



wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 03:43 am BST

PDB ID : 4M8L
Title : crystal structure of RpiA-R5P complex
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Deposited on : 2013-08-13
Resolution : 2.37 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

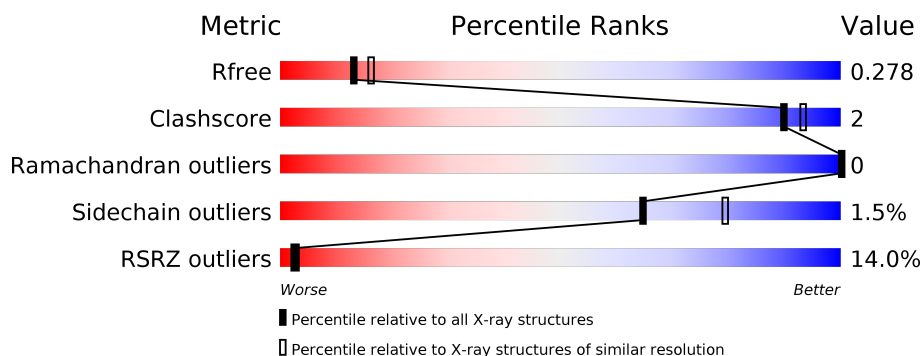
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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}	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 respectively. 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	224	<div> <div>8%</div> <div>97%</div> <div>..</div> </div>
1	B	224	<div> <div>6%</div> <div>93%</div> <div>. .</div> </div>
1	C	224	<div> <div>15%</div> <div>93%</div> <div>. .</div> </div>
1	D	224	<div> <div>25%</div> <div>91%</div> <div>7% .</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	ARF	A	313	-	-	-	X
3	ARF	B	306	-	-	-	X
3	ARF	C	308	-	-	-	X
4	CL	B	315	-	-	X	-

2 Entry composition [i](#)

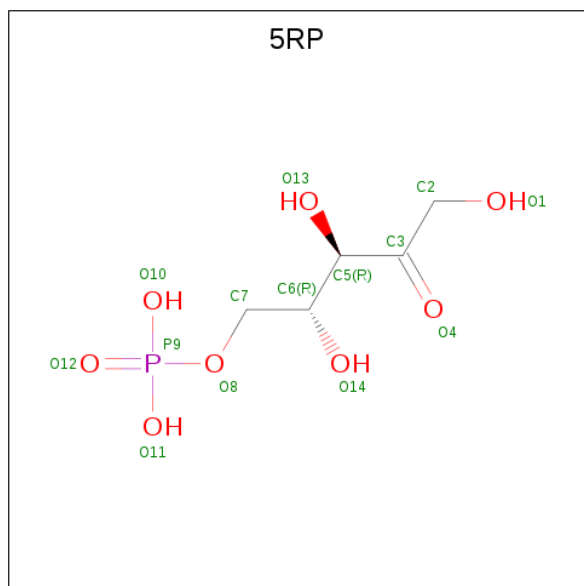
There are 5 unique types of molecules in this entry. The entry contains 7188 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ribose-5-phosphate isomerase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	1	0
			1690	1069	284	332	5			
1	B	217	Total	C	N	O	S	0	2	0
			1675	1059	282	329	5			
1	C	217	Total	C	N	O	S	0	2	0
			1672	1059	279	329	5			
1	D	218	Total	C	N	O	S	0	1	0
			1670	1057	278	330	5			

- Molecule 2 is RIBULOSE-5-PHOSPHATE (three-letter code: 5RP) (formula: C₅H₁₁O₈P).



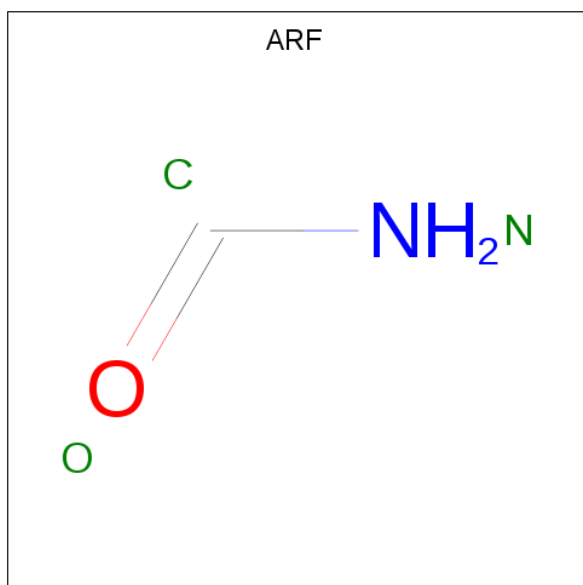
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			14	5	8	1		
2	B	1	Total	C	O	P	0	0
			14	5	8	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	O	P	0	0
			14	5	8	1		
2	D	1	Total	C	O	P	0	0
			14	5	8	1		

- Molecule 3 is FORMAMIDE (three-letter code: ARF) (formula: CH₃NO).



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

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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	C	1	Total 3	C 1	N 1	O 1	0	0
3	D	1	Total 3	C 1	N 1	O 1	0	0
3	D	1	Total 3	C 1	N 1	O 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	104	Total 105	O 105	0	1
5	B	86	Total 88	O 88	0	2
5	C	57	Total 58	O 58	0	1

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	45	Total	O	0	1
			46	46		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.20 Å 151.96 Å 36.29 Å 90.00° 93.98° 90.00°	Depositor
Resolution (Å)	36.75 – 2.37 36.72 – 2.37	Depositor EDS
% Data completeness (in resolution range)	92.9 (36.75-2.37) 93.0 (36.72-2.37)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 2.37 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.237 , 0.274 0.241 , 0.278	Depositor DCC
R_{free} test set	1728 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7188	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 6.32% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: 5RP, ARF, CL

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	0.29	0/1708	0.50	0/2309
1	B	0.30	0/1693	0.52	0/2290
1	C	0.30	0/1690	0.51	0/2287
1	D	0.32	0/1687	0.54	0/2283
All	All	0.30	0/6778	0.52	0/9169

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	104	ALA	Peptide

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	1690	0	1754	3	0
1	B	1675	0	1734	7	0
1	C	1672	0	1732	10	0
1	D	1670	0	1732	10	0
2	A	14	0	9	0	0
2	B	14	0	9	0	0
2	C	14	0	9	0	0
2	D	14	0	9	0	0
3	A	45	0	45	0	0
3	B	39	0	39	1	0
3	C	33	0	33	1	0
3	D	6	0	6	0	0
4	A	1	0	0	0	0
4	B	3	0	0	2	0
4	C	1	0	0	0	0
5	A	105	0	0	0	0
5	B	88	0	0	0	0
5	C	58	0	0	0	0
5	D	46	0	0	0	0
All	All	7188	0	7111	25	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 2.

The worst 5 of 25 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:D:182[B]:ILE:O	1:D:182[B]:ILE:HG13	1.61	0.99
1:C:182[B]:ILE:O	1:C:182[B]:ILE:HG13	1.65	0.93
1:B:49[B]:ARG:HD3	1:B:49[B]:ARG:O	1.77	0.83
1:D:182[B]:ILE:CG1	1:D:182[B]:ILE:O	2.35	0.74
1:A:194:GLN:HE21	1:C:194:GLN:HE21	1.41	0.67

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	219/224 (98%)	213 (97%)	6 (3%)	0	100	100
1	B	217/224 (97%)	211 (97%)	6 (3%)	0	100	100
1	C	217/224 (97%)	210 (97%)	7 (3%)	0	100	100
1	D	217/224 (97%)	211 (97%)	6 (3%)	0	100	100
All	All	870/896 (97%)	845 (97%)	25 (3%)	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	188/191 (98%)	186 (99%)	2 (1%)	73	86
1	B	186/191 (97%)	183 (98%)	3 (2%)	62	78
1	C	186/191 (97%)	185 (100%)	1 (0%)	88	95
1	D	186/191 (97%)	181 (97%)	5 (3%)	44	62
All	All	746/764 (98%)	735 (98%)	11 (2%)	65	79

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	181	LYS
1	C	106	LEU
1	D	106	LEU
1	B	106	LEU
1	D	72	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
1	C	193	ASN
1	C	201	ASN

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 50 ligands modelled in this entry, 5 are monoatomic - leaving 45 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	ARF	B	310	-	2,2,2	0.44	0	1,1,1	0.79	0
3	ARF	B	303	-	2,2,2	0.42	0	1,1,1	0.76	0
3	ARF	B	304	-	2,2,2	0.43	0	1,1,1	0.79	0
3	ARF	C	304	-	2,2,2	0.40	0	1,1,1	0.92	0
3	ARF	B	308	-	2,2,2	0.41	0	1,1,1	0.82	0
3	ARF	C	303	-	2,2,2	0.40	0	1,1,1	0.62	0
3	ARF	C	308	-	2,2,2	0.40	0	1,1,1	0.85	0
3	ARF	A	314	-	2,2,2	0.41	0	1,1,1	0.88	0
3	ARF	A	311	-	2,2,2	0.39	0	1,1,1	0.92	0
3	ARF	D	302	-	2,2,2	0.41	0	1,1,1	0.77	0
3	ARF	C	305	-	2,2,2	0.39	0	1,1,1	0.88	0
3	ARF	A	316	-	2,2,2	0.42	0	1,1,1	0.79	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ARF	B	314	-	2,2,2	0.46	0	1,1,1	0.80	0
3	ARF	A	313	-	2,2,2	0.41	0	1,1,1	0.86	0
3	ARF	A	304	-	2,2,2	0.41	0	1,1,1	0.85	0
3	ARF	A	308	-	2,2,2	0.43	0	1,1,1	0.73	0
3	ARF	C	307	-	2,2,2	0.40	0	1,1,1	0.81	0
3	ARF	C	309	-	2,2,2	0.37	0	1,1,1	0.87	0
3	ARF	B	311	-	2,2,2	0.43	0	1,1,1	0.87	0
3	ARF	C	311	-	2,2,2	0.40	0	1,1,1	0.83	0
3	ARF	A	315	-	2,2,2	0.39	0	1,1,1	0.86	0
3	ARF	A	303	-	2,2,2	0.41	0	1,1,1	0.85	0
2	5RP	B	301	-	11,13,13	0.63	0	11,18,18	0.98	1 (9%)
3	ARF	C	312	-	2,2,2	0.39	0	1,1,1	0.86	0
3	ARF	C	310	-	2,2,2	0.40	0	1,1,1	0.82	0
3	ARF	B	312	-	2,2,2	0.39	0	1,1,1	0.88	0
2	5RP	A	301	-	11,13,13	0.55	0	11,18,18	1.16	0
3	ARF	B	305	-	2,2,2	0.39	0	1,1,1	0.89	0
3	ARF	A	302	-	2,2,2	0.41	0	1,1,1	0.71	0
3	ARF	B	306	-	2,2,2	0.47	0	1,1,1	0.75	0
3	ARF	A	307	-	2,2,2	0.40	0	1,1,1	0.84	0
3	ARF	A	305	-	2,2,2	0.41	0	1,1,1	0.83	0
2	5RP	D	301	-	11,13,13	0.69	0	11,18,18	0.93	0
3	ARF	B	307	-	2,2,2	0.41	0	1,1,1	0.84	0
3	ARF	A	309	-	2,2,2	0.41	0	1,1,1	0.81	0
3	ARF	B	309	-	2,2,2	0.42	0	1,1,1	0.82	0
3	ARF	D	303	-	2,2,2	0.37	0	1,1,1	0.89	0
3	ARF	B	313	-	2,2,2	0.42	0	1,1,1	0.81	0
3	ARF	B	302	-	2,2,2	0.42	0	1,1,1	0.87	0
3	ARF	C	302	-	2,2,2	0.44	0	1,1,1	0.82	0
2	5RP	C	301	-	11,13,13	0.60	0	11,18,18	0.91	0
3	ARF	A	306	-	2,2,2	0.41	0	1,1,1	0.87	0
3	ARF	A	310	-	2,2,2	0.41	0	1,1,1	0.85	0
3	ARF	C	306	-	2,2,2	0.40	0	1,1,1	0.86	0
3	ARF	A	312	-	2,2,2	0.40	0	1,1,1	0.84	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
2	5RP	A	301	-	-	6/16/16/16	-
2	5RP	B	301	-	-	3/16/16/16	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	5RP	C	301	-	-	2/16/16/16	-
2	5RP	D	301	-	-	9/16/16/16	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	5RP	O10-P9-O11	2.01	115.30	107.64

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	5RP	C3-C5-C6-O14
2	A	301	5RP	C3-C5-C6-C7
2	A	301	5RP	O13-C5-C6-O14
2	A	301	5RP	O13-C5-C6-C7
2	D	301	5RP	O1-C2-C3-O4

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	303	ARF	1	0
3	B	314	ARF	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	220/224 (98%)	0.71	18 (8%) 11 12	18, 29, 53, 74	0
1	B	217/224 (96%)	0.73	14 (6%) 18 20	20, 30, 49, 65	0
1	C	217/224 (96%)	0.96	34 (15%) 2 2	22, 39, 72, 96	0
1	D	218/224 (97%)	1.41	56 (25%) 0 0	22, 44, 89, 117	0
All	All	872/896 (97%)	0.95	122 (13%) 2 3	18, 34, 73, 117	0

The worst 5 of 122 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	70	GLY	7.8
1	D	68	ALA	6.8
1	D	66	LEU	6.5
1	D	7	ASN	6.3
1	A	181	LYS	6.1

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. 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	B-factors(\AA^2)	Q<0.9
3	ARF	C	303	3/3	0.50	0.36	38,38,45,46	0
3	ARF	A	313	3/3	0.52	0.53	50,50,54,55	0
3	ARF	A	310	3/3	0.66	0.32	44,44,47,49	0
3	ARF	A	306	3/3	0.67	0.25	54,54,56,56	0
3	ARF	B	313	3/3	0.69	0.32	49,49,54,57	0
3	ARF	D	302	3/3	0.69	0.35	37,37,43,43	0
3	ARF	B	310	3/3	0.69	0.39	38,38,38,41	0
3	ARF	A	304	3/3	0.72	0.25	52,52,53,54	0
3	ARF	A	302	3/3	0.73	0.28	24,24,28,32	0
3	ARF	C	305	3/3	0.74	0.23	56,56,58,59	0
3	ARF	B	311	3/3	0.74	0.27	35,35,40,41	0
3	ARF	C	308	3/3	0.74	0.49	61,61,62,65	0
3	ARF	A	309	3/3	0.75	0.23	49,49,52,53	0
3	ARF	C	310	3/3	0.77	0.23	44,44,45,46	0
3	ARF	B	308	3/3	0.77	0.26	45,45,48,49	0
3	ARF	C	309	3/3	0.78	0.29	38,38,41,42	0
3	ARF	B	306	3/3	0.78	0.42	34,34,36,39	0
3	ARF	C	302	3/3	0.79	0.22	33,33,38,39	0
3	ARF	B	307	3/3	0.80	0.20	49,49,49,52	0
4	CL	B	315	1/1	0.80	0.14	59,59,59,59	0
3	ARF	A	303	3/3	0.81	0.26	40,40,46,47	0
3	ARF	A	308	3/3	0.81	0.27	33,33,39,41	0
3	ARF	B	309	3/3	0.81	0.25	41,41,42,42	0
3	ARF	D	303	3/3	0.81	0.41	32,32,34,37	0
3	ARF	A	307	3/3	0.82	0.29	48,48,52,52	0
2	5RP	D	301	14/14	0.82	0.18	44,51,57,62	0
3	ARF	A	316	3/3	0.82	0.32	38,38,41,41	0
4	CL	B	316	1/1	0.82	0.12	61,61,61,61	0
3	ARF	B	314	3/3	0.82	0.45	43,43,44,49	0
3	ARF	C	306	3/3	0.82	0.23	40,40,45,46	0
3	ARF	B	305	3/3	0.83	0.24	44,44,46,48	0
3	ARF	B	302	3/3	0.83	0.34	36,36,43,44	0
3	ARF	C	312	3/3	0.84	0.14	49,49,53,56	0
3	ARF	C	307	3/3	0.85	0.17	36,36,39,39	0
2	5RP	C	301	14/14	0.86	0.18	35,40,45,46	0
3	ARF	B	303	3/3	0.87	0.22	22,22,25,25	0
3	ARF	C	304	3/3	0.88	0.21	37,37,37,39	0
3	ARF	B	304	3/3	0.91	0.24	40,40,42,44	0
3	ARF	A	314	3/3	0.91	0.17	44,44,46,49	0
4	CL	B	317	1/1	0.91	0.22	60,60,60,60	0
4	CL	C	313	1/1	0.91	0.06	67,67,67,67	0
3	ARF	C	311	3/3	0.92	0.17	52,52,54,57	0
3	ARF	A	311	3/3	0.92	0.25	37,37,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ARF	A	305	3/3	0.92	0.15	37,37,40,40	0
3	ARF	B	312	3/3	0.92	0.17	24,24,26,27	0
4	CL	A	317	1/1	0.94	0.10	44,44,44,44	0
3	ARF	A	312	3/3	0.94	0.16	33,33,36,37	0
2	5RP	A	301	14/14	0.95	0.14	26,31,36,37	0
2	5RP	B	301	14/14	0.95	0.14	29,34,41,42	0
3	ARF	A	315	3/3	0.95	0.18	30,30,34,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.