



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 12, 2024 – 07:43 PM EST

PDB ID : 4HY9  
Title : Crystal structure of the substrate binding domain of E.coli DnaK in complex with pyrrhocoricin\_LYZZ (residues 1 to 11)  
Authors : Zahn, M.; Straeter, N.  
Deposited on : 2012-11-13  
Resolution : 1.55 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

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	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

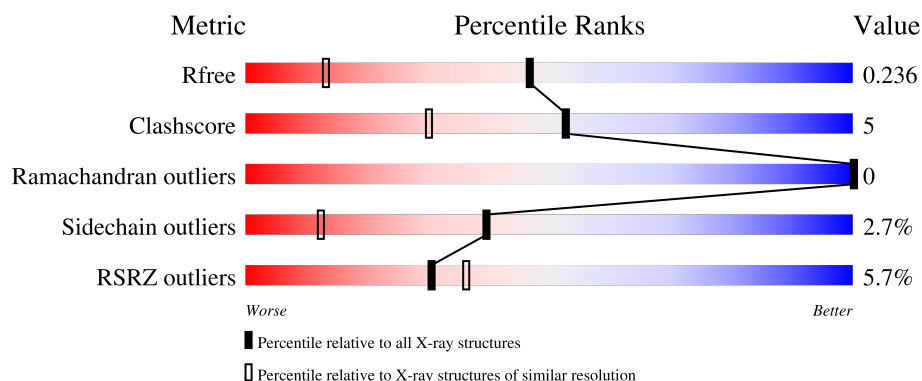
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 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))
$R_{free}$	164625	1935 (1.56-1.56)
Clashscore	180529	2073 (1.56-1.56)
Ramachandran outliers	177936	2037 (1.56-1.56)
Sidechain outliers	177891	2034 (1.56-1.56)
RSRZ outliers	164620	1935 (1.56-1.56)

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	219	<div> <div>5%</div> <div>89%</div> <div>8%</div> <div>.</div> </div>
1	B	219	<div> <div>5%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>
2	C	12	<div> <div>17%</div> <div>58%</div> <div>8%</div> <div>8%</div> <div>25%</div> </div>
2	D	12	<div> <div>17%</div> <div>50%</div> <div>17%</div> <div>8%</div> <div>25%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3846 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 Chaperone protein DnaK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	4	0
			1637	1007	284	340	6			
1	B	215	Total	C	N	O	S	0	4	0
			1653	1017	290	339	7			

- Molecule 2 is a protein called Pyrrhocoricin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	1
			75	53	12	10			
2	D	9	Total	C	N	O	0	0	1
			75	53	12	10			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	LEU	GLY	engineered mutation	UNP P37362
C	5	TYR	SER	engineered mutation	UNP P37362
C	6	ALC	TYR	engineered mutation	UNP P37362
C	7	ALC	LEU	engineered mutation	UNP P37362
C	12	THR	-	expression tag	UNP P37362
D	4	LEU	GLY	engineered mutation	UNP P37362
D	5	TYR	SER	engineered mutation	UNP P37362
D	6	ALC	TYR	engineered mutation	UNP P37362
D	7	ALC	LEU	engineered mutation	UNP P37362
D	12	THR	-	expression tag	UNP P37362

- 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	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

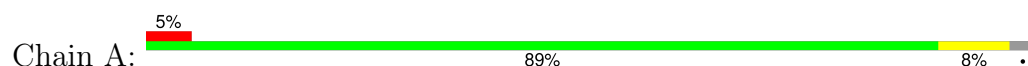
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	192	Total	O	0	0
			192	192		
4	B	177	Total	O	0	0
			177	177		
4	C	14	Total	O	0	0
			14	14		
4	D	3	Total	O	0	0
			3	3		

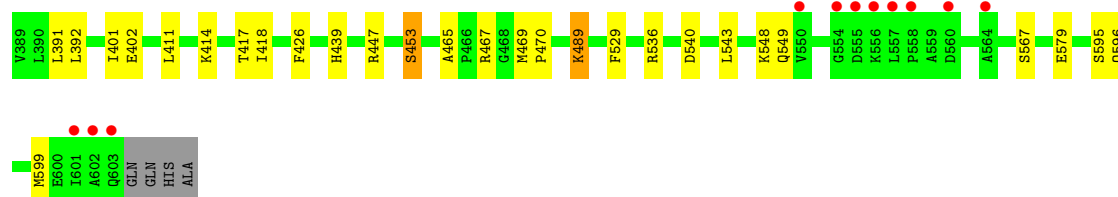
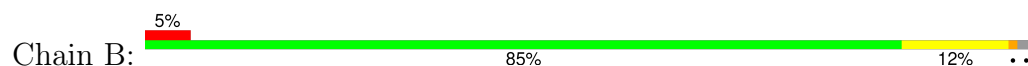
### 3 Residue-property plots [i](#)

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.

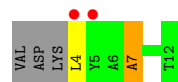
- Molecule 1: Chaperone protein DnaK



- Molecule 1: Chaperone protein DnaK



- Molecule 2: Pyrrhocoricin



- Molecule 2: Pyrrhocoricin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.40Å 162.13Å 45.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.60 – 1.55 24.60 – 1.55	Depositor EDS
% Data completeness (in resolution range)	97.9 (24.60-1.55) 97.9 (24.60-1.55)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.96 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.212 , 0.241 0.209 , 0.236	Depositor DCC
$R_{free}$ test set	1648 reflections (1.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.9	Xtriage
Anisotropy	0.442	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3846	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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: ALC, 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	1.07	1/1663 (0.1%)	1.18	7/2241 (0.3%)
1	B	1.00	1/1677 (0.1%)	1.14	10/2260 (0.4%)
2	C	1.22	0/54	1.33	0/72
2	D	1.02	0/54	0.92	0/72
All	All	1.04	2/3448 (0.1%)	1.16	17/4645 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	444	GLU	CD-OE1	-5.88	1.19	1.25
1	B	414	LYS	C-O	5.28	1.33	1.23

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	467	ARG	NE-CZ-NH1	7.35	123.97	120.30
1	A	491	LYS	CD-CE-NZ	-6.19	97.46	111.70
1	B	418	ILE	CG1-CB-CG2	6.12	124.87	111.40
1	A	532	LEU	CB-CG-CD2	-6.10	100.63	111.00
1	A	577	LYS	CD-CE-NZ	-5.88	98.18	111.70
1	B	529	PHE	CB-CG-CD1	5.84	124.89	120.80
1	A	481	ASP	CB-CG-OD2	-5.72	113.15	118.30
1	B	391	LEU	CB-CG-CD1	-5.72	101.27	111.00
1	B	447	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	A	447	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	431	ASP	CB-CG-OD1	5.39	123.15	118.30
1	B	411	LEU	CA-CB-CG	-5.37	102.95	115.30
1	B	447	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	B	536	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	B	543	LEU	CA-CB-CG	5.04	126.90	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	445	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	B	529	PHE	CB-CG-CD2	-5.00	117.30	120.80

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	1637	0	1654	9	0
1	B	1653	0	1673	25	0
2	C	75	0	81	1	0
2	D	75	0	80	4	0
3	A	5	0	0	0	0
3	B	10	0	0	0	0
3	C	5	0	0	0	0
4	A	192	0	0	3	0
4	B	177	0	0	2	0
4	C	14	0	0	0	0
4	D	3	0	0	0	0
All	All	3846	0	3488	34	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 5.

All (34) 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:439[A]:HIS:NE2	1:B:453[A]:SER:HB3	1.85	0.92
1:A:439:HIS:HE1	4:A:957:HOH:O	1.57	0.88
1:B:465:ALA:HB1	1:B:469[B]:MET:CE	2.06	0.86
1:B:465:ALA:CB	1:B:469[B]:MET:CE	2.54	0.84
1:B:465:ALA:CB	1:B:469[B]:MET:HE1	2.10	0.82
1:B:465:ALA:CB	1:B:469[B]:MET:HE2	2.17	0.75
1:B:401:ILE:HD12	2:D:7:ALC:HE23	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:465:ALA:HB1	1:B:469[B]:MET:HE1	1.68	0.72
1:B:548:LYS:HG3	1:B:549:GLN:N	2.05	0.70
1:A:439:HIS:CE1	4:A:957:HOH:O	2.36	0.69
1:B:465:ALA:HB2	1:B:469[B]:MET:CE	2.25	0.67
1:B:465:ALA:HB2	1:B:469[B]:MET:HE2	1.78	0.65
1:B:439[A]:HIS:NE2	1:B:453[A]:SER:CB	2.61	0.59
1:B:465:ALA:HB1	1:B:469[B]:MET:HE2	1.81	0.57
1:B:401:ILE:HD12	2:D:7:ALC:CZ	2.35	0.56
1:B:401:ILE:HD12	2:D:7:ALC:CE2	2.35	0.56
1:B:548:LYS:HG3	1:B:549:GLN:H	1.69	0.55
1:B:595:SER:O	1:B:599:MET:HG2	2.08	0.54
1:A:517:ARG:NH2	4:A:903:HOH:O	2.42	0.53
1:B:469[B]:MET:HB3	1:B:470:PRO:HD3	1.94	0.50
1:A:562:LYS:O	1:A:566:GLU:HG2	2.13	0.48
1:B:579:GLU:CG	4:B:939:HOH:O	2.61	0.47
1:A:438:ILE:HD12	1:A:438:ILE:N	2.30	0.46
1:A:472:ILE:HD13	2:C:7:ALC:HE23	1.97	0.45
1:A:566:GLU:HA	1:A:566:GLU:OE1	2.16	0.45
1:B:489:LYS:CB	1:B:489:LYS:NZ	2.79	0.45
1:B:392:LEU:HD13	1:B:417:THR:HG22	2.01	0.42
1:A:528:LYS:HD3	1:A:528:LYS:HA	1.79	0.41
1:B:402:GLU:OE1	2:D:8:PRO:HG3	2.19	0.41
1:B:401:ILE:HD11	1:B:426:PHE:CZ	2.55	0.41
1:B:401:ILE:HA	1:B:439[B]:HIS:O	2.21	0.41
1:A:601:ILE:H	1:A:601:ILE:HG13	1.53	0.40
1:B:579:GLU:HG2	4:B:939:HOH:O	2.19	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	215/219 (98%)	214 (100%)	1 (0%)	0	100	100
1	B	217/219 (99%)	215 (99%)	2 (1%)	0	100	100
2	C	5/12 (42%)	4 (80%)	1 (20%)	0	100	100
2	D	5/12 (42%)	5 (100%)	0	0	100	100
All	All	442/462 (96%)	438 (99%)	4 (1%)	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	181/181 (100%)	177 (98%)	4 (2%)	47	19
1	B	182/181 (101%)	176 (97%)	6 (3%)	33	8
2	C	6/10 (60%)	5 (83%)	1 (17%)	2	0
2	D	6/10 (60%)	6 (100%)	0	100	100
All	All	375/382 (98%)	364 (97%)	11 (3%)	40	10

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	392	LEU
1	A	551	GLU
1	A	556	LYS
1	A	601	ILE
1	B	453[A]	SER
1	B	453[B]	SER
1	B	489	LYS
1	B	540	ASP
1	B	567	SER
1	B	596	GLN
2	C	4	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	549	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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$
2	ALC	C	7	2	9,11,12	0.73	0	11,13,15	2.19	5 (45%)
2	ALC	D	7	2	9,11,12	1.47	1 (11%)	11,13,15	3.61	6 (54%)
2	ALC	D	6	2	9,11,12	1.06	0	11,13,15	1.09	1 (9%)
2	ALC	C	6	2	9,11,12	1.23	0	11,13,15	1.10	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	ALC	C	7	2	-	0/5/14/16	0/1/1/1
2	ALC	D	7	2	-	1/5/14/16	0/1/1/1
2	ALC	D	6	2	-	0/5/14/16	0/1/1/1
2	ALC	C	6	2	-	0/5/14/16	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	7	ALC	CB-CG	-3.59	1.47	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	7	ALC	CB-CG-CD2	6.70	127.31	111.71
2	D	7	ALC	CE2-CD2-CG	-5.54	100.43	112.08
2	D	7	ALC	CE1-CD1-CG	-5.18	101.17	112.08
2	D	7	ALC	CB-CA-C	-3.87	105.04	110.99
2	C	7	ALC	CE1-CD1-CG	-3.47	104.78	112.08
2	C	7	ALC	CG-CB-CA	-3.38	109.97	114.52
2	C	7	ALC	CB-CG-CD2	-3.24	104.17	111.71
2	D	7	ALC	CB-CG-CD1	2.99	118.67	111.71
2	D	7	ALC	CG-CB-CA	2.45	117.81	114.52
2	C	7	ALC	CZ-CE1-CD1	2.32	116.18	111.42
2	C	7	ALC	CZ-CE2-CD2	-2.22	106.86	111.42
2	D	6	ALC	CE2-CD2-CG	-2.05	107.78	112.08

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	7	ALC	CA-CB-CG-CD2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	7	ALC	1	0
2	D	7	ALC	3	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

4 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	B	701	-	4,4,4	0.55	0	6,6,6	0.63	0
3	SO4	A	701	-	4,4,4	0.53	0	6,6,6	0.18	0
3	SO4	C	101	-	4,4,4	0.58	0	6,6,6	0.76	0
3	SO4	B	702	-	4,4,4	0.56	0	6,6,6	0.88	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 [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	213/219 (97%)	0.22	10 (4%) 37 44	13, 20, 56, 78	4 (1%)
1	B	215/219 (98%)	0.32	11 (5%) 34 41	9, 23, 58, 71	4 (1%)
2	C	7/12 (58%)	1.31	2 (28%) 1 1	21, 26, 29, 41	0
2	D	7/12 (58%)	1.32	2 (28%) 1 1	25, 27, 38, 44	0
All	All	442/462 (95%)	0.30	25 (5%) 30 36	9, 21, 57, 78	8 (1%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	601	ILE	4.2
2	C	4	LEU	3.7
1	B	602	ALA	3.1
1	A	557	LEU	3.0
1	B	601	ILE	2.9
1	A	550	VAL	2.8
1	B	603	GLN	2.7
1	B	557	LEU	2.6
1	A	599	MET	2.6
1	B	554	GLY	2.5
2	D	12	THR	2.4
1	B	556	LYS	2.4
2	C	5	TYR	2.4
1	A	556	LYS	2.4
1	A	554	GLY	2.3
1	B	560	ASP	2.3
1	B	550	VAL	2.3
1	A	552	GLU	2.3
2	D	9	ARG	2.2
1	A	544	HIS	2.2
1	B	564	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	558	PRO	2.1
1	B	555	ASP	2.0
1	A	594	VAL	2.0
1	A	553	ALA	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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
2	ALC	C	7	11/12	0.94	0.08	16,18,23,25	0
2	ALC	D	6	11/12	0.95	0.08	18,20,22,22	0
2	ALC	D	7	11/12	0.95	0.09	18,19,25,28	0
2	ALC	C	6	11/12	0.97	0.06	16,19,25,26	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 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
3	SO4	B	702	5/5	0.77	0.13	49,57,60,66	0
3	SO4	A	701	5/5	0.80	0.10	74,77,81,82	0
3	SO4	B	701	5/5	0.92	0.08	40,44,49,53	0
3	SO4	C	101	5/5	0.98	0.07	23,25,29,29	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.