



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:14 AM GMT

PDB ID : 3DWI
Title : Crystal structure of Mycobacterium tuberculosis CysM, the cysteine synthase B
Authors : Jurgenson, C.T.; Burns, K.E.; Begley, T.P.; Ealick, S.E.
Deposited on : 2008-07-22
Resolution : 2.81 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

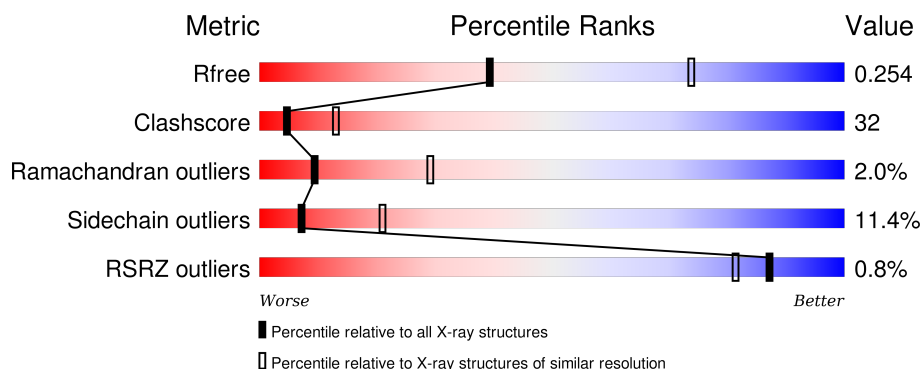
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

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 ≥ 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	323	<div> <div></div> <div>51% 35% 5% 8%</div> </div>
1	B	323	<div> <div></div> <div>46% 42% 8% .</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
2	SO4	A	324	-	-	-	X
3	PLP	A	401	-	X	-	-
3	PLP	B	401	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4646 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 Cysteine synthase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2161	1358	389	407	7			
1	B	309	Total	C	N	O	S	0	0	0
			2280	1435	407	430	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	204	ALA	LYS	ENGINEERED	UNP P63873
B	204	ALA	LYS	ENGINEERED	UNP P63873

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

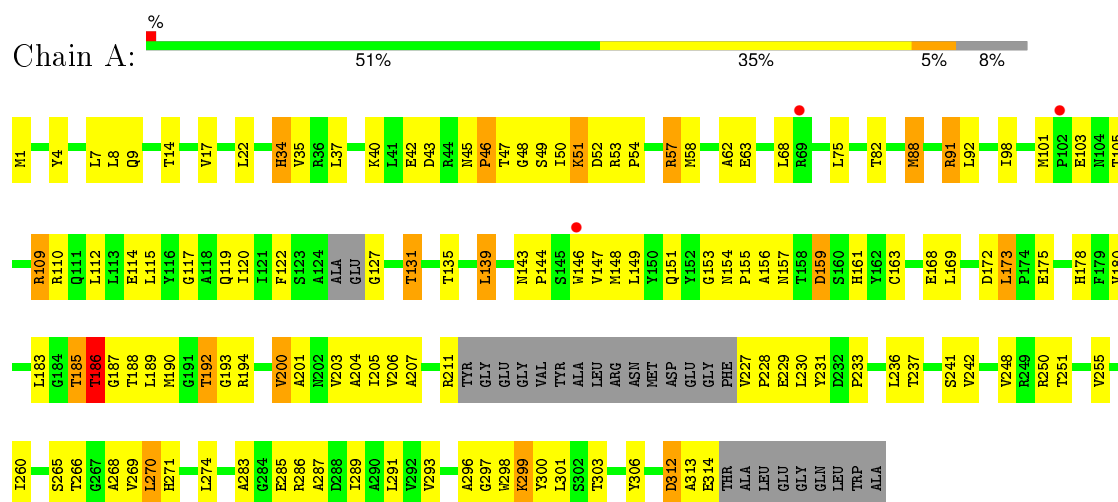
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	72	Total	O	0	0
			72	72		
4	B	88	Total	O	0	0
			88	88		

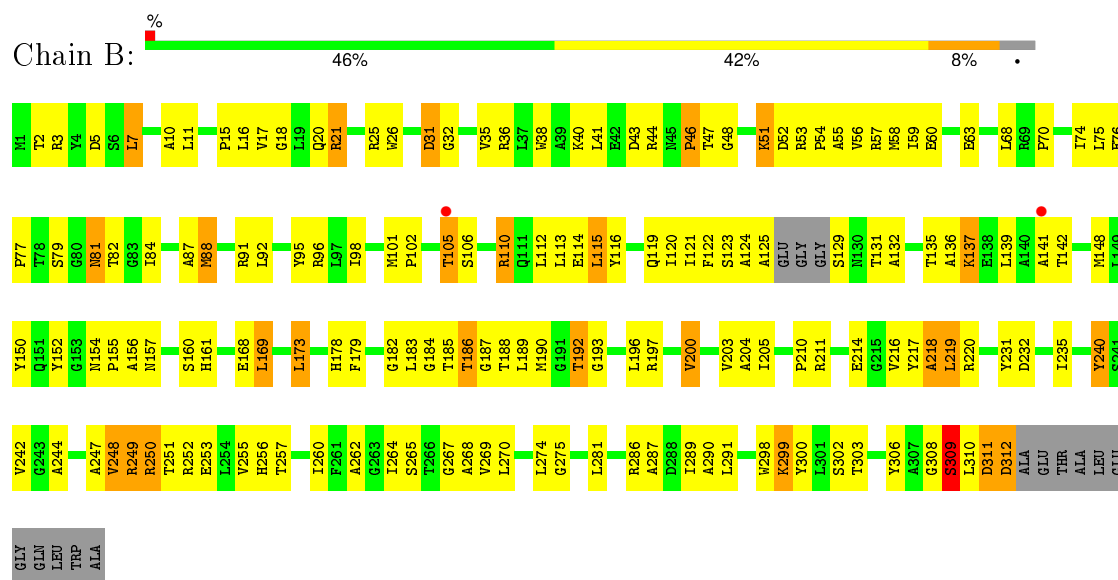
3 Residue-property plots [i](#)

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: Cysteine synthase B



• Molecule 1: Cysteine synthase B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.36Å 85.35Å 98.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.22 – 2.81 48.20 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.22-2.81) 99.4 (48.20-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.49 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.256 0.192 , 0.254	Depositor DCC
R_{free} test set	786 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 15466 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4646	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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.*

¹Intensities estimated from amplitudes.

²Theoretical 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: SO4, PLP

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	0.96	1/2201 (0.0%)	0.97	1/3001 (0.0%)
1	B	1.02	0/2324	1.02	9/3168 (0.3%)
All	All	0.99	1/4525 (0.0%)	1.00	10/6169 (0.2%)

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	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	163	CYS	CB-SG	-6.33	1.71	1.82

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	308	GLY	N-CA-C	7.30	131.35	113.10
1	B	183	LEU	CB-CG-CD2	-6.74	99.55	111.00
1	A	57	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	B	309	SER	N-CA-C	5.67	126.32	111.00
1	B	183	LEU	CA-CB-CG	5.61	128.20	115.30
1	B	5	ASP	CB-CG-OD1	5.45	123.21	118.30
1	B	44	ARG	NE-CZ-NH1	-5.41	117.59	120.30
1	B	169	LEU	CB-CG-CD1	-5.33	101.93	111.00
1	B	81	ASN	CB-CG-OD1	5.09	131.79	121.60
1	B	115	LEU	CA-CB-CG	5.09	127.00	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	MET	Peptide
1	A	312	ASP	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	2161	0	2105	136	0
1	B	2280	0	2240	156	0
2	A	5	0	0	1	0
2	B	10	0	0	0	0
3	A	15	0	6	3	0
3	B	15	0	6	8	0
4	A	72	0	0	9	0
4	B	88	0	0	6	1
All	All	4646	0	4357	282	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 32.

All (282) 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:91:ARG:HB3	1:A:91:ARG:NH1	1.60	1.15
1:A:188:THR:O	1:A:192:THR:HG23	1.48	1.12
1:B:88:MET:CE	1:B:92:LEU:HD12	1.82	1.10
1:A:250:ARG:HE	1:A:271:HIS:CD2	1.75	1.05
1:A:299:LYS:HD2	1:A:299:LYS:H	1.24	1.02
1:A:298:TRP:HB3	1:B:298:TRP:HB3	1.42	1.02
1:B:161:HIS:HB3	1:B:192:THR:HG22	1.43	1.00
1:B:185:THR:N	3:B:401:PLP:O1P	1.95	1.00
1:A:185:THR:O	1:A:186:THR:HB	1.63	0.99
1:B:88:MET:HE2	1:B:92:LEU:HD12	1.43	0.98
1:A:250:ARG:NE	1:A:271:HIS:HD2	1.61	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:VAL:HG13	1:B:287:ALA:HB3	1.42	0.98
1:B:185:THR:O	1:B:186:THR:HB	1.65	0.96
1:A:109:ARG:NH2	4:A:465:HOH:O	2.02	0.93
1:B:81:ASN:ND2	3:B:401:PLP:O3	2.01	0.93
1:A:161:HIS:HB3	1:A:192:THR:HG22	1.48	0.93
1:A:250:ARG:HE	1:A:271:HIS:HD2	0.97	0.92
1:A:131:THR:O	1:A:135:THR:N	2.01	0.92
1:B:186:THR:HG23	1:B:190:MET:HG2	1.54	0.89
1:B:88:MET:HE3	1:B:92:LEU:HD12	1.54	0.86
1:A:200:VAL:HG22	1:A:203:VAL:HB	1.57	0.85
1:B:200:VAL:HG22	1:B:203:VAL:HB	1.56	0.85
3:A:401:PLP:O1P	4:A:425:HOH:O	1.94	0.85
1:A:201:ALA:HB3	4:A:433:HOH:O	1.76	0.84
1:B:51:LYS:HG3	1:B:81:ASN:HB3	1.58	0.83
1:A:91:ARG:CB	1:A:91:ARG:NH1	2.42	0.83
1:A:42:GLU:HG3	1:A:50:ILE:HD13	1.60	0.82
1:A:227:VAL:N	4:A:461:HOH:O	2.11	0.82
1:A:115:LEU:HD11	1:B:306:TYR:HB3	1.60	0.82
1:B:188:THR:O	1:B:192:THR:HG23	1.82	0.80
1:A:185:THR:HG22	1:A:187:GLY:H	1.47	0.79
1:B:88:MET:HE3	1:B:92:LEU:CD1	2.11	0.79
1:A:296:ALA:H	1:A:299:LYS:HZ2	1.29	0.79
1:B:91:ARG:HH11	1:B:91:ARG:HB3	1.45	0.78
1:B:88:MET:CE	1:B:92:LEU:CD1	2.60	0.78
1:A:299:LYS:CD	1:A:299:LYS:H	1.96	0.78
1:B:98:ILE:HG21	1:B:121:ILE:HD12	1.63	0.78
1:A:186:THR:HG23	1:A:190:MET:HG2	1.66	0.77
1:B:252:ARG:HH12	1:B:309:SER:H	1.31	0.77
1:B:186:THR:HG23	1:B:190:MET:CG	2.14	0.77
1:B:75:LEU:HD12	1:B:98:ILE:HB	1.65	0.77
1:A:201:ALA:CB	4:A:433:HOH:O	2.30	0.75
1:B:217:TYR:O	1:B:218:ALA:HB3	1.86	0.75
1:A:91:ARG:CB	1:A:91:ARG:HH11	1.99	0.74
1:A:299:LYS:N	1:A:299:LYS:HD2	2.02	0.74
1:B:244:ALA:O	1:B:248:VAL:HG12	1.88	0.74
1:A:91:ARG:HB3	1:A:91:ARG:HH11	1.47	0.73
1:A:91:ARG:HB3	1:A:91:ARG:CZ	2.18	0.73
1:B:242:VAL:HG11	1:B:268:ALA:HA	1.68	0.73
1:A:101:MET:CE	1:A:110:ARG:HB3	2.19	0.72
1:B:26:TRP:CE3	1:B:36:ARG:HG2	2.23	0.72
1:B:211:ARG:NH2	1:B:214:GLU:OE2	2.21	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:THR:CG2	1:A:187:GLY:H	2.02	0.72
1:B:218:ALA:HB1	3:B:401:PLP:C3	2.20	0.72
1:A:42:GLU:HG3	1:A:50:ILE:CD1	2.20	0.72
1:A:298:TRP:HB3	1:B:298:TRP:CB	2.19	0.71
1:A:63:GLU:OE1	4:A:424:HOH:O	2.09	0.71
1:B:182:GLY:HA3	1:B:219:LEU:HD11	1.72	0.71
1:B:51:LYS:HG3	1:B:81:ASN:CB	2.21	0.70
1:A:296:ALA:O	1:A:299:LYS:HD3	1.91	0.70
1:A:51:LYS:NZ	3:A:401:PLP:O3	1.95	0.70
1:B:68:LEU:HD11	1:B:74:ILE:HD13	1.74	0.69
1:A:189:LEU:O	1:A:193:GLY:HA3	1.93	0.69
1:A:115:LEU:HD11	1:B:306:TYR:CB	2.23	0.68
1:B:218:ALA:HB1	3:B:401:PLP:C2	2.23	0.68
1:B:70:PRO:HA	4:B:467:HOH:O	1.95	0.67
1:A:4:TYR:HE1	1:B:15:PRO:HB3	1.60	0.66
1:B:11:LEU:CD1	1:B:52:ASP:HB3	2.25	0.66
1:B:252:ARG:NH1	1:B:309:SER:H	1.94	0.65
1:B:101:MET:O	1:B:122:PHE:HA	1.95	0.65
1:B:184:GLY:O	1:B:220:ARG:HD3	1.97	0.65
1:B:91:ARG:NH1	1:B:91:ARG:HB3	2.11	0.65
1:A:43:ASP:O	1:A:48:GLY:HA2	1.97	0.65
1:B:87:ALA:O	1:B:91:ARG:HG3	1.98	0.64
1:B:125:ALA:HB3	4:B:498:HOH:O	1.96	0.64
1:B:148:MET:HE3	1:B:150:TYR:HB2	1.81	0.62
1:B:51:LYS:HD3	1:B:82:THR:OG1	2.00	0.62
1:A:185:THR:O	1:A:186:THR:CB	2.41	0.62
1:A:47:THR:CG2	1:A:88:MET:HG2	2.30	0.62
1:B:20:GLN:HG2	1:B:26:TRP:CE2	2.35	0.61
1:B:311:ASP:O	1:B:311:ASP:OD1	2.18	0.61
1:A:186:THR:HG23	1:A:190:MET:CG	2.30	0.61
1:B:75:LEU:HD23	1:B:136:ALA:O	2.01	0.61
1:A:298:TRP:CB	1:B:298:TRP:HB3	2.24	0.61
1:A:101:MET:HE1	1:A:110:ARG:HB3	1.83	0.61
1:B:135:THR:O	1:B:139:LEU:HB2	2.00	0.61
1:B:189:LEU:O	1:B:193:GLY:HA3	2.01	0.60
1:B:16:LEU:HD21	1:B:173:LEU:HG	1.84	0.60
1:A:101:MET:CE	1:A:105:THR:HG21	2.32	0.59
1:B:91:ARG:HD2	1:B:116:TYR:O	2.03	0.59
1:A:299:LYS:HG2	1:A:300:TYR:CE2	2.37	0.59
1:B:217:TYR:O	1:B:218:ALA:CB	2.48	0.59
1:B:75:LEU:O	1:B:148:MET:HA	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:LYS:HA	1:A:54:PRO:HD2	1.85	0.58
1:A:101:MET:HE2	1:A:110:ARG:HB3	1.84	0.58
1:B:102:PRO:O	1:B:105:THR:HG23	2.04	0.58
1:A:35:VAL:HG13	1:A:287:ALA:HB3	1.86	0.58
1:A:306:TYR:HB3	1:B:115:LEU:HD22	1.86	0.58
1:B:53:ARG:NH2	1:B:168:GLU:OE1	2.33	0.58
1:A:101:MET:O	1:A:122:PHE:HA	2.04	0.57
1:B:193:GLY:HA2	1:B:205:ILE:HG12	1.86	0.57
1:A:53:ARG:HB2	1:A:54:PRO:HD3	1.85	0.57
1:A:57:ARG:HG2	1:A:156:ALA:O	2.04	0.57
1:B:81:ASN:HD21	3:B:401:PLP:H2A1	1.68	0.57
1:A:312:ASP:C	1:A:314:GLU:H	2.08	0.56
1:A:172:ASP:O	1:B:3:ARG:HD2	2.05	0.56
1:A:98:ILE:HG12	1:A:119:GLN:HB2	1.88	0.56
1:A:153:GLY:O	1:A:155:PRO:HD3	2.06	0.56
1:B:186:THR:HG21	1:B:231:TYR:HA	1.87	0.56
1:A:47:THR:HG23	1:A:88:MET:HG2	1.88	0.56
1:B:43:ASP:O	1:B:48:GLY:HA2	2.04	0.55
1:B:54:PRO:HG3	4:B:441:HOH:O	2.06	0.55
1:A:231:TYR:CZ	1:A:233:PRO:HA	2.41	0.55
1:A:157:ASN:HD21	1:A:185:THR:HG21	1.72	0.55
1:A:312:ASP:O	1:A:314:GLU:N	2.35	0.55
1:A:101:MET:HE3	1:A:105:THR:CB	2.36	0.55
1:A:114:GLU:HG3	1:A:120:ILE:CD1	2.37	0.55
1:B:11:LEU:HD11	1:B:52:ASP:HB3	1.87	0.55
1:A:62:ALA:HB3	1:A:68:LEU:HD22	1.88	0.55
1:A:185:THR:HG22	1:A:186:THR:N	2.22	0.55
1:A:75:LEU:O	1:A:148:MET:HA	2.07	0.55
1:B:110:ARG:HD2	4:B:485:HOH:O	2.06	0.55
1:B:185:THR:O	1:B:186:THR:CB	2.46	0.55
1:B:197:ARG:HH11	1:B:205:ILE:HD13	1.72	0.54
1:B:47:THR:OG1	1:B:52:ASP:OD1	2.21	0.54
1:A:312:ASP:C	1:A:314:GLU:N	2.60	0.54
1:A:180:VAL:HG12	1:A:269:VAL:HG13	1.90	0.54
1:B:152:TYR:CE1	1:B:185:THR:HG22	2.42	0.54
1:B:211:ARG:HB2	1:B:214:GLU:HG3	1.88	0.54
1:B:7:LEU:HD13	1:B:46:PRO:HB3	1.90	0.54
1:A:178:HIS:HB2	1:A:289:ILE:HG23	1.89	0.54
1:A:110:ARG:NH2	1:A:114:GLU:OE2	2.41	0.53
1:A:291:LEU:HD12	1:A:291:LEU:C	2.29	0.53
1:A:101:MET:HE1	1:A:105:THR:HG21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:ALA:CB	3:B:401:PLP:C2	2.86	0.53
1:B:210:PRO:HD3	1:B:219:LEU:HB3	1.91	0.52
1:A:58:MET:HB3	1:A:149:LEU:HD13	1.90	0.52
1:A:178:HIS:CB	1:A:289:ILE:HG23	2.40	0.52
1:B:79:SER:HA	1:B:113:LEU:HD11	1.92	0.52
1:A:193:GLY:HA2	1:A:205:ILE:HD12	1.92	0.52
1:A:242:VAL:HG11	1:A:268:ALA:HA	1.91	0.51
1:B:265:SER:OG	3:B:401:PLP:N1	2.41	0.51
1:B:31:ASP:CG	1:B:32:GLY:H	2.13	0.51
1:B:186:THR:HG21	1:B:231:TYR:CA	2.40	0.51
1:B:56:VAL:O	1:B:60:GLU:HG3	2.10	0.51
1:B:20:GLN:HG2	1:B:26:TRP:CD2	2.45	0.51
1:B:57:ARG:HG2	1:B:156:ALA:O	2.11	0.50
1:B:63:GLU:HG2	1:B:95:TYR:OH	2.11	0.50
1:A:4:TYR:CE1	1:B:15:PRO:HB3	2.44	0.50
1:A:178:HIS:CD2	1:A:204:ALA:HB3	2.47	0.50
1:A:101:MET:HE3	1:A:105:THR:HB	1.92	0.50
1:A:211:ARG:NH1	4:A:462:HOH:O	2.44	0.50
1:A:231:TYR:OH	1:A:233:PRO:HA	2.12	0.50
1:B:106:SER:HA	4:B:496:HOH:O	2.11	0.50
1:B:51:LYS:C	1:B:54:PRO:HD2	2.31	0.50
1:B:51:LYS:CD	1:B:82:THR:OG1	2.61	0.49
1:B:31:ASP:N	1:B:31:ASP:OD1	2.44	0.49
1:A:188:THR:O	1:A:192:THR:CG2	2.41	0.49
1:A:4:TYR:HB3	1:A:9:GLN:HB2	1.95	0.49
1:B:40:LYS:HD3	1:B:169:LEU:HD21	1.94	0.49
1:B:68:LEU:HD11	1:B:74:ILE:CD1	2.43	0.49
1:A:180:VAL:HG22	1:A:206:VAL:HB	1.94	0.48
1:A:169:LEU:HD23	1:A:169:LEU:HA	1.69	0.48
1:B:54:PRO:O	1:B:58:MET:HG3	2.13	0.48
1:A:183:LEU:HD12	1:A:189:LEU:HD12	1.94	0.48
1:A:34:HIS:O	1:A:34:HIS:CD2	2.67	0.48
1:B:210:PRO:HG3	1:B:216:VAL:HB	1.96	0.48
1:A:185:THR:HB	3:A:401:PLP:O3P	2.13	0.47
1:A:156:ALA:HA	1:A:159:ASP:HB2	1.96	0.47
1:B:25:ARG:HD3	4:B:491:HOH:O	2.14	0.47
1:B:88:MET:HE2	1:B:92:LEU:CD1	2.29	0.47
1:A:127:GLY:HA3	1:A:131:THR:HG21	1.97	0.47
1:B:186:THR:H	3:B:401:PLP:P	2.38	0.47
1:B:54:PRO:HB3	1:B:157:ASN:HA	1.96	0.47
1:A:103:GLU:HA	1:A:122:PHE:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:ILE:HD12	1:B:205:ILE:N	2.30	0.47
1:A:255:VAL:HG21	1:B:115:LEU:HD12	1.96	0.47
1:B:303:THR:HG22	1:B:303:THR:O	2.14	0.47
1:B:51:LYS:HA	1:B:54:PRO:HD2	1.97	0.47
1:B:310:LEU:O	1:B:311:ASP:HB3	2.14	0.46
1:A:303:THR:HB	1:A:314:GLU:HG2	1.97	0.46
1:A:296:ALA:N	1:A:299:LYS:HZ2	2.07	0.46
1:B:210:PRO:CD	1:B:219:LEU:HB3	2.45	0.46
1:A:251:THR:CG2	1:A:306:TYR:HE1	2.29	0.46
1:B:169:LEU:HA	1:B:169:LEU:HD23	1.36	0.46
1:B:247:ALA:O	1:B:267:GLY:HA3	2.16	0.46
1:B:18:GLY:HA3	1:B:26:TRP:CH2	2.50	0.46
1:B:197:ARG:HG3	1:B:205:ILE:CD1	2.45	0.46
1:B:252:ARG:HH12	1:B:309:SER:N	2.06	0.45
1:B:101:MET:CE	1:B:120:ILE:HG23	2.47	0.45
1:B:41:LEU:HD21	1:B:260:ILE:HD13	1.99	0.45
1:A:51:LYS:C	1:A:54:PRO:HD2	2.37	0.45
1:B:311:ASP:C	1:B:312:ASP:CG	2.74	0.45
1:A:286:ARG:HD3	4:A:456:HOH:O	2.16	0.45
1:A:296:ALA:HB1	1:A:298:TRP:CD1	2.50	0.45
1:B:53:ARG:HH22	1:B:168:GLU:CD	2.19	0.45
1:A:227:VAL:HA	1:A:228:PRO:HD3	1.81	0.45
1:A:149:LEU:O	1:A:154:ASN:HB2	2.16	0.45
1:A:270:LEU:HD22	1:A:274:LEU:HG	1.99	0.45
1:A:189:LEU:O	1:A:193:GLY:CA	2.63	0.45
1:A:255:VAL:HA	1:A:260:ILE:O	2.17	0.45
1:A:227:VAL:HG23	4:A:461:HOH:O	2.17	0.45
1:B:249:ARG:HG3	1:B:253:GLU:OE1	2.16	0.45
1:A:271:HIS:HE1	2:A:324:SO4:S	2.40	0.45
1:B:57:ARG:CB	1:B:160:SER:HB3	2.47	0.45
1:B:216:VAL:HG13	1:B:264:ILE:HG22	1.99	0.44
1:A:250:ARG:NE	1:A:271:HIS:CD2	2.51	0.44
1:A:34:HIS:C	1:A:34:HIS:CD2	2.91	0.44
1:A:22:LEU:HA	1:A:22:LEU:HD12	1.64	0.44
1:A:291:LEU:O	1:A:291:LEU:HD12	2.17	0.44
1:A:251:THR:HG22	1:A:306:TYR:HE1	1.82	0.44
1:B:101:MET:HE2	1:B:120:ILE:HG23	1.99	0.44
1:B:10:ALA:O	1:B:46:PRO:HD3	2.17	0.44
1:A:151:GLN:HE21	1:A:151:GLN:HB2	1.63	0.44
1:B:178:HIS:HB2	1:B:289:ILE:HG12	1.98	0.44
1:B:21:ARG:HB2	1:B:257:THR:HG22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:HIS:CD2	1:B:204:ALA:HB3	2.53	0.43
1:B:96:ARG:NH1	1:B:119:GLN:HE22	2.15	0.43
1:B:251:THR:CG2	1:B:262:ALA:O	2.66	0.43
1:B:255:VAL:HG13	1:B:256:HIS:N	2.33	0.43
1:B:112:LEU:O	1:B:115:LEU:HG	2.19	0.43
1:A:297:GLY:O	1:A:300:TYR:N	2.44	0.43
1:B:68:LEU:HD21	1:B:74:ILE:HD11	2.01	0.43
1:B:250:ARG:NH1	1:B:253:GLU:OE2	2.47	0.43
1:A:139:LEU:HB3	1:A:146:TRP:CZ3	2.54	0.43
1:B:274:LEU:HD23	1:B:274:LEU:HA	1.85	0.43
1:B:51:LYS:O	1:B:54:PRO:HD2	2.19	0.43
1:A:91:ARG:CD	1:A:117:GLY:HA3	2.48	0.43
1:B:197:ARG:HG3	1:B:205:ILE:HD11	2.00	0.43
1:A:101:MET:HE3	1:A:105:THR:HG21	2.00	0.42
1:B:154:ASN:HA	1:B:155:PRO:HD2	1.89	0.42
1:B:101:MET:HE3	1:B:120:ILE:HG12	1.99	0.42
1:B:25:ARG:NH1	1:B:32:GLY:HA3	2.33	0.42
1:B:55:ALA:O	1:B:59:ILE:HG13	2.19	0.42
1:B:76:GLU:CD	1:B:77:PRO:HD2	2.39	0.42
1:A:40:LYS:NZ	1:A:168:GLU:OE1	2.47	0.42
1:A:37:LEU:HD11	1:A:274:LEU:CD2	2.49	0.42
1:B:281:LEU:HD12	1:B:281:LEU:HA	1.80	0.42
1:A:92:LEU:HA	1:A:92:LEU:HD23	1.88	0.42
1:A:14:THR:HG21	1:A:42:GLU:HA	2.02	0.42
1:B:11:LEU:HD11	1:B:52:ASP:CB	2.48	0.42
1:B:84:ILE:HG23	1:B:116:TYR:CE2	2.55	0.42
1:B:232:ASP:HB3	1:B:235:ILE:HG12	2.01	0.42
1:B:186:THR:O	1:B:190:MET:HG2	2.20	0.42
1:B:137:LYS:O	1:B:141:ALA:N	2.49	0.42
1:A:45:ASN:HB3	1:A:52:ASP:OD2	2.20	0.42
1:A:157:ASN:ND2	1:A:185:THR:HG21	2.34	0.41
1:B:102:PRO:HD3	1:B:132:ALA:HB2	2.01	0.41
1:A:189:LEU:HB3	1:A:236:LEU:HD21	2.03	0.41
1:A:8:LEU:HD23	1:A:8:LEU:HA	1.79	0.41
1:A:266:THR:HG21	1:A:293:VAL:HG22	2.02	0.41
1:A:51:LYS:CA	1:A:54:PRO:HD2	2.49	0.41
1:A:68:LEU:HD13	1:A:147:VAL:HG11	2.02	0.41
1:B:179:PHE:HB2	1:B:196:LEU:CD1	2.51	0.41
1:B:161:HIS:CG	1:B:188:THR:HA	2.56	0.41
1:A:143:ASN:HA	1:A:144:PRO:HD2	1.80	0.41
1:B:240:TYR:OH	1:B:275:GLY:HA3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:LEU:HD23	1:A:175:GLU:CD	2.41	0.41
1:A:103:GLU:HA	1:A:122:PHE:HD1	1.85	0.41
1:B:219:LEU:HD12	1:B:219:LEU:HA	1.79	0.41
1:B:310:LEU:O	1:B:311:ASP:CB	2.68	0.41
1:A:68:LEU:HD12	1:A:147:VAL:HG21	2.02	0.41
1:B:269:VAL:HG12	1:B:291:LEU:HB2	2.03	0.41
1:A:112:LEU:HD11	1:B:298:TRP:HZ3	1.86	0.41
1:A:43:ASP:C	1:A:45:ASN:H	2.23	0.41
1:B:312:ASP:N	1:B:312:ASP:OD1	2.54	0.41
1:B:110:ARG:HG3	1:B:114:GLU:OE1	2.20	0.41
1:B:299:LYS:HG3	1:B:300:TYR:CE2	2.56	0.41
1:A:82:THR:OG1	1:A:151:GLN:NE2	2.54	0.41
1:B:92:LEU:HD23	1:B:92:LEU:HA	1.87	0.41
1:B:300:TYR:O	1:B:303:THR:HB	2.21	0.41
1:A:112:LEU:O	1:A:115:LEU:HB2	2.21	0.40
1:A:283:ALA:HB3	1:A:285:GLU:HG3	2.03	0.40
1:A:186:THR:HG22	1:A:230:LEU:HB2	2.03	0.40
1:A:207:ALA:HB2	1:A:236:LEU:HD22	2.02	0.40
1:B:38:TRP:CE2	1:B:173:LEU:HD21	2.57	0.40
1:B:38:TRP:NE1	1:B:173:LEU:HD21	2.37	0.40
1:B:40:LYS:HD3	1:B:169:LEU:CD2	2.51	0.40
1:A:45:ASN:O	1:A:48:GLY:N	2.54	0.40
1:B:38:TRP:O	1:B:290:ALA:HA	2.22	0.40
1:B:286:ARG:HD2	1:B:286:ARG:C	2.42	0.40
1:B:157:ASN:OD1	1:B:187:GLY:HA3	2.22	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)
4:B:470:HOH:O	4:B:493:HOH:O[4_566]	2.15	0.05

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	291/323 (90%)	262 (90%)	26 (9%)	3 (1%)	19	51
1	B	305/323 (94%)	278 (91%)	18 (6%)	9 (3%)	5	18
All	All	596/646 (92%)	540 (91%)	44 (7%)	12 (2%)	9	29

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	313	ALA
1	B	309	SER
1	B	311	ASP
1	B	31	ASP
1	B	302	SER
1	B	123	SER
1	B	142	THR
1	B	218	ALA
1	A	186	THR
1	A	46	PRO
1	B	124	ALA
1	B	46	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	206/242 (85%)	180 (87%)	26 (13%)	5	16
1	B	223/242 (92%)	200 (90%)	23 (10%)	9	25
All	All	429/484 (89%)	380 (89%)	49 (11%)	7	21

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU

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Mol	Chain	Res	Type
1	A	17	VAL
1	A	34	HIS
1	A	46	PRO
1	A	49	SER
1	A	51	LYS
1	A	88	MET
1	A	91	ARG
1	A	109	ARG
1	A	131	THR
1	A	139	LEU
1	A	159	ASP
1	A	173	LEU
1	A	185	THR
1	A	186	THR
1	A	192	THR
1	A	194	ARG
1	A	200	VAL
1	A	229	GLU
1	A	237	THR
1	A	241	SER
1	A	248	VAL
1	A	265	SER
1	A	270	LEU
1	A	299	LYS
1	A	301	LEU
1	B	2	THR
1	B	7	LEU
1	B	17	VAL
1	B	21	ARG
1	B	51	LYS
1	B	88	MET
1	B	105	THR
1	B	110	ARG
1	B	129	SER
1	B	131	THR
1	B	137	LYS
1	B	173	LEU
1	B	186	THR
1	B	192	THR
1	B	200	VAL
1	B	219	LEU
1	B	240	TYR

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Mol	Chain	Res	Type
1	B	248	VAL
1	B	249	ARG
1	B	250	ARG
1	B	270	LEU
1	B	299	LYS
1	B	312	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	A	151	GLN
1	A	271	HIS
1	B	61	GLN
1	B	256	HIS

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](#)

5 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
2	SO4	A	324	-	4,4,4	0.23	0	6,6,6	0.35	0
3	PLP	A	401	1	15,15,16	4.58	10 (66%)	21,22,23	3.27	10 (47%)
2	SO4	B	324	-	4,4,4	0.34	0	6,6,6	0.59	0
2	SO4	B	325	-	4,4,4	0.14	0	6,6,6	0.49	0
3	PLP	B	401	1	15,15,16	1.91	2 (13%)	21,22,23	0.97	1 (4%)

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	324	-	-	0/0/0/0	0/0/0/0
3	PLP	A	401	1	-	0/6/6/8	0/1/1/1
2	SO4	B	324	-	-	0/0/0/0	0/0/0/0
2	SO4	B	325	-	-	0/0/0/0	0/0/0/0
3	PLP	B	401	1	-	0/6/6/8	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	PLP	C3-C2	-8.53	1.34	1.40
3	A	401	PLP	P-O2P	-7.65	1.27	1.54
3	A	401	PLP	P-O4P	-6.55	1.38	1.60
3	A	401	PLP	P-O3P	-6.51	1.31	1.54
3	A	401	PLP	C4A-C4	-6.38	1.38	1.51
3	B	401	PLP	C4A-C4	-6.02	1.39	1.51
3	A	401	PLP	P-O1P	-4.47	1.36	1.51
3	A	401	PLP	C3-C4	-2.94	1.32	1.40
3	A	401	PLP	C5-C4	-2.38	1.37	1.40
3	A	401	PLP	O3-C3	-2.37	1.31	1.37
3	A	401	PLP	O4P-C5A	-2.31	1.35	1.44
3	B	401	PLP	C3-C2	-2.15	1.39	1.40

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	PLP	O3P-P-O2P	-4.77	89.21	107.38
3	A	401	PLP	C4A-C4-C3	-4.22	112.72	120.36
3	A	401	PLP	O2P-P-O1P	-3.45	99.48	110.58
3	A	401	PLP	C5A-C5-C6	-2.74	114.10	119.28
3	B	401	PLP	O4P-P-O1P	2.22	112.80	107.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	PLP	O3P-P-O1P	3.06	120.42	110.58
3	A	401	PLP	O3-C3-C2	3.26	123.32	117.66
3	A	401	PLP	O3P-P-O4P	3.35	116.20	106.56
3	A	401	PLP	O4P-P-O1P	4.57	118.77	107.14
3	A	401	PLP	C5A-C5-C4	5.23	128.58	121.65
3	A	401	PLP	C4A-C4-C5	8.34	129.58	120.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	324	SO4	1	0
3	A	401	PLP	3	0
3	B	401	PLP	8	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, 95th 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 ²)	Q<0.9
1	A	297/323 (91%)	-0.08	3 (1%) 84 77	13, 26, 46, 54	0
1	B	309/323 (95%)	0.00	2 (0%) 90 86	12, 26, 46, 66	0
All	All	606/646 (93%)	-0.04	5 (0%) 87 81	12, 26, 46, 66	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	141	ALA	4.3
1	A	102	PRO	2.2
1	A	69	ARG	2.2
1	A	146	TRP	2.1
1	B	105	THR	2.1

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, 95th 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(\AA^2)	Q<0.9
2	SO4	A	324	5/5	0.82	0.28	2.47	81,81,82,82	0
3	PLP	B	401	15/16	0.90	0.23	1.64	7,20,25,26	0
3	PLP	A	401	15/16	0.92	0.20	1.27	12,17,21,24	0
2	SO4	B	324	5/5	0.87	0.22	0.07	66,68,68,68	0
2	SO4	B	325	5/5	0.94	0.32	-	52,53,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.