



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 22, 2016 – 01:38 PM EST

PDB ID : 5CX9  
Title : Crystal structure of CK2alpha with (methyl 4-((3-(3-chloro-4-(phenyl)benzyl amino)propyl)amino)-4-oxobutanoate bound  
Authors : Brear, P.; De Fusco, C.; Georgiou, K.H.; Spring, D.; Hyvonen, M.  
Deposited on : 2015-07-28  
Resolution : 1.73 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-20028320  
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-20028320

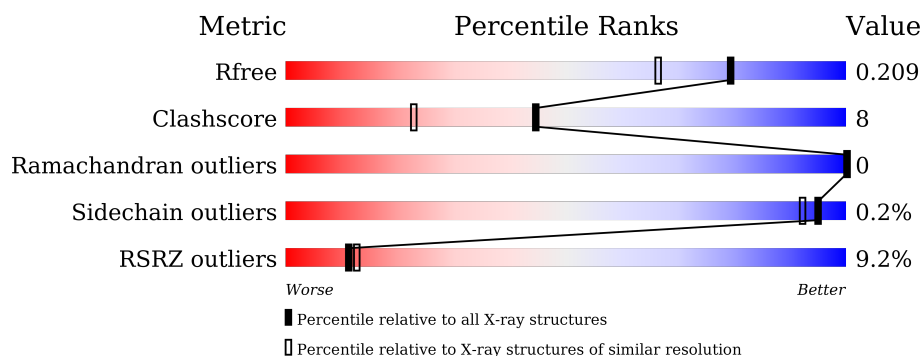
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.73 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	2417 (1.76-1.72)
Clashscore	102246	2570 (1.76-1.72)
Ramachandran outliers	100387	2544 (1.76-1.72)
Sidechain outliers	100360	2544 (1.76-1.72)
RSRZ outliers	91569	2420 (1.76-1.72)

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	352	<div> <div>5%</div> <div>85%</div> <div>8%</div> <div>7%</div> </div>
1	B	352	<div> <div>13%</div> <div>84%</div> <div>8%</div> <div>8%</div> </div>

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
2	ACT	B	402	-	-	-	X
3	551	A	406	-	-	-	X
3	551	B	404	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6033 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
1	A	328	Total	C	N	O	S	0	7	0
			2821	1804	495	511	11			
1	B	325	Total	C	N	O	S	0	4	0
			2773	1775	485	502	11			

There are 50 discrepancies between the modelled and reference sequences:

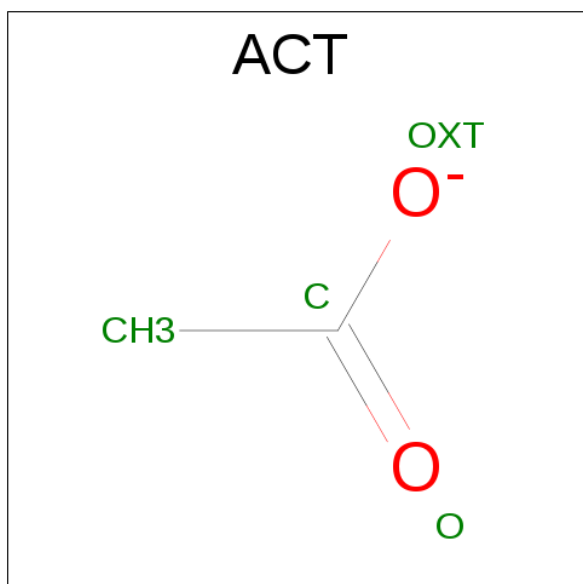
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	GLY	-	expression tag	UNP P68400
A	-21	SER	-	expression tag	UNP P68400
A	-20	MET	-	expression tag	UNP P68400
A	-19	ASP	-	expression tag	UNP P68400
A	-18	ILE	-	expression tag	UNP P68400
A	-17	GLU	-	expression tag	UNP P68400
A	-16	PHE	-	expression tag	UNP P68400
A	-15	ASP	-	expression tag	UNP P68400
A	-14	ASP	-	expression tag	UNP P68400
A	-13	ASP	-	expression tag	UNP P68400
A	-12	ALA	-	expression tag	UNP P68400
A	-11	ASP	-	expression tag	UNP P68400
A	-10	ASP	-	expression tag	UNP P68400
A	-9	ASP	-	expression tag	UNP P68400
A	-8	GLY	-	expression tag	UNP P68400
A	-7	SER	-	expression tag	UNP P68400
A	-6	GLY	-	expression tag	UNP P68400
A	-5	SER	-	expression tag	UNP P68400
A	-4	GLY	-	expression tag	UNP P68400
A	-3	SER	-	expression tag	UNP P68400
A	-2	GLY	-	expression tag	UNP P68400
A	-1	SER	-	expression tag	UNP P68400
A	0	GLY	-	expression tag	UNP P68400
A	1	SER	-	expression tag	UNP P68400
A	21	SER	ARG	engineered mutation	UNP P68400

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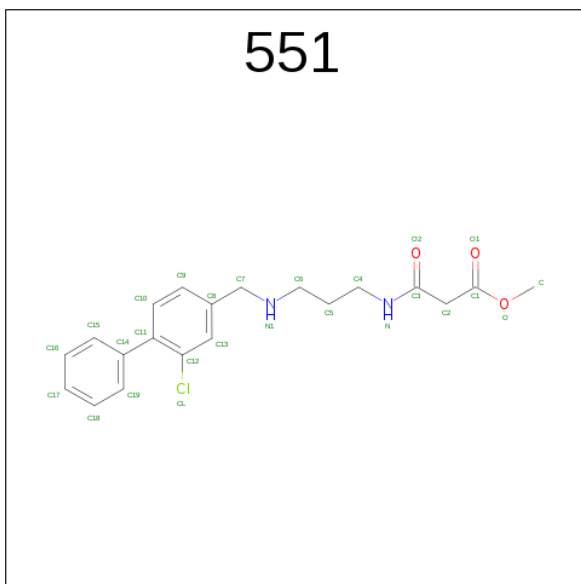
Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	GLY	-	expression tag	UNP P68400
B	-21	SER	-	expression tag	UNP P68400
B	-20	MET	-	expression tag	UNP P68400
B	-19	ASP	-	expression tag	UNP P68400
B	-18	ILE	-	expression tag	UNP P68400
B	-17	GLU	-	expression tag	UNP P68400
B	-16	PHE	-	expression tag	UNP P68400
B	-15	ASP	-	expression tag	UNP P68400
B	-14	ASP	-	expression tag	UNP P68400
B	-13	ASP	-	expression tag	UNP P68400
B	-12	ALA	-	expression tag	UNP P68400
B	-11	ASP	-	expression tag	UNP P68400
B	-10	ASP	-	expression tag	UNP P68400
B	-9	ASP	-	expression tag	UNP P68400
B	-8	GLY	-	expression tag	UNP P68400
B	-7	SER	-	expression tag	UNP P68400
B	-6	GLY	-	expression tag	UNP P68400
B	-5	SER	-	expression tag	UNP P68400
B	-4	GLY	-	expression tag	UNP P68400
B	-3	SER	-	expression tag	UNP P68400
B	-2	GLY	-	expression tag	UNP P68400
B	-1	SER	-	expression tag	UNP P68400
B	0	GLY	-	expression tag	UNP P68400
B	1	SER	-	expression tag	UNP P68400
B	21	SER	ARG	engineered mutation	UNP P68400

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula:  $\text{C}_2\text{H}_3\text{O}_2$ ).



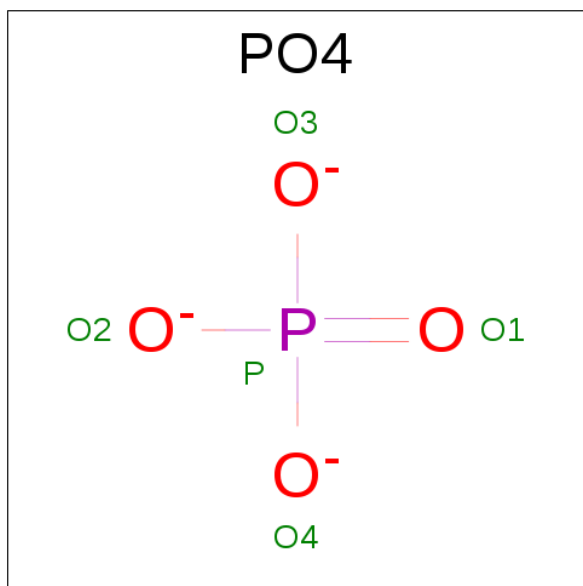
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0

- Molecule 3 is methyl 3-[(3-[(2-chlorobiphenyl-4-yl)methyl]amino}propyl)amino]-3-oxopropionate (three-letter code: 551) (formula: C<sub>20</sub>H<sub>23</sub>ClN<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C Cl N O 26 20 1 2 3	0	0
3	A	1	Total C Cl N 15 13 1 1	0	0
3	A	1	Total C Cl N 15 13 1 1	0	0
3	B	1	Total C Cl N O 26 20 1 2 3	0	0
3	B	1	Total C Cl N 15 13 1 1	0	0

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	227	Total	O	0	0
			227	227		
5	B	86	Total	O	0	0
			86	86		





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.68Å 68.13Å 334.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.46 – 1.73 46.46 – 1.73	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.46-1.73) 99.7 (46.46-1.73)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.22 (at 1.73Å)	Xtriage
Refinement program	BUSTER-TNT	Depositor
R, $R_{free}$	0.189 , 0.198 0.195 , 0.209	Depositor DCC
$R_{free}$ test set	3904 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.9	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 48.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.066 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6033	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.61% 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, 0.333 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: PO4, 551, ACT

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	0.70	2/2896 (0.1%)	0.67	1/3916 (0.0%)
1	B	0.57	0/2848	0.65	0/3854
All	All	0.64	2/5744 (0.0%)	0.66	1/7770 (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
1	A	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	296	GLU	CD-OE1	-5.55	1.19	1.25
1	A	180	GLU	CD-OE1	-5.52	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	C-N-CD	5.40	139.73	128.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	18	HIS	Mainchain

## 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	2821	0	2755	32	0
1	B	2773	0	2701	58	0
2	A	16	0	12	0	0
2	B	8	0	6	0	0
3	A	56	0	0	6	0
3	B	41	0	0	6	0
4	A	5	0	0	0	0
5	A	227	0	0	0	0
5	B	86	0	0	1	0
All	All	6033	0	5474	88	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 8.

All (88) 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:36:GLN:HE21	1:B:103[B]:ASP:CA	1.42	1.31
1:B:36:GLN:NE2	1:B:103[A]:ASP:OD1	1.65	1.30
1:B:36:GLN:NE2	1:B:103[A]:ASP:HA	1.50	1.25
1:B:36:GLN:NE2	1:B:103[B]:ASP:CB	2.01	1.23
1:B:258:ILE:HD13	1:B:265:LEU:CD1	1.73	1.18
1:B:36:GLN:NE2	1:B:103[B]:ASP:HA	1.62	1.15
1:B:36:GLN:NE2	1:B:103[B]:ASP:HB2	1.61	1.12
1:A:221:MET:HE1	3:A:405:551:CL	1.87	1.11
1:B:258:ILE:HD13	1:B:265:LEU:HD13	1.36	1.05
1:B:36:GLN:HE22	1:B:103[B]:ASP:HB2	1.13	1.03
1:A:221:MET:CE	3:A:405:551:CL	2.44	1.02
1:A:36:GLN:HG2	1:A:104:PRO:HD3	1.38	1.02
1:B:36:GLN:HE21	1:B:103[B]:ASP:HA	0.88	1.01
1:B:224:SER:HB2	1:B:230:GLU:HG2	1.47	0.96
1:A:36:GLN:HE21	1:A:103:ASP:HA	1.32	0.95
1:B:36:GLN:CG	1:B:104:PRO:HD3	1.96	0.95
1:A:33:TRP:CE3	1:A:100:ILE:HG22	2.05	0.91
1:B:258:ILE:HD13	1:B:265:LEU:HD11	1.50	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:GLN:HE21	1:B:103[A]:ASP:HA	0.81	0.89
1:B:162:VAL:HG11	3:B:403:551:CL	2.11	0.87
1:B:36:GLN:NE2	1:B:103[B]:ASP:CA	2.17	0.86
1:A:36:GLN:CG	1:A:104:PRO:HD3	2.05	0.85
1:A:36:GLN:NE2	1:A:103:ASP:HA	1.92	0.85
1:B:258:ILE:CD1	1:B:265:LEU:HD13	2.09	0.83
1:B:36:GLN:HG2	1:B:104:PRO:HD3	1.59	0.82
1:A:36:GLN:NE2	3:A:406:551:C10	2.39	0.81
1:B:10:ARG:NH2	1:B:317:GLU:OE2	2.14	0.81
1:B:17:THR:HG22	5:B:536:HOH:O	1.81	0.80
1:B:162:VAL:CG1	3:B:403:551:CL	2.72	0.75
1:B:36:GLN:HE22	1:B:103[A]:ASP:CG	1.86	0.74
1:A:225:MET:HG3	3:A:405:551:C16	2.18	0.73
1:B:258:ILE:CD1	1:B:265:LEU:CD1	2.61	0.72
1:A:121:PHE:HD2	1:A:125:TYR:CE2	2.08	0.71
1:B:36:GLN:NE2	1:B:104:PRO:CD	2.55	0.70
1:A:36:GLN:HE21	1:A:104:PRO:CD	2.05	0.70
1:A:36:GLN:NE2	1:A:104:PRO:CD	2.56	0.68
1:B:258:ILE:HG21	1:B:265:LEU:HD11	1.74	0.68
1:B:103[A]:ASP:HB3	1:B:106[A]:SER:HB3	1.77	0.66
1:B:36:GLN:HE22	1:B:103[B]:ASP:CB	1.85	0.63
1:A:102:LYS:HD3	1:A:107:ARG:O	2.00	0.61
1:A:36:GLN:NE2	1:A:104:PRO:HD2	2.17	0.59
1:B:225:MET:HG3	3:B:403:551:C16	2.32	0.59
1:A:33:TRP:CZ3	1:A:100:ILE:CG2	2.85	0.59
1:A:33:TRP:CE3	1:A:100:ILE:CG2	2.84	0.58
1:A:36:GLN:HE21	1:A:103:ASP:CA	2.11	0.57
1:B:103[A]:ASP:O	1:B:107:ARG:N	2.36	0.57
1:B:10:ARG:HH22	1:B:317:GLU:CD	2.08	0.57
1:B:103[A]:ASP:OD2	1:B:106[A]:SER:HB2	2.05	0.56
1:A:121:PHE:N	1:A:121:PHE:CD1	2.73	0.56
1:A:33:TRP:CZ3	1:A:100:ILE:HG22	2.42	0.55
1:A:47:ARG:HG2	1:A:52:GLU:HG3	1.90	0.54
1:B:36:GLN:CD	1:B:104:PRO:HD3	2.29	0.53
1:B:36:GLN:NE2	1:B:103[B]:ASP:HB3	2.12	0.53
1:A:23:TYR:HB2	1:B:105:VAL:HG22	1.91	0.53
1:B:36:GLN:NE2	1:B:104:PRO:HD2	2.24	0.52
1:B:36:GLN:NE2	1:B:103[A]:ASP:CA	2.32	0.52
1:B:36:GLN:HE21	1:B:104:PRO:CD	2.22	0.51
1:B:103[A]:ASP:CB	1:B:106[A]:SER:HB3	2.40	0.51
1:B:102:LYS:HD3	1:B:107:ARG:O	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:PRO:HD2	1:B:105:VAL:HG21	1.93	0.50
1:B:36:GLN:HG2	1:B:103[B]:ASP:HA	1.94	0.49
1:B:118:ASN:HD22	1:B:164:ILE:H	1.61	0.49
1:A:36:GLN:HE22	3:A:406:551:C10	2.21	0.48
1:A:221:MET:HE2	3:A:405:551:CL	2.44	0.48
1:A:36:GLN:NE2	1:A:104:PRO:HD3	2.29	0.47
1:B:239:TYR:CZ	1:B:268:ARG:HD3	2.49	0.47
1:A:239:TYR:CZ	1:A:268:ARG:HD3	2.50	0.47
1:B:36:GLN:NE2	1:B:104:PRO:HD3	2.30	0.47
1:B:162:VAL:CB	3:B:403:551:CL	3.02	0.45
1:B:162:VAL:HB	3:B:403:551:CL	2.54	0.45
1:B:33:TRP:NE1	1:B:100:ILE:HG22	2.32	0.45
1:B:36:GLN:CD	1:B:104:PRO:CD	2.86	0.44
1:B:67:VAL:HG22	1:B:112:VAL:HG22	1.99	0.44
1:A:121:PHE:O	1:A:124:LEU:N	2.46	0.43
1:B:102:LYS:HG2	1:B:107:ARG:O	2.19	0.43
1:B:36:GLN:NE2	3:B:404:551:C10	2.79	0.43
1:B:103[B]:ASP:O	1:B:107:ARG:N	2.47	0.42
1:B:36:GLN:HG3	1:B:104:PRO:HD3	1.94	0.42
1:B:5:VAL:HB	1:B:261:TYR:HA	2.01	0.42
1:B:36:GLN:NE2	1:B:103[A]:ASP:CG	2.46	0.41
1:B:103[A]:ASP:HB3	1:B:106[A]:SER:CB	2.48	0.41
1:B:106[A]:SER:O	1:B:107:ARG:HB2	2.20	0.41
1:A:121:PHE:HD2	1:A:125:TYR:CZ	2.38	0.41
1:A:96:THR:H	1:A:114[A]:GLU:HG2	1.86	0.41
1:A:5:VAL:HB	1:A:261:TYR:HA	2.03	0.41
1:A:258:ILE:HG23	1:A:263:ILE:O	2.21	0.40
1:A:124:LEU:HD23	1:A:127:THR:OG1	2.22	0.40
1:B:95:ILE:HB	1:B:174:ILE:HG22	2.03	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	333/352 (95%)	326 (98%)	7 (2%)	0	100	100
1	B	327/352 (93%)	320 (98%)	7 (2%)	0	100	100
All	All	660/704 (94%)	646 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	309/319 (97%)	308 (100%)	1 (0%)	94	91
1	B	303/319 (95%)	303 (100%)	0	100	100
All	All	612/638 (96%)	611 (100%)	1 (0%)	95	92

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	118	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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

12 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$
2	ACT	A	401	-	0,3,3	0.00	-	0,3,3	0.00	-
2	ACT	A	402	-	0,3,3	0.00	-	0,3,3	0.00	-
2	ACT	A	403	-	0,3,3	0.00	-	0,3,3	0.00	-
2	ACT	A	404	-	0,3,3	0.00	-	0,3,3	0.00	-
3	551	A	405	-	27,27,27	0.18	0	32,34,34	0.19	0
3	551	A	406	1	16,16,27	0.20	0	20,21,34	0.24	0
3	551	A	407	-	16,16,27	0.18	0	20,21,34	0.25	0
4	PO4	A	408	-	4,4,4	1.88	2 (50%)	6,6,6	0.26	0
2	ACT	B	401	-	0,3,3	0.00	-	0,3,3	0.00	-
2	ACT	B	402	-	0,3,3	0.00	-	0,3,3	0.00	-
3	551	B	403	-	27,27,27	0.19	0	32,34,34	0.20	0
3	551	B	404	-	16,16,27	0.19	0	20,21,34	0.22	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
2	ACT	A	401	-	-	0/0/0/0	0/0/0/0
2	ACT	A	402	-	-	0/0/0/0	0/0/0/0
2	ACT	A	403	-	-	0/0/0/0	0/0/0/0
2	ACT	A	404	-	-	0/0/0/0	0/0/0/0
3	551	A	405	-	-	0/19/19/19	0/2/2/2
3	551	A	406	1	-	0/6/6/19	0/2/2/2
3	551	A	407	-	-	0/6/6/19	0/2/2/2
4	PO4	A	408	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACT	B	401	-	-	0/0/0/0	0/0/0/0
2	ACT	B	402	-	-	0/0/0/0	0/0/0/0
3	551	B	403	-	-	0/19/19/19	0/2/2/2
3	551	B	404	-	-	0/6/6/19	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	408	PO4	P-O4	2.16	1.60	1.53
4	A	408	PO4	P-O3	2.33	1.61	1.53

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	405	551	4	0
3	A	406	551	2	0
3	B	403	551	5	0
3	B	404	551	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	328/352 (93%)	0.28	16 (4%) 33 38	15, 27, 58, 101	0
1	B	325/352 (92%)	0.81	44 (13%) 4 5	25, 54, 90, 109	0
All	All	653/704 (92%)	0.55	60 (9%) 11 13	15, 39, 83, 109	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	TYR	13.0
1	B	50	TYR	10.3
1	B	122	LYS	8.7
1	A	124	LEU	7.2
1	A	121	PHE	7.0
1	A	125	TYR	7.0
1	B	105	VAL	6.6
1	B	123	GLN	6.5
1	A	123	GLN	6.5
1	B	121	PHE	6.1
1	B	126	GLN	5.5
1	B	49	LYS	5.4
1	B	124	LEU	5.4
1	A	122	LYS	5.2
1	B	104	PRO	5.0
1	A	50	TYR	5.0
1	B	268	ARG	5.0
1	B	48	GLY	4.9
1	B	270	ASN	4.6
1	B	288	GLU	4.4
1	B	127	THR	4.3
1	B	33	TRP	4.0
1	A	49	LYS	4.0
1	A	2	SER	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	33	TRP	3.6
1	B	255	TYR	3.4
1	B	47	ARG	3.3
1	A	126	GLN	3.3
1	A	104	PRO	3.2
1	A	105	VAL	3.2
1	A	329	LYS	3.2
1	B	107	ARG	3.2
1	B	257	TYR	3.1
1	B	36	GLN	3.0
1	B	271	ASP	3.0
1	B	3	GLY	2.9
1	B	59	ILE	2.9
1	B	280	ARG	2.8
1	B	252	GLU	2.7
1	A	107	ARG	2.7
1	B	272	ILE	2.6
1	B	273	LEU	2.5
1	B	284	PHE	2.4
1	B	269	PHE	2.3
1	B	265	LEU	2.3
1	B	251	THR	2.2
1	B	227	PHE	2.2
1	B	267	PRO	2.2
1	B	239	TYR	2.2
1	B	57	ILE	2.2
1	B	117	ASN	2.2
1	B	307	TYR	2.2
1	B	21	SER	2.2
1	B	73	VAL	2.2
1	B	264	GLU	2.1
1	B	166	HIS	2.1
1	A	117	ASN	2.1
1	B	40[A]	GLN	2.0
1	B	327	VAL	2.0
1	A	35	ASN	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	551	A	406	15/26	0.57	0.32	3.69	47,56,64,65	0
2	ACT	B	402	4/4	0.90	0.14	3.53	35,38,42,44	0
3	551	B	404	15/26	0.47	0.37	3.02	32,32,33,33	15
3	551	B	403	26/26	0.65	0.22	1.50	56,75,80,87	0
3	551	A	407	15/26	0.90	0.13	1.33	63,64,66,69	0
3	551	A	405	26/26	0.58	0.22	1.27	55,57,73,74	0
2	ACT	B	401	4/4	0.92	0.16	1.01	47,49,49,49	0
2	ACT	A	402	4/4	0.97	0.09	-0.38	30,33,34,35	0
2	ACT	A	401	4/4	0.95	0.09	-0.89	26,27,30,30	0
4	PO4	A	408	5/5	0.95	0.07	-1.70	43,44,50,54	0
2	ACT	A	404	4/4	0.60	0.33	-	53,55,55,55	0
2	ACT	A	403	4/4	0.94	0.17	-	42,42,43,46	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.