530:HOH:O	2.12	0.48
1:A:276:ARG:NE	5:A:547:HOH:O	2.36	0.48

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:SD	1:A:44:GLU:HG3	2.53	0.47
2:B:2:ARG:HG2	2:B:289:LEU:HD23	1.96	0.47
1:A:210:GLU:OE1	1:A:234:SER:OG	2.28	0.47
1:A:95:ASN:HB3	1:A:153:ASP:O	2.15	0.47
1:A:59:ASP:OD2	1:A:304:THR:CG2	2.61	0.47
1:A:60:ILE:HA	1:A:91:VAL:HB	1.97	0.47
1:A:13:ASP:HB2	1:A:367:LEU:HD12	1.96	0.47
1:A:410:ILE:HA	1:A:422:ASP:HB2	1.96	0.46
2:B:2:ARG:CG	2:B:289:LEU:HD23	2.46	0.46
1:A:194:GLY:HA2	2:B:187:ASP:OD2	2.15	0.46
2:B:267:GLU:HB2	5:B:416:HOH:O	2.16	0.45
1:A:180:HIS:ND1	5:A:577:HOH:O	2.36	0.45
2:B:126:HIS:CD2	5:B:427:HOH:O	2.61	0.43
1:A:126:THR:HG21	1:A:131:ARG:HG3	1.99	0.43
1:A:208:GLU:HG3	1:A:208:GLU:H	1.32	0.43
2:B:94:VAL:HB	2:B:116:LEU:HD23	1.99	0.43
2:B:170:MET:HE2	2:B:170:MET:HB3	1.99	0.42
1:A:126:THR:HA	1:A:151:THR:O	2.20	0.42
1:A:284:LYS:HE2	1:A:287:ARG:HH11	1.84	0.42
2:B:113:ASN:HB2	5:B:482:HOH:O	2.21	0.41
2:B:201:LYS:HZ3	2:B:201:LYS:HG3	1.75	0.41
2:B:1:MET:CE	5:B:425:HOH:O	2.45	0.40
1:A:171:SER:HB3	1:A:226:GLY:HA2	2.04	0.40

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)
5:B:487:HOH:O	5:B:492:HOH:O[17_555]	2.11	0.09

## 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	420/422 (100%)	398 (95%)	19 (4%)	3 (1%)	24	29
2	B	289/291 (99%)	276 (96%)	11 (4%)	2 (1%)	24	29
All	All	709/713 (99%)	674 (95%)	30 (4%)	5 (1%)	24	29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	72	SER
1	A	37	ASN
1	A	306	HIS
2	B	250	PRO
1	A	326	GLY

### 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	362/362 (100%)	347 (96%)	15 (4%)	33	46
2	B	262/262 (100%)	252 (96%)	10 (4%)	36	50
All	All	624/624 (100%)	599 (96%)	25 (4%)	34	48

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	ILE
1	A	36	LYS
1	A	37	ASN
1	A	120	ARG
1	A	208	GLU
1	A	232	HIS
1	A	276	ARG
1	A	329	THR
1	A	347	LYS
1	A	368	LYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	369	LEU
1	A	389	GLU
1	A	396	ARG
1	A	415	ASP
2	B	2	ARG
2	B	29	LYS
2	B	39	LEU
2	B	73	THR
2	B	106	LYS
2	B	123	THR
2	B	143	GLU
2	B	189	GLU
2	B	193	VAL
2	B	220	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	126	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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
4	FLC	A	424	3	3,12,12	1.35	1 (33%)	3,17,17	1.44	0
4	FLC	B	292	-	3,12,12	0.71	0	3,17,17	1.77	1 (33%)

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
4	FLC	A	424	3	-	0/6/16/16	0/0/0/0
4	FLC	B	292	-	-	0/6/16/16	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	424	FLC	CA-CB	2.12	1.57	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	292	FLC	CB-CG-CGC	-2.74	110.84	114.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	424	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	422/422 (100%)	0.37	46 (10%) <b>5</b> <b>8</b>	24, 38, 49, 68	0
2	B	291/291 (100%)	-0.41	4 (1%) <b>75</b> <b>80</b>	18, 26, 44, 78	0
All	All	713/713 (100%)	0.05	50 (7%) <b>16</b> <b>22</b>	18, 32, 48, 78	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	12.9
1	A	41	PRO	7.0
1	A	43	ALA	6.4
1	A	42	GLU	6.2
1	A	26	VAL	5.7
1	A	2	LEU	5.2
1	A	39	LEU	5.2
1	A	38	ILE	5.1
1	A	4	LEU	5.1
1	A	6	VAL	5.0
1	A	40	VAL	4.8
1	A	3	LYS	4.6
2	B	73	THR	4.6
1	A	5	ILE	4.5
1	A	45	ILE	4.4
1	A	44	GLU	4.4
1	A	46	ILE	4.3
1	A	25	LEU	4.0
1	A	47	ASP	3.9
1	A	24	ILE	3.7
1	A	35	ASP	3.7
1	A	36	LYS	3.7
1	A	27	GLU	3.6
1	A	10	TYR	3.5

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	11	VAL	3.5
1	A	49	LYS	3.5
1	A	37	ASN	3.4
1	A	384	TRP	3.3
1	A	8	ASN	3.2
1	A	7	LYS	3.2
1	A	34	ILE	3.1
1	A	33	LYS	3.0
1	A	411	TYR	3.0
1	A	50	GLY	2.9
1	A	28	ASN	2.9
1	A	31	ILE	2.9
1	A	9	GLY	2.9
1	A	376	THR	2.8
1	A	32	LYS	2.8
1	A	48	ALA	2.7
2	B	291	THR	2.6
1	A	21	GLU	2.6
1	A	410	ILE	2.5
1	A	23	ASP	2.4
1	A	22	PHE	2.3
1	A	51	LEU	2.3
2	B	75	LYS	2.3
1	A	385	ILE	2.2
1	A	422	ASP	2.2
2	B	2	ARG	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 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
4	FLC	A	424	13/13	0.67	0.41	31,38,46,50	0
4	FLC	B	292	13/13	0.82	0.36	70,72,75,75	0
3	ZN	A	423	1/1	0.99	0.10	54,54,54,54	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.