



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:33 AM UTC

PDB ID : 3AXW / pdb_00003axw
Title : Crystal structure of human CK2alpha complexed with a potent inhibitor
Authors : Kinoshita, T.; Nakanishi, I.
Deposited on : 2011-04-18
Resolution : 2.50 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

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
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

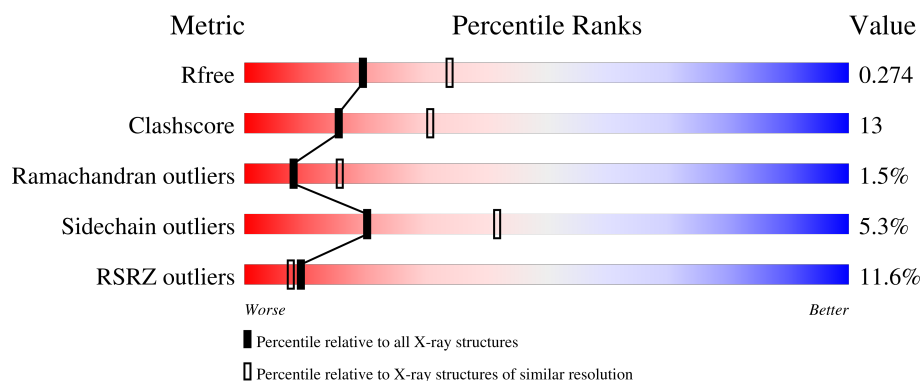
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	340	11% 69% 22% ...

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2914 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.

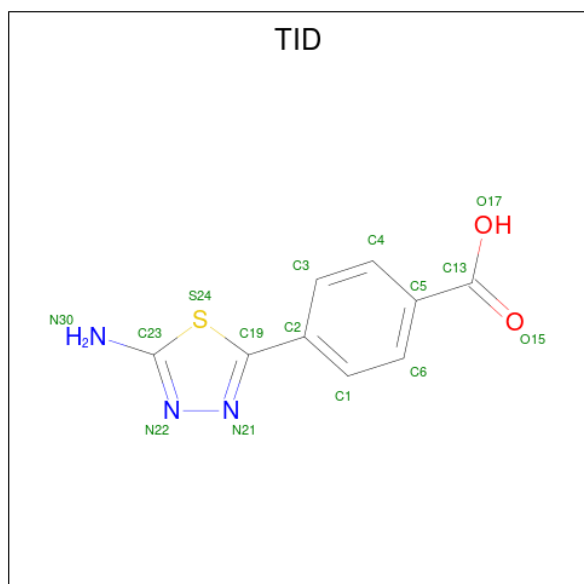
- Molecule 1 is a protein called Casein kinase II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2779	1781	488	499	11	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P68400
A	-3	PRO	-	expression tag	UNP P68400
A	-2	LEU	-	expression tag	UNP P68400
A	-1	GLY	-	expression tag	UNP P68400
A	0	SER	-	expression tag	UNP P68400

- Molecule 2 is 4-(5-amino-1,3,4-thiadiazol-2-yl)benzoic acid (CCD ID: TID) (formula: $C_9H_7N_3O_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	9	3	2	1		

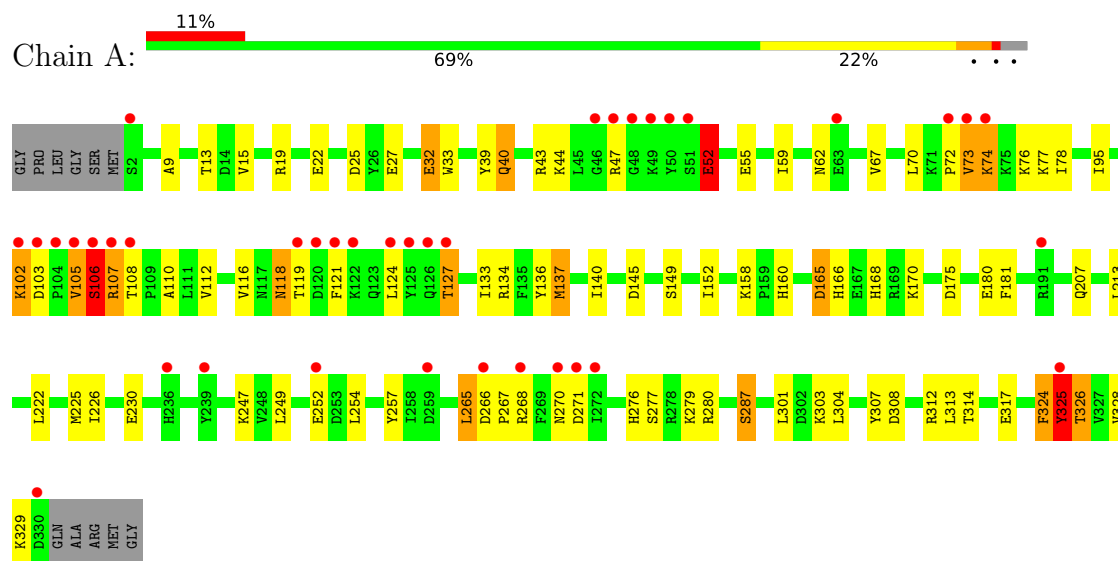
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	120	Total	O	0	0
			120	120		

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: Casein kinase II subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.19Å 80.22Å 83.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 95.9 (50.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.51Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.268 , 0.295 0.238 , 0.274	Depositor DCC
R_{free} test set	530 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å ²)	37.3	Xtriage
Anisotropy	0.428	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,l,k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2914	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.38% 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: TID

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	0.59	4/2854 (0.1%)	1.16	28/3860 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	GLN	CD-NE2	-12.76	1.06	1.33
1	A	325	TYR	CD2-CE2	11.88	1.74	1.38
1	A	325	TYR	CE2-CZ	-6.86	1.21	1.38
1	A	325	TYR	CG-CD2	-6.28	1.26	1.39

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	ASN	N-CA-C	9.86	124.09	109.59
1	A	40	GLN	CG-CD-NE2	-9.76	101.76	116.40
1	A	55	GLU	N-CA-C	-9.48	95.22	110.20
1	A	326	THR	N-CA-C	-9.08	101.38	111.28
1	A	207	GLN	N-CA-C	7.76	122.01	112.23
1	A	280	ARG	CG-CD-NE	-7.65	95.18	112.00
1	A	119	THR	N-CA-C	-7.64	96.78	109.46
1	A	134	ARG	NE-CZ-NH2	7.47	125.92	119.20
1	A	134	ARG	NE-CZ-NH1	-7.46	114.05	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	324	PHE	N-CA-C	6.97	121.10	112.59
1	A	13	THR	N-CA-C	6.50	118.36	111.28
1	A	152	ILE	N-CA-C	6.43	117.38	108.12
1	A	32	GLU	N-CA-C	-6.28	98.96	109.07
1	A	52	GLU	N-CA-C	-6.17	98.22	108.34
1	A	213	LEU	N-CA-C	-6.14	104.14	111.69
1	A	33	TRP	N-CA-C	6.09	119.97	110.17
1	A	134	ARG	CG-CD-NE	-6.09	98.61	112.00
1	A	287	SER	N-CA-C	6.07	120.30	113.01
1	A	39	TYR	N-CA-C	6.07	119.30	109.40
1	A	280	ARG	CD-NE-CZ	6.03	132.84	124.40
1	A	280	ARG	NE-CZ-NH1	-5.87	115.63	121.50
1	A	137	MET	N-CA-C	-5.87	104.97	111.36
1	A	9	ALA	N-CA-C	-5.57	102.15	110.23
1	A	165	ASP	N-CA-C	-5.52	98.25	108.24
1	A	280	ARG	NE-CZ-NH2	5.39	124.05	119.20
1	A	59	ILE	N-CA-C	5.30	120.36	109.34
1	A	166	HIS	N-CA-C	5.22	119.99	111.37
1	A	271	ASP	N-CA-C	5.21	117.86	111.82

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	325	TYR	Sidechain
1	A	40	GLN	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	2779	0	2722	74	0
2	A	15	0	6	0	0
3	A	120	0	0	9	0
All	All	2914	0	2728	74	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 13.

All (74) 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:325:TYR:CE1	1:A:329:LYS:HB2	1.65	1.30
1:A:121:PHE:HE2	1:A:160:HIS:CD2	1.56	1.24
1:A:121:PHE:CE2	1:A:160:HIS:CD2	2.28	1.21
1:A:325:TYR:HE1	1:A:329:LYS:HD2	1.10	1.14
1:A:325:TYR:HE1	1:A:329:LYS:CD	1.65	1.07
1:A:121:PHE:HE2	1:A:160:HIS:NE2	1.52	1.07
1:A:325:TYR:CE1	1:A:329:LYS:HD2	1.89	1.06
1:A:325:TYR:CE1	1:A:329:LYS:CB	2.44	1.00
1:A:314:THR:OG1	1:A:317:GLU:HG3	1.67	0.95
1:A:25:ASP:HB3	3:A:440:HOH:O	1.66	0.95
1:A:325:TYR:CD1	1:A:325:TYR:O	2.24	0.89
1:A:22:GLU:HA	3:A:440:HOH:O	1.72	0.88
1:A:121:PHE:CE2	1:A:160:HIS:HD2	1.87	0.87
1:A:106:SER:C	1:A:108:THR:H	1.85	0.85
1:A:121:PHE:CE2	1:A:160:HIS:NE2	2.38	0.84
1:A:325:TYR:CD1	1:A:325:TYR:C	2.55	0.83
1:A:325:TYR:CZ	1:A:329:LYS:HB2	2.12	0.83
1:A:325:TYR:CE1	1:A:329:LYS:CD	2.57	0.82
1:A:165:ASP:OD2	1:A:168:HIS:HD2	1.68	0.77
1:A:73:VAL:HG13	1:A:73:VAL:O	1.87	0.73
1:A:265:LEU:HD13	3:A:383:HOH:O	1.88	0.73
1:A:102:LYS:NZ	1:A:107:ARG:HB3	2.05	0.71
1:A:44:LYS:NZ	1:A:52:GLU:OE2	2.26	0.69
1:A:102:LYS:HZ2	1:A:107:ARG:HB3	1.59	0.67
1:A:165:ASP:HB3	1:A:170:LYS:HB3	1.76	0.67
1:A:106:SER:C	1:A:108:THR:N	2.53	0.64
1:A:106:SER:O	1:A:108:THR:N	2.34	0.61
1:A:136:TYR:O	1:A:140:ILE:HG13	1.98	0.61
1:A:107:ARG:HG2	3:A:352:HOH:O	2.03	0.57
1:A:27:GLU:OE2	1:A:76:LYS:HE2	2.05	0.57
1:A:325:TYR:CE1	1:A:329:LYS:CG	2.87	0.57
1:A:70:LEU:HD13	1:A:78:ILE:HG12	1.86	0.57
1:A:165:ASP:OD2	1:A:168:HIS:CD2	2.56	0.57
1:A:15:VAL:O	1:A:19:ARG:HG2	2.04	0.57
1:A:267:PRO:HA	1:A:270:ASN:ND2	2.20	0.56
1:A:314:THR:HG1	1:A:317:GLU:HG3	1.68	0.55
1:A:325:TYR:C	1:A:325:TYR:HD1	2.11	0.55
1:A:133:ILE:HG12	1:A:225:MET:HE2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:TYR:HE1	1:A:329:LYS:CG	2.19	0.52
1:A:254:LEU:HD13	1:A:307:TYR:CE1	2.45	0.52
1:A:124:LEU:HA	3:A:405:HOH:O	2.11	0.51
1:A:74:LYS:HG2	1:A:76:LYS:H	1.76	0.51
1:A:267:PRO:HA	1:A:270:ASN:HD21	1.75	0.50
1:A:257:TYR:CE1	1:A:308:ASP:HA	2.47	0.49
1:A:324:PHE:O	1:A:328:VAL:HG23	2.12	0.49
1:A:73:VAL:O	1:A:73:VAL:CG1	2.60	0.48
1:A:226:ILE:HD11	1:A:301:LEU:HD22	1.95	0.48
1:A:137:MET:HE2	1:A:222:LEU:HD13	1.95	0.47
1:A:180:GLU:HG3	1:A:181:PHE:N	2.29	0.47
1:A:73:VAL:O	1:A:74:LYS:O	2.32	0.47
1:A:72:PRO:HA	3:A:384:HOH:O	2.15	0.47
1:A:325:TYR:O	1:A:329:LYS:HB3	2.15	0.47
1:A:103:ASP:HB2	1:A:110:ALA:HB2	1.97	0.46
1:A:116:VAL:O	1:A:118:ASN:N	2.48	0.46
1:A:108:THR:HG22	3:A:384:HOH:O	2.15	0.46
1:A:67:VAL:HG22	1:A:112:VAL:HG22	1.98	0.46
1:A:127:THR:HA	3:A:375:HOH:O	2.15	0.46
1:A:303:LYS:HB3	1:A:313:LEU:HG	1.98	0.45
1:A:95:ILE:O	1:A:95:ILE:HG23	2.18	0.43
1:A:121:PHE:CD1	1:A:121:PHE:C	2.96	0.43
1:A:268:ARG:HG3	1:A:268:ARG:HH21	1.84	0.43
1:A:145:ASP:O	1:A:149:SER:HB3	2.18	0.43
1:A:247:LYS:HG2	1:A:276:HIS:CE1	2.53	0.42
1:A:70:LEU:O	1:A:108:THR:HG21	2.19	0.42
1:A:72:PRO:O	1:A:73:VAL:HB	2.20	0.42
1:A:52:GLU:HG2	3:A:453:HOH:O	2.20	0.42
1:A:304:LEU:O	1:A:312:ARG:HG2	2.20	0.42
1:A:158:LYS:HE3	1:A:160:HIS:HB2	2.01	0.42
1:A:62:ASN:HD22	1:A:62:ASN:HA	1.68	0.41
1:A:103:ASP:OD1	1:A:105:VAL:HG23	2.21	0.41
1:A:277:SER:O	1:A:279:LYS:NZ	2.48	0.41
1:A:267:PRO:O	1:A:270:ASN:ND2	2.53	0.41
1:A:313:LEU:HA	1:A:317:GLU:OE1	2.20	0.40
1:A:254:LEU:HB2	1:A:307:TYR:CZ	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	Percentiles	
1	A	327/340 (96%)	303 (93%)	19 (6%)	5 (2%)	8	16

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	VAL
1	A	74	LYS
1	A	106	SER
1	A	107	ARG
1	A	175	ASP

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	303/310 (98%)	287 (95%)	16 (5%)	20	42

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	GLU
1	A	43	ARG
1	A	47	ARG
1	A	52	GLU
1	A	77	LYS
1	A	102	LYS

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Mol	Chain	Res	Type
1	A	105	VAL
1	A	106	SER
1	A	127	THR
1	A	230	GLU
1	A	249	LEU
1	A	252	GLU
1	A	265	LEU
1	A	266	ASP
1	A	287	SER
1	A	326	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	40	GLN
1	A	62	ASN
1	A	118	ASN
1	A	160	HIS
1	A	168	HIS
1	A	262	ASN
1	A	270	ASN
1	A	276	HIS
1	A	310	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TID	A	336	-	16,16,16	2.91	7 (43%)	22,22,22	5.92	8 (36%)

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
2	TID	A	336	-	-	0/8/8/8	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	336	TID	C23-N30	6.87	1.48	1.34
2	A	336	TID	C23-N22	5.11	1.38	1.31
2	A	336	TID	C2-C19	-4.80	1.40	1.47
2	A	336	TID	C23-S24	-2.92	1.70	1.74
2	A	336	TID	C19-N21	2.50	1.36	1.30
2	A	336	TID	C1-C2	2.16	1.42	1.39
2	A	336	TID	C6-C5	2.10	1.42	1.39

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	336	TID	C23-S24-C19	22.48	94.10	86.98
2	A	336	TID	N30-C23-N22	11.47	133.50	123.83
2	A	336	TID	S24-C23-N22	-7.53	108.39	113.79
2	A	336	TID	S24-C19-N21	-5.67	107.52	113.58
2	A	336	TID	S24-C23-N30	-2.97	118.11	121.97
2	A	336	TID	C2-C19-N21	2.88	128.46	122.80
2	A	336	TID	C23-N22-N21	2.51	114.40	112.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	336	TID	C19-N21-N22	2.41	115.58	113.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	329/340 (96%)	0.43	38 (11%) 9 8	9, 28, 71, 90	3 (0%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	105	VAL	7.9
1	A	104	PRO	7.9
1	A	103	ASP	7.4
1	A	325	TYR	6.2
1	A	108	THR	6.2
1	A	107	ARG	5.4
1	A	106	SER	5.1
1	A	50	TYR	4.2
1	A	2	SER	3.9
1	A	49	LYS	3.8
1	A	125	TYR	3.7
1	A	51	SER	3.6
1	A	102	LYS	3.5
1	A	72	PRO	3.4
1	A	46	GLY	3.1
1	A	126	GLN	2.9
1	A	121	PHE	2.8
1	A	48	GLY	2.8
1	A	73	VAL	2.8
1	A	74	LYS	2.8
1	A	127	THR	2.7
1	A	271	ASP	2.6
1	A	236	HIS	2.6
1	A	259	ASP	2.5
1	A	266	ASP	2.5
1	A	268	ARG	2.4
1	A	119	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	330	ASP	2.4
1	A	270	ASN	2.3
1	A	122	LYS	2.3
1	A	124	LEU	2.3
1	A	63	GLU	2.3
1	A	120	ASP	2.3
1	A	47	ARG	2.1
1	A	191	ARG	2.1
1	A	252	GLU	2.1
1	A	272	ILE	2.1
1	A	239	TYR	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($\text{\AA}^2$)	Q<0.9
2	TID	A	336	15/15	0.87	0.11	35,38,40,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.