



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:44 AM GMT

PDB ID : 3C0G
Title : CASK CaM-Kinase Domain- 3'-AMP complex, P1 form
Authors : Wahl, M.C.
Deposited on : 2008-01-20
Resolution : 2.19 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

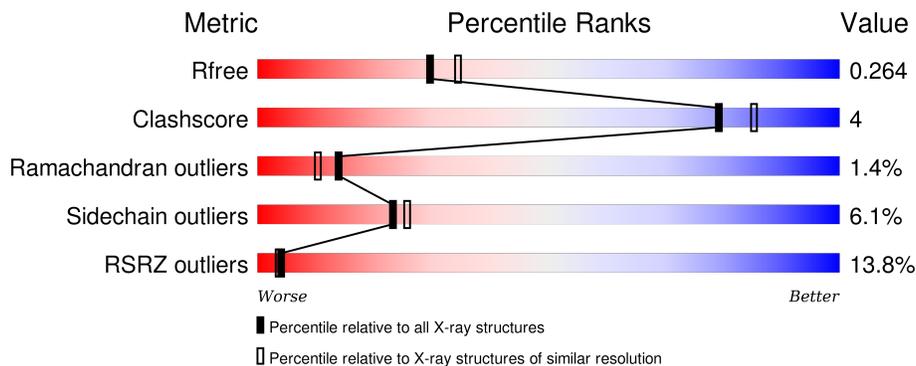
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 ≥ 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	351	
1	B	351	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5363 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 Peripheral plasma membrane protein CASK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	309	2500	1609	436	438	17	0	4	0
1	B	320	2563	1646	446	454	17	0	2	0

There are 28 discrepancies between the modelled and reference sequences:

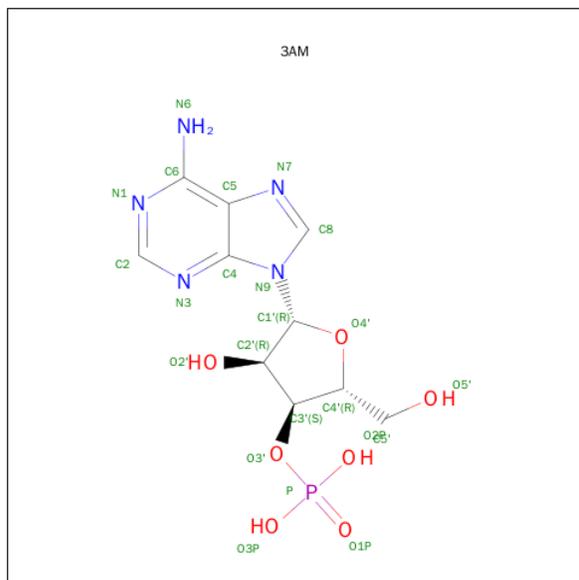
Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	GLY	-	EXPRESSION TAG	UNP O14936
A	-12	SER	-	EXPRESSION TAG	UNP O14936
A	-11	PRO	-	EXPRESSION TAG	UNP O14936
A	-10	GLY	-	EXPRESSION TAG	UNP O14936
A	-9	ILE	-	EXPRESSION TAG	UNP O14936
A	-8	SER	-	EXPRESSION TAG	UNP O14936
A	-7	GLY	-	EXPRESSION TAG	UNP O14936
A	-6	GLY	-	EXPRESSION TAG	UNP O14936
A	-5	GLY	-	EXPRESSION TAG	UNP O14936
A	-4	GLY	-	EXPRESSION TAG	UNP O14936
A	-3	GLY	-	EXPRESSION TAG	UNP O14936
A	-2	ILE	-	EXPRESSION TAG	UNP O14936
A	-1	LEU	-	EXPRESSION TAG	UNP O14936
A	0	ASP	-	EXPRESSION TAG	UNP O14936
B	-13	GLY	-	EXPRESSION TAG	UNP O14936
B	-12	SER	-	EXPRESSION TAG	UNP O14936
B	-11	PRO	-	EXPRESSION TAG	UNP O14936
B	-10	GLY	-	EXPRESSION TAG	UNP O14936
B	-9	ILE	-	EXPRESSION TAG	UNP O14936
B	-8	SER	-	EXPRESSION TAG	UNP O14936
B	-7	GLY	-	EXPRESSION TAG	UNP O14936
B	-6	GLY	-	EXPRESSION TAG	UNP O14936
B	-5	GLY	-	EXPRESSION TAG	UNP O14936
B	-4	GLY	-	EXPRESSION TAG	UNP O14936
B	-3	GLY	-	EXPRESSION TAG	UNP O14936

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	ILE	-	EXPRESSION TAG	UNP O14936
B	-1	LEU	-	EXPRESSION TAG	UNP O14936
B	0	ASP	-	EXPRESSION TAG	UNP O14936

- Molecule 2 is [(2R,3S,4R,5R)-5-(6-AMINOPURIN-9-YL)-4-HYDROXY-2-(HYDROXYMETHYL)OXOLAN-3-YL] DIHYDROGEN PHOSPHATE (three-letter code: 3AM) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

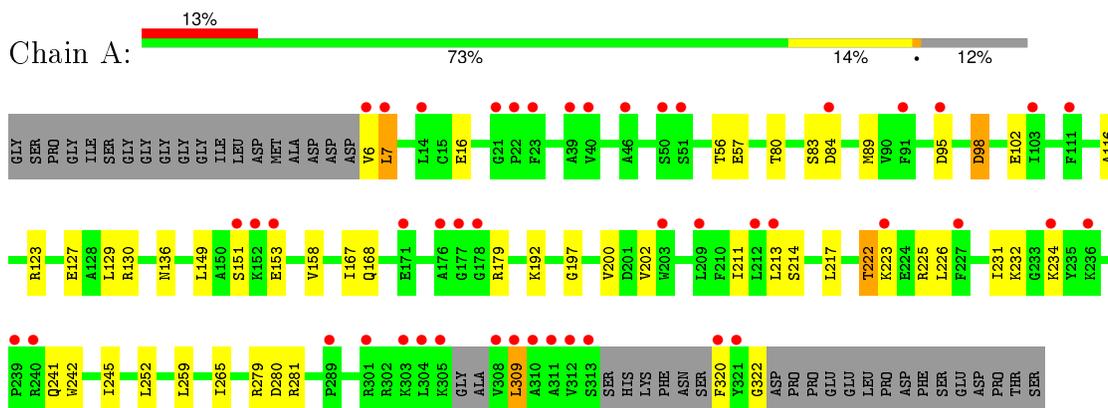
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	133	Total	O	0	1
			134	134		
3	B	119	Total	O	0	1
			120	120		

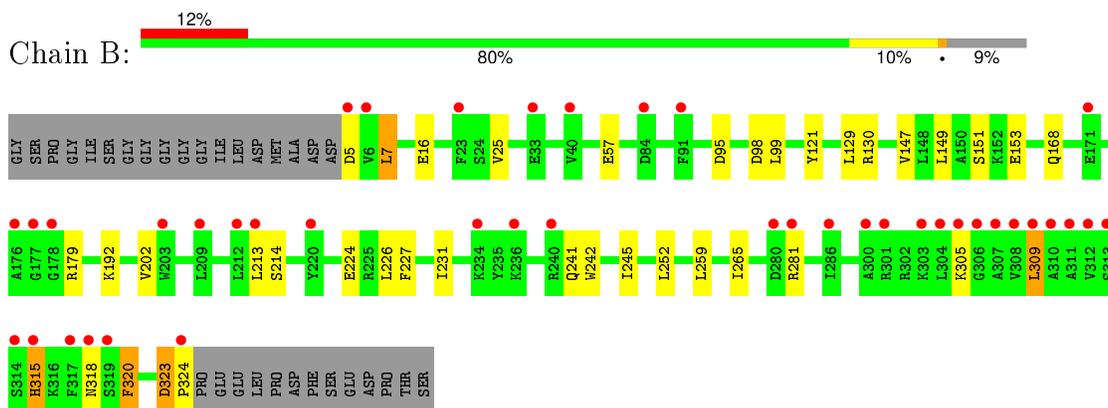
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Peripheral plasma membrane protein CASK



- Molecule 1: Peripheral plasma membrane protein CASK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.33Å 60.27Å 60.55Å 90.00° 106.49° 105.80°	Depositor
Resolution (Å)	20.00 – 2.19 19.90 – 2.19	Depositor EDS
% Data completeness (in resolution range)	86.3 (20.00-2.19) 84.5 (19.90-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.208 , 0.270 0.204 , 0.264	Depositor DCC
R_{free} test set	1734 reflections (5.70%)	DCC
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.555	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.8	EDS
Estimated twinning fraction	0.017 for -h,-l,-k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 34845 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5363	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.00% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 3AM

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	0.39	0/2557	0.55	1/3445 (0.0%)
1	B	0.39	0/2624	0.57	2/3540 (0.1%)
All	All	0.39	0/5181	0.56	3/6985 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	7	LEU	CA-CB-CG	6.06	129.23	115.30
1	B	309	LEU	CA-CB-CG	5.76	128.54	115.30
1	A	309	LEU	CA-CB-CG	5.48	127.90	115.30

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	A	2500	0	2527	24	0
1	B	2563	0	2579	12	0
2	A	23	0	12	0	0
2	B	23	0	12	0	0
3	A	134	0	0	2	0
3	B	120	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5363	0	5130	36	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 (36) 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:323:ASP:HB2	1:B:324:PRO:HD3	1.53	0.88
1:A:179:ARG:HD2	1:A:320:PHE:HB2	1.59	0.82
1:A:232:LYS:HB3	1:A:234:LYS:HG2	1.68	0.74
1:A:223[A]:LYS:HA	1:A:322:GLY:HA3	1.75	0.68
1:A:223[B]:LYS:HA	1:A:322:GLY:HA3	1.77	0.66
1:B:99:LEU:HD22	1:B:147:VAL:HG11	1.81	0.63
1:A:102:GLU:HG3	1:A:149:LEU:HD12	1.83	0.61
1:A:95:ASP:HB3	1:A:151:SER:HA	1.81	0.60
1:A:222:THR:HG22	1:A:225:ARG:HB2	1.84	0.59
1:A:214:SER:HB2	1:A:241:GLN:HG3	1.88	0.55
1:B:179:ARG:HD2	1:B:320:PHE:HB2	1.90	0.53
1:A:202:VAL:HG11	1:A:265:ILE:O	2.09	0.52
1:B:224:GLU:HA	1:B:227:PHE:HB3	1.91	0.51
1:B:323:ASP:HB2	1:B:324:PRO:CD	2.35	0.48
1:A:192:LYS:HE3	1:A:231:ILE:HG12	1.96	0.48
1:A:179:ARG:CD	1:A:320:PHE:HB2	2.37	0.48
1:B:242:TRP:HA	1:B:245:ILE:HD12	1.95	0.47
1:A:242:TRP:HA	1:A:245:ILE:HD12	1.97	0.47
1:B:99:LEU:HD11	1:B:121:TYR:CD1	2.50	0.46
1:A:102:GLU:CG	1:A:149:LEU:HD12	2.45	0.46
1:B:95:ASP:HB3	1:B:151:SER:HA	1.96	0.46
1:A:116:ALA:O	1:A:279:ARG:HD3	2.15	0.45
1:A:123:ARG:O	1:A:127:GLU:HG3	2.17	0.45
1:B:192:LYS:HE3	1:B:231:ILE:HG12	1.98	0.44
1:B:213:LEU:O	1:B:245:ILE:HD11	2.16	0.44
1:B:202:VAL:HG11	1:B:265:ILE:O	2.18	0.43
1:A:232:LYS:HB3	1:A:234:LYS:CG	2.43	0.43
1:A:98:ASP:C	1:A:98:ASP:OD2	2.57	0.43
1:A:213:LEU:O	1:A:245:ILE:HD11	2.20	0.42
1:A:211:ILE:HD11	1:A:217:LEU:HD23	2.02	0.41
1:A:197:GLY:O	1:A:200:VAL:HG22	2.20	0.41
1:A:7:LEU:HD22	1:A:83:SER:HB3	2.03	0.41
1:A:280:ASP:HB2	3:A:395:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:THR:HG22	1:A:89:MET:HG2	2.03	0.41
1:A:136:ASN:ND2	3:A:421:HOH:O	2.54	0.41
1:B:214:SER:HB3	1:B:242:TRP:CE2	2.56	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	307/351 (88%)	297 (97%)	7 (2%)	3 (1%)	19 16
1	B	320/351 (91%)	303 (95%)	11 (3%)	6 (2%)	10 6
All	All	627/702 (89%)	600 (96%)	18 (3%)	9 (1%)	14 10

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	222	THR
1	B	305	LYS
1	B	315	HIS
1	B	323	ASP
1	B	7	LEU
1	B	318	ASN
1	B	320	PHE
1	A	84	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	269/297 (91%)	253 (94%)	16 (6%)	24	27
1	B	276/297 (93%)	259 (94%)	17 (6%)	23	25
All	All	545/594 (92%)	512 (94%)	33 (6%)	23	26

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	VAL
1	A	16	GLU
1	A	56	THR
1	A	57	GLU
1	A	98	ASP
1	A	129	LEU
1	A	130	ARG
1	A	153	GLU
1	A	158	VAL
1	A	167	ILE
1	A	168	GLN
1	A	226	LEU
1	A	252	LEU
1	A	259	LEU
1	A	281	ARG
1	A	309	LEU
1	B	5	ASP
1	B	16	GLU
1	B	25	VAL
1	B	57	GLU
1	B	98	ASP
1	B	129	LEU
1	B	130	ARG
1	B	149	LEU
1	B	153	GLU
1	B	168	GLN
1	B	226	LEU
1	B	241	GLN
1	B	252	LEU
1	B	259	LEU
1	B	281	ARG
1	B	309	LEU
1	B	315	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	37	GLN
1	A	168	GLN
1	B	36	GLN
1	B	168	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	3AM	A	338	-	20,25,25	1.23	2 (10%)	22,38,38	2.69	1 (4%)
2	3AM	B	338	-	20,25,25	1.20	2 (10%)	22,38,38	2.60	1 (4%)

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	3AM	A	338	-	-	0/7/27/27	0/3/3/3
2	3AM	B	338	-	-	0/7/27/27	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	338	3AM	C2-N1	2.59	1.38	1.33
2	A	338	3AM	C2-N1	2.65	1.38	1.33
2	B	338	3AM	C2-N3	3.68	1.38	1.32
2	A	338	3AM	C2-N3	3.87	1.39	1.32

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	338	3AM	N3-C2-N1	-12.07	119.66	128.89
2	B	338	3AM	N3-C2-N1	-11.58	120.02	128.89

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	309/351 (88%)	0.88	46 (14%) 3 3	32, 45, 69, 128	1 (0%)
1	B	320/351 (91%)	0.82	41 (12%) 5 4	32, 46, 88, 107	0
All	All	629/702 (89%)	0.85	87 (13%) 4 3	32, 45, 76, 128	1 (0%)

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	308	VAL	10.6
1	B	306	GLY	9.2
1	A	6	VAL	9.1
1	B	308	VAL	8.8
1	A	304	LEU	8.5
1	A	311	ALA	8.3
1	B	6	VAL	8.0
1	B	307	ALA	7.8
1	B	311	ALA	7.6
1	A	309	LEU	7.5
1	A	312	VAL	7.2
1	A	178	GLY	7.0
1	B	312	VAL	6.9
1	A	305	LYS	6.7
1	A	313	SER	6.6
1	A	310	ALA	6.3
1	B	305	LYS	5.9
1	B	319	SER	5.8
1	B	240	ARG	5.2
1	B	317	PHE	4.7
1	B	178	GLY	4.6
1	A	176	ALA	4.6
1	A	7	LEU	4.4
1	B	314	SER	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	324	PRO	4.2
1	A	321	TYR	4.1
1	B	315	HIS	4.0
1	B	280	ASP	3.9
1	A	40	VAL	3.8
1	B	176	ALA	3.8
1	A	23	PHE	3.7
1	B	304	LEU	3.7
1	A	236	LYS	3.7
1	A	84	ASP	3.6
1	B	309	LEU	3.6
1	A	22	PRO	3.4
1	B	212[A]	LEU	3.4
1	A	50	SER	3.3
1	A	213	LEU	3.3
1	A	227	PHE	3.2
1	A	320	PHE	3.2
1	B	301	ARG	3.2
1	B	209	LEU	3.2
1	B	40	VAL	3.1
1	B	318	ASN	3.0
1	B	220	TYR	3.0
1	B	313	SER	2.9
1	A	152	LYS	2.9
1	A	171	GLU	2.9
1	B	203	TRP	2.8
1	A	95	ASP	2.8
1	B	310	ALA	2.8
1	A	51	SER	2.7
1	A	223[A]	LYS	2.7
1	A	209	LEU	2.6
1	A	46	ALA	2.6
1	B	5	ASP	2.6
1	A	177	GLY	2.6
1	B	300	ALA	2.6
1	A	212[A]	LEU	2.6
1	B	177	GLY	2.6
1	A	234	LYS	2.6
1	B	91	PHE	2.5
1	B	171	GLU	2.5
1	B	213	LEU	2.5
1	A	239	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	240	ARG	2.4
1	A	151	SER	2.4
1	B	234	LYS	2.4
1	A	303	LYS	2.4
1	A	91	PHE	2.4
1	B	23	PHE	2.3
1	B	33	GLU	2.3
1	B	84	ASP	2.3
1	A	111	PHE	2.3
1	A	203	TRP	2.3
1	A	289	PRO	2.3
1	A	21	GLY	2.2
1	A	39	ALA	2.2
1	A	153	GLU	2.2
1	A	14	LEU	2.2
1	B	303	LYS	2.2
1	A	103	ILE	2.1
1	B	236	LYS	2.1
1	A	301	ARG	2.1
1	B	281	ARG	2.1
1	B	286	ILE	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 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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	3AM	A	338	23/23	0.89	0.17	0.73	62,66,81,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	3AM	B	338	23/23	0.94	0.16	0.55	45,53,71,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.