



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 05:01 AM GMT

PDB ID : 2XMR
Title : CRYSTAL STRUCTURE OF HUMAN NDRG2 PROTEIN PROVIDES INSIGHT INTO ITS ROLE AS A TUMOR SUPPRESSOR
Authors : Hwang, J.; Kim, Y.; Lee, H.; Kim, M.H.
Deposited on : 2010-07-29
Resolution : 2.00 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>

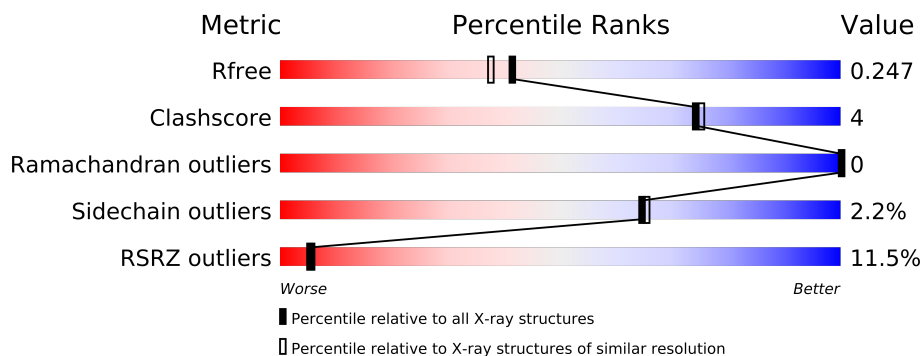
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.00 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	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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	281	
1	B	281	
1	C	281	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7032 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 PROTEIN NDRG2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2197	1405	363	416	13			
1	B	281	Total	C	N	O	S	0	0	0
			2197	1405	363	416	13			
1	C	281	Total	C	N	O	S	0	1	0
			2208	1411	367	417	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36
A	47	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36
B	45	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36
B	47	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36
C	45	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36
C	47	ALA	LYS	ENGINEERED MUTATION	UNP Q9UN36

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

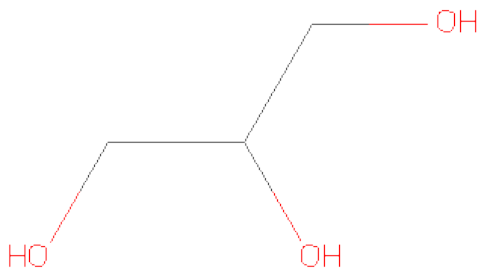
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

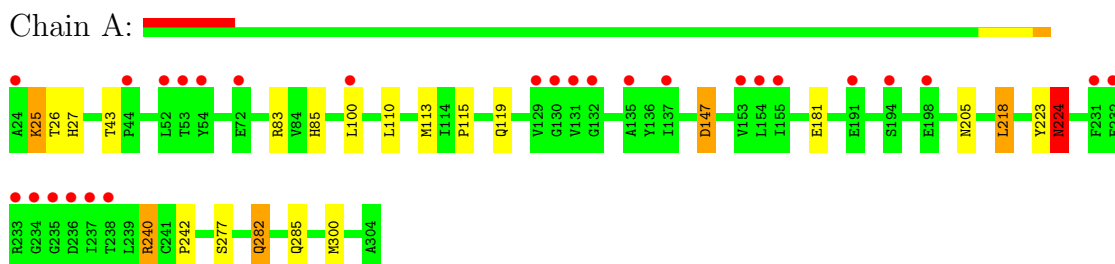
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	132	Total	O	0	0
			132	132		
5	B	152	Total	O	0	0
			152	152		
5	C	115	Total	O	0	0
			115	115		

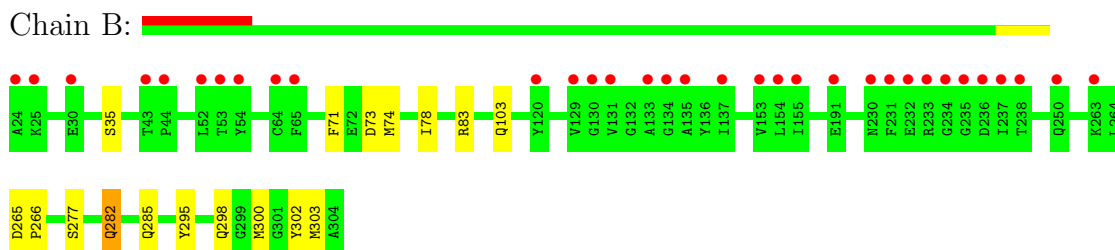
3 Residue-property plots

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.

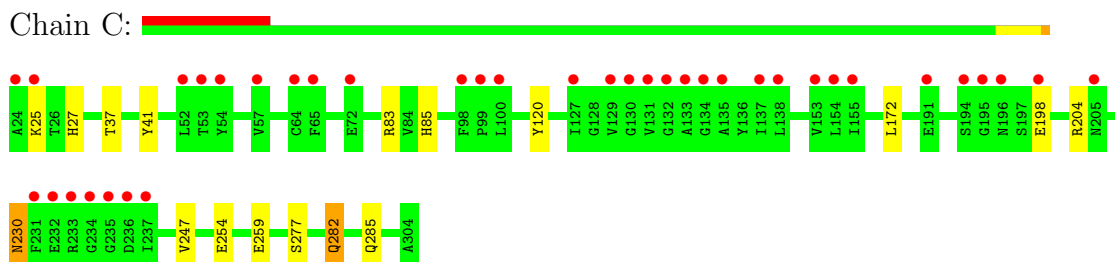
• Molecule 1: PROTEIN NDRG2



• Molecule 1: PROTEIN NDRG2



• Molecule 1: PROTEIN NDRG2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.23Å 88.05Å 126.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 29.93 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.5 (30.00-2.00) 96.4 (29.93-2.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.180 , 0.232 0.196 , 0.247	Depositor DCC
R_{free} test set	3217 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 40.4	EDS
Estimated twinning fraction	0.019 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 63288 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7032	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: GOL, CA, 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	0.60	1/2255 (0.0%)	0.67	2/3072 (0.1%)
1	B	0.57	0/2255	0.64	0/3072
1	C	0.53	0/2266	0.61	0/3086
All	All	0.57	1/6776 (0.0%)	0.64	2/9230 (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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	218	LEU	CG-CD1	-7.71	1.23	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	218	LEU	CB-CG-CD1	6.23	121.60	111.00
1	A	224	ASN	N-CA-C	-5.49	96.17	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	223	TYR	Peptide

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Mol	Chain	Res	Type	Group
1	A	224	ASN	Peptide

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	2197	0	2131	26	0
1	B	2197	0	2131	16	0
1	C	2208	0	2143	17	0
2	A	1	0	0	0	0
3	A	4	0	3	4	0
3	B	4	0	3	2	0
3	C	4	0	3	5	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
4	C	6	0	8	0	0
5	A	132	0	0	2	0
5	B	152	0	0	0	0
5	C	115	0	0	3	0
All	All	7032	0	6438	54	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 4.

All (54) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:83:ARG:HE	3:B:1305:ACT:H1	1.00	1.09
1:A:83:ARG:HE	3:A:1306:ACT:H1	1.18	1.01
1:C:85:HIS:NE2	3:C:1305:ACT:H1	1.80	0.97
1:A:25:LYS:HZ2	1:A:26:THR:H	1.13	0.94
1:A:282:GLN:H	1:A:282:GLN:HE21	1.16	0.93
1:A:119:GLN:HE22	1:B:298:GLN:HE22	1.10	0.93
1:B:83:ARG:NE	3:B:1305:ACT:H1	1.86	0.90
1:C:282:GLN:H	1:C:282:GLN:HE21	1.20	0.84
1:B:282:GLN:H	1:B:282:GLN:HE21	1.33	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:83:ARG:HE	3:C:1305:ACT:CH3	1.99	0.75
1:A:240:ARG:HG2	1:A:240:ARG:HH11	1.53	0.73
1:A:119:GLN:NE2	1:B:298:GLN:HE22	1.86	0.73
1:A:83:ARG:NE	3:A:1306:ACT:H1	2.01	0.71
1:A:282:GLN:NE2	1:A:282:GLN:H	1.91	0.66
1:C:282:GLN:H	1:C:282:GLN:NE2	1.92	0.66
1:C:83:ARG:HE	3:C:1305:ACT:H2	1.61	0.66
1:A:240:ARG:CG	1:A:240:ARG:HH11	2.09	0.65
1:A:85:HIS:NE2	3:A:1306:ACT:H2	2.15	0.61
1:C:37:THR:HG22	5:C:2025:HOH:O	2.04	0.57
1:C:83:ARG:HE	3:C:1305:ACT:H3	1.69	0.57
1:A:85:HIS:NE2	3:A:1306:ACT:CH3	2.68	0.56
1:A:282:GLN:N	1:A:282:GLN:HE21	1.96	0.56
1:B:282:GLN:H	1:B:282:GLN:NE2	2.00	0.56
1:C:25:LYS:HG3	1:C:27:HIS:CE1	2.41	0.56
1:B:300:MET:HE2	1:B:302:TYR:CE2	2.42	0.55
1:B:300:MET:HE2	1:B:302:TYR:HE2	1.72	0.55
1:A:110:LEU:HD23	1:A:113:MET:HE1	1.88	0.54
1:A:25:LYS:HE3	5:C:2036:HOH:O	2.08	0.52
1:C:25:LYS:HG3	1:C:27:HIS:HE1	1.75	0.51
1:A:119:GLN:HE22	1:B:298:GLN:NE2	1.92	0.50
1:C:282:GLN:HE21	1:C:282:GLN:N	2.00	0.49
1:A:240:ARG:NH1	1:A:240:ARG:CG	2.72	0.47
1:B:300:MET:CE	1:B:302:TYR:HE2	2.26	0.47
1:A:25:LYS:NZ	1:A:26:THR:H	1.97	0.46
1:A:110:LEU:HD23	1:A:113:MET:CE	2.47	0.45
1:A:242:PRO:HG2	1:A:300:MET:HG2	1.99	0.45
1:C:230:ASN:HA	1:C:230:ASN:HD22	1.67	0.44
1:B:71:PHE:CE2	1:B:73:ASP:HB2	2.53	0.44
1:A:115:PRO:HB2	1:B:303:MET:CE	2.48	0.44
1:B:277:SER:HB3	1:B:285:GLN:HE21	1.83	0.43
1:C:277:SER:HB3	1:C:285:GLN:OE1	2.19	0.43
1:A:25:LYS:HA	1:A:25:LYS:HZ3	1.84	0.43
1:B:74:MET:O	1:B:78:ILE:HG12	2.18	0.42
1:A:25:LYS:HB3	1:A:27:HIS:CE1	2.53	0.42
1:B:265:ASP:HA	1:B:266:PRO:HD2	1.86	0.42
1:B:295:TYR:HB2	1:C:172:LEU:HG	2.00	0.42
1:A:181:GLU:HG2	5:A:2088:HOH:O	2.19	0.42
1:C:27:HIS:HD2	1:C:120:TYR:OH	2.03	0.41
1:C:41:TYR:HB2	3:C:1305:ACT:H3	2.02	0.41
1:A:115:PRO:O	1:A:119:GLN:HG3	2.19	0.41
1:A:277:SER:CB	1:A:285:GLN:HG3	2.51	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:247:VAL:HG21	1:C:254:GLU:HA	2.02	0.41
1:C:204:ARG:HD2	5:C:2077:HOH:O	2.19	0.41
1:A:147:ASP:CG	5:A:2062:HOH:O	2.59	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	279/281 (99%)	276 (99%)	3 (1%)	0	100	100
1	B	279/281 (99%)	278 (100%)	1 (0%)	0	100	100
1	C	280/281 (100%)	280 (100%)	0	0	100	100
All	All	838/843 (99%)	834 (100%)	4 (0%)	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	238/238 (100%)	229 (96%)	9 (4%)	44	39
1	B	238/238 (100%)	235 (99%)	3 (1%)	80	82
1	C	239/238 (100%)	235 (98%)	4 (2%)	73	75
All	All	715/714 (100%)	699 (98%)	16 (2%)	64	65

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

Mol	Chain	Res	Type
1	A	25	LYS
1	A	43	THR
1	A	100	LEU
1	A	147	ASP
1	A	205	ASN
1	A	218	LEU
1	A	224	ASN
1	A	240	ARG
1	A	282	GLN
1	B	35	SER
1	B	103	GLN
1	B	282	GLN
1	C	198	GLU
1	C	230	ASN
1	C	259	GLU
1	C	282	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	75	GLN
1	A	119	GLN
1	A	224	ASN
1	A	280	GLN
1	A	282	GLN
1	A	298	GLN
1	B	27	HIS
1	B	75	GLN
1	B	103	GLN
1	B	230	ASN
1	B	280	GLN
1	B	282	GLN
1	B	285	GLN
1	C	27	HIS
1	C	75	GLN
1	C	103	GLN
1	C	224	ASN
1	C	230	ASN
1	C	280	GLN
1	C	282	GLN
1	C	298	GLN

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 ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
3	ACT	A	1306	-	1,3,3	1.48	0	0,3,3	0.00	-
4	GOL	A	1307	-	5,5,5	0.28	0	5,5,5	0.50	0
3	ACT	B	1305	-	1,3,3	0.68	0	0,3,3	0.00	-
4	GOL	B	1306	-	5,5,5	0.33	0	5,5,5	0.24	0
3	ACT	C	1305	-	1,3,3	0.40	0	0,3,3	0.00	-
4	GOL	C	1306	-	5,5,5	0.28	0	5,5,5	0.39	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
3	ACT	A	1306	-	-	0/0/0/0	0/0/0/0
4	GOL	A	1307	-	-	0/4/4/4	0/0/0/0
3	ACT	B	1305	-	-	0/0/0/0	0/0/0/0
4	GOL	B	1306	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ACT	C	1305	-	-	0/0/0/0	0/0/0/0
4	GOL	C	1306	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

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	281/281 (100%)	0.36	27 (9%) 8 8	17, 25, 42, 61	1 (0%)
1	B	281/281 (100%)	0.43	33 (11%) 5 5	18, 25, 42, 60	1 (0%)
1	C	281/281 (100%)	0.51	38 (13%) 4 4	17, 27, 42, 59	1 (0%)
All	All	843/843 (100%)	0.43	98 (11%) 5 5	17, 26, 43, 61	3 (0%)

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	ASP	6.4
1	C	234	GLY	5.9
1	B	234	GLY	5.8
1	C	237	ILE	5.6
1	B	232	GLU	5.5
1	B	24	ALA	5.4
1	C	236	ASP	5.4
1	A	234	GLY	5.1
1	A	233	ARG	4.5
1	B	235	GLY	4.4
1	A	236	ASP	4.3
1	C	131	VAL	4.2
1	C	25	LYS	4.2
1	C	191	GLU	4.1
1	B	237	ILE	3.9
1	C	232	GLU	3.6
1	A	135	ALA	3.6
1	C	24	ALA	3.5
1	A	232	GLU	3.5
1	B	233	ARG	3.5
1	C	134	GLY	3.4
1	B	129	VAL	3.4
1	A	129	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	24	ALA	3.3
1	A	130	GLY	3.3
1	B	153	VAL	3.3
1	C	235	GLY	3.3
1	B	231	PHE	3.3
1	B	238	THR	3.3
1	A	237	ILE	3.2
1	C	153	VAL	3.2
1	C	133	ALA	3.0
1	B	154	LEU	3.0
1	C	100	LEU	3.0
1	C	233	ARG	3.0
1	B	43	THR	3.0
1	C	194	SER	3.0
1	C	129	VAL	3.0
1	A	100	LEU	2.8
1	A	131	VAL	2.8
1	A	238	THR	2.8
1	A	194	SER	2.7
1	C	135	ALA	2.7
1	C	64	CYS	2.7
1	B	65	PHE	2.7
1	B	134	GLY	2.7
1	B	135	ALA	2.6
1	C	127	ILE	2.6
1	C	57	VAL	2.6
1	B	130	GLY	2.6
1	A	153	VAL	2.6
1	A	155	ILE	2.6
1	B	155	ILE	2.6
1	C	155	ILE	2.5
1	C	132	GLY	2.5
1	C	205	ASN	2.5
1	B	131	VAL	2.5
1	B	30	GLU	2.5
1	C	65	PHE	2.5
1	A	154	LEU	2.4
1	C	137	ILE	2.4
1	A	44	PRO	2.4
1	C	196	ASN	2.4
1	B	263	LYS	2.4
1	B	44	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	191	GLU	2.3
1	A	132	GLY	2.3
1	A	231	PHE	2.3
1	B	54	TYR	2.3
1	B	25	LYS	2.3
1	C	195	GLY	2.3
1	C	130	GLY	2.3
1	B	64	CYS	2.3
1	A	54	TYR	2.3
1	C	54	TYR	2.2
1	B	250	GLN	2.2
1	C	72	GLU	2.2
1	B	53	THR	2.2
1	B	52	LEU	2.2
1	C	53	THR	2.2
1	A	72	GLU	2.2
1	A	198	GLU	2.2
1	B	230	ASN	2.2
1	A	53	THR	2.2
1	C	98	PHE	2.1
1	A	235	GLY	2.1
1	A	191	GLU	2.1
1	B	120	TYR	2.1
1	A	137	ILE	2.1
1	B	133	ALA	2.1
1	C	198	GLU	2.1
1	C	52	LEU	2.1
1	C	154	LEU	2.1
1	C	231	PHE	2.1
1	A	52	LEU	2.1
1	B	137	ILE	2.0
1	C	99	PRO	2.0
1	C	138	LEU	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
3	ACT	A	1306	4/4	0.14	1.74	16,20,21,22	0
3	ACT	C	1305	4/4	0.16	1.58	27,29,30,30	0
4	GOL	A	1307	6/6	0.10	-0.55	28,30,31,33	0
4	GOL	C	1306	6/6	0.12	-0.62	33,37,38,40	0
4	GOL	B	1306	6/6	0.10	-0.72	24,28,31,35	0
3	ACT	B	1305	4/4	0.10	-1.39	18,20,21,22	0
2	CA	A	1305	1/1	0.07	-1.57	26,26,26,26	0

6.5 Other polymers ⓘ

There are no such residues in this entry.