



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 09:40 PM GMT

PDB ID : 2PGL
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 : 1.76 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>

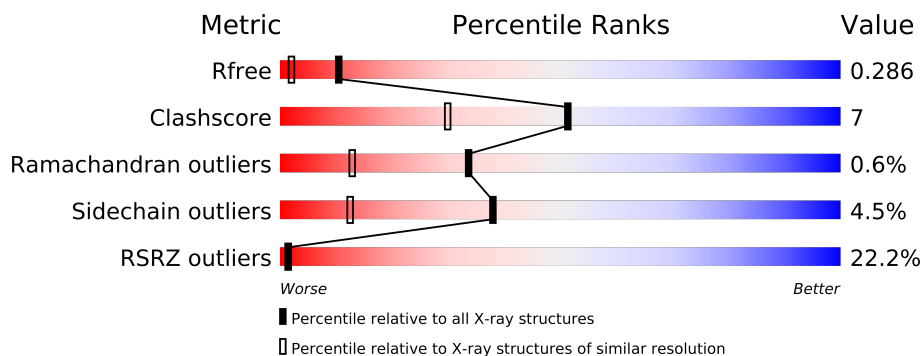
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.76 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	1134 (1.76-1.76)
Clashscore	79885	1304 (1.76-1.76)
Ramachandran outliers	78287	1288 (1.76-1.76)
Sidechain outliers	78261	1288 (1.76-1.76)
RSRZ outliers	66119	1135 (1.76-1.76)

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 4520 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.

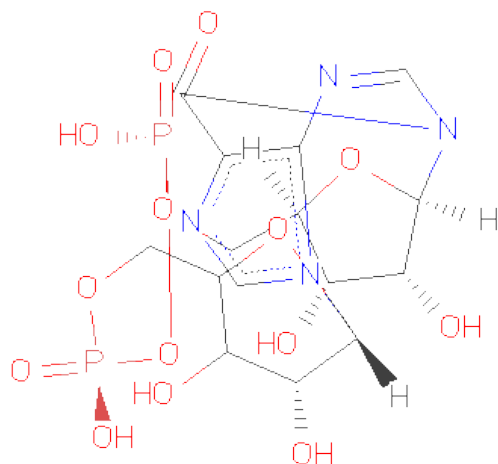
- Molecule 1 is a protein called ADP-ribosyl cyclase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			2050	1290	359	385	16			
1	B	252	Total	C	N	O	S	0	0	0
			2050	1290	359	385	16			

There are 24 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
A	226	GLN	GLU	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
B	226	GLN	GLU	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
			Total	C	N	O	P		
2	A	1	35	15	4	14	2	0	0

- Molecule 3 is water.

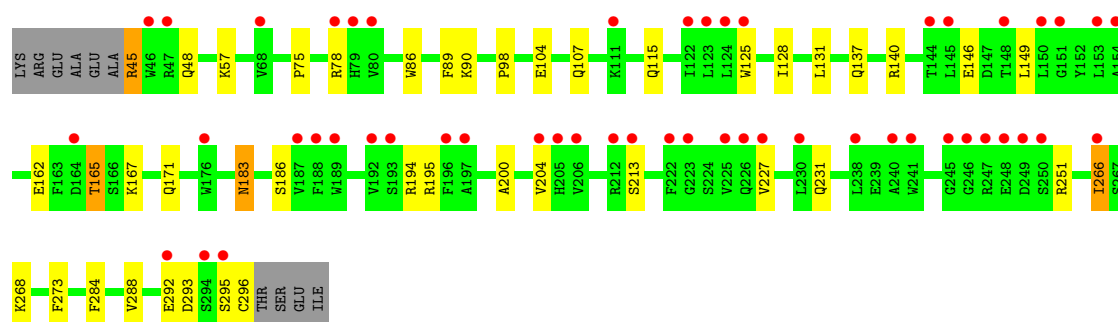
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	198	Total	O	0	0
			198	198		
3	B	187	Total	O	0	0
			187	187		

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.

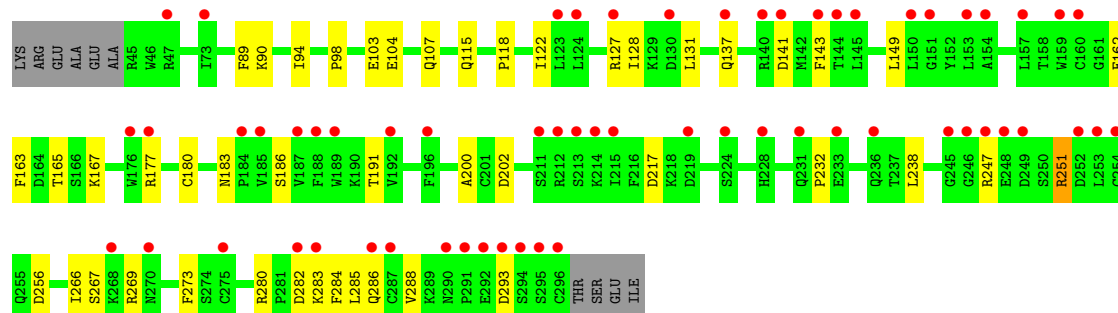
• Molecule 1: ADP-ribosyl cyclase 1

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	41.87Å 53.20Å 65.71Å 105.79° 92.07° 95.20°	Depositor
Resolution (Å)	20.00 – 1.76 26.26 – 1.76	Depositor EDS
% Data completeness (in resolution range)	95.5 (20.00-1.76) 95.5 (26.26-1.76)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 1.76Å)	Xtriage
Refinement program	REFMAC 5.3.0021	Depositor
R, R_{free}	0.210 , 0.245 0.254 , 0.286	Depositor DCC
R_{free} test set	2558 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 51403 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4520	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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	0.63	0/2101	0.68	0/2846
1	B	0.62	0/2101	0.64	0/2846
All	All	0.62	0/4202	0.66	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	1978	30	0
1	B	2050	0	1978	29	0
2	A	35	0	15	4	0
3	A	198	0	0	7	0
3	B	187	0	0	2	0
All	All	4520	0	3971	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:A:165:THR:HG23	1:A:167:LYS:H	1.40	0.86
1:A:115:GLN:HE22	1:A:149:LEU:H	1.25	0.84
1:B:115:GLN:HE22	1:B:149:LEU:H	1.24	0.82
1:A:266:ILE:HD11	1:A:273:PHE:HB2	1.64	0.78
1:A:231:GLN:HG3	3:A:498:HOH:O	1.84	0.76
1:A:268:LYS:HD3	1:B:163:PHE:HE1	1.52	0.74
1:B:165:THR:HG23	1:B:167:LYS:H	1.53	0.72
1:A:162:GLU:OE2	1:A:165:THR:HG21	1.91	0.70
1:A:268:LYS:HD3	1:B:163:PHE:CE1	2.27	0.70
1:B:103:GLU:HG2	1:B:191:THR:OG1	1.91	0.69
1:B:183:ASN:ND2	1:B:186:SER:H	1.94	0.66
1:A:183:ASN:ND2	1:A:186:SER:H	1.95	0.65
1:B:127:ARG:HH12	1:B:217:ASP:HB3	1.64	0.62
1:B:266:ILE:HD11	1:B:273:PHE:HB2	1.81	0.62
1:A:75:PRO:O	1:A:78:ARG:HB2	2.00	0.61
1:A:183:ASN:HD21	1:A:186:SER:H	1.48	0.61
1:A:194:ARG:NH2	3:A:326:HOH:O	2.36	0.59
1:B:232:PRO:HG3	1:B:269:ARG:O	2.02	0.58
1:B:162:GLU:HB2	1:B:165:THR:HG22	1.85	0.58
1:A:293:ASP:HB3	1:A:296:CYS:SG	2.43	0.58
1:A:140:ARG:NH2	3:A:388:HOH:O	2.37	0.57
1:B:137:GLN:HG2	3:B:474:HOH:O	2.03	0.57
2:A:301:N1C:H2D	2:A:301:N1C:O6	2.04	0.57
1:B:284:PHE:O	1:B:288:VAL:HG23	2.05	0.56
1:A:251:ARG:HD3	1:A:251:ARG:H	1.71	0.56
1:A:146:GLU:OE2	2:A:301:N1C:O6	2.23	0.56
1:B:122:ILE:HD12	1:B:200:ALA:HA	1.90	0.52
1:A:98:PRO:O	1:A:183:ASN:HA	2.10	0.52
1:B:180:CYS:HB2	3:B:324:HOH:O	2.10	0.52
1:B:104:GLU:HA	1:B:107:GLN:HG2	1.92	0.51
1:A:125:TRP:O	2:A:301:N1C:H5D1	2.11	0.51
1:B:183:ASN:HD21	1:B:186:SER:H	1.60	0.49
1:B:90:LYS:HG2	1:B:94:ILE:HG13	1.95	0.49
1:B:115:GLN:NE2	1:B:149:LEU:H	2.03	0.48
1:B:127:ARG:HD2	1:B:127:ARG:N	2.27	0.48
1:A:137:GLN:NE2	3:A:419:HOH:O	2.47	0.47
1:A:200:ALA:HB1	1:A:204:VAL:HG22	1.97	0.47
1:A:266:ILE:CD1	1:A:273:PHE:HB2	2.41	0.47
1:A:200:ALA:HB1	1:A:204:VAL:CG2	2.46	0.46
1:A:162:GLU:HB2	1:A:165:THR:HG22	1.97	0.45
1:A:266:ILE:HD11	1:A:273:PHE:CB	2.41	0.45
1:B:266:ILE:HD11	1:B:273:PHE:CB	2.46	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:86:TRP:CE2	1:A:90:LYS:HG3	2.51	0.45
1:B:280:ARG:NH1	1:B:283:LYS:HE3	2.32	0.44
1:A:165:THR:CG2	1:A:167:LYS:H	2.20	0.44
1:B:98:PRO:O	1:B:183:ASN:HA	2.18	0.43
1:B:118:PRO:HD2	1:B:143:PHE:CE2	2.54	0.43
1:B:90:LYS:CG	1:B:94:ILE:HG13	2.48	0.43
1:B:238:LEU:HB3	1:B:266:ILE:HD13	2.01	0.42
3:A:398:HOH:O	1:B:177:ARG:HD2	2.18	0.42
1:A:104:GLU:HA	1:A:107:GLN:HG2	2.00	0.42
2:A:301:N1C:H8	3:A:334:HOH:O	2.19	0.42
1:B:266:ILE:HG13	1:B:267:SER:N	2.34	0.41
1:A:48:GLN:NE2	1:A:171:GLN:HB3	2.35	0.41
1:A:284:PHE:O	1:A:288:VAL:HG23	2.20	0.41
1:B:162:GLU:HB2	1:B:165:THR:CG2	2.50	0.40
1:A:195:ARG:HG3	3:A:326:HOH:O	2.20	0.40
1:B:251:ARG:HG3	1:B:251:ARG:H	1.54	0.40
1:A:45:ARG:HH11	1:A:45:ARG:HG2	1.86	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%)	10 (4%)	1 (0%)	43	21
1	B	250/262 (95%)	234 (94%)	14 (6%)	2 (1%)	27	9
All	All	500/524 (95%)	473 (95%)	24 (5%)	3 (1%)	33	13

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	247	ARG
1	B	293	ASP
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%)	222 (95%)	11 (5%)	36	11
1	B	233/241 (97%)	223 (96%)	10 (4%)	40	13
All	All	466/482 (97%)	445 (96%)	21 (4%)	38	12

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ARG
1	A	57	LYS
1	A	89	PHE
1	A	131	LEU
1	A	165	THR
1	A	183	ASN
1	A	213	SER
1	A	227	VAL
1	A	266	ILE
1	A	292	GLU
1	A	295	SER
1	B	89	PHE
1	B	128	ILE
1	B	131	LEU
1	B	141	ASP
1	B	202	ASP
1	B	251	ARG
1	B	256	ASP
1	B	282	ASP
1	B	285	LEU
1	B	286	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	GLN
1	A	115	GLN
1	A	134	GLN

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Mol	Chain	Res	Type
1	A	137	GLN
1	A	139	GLN
1	A	183	ASN
1	B	115	GLN
1	B	183	ASN

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	1.91	5 (12%)	60,62,62	2.57	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 (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	N1C	C3'-C4'	-7.60	1.32	1.53
2	A	301	N1C	C6-C5	5.32	1.49	1.41
2	A	301	N1C	O3'-C3'	-4.03	1.33	1.43
2	A	301	N1C	C2-N3	3.21	1.35	1.30
2	A	301	N1C	C6-N1	2.30	1.41	1.35

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	N1C	C6-C5-N7	-11.65	133.62	134.24
2	A	301	N1C	N3-C4-N9	5.02	134.50	125.43
2	A	301	N1C	O3'-C3'-C4'	4.35	123.91	111.08
2	A	301	N1C	C2'-C3'-C4'	4.31	111.25	102.65
2	A	301	N1C	C4'-O4'-C1'	-4.11	105.29	109.75
2	A	301	N1C	O4'-C4'-C5'	4.04	123.79	109.36
2	A	301	N1C	PA-O3A-PB	-4.03	119.86	131.68
2	A	301	N1C	O3'-C3'-C2'	4.00	124.84	111.83
2	A	301	N1C	O4'-C1'-N9	-3.71	104.98	108.44
2	A	301	N1C	C3D-C2D-C1D	3.65	106.61	100.91
2	A	301	N1C	C2-N3-C4	3.58	122.34	116.23
2	A	301	N1C	O4'-C4'-C3'	3.08	111.42	105.17
2	A	301	N1C	C5-C4-N3	-3.05	119.05	125.70
2	A	301	N1C	C8-N9-C4	2.90	109.11	106.90
2	A	301	N1C	C4D-O4D-C1D	2.66	112.64	109.75
2	A	301	N1C	O4D-C1D-C2D	-2.53	102.90	106.77
2	A	301	N1C	C5'-C4'-C3'	2.39	124.79	115.21
2	A	301	N1C	C6-C5-C4	-2.01	118.48	119.92

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%)	1.17	51 (20%) 2 2	26, 33, 47, 58	0
1	B	252/262 (96%)	1.41	61 (24%) 1 1	27, 35, 53, 64	0
All	All	504/524 (96%)	1.29	112 (22%) 1 1	26, 34, 50, 64	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	246	GLY	12.8
1	A	249	ASP	8.7
1	B	249	ASP	7.9
1	A	246	GLY	7.7
1	B	292	GLU	7.6
1	B	245	GLY	7.5
1	B	295	SER	7.4
1	B	291	PRO	7.3
1	A	247	ARG	7.2
1	B	293	ASP	6.8
1	B	213	SER	6.4
1	B	294	SER	6.4
1	B	248	GLU	6.3
1	B	290	ASN	6.1
1	B	296	CYS	5.9
1	A	248	GLU	5.7
1	B	212	ARG	5.1
1	A	124	LEU	5.1
1	B	247	ARG	5.0
1	B	270	ASN	4.8
1	A	292	GLU	4.6
1	A	245	GLY	4.5
1	A	213	SER	4.5
1	A	47	ARG	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	159	TRP	4.2
1	A	46	TRP	4.1
1	A	145	LEU	4.0
1	A	123	LEU	3.8
1	A	295	SER	3.8
1	A	78	ARG	3.7
1	A	212	ARG	3.6
1	B	189	TRP	3.5
1	B	127	ARG	3.5
1	A	222	PHE	3.5
1	B	283	LYS	3.5
1	B	211	SER	3.4
1	A	238	LEU	3.4
1	A	227	VAL	3.4
1	B	188	PHE	3.3
1	B	143	PHE	3.3
1	B	282	ASP	3.3
1	A	79	HIS	3.3
1	B	286	GLN	3.3
1	A	206	VAL	3.2
1	A	204	VAL	3.2
1	B	287	CYS	3.1
1	B	145	LEU	3.1
1	A	189	TRP	3.0
1	A	225	VAL	3.0
1	B	254	CYS	3.0
1	A	80	VAL	3.0
1	A	164	ASP	3.0
1	B	137	GLN	2.8
1	B	144	THR	2.8
1	B	185	VAL	2.8
1	A	144	THR	2.7
1	A	230	LEU	2.7
1	A	192	VAL	2.7
1	A	294	SER	2.7
1	B	153	LEU	2.7
1	A	154	ALA	2.7
1	A	197	ALA	2.7
1	B	214	LYS	2.6
1	A	125	TRP	2.6
1	B	228	HIS	2.6
1	B	219	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	160	CYS	2.6
1	B	192	VAL	2.6
1	A	196	PHE	2.6
1	B	157	LEU	2.6
1	B	140	ARG	2.6
1	A	188	PHE	2.6
1	B	184	PRO	2.5
1	B	176	TRP	2.5
1	A	148	THR	2.5
1	B	123	LEU	2.5
1	B	141	ASP	2.5
1	B	224	SER	2.4
1	A	122	ILE	2.4
1	A	150	LEU	2.4
1	B	124	LEU	2.4
1	B	236	GLN	2.4
1	B	215	ILE	2.4
1	A	111	LYS	2.3
1	B	268	LYS	2.3
1	A	187	VAL	2.3
1	A	193	SER	2.3
1	A	151	GLY	2.3
1	A	223	GLY	2.3
1	B	150	LEU	2.2
1	A	240	ALA	2.2
1	B	196	PHE	2.2
1	A	266	ILE	2.2
1	B	177	ARG	2.2
1	B	252	ASP	2.2
1	A	176	TRP	2.1
1	B	73	ILE	2.1
1	B	233	GLU	2.1
1	A	153	LEU	2.1
1	B	253	LEU	2.1
1	A	226	GLN	2.1
1	A	250	SER	2.1
1	B	187	VAL	2.1
1	A	68	VAL	2.1
1	A	241	TRP	2.1
1	B	231	GLN	2.1
1	B	130	ASP	2.0
1	B	151	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	47	ARG	2.0
1	A	205	HIS	2.0
1	B	275	CYS	2.0
1	B	154	ALA	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(Å ²)	Q<0.9
2	N1C	A	301	35/35	0.17	-0.36	25,44,50,52	0

6.5 Other polymers ⓘ

There are no such residues in this entry.