



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 24, 2024 – 03:33 AM EDT

PDB ID : 6QH0  
Title : The complex structure of hsRosR-S5 (VNG0258H/RosR-S5)  
Authors : Shaanan, B.; Kutnowski, N.  
Deposited on : 2019-01-14  
Resolution : 2.44 Å(reported)

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We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

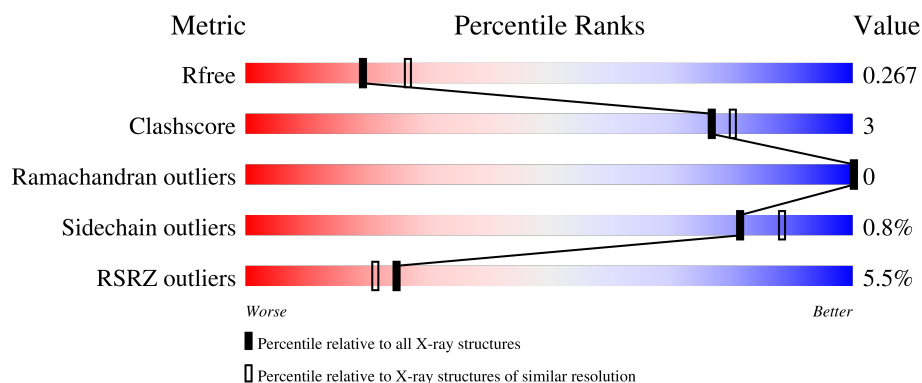
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.44 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	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	116	<div> <div>7%</div> <div>88%</div> <div>11%</div> <div>.</div> </div>
1	B	116	<div> <div>9%</div> <div>91%</div> <div>9%</div> <div>.</div> </div>
1	C	116	<div> <div>2%</div> <div>84%</div> <div>9%</div> <div>5%</div> <div>.</div> </div>
1	D	116	<div> <div>4%</div> <div>91%</div> <div>7%</div> <div>.</div> </div>
2	E	28	<div> <div>7%</div> <div>96%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
2	G	28	<div> <div>4%</div> <div>96%</div> <div>.</div> </div>
3	F	28	<div> <div>4%</div> <div>96%</div> <div>.</div> </div>
3	H	28	<div> <div>4%</div> <div>96%</div> <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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	A	204	-	-	-	X
4	SO4	A	205	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10940 atoms, of which 4780 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 hsRosR DNA binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	115	Total	C	H	N	O	0	0	0
			1805	571	884	158	192			
1	B	116	Total	C	H	N	O	0	2	0
			1833	580	897	160	196			
1	C	110	Total	C	H	N	O	0	0	0
			1736	551	853	150	182			
1	D	114	Total	C	H	N	O	0	0	0
			1793	567	880	157	189			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	ALA	-	expression tag	UNP Q9HSF4
A	118	ALA	-	expression tag	UNP Q9HSF4
A	119	ALA	-	expression tag	UNP Q9HSF4
A	120	LEU	-	expression tag	UNP Q9HSF4
A	121	GLU	-	expression tag	UNP Q9HSF4
B	117	ALA	-	expression tag	UNP Q9HSF4
B	118	ALA	-	expression tag	UNP Q9HSF4
B	119	ALA	-	expression tag	UNP Q9HSF4
B	120	LEU	-	expression tag	UNP Q9HSF4
B	121	GLU	-	expression tag	UNP Q9HSF4
C	117	ALA	-	expression tag	UNP Q9HSF4
C	118	ALA	-	expression tag	UNP Q9HSF4
C	119	ALA	-	expression tag	UNP Q9HSF4
C	120	LEU	-	expression tag	UNP Q9HSF4
C	121	GLU	-	expression tag	UNP Q9HSF4
D	117	ALA	-	expression tag	UNP Q9HSF4
D	118	ALA	-	expression tag	UNP Q9HSF4
D	119	ALA	-	expression tag	UNP Q9HSF4
D	120	LEU	-	expression tag	UNP Q9HSF4
D	121	GLU	-	expression tag	UNP Q9HSF4

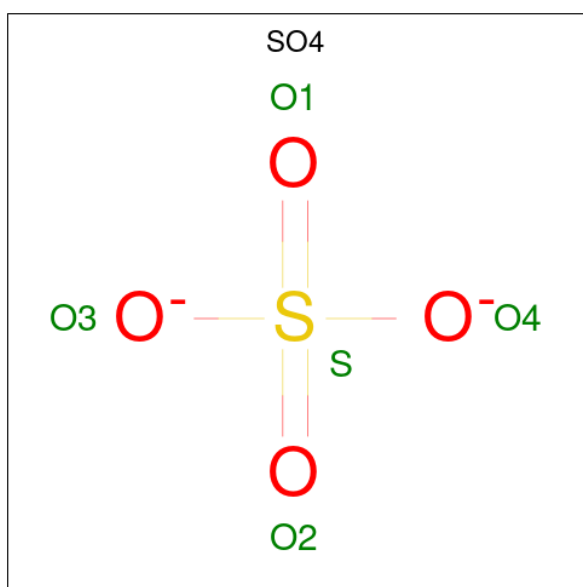
- Molecule 2 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	28	Total	C	H	N	O	P	0	0	0
			895	275	315	112	165	28			
2	G	28	Total	C	H	N	O	P	0	0	0
			895	275	315	112	165	28			

- Molecule 3 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	F	28	Total	C	H	N	O	P	0	0	0
			886	272	318	97	171	28			
3	H	28	Total	C	H	N	O	P	0	0	0
			886	272	318	97	171	28			

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



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total 1	Mn 1	0	0
5	C	2	Total 2	Mn 2	0	0
5	D	1	Total 1	Mn 1	0	0

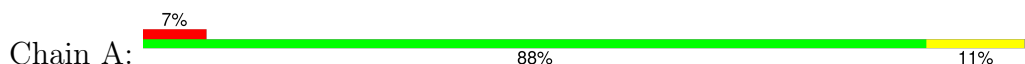
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	13	Total 13	O 13	0	0
6	B	14	Total 14	O 14	0	0
6	C	16	Total 16	O 16	0	0
6	D	14	Total 14	O 14	0	0
6	E	6	Total 6	O 6	0	0
6	F	6	Total 6	O 6	0	0
6	G	10	Total 10	O 10	0	0
6	H	12	Total 12	O 12	0	0

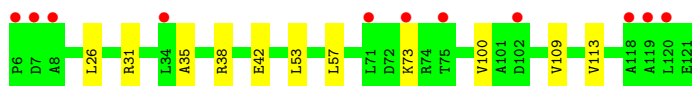
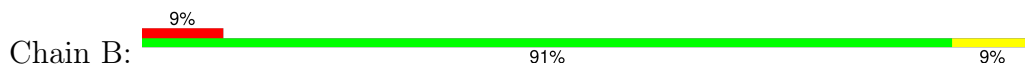
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

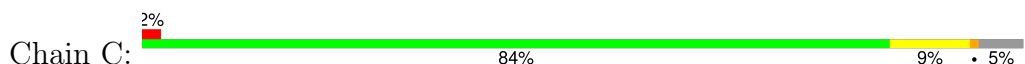
- Molecule 1: hsRosR DNA binding protein



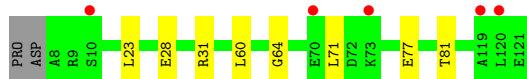
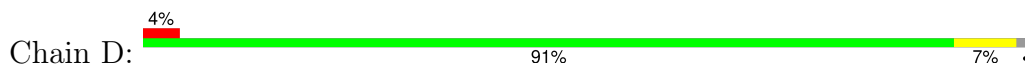
- Molecule 1: hsRosR DNA binding protein



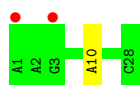
- Molecule 1: hsRosR DNA binding protein



- Molecule 1: hsRosR DNA binding protein

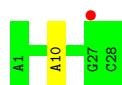


- Molecule 2: DNA (28-MER)

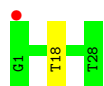


- Molecule 2: DNA (28-MER)

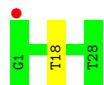




- Molecule 3: DNA (28-MER)



- Molecule 3: DNA (28-MER)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.55Å 83.11Å 89.60Å 95.81° 95.50° 105.11°	Depositor
Resolution (Å)	44.57 – 2.44 44.57 – 2.44	Depositor EDS
% Data completeness (in resolution range)	57.5 (44.57-2.44) 57.5 (44.57-2.44)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.18	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.31 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.239 , 0.268 0.239 , 0.267	Depositor DCC
$R_{free}$ test set	1356 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 58.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	10940	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, MN

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	0.26	0/932	0.43	0/1260
1	B	0.25	0/955	0.42	0/1292
1	C	0.26	0/894	0.41	0/1209
1	D	0.26	0/924	0.41	0/1249
2	E	0.53	0/652	0.89	0/1005
2	G	0.54	0/652	0.90	0/1005
3	F	0.54	0/634	0.95	0/975
3	H	0.52	0/634	0.95	0/975
All	All	0.39	0/6277	0.69	0/8970

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	921	884	884	6	0
1	B	936	897	888	10	0
1	C	883	853	853	7	0
1	D	913	880	880	5	0
2	E	580	315	315	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	580	315	315	1	0
3	F	568	318	318	1	0
3	H	568	318	318	1	0
4	A	25	0	0	1	0
4	B	30	0	0	1	0
4	C	15	0	0	0	0
4	D	25	0	0	0	0
4	E	5	0	0	0	0
4	F	5	0	0	0	0
4	G	5	0	0	0	0
4	H	5	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
5	D	1	0	0	0	0
6	A	13	0	0	0	0
6	B	14	0	0	1	0
6	C	16	0	0	1	0
6	D	14	0	0	1	0
6	E	6	0	0	0	0
6	F	6	0	0	0	0
6	G	10	0	0	0	0
6	H	12	0	0	0	0
All	All	6160	4780	4771	30	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.

The worst 5 of 30 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:23:LEU:HD21	1:D:60:LEU:HD22	1.53	0.91
1:A:26:LEU:HD21	1:A:36:ILE:HD11	1.64	0.79
1:A:67:GLU:HG2	1:A:81:THR:HG22	1.79	0.64
1:D:28:GLU:OE2	1:D:31:ARG:NH1	2.30	0.64
1:B:38:ARG:NH1	4:B:201:SO4:O3	2.30	0.63

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	113/116 (97%)	109 (96%)	4 (4%)	0	100	100
1	B	116/116 (100%)	112 (97%)	4 (3%)	0	100	100
1	C	108/116 (93%)	103 (95%)	5 (5%)	0	100	100
1	D	112/116 (97%)	108 (96%)	4 (4%)	0	100	100
All	All	449/464 (97%)	432 (96%)	17 (4%)	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	96/97 (99%)	94 (98%)	2 (2%)	53	66
1	B	99/97 (102%)	99 (100%)	0	100	100
1	C	92/97 (95%)	91 (99%)	1 (1%)	73	83
1	D	95/97 (98%)	95 (100%)	0	100	100
All	All	382/388 (98%)	379 (99%)	3 (1%)	81	88

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	106	ARG
1	C	77	GLU

Sometimes 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 5 are monoatomic - leaving 23 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	$\# Z  > 2$
4	SO4	A	201	-	4,4,4	0.23	0	6,6,6	0.06	0
4	SO4	D	201	-	4,4,4	0.24	0	6,6,6	0.09	0
4	SO4	C	202	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	D	203	-	4,4,4	0.24	0	6,6,6	0.06	0
4	SO4	B	202	-	4,4,4	0.24	0	6,6,6	0.07	0
4	SO4	D	205	-	4,4,4	0.23	0	6,6,6	0.07	0
4	SO4	B	201	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	F	101	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	D	202	-	4,4,4	0.25	0	6,6,6	0.07	0
4	SO4	A	203	-	4,4,4	0.24	0	6,6,6	0.07	0
4	SO4	H	101	-	4,4,4	0.23	0	6,6,6	0.09	0
4	SO4	B	203	-	4,4,4	0.23	0	6,6,6	0.12	0
4	SO4	B	206	-	4,4,4	0.25	0	6,6,6	0.07	0
4	SO4	B	205	-	4,4,4	0.24	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	SO4	A	202	-	4,4,4	0.24	0	6,6,6	0.07	0
4	SO4	C	201	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	E	101	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	B	204	-	4,4,4	0.23	0	6,6,6	0.11	0
4	SO4	C	203	-	4,4,4	0.23	0	6,6,6	0.07	0
4	SO4	A	204	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	A	205	-	4,4,4	0.24	0	6,6,6	0.06	0
4	SO4	D	204	-	4,4,4	0.24	0	6,6,6	0.06	0
4	SO4	G	101	-	4,4,4	0.24	0	6,6,6	0.07	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	201	SO4	1	0
4	A	203	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 [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, 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	115/116 (99%)	0.67	8 (6%) 16 12	14, 36, 71, 89	0
1	B	116/116 (100%)	0.77	11 (9%) 8 5	19, 44, 98, 158	0
1	C	110/116 (94%)	0.60	2 (1%) 68 64	16, 35, 62, 78	0
1	D	114/116 (98%)	0.75	5 (4%) 34 32	18, 45, 84, 107	0
2	E	28/28 (100%)	0.73	2 (7%) 16 12	49, 81, 114, 128	0
2	G	28/28 (100%)	0.40	1 (3%) 42 39	43, 69, 88, 105	0
3	F	28/28 (100%)	0.66	1 (3%) 42 39	49, 83, 91, 100	0
3	H	28/28 (100%)	0.34	1 (3%) 42 39	39, 69, 84, 98	0
All	All	567/576 (98%)	0.67	31 (5%) 25 21	14, 45, 89, 158	0

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	71	LEU	5.3
2	E	1	DA	3.9
1	B	7	ASP	3.5
1	A	71	LEU	3.3
1	B	6	PRO	3.2

### 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 monosaccharides 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
4	SO4	A	205	5/5	0.74	0.64	84,85,85,85	5
4	SO4	D	201	5/5	0.78	0.32	126,126,126,126	0
4	SO4	A	203	5/5	0.79	0.19	105,105,106,106	0
4	SO4	A	204	5/5	0.79	0.59	134,135,135,135	0
4	SO4	B	201	5/5	0.80	0.21	124,124,125,125	0
4	SO4	B	203	5/5	0.83	0.30	106,106,106,107	0
4	SO4	A	201	5/5	0.84	0.30	140,140,140,140	0
4	SO4	B	202	5/5	0.85	0.22	104,104,104,105	0
4	SO4	H	101	5/5	0.85	0.22	97,97,98,98	0
4	SO4	B	204	5/5	0.87	0.24	110,110,110,110	0
4	SO4	C	203	5/5	0.87	0.26	110,110,111,111	0
4	SO4	B	206	5/5	0.88	0.22	57,57,57,58	5
4	SO4	D	204	5/5	0.90	0.21	110,110,110,110	0
5	MN	A	206	1/1	0.90	0.16	67,67,67,67	0
4	SO4	C	202	5/5	0.91	0.21	117,117,117,117	0
4	SO4	B	205	5/5	0.91	0.24	110,110,111,111	0
4	SO4	C	201	5/5	0.91	0.24	96,96,97,97	0
4	SO4	A	202	5/5	0.92	0.20	95,95,96,96	0
5	MN	C	204	1/1	0.92	0.12	81,81,81,81	1
4	SO4	D	202	5/5	0.93	0.18	57,57,58,58	5
4	SO4	D	203	5/5	0.95	0.20	44,44,44,45	5
4	SO4	D	205	5/5	0.95	0.15	94,95,95,95	0
4	SO4	G	101	5/5	0.95	0.13	71,71,71,72	0
4	SO4	E	101	5/5	0.96	0.08	78,78,79,79	0
5	MN	B	207	1/1	0.96	0.08	25,25,25,25	1
4	SO4	F	101	5/5	0.96	0.21	83,83,84,84	0
5	MN	C	205	1/1	0.97	0.14	58,58,58,58	1
5	MN	D	206	1/1	0.98	0.05	19,19,19,19	1

## 6.5 Other polymers ⓘ

There are no such residues in this entry.