



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:26 AM UTC

PDB ID : 2RKU / pdb_00002rku
Title : Structure of PLK1 in complex with BI2536
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Deposited on : 2007-10-17
Resolution : 1.95 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

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>.

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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

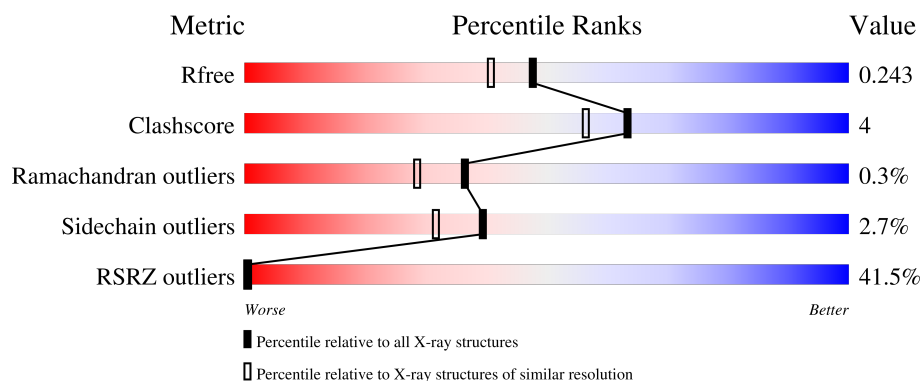
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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 ≥ 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	294	41% 88% 11%

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
5	SRT	A	505	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	SRT	A	507	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2785 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 Serine/threonine-protein kinase PLK1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	5	0
			2411	1546	428	425	12			

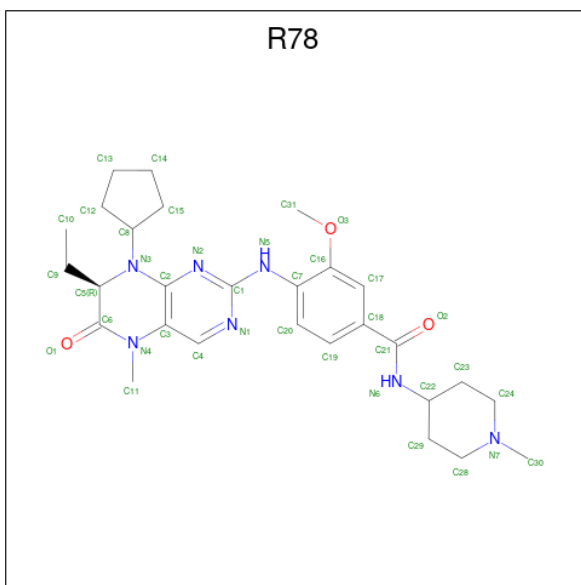
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	210	VAL	THR	engineered mutation	UNP P53350

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

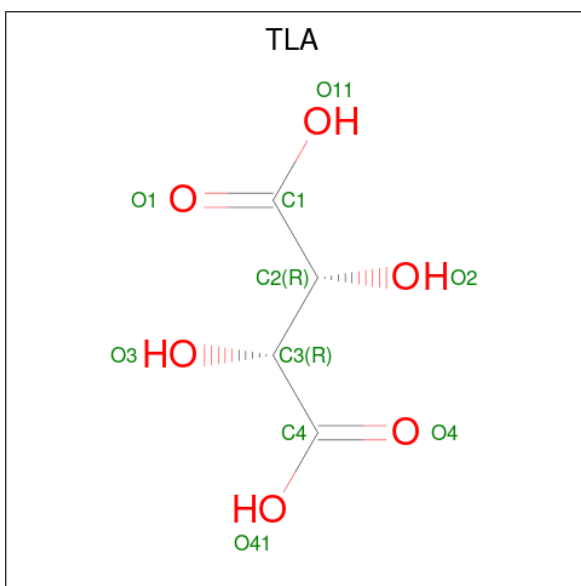
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 4-[[[(7R)-8-cyclopentyl-7-ethyl-5-methyl-6-oxo-5,6,7,8-tetrahydropteridin-2-yl]amino}-3-methoxy-N-(1-methylpiperidin-4-yl)benzamide (CCD ID: R78) (formula: C₂₈H₃₉N₇O₃).



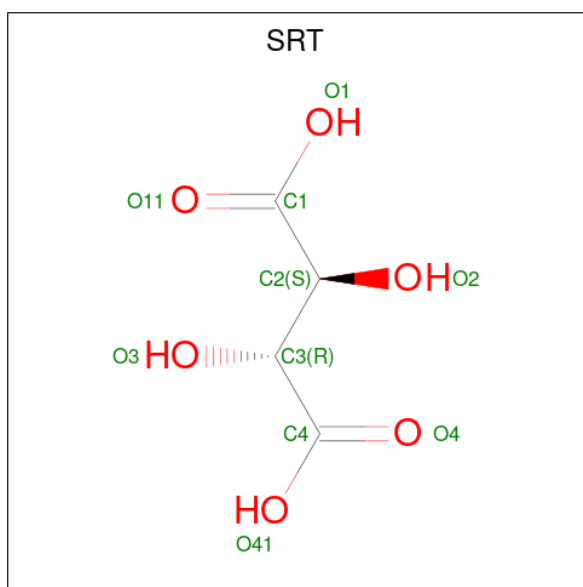
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			38	28	7	3		

- Molecule 4 is L(+)-TARTARIC ACID (CCD ID: TLA) (formula: $C_4H_6O_6$).



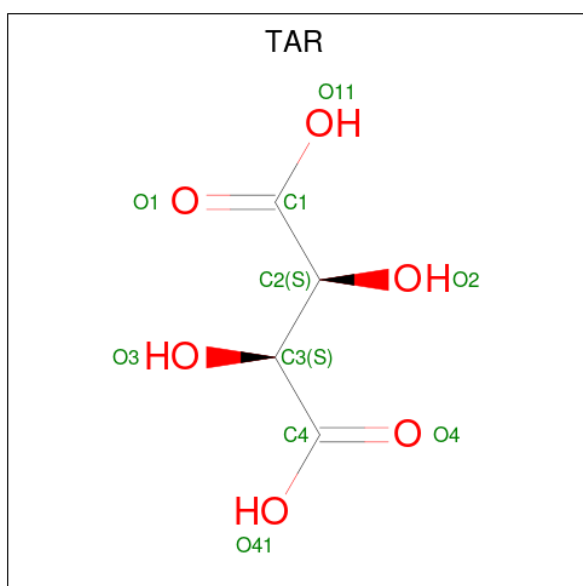
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	4	6		
4	A	1	Total	C	O	0	0
			10	4	6		

- Molecule 5 is S,R MESO-TARTARIC ACID (CCD ID: SRT) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	4	6		
5	A	1	Total	C	O	0	0
			10	4	6		

- Molecule 6 is D(-)-TARTARIC ACID (CCD ID: TAR) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	4	6		

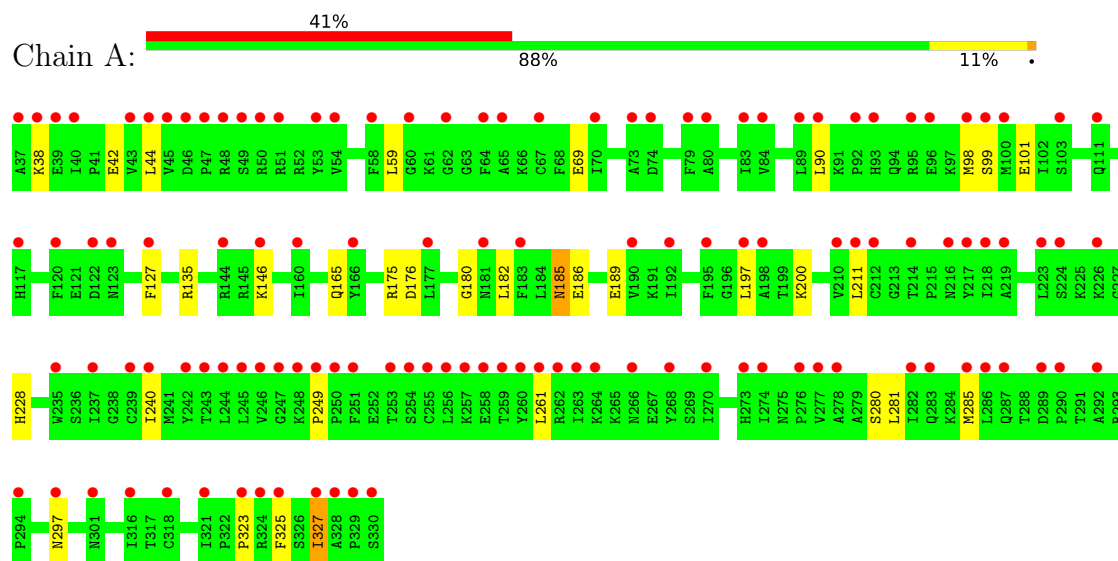
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	285	Total 285	O 285	0	0

3 Residue-property plots

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: Serine/threonine-protein kinase PLK1



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	66.59Å 66.59Å 154.11Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.41 – 1.95 33.41 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.4 (33.41-1.95) 99.4 (33.41-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.241 0.205 , 0.243	Depositor DCC
R_{free} test set	1494 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2785	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: R78, SRT, TLA, TAR, ZN

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	0.77	1/2466 (0.0%)	0.88	1/3327 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	GLU	CD-OE1	5.67	1.36	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	PRO	O-C-N	5.75	123.85	121.15

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	2411	0	2465	18	1
2	A	1	0	0	0	0
3	A	38	0	39	1	0
4	A	20	0	8	0	0
5	A	20	0	8	2	0
6	A	10	0	4	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	285	0	0	3	0
All	All	2785	0	2524	20	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 4.

All (20) 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:176:ASP:OD2	7:A:513:HOH:O	2.03	0.74
1:A:281:LEU:HG	1:A:285:MET:HE2	1.81	0.63
1:A:186:GLU:HB2	7:A:664:HOH:O	2.05	0.55
1:A:175:ARG:HH11	1:A:228:HIS:HE1	1.55	0.53
1:A:180:GLY:N	6:A:503:TAR:O4	2.41	0.53
3:A:500:R78:H2O	3:A:500:R78:N2	2.25	0.51
1:A:185:ASN:ND2	1:A:189:GLU:H	2.09	0.50
1:A:185:ASN:HD22	1:A:185:ASN:C	2.20	0.48
1:A:59:LEU:HD21	1:A:69:GLU:HG3	1.99	0.45
1:A:182:LEU:HD11	1:A:240:ILE:HD13	2.00	0.44
1:A:101:GLU:OE2	7:A:704:HOH:O	2.21	0.44
1:A:165:GLN:HG2	1:A:297:ASN:HD22	1.83	0.44
1:A:175:ARG:HH11	1:A:228:HIS:CE1	2.35	0.43
1:A:146:LYS:HA	1:A:146:LYS:HD3	1.81	0.42
1:A:197:LEU:HD13	1:A:211:LEU:HD21	2.02	0.42
1:A:165:GLN:HG2	1:A:297:ASN:ND2	2.35	0.41
1:A:44:LEU:HD22	1:A:127:PHE:CE2	2.55	0.41
1:A:135:ARG:NH2	5:A:507:SRT:O4	2.51	0.41
1:A:323:PRO:HG2	1:A:325:PHE:CE1	2.55	0.41
5:A:507:SRT:O11	5:A:507:SRT:O3	2.33	0.41

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)
1:A:99[B]:SER:OG	1:A:327:ILE:O[2_564]	2.09	0.11

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	297/294 (101%)	286 (96%)	10 (3%)	1 (0%)	36	28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	LYS

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	268/263 (102%)	261 (97%)	7 (3%)	40	33

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	A	98	MET
1	A	185	ASN
1	A	200	LYS
1	A	261	LEU
1	A	280	SER
1	A	327	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	ASN
1	A	165	GLN
1	A	185	ASN
1	A	228	HIS
1	A	266	ASN
1	A	297	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z > 2$
3	R78	A	500	-	40,42,42	0.93	2 (5%)	50,60,60	2.18	12 (24%)
5	SRT	A	505	-	9,9,9	1.17	0	12,12,12	0.95	0
4	TLA	A	502	2	9,9,9	1.08	0	12,12,12	1.21	2 (16%)
6	TAR	A	503	-	9,9,9	1.11	0	12,12,12	1.38	1 (8%)
4	TLA	A	504	-	9,9,9	1.14	0	12,12,12	1.04	1 (8%)
5	SRT	A	507	-	9,9,9	1.22	0	12,12,12	1.13	1 (8%)

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
5	SRT	A	507	-	2/2/4/4	12/12/12/12	-
3	R78	A	500	-	-	0/20/57/57	0/5/5/5
5	SRT	A	505	-	2/2/4/4	5/12/12/12	-
4	TLA	A	502	2	-	0/12/12/12	-
6	TAR	A	503	-	-	10/12/12/12	-
4	TLA	A	504	-	-	0/12/12/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	R78	O1-C6	2.13	1.26	1.22
3	A	500	R78	C5-C6	-2.04	1.48	1.51

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	R78	N1-C1-N2	-6.21	120.42	126.42
3	A	500	R78	C11-N4-C6	4.88	122.32	117.69
3	A	500	R78	C31-O3-C16	4.78	124.53	117.51
3	A	500	R78	C4-N1-C1	4.48	121.96	115.81
3	A	500	R78	C9-C5-C6	-4.11	101.46	111.20
3	A	500	R78	O3-C16-C7	3.11	118.69	114.81
3	A	500	R78	C5-C6-N4	3.11	121.27	116.28
6	A	503	TAR	O41-C4-C3	2.92	121.42	113.31
3	A	500	R78	C10-C9-C5	2.85	117.88	112.52
3	A	500	R78	C23-C22-N6	-2.74	105.10	110.57
3	A	500	R78	C29-C22-N6	-2.53	105.52	110.57
3	A	500	R78	C20-C19-C18	-2.52	118.11	120.80
4	A	502	TLA	O11-C1-C2	2.44	120.08	113.31
3	A	500	R78	C30-N7-C24	-2.20	106.44	110.63
5	A	507	SRT	C3-C2-C1	2.16	114.62	109.82
4	A	502	TLA	O41-C4-C3	2.06	119.03	113.31
4	A	504	TLA	O11-C1-C2	2.05	119.01	113.31

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	505	SRT	C3

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Mol	Chain	Res	Type	Atom
5	A	505	SRT	C2
5	A	507	SRT	C3
5	A	507	SRT	C2

All (27) torsion outliers are listed below:

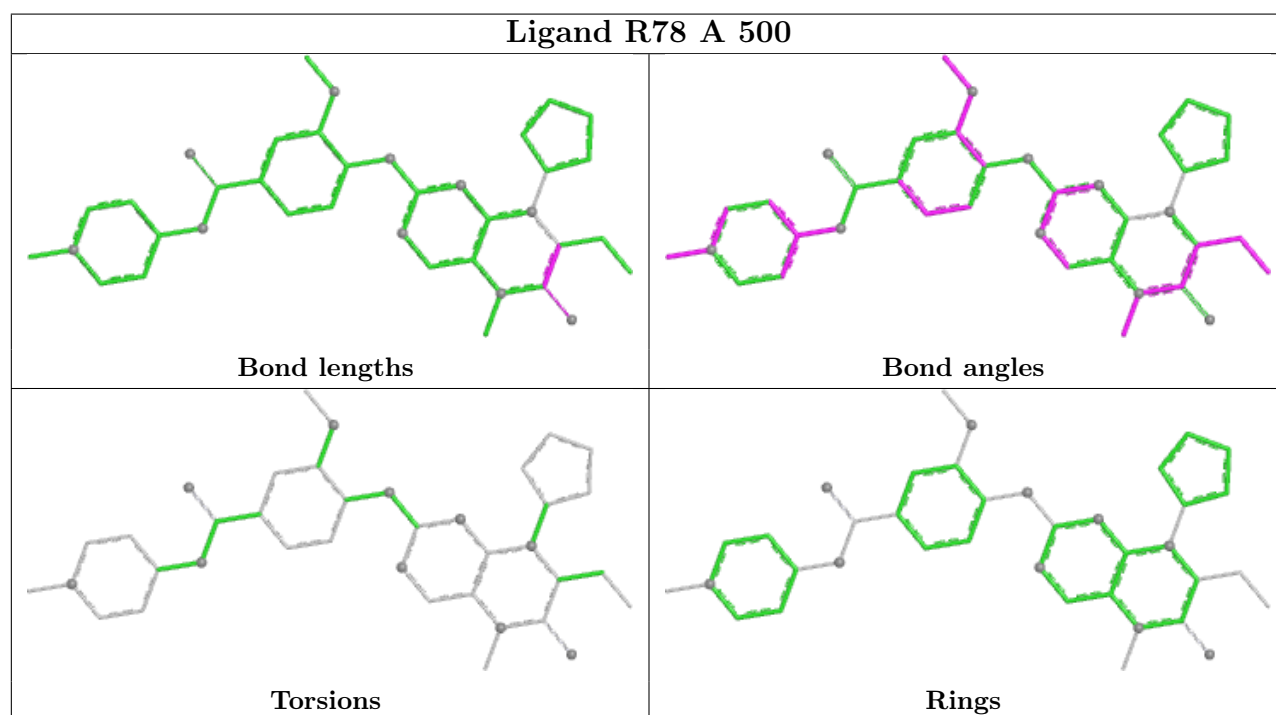
Mol	Chain	Res	Type	Atoms
5	A	505	SRT	O1-C1-C2-O2
5	A	505	SRT	O11-C1-C2-O2
5	A	507	SRT	O1-C1-C2-C3
5	A	507	SRT	O2-C2-C3-O3
6	A	503	TAR	O1-C1-C2-O2
6	A	503	TAR	O11-C1-C2-O2
6	A	503	TAR	C1-C2-C3-O3
5	A	507	SRT	O2-C2-C3-C4
6	A	503	TAR	O2-C2-C3-O3
5	A	507	SRT	C1-C2-C3-C4
5	A	505	SRT	O1-C1-C2-C3
5	A	505	SRT	O11-C1-C2-C3
5	A	507	SRT	O11-C1-C2-C3
5	A	507	SRT	C2-C3-C4-O4
5	A	507	SRT	C2-C3-C4-O41
6	A	503	TAR	O2-C2-C3-C4
6	A	503	TAR	C1-C2-C3-C4
5	A	507	SRT	O3-C3-C4-O4
5	A	507	SRT	O3-C3-C4-O41
5	A	507	SRT	C1-C2-C3-O3
5	A	507	SRT	O1-C1-C2-O2
5	A	507	SRT	O11-C1-C2-O2
6	A	503	TAR	C2-C3-C4-O41
6	A	503	TAR	C2-C3-C4-O4
5	A	505	SRT	O2-C2-C3-O3
6	A	503	TAR	O3-C3-C4-O4
6	A	503	TAR	O3-C3-C4-O41

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	R78	1	0
6	A	503	TAR	1	0
5	A	507	SRT	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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, 95th 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 ²)	Q<0.9
1	A	294/294 (100%)	1.93	122 (41%) 0 0	18, 45, 56, 65	5 (1%)

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	PRO	6.0
1	A	48	ARG	5.7
1	A	47	PRO	4.9
1	A	217	TYR	4.7
1	A	45	VAL	4.6
1	A	256	LEU	4.3
1	A	64	PHE	4.3
1	A	244	LEU	4.0
1	A	100	MET	3.8
1	A	99[A]	SER	3.7
1	A	40	ILE	3.7
1	A	327	ILE	3.7
1	A	263	ILE	3.7
1	A	44	LEU	3.7
1	A	328	ALA	3.6
1	A	324	ARG	3.6
1	A	255	CYS	3.6
1	A	245	LEU	3.6
1	A	254[A]	SER	3.5
1	A	243	THR	3.5
1	A	259	THR	3.5
1	A	98	MET	3.5
1	A	211	LEU	3.5
1	A	49	SER	3.4
1	A	292	ALA	3.4
1	A	246	VAL	3.4
1	A	251	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	286	LEU	3.3
1	A	260	TYR	3.3
1	A	123	ASN	3.3
1	A	318[A]	CYS	3.3
1	A	240	ILE	3.2
1	A	250	PRO	3.1
1	A	43	VAL	3.0
1	A	277	VAL	3.0
1	A	237	ILE	3.0
1	A	330	SER	3.0
1	A	261	LEU	2.9
1	A	242	TYR	2.9
1	A	226	LYS	2.9
1	A	247	GLY	2.9
1	A	79	PHE	2.9
1	A	325	PHE	2.9
1	A	274	ILE	2.9
1	A	253	THR	2.8
1	A	53	TYR	2.8
1	A	92	PRO	2.8
1	A	290	PRO	2.8
1	A	74	ASP	2.8
1	A	224	SER	2.8
1	A	117	HIS	2.8
1	A	127	PHE	2.8
1	A	50	ARG	2.8
1	A	111	GLN	2.8
1	A	212	CYS	2.7
1	A	216	ASN	2.7
1	A	266	ASN	2.7
1	A	58	PHE	2.7
1	A	278	ALA	2.6
1	A	46	ASP	2.6
1	A	89	LEU	2.6
1	A	297	ASN	2.6
1	A	80	ALA	2.6
1	A	210	VAL	2.6
1	A	257	LYS	2.6
1	A	197	LEU	2.6
1	A	37	ALA	2.6
1	A	103	SER	2.6
1	A	93	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	268	TYR	2.5
1	A	177	LEU	2.5
1	A	39	GLU	2.5
1	A	223	LEU	2.5
1	A	84	VAL	2.5
1	A	120	PHE	2.5
1	A	62	GLY	2.4
1	A	65	ALA	2.4
1	A	276	PRO	2.4
1	A	301	ASN	2.4
1	A	248	LYS	2.4
1	A	144	ARG	2.4
1	A	90	LEU	2.4
1	A	270	ILE	2.4
1	A	321	ILE	2.4
1	A	258	GLU	2.3
1	A	122	ASP	2.3
1	A	214	THR	2.3
1	A	264	LYS	2.3
1	A	239	CYS	2.3
1	A	235	TRP	2.3
1	A	38	LYS	2.3
1	A	183	PHE	2.3
1	A	323	PRO	2.2
1	A	83	ILE	2.2
1	A	146	LYS	2.2
1	A	96	GLU	2.2
1	A	219	ALA	2.2
1	A	218	ILE	2.2
1	A	316	ILE	2.2
1	A	181[A]	ASN	2.2
1	A	70	ILE	2.2
1	A	249	PRO	2.2
1	A	95	ARG	2.2
1	A	262	ARG	2.2
1	A	289	ASP	2.1
1	A	54	VAL	2.1
1	A	198	ALA	2.1
1	A	190	VAL	2.1
1	A	67	CYS	2.1
1	A	273	HIS	2.1
1	A	73	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	283	GLN	2.1
1	A	195	PHE	2.1
1	A	166	TYR	2.1
1	A	282	ILE	2.1
1	A	285	MET	2.1
1	A	287	GLN	2.1
1	A	192	ILE	2.0
1	A	294	PRO	2.0
1	A	51	ARG	2.0
1	A	160	ILE	2.0
1	A	60	GLY	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

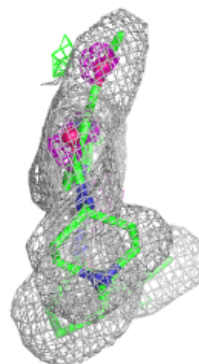
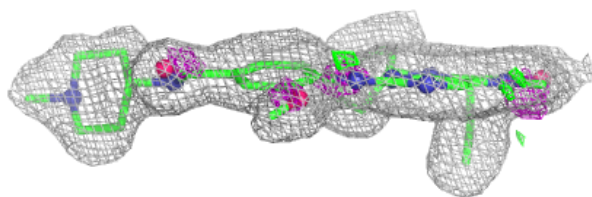
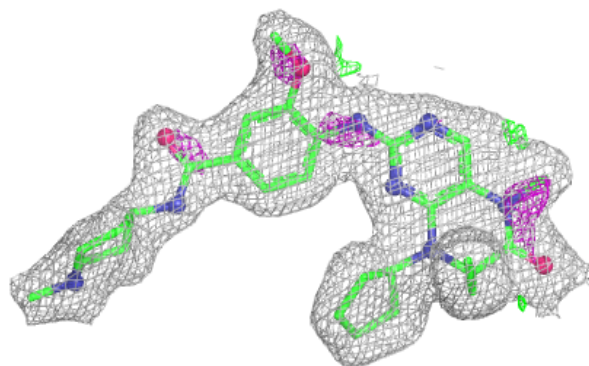
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	TLA	A	502	10/10	0.59	0.31	35,39,43,44	0
4	TLA	A	504	10/10	0.64	0.22	56,57,58,58	0
6	TAR	A	503	10/10	0.65	0.17	56,65,68,69	0
5	SRT	A	505	10/10	0.70	0.22	86,87,88,88	0
5	SRT	A	507	10/10	0.80	0.19	78,81,82,82	0
2	ZN	A	501	1/1	0.81	0.32	34,34,34,34	0
3	R78	A	500	38/38	0.91	0.10	24,28,47,49	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around R78 A 500:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.