

Full wwPDB X-ray Structure Validation Report (i)

Aug 16, 2023 - 06:16 PM EDT

PDB ID	:	2F57
Title	:	Crystal Structure Of The Human P21-Activated Kinase 5
Authors	:	Eswaran, J.; Turnbull, A.; Ugochukwu, E.; Papagrigoriou, E.; Bray, J.; Das,
		S.; Savitsky, P.; Smee, C.; Burgess, N.; Fedorov, O.; Filippakopoulos, P.; von
		Delft, F.; Arrowsmith, C.; Weigelt, J.; Sundstrom, M.; Edwards, A.; Knapp,
		S.; Structural Genomics Consortium (SGC)
Deposited on	:	2005-11-25
Resolution	:	1.80 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 (i) 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report		
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

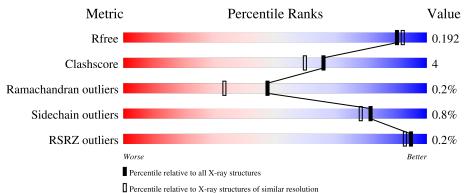


1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	А	317	90%	•	6%
1	В	317	86%	8%	5%

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	23D	В	1	-	-	Х	-



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5389 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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace		
1	Δ	298	Total	С	Ν	0	Р	\mathbf{S}	0	2	0
			2313	1471	399	426	1	16	0	2	0
1	D	200	Total	С	Ν	0	Р	S	0	0	0
	I B	300	2418	1539	421	441	1	16	0	0	U

• Molecule 1 is a protein called Serine/threonine-protein kinase PAK 7.

Chain	Residue	Modelled	Actual	Comment	Reference
А	-21	MET	-	cloning artifact	UNP Q9P286
А	-20	GLY	-	cloning artifact	UNP Q9P286
А	-19	SER	-	cloning artifact	UNP Q9P286
А	-18	SER	-	cloning artifact	UNP Q9P286
А	-17	HIS	-	expression tag	UNP Q9P286
А	-16	HIS	-	expression tag	UNP Q9P286
А	-15	HIS	-	expression tag	UNP Q9P286
А	-14	HIS	-	expression tag	UNP Q9P286
А	-13	HIS	-	expression tag	UNP Q9P286
A	-12	HIS	-	expression tag	UNP Q9P286
A	-11	SER	-	cloning artifact	UNP Q9P286
А	-10	SER	-	cloning artifact	UNP Q9P286
A	-9	GLY	-	cloning artifact	UNP Q9P286
А	-8	ARG	-	cloning artifact	UNP Q9P286
A	-7	GLU	-	cloning artifact	UNP Q9P286
А	-6	ASN	-	cloning artifact	UNP Q9P286
А	-5	LEU	-	cloning artifact	UNP Q9P286
А	-4	TYR	-	cloning artifact	UNP Q9P286
А	-3	PHE	-	cloning artifact	UNP Q9P286
А	-2	GLN	-	cloning artifact	UNP Q9P286
А	-1	SER	-	cloning artifact	UNP Q9P286
А	0	MET	-	cloning artifact	UNP Q9P286
А	602	SEP	SER	modified residue	UNP Q9P286
В	-21	MET	-	cloning artifact	UNP Q9P286
В	-20	GLY	-	cloning artifact	UNP Q9P286

There are 46 discrepancies between the modelled and reference sequences:

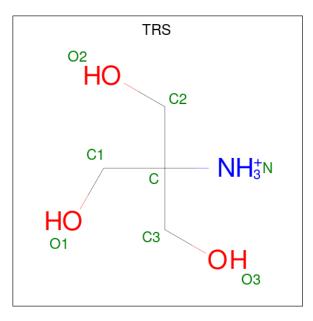
Continued on next page...



Chain	Residue	Modelled	Actual	Comment	Reference
В	-19	SER	-	cloning artifact	UNP Q9P286
В	-18	SER	-	cloning artifact	UNP Q9P286
В	-17	HIS	-	expression tag	UNP Q9P286
В	-16	HIS	-	expression tag	UNP Q9P286
В	-15	HIS	-	expression tag	UNP Q9P286
В	-14	HIS	-	expression tag	UNP Q9P286
В	-13	HIS	-	expression tag	UNP Q9P286
В	-12	HIS	-	expression tag	UNP Q9P286
В	-11	SER	-	cloning artifact	UNP Q9P286
В	-10	SER	-	cloning artifact	UNP Q9P286
В	-9	GLY	-	cloning artifact	UNP Q9P286
В	-8	ARG	-	cloning artifact	UNP Q9P286
В	-7	GLU	-	cloning artifact	UNP Q9P286
В	-6	ASN	-	cloning artifact	UNP Q9P286
В	-5	LEU	-	cloning artifact	UNP Q9P286
В	-4	TYR	-	cloning artifact	UNP Q9P286
В	-3	PHE	-	cloning artifact	UNP Q9P286
В	-2	GLN	-	cloning artifact	UNP Q9P286
В	-1	SER	-	cloning artifact	UNP Q9P286
В	0	MET	-	cloning artifact	UNP Q9P286
В	602	SEP	SER	modified residue	UNP Q9P286

Continued from previous page...

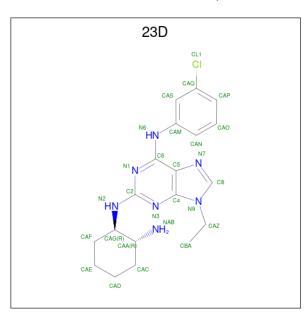
• Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{N} & \text{O} \\ 8 & 4 & 1 & 3 \end{array}$	0	0
2	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{N} & \text{O} \\ 8 & 4 & 1 & 3 \end{array}$	0	0

• Molecule 3 is N2-[(1R,2S)-2-AMINOCYCLOHEXYL]-N6-(3-CHLOROPHENYL)-9-ETHY L-9H-PURINE-2,6-DIAMINE (three-letter code: 23D) (formula: C₁₉H₂₄ClN₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	D	1	Total	С	Cl	Ν	0	0
0	3 B		27	19	1	7	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	239	Total O 240 240	0	1
4	В	374	Total O 375 375	0	1



3 Residue-property plots (i)

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 PAK 7

Chain A:	90%	• 6%
MET GLY SER HIS HIS HIS HIS HIS HIS RIS SER SER SER ARC CLY ARC CLY ARC CLY ARC CLY M-6 1465 1465 1465	1538 R567 R567 R567 R567 L603 V604 V604 L649 M555 M555 M555 H15 H15 H15 H15 H15 H15 H15	
• Molecule 1: Serine/threonine-pro	otein kinase PAK 7	
Chain B:	86%	8% 5%
MET SER SER HIS HIS HIS HIS HIS HIS SER SER SER SER SER SER SER SER SER SE	K478 K479 M480 F492 F492 F492 F525 F525 E523 E524 F525 E525 E526 E526 E526 E526 E526 E527 E527 E527 F557 S578 S578	1582 K583 L584 L584 R600 8601 S602 L603 V604
P607 1655 P660 R683 M713 R714 H18 H18		



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	98.66Å 56.60 Å 120.76 Å	Deperitor
a, b, c, α , β , γ	90.00° 103.13° 90.00°	Depositor
Resolution (Å)	50.00 - 1.80	Depositor
Resolution (A)	48.77 - 1.80	EDS
% Data completeness	99.4 (50.00-1.80)	Depositor
(in resolution range)	99.3(48.77-1.80)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.45 (at 1.79 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D	0.154 , 0.185	Depositor
R, R_{free}	0.163 , 0.192	DCC
R_{free} test set	3031 reflections $(5.05%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	22.0	Xtriage
Anisotropy	0.372	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.34 , 54.3	EDS
L-test for twinning ²	$< L > = 0.50, < L^2 > = 0.34$	Xtriage
	0.012 for 1/2*h-3/2*k,-1/2*h-1/2*k,-1/2*h	
Estimated twinning fraction	$+1/2^{*}$ k-l 0.011 for $1/2^{*}$ h+ $3/2^{*}$ k, $1/2^{*}$ h- $1/2^{*}$ k, $-1/2^{*}$ h-	Xtriage
	$1/2^*$ k-l	
F_o, F_c correlation	0.97	EDS
Total number of atoms	5389	wwPDB-VP
Average B, all atoms $(Å^2)$	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.42% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: 23D, TRS, SEP

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	Bo	nd angles
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.39	0/2356	0.58	1/3196~(0.0%)
1	В	0.52	0/2483	0.65	0/3361
All	All	0.46	0/4839	0.62	1/6557~(0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	539	ARG	NE-CZ-NH1	-5.94	117.33	120.30

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	А	2313	0	2285	9	0
1	В	2418	0	2445	27	0
2	А	16	0	24	0	0
3	В	27	0	24	10	0
4	А	240	0	0	1	0
4	В	375	0	0	3	0
All	All	5389	0	4778	38	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 4.

All (38) 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:B:523:MET:HE3	3:B:1:23D:CAZ	2.08	0.82
1:A:649:LEU:HD23	1:A:652:MET:HE3	1.63	0.78
1:B:524:GLU:O	3:B:1:23D:H8	1.85	0.76
1:A:649:LEU:HA	1:A:652:MET:HE3	1.68	0.76
1:B:523:MET:CE	3:B:1:23D:HAZ1	2.17	0.75
1:B:523:MET:HE3	3:B:1:23D:HAZ2	1.69	0.73
1:B:523:MET:CE	3:B:1:23D:CAZ	2.70	0.68
1:B:567:ARG:CZ	1:B:604[A]:VAL:HG21	2.25	0.67
1:A:649:LEU:HD23	1:A:652:MET:CE	2.27	0.64
1:B:567:ARG:CZ	1:B:604[B]:VAL:HG11	2.29	0.63
1:A:607:PRO:HB3	1:A:652:MET:HE2	1.82	0.62
1:A:567:ARG:CZ	1:A:604:VAL:HG21	2.29	0.62
1:B:478:LYS:HE3	1:B:523:MET:HE1	1.84	0.60
1:B:508:ASP:HB2	1:B:524:GLU:HB2	1.84	0.59
1:A:652:MET:HA	1:A:655:ILE:HD12	1.86	0.57
3:B:1:23D:HBA3	4:B:1090:HOH:O	2.05	0.57
1:A:655:ILE:HG12	4:A:907:HOH:O	2.04	0.57
1:B:431:GLN:OE1	4:B:1072:HOH:O	2.18	0.56
1:B:714[A]:ARG:CZ	1:B:714[A]:ARG:HB2	2.35	0.55
1:B:655:ILE:HG12	4:B:1085:HOH:O	2.06	0.54
1:B:523:MET:HE3	3:B:1:23D:HAZ1	1.80	0.54
1:B:436:LEU:HD23	1:B:492:PHE:CE1	2.43	0.53
1:B:567:ARG:NE	1:B:604[A]:VAL:HG21	2.24	0.53
3:B:1:23D:HAS	3:B:1:23D:N1	2.24	0.52
1:A:462:ILE:HG22	1:A:479:LYS:HB2	1.91	0.51
1:B:567:ARG:NH2	1:B:604[B]:VAL:HG11	2.26	0.51
1:B:436:LEU:HD23	1:B:492:PHE:CD1	2.48	0.49
1:B:602:SEP:HB2	1:B:604[A]:VAL:HG13	1.94	0.49
1:B:660:PRO:HG3	1:B:683:ARG:HG3	1.97	0.47
1:A:455:ILE:HD13	1:A:465:ILE:CG1	2.45	0.46
1:B:523:MET:HE1	3:B:1:23D:HAZ1	1.94	0.46
1:B:456:GLY:HA3	3:B:1:23D:HAE1	1.97	0.46
1:B:569:ILE:HD13	1:B:584:LEU:HD11	2.00	0.43
1:B:600:ARG:HB3	1:B:602:SEP:O1P	2.18	0.43
1:B:576:LEU:HG	1:B:582:ILE:HD13	2.00	0.43
1:B:578:SER:O	1:B:713:MET:HE1	2.21	0.41
1:B:480:MET:HG2	1:B:488:ARG:HA	2.04	0.41

Continued on next page...



α \cdot \cdot \cdot	C		
Continued	trom	previous	<i>paae</i>
00100000000	J. 01.0	proceedae	P ~ 9 ~

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:525:PHE:CE2	1:B:527:GLU:HG2	2.55	0.40

There are no symmetry-related clashes.

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	Perce	ntiles
1	А	297/317~(94%)	291~(98%)	5(2%)	1 (0%)	41	27
1	В	305/317~(96%)	298~(98%)	7~(2%)	0	100	100
All	All	602/634~(95%)	589~(98%)	12 (2%)	1 (0%)	47	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	607	PRO

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	А	248/282~(88%)	245~(99%)	3~(1%)	71 65
1	В	269/282~(95%)	268 (100%)	1 (0%)	91 89
All	All	517/564~(92%)	513~(99%)	4 (1%)	81 78



All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	538	THR
1	А	539	ARG
1	А	607	PRO
1	В	607	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	650	GLN
1	В	537	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)

2 non-standard protein/DNA/RNA residues 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	B	Bond lengths			Bond angles		
IVIOI	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
1	SEP	В	602	1	8,9,10	1.31	1 (12%)	8,12,14	2.08	4 (50%)	
1	SEP	А	602	1	8,9,10	1.54	1 (12%)	8,12,14	1.23	1 (12%)	

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
1	SEP	В	602	1	-	1/5/8/10	-
1	SEP	А	602	1	-	0/5/8/10	-



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	602	SEP	P-O1P	3.47	1.61	1.50
1	В	602	SEP	P-O1P	2.38	1.58	1.50

All (2) bond length outliers are listed below:

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	602	SEP	OG-P-O1P	3.12	115.23	106.47
1	В	602	SEP	OG-CB-CA	2.67	110.75	108.14
1	В	602	SEP	O3P-P-OG	-2.49	100.12	106.73
1	В	602	SEP	O3P-P-O2P	2.19	116.00	107.64
1	А	602	SEP	OG-P-O1P	2.13	112.44	106.47

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	В	602	SEP	CB-OG-P-O3P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	В	602	SEP	2	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	Turne	Chain	Res	Link	Bo	Bond lengths			Bond angles		
INIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
3	23D	В	1	-	26,30,30	1.07	3 (11%)	29,42,42	1.91	6 (20%)	
2	TRS	А	802	-	7,7,7	0.21	0	9,9,9	0.70	0	
2	TRS	А	801	-	7,7,7	0.34	0	9,9,9	0.32	0	

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	23D	В	1	-	-	3/10/21/21	0/4/4/4
2	TRS	А	802	-	-	0/9/9/9	-
2	TRS	А	801	-	_	0/9/9/9	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	В	1	23D	CAZ-N9	-2.62	1.47	1.49
3	В	1	23D	CAM-N6	-2.46	1.35	1.40
3	В	1	23D	C6-N1	2.40	1.36	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	В	1	23D	N6-C6-N1	6.33	124.60	118.66
3	В	1	23D	C2-N3-C4	4.19	120.03	115.28
3	В	1	23D	N3-C2-N1	-3.35	120.92	126.23
3	В	1	23D	C5-C6-N1	-3.24	118.12	120.81
3	В	1	23D	N2-C2-N1	2.65	121.15	117.18
3	В	1	23D	C2-N1-C6	2.16	121.33	116.39

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	В	1	23D	CBA-CAZ-N9-C4
3	В	1	23D	CAF-CAG-N2-C2
3	В	1	23D	CAA-CAG-N2-C2

There are no ring outliers.

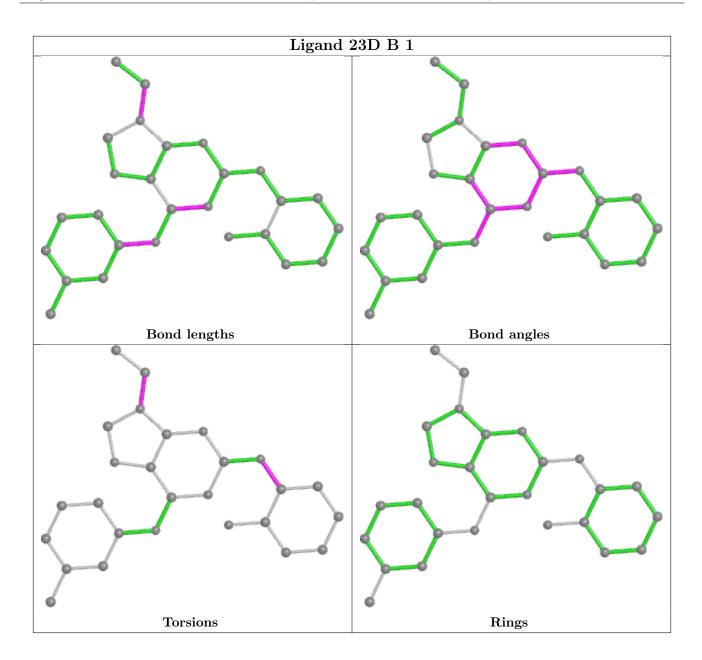
1 monomer is involved in 10 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	1	23D	10	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 then 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, 95^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	297/317~(93%)	-0.44	0 100 100	11, 18, 26, 42	0
1	В	299/317~(94%)	-0.65	1 (0%) 94 92	13, 17, 27, 44	0
All	All	596/634~(94%)	-0.55	1 (0%) 95 93	11, 17, 27, 44	0

All (1) RSRZ outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	RSRZ
1	В	425	SER	2.4

6.2 Non-standard residues in protein, DNA, RNA chains (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^{th} 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
1	SEP	А	602	10/11	0.96	0.09	18,23,30,38	0
1	SEP	В	602	10/11	0.97	0.09	13,20,30,41	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

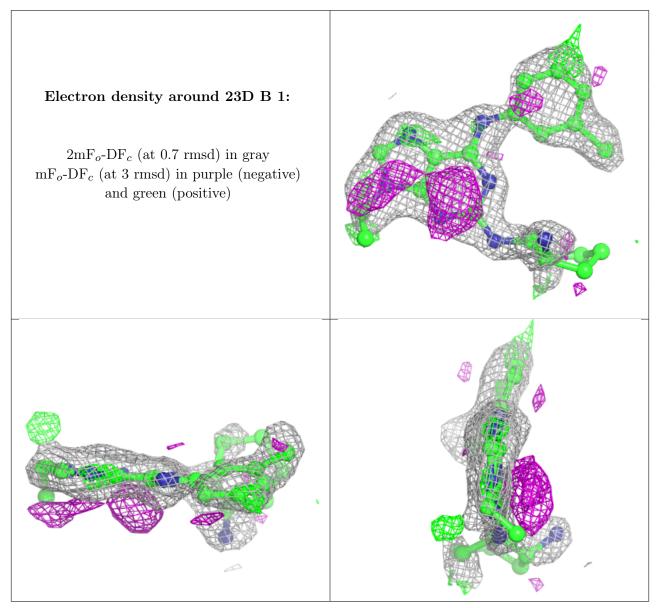
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q < 0.9
3	23D	В	1	27/27	0.75	0.24	$27,\!53,\!77,\!80$	0
2	TRS	А	802	8/8	0.83	0.12	33,36,43,47	0
2	TRS	А	801	8/8	0.86	0.14	45,51,55,66	0

median, 95^{th} 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.

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 (i)

There are no such residues in this entry.

