



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:28 PM UTC

PDB ID : 1O5C / pdb_00001o5c
Title : Dissecting and Designing Inhibitor Selectivity Determinants at the S1 site
Using an Artificial Ala190 Protease (Ala190 uPA)
Authors : Katz, B.A.; Luong, C.; Ho, J.D.; Somoza, J.R.; Gjerstad, E.; Tang, J.;
Williams, S.R.; Verner, E.; Mackman, R.L.; Young, W.B.; Sprengeler, P.A.;
Chan, H.; Mortara, K.; Janc, J.W.; McGrath, M.E.
Deposited on : 2003-09-09
Resolution : 1.63 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](#) ①) 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

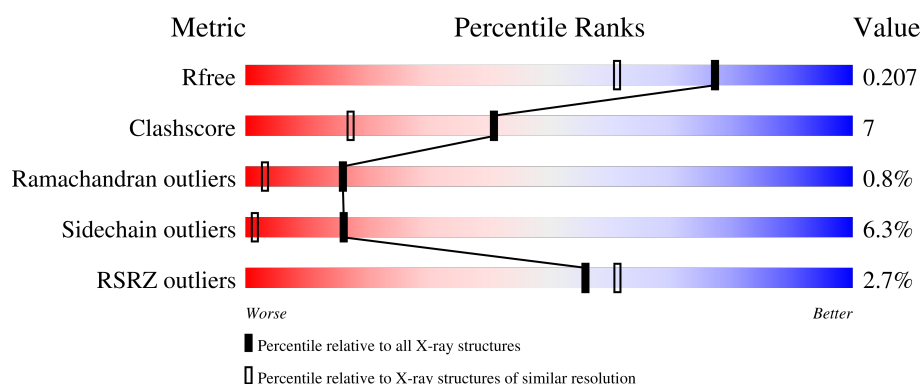
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	23	9% 22% 17% 61%
2	B	253	2% 64% 28% . .

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	CR9	B	251	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4786 atoms, of which 2479 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 Urokinase-type plasminogen activator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	9	Total	C	H	N	O	S	0	0	0
			147	46	75	13	12	1			

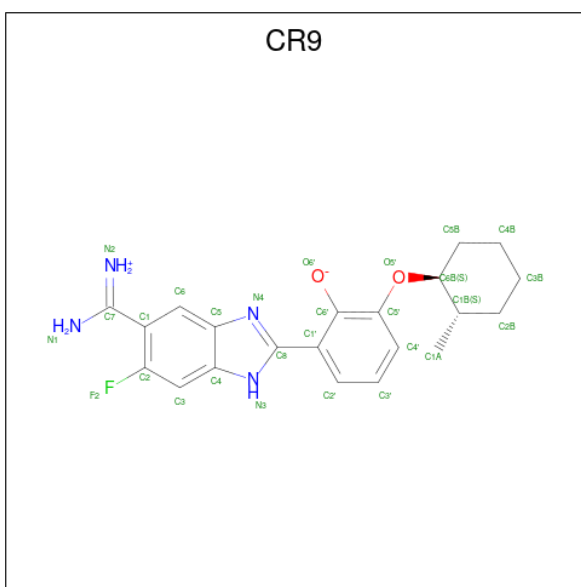
- Molecule 2 is a protein called Urokinase-type plasminogen activator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	246	Total	C	H	N	O	S	0	11	0
			3934	1252	1959	340	365	18			

There are 2 discrepancies between the modelled and reference sequences:

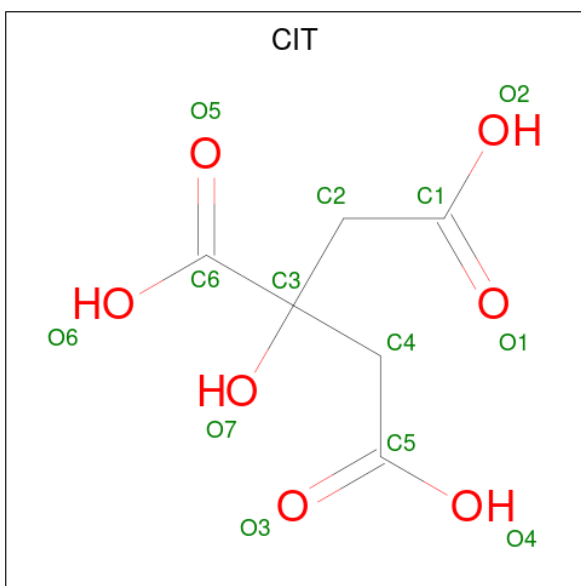
Chain	Residue	Modelled	Actual	Comment	Reference
B	145	ALA	ASN	engineered mutation	UNP P00749
B	190	ALA	SER	engineered mutation	UNP P00749

- Molecule 3 is 2-{5-[AMINO(IMINIO)METHYL]-6-FLUORO-1H-BENZIMIDAZOL-2-YL}-6-[(2-METHYLCYCLOHEXYL)OXY]BENZENOLATE (CCD ID: CR9) (formula: C₂₁H₂₃FN₄O₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	F	H	N	O	0	0
			51	21	1	23	4	2		

- Molecule 4 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			18	6	5	7		
4	B	1	Total	C	H	O	0	0
			18	6	5	7		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	8	Total 24	H 16	O 8	0	0
5	B	197	Total 594	H 396	O 198	0	1

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	81.71Å 50.28Å 66.20Å 90.00° 113.04° 90.00°	Depositor
Resolution (Å)	7.00 – 1.63 7.00 – 1.63	Depositor EDS
% Data completeness (in resolution range)	78.6 (7.00-1.63) 77.6 (7.00-1.63)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.61 (at 1.63Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.191 , 0.219 (Not available) , 0.207	Depositor DCC
R_{free} test set	2410 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 73.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4786	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.57% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CR9, CIT

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	1.65	0/72	1.84	2/93 (2.2%)
2	B	1.79	22/2061 (1.1%)	1.75	60/2792 (2.1%)
All	All	1.79	22/2133 (1.0%)	1.76	62/2885 (2.1%)

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	8

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	241	HIS	CG-ND1	-7.68	1.29	1.38
2	B	170(B)	HIS	CG-ND1	-6.79	1.30	1.38
2	B	37	HIS	CG-ND1	-6.72	1.30	1.38
2	B	37	HIS	CD2-NE2	-6.38	1.30	1.37
2	B	57	HIS	CG-ND1	-6.19	1.31	1.38
2	B	170(B)	HIS	CD2-NE2	-6.06	1.31	1.37
2	B	165	HIS	CG-ND1	-5.99	1.31	1.38
2	B	100	HIS	CG-ND1	-5.86	1.31	1.38
2	B	91	HIS	CG-ND1	-5.81	1.31	1.38
2	B	99	HIS	CD2-NE2	-5.79	1.31	1.37
2	B	99	HIS	CG-ND1	-5.73	1.31	1.38
2	B	165	HIS	CG-CD2	5.55	1.42	1.35
2	B	91	HIS	CD2-NE2	-5.53	1.31	1.37
2	B	57	HIS	CD2-NE2	-5.47	1.31	1.37
2	B	165	HIS	CD2-NE2	-5.46	1.31	1.37
2	B	233	HIS	CG-CD2	5.34	1.41	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	233	HIS	CG-ND1	-5.33	1.32	1.38
2	B	100	HIS	CD2-NE2	-5.12	1.32	1.37
2	B	51	TRP	CG-CD2	-5.06	1.34	1.43
2	B	233	HIS	CD2-NE2	-5.04	1.32	1.37
2	B	100	HIS	CG-CD2	5.03	1.41	1.35
2	B	241	HIS	CD2-NE2	-5.00	1.32	1.37

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	207	MET	N-CA-C	-9.55	94.64	109.25
2	B	37(A)	ARG	N-CA-C	-8.79	94.86	109.46
2	B	199	LEU	N-CA-C	-8.54	93.07	108.02
2	B	92	LYS	N-CA-C	-8.52	96.22	109.76
2	B	21	PHE	N-CA-C	-7.99	97.62	110.32
1	A	16	LYS	N-CA-C	-7.66	97.84	110.17
2	B	100	HIS	CA-CB-CG	-7.51	106.29	113.80
2	B	185	ASP	N-CA-C	-7.25	100.66	109.72
2	B	26	ASN	CA-CB-CG	-7.19	105.41	112.60
2	B	133	GLY	N-CA-C	-7.03	106.32	114.69
2	B	84[A]	GLU	N-CA-C	-7.01	99.56	110.42
2	B	84[B]	GLU	N-CA-C	-7.01	99.56	110.42
2	B	171	TYR	N-CA-C	-6.98	97.02	108.12
2	B	190	ALA	N-CA-C	-6.95	100.51	110.59
2	B	226	GLY	N-CA-C	-6.90	100.18	112.54
2	B	73	LEU	N-CA-C	6.85	118.53	111.14
2	B	51	TRP	NE1-CE2-CZ2	6.75	140.22	130.10
2	B	57	HIS	CA-CB-CG	6.51	120.31	113.80
2	B	125	SER	N-CA-C	-6.45	100.97	110.52
2	B	170	GLN	N-CA-C	-6.43	101.66	109.83
2	B	107	LYS	N-CA-C	-6.41	98.80	109.24
2	B	48	SER	N-CA-C	-6.39	102.38	108.07
2	B	237	TRP	NE1-CE2-CZ2	6.39	139.68	130.10
2	B	29	TRP	NE1-CE2-CZ2	6.24	139.46	130.10
2	B	193	GLY	N-CA-C	-6.18	107.34	114.69
2	B	202	SER	N-CA-C	-6.16	99.61	109.96
2	B	229	THR	N-CA-C	-6.13	101.21	110.28
2	B	99	HIS	N-CA-C	-6.06	98.52	108.76
2	B	37	HIS	CA-CB-CG	-5.93	107.87	113.80
2	B	186	TRP	CG-CD1-NE1	-5.93	102.49	110.20
1	A	10	LYS	N-CA-C	-5.88	100.38	109.14
2	B	123	LEU	N-CA-C	-5.87	102.76	110.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	165	HIS	N-CA-C	-5.87	104.96	111.36
2	B	68	LEU	N-CA-C	-5.79	100.46	109.72
2	B	94	TYR	N-CA-C	-5.77	101.74	110.28
2	B	29	TRP	CG-CD1-NE1	-5.65	102.86	110.20
2	B	35	ARG	CD-NE-CZ	-5.62	116.53	124.40
2	B	38	VAL	N-CA-C	-5.58	100.31	108.11
2	B	121	ILE	N-CA-C	-5.54	100.62	109.20
2	B	43	GLY	N-CA-C	-5.50	103.11	112.51
2	B	158	THR	CA-CB-OG1	-5.50	101.36	109.60
2	B	189	ASP	CA-CB-CG	5.49	118.09	112.60
2	B	237	TRP	CG-CD1-NE1	-5.48	103.07	110.20
2	B	75	SER	N-CA-C	-5.48	101.20	109.85
2	B	186	TRP	NE1-CE2-CZ2	5.48	138.31	130.10
2	B	215	TRP	NE1-CE2-CZ2	5.45	138.27	130.10
2	B	51	TRP	NE1-CE2-CD2	-5.41	100.36	107.40
2	B	51	TRP	CG-CD1-NE1	-5.41	103.17	110.20
2	B	17	ILE	N-CA-C	-5.39	100.66	108.36
2	B	237	TRP	NE1-CE2-CD2	-5.33	100.47	107.40
2	B	234	PHE	CA-CB-CG	-5.30	108.50	113.80
2	B	98	ALA	N-CA-C	5.30	117.71	110.55
2	B	160	VAL	N-CA-CB	-5.21	102.72	112.36
2	B	51	TRP	CG-CD2-CE3	-5.20	128.70	133.90
2	B	74	ASN	OD1-CG-ND2	-5.20	117.40	122.60
2	B	186	TRP	CD1-NE1-CE2	5.15	118.17	108.90
2	B	153	GLU	N-CA-C	-5.15	106.16	112.90
2	B	215	TRP	CG-CD1-NE1	-5.10	103.58	110.20
2	B	241	HIS	CE1-NE2-CD2	-5.08	103.92	109.00
2	B	170	GLN	OE1-CD-NE2	-5.05	117.55	122.60
2	B	51	TRP	N-CA-C	5.03	117.11	108.90
2	B	231	VAL	N-CA-C	5.01	116.33	110.62

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	109	ARG	Sidechain
2	B	110(D)	ARG	Sidechain
2	B	116	ARG	Sidechain
2	B	206	ARG	Sidechain
2	B	217	ARG	Sidechain
2	B	230	ARG	Sidechain
2	B	239	ARG	Sidechain

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Mol	Chain	Res	Type	Group
2	B	35	ARG	Sidechain

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	72	75	75	1	0
2	B	1975	1959	1928	22	1
3	B	28	23	23	3	0
4	B	26	10	10	5	0
5	A	8	16	0	0	0
5	B	198	396	0	3	1
All	All	2307	2479	2036	27	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 7.

All (27) 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:B:195[A]:SER:HG	3:B:251:CR9:HN3	1.25	0.84
2:B:92:LYS:H	2:B:92:LYS:HD2	1.57	0.69
4:B:1:CIT:O7	4:B:2:CIT:O7	2.13	0.67
2:B:92:LYS:H	2:B:92:LYS:CD	2.08	0.66
2:B:195[A]:SER:OG	3:B:251:CR9:N3	2.29	0.65
2:B:27:GLN:HG3	2:B:139[A]:THR:HG21	1.79	0.65
2:B:209[B]:LEU:CD1	2:B:231:VAL:HG21	2.33	0.59
2:B:170(B):HIS:H	2:B:170(B):HIS:CD2	2.28	0.52
2:B:47[A]:MET:HE2	2:B:123:LEU:CD1	2.41	0.51
2:B:195[A]:SER:OG	5:B:300:HOH:O	2.18	0.51
4:B:1:CIT:C5	4:B:2:CIT:O6	2.58	0.51
2:B:138:ILE:C	2:B:139[A]:THR:HG23	2.36	0.50
2:B:47[A]:MET:HE1	2:B:121:ILE:HG23	1.95	0.48
2:B:126:MET:HA	2:B:235[A]:LEU:HD12	1.95	0.48
2:B:27:GLN:CG	2:B:139[A]:THR:HG21	2.44	0.48
2:B:92:LYS:O	2:B:93:ASP:CB	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1:CIT:HO7	4:B:2:CIT:HO7	1.49	0.47
2:B:195[A]:SER:CB	5:B:300:HOH:O	2.63	0.46
3:B:251:CR9:N3	5:B:300:HOH:O	2.35	0.46
2:B:47[B]:MET:SD	2:B:53:ILE:HB	2.56	0.45
2:B:36:ARG:HH22	4:B:2:CIT:C5	2.30	0.44
2:B:209[B]:LEU:HD12	2:B:231:VAL:HG21	2.01	0.43
2:B:110:SER:OG	2:B:110(D):ARG:HG2	2.18	0.43
2:B:97(A):THR:O	2:B:97(B):LEU:HB2	2.19	0.42
4:B:1:CIT:O6	4:B:2:CIT:O7	2.36	0.42
2:B:47[A]:MET:HE2	2:B:123:LEU:HD13	2.01	0.41
1:A:11:PHE:CD2	2:B:120:THR:HG23	2.56	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)
2:B:76:ASN:OD1	5:B:259:HOH:H1[2_656]	1.58	0.02

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	7/23 (30%)	7 (100%)	0	0	100	100
2	B	252/253 (100%)	238 (94%)	12 (5%)	2 (1%)	16	3
All	All	259/276 (94%)	245 (95%)	12 (5%)	2 (1%)	16	3

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	93	ASP
2	B	97(B)	LEU

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	8/22 (36%)	7 (88%)	1 (12%)	4	0
2	B	222/218 (102%)	208 (94%)	14 (6%)	16	1
All	All	230/240 (96%)	215 (94%)	15 (6%)	16	1

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
2	B	20	GLU
2	B	60	ILE
2	B	60(A)	ASP
2	B	60(B)	TYR
2	B	60(C)	PRO
2	B	70	ARG
2	B	92	LYS
2	B	116	ARG
2	B	150	LEU
2	B	179	LYS
2	B	202	SER
2	B	203	LEU
2	B	209[A]	LEU
2	B	209[B]	LEU

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

Mol	Chain	Res	Type
1	A	15	GLN
2	B	78	GLN
2	B	91	HIS
2	B	101	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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
4	CIT	B	1	-	12,12,12	2.06	5 (41%)	17,17,17	1.59	3 (17%)
4	CIT	B	2	-	12,12,12	2.30	5 (41%)	17,17,17	1.34	3 (17%)
3	CR9	B	251	-	31,31,31	2.15	8 (25%)	43,45,45	1.68	10 (23%)

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	CIT	B	1	-	-	0/16/16/16	-
4	CIT	B	2	-	-	2/16/16/16	-
3	CR9	B	251	-	1/1/3/3	2/11/23/23	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	251	CR9	C4-N3	-5.68	1.28	1.38
4	B	2	CIT	C4-C3	4.53	1.59	1.54
3	B	251	CR9	C1-C7	-4.29	1.41	1.48
4	B	1	CIT	C2-C3	4.25	1.59	1.54
3	B	251	CR9	C1-C2	3.90	1.44	1.38
3	B	251	CR9	C4-C5	-3.83	1.34	1.40
3	B	251	CR9	C5-N4	-3.83	1.32	1.39
4	B	2	CIT	C2-C3	3.58	1.58	1.54
3	B	251	CR9	C1'-C8	-3.51	1.41	1.47
3	B	251	CR9	C8-N3	-2.93	1.30	1.36
3	B	251	CR9	C3-C4	-2.90	1.35	1.39
4	B	2	CIT	C4-C5	2.76	1.59	1.50
4	B	1	CIT	C4-C3	2.73	1.57	1.54
4	B	2	CIT	O2-C1	-2.65	1.22	1.30
4	B	2	CIT	O4-C5	-2.51	1.22	1.30
4	B	1	CIT	O4-C5	-2.43	1.22	1.30
4	B	1	CIT	O2-C1	-2.35	1.23	1.30
4	B	1	CIT	O6-C6	-2.33	1.22	1.30

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	251	CR9	C5-N4-C8	-5.81	98.39	105.17
4	B	1	CIT	C2-C3-C6	3.47	117.71	110.03
4	B	1	CIT	O7-C3-C6	-3.25	104.35	108.96
3	B	251	CR9	C4-N3-C8	-3.01	104.04	107.06
3	B	251	CR9	N3-C8-N4	2.75	117.13	112.44
4	B	2	CIT	O5-C6-C3	-2.69	116.89	122.09
3	B	251	CR9	C4-C5-N4	2.55	112.91	109.42
3	B	251	CR9	C1A-C1B-C6B	-2.50	108.42	112.62
3	B	251	CR9	C1'-C6'-C5'	2.43	121.54	119.56
4	B	2	CIT	O7-C3-C4	-2.31	104.11	109.38
3	B	251	CR9	C5'-O5'-C6B	-2.21	114.23	119.00
3	B	251	CR9	C6-C5-N4	-2.19	124.58	129.31
4	B	2	CIT	C2-C3-C6	2.18	114.85	110.03
4	B	1	CIT	O3-C5-C4	-2.16	116.84	122.95
3	B	251	CR9	C2'-C1'-C6'	-2.13	116.70	118.76
3	B	251	CR9	C6'-C1'-C8	2.02	121.97	119.55

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	251	CR9	C1B

All (4) torsion outliers are listed below:

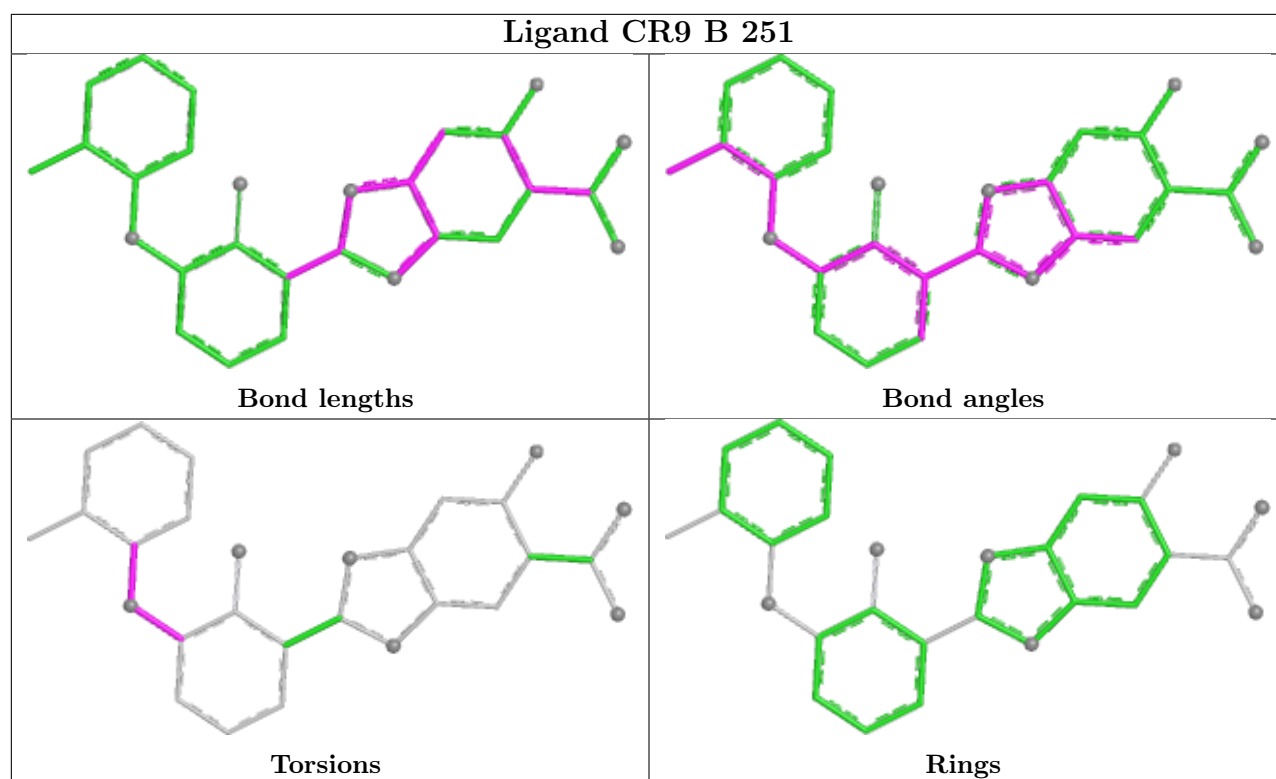
Mol	Chain	Res	Type	Atoms
3	B	251	CR9	C5B-C6B-O5'-C5'
4	B	2	CIT	C4-C3-C6-O5
4	B	2	CIT	C4-C3-C6-O6
3	B	251	CR9	C6'-C5'-O5'-C6B

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1	CIT	4	0
4	B	2	CIT	5	0
3	B	251	CR9	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	9/23 (39%)	0.50	2 (22%) 2 2	10, 19, 38, 41	3 (33%)
2	B	246/253 (97%)	-0.49	5 (2%) 65 71	5, 17, 33, 55	29 (11%)
All	All	255/276 (92%)	-0.46	7 (2%) 56 61	5, 18, 33, 55	32 (12%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	17	THR	3.3
2	B	243	LYS	3.3
1	A	9	LEU	3.0
2	B	97(A)	THR	2.8
2	B	97	ASP	2.5
2	B	242	THR	2.4
2	B	37(A)	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](#)

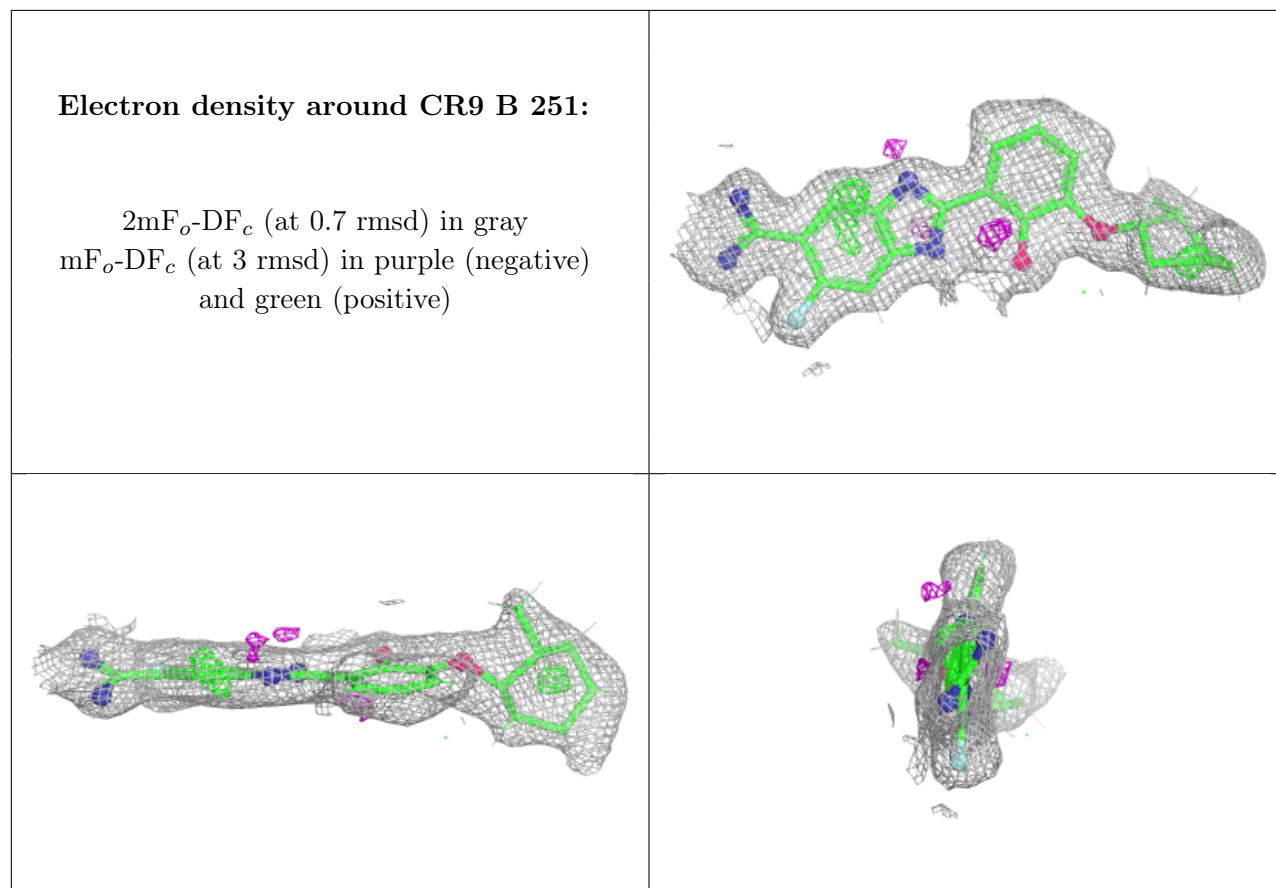
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
4	CIT	B	2	13/13	0.78	0.12	26,32,38,43	0
3	CR9	B	251	28/28	0.89	0.07	18,24,38,39	0
4	CIT	B	1	13/13	0.90	0.07	11,18,22,24	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.