



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:16 AM GMT

PDB ID : 3OIY  
Title : Helicase domain of reverse gyrase from *Thermotoga maritima*  
Authors : Rudolph, M.G.; Klostermeier, D.  
Deposited on : 2010-08-20  
Resolution : 2.35 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](mailto: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.

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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

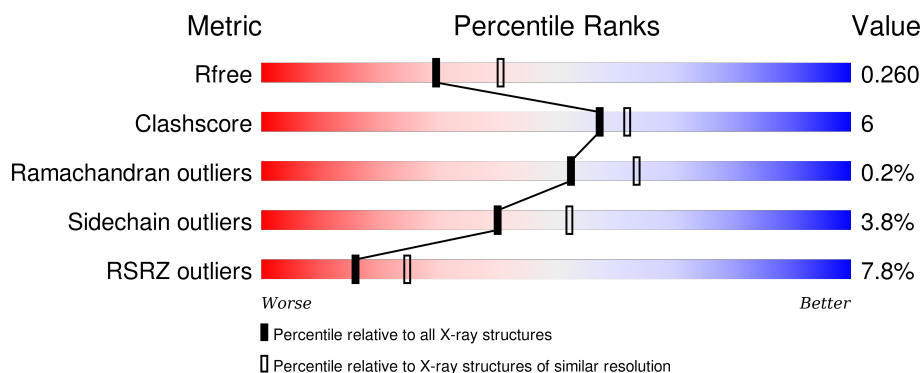
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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  $\geq 3$ , 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	414	<div> <div>5%</div> <div>83%</div> <div>14%</div> <div>••</div> </div>
1	B	414	<div> <div>10%</div> <div>81%</div> <div>14%</div> <div>••</div> </div>

## 2 Entry composition [i](#)

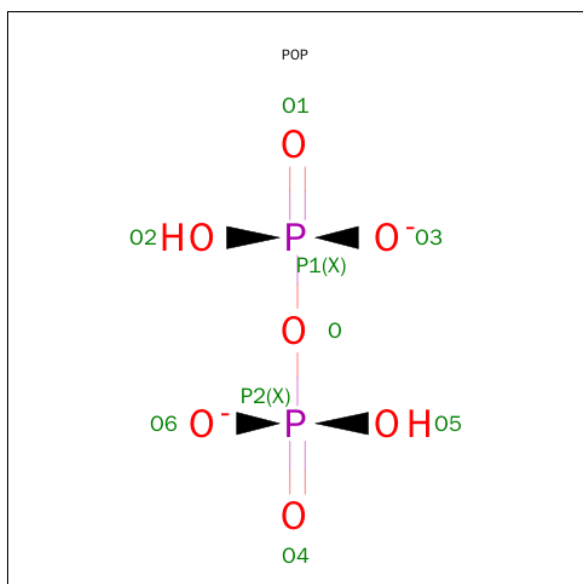
There are 4 unique types of molecules in this entry. The entry contains 6742 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 reverse gyrase helicase domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	0	0
			3343	2164	566	607	6			
1	B	400	Total	C	N	O	S	0	0	0
			3290	2131	555	599	5			

- Molecule 2 is PYROPHOSPHATE 2- (three-letter code: POP) (formula:  $\text{H}_2\text{O}_7\text{P}_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			9	7	2		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula:  $\text{Cl}$ ).

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

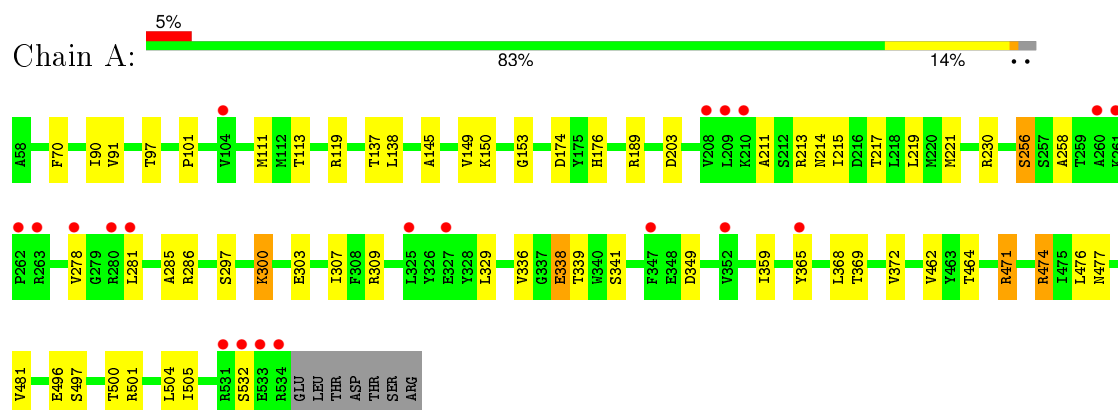
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total 56	O 56	0	0
4	B	43	Total 43	O 43	0	0

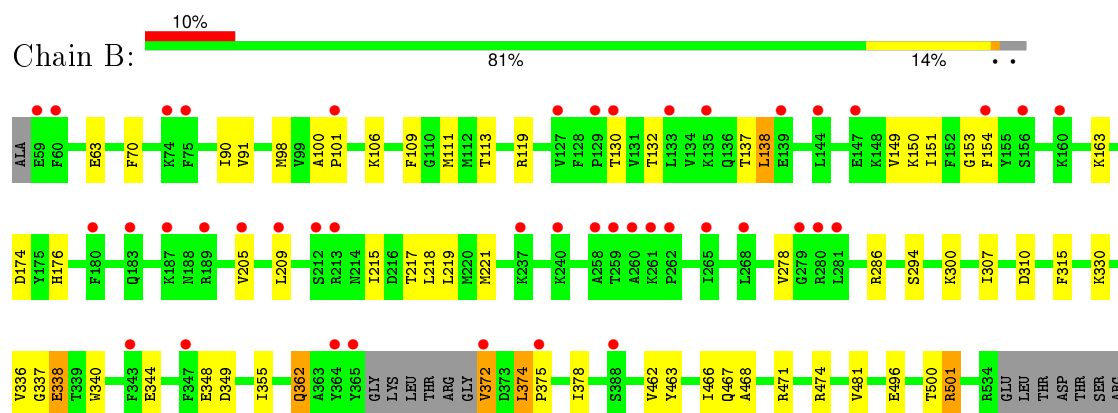
### 3 Residue-property plots

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: reverse gyrase helicase domain



- Molecule 1: reverse gyrase helicase domain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.65Å 126.51Å 132.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.36 – 2.35 44.36 – 2.35	Depositor EDS
% Data completeness (in resolution range)	90.4 (44.36-2.35) 94.1 (44.36-2.35)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.30 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.208 , 0.251 0.220 , 0.260	Depositor DCC
$R_{free}$ test set	2151 reflections (5.65%)	DCC
Wilson B-factor (Å <sup>2</sup> )	54.5	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 47.1	EDS
Estimated twinning fraction	0.014 for -h,l,k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 42481 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6742	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: POP, 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	0.51	0/3409	0.68	0/4576
1	B	0.49	0/3355	0.66	0/4504
All	All	0.50	0/6764	0.67	0/9080

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

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	3343	0	3425	40	0
1	B	3290	0	3355	44	0
2	A	9	0	0	0	0
3	A	1	0	0	0	0
4	A	56	0	0	1	0
4	B	43	0	0	3	0
All	All	6742	0	6780	80	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 6.

All (80) 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:B:374:LEU:H	1:B:374:LEU:HD12	1.16	1.07
1:A:138:LEU:HD11	1:A:153:GLY:HA3	1.48	0.94
1:B:138:LEU:HD11	1:B:153:GLY:HA3	1.55	0.89
1:B:374:LEU:N	1:B:374:LEU:HD12	1.94	0.83
1:B:154:PHE:CZ	1:B:163:LYS:HE3	2.16	0.80
1:B:90:ILE:HD13	1:B:113:THR:HG21	1.63	0.79
1:A:477:ASN:HB2	1:A:532:SER:O	1.83	0.79
1:A:462:VAL:HG11	1:A:501:ARG:HB3	1.65	0.78
1:B:149:VAL:HG12	1:B:151:ILE:HG13	1.66	0.78
1:B:462:VAL:HG21	1:B:501:ARG:HB3	1.66	0.76
1:A:462:VAL:HG11	1:A:501:ARG:CB	2.16	0.76
1:A:281:LEU:O	1:A:281:LEU:HD12	1.86	0.75
1:A:90:ILE:HG21	1:A:113:THR:HG23	1.74	0.70
1:B:474:ARG:O	1:B:481:VAL:HG22	1.93	0.69
1:B:90:ILE:HG21	1:B:113:THR:HG23	1.79	0.65
1:B:374:LEU:H	1:B:374:LEU:CD1	1.87	0.63
1:A:150:LYS:H	1:A:176:HIS:HD2	1.45	0.63
1:A:462:VAL:HG21	1:A:505:ILE:CD1	2.29	0.62
1:B:154:PHE:CE2	1:B:163:LYS:HE3	2.34	0.61
1:B:150:LYS:H	1:B:176:HIS:HD2	1.47	0.61
1:A:303:GLU:HG3	4:A:560:HOH:O	2.01	0.60
1:A:111:MET:HE2	1:A:137:THR:HG23	1.85	0.59
1:B:315:PHE:HB3	1:B:362:GLN:HG3	1.84	0.59
1:B:174:ASP:HB3	4:B:554:HOH:O	2.03	0.58
1:B:111:MET:HE2	1:B:137:THR:HG23	1.85	0.57
1:A:474:ARG:O	1:A:481:VAL:HG22	2.05	0.57
1:B:337:GLY:HA3	1:B:355:ILE:HD13	1.86	0.56
1:A:189:ARG:HD2	1:A:213:ARG:HD3	1.88	0.55
1:A:505:ILE:HD11	1:B:463:TYR:HD2	1.71	0.55
1:A:278:VAL:O	1:A:278:VAL:HG12	2.06	0.55
1:A:462:VAL:HG11	1:A:501:ARG:HB2	1.89	0.54
1:A:462:VAL:HG21	1:A:505:ILE:HD12	1.89	0.54
1:B:215:ILE:O	1:B:219:LEU:HG	2.09	0.53
1:A:372:VAL:HG23	1:A:471:ARG:HG3	1.91	0.53
1:A:211:ALA:O	1:A:214:ASN:HB2	2.09	0.52
1:A:339:THR:HG21	1:A:368:LEU:HD13	1.92	0.52
1:A:505:ILE:HD11	1:B:463:TYR:CD2	2.44	0.52
1:B:468:ALA:HA	1:B:471:ARG:HD2	1.92	0.51
1:A:215:ILE:O	1:A:219:LEU:HG	2.11	0.50
1:A:90:ILE:HG21	1:A:113:THR:CG2	2.41	0.50
1:A:497:SER:O	1:A:501:ARG:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:THR:HG22	1:A:221:MET:HE2	1.93	0.49
1:A:285:ALA:HB3	1:B:101:PRO:HB3	1.95	0.49
1:B:90:ILE:HD13	1:B:113:THR:CG2	2.38	0.49
1:B:286:ARG:HG2	4:B:557:HOH:O	2.13	0.48
1:A:329:LEU:HD12	1:A:359:ILE:HD11	1.96	0.46
1:A:217:THR:HG22	1:A:221:MET:CE	2.45	0.46
1:B:344:GLU:O	1:B:348:GLU:HG2	2.14	0.46
1:B:496:GLU:O	1:B:500:THR:HG23	2.15	0.46
1:B:374:LEU:HA	1:B:375:PRO:HD3	1.72	0.46
1:B:462:VAL:HG12	1:B:466:ILE:HD12	1.98	0.46
1:A:369:THR:HG21	1:A:464:THR:HG23	1.98	0.45
1:B:90:ILE:HD12	1:B:109:PHE:CE2	2.51	0.45
1:B:294:SER:O	1:B:294:SER:OG	2.31	0.45
1:A:462:VAL:CG1	1:A:501:ARG:CB	2.93	0.45
1:A:496:GLU:O	1:A:500:THR:HG23	2.16	0.45
1:B:149:VAL:HG11	1:B:151:ILE:HD11	1.99	0.45
1:A:297:SER:HB3	1:A:300:LYS:HB2	1.99	0.45
1:A:145:ALA:HB1	1:A:149:VAL:HG13	1.99	0.44
1:A:476:LEU:HD23	1:A:532:SER:OG	2.18	0.43
1:A:462:VAL:CG1	1:A:501:ARG:HB2	2.47	0.43
1:A:97:THR:HG21	1:A:258:ALA:HB2	2.00	0.43
1:B:209:LEU:HD11	1:B:278:VAL:HG11	2.00	0.43
1:B:130:THR:HG22	1:B:132:THR:H	1.83	0.43
1:B:217:THR:HG22	1:B:221:MET:HE2	2.00	0.43
1:B:174:ASP:CB	4:B:554:HOH:O	2.66	0.42
1:B:462:VAL:HG12	1:B:466:ILE:CD1	2.49	0.42
1:B:330:LYS:HD2	1:B:336:VAL:CG2	2.50	0.42
1:B:338:GLU:HG3	1:B:340:TRP:CE2	2.54	0.42
1:A:329:LEU:HB3	1:A:336:VAL:HG21	2.02	0.42
1:B:217:THR:HG22	1:B:221:MET:CE	2.50	0.41
1:A:203:ASP:HA	1:A:256:SER:HB2	2.02	0.41
1:B:100:ALA:HB3	1:B:106:LYS:HE3	2.01	0.41
1:A:501:ARG:HG2	1:B:463:TYR:CZ	2.55	0.41
1:B:372:VAL:HG12	1:B:378:ILE:HD11	2.02	0.41
1:B:463:TYR:O	1:B:467:GLN:HG2	2.20	0.41
1:A:70:PHE:CE1	1:A:119:ARG:HD2	2.56	0.41
1:B:70:PHE:CE1	1:B:119:ARG:HD2	2.55	0.41
1:B:98:MET:HB3	1:B:106:LYS:HE2	2.04	0.40
1:A:338:GLU:OE1	1:A:341:SER:HB3	2.21	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	405/414 (98%)	393 (97%)	11 (3%)	1 (0%)	52	63
1	B	396/414 (96%)	383 (97%)	12 (3%)	1 (0%)	46	55
All	All	801/828 (97%)	776 (97%)	23 (3%)	2 (0%)	52	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	362	GLN
1	A	101	PRO

### 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	363/370 (98%)	349 (96%)	14 (4%)	39	51
1	B	357/370 (96%)	344 (96%)	13 (4%)	42	55
All	All	720/740 (97%)	693 (96%)	27 (4%)	40	52

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	VAL
1	A	174	ASP
1	A	230	ARG
1	A	256	SER

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Mol	Chain	Res	Type
1	A	286	ARG
1	A	300	LYS
1	A	307	ILE
1	A	309	ARG
1	A	338	GLU
1	A	349	ASP
1	A	365	TYR
1	A	471	ARG
1	A	474	ARG
1	A	504	LEU
1	B	63	GLU
1	B	91	VAL
1	B	138	LEU
1	B	205	VAL
1	B	218	LEU
1	B	300	LYS
1	B	307	ILE
1	B	310	ASP
1	B	338	GLU
1	B	349	ASP
1	B	372	VAL
1	B	374	LEU
1	B	501	ARG

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	92	GLN
1	A	176	HIS
1	A	183	GLN
1	A	188	ASN
1	B	92	GLN
1	B	176	HIS
1	B	188	ASN
1	B	214	ASN

### 5.3.3 RNA ⓘ

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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z  > 2$
2	POP	A	1	-	8,8,8	1.60	2 (25%)	13,13,13	1.22	1 (7%)

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	POP	A	1	-	-	0/6/6/6	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	POP	P1-O	2.18	1.66	1.61
2	A	1	POP	P2-O	2.61	1.67	1.61

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	POP	O5-P2-O	2.06	114.14	104.65

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	407/414 (98%)	0.28	20 (4%) 33 48	34, 64, 117, 167	0
1	B	400/414 (96%)	0.69	43 (10%) 8 13	30, 80, 137, 181	0
All	All	807/828 (97%)	0.48	63 (7%) 16 25	30, 72, 131, 181	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	ARG	7.0
1	A	260	ALA	6.2
1	B	144	LEU	5.3
1	B	265	ILE	5.2
1	B	154	PHE	5.1
1	A	532	SER	4.6
1	A	261	LYS	4.5
1	B	213	ARG	4.2
1	A	531	ARG	4.1
1	B	133	LEU	4.0
1	A	278	VAL	3.9
1	A	365	TYR	3.9
1	A	280	ARG	3.9
1	B	129	PRO	3.9
1	B	364	TYR	3.9
1	B	281	LEU	3.8
1	B	262	PRO	3.7
1	B	156	SER	3.7
1	A	533	GLU	3.6
1	B	365	TYR	3.5
1	B	258	ALA	3.5
1	B	280	ARG	3.5
1	A	208	VAL	3.4
1	B	259	THR	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	268	LEU	3.4
1	A	210	LYS	3.3
1	B	375	PRO	3.2
1	A	352	VAL	3.1
1	B	101	PRO	3.0
1	B	372	VAL	3.0
1	B	74	LYS	2.9
1	A	104	VAL	2.9
1	B	212	SER	2.9
1	B	343	PHE	2.8
1	A	325	LEU	2.8
1	B	187	LYS	2.8
1	B	75	PHE	2.7
1	B	189	ARG	2.7
1	B	261	LYS	2.7
1	A	262	PRO	2.7
1	B	183	GLN	2.6
1	A	281	LEU	2.4
1	B	260	ALA	2.4
1	B	147	GLU	2.3
1	B	135	LYS	2.3
1	B	130	THR	2.3
1	B	279	GLY	2.2
1	B	240	LYS	2.2
1	A	347	PHE	2.2
1	B	127	VAL	2.2
1	B	347	PHE	2.2
1	B	205	VAL	2.2
1	B	237	LYS	2.1
1	A	327	GLU	2.1
1	B	209	LEU	2.1
1	B	59	GLU	2.1
1	B	60	PHE	2.1
1	A	209	LEU	2.1
1	B	160	LYS	2.1
1	B	139	GLU	2.1
1	B	180	PHE	2.1
1	A	534	ARG	2.1
1	B	388	SER	2.0

## 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. 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	POP	A	1	9/9	0.93	0.12	-0.64	58,92,139,142	0
3	CL	A	2	1/1	0.89	0.13	-	79,79,79,79	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.