



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:55 pm GMT

PDB ID : 5DA7  
Title : monomeric PCNA bound to a small protein inhibitor  
Authors : Ladner, J.E.; Altieri, A.S.; Kelman, Z.  
Deposited on : 2015-08-19  
Resolution : 2.80 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

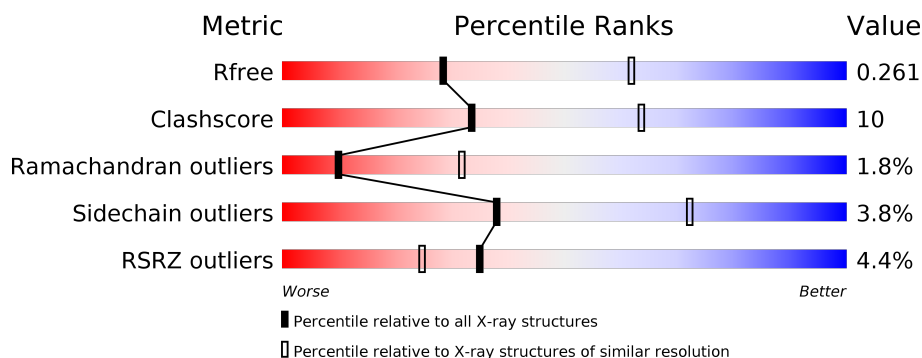
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	255	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>18%</div> <div>••</div> </div> </div>
1	D	255	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>27%</div> <div>••</div> </div> </div>
2	B	64	<div> <div>3%</div> <div> <div></div> <div>44%</div> <div>9%</div> <div>•</div> <div>45%</div> </div> </div>
2	E	64	<div> <div>6%</div> <div> <div></div> <div>36%</div> <div>14%</div> <div>5%</div> <div>45%</div> </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	SO4	A	304	-	-	X	X
3	SO4	A	307	-	-	-	X
3	SO4	E	102	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4598 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 DNA polymerase sliding clamp 1, Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1965	1258	311	388	8			
1	D	247	Total	C	N	O	S	0	0	0
			1965	1258	311	388	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	250	HIS	-	expression tag	UNP Q5JF32
A	251	HIS	-	expression tag	UNP Q5JF32
A	252	HIS	-	expression tag	UNP Q5JF32
A	253	HIS	-	expression tag	UNP Q5JF32
A	254	HIS	-	expression tag	UNP Q5JF32
A	255	HIS	-	expression tag	UNP Q5JF32
D	250	HIS	-	expression tag	UNP Q5JF32
D	251	HIS	-	expression tag	UNP Q5JF32
D	252	HIS	-	expression tag	UNP Q5JF32
D	253	HIS	-	expression tag	UNP Q5JF32
D	254	HIS	-	expression tag	UNP Q5JF32
D	255	HIS	-	expression tag	UNP Q5JF32

- Molecule 2 is a protein called Thermococcales inhibitor of PCNA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	35	Total	C	N	O	0	0	0
			289	185	52	52			
2	E	35	Total	C	N	O	0	0	0
			289	185	52	52			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		



## 4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.34Å 184.08Å 330.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.96 – 2.80 29.96 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.7 (29.96-2.80) 95.7 (29.96-2.80)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.06 (at 2.80Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.201 , 0.261 0.203 , 0.261	Depositor DCC
$R_{free}$ test set	1503 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	78.4	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 64.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4598	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.01% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: SO4

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	0/1992	0.66	1/2679 (0.0%)
1	D	0.49	0/1992	0.69	3/2679 (0.1%)
2	B	0.55	0/294	0.67	0/393
2	E	0.48	0/294	0.72	0/393
All	All	0.50	0/4572	0.68	4/6144 (0.1%)

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.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	246	ARG	O-C-N	7.88	135.30	122.70
1	A	82	ARG	NE-CZ-NH1	-5.95	117.33	120.30
1	D	246	ARG	CA-C-N	-5.34	105.45	117.20
1	D	246	ARG	C-N-CA	-5.04	109.11	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	122	GLU	Peptide

## 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	1965	0	1993	35	0
1	D	1965	0	1993	48	1
2	B	289	0	296	4	0
2	E	289	0	296	7	0
3	A	50	0	0	4	0
3	B	10	0	0	0	0
3	D	20	0	0	2	0
3	E	10	0	0	1	0
All	All	4598	0	4578	89	1

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 10.

All (89) 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:A:4:GLU:OE1	1:A:60:LYS:NZ	2.25	0.68
1:A:213:LYS:HD2	1:A:213:LYS:H	1.59	0.67
2:E:30:LYS:HG3	2:E:31:PRO:HD2	1.76	0.67
1:D:230:GLU:OE1	1:D:238:ARG:NH1	2.29	0.65
1:D:160:LYS:NZ	3:D:303:SO4:O2	2.31	0.64
1:A:39:ARG:NH2	1:A:50:ASP:OD1	2.30	0.63
1:D:87:ASP:HB2	1:D:104:GLU:O	1.99	0.62
1:D:11:LYS:HG2	1:D:84:LYS:O	2.00	0.61
1:A:246:ARG:NE	3:A:304:SO4:O3	2.24	0.61
1:D:109:ARG:HG3	1:D:111:PHE:CE2	2.36	0.60
1:D:39:ARG:NH2	1:D:123:LEU:HD23	2.16	0.60
1:D:11:LYS:HE2	1:D:85:ALA:N	2.16	0.60
1:D:106:THR:OG1	1:D:107:ALA:N	2.36	0.59
1:A:55:GLU:HG2	1:A:56:SER:N	2.18	0.59
1:D:82:ARG:HH11	1:D:111:PHE:HZ	1.50	0.59
1:D:234:ARG:O	1:D:236:GLU:HG2	2.02	0.58
1:A:2:PRO:O	1:A:64:GLU:HG3	2.04	0.57
1:A:112:ARG:NH2	1:D:119:GLU:OE2	2.37	0.57
1:A:155:ILE:HB	1:A:201:ILE:HD11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:LEU:HD23	1:A:16:LEU:C	2.26	0.55
1:A:236:GLU:OE1	1:A:236:GLU:HA	2.06	0.55
1:D:247:VAL:HG22	1:D:248:GLU:H	1.72	0.54
1:D:140:VAL:HG11	1:D:179:LEU:HD11	1.90	0.53
1:D:246:ARG:NE	3:D:304:SO4:O1	2.40	0.53
2:E:54:LYS:NZ	3:E:101:SO4:O3	2.41	0.53
1:D:9:GLY:O	1:D:12:GLU:HB3	2.09	0.53
1:A:22:ASN:HD22	1:A:209:LYS:HE2	1.72	0.53
1:A:30:LYS:HD3	1:A:123:LEU:HD23	1.90	0.53
1:D:109:ARG:CG	1:D:111:PHE:CE2	2.92	0.53
1:A:45:ARG:NH2	3:A:301:SO4:O4	2.41	0.53
1:D:72:ASN:OD1	1:D:75:GLN:HG3	2.09	0.53
1:A:234:ARG:O	1:A:234:ARG:HG3	2.10	0.52
1:D:90:ILE:O	1:D:91:LEU:HD23	2.10	0.52
1:A:90:ILE:HB	1:A:102:THR:HB	1.91	0.51
1:A:94:GLY:HA3	3:A:302:SO4:O2	2.10	0.51
1:D:5:VAL:HG22	1:D:91:LEU:HB2	1.92	0.51
1:D:233:ILE:O	1:D:235:ASP:N	2.44	0.50
2:E:61:ILE:HD12	2:E:61:ILE:H	1.76	0.50
1:D:190:GLU:N	1:D:190:GLU:OE1	2.45	0.50
1:D:19:THR:HG21	1:D:210:GLY:HA3	1.94	0.49
1:D:234:ARG:HB3	1:D:234:ARG:NH1	2.28	0.49
1:A:39:ARG:HH12	1:A:124:GLU:H	1.59	0.49
1:D:213:LYS:HD3	1:D:213:LYS:N	2.27	0.48
1:D:65:GLU:HG2	1:D:66:PRO:HD2	1.95	0.48
1:A:247:VAL:HG22	1:A:248:GLU:H	1.78	0.48
1:D:71:ILE:HG23	1:D:113:LEU:HD21	1.96	0.47
1:D:233:ILE:HG22	1:D:237:GLY:O	2.14	0.47
1:D:233:ILE:O	1:D:233:ILE:HG23	2.14	0.47
1:A:39:ARG:NH1	1:A:124:GLU:H	2.13	0.47
1:A:5:VAL:CG2	1:A:91:LEU:HB2	2.45	0.47
2:E:56:ILE:O	2:E:58:ASN:N	2.47	0.47
1:A:82:ARG:HH22	1:A:109:ARG:NH1	2.13	0.47
1:D:226:PRO:HA	1:D:245:PRO:HD3	1.96	0.47
1:A:167:LYS:HE3	1:A:176:GLU:CD	2.35	0.47
1:D:26:GLU:HG3	2:E:49:LEU:HG	1.97	0.46
1:A:39:ARG:HG2	2:B:46:PHE:HZ	1.79	0.46
1:D:104:GLU:HG3	1:D:108:LYS:HE3	1.98	0.46
1:A:119:GLU:OE2	1:D:112:ARG:NH2	2.49	0.46
1:A:29:PHE:CD2	1:A:38:MET:HB2	2.50	0.46
1:A:46:VAL:HG11	3:A:304:SO4:O2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:103:PHE:O	1:D:108:LYS:HA	2.16	0.45
1:D:12:GLU:CD	1:D:234:ARG:NH1	2.70	0.45
2:E:53:SER:O	2:E:56:ILE:HG22	2.17	0.45
1:A:141:LEU:O	1:A:145:ILE:HD12	2.16	0.45
1:A:41:MET:SD	1:A:45:ARG:HG2	2.57	0.45
1:A:39:ARG:NH1	1:A:123:LEU:HA	2.32	0.44
1:A:87:ASP:N	1:A:87:ASP:OD1	2.38	0.44
1:A:52:ASN:O	1:A:53:LEU:HD23	2.17	0.44
1:D:54:PRO:HD2	1:D:57:ILE:HD11	1.98	0.44
1:D:199:TYR:CZ	1:D:227:LEU:HB2	2.52	0.44
1:D:84:LYS:HB2	1:D:87:ASP:CG	2.38	0.44
1:D:179:LEU:HD13	1:D:186:LEU:HB2	2.00	0.44
1:D:207:MET:SD	1:D:243:LEU:HD22	2.58	0.43
1:D:109:ARG:HG2	1:D:111:PHE:CZ	2.53	0.43
1:D:155:ILE:HD11	1:D:166:MET:HB3	2.00	0.43
1:D:230:GLU:HG2	1:D:240:THR:HG23	1.99	0.43
1:D:233:ILE:C	1:D:235:ASP:N	2.73	0.42
1:A:16:LEU:HD23	1:A:16:LEU:O	2.19	0.42
1:A:22:ASN:ND2	1:A:209:LYS:HE2	2.34	0.42
1:D:59:SER:HB3	1:D:60:LYS:HE2	2.01	0.42
1:D:30:LYS:HD3	1:D:123:LEU:HD12	2.02	0.41
2:B:57:ARG:HG3	2:B:58:ASN:OD1	2.20	0.41
2:B:30:LYS:HA	2:B:31:PRO:HD3	1.92	0.41
1:A:112:ARG:H	2:B:61:ILE:HG22	1.86	0.41
1:D:197:SER:HB3	1:D:198:ALA:H	1.73	0.41
2:E:60:LYS:HG2	2:E:60:LYS:O	2.20	0.41
1:A:16:LEU:C	1:A:16:LEU:CD2	2.89	0.41
1:D:82:ARG:NH1	1:D:111:PHE:HZ	2.19	0.41
1:D:181:LEU:HD12	1:D:181:LEU:H	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:GLU:OE1	1:D:235:ASP:OD2[2_555]	1.76	0.44

## 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	245/255 (96%)	231 (94%)	11 (4%)	3 (1%)	15	44
1	D	245/255 (96%)	227 (93%)	14 (6%)	4 (2%)	11	36
2	B	33/64 (52%)	32 (97%)	1 (3%)	0	100	100
2	E	33/64 (52%)	30 (91%)	0	3 (9%)	1	1
All	All	556/638 (87%)	520 (94%)	26 (5%)	10 (2%)	10	32

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	64	GLU
1	D	123	LEU
1	D	124	GLU
2	E	57	ARG
2	E	61	ILE
1	A	127	GLU
1	D	234	ARG
1	A	59	SER
1	A	64	GLU
2	E	59	ALA

### 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	215/223 (96%)	211 (98%)	4 (2%)	62	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	215/223 (96%)	205 (95%)	10 (5%)	30	64
2	B	32/60 (53%)	30 (94%)	2 (6%)	21	51
2	E	32/60 (53%)	29 (91%)	3 (9%)	10	29
All	All	494/566 (87%)	475 (96%)	19 (4%)	38	72

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	GLU
1	A	84	LYS
1	A	95	ASP
1	A	143	GLU
1	D	8	ASP
1	D	44	SER
1	D	56	SER
1	D	88	THR
1	D	131	THR
1	D	160	LYS
1	D	163	GLU
1	D	196	LYS
1	D	202	ARG
1	D	247	VAL
2	B	60	LYS
2	B	61	ILE
2	E	42	HIS
2	E	54	LYS
2	E	60	LYS

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

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

18 ligands are 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	$\# Z  > 2$
3	SO4	A	301	-	4,4,4	0.16	0	6,6,6	0.19	0
3	SO4	A	302	-	4,4,4	0.20	0	6,6,6	0.35	0
3	SO4	A	303	-	4,4,4	0.21	0	6,6,6	0.18	0
3	SO4	A	304	-	4,4,4	0.32	0	6,6,6	0.57	0
3	SO4	A	305	-	4,4,4	0.28	0	6,6,6	0.44	0
3	SO4	A	306	-	4,4,4	0.21	0	6,6,6	0.24	0
3	SO4	A	307	-	4,4,4	0.23	0	6,6,6	0.30	0
3	SO4	A	308	-	4,4,4	0.25	0	6,6,6	0.43	0
3	SO4	A	309	-	4,4,4	0.19	0	6,6,6	0.13	0
3	SO4	A	310	-	4,4,4	0.24	0	6,6,6	0.10	0
3	SO4	B	101	-	4,4,4	0.15	0	6,6,6	0.24	0
3	SO4	B	102	-	4,4,4	0.19	0	6,6,6	0.24	0
3	SO4	D	301	-	4,4,4	0.20	0	6,6,6	0.11	0
3	SO4	D	302	-	4,4,4	0.20	0	6,6,6	0.30	0
3	SO4	D	303	-	4,4,4	0.18	0	6,6,6	0.11	0
3	SO4	D	304	-	4,4,4	0.25	0	6,6,6	0.14	0
3	SO4	E	101	-	4,4,4	0.22	0	6,6,6	0.11	0
3	SO4	E	102	-	4,4,4	0.20	0	6,6,6	0.09	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	SO4	A	301	-	-	0/0/0/0	0/0/0/0
3	SO4	A	302	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	303	-	-	0/0/0/0	0/0/0/0
3	SO4	A	304	-	-	0/0/0/0	0/0/0/0
3	SO4	A	305	-	-	0/0/0/0	0/0/0/0
3	SO4	A	306	-	-	0/0/0/0	0/0/0/0
3	SO4	A	307	-	-	0/0/0/0	0/0/0/0
3	SO4	A	308	-	-	0/0/0/0	0/0/0/0
3	SO4	A	309	-	-	0/0/0/0	0/0/0/0
3	SO4	A	310	-	-	0/0/0/0	0/0/0/0
3	SO4	B	101	-	-	0/0/0/0	0/0/0/0
3	SO4	B	102	-	-	0/0/0/0	0/0/0/0
3	SO4	D	301	-	-	0/0/0/0	0/0/0/0
3	SO4	D	302	-	-	0/0/0/0	0/0/0/0
3	SO4	D	303	-	-	0/0/0/0	0/0/0/0
3	SO4	D	304	-	-	0/0/0/0	0/0/0/0
3	SO4	E	101	-	-	0/0/0/0	0/0/0/0
3	SO4	E	102	-	-	0/0/0/0	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.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	SO4	1	0
3	A	302	SO4	1	0
3	A	304	SO4	2	0
3	D	303	SO4	1	0
3	D	304	SO4	1	0
3	E	101	SO4	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

### 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	247/255 (96%)	0.05	8 (3%) 48 37	56, 76, 114, 131	0
1	D	247/255 (96%)	0.17	11 (4%) 34 24	68, 90, 117, 130	0
2	B	35/64 (54%)	0.36	2 (5%) 24 16	63, 80, 113, 115	0
2	E	35/64 (54%)	0.12	4 (11%) 6 3	79, 100, 125, 132	0
All	All	564/638 (88%)	0.12	25 (4%) 35 25	56, 86, 117, 132	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	109	ARG	4.5
1	D	107	ALA	3.3
1	D	86	LYS	2.6
2	B	61	ILE	2.6
1	A	104	GLU	2.6
1	D	108	LYS	2.6
2	E	30	LYS	2.5
1	A	126	PRO	2.4
1	A	65	GLU	2.4
1	A	70	GLY	2.4
2	B	57	ARG	2.4
2	E	28	ARG	2.4
1	D	106	THR	2.4
1	D	52	ASN	2.4
1	A	106	THR	2.4
1	D	161	GLU	2.3
1	D	65	GLU	2.2
1	A	235	ASP	2.2
1	D	85	ALA	2.2
1	D	64	GLU	2.2
1	D	182	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	E	42	HIS	2.1
1	A	107	ALA	2.1
2	E	29	LEU	2.0
1	A	105	GLY	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(Å <sup>2</sup> )	Q<0.9
3	SO4	E	102	5/5	0.92	0.35	6.93	107,111,116,120	5
3	SO4	A	304	5/5	0.79	0.34	3.75	64,67,85,88	5
3	SO4	A	307	5/5	0.84	0.34	2.08	94,98,102,121	5
3	SO4	D	304	5/5	0.86	0.21	0.63	97,103,111,113	5
3	SO4	A	305	5/5	0.91	0.22	0.07	78,79,96,104	5
3	SO4	D	302	5/5	0.96	0.10	-1.64	80,85,97,99	0
3	SO4	A	302	5/5	0.96	0.09	-2.62	75,92,93,96	0
3	SO4	E	101	5/5	0.94	0.24	-	110,114,116,122	0
3	SO4	B	102	5/5	0.95	0.23	-	93,94,105,109	0
3	SO4	B	101	5/5	0.92	0.20	-	80,81,90,94	5
3	SO4	A	306	5/5	0.81	0.36	-	98,100,116,118	5
3	SO4	A	301	5/5	0.96	0.23	-	108,110,116,118	0
3	SO4	A	308	5/5	0.90	0.16	-	76,85,96,102	5
3	SO4	A	310	5/5	0.81	0.49	-	78,90,94,99	5
3	SO4	A	303	5/5	0.94	0.28	-	82,88,92,99	5
3	SO4	D	303	5/5	0.91	0.42	-	138,141,151,152	0
3	SO4	A	309	5/5	0.83	0.30	-	113,119,129,131	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SO4	D	301	5/5	0.88	0.29	-	124,124,136,137	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.