



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 09:40 PM GMT

PDB ID : 2PGJ
Title : Catalysis associated conformational changes revealed by human cd38 complexed with a non-hydrolyzable substrate analog
Authors : Liu, Q.; Kriksunov, I.A.; Moreau, C.; Graeff, R.; Potter, B.V.L.; Lee, H.C.; Hao, Q.
Deposited on : 2007-04-09
Resolution : 1.71 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

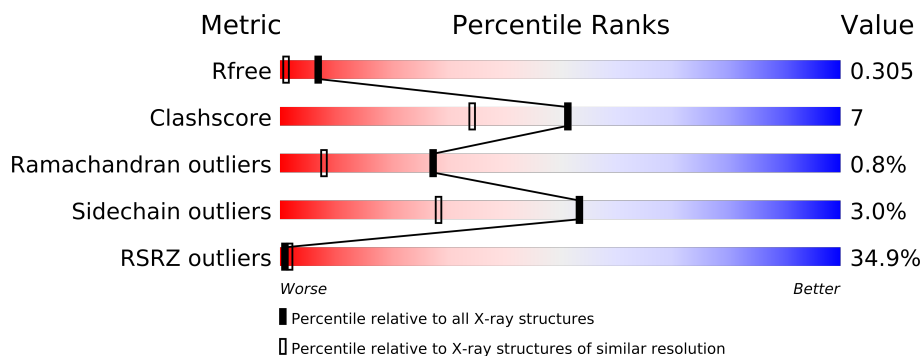
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.71 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}	66092	2979 (1.74-1.70)
Clashscore	79885	3506 (1.74-1.70)
Ramachandran outliers	78287	3449 (1.74-1.70)
Sidechain outliers	78261	3449 (1.74-1.70)
RSRZ outliers	66119	2979 (1.74-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	262	
1	B	262	

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 4535 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

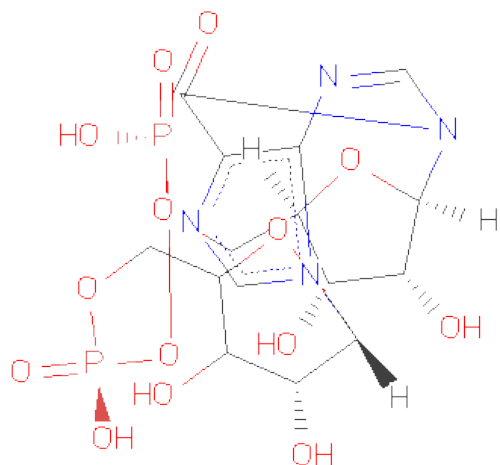
- Molecule 1 is a protein called ADP-ribosyl cyclase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			2050	1290	358	386	16			
1	B	252	Total	C	N	O	S	0	0	0
			2050	1290	358	386	16			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	LYS	-	CLONING ARTIFACT	UNP P28907
A	40	ARG	-	CLONING ARTIFACT	UNP P28907
A	41	GLU	-	CLONING ARTIFACT	UNP P28907
A	42	ALA	-	CLONING ARTIFACT	UNP P28907
A	43	GLU	-	CLONING ARTIFACT	UNP P28907
A	44	ALA	-	CLONING ARTIFACT	UNP P28907
A	49	THR	GLN	ENGINEERED	UNP P28907
A	100	ASP	ASN	ENGINEERED	UNP P28907
A	164	ASP	ASN	ENGINEERED	UNP P28907
A	209	ASP	ASN	ENGINEERED	UNP P28907
A	219	ASP	ASN	ENGINEERED	UNP P28907
B	39	LYS	-	CLONING ARTIFACT	UNP P28907
B	40	ARG	-	CLONING ARTIFACT	UNP P28907
B	41	GLU	-	CLONING ARTIFACT	UNP P28907
B	42	ALA	-	CLONING ARTIFACT	UNP P28907
B	43	GLU	-	CLONING ARTIFACT	UNP P28907
B	44	ALA	-	CLONING ARTIFACT	UNP P28907
B	49	THR	GLN	ENGINEERED	UNP P28907
B	100	ASP	ASN	ENGINEERED	UNP P28907
B	164	ASP	ASN	ENGINEERED	UNP P28907
B	209	ASP	ASN	ENGINEERED	UNP P28907
B	219	ASP	ASN	ENGINEERED	UNP P28907

- Molecule 2 is N1-CYCLIC INOSINE 5'-DIPHOSPHORIBOSE (three-letter code: N1C) (formula: $C_{15}H_{20}N_4O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			35	15	4	14	2		

- Molecule 3 is water.

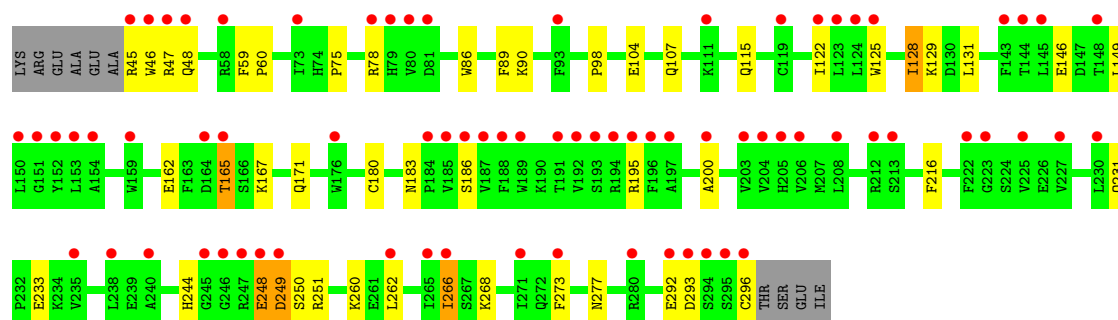
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	210	Total	O	0	0
			210	210		
3	B	190	Total	O	0	0
			190	190		

3 Residue-property plots

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.

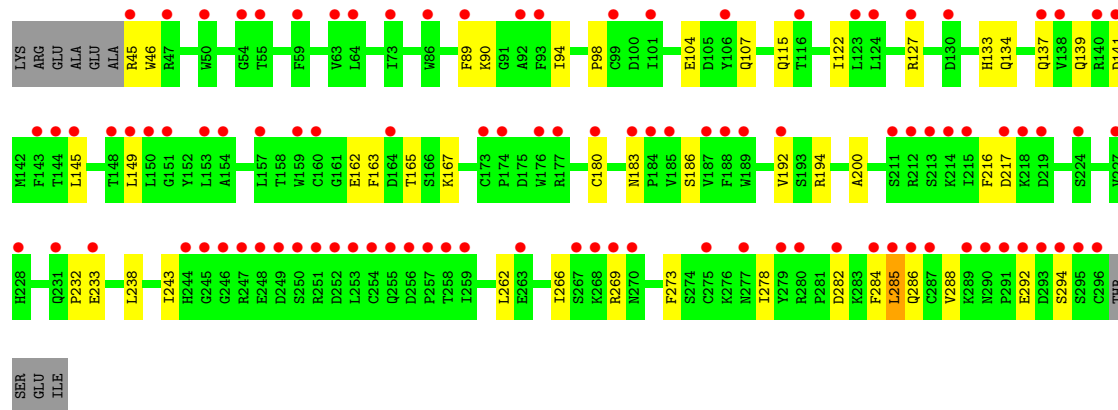
• Molecule 1: ADP-ribosyl cyclase 1

Chain A: 



• Molecule 1: ADP-ribosyl cyclase 1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	41.89Å 53.20Å 65.64Å 105.34° 91.96° 94.94°	Depositor
Resolution (Å)	28.41 – 1.71 28.41 – 1.71	Depositor EDS
% Data completeness (in resolution range)	100.0 (28.41-1.71) 91.9 (28.41-1.71)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.21 (at 1.71Å)	Xtriage
Refinement program	REFMAC 5.3.0021	Depositor
R, R_{free}	0.191 , 0.224 0.281 , 0.305	Depositor DCC
R_{free} test set	2762 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 57.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 54284 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4535	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: N1C

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	0.44	0/2101	0.57	0/2846
1	B	0.46	0/2101	0.55	0/2846
All	All	0.45	0/4202	0.56	0/5692

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2050	0	1976	34	0
1	B	2050	0	1976	26	0
2	A	35	0	15	3	0
3	A	210	0	0	5	0
3	B	190	0	0	1	0
All	All	4535	0	3967	59	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (59) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:115:GLN:HE22	1:B:149:LEU:H	1.20	0.88
1:A:115:GLN:HE22	1:A:149:LEU:H	1.23	0.86
1:B:165:THR:HG23	1:B:167:LYS:H	1.46	0.80
1:A:268:LYS:HD3	1:B:163:PHE:CE1	2.17	0.79
1:A:165:THR:HG23	1:A:167:LYS:H	1.47	0.78
1:A:268:LYS:HD3	1:B:163:PHE:HE1	1.49	0.77
2:A:301:N1C:H2D	2:A:301:N1C:O6	1.94	0.67
1:B:162:GLU:HB2	1:B:165:THR:HG22	1.79	0.64
1:A:180:CYS:HB2	3:A:432:HOH:O	2.00	0.61
1:A:183:ASN:ND2	1:A:186:SER:H	2.00	0.60
1:A:162:GLU:HB2	1:A:165:THR:HG22	1.84	0.59
1:A:266:ILE:HD11	1:A:273:PHE:HB2	1.86	0.57
1:A:75:PRO:O	1:A:78:ARG:HB2	2.03	0.57
1:B:266:ILE:HD11	1:B:273:PHE:HB2	1.86	0.57
1:B:127:ARG:HH12	1:B:217:ASP:HB3	1.68	0.57
1:A:104:GLU:HA	1:A:107:GLN:HG2	1.87	0.57
1:B:104:GLU:HA	1:B:107:GLN:HG2	1.86	0.56
1:B:133:HIS:O	1:B:137:GLN:HG3	2.06	0.56
1:A:231:GLN:HG3	3:A:510:HOH:O	2.05	0.56
1:A:115:GLN:NE2	1:A:149:LEU:H	2.01	0.56
1:A:146:GLU:OE2	2:A:301:N1C:O6	2.24	0.55
1:B:238:LEU:HB3	1:B:266:ILE:HD13	1.89	0.54
1:A:183:ASN:HD21	1:A:186:SER:H	1.54	0.54
1:B:115:GLN:NE2	1:B:149:LEU:H	1.98	0.53
1:B:284:PHE:O	1:B:288:VAL:HG23	2.08	0.53
1:A:293:ASP:HB3	1:A:296:CYS:SG	2.50	0.52
1:B:139:GLN:C	1:B:141:ASP:H	2.13	0.51
1:A:195:ARG:HG3	3:A:326:HOH:O	2.11	0.50
1:B:183:ASN:ND2	1:B:186:SER:H	2.10	0.49
1:B:145:LEU:HD21	1:B:192:VAL:HG23	1.95	0.49
1:A:216:PHE:CE1	1:A:262:LEU:HD13	2.48	0.49
1:A:48:GLN:NE2	1:A:171:GLN:HB3	2.28	0.48
1:B:45:ARG:HG2	1:B:46:TRP:H	1.79	0.48
1:B:134:GLN:HB3	1:B:285:LEU:HD11	1.96	0.47
1:A:86:TRP:CE2	1:A:90:LYS:HG3	2.49	0.47
1:B:90:LYS:HG3	1:B:94:ILE:HG13	1.96	0.45
1:A:125:TRP:CH2	1:A:129:LYS:HB3	2.51	0.45
1:A:125:TRP:CZ3	1:A:129:LYS:HB3	2.52	0.44
1:A:260:LYS:NZ	3:A:446:HOH:O	2.49	0.44
1:A:125:TRP:O	2:A:301:N1C:H5D1	2.16	0.44
1:A:59:PHE:HB3	1:A:60:PRO:HD3	2.00	0.44
1:B:122:ILE:HD12	1:B:200:ALA:HA	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:78:ARG:NH1	3:A:382:HOH:O	2.51	0.43
1:B:216:PHE:CE1	1:B:262:LEU:HD13	2.53	0.43
1:B:232:PRO:HG3	1:B:269:ARG:O	2.18	0.43
1:B:180:CYS:HB2	3:B:324:HOH:O	2.18	0.43
1:A:46:TRP:HE1	1:A:47:ARG:NH1	2.17	0.42
1:A:162:GLU:OE2	1:A:165:THR:HG21	2.19	0.42
1:B:233:GLU:CD	1:B:233:GLU:H	2.22	0.42
1:A:98:PRO:O	1:A:183:ASN:HA	2.20	0.42
1:B:98:PRO:O	1:B:183:ASN:HA	2.18	0.42
1:A:248:GLU:O	1:A:249:ASP:CB	2.67	0.42
1:B:243:ILE:CD1	1:B:278:ILE:HD12	2.50	0.42
1:A:233:GLU:H	1:A:233:GLU:CD	2.22	0.42
1:A:244:HIS:HD2	1:A:250:SER:HB3	1.85	0.41
1:A:122:ILE:HD12	1:A:200:ALA:HA	2.03	0.41
1:A:248:GLU:O	1:A:249:ASP:HB2	2.20	0.41
1:B:139:GLN:C	1:B:141:ASP:N	2.74	0.40
1:A:244:HIS:HE1	1:A:277:ASN:OD1	2.04	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	250/262 (95%)	239 (96%)	8 (3%)	3 (1%)	19	4
1	B	250/262 (95%)	237 (95%)	12 (5%)	1 (0%)	43	22
All	All	500/524 (95%)	476 (95%)	20 (4%)	4 (1%)	27	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	ASP
1	A	248	GLU
1	B	292	GLU

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Mol	Chain	Res	Type
1	A	128	ILE

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	233/241 (97%)	225 (97%)	8 (3%)	49	23
1	B	233/241 (97%)	227 (97%)	6 (3%)	59	35
All	All	466/482 (97%)	452 (97%)	14 (3%)	53	28

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ARG
1	A	89	PHE
1	A	128	ILE
1	A	131	LEU
1	A	165	THR
1	A	251	ARG
1	A	266	ILE
1	A	292	GLU
1	B	89	PHE
1	B	194	ARG
1	B	282	ASP
1	B	285	LEU
1	B	286	GLN
1	B	294	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	GLN
1	A	115	GLN
1	A	134	GLN
1	A	139	GLN
1	A	183	ASN

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Mol	Chain	Res	Type
1	A	244	HIS
1	A	290	ASN
1	B	115	GLN
1	B	137	GLN
1	B	171	GLN
1	B	183	ASN
1	B	244	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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	N1C	A	301	-	39,39,39	1.81	6 (15%)	60,62,62	3.50	18 (30%)

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	N1C	A	301	-	2/2/10/10	0/24/58/58	0/0/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	N1C	C3'-C4'	-7.52	1.32	1.53
2	A	301	N1C	C6-C5	4.52	1.48	1.41
2	A	301	N1C	O3'-C3'	-3.92	1.33	1.43
2	A	301	N1C	C2-N3	2.68	1.34	1.30
2	A	301	N1C	C6-N1	2.49	1.42	1.35
2	A	301	N1C	C4-N9	-2.32	1.34	1.37

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	N1C	C6-C5-N7	-22.28	133.05	134.24
2	A	301	N1C	O3'-C3'-C4'	4.78	125.17	111.08
2	A	301	N1C	N3-C4-N9	4.44	133.44	125.43
2	A	301	N1C	O4'-C1'-N9	-4.31	104.43	108.44
2	A	301	N1C	C2'-C3'-C4'	4.13	110.89	102.65
2	A	301	N1C	C3D-C2D-C1D	4.06	107.26	100.91
2	A	301	N1C	O4'-C4'-C5'	3.92	123.35	109.36
2	A	301	N1C	O3'-C3'-C2'	3.71	123.91	111.83
2	A	301	N1C	PA-O3A-PB	-3.48	121.48	131.68
2	A	301	N1C	C2-N3-C4	3.35	121.95	116.23
2	A	301	N1C	C4'-O4'-C1'	-2.98	106.51	109.75
2	A	301	N1C	O4'-C4'-C3'	2.88	111.00	105.17
2	A	301	N1C	C5-C4-N3	-2.71	119.79	125.70
2	A	301	N1C	C4D-O4D-C1D	2.61	112.58	109.75
2	A	301	N1C	C5'-C4'-C3'	2.61	125.65	115.21
2	A	301	N1C	C5-C6-N1	-2.55	117.84	120.16
2	A	301	N1C	C8-N9-C4	2.51	108.81	106.90
2	A	301	N1C	O4D-C1D-C2D	-2.50	102.94	106.77

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	301	N1C	C4'
2	A	301	N1C	C3'

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	252/262 (96%)	1.70	75 (29%) 1 2	23, 29, 36, 43	0
1	B	252/262 (96%)	2.07	101 (40%) 1 1	23, 29, 35, 40	0
All	All	504/524 (96%)	1.88	176 (34%) 1 2	23, 29, 36, 43	0

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	CYS	10.8
1	A	246	GLY	10.8
1	B	246	GLY	10.6
1	B	213	SER	9.2
1	B	248	GLU	9.1
1	B	295	SER	8.9
1	B	245	GLY	8.9
1	A	249	ASP	8.7
1	B	291	PRO	8.7
1	B	249	ASP	8.3
1	B	247	ARG	7.6
1	B	292	GLU	7.6
1	B	287	CYS	7.0
1	B	290	ASN	6.6
1	A	296	CYS	6.2
1	A	292	GLU	6.1
1	A	47	ARG	6.1
1	A	247	ARG	6.1
1	A	294	SER	6.1
1	A	295	SER	6.0
1	B	294	SER	5.9
1	B	293	ASP	5.6
1	B	212	ARG	5.4
1	A	248	GLU	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	143	PHE	5.2
1	A	145	LEU	5.1
1	B	159	TRP	5.1
1	B	254	CYS	5.0
1	A	124	LEU	5.0
1	A	189	TRP	4.7
1	B	219	ASP	4.7
1	A	188	PHE	4.6
1	B	141	ASP	4.5
1	A	123	LEU	4.5
1	A	222	PHE	4.5
1	A	46	TRP	4.5
1	A	238	LEU	4.5
1	B	153	LEU	4.5
1	B	270	ASN	4.4
1	A	230	LEU	4.4
1	B	188	PHE	4.4
1	B	286	GLN	4.3
1	A	213	SER	4.3
1	B	185	VAL	4.2
1	B	214	LYS	4.2
1	B	211	SER	4.1
1	A	187	VAL	4.0
1	B	257	PRO	4.0
1	B	215	ILE	4.0
1	A	192	VAL	4.0
1	A	206	VAL	3.9
1	B	176	TRP	3.8
1	A	204	VAL	3.7
1	B	140	ARG	3.7
1	B	252	ASP	3.7
1	A	266	ILE	3.7
1	B	259	ILE	3.7
1	A	79	HIS	3.7
1	A	164	ASP	3.6
1	A	212	ARG	3.6
1	B	157	LEU	3.6
1	A	245	GLY	3.6
1	B	184	PRO	3.6
1	B	160	CYS	3.6
1	A	154	ALA	3.6
1	B	116	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	165	THR	3.5
1	A	227	VAL	3.5
1	B	154	ALA	3.5
1	B	127	ARG	3.5
1	B	189	TRP	3.4
1	B	279	TYR	3.4
1	B	285	LEU	3.4
1	A	196	PHE	3.4
1	B	282	ASP	3.4
1	B	89	PHE	3.3
1	A	176	TRP	3.3
1	B	268	LYS	3.3
1	A	125	TRP	3.3
1	B	145	LEU	3.3
1	B	218	LYS	3.3
1	B	138	VAL	3.3
1	A	122	ILE	3.3
1	A	185	VAL	3.3
1	A	235	VAL	3.2
1	A	159	TRP	3.2
1	A	150	LEU	3.2
1	B	73	ILE	3.2
1	A	293	ASP	3.2
1	A	203	VAL	3.2
1	B	255	GLN	3.1
1	A	197	ALA	3.1
1	B	231	GLN	3.1
1	A	262	LEU	3.1
1	A	73	ILE	3.0
1	B	289	LYS	3.0
1	B	130	ASP	3.0
1	A	144	THR	2.9
1	B	149	LEU	2.9
1	B	150	LEU	2.9
1	A	80	VAL	2.9
1	A	271	ILE	2.9
1	B	93	PHE	2.9
1	B	180	CYS	2.8
1	A	265	ILE	2.8
1	A	148	THR	2.8
1	B	55	THR	2.8
1	B	224	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	250	SER	2.8
1	B	233	GLU	2.8
1	A	93	PHE	2.8
1	A	48	GLN	2.7
1	B	217	ASP	2.7
1	B	144	THR	2.7
1	A	194	ARG	2.7
1	B	253	LEU	2.7
1	A	225	VAL	2.7
1	B	143	PHE	2.7
1	B	63	VAL	2.7
1	B	251	ARG	2.7
1	B	124	LEU	2.6
1	B	228	HIS	2.6
1	A	111	LYS	2.6
1	B	269	ARG	2.6
1	B	227	VAL	2.5
1	B	50	TRP	2.5
1	A	273	PHE	2.5
1	A	151	GLY	2.5
1	B	123	LEU	2.5
1	B	244	HIS	2.5
1	B	59	PHE	2.5
1	A	45	ARG	2.5
1	A	152	TYR	2.5
1	B	187	VAL	2.4
1	B	101	ILE	2.4
1	B	284	PHE	2.4
1	A	208	LEU	2.4
1	B	137	GLN	2.4
1	B	45	ARG	2.4
1	B	54	GLY	2.4
1	B	106	TYR	2.4
1	B	280	ARG	2.4
1	B	151	GLY	2.4
1	A	205	HIS	2.4
1	A	78	ARG	2.4
1	B	99	CYS	2.3
1	A	191	THR	2.3
1	A	58	ARG	2.3
1	A	223	GLY	2.3
1	A	280	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	153	LEU	2.3
1	A	81	ASP	2.3
1	B	192	VAL	2.2
1	B	173	CYS	2.2
1	A	193	SER	2.2
1	A	195	ARG	2.2
1	B	256	ASP	2.2
1	B	177	ARG	2.2
1	B	267	SER	2.2
1	B	47	ARG	2.2
1	A	200	ALA	2.1
1	A	240	ALA	2.1
1	B	275	CYS	2.1
1	B	174	PRO	2.1
1	B	183	ASN	2.1
1	B	86	TRP	2.1
1	B	148	THR	2.1
1	B	92	ALA	2.1
1	B	164	ASP	2.1
1	A	184	PRO	2.1
1	A	186	SER	2.1
1	A	119	CYS	2.1
1	B	263	GLU	2.1
1	B	64	LEU	2.1
1	B	277	ASN	2.0
1	B	258	THR	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	N1C	A	301	35/35	0.13	-1.25	25,33,38,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.