



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 7, 2017 – 04:06 AM EST

PDB ID : 4H4O  
Title : Crystal Structure of HIV-1 Reverse Transcriptase (RT) in Complex with (E)-3-(3-(2-(2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy)-4-fluorophenoxy)-5-fluorophenyl)acrylonitrile (JLJ506), A Non-nucleoside inhibitor  
Authors : Frey, K.M.; Anderson, K.S.  
Deposited on : unknown  
Resolution : 2.90 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

<http://wwpdb.org/validation/2016/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  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

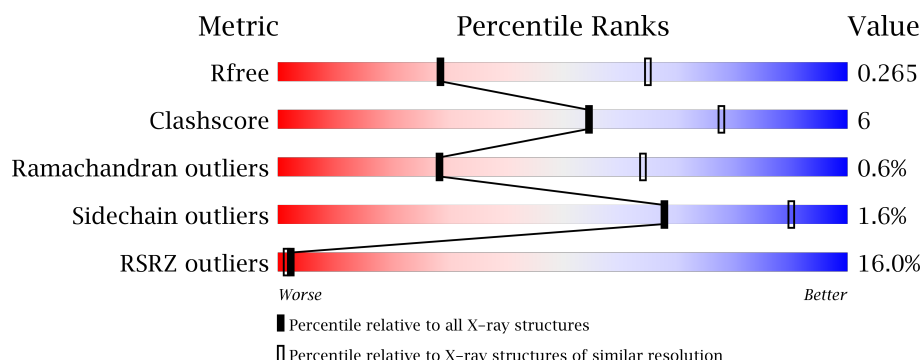
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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	557	<div> <div>16%</div> <div>85%</div> <div>13%</div> <div>.</div> </div>
2	B	428	<div> <div>16%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8082 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 Reverse transcriptase/ribonuclease H, Exoribonuclease H, p66 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4465	2891	742	825	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	INITIATING METHIONINE	UNP P03366
A	0	VAL	-	EXPRESSION TAG	UNP P03366
A	172	ALA	LYS	ENGINEERED MUTATION	UNP P03366
A	173	ALA	LYS	ENGINEERED MUTATION	UNP P03366
A	280	SER	CYS	ENGINEERED MUTATION	UNP P03366

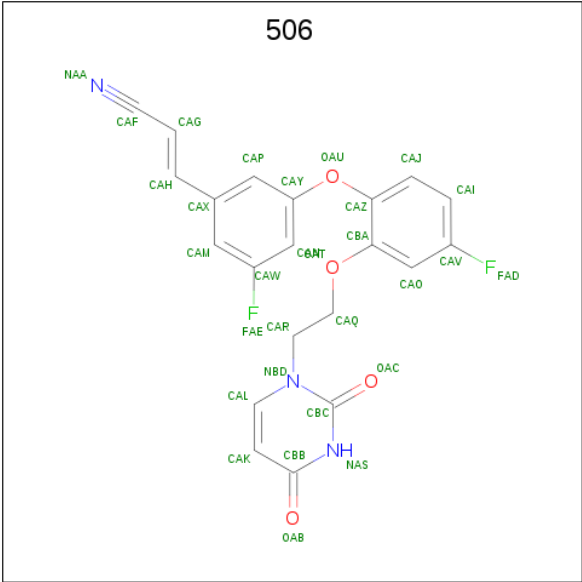
- Molecule 2 is a protein called Reverse transcriptase/ribonuclease H, Exoribonuclease H, p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	428	Total	C	N	O	S	0	0	0
			3538	2305	586	640	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	ENGINEERED MUTATION	UNP P03366

- Molecule 3 is (2E)-3-(3-{2-[2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy]-4-fluorophenoxy}-5-fluorophenyl)prop-2-enenitrile (three-letter code: 506) (formula: C<sub>21</sub>H<sub>15</sub>F<sub>2</sub>N<sub>3</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			30	21	2	3	4		

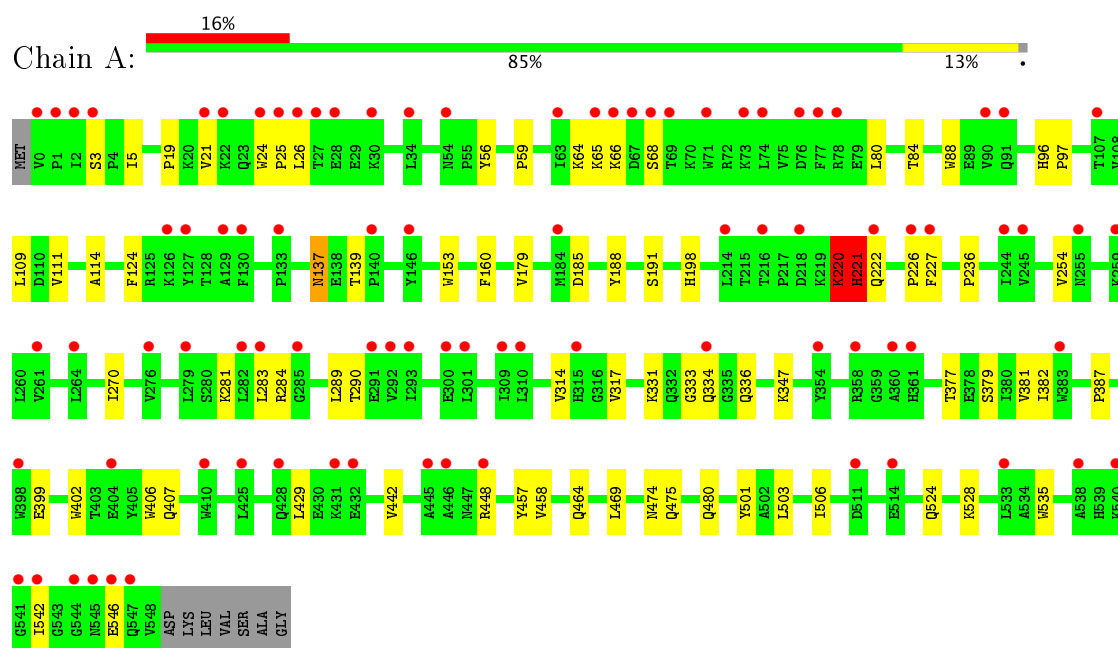
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	27	Total	O	0	0
			27	27		

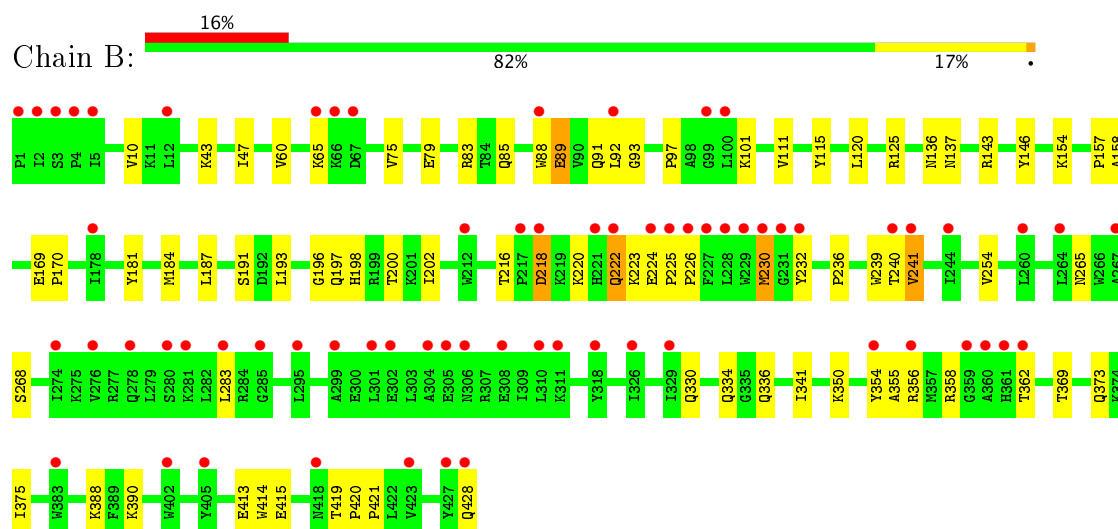
### 3 Residue-property plots [i](#)

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: Reverse transcriptase/ribonuclease H, Exoribonuclease H, p66 RT



- Molecule 2: Reverse transcriptase/ribonuclease H, Exoribonuclease H, p51 RT



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	225.45Å 69.58Å 104.40Å 90.00° 106.21° 90.00°	Depositor
Resolution (Å)	41.72 – 2.90 41.72 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (41.72-2.90) 99.8 (41.72-2.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.27 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, $R_{free}$	0.232 , 0.265 0.232 , 0.265	Depositor DCC
$R_{free}$ test set	1997 reflections (5.75%)	DCC
Wilson B-factor (Å <sup>2</sup> )	79.0	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 113.0	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	0.93	EDS
Total number of atoms	8082	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	164.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: 506

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	0.23	0/4583	0.45	0/6231
2	B	0.24	0/3643	0.48	1/4952 (0.0%)
All	All	0.24	0/8226	0.46	1/11183 (0.0%)

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	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	89	GLU	N-CA-C	-5.05	97.36	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	220	LYS	Peptide
2	B	419	THR	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	4465	0	4516	49	0
2	B	3538	0	3576	52	0
3	A	30	0	15	2	0
4	A	22	0	0	0	1
4	B	27	0	0	1	0
All	All	8082	0	8107	96	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 (96) 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:220:LYS:HB3	1:A:222:GLN:H	1.51	0.76
1:A:331:LYS:HE2	1:A:333:GLY:HA2	1.68	0.76
1:A:191:SER:OG	1:A:198:HIS:ND1	2.23	0.68
1:A:64:LYS:HD2	1:A:68:SER:HA	1.75	0.67
1:A:226:PRO:HA	1:A:227:PHE:HB2	1.77	0.66
1:A:317:VAL:HG11	1:A:347:LYS:HD3	1.79	0.65
2:B:137:ASN:O	4:B:505:HOH:O	2.15	0.64
2:B:336:GLN:HG2	2:B:355:ALA:HB2	1.80	0.63
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.61	0.63
2:B:390:LYS:NZ	2:B:415:GLU:OE2	2.34	0.61
1:A:88:TRP:CE2	2:B:143:ARG:HD2	2.37	0.59
1:A:334:GLN:HB2	1:A:336:GLN:HE21	1.68	0.58
2:B:157:PRO:HG3	2:B:184:MET:HA	1.86	0.58
2:B:354:TYR:HB3	2:B:356:ARG:HH12	1.71	0.56
1:A:220:LYS:O	1:A:221:HIS:HB2	2.04	0.55
2:B:191:SER:OG	2:B:198:HIS:ND1	2.30	0.54
1:A:220:LYS:CB	1:A:222:GLN:H	2.20	0.54
1:A:270:ILE:HG13	1:A:314:VAL:HG13	1.90	0.54
2:B:91:GLN:HG2	2:B:92:LEU:HD12	1.88	0.53
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.90	0.52
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.45	0.52
1:A:84:THR:HG22	1:A:124:PHE:HZ	1.75	0.52

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:236:PRO:HA	2:B:239:TRP:CG	2.46	0.51
1:A:88:TRP:CD2	2:B:143:ARG:HD2	2.46	0.51
2:B:79:GLU:OE2	2:B:83:ARG:NH1	2.44	0.50
1:A:281:LYS:HG3	1:A:284:ARG:HE	1.76	0.50
2:B:89:GLU:O	2:B:93:GLY:HA3	2.12	0.50
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.93	0.50
2:B:97:PRO:HG3	2:B:181:TYR:HB2	1.94	0.49
2:B:239:TRP:O	2:B:350:LYS:HE2	2.13	0.49
2:B:92:LEU:HB3	2:B:158:ALA:HB1	1.95	0.48
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.49	0.48
1:A:220:LYS:HD2	1:A:220:LYS:HA	1.55	0.48
2:B:336:GLN:OE1	2:B:428:GLN:NE2	2.46	0.48
2:B:240:THR:OG1	2:B:241:VAL:N	2.46	0.48
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.96	0.48
2:B:334:GLN:O	2:B:358:ARG:NH1	2.47	0.47
1:A:19:PRO:O	1:A:56:TYR:HB3	2.14	0.47
1:A:254:VAL:HG13	1:A:283:LEU:HD11	1.96	0.47
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.49	0.47
1:A:236:PRO:HA	3:A:601:506:CBB	2.44	0.47
1:A:399:GLU:HA	1:A:402:TRP:NE1	2.29	0.47
2:B:120:LEU:HD23	2:B:125:ARG:HG2	1.98	0.46
2:B:88:TRP:N	2:B:88:TRP:CD1	2.84	0.46
2:B:354:TYR:HB3	2:B:356:ARG:NH1	2.29	0.46
1:A:111:VAL:HG23	1:A:114:ALA:HB2	1.98	0.46
1:A:65:LYS:HG2	1:A:66:LYS:H	1.80	0.46
2:B:193:LEU:HB3	2:B:197:GLN:HG3	1.98	0.45
2:B:222:GLN:HB3	2:B:223:LYS:H	1.59	0.45
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.98	0.45
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.65	0.45
1:A:448:ARG:HE	1:A:474:ASN:ND2	2.15	0.45
1:A:25:PRO:HA	1:A:26:LEU:HA	1.72	0.45
1:A:442:VAL:HG12	1:A:457:TYR:HB3	1.99	0.45
1:A:406:TRP:CH2	2:B:420:PRO:HD2	2.52	0.45
1:A:406:TRP:CZ3	1:A:407:GLN:HG3	2.52	0.45
2:B:115:TYR:OH	2:B:184:MET:O	2.33	0.45
1:A:379:SER:CB	1:A:387:PRO:HD3	2.47	0.44
1:A:377:THR:O	1:A:381:VAL:HG23	2.17	0.44
1:A:111:VAL:HG22	1:A:185:ASP:O	2.17	0.44
1:A:254:VAL:HB	1:A:289:LEU:O	2.17	0.44
2:B:220:LYS:HD3	2:B:220:LYS:HA	1.68	0.44
2:B:265:ASN:O	2:B:268:SER:OG	2.27	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.98	0.44
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.53	0.44
2:B:10:VAL:HG22	2:B:88:TRP:CH2	2.52	0.44
2:B:220:LYS:C	2:B:222:GLN:H	2.21	0.44
2:B:101:LYS:O	2:B:236:PRO:HB2	2.17	0.44
2:B:43:LYS:HE3	2:B:43:LYS:HB2	1.86	0.44
1:A:3:SER:HB3	1:A:5:ILE:HG13	2.00	0.43
1:A:188:TYR:CZ	3:A:601:506:H14	2.53	0.43
1:A:96:HIS:HA	1:A:97:PRO:HD3	1.92	0.43
2:B:341:ILE:HD11	2:B:375:ILE:HG23	2.00	0.43
1:A:469:LEU:HD11	1:A:480:GLN:HG2	2.00	0.43
2:B:369:THR:HG22	2:B:373:GLN:NE2	2.34	0.43
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.99	0.43
1:A:524:GLN:O	1:A:528:LYS:HG2	2.19	0.42
1:A:458:VAL:HA	1:A:464:GLN:HB3	2.01	0.42
2:B:198:HIS:O	2:B:202:ILE:HG12	2.20	0.42
1:A:429:LEU:HD11	1:A:506:ILE:HG22	2.02	0.42
1:A:24:TRP:N	1:A:25:PRO:HD2	2.35	0.42
1:A:503:LEU:HD22	1:A:535:TRP:HB2	2.01	0.42
2:B:224:GLU:HA	2:B:225:PRO:HD2	1.96	0.42
1:A:382:ILE:O	2:B:136:ASN:HB2	2.20	0.41
2:B:187:LEU:HA	2:B:187:LEU:HD12	1.90	0.41
1:A:458:VAL:HG22	1:A:464:GLN:HB3	2.01	0.41
1:A:406:TRP:CZ3	2:B:420:PRO:HD2	2.56	0.41
2:B:85:GLN:HG3	2:B:154:LYS:HB3	2.02	0.41
2:B:388:LYS:HD2	2:B:413:GLU:HB2	2.02	0.41
2:B:216:THR:O	2:B:218:ASP:N	2.49	0.41
2:B:169:GLU:HB3	2:B:170:PRO:HD3	2.03	0.41
2:B:196:GLY:O	2:B:200:THR:HG23	2.21	0.40
2:B:236:PRO:O	2:B:239:TRP:HB2	2.21	0.40
1:A:109:LEU:HD21	1:A:221:HIS:HE1	1.87	0.40
2:B:369:THR:HG22	2:B:373:GLN:HE21	1.85	0.40
2:B:220:LYS:HG3	2:B:230:MET:HG3	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)
4:A:718:HOH:O	4:A:718:HOH:O[2_756]	2.16	0.04

## 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	547/557 (98%)	519 (95%)	24 (4%)	4 (1%)	25	60
2	B	426/428 (100%)	397 (93%)	27 (6%)	2 (0%)	32	68
All	All	973/985 (99%)	916 (94%)	51 (5%)	6 (1%)	28	64

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	HIS
1	A	220	LYS
1	A	137	ASN
2	B	226	PRO
1	A	542	ILE
2	B	65	LYS

### 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	489/495 (99%)	483 (99%)	6 (1%)	75	93
2	B	390/390 (100%)	382 (98%)	8 (2%)	59	86
All	All	879/885 (99%)	865 (98%)	14 (2%)	68	90

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	139	THR
1	A	179	VAL
1	A	221	HIS
1	A	290	THR
1	A	546	GLU
2	B	218	ASP
2	B	222	GLN
2	B	230	MET
2	B	232	TYR
2	B	241	VAL
2	B	330	GLN
2	B	362	THR
2	B	414	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	361	HIS
1	A	474	ASN
2	B	197	GLN
2	B	278	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	$\# Z  > 2$
3	506	A	601	-	29,32,32	1.69	5 (17%)	35,43,43	2.69	10 (28%)

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	506	A	601	-	-	0/13/14/14	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	506	CAM-CAW	2.17	1.41	1.37
3	A	601	506	CAR-CAQ	2.37	1.56	1.51
3	A	601	506	CAZ-CBA	2.47	1.45	1.40
3	A	601	506	CAL-NBD	2.73	1.40	1.36
3	A	601	506	CBB-NAS	5.03	1.42	1.33

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	506	OAT-CBA-CAO	-3.76	114.45	123.61
3	A	601	506	CAK-CBB-NAS	-3.54	114.66	123.12
3	A	601	506	CAL-NBD-CBC	-2.18	118.62	122.14
3	A	601	506	CAI-CAV-CAO	-2.06	120.59	123.29
3	A	601	506	CAI-CAJ-CAZ	2.07	124.00	120.06
3	A	601	506	OAU-CAZ-CBA	2.21	124.08	117.39
3	A	601	506	CBA-CAO-CAV	2.89	122.31	116.51
3	A	601	506	CAX-CAP-CAY	3.26	122.62	119.88
3	A	601	506	OAT-CBA-CAZ	5.74	127.74	115.76
3	A	601	506	CBB-NAS-CBC	11.52	124.02	114.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	506	2	0

## 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 ⓘ

### 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	549/557 (98%)	1.04	89 (16%) 2 1	112, 157, 242, 303	1 (0%)
2	B	428/428 (100%)	1.23	67 (15%) 2 1	114, 148, 248, 369	1 (0%)
All	All	977/985 (99%)	1.12	156 (15%) 2 1	112, 154, 246, 369	2 (0%)

All (156) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	PRO	39.7
2	B	2	ILE	18.7
1	A	541	GLY	18.0
2	B	224	GLU	16.4
2	B	231	GLY	16.0
1	A	25	PRO	14.8
2	B	4	PRO	12.5
1	A	24	TRP	12.5
1	A	545	ASN	12.0
2	B	67	ASP	10.7
1	A	67	ASP	9.8
1	A	2	ILE	9.0
1	A	542	ILE	8.8
2	B	226	PRO	8.7
2	B	217	PRO	8.1
2	B	225	PRO	7.5
2	B	3	SER	7.4
2	B	229	TRP	7.1
2	B	221	HIS	7.0
2	B	66	LYS	7.0
2	B	228	LEU	6.8
2	B	227	PHE	6.6
2	B	361	HIS	6.3
2	B	428	GLN	6.3

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	B	360	ALA	6.1
1	A	448	ARG	6.0
2	B	418	ASN	5.9
1	A	546	GLU	5.8
2	B	230	MET	5.7
1	A	26	LEU	5.6
2	B	310	LEU	5.6
2	B	301	LEU	5.4
2	B	354	TYR	5.3
1	A	22	LYS	5.0
1	A	264	LEU	5.0
1	A	68	SER	4.9
1	A	107	THR	4.8
2	B	5	ILE	4.7
2	B	88	TRP	4.7
1	A	28	GLU	4.7
1	A	227	PHE	4.6
1	A	360	ALA	4.6
2	B	299	ALA	4.6
1	A	21	VAL	4.5
2	B	283	LEU	4.5
2	B	218	ASP	4.5
1	A	226	PRO	4.4
1	A	354	TYR	4.4
1	A	54	ASN	4.4
1	A	129	ALA	4.4
2	B	311	LYS	4.4
1	A	133	PRO	4.4
1	A	283	LEU	4.3
1	A	77	PHE	4.3
1	A	63	ILE	4.3
1	A	255	ASN	4.3
1	A	130	PHE	4.3
1	A	0	VAL	4.2
1	A	358	ARG	4.1
1	A	261	VAL	4.0
2	B	304	ALA	3.9
1	A	309	ILE	3.9
1	A	514	GLU	3.9
2	B	326	ILE	3.8
2	B	295	LEU	3.8
1	A	540	LYS	3.6

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	74	LEU	3.6
1	A	291	GLU	3.6
1	A	446	ALA	3.6
2	B	285	GLY	3.5
1	A	71	TRP	3.5
2	B	329	ILE	3.4
1	A	361	HIS	3.4
2	B	318	TYR	3.4
2	B	65	LYS	3.4
2	B	302	GLU	3.4
1	A	222	GLN	3.3
2	B	305	GLU	3.3
2	B	306	ASN	3.3
1	A	282	LEU	3.3
1	A	216	THR	3.3
1	A	245	VAL	3.2
1	A	30	LYS	3.2
1	A	78	ARG	3.2
2	B	232	TYR	3.2
1	A	431	LYS	3.2
2	B	359	GLY	3.1
1	A	91	GLN	3.1
1	A	538	ALA	3.0
1	A	445	ALA	2.9
1	A	293	ILE	2.9
1	A	140	PRO	2.9
1	A	3	SER	2.9
1	A	410	TRP	2.9
1	A	544	GLY	2.9
1	A	127	TYR	2.8
2	B	427	TYR	2.8
1	A	334	GLN	2.8
1	A	218	ASP	2.8
2	B	264	LEU	2.8
2	B	356	ARG	2.8
2	B	362	THR	2.7
2	B	423	VAL	2.7
1	A	300	GLU	2.7
1	A	90	VAL	2.7
1	A	279	LEU	2.6
1	A	126	LYS	2.6
2	B	405	TYR	2.6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	B	276	VAL	2.6
1	A	34	LEU	2.6
2	B	100	LEU	2.6
1	A	404	GLU	2.6
1	A	292	VAL	2.6
2	B	241	VAL	2.6
1	A	285	GLY	2.6
1	A	511	ASP	2.6
1	A	69	THR	2.6
1	A	76	ASP	2.6
1	A	315	HIS	2.5
1	A	244	ILE	2.5
1	A	276	VAL	2.5
1	A	398	TRP	2.5
2	B	281	LYS	2.5
1	A	259	LYS	2.5
2	B	280	SER	2.4
1	A	428	GLN	2.4
1	A	1	PRO	2.4
1	A	73	LYS	2.4
2	B	308	GLU	2.3
1	A	310	LEU	2.3
2	B	383	TRP	2.3
2	B	278	GLN	2.3
1	A	533	LEU	2.3
2	B	274	ILE	2.3
1	A	432	GLU	2.3
1	A	27	THR	2.2
2	B	222	GLN	2.2
1	A	65	LYS	2.2
1	A	425	LEU	2.2
2	B	178	ILE	2.2
1	A	66	LYS	2.2
2	B	12	LEU	2.1
1	A	184	MET	2.1
1	A	214	LEU	2.1
2	B	99	GLY	2.1
1	A	146	TYR	2.1
1	A	383	TRP	2.1
2	B	92	LEU	2.1
2	B	260	LEU	2.0
2	B	212	TRP	2.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	B	402	TRP	2.0
2	B	240	THR	2.0
1	A	301	LEU	2.0
2	B	267	ALA	2.0
1	A	547	GLN	2.0
2	B	244	ILE	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	506	A	601	30/30	0.79	0.39	1.47	125,145,153,160	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.