



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:17 PM GMT

PDB ID : 4OF8
Title : Crystal Structure of Rst D1-D2
Authors : Ozkan, E.; Garcia, K.C.
Deposited on : 2014-01-14
Resolution : 1.90 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

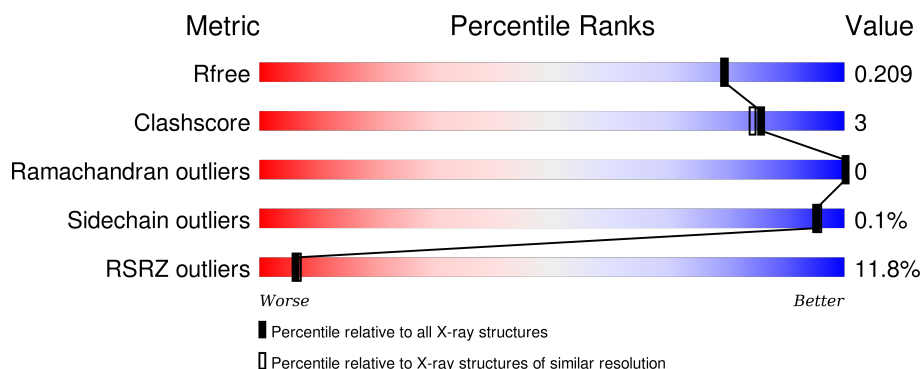
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	228	<div> <div>4%</div> <div>89%</div> <div>7%</div> </div>
1	B	228	<div> <div>8%</div> <div>81%</div> <div>11%</div> <div>8%</div> </div>
1	C	228	<div> <div>12%</div> <div>88%</div> <div>8%</div> </div>
1	D	228	<div> <div>18%</div> <div>85%</div> <div>7%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	302	-	-	-	X
3	GOL	B	302	-	-	-	X
3	GOL	D	302	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7448 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 Irregular chiasm C-roughest protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1647	1027	289	324	7			
1	B	209	Total	C	N	O	S	0	1	0
			1639	1021	290	321	7			
1	C	210	Total	C	N	O	S	0	1	0
			1640	1021	288	324	7			
1	D	209	Total	C	N	O	S	0	1	0
			1639	1021	290	321	7			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	ALA	-	EXPRESSION TAG	UNP Q08180
A	18	ASP	-	EXPRESSION TAG	UNP Q08180
A	19	PRO	-	EXPRESSION TAG	UNP Q08180
A	238	HIS	-	EXPRESSION TAG	UNP Q08180
A	239	HIS	-	EXPRESSION TAG	UNP Q08180
A	240	HIS	-	EXPRESSION TAG	UNP Q08180
A	241	HIS	-	EXPRESSION TAG	UNP Q08180
A	242	HIS	-	EXPRESSION TAG	UNP Q08180
A	243	HIS	-	EXPRESSION TAG	UNP Q08180
A	244	HIS	-	EXPRESSION TAG	UNP Q08180
B	17	ALA	-	EXPRESSION TAG	UNP Q08180
B	18	ASP	-	EXPRESSION TAG	UNP Q08180
B	19	PRO	-	EXPRESSION TAG	UNP Q08180
B	238	HIS	-	EXPRESSION TAG	UNP Q08180
B	239	HIS	-	EXPRESSION TAG	UNP Q08180
B	240	HIS	-	EXPRESSION TAG	UNP Q08180
B	241	HIS	-	EXPRESSION TAG	UNP Q08180
B	242	HIS	-	EXPRESSION TAG	UNP Q08180
B	243	HIS	-	EXPRESSION TAG	UNP Q08180
B	244	HIS	-	EXPRESSION TAG	UNP Q08180
C	17	ALA	-	EXPRESSION TAG	UNP Q08180

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Chain	Residue	Modelled	Actual	Comment	Reference
C	18	ASP	-	EXPRESSION TAG	UNP Q08180
C	19	PRO	-	EXPRESSION TAG	UNP Q08180
C	238	HIS	-	EXPRESSION TAG	UNP Q08180
C	239	HIS	-	EXPRESSION TAG	UNP Q08180
C	240	HIS	-	EXPRESSION TAG	UNP Q08180
C	241	HIS	-	EXPRESSION TAG	UNP Q08180
C	242	HIS	-	EXPRESSION TAG	UNP Q08180
C	243	HIS	-	EXPRESSION TAG	UNP Q08180
C	244	HIS	-	EXPRESSION TAG	UNP Q08180
D	17	ALA	-	EXPRESSION TAG	UNP Q08180
D	18	ASP	-	EXPRESSION TAG	UNP Q08180
D	19	PRO	-	EXPRESSION TAG	UNP Q08180
D	238	HIS	-	EXPRESSION TAG	UNP Q08180
D	239	HIS	-	EXPRESSION TAG	UNP Q08180
D	240	HIS	-	EXPRESSION TAG	UNP Q08180
D	241	HIS	-	EXPRESSION TAG	UNP Q08180
D	242	HIS	-	EXPRESSION TAG	UNP Q08180
D	243	HIS	-	EXPRESSION TAG	UNP Q08180
D	244	HIS	-	EXPRESSION TAG	UNP Q08180

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Na 1 1	0	0
2	A	1	Total Na 1 1	0	0
2	D	1	Total Na 1 1	0	0
2	C	1	Total Na 1 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

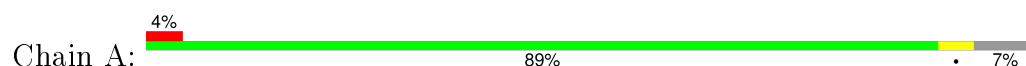
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	225	Total	O	0	0
			225	225		
4	B	188	Total	O	0	0
			188	188		
4	C	229	Total	O	0	0
			229	229		
4	D	201	Total	O	0	0
			201	201		

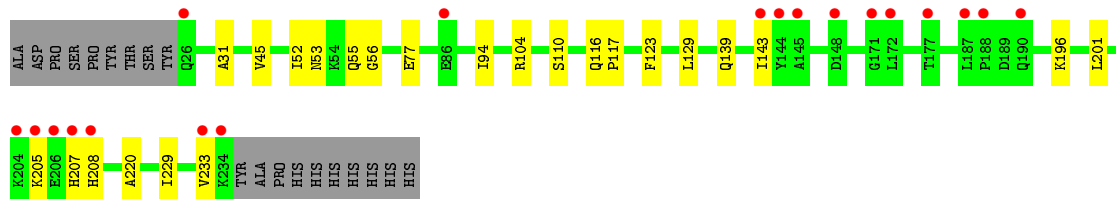
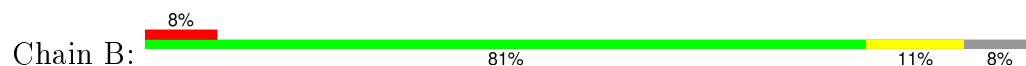
3 Residue-property plots

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.

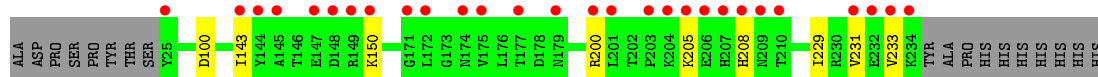
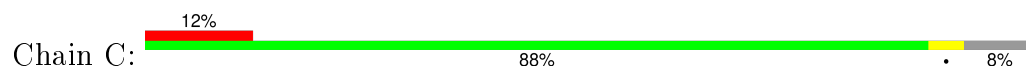
- Molecule 1: Irregular chiasm C-roughest protein



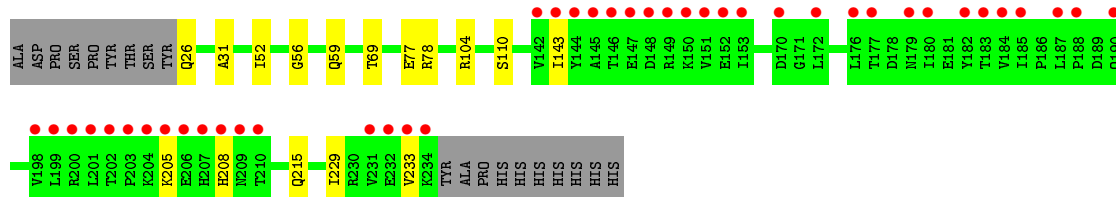
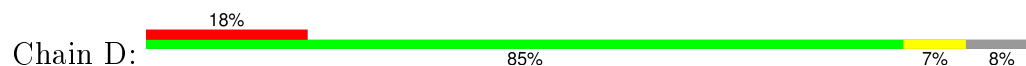
- Molecule 1: Irregular chiasm C-roughest protein



- Molecule 1: Irregular chiasm C-roughest protein



- Molecule 1: Irregular chiasm C-roughest protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.94Å 77.94Å 121.10Å 90.00° 95.44° 90.00°	Depositor
Resolution (Å)	44.57 – 1.90 44.57 – 1.90	Depositor EDS
% Data completeness (in resolution range)	88.5 (44.57-1.90) 88.5 (44.57-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 1.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1593)	Depositor
R, R_{free}	0.173 , 0.208 0.174 , 0.209	Depositor DCC
R_{free} test set	3889 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	1.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	7 of 77221 reflections (0.009%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7448	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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: GOL, NA

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	0.38	0/1679	0.51	0/2281
1	B	0.35	0/1673	0.51	0/2272
1	C	0.37	0/1674	0.51	0/2274
1	D	0.34	0/1673	0.51	0/2272
All	All	0.36	0/6699	0.51	0/9099

There are no bond length outliers.

There are no bond angle outliers.

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	1647	0	1609	5	0
1	B	1639	0	1612	15	0
1	C	1640	0	1604	5	0
1	D	1639	0	1612	12	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	12	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	12	0	16	0	0
3	D	12	0	16	0	0
4	A	225	0	0	0	0
4	B	188	0	0	2	0
4	C	229	0	0	0	0
4	D	201	0	0	5	0
All	All	7448	0	6485	37	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 3.

All (37) 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:C:208:HIS:HB2	1:C:233:VAL:HG12	1.72	0.70
1:D:26:GLN:NE2	4:D:564:HOH:O	2.25	0.69
1:D:26:GLN:N	4:D:560:HOH:O	2.28	0.65
1:B:53:ASN:OD1	1:B:55:GLN:NE2	2.34	0.61
1:D:205:LYS:HA	1:D:233:VAL:HG11	1.84	0.59
1:B:208:HIS:HB2	1:B:233:VAL:HG12	1.84	0.59
1:D:104[B]:ARG:NH2	4:D:569:HOH:O	2.37	0.57
1:D:143:ILE:HG13	1:D:229:ILE:HD11	1.89	0.54
1:C:150:LYS:HE3	1:C:200:ARG:HB3	1.88	0.54
1:D:104[B]:ARG:NH1	4:D:584:HOH:O	2.33	0.53
1:C:205:LYS:HA	1:C:233:VAL:HG11	1.90	0.53
1:B:77:GLU:OE2	4:B:529:HOH:O	2.18	0.53
1:B:205:LYS:HA	1:B:233:VAL:HG11	1.90	0.53
1:D:215:GLN:NE2	4:D:567:HOH:O	2.43	0.52
1:B:139:GLN:HG2	1:B:143:ILE:HD11	1.92	0.51
1:B:143:ILE:HG13	1:B:229:ILE:HD11	1.93	0.51
1:B:201:LEU:HD11	1:B:207:HIS:CE1	2.47	0.50
1:B:45:VAL:HG12	1:B:94:ILE:HB	1.93	0.49
1:B:104[B]:ARG:HB3	1:B:104[B]:ARG:HH21	1.77	0.49
1:B:196:LYS:NZ	4:B:561:HOH:O	2.43	0.47
1:D:208:HIS:HB2	1:D:233:VAL:HG12	1.97	0.47
1:A:35:GLN:N	1:A:35:GLN:OE1	2.48	0.47
1:C:143:ILE:HG13	1:C:229:ILE:HD11	1.96	0.47
1:A:104:ARG:HD2	1:A:123:PHE:CD1	2.51	0.45
1:B:104[B]:ARG:HD2	1:B:123:PHE:CD1	2.52	0.45
1:D:59:GLN:HB2	1:D:69:THR:HG22	1.99	0.45
1:A:116:GLN:HA	1:A:117:PRO:HD3	1.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ALA:HB2	1:B:52:ILE:HG13	2.00	0.44
1:D:56:GLY:HA3	1:D:110:SER:O	2.18	0.43
1:B:129:LEU:HB3	1:B:220:ALA:HB2	2.00	0.43
1:A:77:GLU:HG2	1:A:78:ARG:HG3	2.00	0.43
1:A:139:GLN:HG2	1:A:143:ILE:HD12	2.02	0.41
1:B:116:GLN:HA	1:B:117:PRO:HD3	1.93	0.41
1:D:77:GLU:HG3	1:D:78:ARG:HG3	2.03	0.41
1:B:56:GLY:HA3	1:B:110:SER:O	2.21	0.41
1:D:31:ALA:HB2	1:D:52:ILE:HG13	2.02	0.41

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	209/228 (92%)	205 (98%)	4 (2%)	0	100	100
1	B	208/228 (91%)	201 (97%)	7 (3%)	0	100	100
1	C	209/228 (92%)	208 (100%)	1 (0%)	0	100	100
1	D	208/228 (91%)	201 (97%)	7 (3%)	0	100	100
All	All	834/912 (91%)	815 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	180/196 (92%)	180 (100%)	0	100	100
1	B	180/196 (92%)	180 (100%)	0	100	100
1	C	180/196 (92%)	179 (99%)	1 (1%)	90	90
1	D	180/196 (92%)	180 (100%)	0	100	100
All	All	720/784 (92%)	719 (100%)	1 (0%)	95	95

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

Mol	Chain	Res	Type
1	C	231	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 10 ligands modelled in this entry, 4 are 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	GOL	A	302	-	5,5,5	0.35	0	5,5,5	0.23	0
3	GOL	A	303	-	5,5,5	0.26	0	5,5,5	0.43	0
3	GOL	B	302	-	5,5,5	0.25	0	5,5,5	0.35	0
3	GOL	B	303	-	5,5,5	0.33	0	5,5,5	0.33	0
3	GOL	D	302	-	5,5,5	0.18	0	5,5,5	0.78	0
3	GOL	D	303	-	5,5,5	0.35	0	5,5,5	0.40	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	GOL	A	302	-	-	0/4/4/4	0/0/0/0
3	GOL	A	303	-	-	0/4/4/4	0/0/0/0
3	GOL	B	302	-	-	0/4/4/4	0/0/0/0
3	GOL	B	303	-	-	0/4/4/4	0/0/0/0
3	GOL	D	302	-	-	0/4/4/4	0/0/0/0
3	GOL	D	303	-	-	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.

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	211/228 (92%)	0.32	10 (4%) 35 38	12, 26, 74, 112	0
1	B	209/228 (91%)	0.41	19 (9%) 11 13	15, 30, 82, 109	0
1	C	210/228 (92%)	0.63	28 (13%) 4 5	16, 28, 93, 129	0
1	D	209/228 (91%)	0.88	42 (20%) 1 1	14, 31, 101, 129	0
All	All	839/912 (91%)	0.56	99 (11%) 6 6	12, 28, 90, 129	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	233	VAL	7.4
1	D	207	HIS	6.9
1	D	204	LYS	6.3
1	D	233	VAL	6.2
1	B	204	LYS	6.1
1	D	201	LEU	5.7
1	D	172	LEU	5.6
1	B	207	HIS	5.5
1	A	172	LEU	5.5
1	C	234	LYS	5.4
1	D	147	GLU	5.4
1	B	234	LYS	5.4
1	C	207	HIS	5.4
1	D	203	PRO	5.3
1	C	205	LYS	5.2
1	D	202	THR	5.1
1	D	206	GLU	5.1
1	A	204	LYS	5.0
1	D	146	THR	4.9
1	D	148	ASP	4.8
1	C	172	LEU	4.7

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Mol	Chain	Res	Type	RSRZ
1	D	143	ILE	4.6
1	C	206	GLU	4.5
1	D	205	LYS	4.4
1	A	205	LYS	4.4
1	C	149	ARG	4.4
1	B	171	GLY	4.2
1	D	188	PRO	4.2
1	C	233	VAL	3.9
1	C	208	HIS	3.8
1	B	148	ASP	3.8
1	D	208	HIS	3.8
1	C	209	ASN	3.8
1	C	175	VAL	3.7
1	C	204	LYS	3.7
1	D	149	ARG	3.7
1	D	200	ARG	3.6
1	B	172	LEU	3.5
1	A	206	GLU	3.5
1	B	145	ALA	3.5
1	C	148	ASP	3.5
1	D	176	LEU	3.5
1	D	187	LEU	3.5
1	B	188	PRO	3.5
1	D	144	TYR	3.5
1	B	206	GLU	3.5
1	D	210	THR	3.4
1	C	143	ILE	3.4
1	A	207	HIS	3.3
1	D	199	LEU	3.2
1	D	153	ILE	3.2
1	D	151	VAL	3.1
1	C	201	LEU	3.1
1	D	234	LYS	3.0
1	A	177	THR	3.0
1	D	209	ASN	3.0
1	C	25	TYR	3.0
1	C	144	TYR	3.0
1	D	179	ASN	3.0
1	B	144	TYR	3.0
1	D	184	VAL	3.0
1	B	26	GLN	3.0
1	D	170	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	235	TYR	2.9
1	C	232	GLU	2.8
1	D	198	VAL	2.8
1	B	187	LEU	2.8
1	D	190	GLN	2.8
1	C	231	VAL	2.8
1	D	177	THR	2.7
1	C	171	GLY	2.6
1	C	210	THR	2.6
1	A	233	VAL	2.6
1	B	190	GLN	2.5
1	D	152	GLU	2.5
1	A	148	ASP	2.5
1	D	145	ALA	2.5
1	A	176	LEU	2.4
1	C	147	GLU	2.4
1	C	150	LYS	2.4
1	C	179	ASN	2.4
1	D	183	THR	2.4
1	D	182	TYR	2.4
1	B	205	LYS	2.4
1	B	208	HIS	2.4
1	D	142	VAL	2.4
1	C	200	ARG	2.3
1	D	231	VAL	2.3
1	C	177	THR	2.3
1	D	232	GLU	2.2
1	C	174	ASN	2.2
1	B	86	GLU	2.1
1	C	203	PRO	2.1
1	D	180	ILE	2.1
1	B	177	THR	2.1
1	C	145	ALA	2.1
1	D	150	LYS	2.1
1	B	143	ILE	2.0
1	D	185	ILE	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, 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(\AA^2)	Q<0.9
3	GOL	B	302	6/6	0.80	0.23	7.04	44,50,51,57	0
3	GOL	D	302	6/6	0.75	0.23	6.64	34,40,43,44	0
3	GOL	A	302	6/6	0.77	0.18	2.44	50,51,53,54	0
3	GOL	B	303	6/6	0.82	0.20	0.87	81,82,85,87	0
3	GOL	D	303	6/6	0.90	0.14	0.23	58,61,62,64	0
3	GOL	A	303	6/6	0.86	0.12	0.04	33,38,42,43	0
2	NA	B	301	1/1	0.98	0.06	-2.42	19,19,19,19	0
2	NA	C	301	1/1	0.99	0.06	-3.27	23,23,23,23	0
2	NA	A	301	1/1	0.98	0.06	-3.29	20,20,20,20	0
2	NA	D	301	1/1	0.99	0.05	-6.82	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.