



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:58 PM GMT

PDB ID : 5C42  
Title : Crystal Structure of HIV-1 Reverse Transcriptase (K101P) Variant in Complex with 8-(2-(2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy)phenoxy)indoline-2-carbonitrile (JLJ555), a non-nucleoside inhibitor  
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Deposited on : 2015-06-17  
Resolution : 3.50 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.

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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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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-20026982

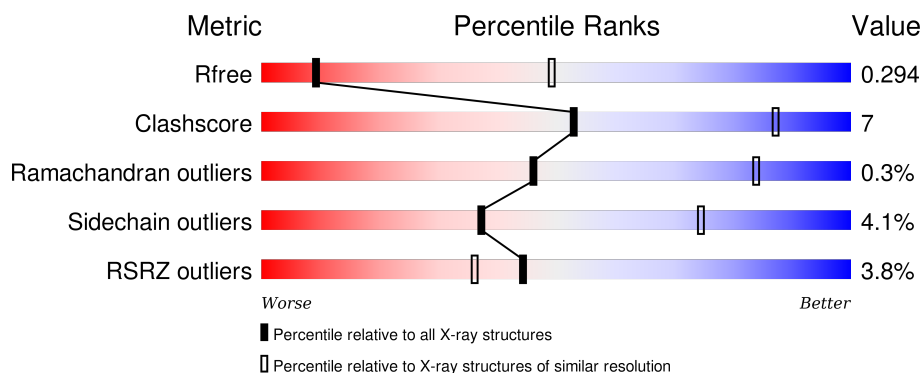
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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))
$R_{free}$	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	 3% 76% 21% ..
2	B	428	 5% 82% 16% ..

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
3	29T	A	601	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7923 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 HIV-1 Reverse Transcriptase, p66 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	544	Total	C	N	O	S	0	0	0
			4391	2842	727	815	7			

There are 6 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	101	PRO	LYS	engineered mutation	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

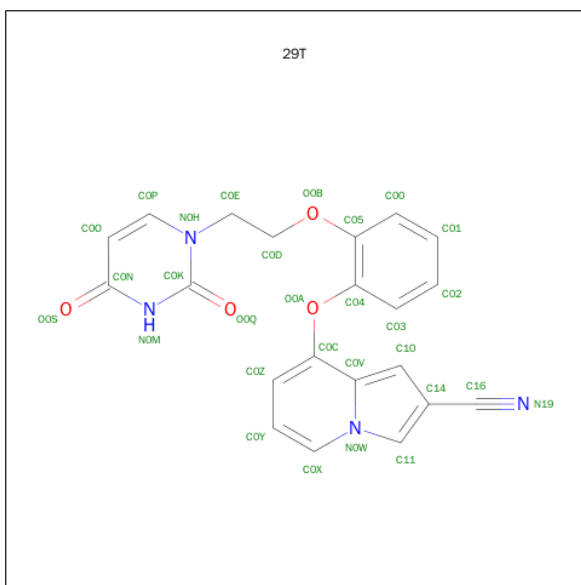
- Molecule 2 is a protein called HIV-1 Reverse Transcriptase, p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	424	Total	C	N	O	S	0	0	0
			3503	2281	582	633	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 8-{2-[2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy]phenoxy}indolizine-2-carbonitrile (three-letter code: 29T) (formula: C<sub>21</sub>H<sub>16</sub>N<sub>4</sub>O<sub>4</sub>).

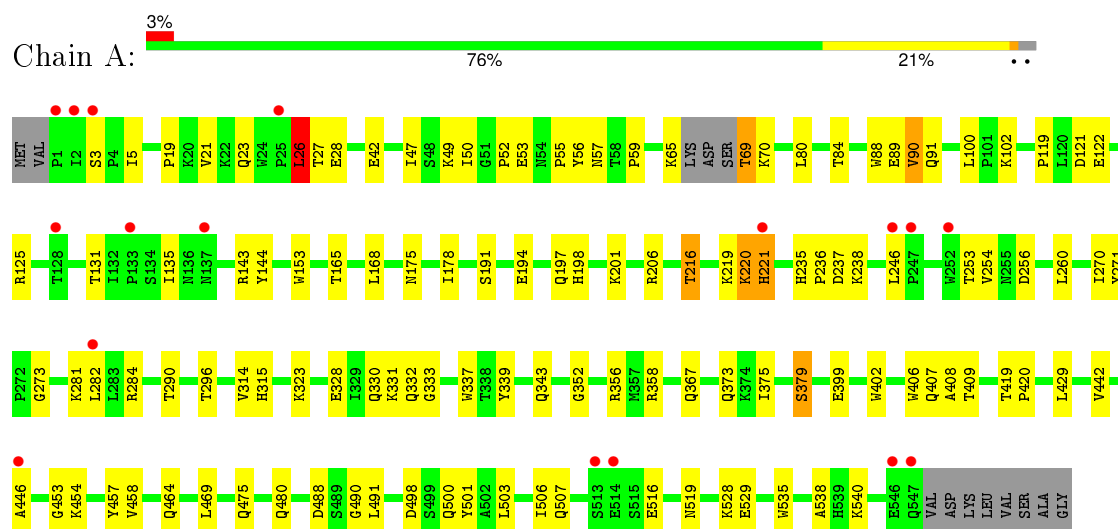


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	21	4	4		

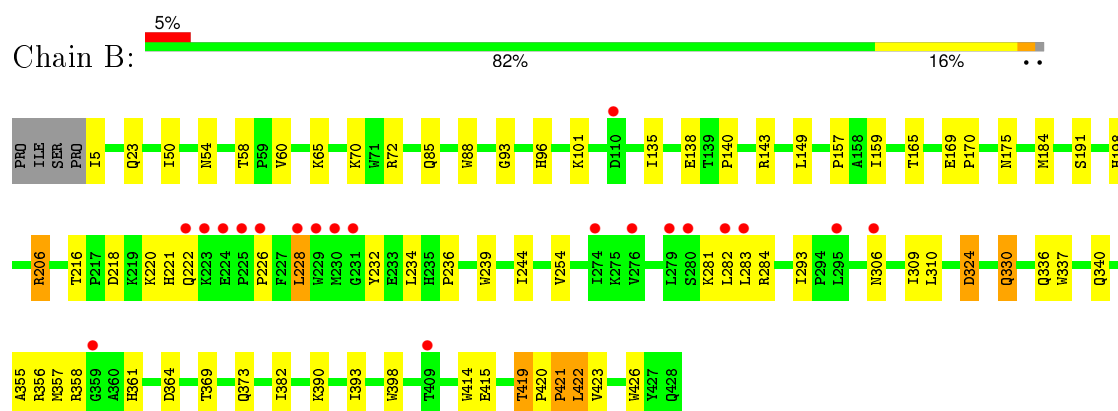
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: HIV-1 Reverse Transcriptase, p66 subunit



- Molecule 2: HIV-1 Reverse Transcriptase, p51 subunit



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	224.20 Å 69.01 Å 104.28 Å 90.00° 106.49° 90.00°	Depositor
Resolution (Å)	38.22 – 3.50 38.22 – 3.23	Depositor EDS
% Data completeness (in resolution range)	99.3 (38.22-3.50) 86.4 (38.22-3.23)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.31 (at 3.25 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.244 , 0.282 0.263 , 0.294	Depositor DCC
$R_{free}$ test set	1638 reflections (9.70%)	DCC
Wilson B-factor (Å <sup>2</sup> )	124.0	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 54.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	1 of 23862 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	7923	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 29T

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	0/4507	0.45	1/6134 (0.0%)
2	B	0.23	0/3606	0.46	0/4901
All	All	0.23	0/8113	0.46	1/11035 (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
2	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	26	LEU	CA-CB-CG	6.48	130.20	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
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	4391	0	4402	68	0
2	B	3503	0	3530	41	0
3	A	29	0	16	1	0
All	All	7923	0	7948	105	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 7.

All (105) 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:281:LYS:HG2	1:A:284:ARG:HE	1.49	0.76
1:A:56:TYR:O	1:A:143:ARG:NH2	2.20	0.74
1:A:23:GLN:NE2	1:A:131:THR:O	2.22	0.72
2:B:157:PRO:HG3	2:B:184:MET:HA	1.75	0.69
2:B:221:HIS:HE1	2:B:228:LEU:HD13	1.58	0.68
1:A:26:LEU:HG	1:A:27:THR:HG23	1.76	0.67
1:A:408:ALA:HB2	2:B:337:TRP:HH2	1.60	0.66
1:A:331:LYS:HE2	1:A:333:GLY:HA2	1.79	0.65
2:B:356:ARG:HE	2:B:357:MET:H	1.46	0.63
1:A:3:SER:HB3	1:A:5:ILE:HG13	1.81	0.63
1:A:100:LEU:HB3	3:A:601:29T:H11	1.81	0.62
2:B:330:GLN:OE1	2:B:340:GLN:NE2	2.33	0.61
2:B:54:ASN:O	2:B:143:ARG:NH2	2.36	0.59
1:A:469:LEU:HD11	1:A:480:GLN:HG2	1.85	0.59
2:B:390:LYS:NZ	2:B:415:GLU:OE2	2.35	0.59
2:B:191:SER:OG	2:B:198:HIS:ND1	2.28	0.58
2:B:206:ARG:HD2	2:B:218:ASP:HB3	1.86	0.57
2:B:206:ARG:NH1	2:B:216:THR:OG1	2.37	0.57
2:B:306:ASN:HA	2:B:309:ILE:HG22	1.86	0.56
2:B:282:LEU:HG	2:B:293:ILE:HD11	1.87	0.56
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.71	0.55
1:A:498:ASP:HB2	1:A:538:ALA:HB2	1.87	0.54
1:A:282:LEU:HD21	1:A:296:THR:HG23	1.88	0.54
1:A:281:LYS:HA	1:A:284:ARG:HG3	1.88	0.54
1:A:50:ILE:O	1:A:143:ARG:HB2	2.07	0.54
2:B:423:VAL:HG12	2:B:426:TRP:CE3	2.42	0.54
1:A:3:SER:HB2	1:A:119:PRO:HD3	1.89	0.53
1:A:490:GLY:O	1:A:528:LYS:NZ	2.32	0.53
1:A:406:TRP:CH2	1:A:407:GLN:HG3	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:VAL:HG12	1:A:457:TYR:HB3	1.91	0.52
2:B:149:LEU:HD21	2:B:159:ILE:HD12	1.91	0.52
1:A:42:GLU:OE1	1:A:49:LYS:NZ	2.42	0.52
2:B:65:LYS:HA	2:B:72:ARG:HD2	1.92	0.51
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.92	0.51
1:A:356:ARG:NE	1:A:358:ARG:O	2.44	0.51
1:A:399:GLU:HA	1:A:402:TRP:NE1	2.25	0.51
2:B:324:ASP:OD1	2:B:324:ASP:N	2.45	0.50
1:A:102:LYS:NZ	1:A:236:PRO:O	2.28	0.49
1:A:253:THR:HG23	1:A:256:ASP:H	1.77	0.49
1:A:52:PRO:HA	1:A:53:GLU:C	2.32	0.49
1:A:328:GLU:HG2	1:A:330:GLN:HE22	1.77	0.49
1:A:337:TRP:HE1	1:A:367:GLN:NE2	2.10	0.49
1:A:69:THR:HG23	1:A:70:LYS:HG3	1.95	0.48
1:A:28:GLU:HG3	1:A:135:ILE:HD13	1.96	0.48
1:A:88:TRP:CH2	2:B:140:PRO:HB2	2.49	0.48
2:B:423:VAL:HG12	2:B:426:TRP:HE3	1.77	0.47
1:A:191:SER:HG	1:A:198:HIS:CE1	2.33	0.47
1:A:19:PRO:HD3	1:A:80:LEU:HD13	1.96	0.47
1:A:406:TRP:CH2	2:B:420:PRO:HD2	2.49	0.47
2:B:281:LYS:O	2:B:284:ARG:HB2	2.15	0.47
1:A:121:ASP:O	1:A:125:ARG:HG3	2.15	0.46
2:B:50:ILE:HG12	2:B:143:ARG:HB3	1.98	0.46
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.97	0.46
2:B:220:LYS:C	2:B:222:GLN:H	2.17	0.46
1:A:429:LEU:HD11	1:A:506:ILE:HG22	1.98	0.46
2:B:96:HIS:NE2	2:B:382:ILE:O	2.41	0.46
2:B:85:GLN:HA	2:B:88:TRP:CE2	2.51	0.45
1:A:446:ALA:HA	1:A:453:GLY:HA3	1.97	0.45
1:A:458:VAL:HG22	1:A:464:GLN:HG2	1.98	0.45
1:A:57:ASN:OD1	1:A:131:THR:OG1	2.25	0.45
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.52	0.45
1:A:191:SER:OG	1:A:198:HIS:ND1	2.49	0.45
2:B:244:ILE:HD12	2:B:310:LEU:HD11	1.98	0.45
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.99	0.45
1:A:337:TRP:HE1	1:A:367:GLN:HE21	1.64	0.45
1:A:122:GLU:OE1	1:A:125:ARG:NH1	2.50	0.44
1:A:194:GLU:HB3	1:A:197:GLN:HG3	1.99	0.44
2:B:88:TRP:O	2:B:93:GLY:HA2	2.18	0.44
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.52	0.44
2:B:393:ILE:HD13	2:B:398:TRP:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.53	0.44
1:A:457:TYR:OH	1:A:488:ASP:OD2	2.25	0.44
1:A:253:THR:OG1	1:A:254:VAL:N	2.50	0.44
1:A:55:PRO:HB2	1:A:143:ARG:NH2	2.33	0.43
2:B:369:THR:HG22	2:B:373:GLN:HE21	1.83	0.43
2:B:422:LEU:O	2:B:422:LEU:HD23	2.17	0.43
1:A:201:LYS:HD3	1:A:201:LYS:HA	1.86	0.43
1:A:90:VAL:HG23	1:A:91:GLN:H	1.82	0.43
1:A:246:LEU:HD22	1:A:260:LEU:HD12	2.00	0.43
1:A:323:LYS:HB2	1:A:343:GLN:NE2	2.34	0.43
2:B:336:GLN:HE21	2:B:358:ARG:NH2	2.17	0.43
1:A:220:LYS:HB3	1:A:221:HIS:H	1.68	0.42
1:A:271:TYR:CD2	1:A:314:VAL:HG12	2.54	0.42
1:A:375:ILE:O	1:A:379:SER:HB2	2.20	0.42
1:A:270:ILE:HG23	1:A:271:TYR:H	1.83	0.42
1:A:540:LYS:HA	1:A:540:LYS:HD3	1.83	0.42
2:B:169:GLU:HB3	2:B:170:PRO:HD3	2.00	0.42
1:A:408:ALA:HB1	2:B:364:ASP:HB3	2.02	0.42
2:B:220:LYS:HB3	2:B:221:HIS:H	1.73	0.42
1:A:131:THR:HG23	1:A:143:ARG:HD2	2.01	0.41
1:A:235:HIS:HB2	1:A:238:LYS:O	2.20	0.41
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.62	0.41
1:A:19:PRO:O	1:A:56:TYR:HB3	2.21	0.41
2:B:101:LYS:HD3	2:B:382:ILE:HG23	2.03	0.41
1:A:516:GLU:HA	1:A:519:ASN:HD22	1.86	0.41
1:A:491:LEU:HB3	1:A:529:GLU:HG3	2.02	0.41
1:A:206:ARG:HE	1:A:216:THR:HB	1.85	0.41
2:B:135:ILE:O	2:B:138:GLU:HG3	2.19	0.41
1:A:238:LYS:HD2	1:A:315:HIS:NE2	2.36	0.41
1:A:47:ILE:HD12	1:A:144:TYR:CD1	2.56	0.41
1:A:273:GLY:HA2	1:A:332:GLN:NE2	2.36	0.41
1:A:419:THR:HA	1:A:420:PRO:HD3	1.84	0.41
2:B:23:GLN:HE22	2:B:60:VAL:H	1.68	0.41
2:B:336:GLN:HG2	2:B:355:ALA:HB2	2.03	0.40
2:B:356:ARG:HA	2:B:356:ARG:HD2	1.82	0.40

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	540/557 (97%)	513 (95%)	27 (5%)	0	100	100
2	B	422/428 (99%)	394 (93%)	25 (6%)	3 (1%)	26	72
All	All	962/985 (98%)	907 (94%)	52 (5%)	3 (0%)	46	84

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	422	LEU
2	B	226	PRO
2	B	421	PRO

### 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	477/495 (96%)	456 (96%)	21 (4%)	35	73
2	B	384/390 (98%)	370 (96%)	14 (4%)	42	77
All	All	861/885 (97%)	826 (96%)	35 (4%)	37	74

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	65	LYS
1	A	69	THR

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Mol	Chain	Res	Type
1	A	89	GLU
1	A	90	VAL
1	A	165	THR
1	A	168	LEU
1	A	175	ASN
1	A	178	ILE
1	A	216	THR
1	A	219	LYS
1	A	220	LYS
1	A	221	HIS
1	A	237	ASP
1	A	290	THR
1	A	373	GLN
1	A	379	SER
1	A	409	THR
1	A	454	LYS
1	A	500	GLN
1	A	507	GLN
2	B	5	ILE
2	B	58	THR
2	B	70	LYS
2	B	165	THR
2	B	175	ASN
2	B	206	ARG
2	B	228	LEU
2	B	232	TYR
2	B	234	LEU
2	B	324	ASP
2	B	330	GLN
2	B	361	HIS
2	B	414	TRP
2	B	419	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	367	GLN
1	A	394	GLN
1	A	428	GLN
1	A	519	ASN
1	A	520	GLN

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Mol	Chain	Res	Type
2	B	23	GLN
2	B	161	GLN
2	B	278	GLN
2	B	330	GLN
2	B	340	GLN
2	B	373	GLN

### 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](#)

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	29T	A	601	-	26,32,32	1.40	3 (11%)	29,44,44	2.89	4 (13%)

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	29T	A	601	-	-	0/10/12/12	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	29T	C0Z-C0C	2.36	1.43	1.38
3	A	601	29T	C0Y-C0Z	3.30	1.45	1.38
3	A	601	29T	C0P-N0H	3.90	1.42	1.36

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	29T	C10-C14-C16	-3.62	121.87	126.72
3	A	601	29T	C0O-C0N-N0M	-3.52	114.63	123.28
3	A	601	29T	C0D-C0E-N0H	3.62	115.91	110.94
3	A	601	29T	C0N-N0M-C0K	13.34	128.27	114.21

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
3	A	601	29T	1	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	544/557 (97%)	-0.17	17 (3%)	52 43	98, 143, 187, 263	1 (0%)
2	B	424/428 (99%)	-0.06	20 (4%)	35 28	91, 132, 216, 290	0
All	All	968/985 (98%)	-0.12	37 (3%)	44 36	91, 139, 208, 290	1 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	229	TRP	6.8
2	B	225	PRO	5.7
1	A	133	PRO	5.3
2	B	409	THR	4.5
2	B	226	PRO	4.2
1	A	514	GLU	4.2
2	B	279	LEU	3.6
2	B	222	GLN	3.3
1	A	246	LEU	3.2
2	B	110	ASP	3.1
2	B	359	GLY	3.0
2	B	282	LEU	3.0
1	A	546	GLU	2.9
2	B	224	GLU	2.9
1	A	137	ASN	2.9
2	B	283	LEU	2.8
1	A	247	PRO	2.8
2	B	231	GLY	2.8
2	B	223	LYS	2.7
2	B	228	LEU	2.7
2	B	274	ILE	2.7
2	B	306	ASN	2.6
1	A	252	TRP	2.5
2	B	276	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	547	GLN	2.5
1	A	221	HIS	2.5
2	B	280	SER	2.4
1	A	446	ALA	2.3
2	B	230	MET	2.3
1	A	2	ILE	2.2
1	A	513	SER	2.2
1	A	1	PRO	2.2
2	B	295	LEU	2.1
1	A	128	THR	2.0
1	A	282	LEU	2.0
1	A	3	SER	2.0
1	A	25	PRO	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(Å <sup>2</sup> )	Q<0.9
3	29T	A	601	29/29	0.88	0.44	2.35	112,132,151,153	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.