



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:34 AM GMT

PDB ID : 2HRO  
Title : Structure of the full-length Enzyme I of the PTS system from *Staphylococcus carnosus*  
Authors : Marquez, J.A.; Reinelt, S.; Koch, B.; Engelman, R.; Hengstenberg, W.; Schefzek, K.  
Deposited on : 2006-07-20  
Resolution : 2.50 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

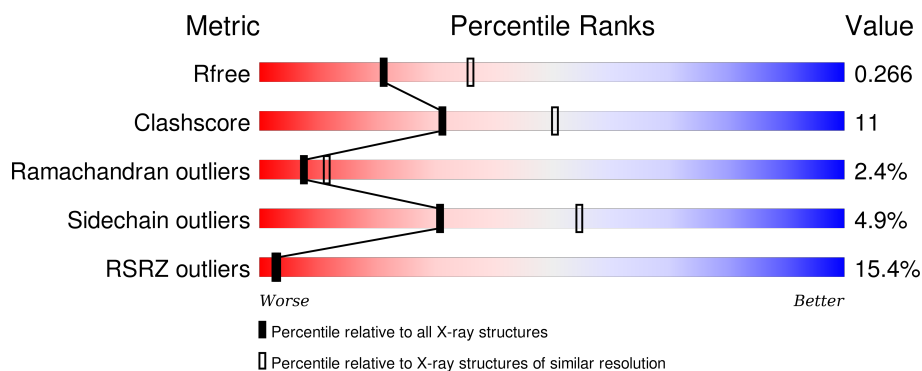
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	573	<div> <div>15%</div> <div>70%</div> <div>24%</div> <div>• •</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4247 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 Phosphoenolpyruvate-protein phosphotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	0
			4070	2551	697	803	19			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	MET	ILE	SEE REMARK 999	UNP P23533
A	?	-	LYS	SEE REMARK 999	UNP P23533
A	406	LEU	PHE	SEE REMARK 999	UNP P23533
A	408	GLU	LYS	SEE REMARK 999	UNP P23533
A	409	GLU	LYS	SEE REMARK 999	UNP P23533
A	410	GLU	ASN	SEE REMARK 999	UNP P23533
A	411	ARG	VAL	SEE REMARK 999	UNP P23533
A	412	ALA	LEU	SEE REMARK 999	UNP P23533
A	413	ASN	THR	SEE REMARK 999	UNP P23533
A	416	ASN	MET	SEE REMARK 999	UNP P23533
A	417	GLU	LYS	SEE REMARK 999	UNP P23533
A	418	GLY	ALA	SEE REMARK 999	UNP P23533
A	419	TYR	MET	SEE REMARK 999	UNP P23533
A	420	GLU	LYS	SEE REMARK 999	UNP P23533
A	479	ASN	ILE	SEE REMARK 999	UNP P23533
A	480	PRO	SER	SEE REMARK 999	UNP P23533
A	481	ALA	ASN	SEE REMARK 999	UNP P23533
A	482	ILE	PHE	SEE REMARK 999	UNP P23533
A	483	LEU	SER	SEE REMARK 999	UNP P23533
A	484	ARG	PHE	SEE REMARK 999	UNP P23533
A	536	ARG	VAL	SEE REMARK 999	UNP P23533

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

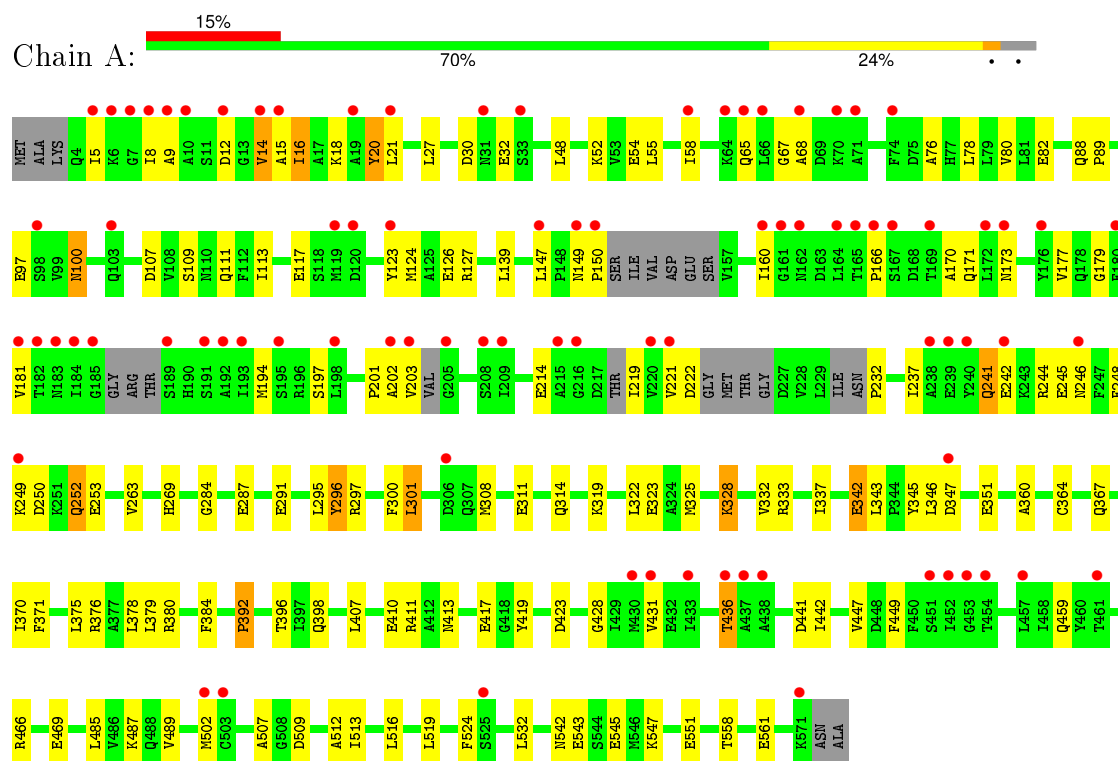
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	172	Total	O	0	0
			172	172		

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

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: Phosphoenolpyruvate-protein phosphotransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	173.36Å 46.85Å 85.30Å 90.00° 101.46° 90.00°	Depositor
Resolution (Å)	34.86 – 2.50 34.86 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (34.86-2.50) 98.9 (34.86-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.15 (at 2.39Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.215 , 0.267 0.216 , 0.266	Depositor DCC
$R_{free}$ test set	1179 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	43.0	Xtriage
Anisotropy	0.972	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 50.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 26420 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4247	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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 [i](#)

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

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	0.35	0/4119	0.59	3/5572 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	201	PRO	N-CA-CB	5.74	110.18	103.30
1	A	166	PRO	N-CA-CB	5.62	110.04	103.30
1	A	232	PRO	N-CA-CB	5.59	110.02	103.30

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	4070	0	3828	85	0
2	A	5	0	0	0	0
3	A	172	0	0	0	0
All	All	4247	0	3828	85	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 11.

All (85) 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:396:THR:HG22	1:A:398:GLN:H	1.42	0.85
1:A:345:TYR:O	1:A:346:LEU:HB2	1.80	0.79
1:A:558:THR:OG1	1:A:561:GLU:HG3	1.81	0.78
1:A:428:GLY:HA2	1:A:447:VAL:HG13	1.68	0.76
1:A:54:GLU:O	1:A:58:ILE:HG12	1.92	0.69
1:A:547:LYS:O	1:A:551:GLU:HG3	1.94	0.68
1:A:411:ARG:NH2	1:A:423:ASP:HA	2.08	0.68
1:A:65:GLN:HE22	1:A:173:ASN:HA	1.60	0.66
1:A:431:VAL:HG13	1:A:436:THR:CG2	2.26	0.66
1:A:411:ARG:HH21	1:A:423:ASP:HA	1.60	0.65
1:A:246:ASN:HA	1:A:249:LYS:HE3	1.79	0.64
1:A:367:GLN:O	1:A:370:ILE:HG22	1.99	0.62
1:A:100:ASN:C	1:A:100:ASN:HD22	2.04	0.60
1:A:149:ASN:HB2	1:A:150:PRO:HD2	1.86	0.58
1:A:8:ILE:HA	1:A:203:VAL:HA	1.85	0.58
1:A:447:VAL:HG12	1:A:449:PHE:H	1.69	0.58
1:A:345:TYR:HD1	1:A:346:LEU:HD13	1.69	0.57
1:A:428:GLY:HA2	1:A:447:VAL:CG1	2.36	0.56
1:A:308:MET:SD	1:A:370:ILE:HD12	2.46	0.56
1:A:396:THR:HG22	1:A:398:GLN:N	2.16	0.55
1:A:14:VAL:H	1:A:222:ASP:HA	1.72	0.55
1:A:509:ASP:O	1:A:513:ILE:HG13	2.07	0.54
1:A:297:ARG:HG2	1:A:333:ARG:NH2	2.23	0.53
1:A:487:LYS:HB2	1:A:519:LEU:HD22	1.90	0.53
1:A:325:MET:O	1:A:328:LYS:HB2	2.10	0.52
1:A:380:ARG:HD3	1:A:419:TYR:CD2	2.45	0.51
1:A:78:LEU:O	1:A:82:GLU:HG2	2.11	0.50
1:A:300:PHE:CE2	1:A:301:LEU:HD22	2.46	0.50
1:A:345:TYR:CD1	1:A:346:LEU:HD13	2.46	0.50
1:A:150:PRO:HB3	1:A:171:GLN:O	2.10	0.50
1:A:30:ASP:O	1:A:32:GLU:HG3	2.13	0.49
1:A:21:LEU:HA	1:A:160:ILE:O	2.12	0.49
1:A:301:LEU:HG	1:A:314:GLN:HG2	1.94	0.49
1:A:337:ILE:HD13	1:A:343:LEU:HD11	1.93	0.49
1:A:332:VAL:HG11	1:A:378:LEU:HD21	1.95	0.49
1:A:67:GLY:O	1:A:68:ALA:HB3	2.13	0.48
1:A:113:ILE:O	1:A:117:GLU:HG2	2.13	0.48
1:A:237:ILE:O	1:A:241:GLN:HB2	2.14	0.47
1:A:14:VAL:H	1:A:222:ASP:CA	2.26	0.47
1:A:109:SER:O	1:A:113:ILE:HG13	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ASN:C	1:A:100:ASN:ND2	2.68	0.47
1:A:76:ALA:O	1:A:80:VAL:HG23	2.14	0.46
1:A:245:GLU:O	1:A:249:LYS:HB2	2.15	0.46
1:A:88:GLN:HB2	1:A:89:PRO:HD3	1.97	0.46
1:A:319:LYS:HE2	1:A:384:PHE:CE2	2.51	0.46
1:A:542:ASN:HB3	1:A:545:GLU:HB2	1.97	0.46
1:A:431:VAL:HG13	1:A:436:THR:HG22	1.97	0.45
1:A:20:TYR:H	1:A:21:LEU:HD12	1.82	0.45
1:A:328:LYS:HA	1:A:328:LYS:HE2	1.99	0.45
1:A:459:GLN:OE1	1:A:466:ARG:HG2	2.16	0.45
1:A:248:PHE:O	1:A:252:GLN:HB2	2.17	0.45
1:A:65:GLN:NE2	1:A:173:ASN:HA	2.31	0.45
1:A:15:ALA:HB1	1:A:179:GLY:HA2	1.99	0.44
1:A:244:ARG:C	1:A:246:ASN:H	2.21	0.44
1:A:55:LEU:HB2	1:A:78:LEU:HD21	1.99	0.44
1:A:364:CYS:HB3	1:A:371:PHE:CG	2.52	0.44
1:A:124:MET:SD	1:A:127:ARG:HD3	2.57	0.44
1:A:194:MET:HA	1:A:197:SER:CB	2.48	0.44
1:A:502:MET:HG2	1:A:507:ALA:HB2	1.99	0.44
1:A:360:ALA:HB1	1:A:392:PRO:HG2	2.00	0.44
1:A:263:VAL:HG22	1:A:269:HIS:CD2	2.53	0.43
1:A:48:LEU:HD11	1:A:52:LYS:HE3	2.00	0.43
1:A:48:LEU:CD1	1:A:52:LYS:HE3	2.48	0.43
1:A:323:GLU:HG3	1:A:384:PHE:HB3	2.00	0.43
1:A:16:ILE:HA	1:A:219:ILE:O	2.17	0.43
1:A:27:LEU:HB3	1:A:139:LEU:CD1	2.49	0.42
1:A:447:VAL:HG12	1:A:449:PHE:N	2.32	0.42
1:A:123:TYR:O	1:A:126:GLU:HB3	2.19	0.42
1:A:311:GLU:OE1	1:A:376:ARG:HD3	2.20	0.42
1:A:512:ALA:O	1:A:516:LEU:HG	2.19	0.42
1:A:21:LEU:HD12	1:A:21:LEU:N	2.35	0.42
1:A:107:ASP:O	1:A:111:GLN:HG2	2.20	0.41
1:A:547:LYS:HE3	1:A:547:LYS:HB2	1.87	0.41
1:A:524:PHE:CD1	1:A:524:PHE:N	2.88	0.41
1:A:296:TYR:HB3	1:A:332:VAL:HA	2.01	0.41
1:A:469:GLU:OE1	1:A:469:GLU:HA	2.20	0.41
1:A:249:LYS:O	1:A:253:GLU:HG3	2.20	0.41
1:A:485:LEU:O	1:A:489:VAL:HG23	2.21	0.41
1:A:291:GLU:O	1:A:328:LYS:HG3	2.20	0.41
1:A:9:ALA:O	1:A:202:ALA:HB3	2.21	0.41
1:A:413:ASN:O	1:A:417:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:GLY:HA2	1:A:287:GLU:HG2	2.03	0.40
1:A:375:LEU:HD22	1:A:407:LEU:CD1	2.51	0.40
1:A:410:GLU:HA	1:A:410:GLU:OE1	2.21	0.40
1:A:441:ASP:OD2	1:A:442:ILE:N	2.54	0.40

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	537/573 (94%)	460 (86%)	64 (12%)	13 (2%)	<b>7</b>	<b>11</b>

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	VAL
1	A	342	GLU
1	A	12	ASP
1	A	18	LYS
1	A	20	TYR
1	A	170	ALA
1	A	351	GLU
1	A	14	VAL
1	A	177	VAL
1	A	181	VAL
1	A	16	ILE
1	A	214	GLU
1	A	5	ILE

### 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	391/481 (81%)	372 (95%)	19 (5%)	31 55

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

Mol	Chain	Res	Type
1	A	97	GLU
1	A	100	ASN
1	A	147	LEU
1	A	241	GLN
1	A	242	GLU
1	A	250	ASP
1	A	252	GLN
1	A	295	LEU
1	A	296	TYR
1	A	301	LEU
1	A	322	LEU
1	A	328	LYS
1	A	342	GLU
1	A	347	ASP
1	A	379	LEU
1	A	392	PRO
1	A	436	THR
1	A	532	LEU
1	A	543	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	65	GLN
1	A	100	ASN
1	A	103	GLN
1	A	121	ASN
1	A	353	ASN

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Mol	Chain	Res	Type
1	A	416	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](#)

1 ligand is 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	$\# Z  > 2$
2	SO4	A	900	-	4,4,4	0.13	0	6,6,6	0.10	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	SO4	A	900	-	-	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.

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	551/573 (96%)	0.76	85 (15%) 3 3	29, 58, 132, 139	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	15	ALA	8.5
1	A	161	GLY	7.5
1	A	123	TYR	5.9
1	A	202	ALA	5.9
1	A	191	SER	5.8
1	A	9	ALA	5.8
1	A	203	VAL	5.7
1	A	180	PHE	5.5
1	A	215	ALA	5.0
1	A	5	ILE	5.0
1	A	164	LEU	4.8
1	A	192	ALA	4.7
1	A	169	THR	4.6
1	A	10	ALA	4.5
1	A	65	GLN	4.4
1	A	64	LYS	4.4
1	A	7	GLY	4.3
1	A	19	ALA	4.0
1	A	165	THR	4.0
1	A	167	SER	4.0
1	A	31	ASN	3.9
1	A	185	GLY	3.9
1	A	21	LEU	3.9
1	A	246	ASN	3.6
1	A	208	SER	3.6
1	A	431	VAL	3.6
1	A	216	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	240	TYR	3.4
1	A	181	VAL	3.3
1	A	184	ILE	3.3
1	A	239	GLU	3.3
1	A	58	ILE	3.3
1	A	430	MET	3.2
1	A	172	LEU	3.2
1	A	193	ILE	3.2
1	A	71	ALA	3.1
1	A	238	ALA	3.1
1	A	249	LYS	3.1
1	A	74	PHE	3.1
1	A	8	ILE	3.1
1	A	66	LEU	3.0
1	A	221	VAL	3.0
1	A	453	GLY	2.9
1	A	452	ILE	2.9
1	A	68	ALA	2.8
1	A	14	VAL	2.8
1	A	205	GLY	2.8
1	A	242	GLU	2.7
1	A	198	LEU	2.7
1	A	457	LEU	2.7
1	A	182	THR	2.7
1	A	183	ASN	2.7
1	A	70	LYS	2.7
1	A	162	ASN	2.7
1	A	166	PRO	2.6
1	A	120	ASP	2.6
1	A	173	ASN	2.5
1	A	12	ASP	2.5
1	A	195	SER	2.5
1	A	437	ALA	2.5
1	A	502	MET	2.5
1	A	147	LEU	2.5
1	A	160	ILE	2.5
1	A	149	ASN	2.5
1	A	189	SER	2.5
1	A	438	ALA	2.5
1	A	150	PRO	2.4
1	A	98	SER	2.4
1	A	454	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	306	ASP	2.3
1	A	436	THR	2.2
1	A	433	ILE	2.2
1	A	571	LYS	2.2
1	A	33	SER	2.1
1	A	176	TYR	2.1
1	A	6	LYS	2.1
1	A	525	SER	2.1
1	A	220	VAL	2.1
1	A	119	MET	2.1
1	A	209	ILE	2.1
1	A	461	THR	2.1
1	A	451	SER	2.1
1	A	347	ASP	2.1
1	A	103	GLN	2.0
1	A	503	CYS	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	SO4	A	900	5/5	0.96	0.29	1.02	63,64,66,66	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.