



# Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 06:38 AM GMT

PDB ID : 3VW2  
Title : Crystal Structure of The Berberine-bound Form of RamR (Transcriptional Regulator of TetR Family) from Salmonella Typhimurium  
Authors : Sakurai, K.; Nikaido, E.; Nakashima, R.; Yamasaki, S.; Yamaguchi, A.; Nishino, K.  
Deposited on : 2012-07-30  
Resolution : 2.34 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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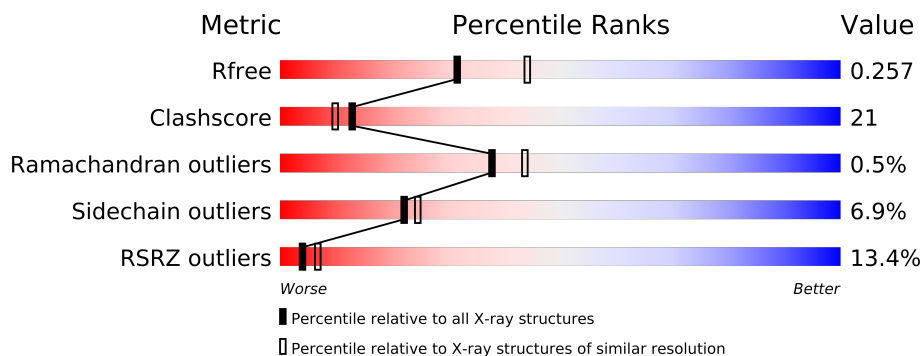
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.34 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))
$R_{free}$	66092	4049 (2.38-2.30)
Clashscore	79885	1094 (2.36-2.32)
Ramachandran outliers	78287	1080 (2.36-2.32)
Sidechain outliers	78261	1081 (2.36-2.32)
RSRZ outliers	66119	4050 (2.38-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	194	
1	B	194	
1	C	194	
1	D	194	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	BER	A	301	X	-
2	BER	B	301	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6032 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

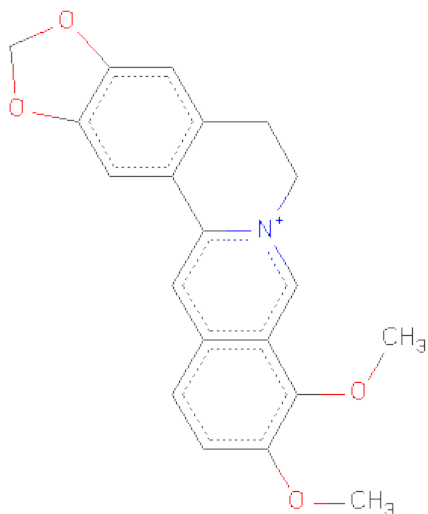
- Molecule 1 is a protein called Putative regulatory protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1462	922	258	271	11			
1	B	184	Total	C	N	O	S	0	0	0
			1462	922	258	271	11			
1	C	185	Total	C	N	O	S	0	0	0
			1468	926	259	272	11			
1	D	183	Total	C	N	O	S	0	0	0
			1453	917	257	268	11			

There are 8 discrepancies between the modelled and reference sequences:

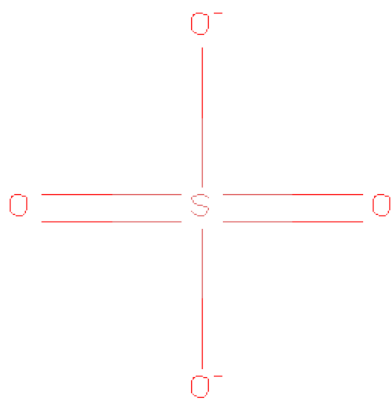
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	EXPRESSION TAG	UNP D0ZP76
A	1	VAL	-	EXPRESSION TAG	UNP D0ZP76
B	0	MET	-	EXPRESSION TAG	UNP D0ZP76
B	1	VAL	-	EXPRESSION TAG	UNP D0ZP76
C	0	MET	-	EXPRESSION TAG	UNP D0ZP76
C	1	VAL	-	EXPRESSION TAG	UNP D0ZP76
D	0	MET	-	EXPRESSION TAG	UNP D0ZP76
D	1	VAL	-	EXPRESSION TAG	UNP D0ZP76

- Molecule 2 is BERBERINE (three-letter code: BER) (formula: C<sub>20</sub>H<sub>18</sub>NO<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	20	1	4		
2	B	1	Total	C	N	O	0	0
			25	20	1	4		
2	C	1	Total	C	N	O	0	0
			25	20	1	4		
2	D	1	Total	C	N	O	0	0
			25	20	1	4		

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

- Molecule 4 is water.

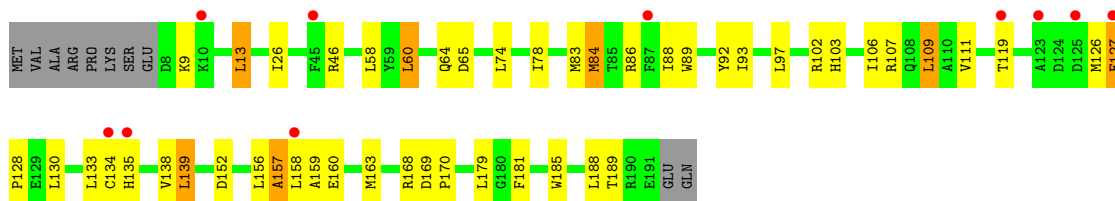
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	16	Total	O	0	0
			16	16		
4	C	16	Total	O	0	0
			16	16		
4	D	13	Total	O	0	0
			13	13		

### 3 Residue-property plots

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.

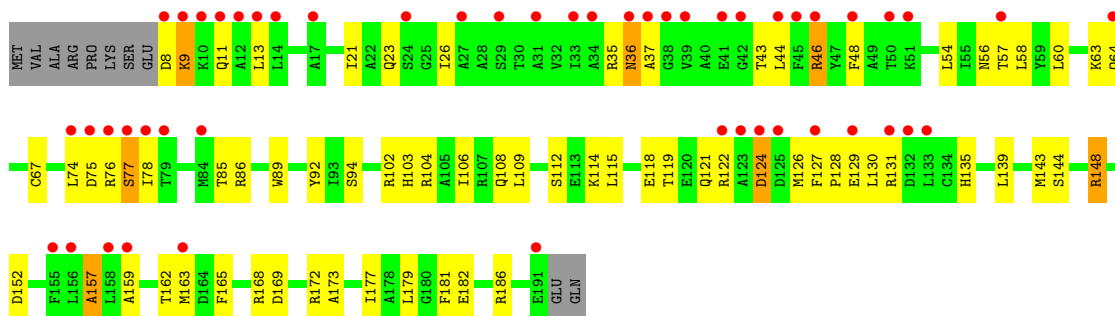
- Molecule 1: Putative regulatory protein

Chain A: 



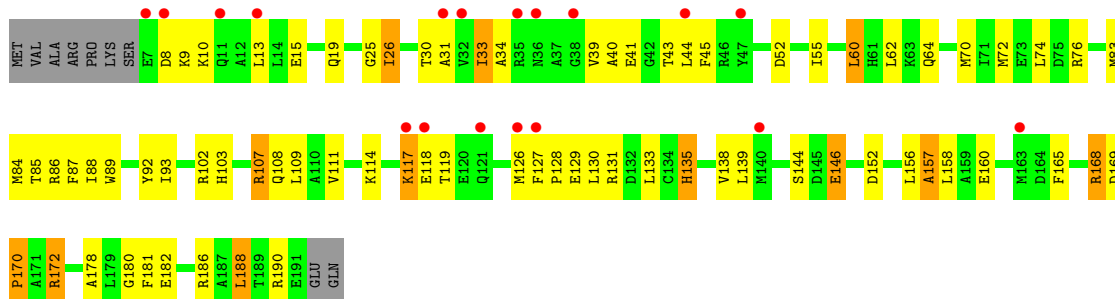
- Molecule 1: Putative regulatory protein

Chain B: 



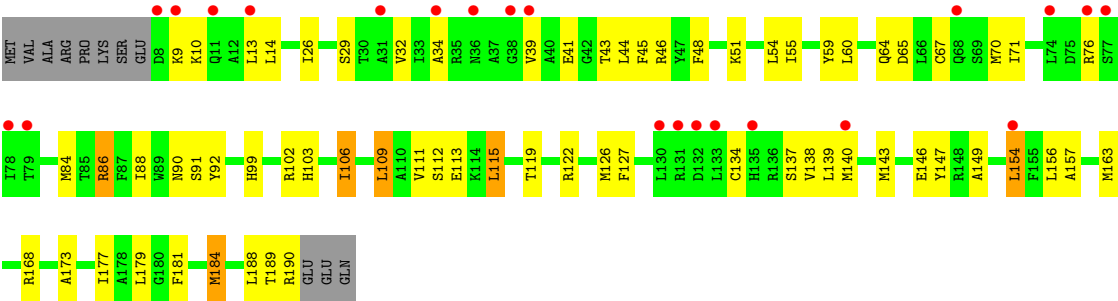
- Molecule 1: Putative regulatory protein

Chain C: 



- Molecule 1: Putative regulatory protein

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.82Å 54.62Å 92.29Å 104.97° 97.74° 90.04°	Depositor
Resolution (Å)	41.29 – 2.34 41.29 – 2.34	Depositor EDS
% Data completeness (in resolution range)	94.6 (41.29-2.34) 94.6 (41.29-2.34)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.78 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.225 , 0.261 0.222 , 0.257	Depositor DCC
$R_{free}$ test set	1693 reflections (5.44%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.4	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 47.4	EDS
Estimated twinning fraction	0.026 for -h,k,-k-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32793 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6032	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BER, 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	1.14	1/1488 (0.1%)	0.83	0/2006
1	B	1.12	3/1488 (0.2%)	0.79	0/2006
1	C	1.18	4/1494 (0.3%)	0.87	1/2014 (0.0%)
1	D	1.03	1/1479 (0.1%)	0.84	1/1994 (0.1%)
All	All	1.12	9/5949 (0.2%)	0.83	2/8020 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	159	ALA	CA-CB	5.71	1.64	1.52
1	A	157	ALA	CA-CB	5.68	1.64	1.52
1	B	157	ALA	CA-CB	5.59	1.64	1.52
1	C	180	GLY	C-O	5.44	1.32	1.23
1	C	157	ALA	CA-CB	5.29	1.63	1.52
1	D	149	ALA	CA-CB	5.18	1.63	1.52
1	C	129	GLU	CD-OE1	5.17	1.31	1.25
1	C	146	GLU	CG-CD	5.10	1.59	1.51
1	B	94	SER	CB-OG	-5.09	1.35	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	172	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	D	154	LEU	CB-CG-CD1	5.26	119.95	111.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1462	0	1450	40	0
1	B	1462	0	1450	61	0
1	C	1468	0	1454	70	0
1	D	1453	0	1444	69	0
2	A	25	0	18	5	0
2	B	25	0	18	9	0
2	C	25	0	18	3	0
2	D	25	0	18	10	0
3	A	5	0	0	0	0
3	B	5	0	0	1	0
3	C	5	0	0	1	0
3	D	5	0	0	0	0
4	A	22	0	0	1	0
4	B	16	0	0	4	0
4	C	16	0	0	3	0
4	D	13	0	0	1	0
All	All	6032	0	5870	246	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 21.

All (246) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:301:BER:C19	2:D:301:BER:H61	1.61	1.27
1:B:75:ASP:HB3	1:B:78:ILE:HD11	1.17	1.16
2:D:301:BER:C6	2:D:301:BER:H193	1.72	1.15
2:B:301:BER:C19	2:B:301:BER:H61	1.75	1.14
2:B:301:BER:H193	2:B:301:BER:H61	1.28	1.08
2:A:301:BER:O4	2:A:301:BER:H192	1.53	1.07
2:B:301:BER:C6	2:B:301:BER:H193	1.86	1.02
1:D:60:LEU:HD21	1:D:119:THR:HG23	1.40	1.02
2:D:301:BER:H61	2:D:301:BER:H193	1.04	1.01
1:C:33:ILE:O	1:C:33:ILE:HG12	1.58	1.01
2:B:301:BER:O3	2:B:301:BER:H203	1.63	0.97

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:60:LEU:HD11	1:B:119:THR:HG23	1.44	0.95
1:B:118:GLU:HB3	4:B:405:HOH:O	1.64	0.95
1:D:60:LEU:CD2	1:D:119:THR:HG23	2.00	0.91
1:A:92:TYR:OH	1:A:103:HIS:CD2	2.24	0.91
2:D:301:BER:C19	2:D:301:BER:C6	2.30	0.90
2:B:301:BER:C6	2:B:301:BER:C19	2.43	0.90
1:D:86:ARG:HH12	1:D:90:ASN:HD21	1.17	0.89
1:D:86:ARG:NH1	1:D:90:ASN:HD21	1.72	0.88
1:C:165:PHE:CZ	1:D:111:VAL:HG21	2.09	0.86
1:A:92:TYR:OH	1:A:103:HIS:HD2	1.58	0.85
2:D:301:BER:H61	2:D:301:BER:H192	1.60	0.83
1:D:115:LEU:HD12	1:D:115:LEU:H	1.43	0.83
1:C:168:ARG:NH2	1:D:111:VAL:HG22	1.92	0.83
1:D:106:ILE:HD12	2:D:301:BER:H201	1.60	0.82
1:D:13:LEU:HD21	1:D:44:LEU:HA	1.61	0.82
1:D:86:ARG:HG3	1:D:86:ARG:HH11	1.45	0.82
2:B:301:BER:O3	2:B:301:BER:C20	2.30	0.80
2:A:301:BER:O4	2:A:301:BER:C19	2.30	0.80
1:D:106:ILE:CD1	2:D:301:BER:H201	2.12	0.79
1:C:92:TYR:OH	1:C:103:HIS:CD2	2.35	0.79
2:B:301:BER:H192	2:B:301:BER:H61	1.63	0.79
1:D:70:MET:HG2	1:D:91:SER:HB3	1.65	0.78
1:C:70:MET:HE1	1:C:130:LEU:HD22	1.65	0.78
1:B:43:THR:O	1:B:46:ARG:HG3	1.84	0.78
1:B:9:LYS:HD3	1:B:9:LYS:H	1.47	0.77
1:C:13:LEU:HD21	1:C:44:LEU:HA	1.66	0.77
1:D:29:SER:OG	1:D:32:VAL:HG23	1.85	0.77
1:B:60:LEU:CD1	1:B:119:THR:HG23	2.15	0.75
1:D:106:ILE:HD13	1:D:106:ILE:O	1.85	0.74
1:D:76:ARG:HB2	4:D:408:HOH:O	1.87	0.74
1:C:55:ILE:HG23	1:C:109:LEU:HD22	1.69	0.73
1:C:84:MET:O	1:C:88:ILE:HG12	1.88	0.73
1:A:134:CYS:O	1:A:138:VAL:HG23	1.87	0.73
1:A:60:LEU:HD13	1:A:119:THR:HG23	1.69	0.73
1:D:92:TYR:OH	1:D:103:HIS:HD2	1.73	0.72
1:C:128:PRO:HB3	1:C:131:ARG:NH2	2.06	0.70
1:D:99:HIS:HB3	1:D:102:ARG:HG3	1.73	0.70
1:C:152:ASP:OD1	2:C:301:BER:H61	1.93	0.68
1:D:71:ILE:HD11	1:D:127:PHE:CD2	2.29	0.68
1:A:88:ILE:HD13	1:A:130:LEU:HD13	1.76	0.67
1:D:106:ILE:HD12	2:D:301:BER:C20	2.25	0.66
1:C:92:TYR:OH	1:C:103:HIS:HD2	1.77	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:172:ARG:HE	1:D:146:GLU:CD	1.99	0.66
1:C:117:LYS:HB2	1:C:117:LYS:NZ	2.11	0.66
1:D:13:LEU:HD23	1:D:48:PHE:CE2	2.31	0.65
1:B:124:ASP:HB3	1:B:131:ARG:HD3	1.77	0.65
1:C:60:LEU:HD13	1:C:119:THR:HG23	1.80	0.64
1:C:92:TYR:HH	1:C:103:HIS:CD2	2.15	0.64
1:A:64:GLN:HB2	1:A:126:MET:HE1	1.80	0.64
1:D:109:LEU:O	1:D:115:LEU:HD11	1.98	0.64
1:C:168:ARG:NH2	1:D:111:VAL:O	2.31	0.64
1:B:163:MET:HE3	4:B:407:HOH:O	1.98	0.63
1:A:92:TYR:HH	1:A:103:HIS:CD2	2.14	0.63
1:C:139:LEU:HD13	1:C:190:ARG:HA	1.79	0.63
1:A:74:LEU:HD22	1:A:127:PHE:HZ	1.64	0.63
1:C:168:ARG:HH22	1:D:111:VAL:HG22	1.62	0.63
1:A:111:VAL:O	1:A:111:VAL:HG22	1.99	0.62
1:B:148:ARG:HH11	1:B:148:ARG:HG3	1.65	0.62
1:D:111:VAL:O	1:D:111:VAL:HG22	1.99	0.61
1:D:13:LEU:HD23	1:D:48:PHE:HE2	1.65	0.61
1:C:107:ARG:NH2	1:C:160:GLU:OE1	2.33	0.61
1:C:25:GLY:HA2	1:C:108:GLN:HG2	1.83	0.60
1:D:60:LEU:HD21	1:D:119:THR:CG2	2.24	0.60
1:D:122:ARG:O	1:D:126:MET:HG3	2.03	0.59
1:B:104:ARG:O	1:B:108:GLN:HG3	2.02	0.59
1:C:30:THR:HG22	4:C:410:HOH:O	2.02	0.59
1:A:152:ASP:OD1	2:A:301:BER:H61	2.03	0.59
1:B:148:ARG:HH11	1:B:148:ARG:CG	2.15	0.58
1:D:70:MET:HG2	1:D:91:SER:CB	2.33	0.58
1:C:111:VAL:HG12	1:C:111:VAL:O	2.03	0.58
1:B:48:PHE:CE1	1:B:57:THR:HG21	2.38	0.58
1:B:75:ASP:HB3	1:B:78:ILE:CD1	2.11	0.58
1:A:86:ARG:HD2	1:A:181:PHE:CD2	2.38	0.58
1:C:107:ARG:NH1	1:C:157:ALA:HA	2.19	0.57
1:A:156:LEU:HD11	2:A:301:BER:H193	1.86	0.57
1:A:127:PHE:CG	1:A:130:LEU:HB3	2.40	0.56
1:D:184:MET:O	1:D:184:MET:HE2	2.05	0.56
1:B:23:GLN:HG3	4:B:410:HOH:O	2.04	0.56
2:D:301:BER:C6	2:D:301:BER:H192	2.28	0.56
1:C:144:SER:HB2	1:C:146:GLU:OE2	2.06	0.56
1:C:118:GLU:H	1:C:118:GLU:CD	2.08	0.56
1:C:31:ALA:O	1:C:34:ALA:HB3	2.06	0.55
1:D:106:ILE:HD13	1:D:106:ILE:C	2.26	0.55
1:C:139:LEU:CD1	1:C:190:ARG:HA	2.36	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:54:LEU:HA	1:B:57:THR:HG22	1.88	0.55
1:C:30:THR:HG21	1:C:45:PHE:HZ	1.72	0.55
1:C:102:ARG:HG3	3:C:302:SO4:O3	2.07	0.54
1:C:70:MET:CE	1:C:130:LEU:HD22	2.36	0.54
1:C:30:THR:HG21	1:C:45:PHE:CZ	2.43	0.54
1:B:56:ASN:O	1:B:60:LEU:HD13	2.07	0.54
1:D:86:ARG:HD2	1:D:181:PHE:CD2	2.43	0.54
1:C:13:LEU:HD13	1:C:39:VAL:HG21	1.89	0.54
1:A:111:VAL:HG23	1:B:165:PHE:CZ	2.44	0.53
1:D:92:TYR:HE2	1:D:163:MET:HE1	1.73	0.53
1:C:74:LEU:HD12	1:C:84:MET:HE1	1.89	0.53
1:D:84:MET:O	1:D:88:ILE:HG12	2.08	0.53
1:C:33:ILE:HD13	1:C:44:LEU:HD11	1.91	0.53
1:A:157:ALA:HB1	1:B:157:ALA:HB1	1.90	0.53
1:D:86:ARG:NH1	1:D:86:ARG:HG3	2.19	0.53
1:B:89:TRP:CD1	1:B:181:PHE:HB2	2.44	0.53
1:B:13:LEU:HD21	1:B:44:LEU:HA	1.90	0.52
1:B:121:GLN:HE22	1:D:10:LYS:NZ	2.07	0.52
1:A:60:LEU:CD1	1:A:119:THR:HG23	2.38	0.51
1:C:107:ARG:HD3	1:C:156:LEU:HB3	1.92	0.51
1:A:78:ILE:CG2	1:A:83:MET:HG2	2.41	0.51
1:A:84:MET:HG2	1:A:133:LEU:HD11	1.93	0.51
1:A:9:LYS:O	1:A:13:LEU:HD22	2.11	0.51
1:C:117:LYS:HZ2	1:C:117:LYS:HB2	1.75	0.50
1:B:76:ARG:O	1:B:77:SER:HB3	2.12	0.50
1:B:64:GLN:HB2	1:B:126:MET:HE1	1.93	0.50
1:C:152:ASP:HA	2:C:301:BER:H72	1.94	0.50
1:B:148:ARG:NH1	1:B:148:ARG:CG	2.74	0.50
1:B:56:ASN:OD1	1:B:115:LEU:HA	2.12	0.50
1:C:138:VAL:HG12	1:C:139:LEU:O	2.13	0.49
1:C:52:ASP:OD1	1:C:114:LYS:HD2	2.13	0.49
1:A:103:HIS:CE1	1:A:160:GLU:OE1	2.66	0.49
1:D:64:GLN:HB2	1:D:126:MET:CE	2.42	0.49
1:D:184:MET:CE	1:D:188:LEU:HB2	2.42	0.49
1:A:84:MET:HG3	4:A:413:HOH:O	2.11	0.49
1:D:67:CYS:SG	1:D:127:PHE:HB2	2.53	0.49
1:B:182:GLU:HG3	1:B:186:ARG:NH1	2.28	0.49
1:D:54:LEU:C	1:D:54:LEU:HD23	2.32	0.49
1:C:70:MET:HE1	1:C:130:LEU:CD2	2.40	0.49
1:D:29:SER:HG	1:D:32:VAL:HG23	1.77	0.49
1:C:30:THR:HG23	1:C:41:GLU:OE1	2.12	0.49
1:B:92:TYR:OH	1:B:103:HIS:HD2	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:78:ILE:HG23	1:A:83:MET:HG2	1.93	0.48
1:B:168:ARG:HG3	1:B:169:ASP:N	2.27	0.48
1:B:26:ILE:HG21	1:B:112:SER:HB2	1.95	0.48
1:A:169:ASP:N	1:A:170:PRO:HD3	2.28	0.48
1:B:75:ASP:HA	4:B:413:HOH:O	2.13	0.48
1:D:9:LYS:HD3	1:D:39:VAL:CG1	2.44	0.48
1:C:182:GLU:HB3	1:C:186:ARG:NH1	2.28	0.48
1:A:64:GLN:CB	1:A:126:MET:HE1	2.44	0.47
1:D:45:PHE:HA	1:D:48:PHE:O	2.14	0.47
1:B:85:THR:HG22	1:B:181:PHE:HE1	1.79	0.47
1:A:46:ARG:HG2	1:A:46:ARG:HH11	1.79	0.47
1:B:26:ILE:HD12	1:B:114:LYS:NZ	2.29	0.47
1:B:35:ARG:C	1:B:37:ALA:H	2.18	0.47
1:D:71:ILE:HG12	1:D:127:PHE:CE2	2.49	0.47
1:D:67:CYS:HA	1:D:70:MET:HE2	1.96	0.47
1:D:137:SER:N	1:D:190:ARG:HH21	2.13	0.47
1:B:152:ASP:HA	2:B:301:BER:H72	1.96	0.46
1:D:139:LEU:HD21	1:D:189:THR:O	2.15	0.46
2:A:301:BER:H51	2:A:301:BER:H31	1.69	0.46
1:C:182:GLU:HB3	1:C:186:ARG:HH12	1.81	0.46
1:B:8:ASP:OD2	1:B:9:LYS:NZ	2.49	0.46
1:C:146:GLU:H	1:C:146:GLU:CD	2.18	0.46
1:D:43:THR:O	1:D:46:ARG:HB3	2.16	0.46
1:C:64:GLN:HB2	1:C:126:MET:HE1	1.97	0.46
1:A:89:TRP:CD1	1:A:181:PHE:HB2	2.51	0.46
1:C:131:ARG:HG2	1:C:131:ARG:HH11	1.80	0.45
1:C:83:MET:HG2	1:C:87:PHE:CE2	2.51	0.45
1:B:173:ALA:O	1:B:177:ILE:HG13	2.16	0.45
1:D:134:CYS:O	1:D:143:MET:HE1	2.16	0.45
1:A:103:HIS:O	1:A:106:ILE:HG22	2.17	0.45
1:C:10:LYS:O	1:C:13:LEU:N	2.49	0.45
1:D:146:GLU:HG2	1:D:147:TYR:CE2	2.52	0.45
1:A:111:VAL:CG2	1:B:165:PHE:CZ	3.00	0.45
1:A:89:TRP:O	1:A:93:ILE:HG12	2.16	0.45
1:A:139:LEU:HD13	1:A:189:THR:O	2.16	0.45
1:D:9:LYS:HD3	1:D:39:VAL:HG11	1.99	0.45
1:B:21:ILE:HD13	1:B:109:LEU:HD11	1.98	0.45
1:D:92:TYR:HE2	1:D:163:MET:CE	2.30	0.45
1:C:76:ARG:NH2	4:C:413:HOH:O	2.50	0.45
1:C:40:ALA:HB3	1:C:43:THR:OG1	2.17	0.45
1:B:102:ARG:HG3	3:B:302:SO4:O4	2.17	0.45
1:C:157:ALA:HB1	1:D:157:ALA:HB1	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:54:LEU:O	1:B:57:THR:HG22	2.16	0.44
1:B:63:LYS:HE3	1:B:67:CYS:SG	2.57	0.44
1:B:74:LEU:HD11	1:B:127:PHE:HE1	1.82	0.44
1:B:122:ARG:O	1:B:126:MET:HG3	2.17	0.44
1:D:86:ARG:HH12	1:D:90:ASN:ND2	1.99	0.44
1:C:86:ARG:HH21	1:C:178:ALA:HB1	1.83	0.44
1:D:65:ASP:OD2	1:D:102:ARG:NH1	2.51	0.44
1:C:111:VAL:CG1	1:C:111:VAL:O	2.65	0.44
1:A:168:ARG:C	1:A:170:PRO:HD3	2.38	0.44
1:B:127:PHE:CD2	1:B:130:LEU:HB2	2.52	0.44
1:D:51:LYS:O	1:D:55:ILE:HG12	2.17	0.44
1:C:127:PHE:HD2	1:C:130:LEU:HB3	1.81	0.44
1:D:92:TYR:HE1	2:D:301:BER:H202	1.83	0.43
1:C:86:ARG:HA	1:C:181:PHE:CE1	2.53	0.43
1:B:144:SER:O	1:B:148:ARG:HB2	2.18	0.43
1:A:107:ARG:O	1:A:111:VAL:HG12	2.18	0.43
1:C:62:LEU:HD22	1:C:102:ARG:HB3	2.01	0.43
1:C:89:TRP:O	1:C:93:ILE:HG12	2.18	0.43
1:D:64:GLN:HB2	1:D:126:MET:HE2	2.00	0.43
1:A:58:LEU:HD23	1:A:109:LEU:HD21	2.01	0.43
1:C:26:ILE:HD13	4:C:404:HOH:O	2.18	0.43
1:C:85:THR:HG22	1:C:181:PHE:HE1	1.84	0.43
1:B:162:THR:HG21	1:B:177:ILE:HA	2.00	0.43
1:A:159:ALA:O	1:A:163:MET:HG3	2.19	0.42
1:D:86:ARG:NH1	1:D:90:ASN:ND2	2.54	0.42
1:D:92:TYR:CE2	1:D:163:MET:CE	3.02	0.42
1:B:26:ILE:HG21	1:B:112:SER:CB	2.49	0.42
1:B:127:PHE:N	1:B:128:PRO:HD3	2.34	0.42
1:D:59:TYR:CD2	1:D:115:LEU:HD23	2.54	0.42
1:C:83:MET:HG2	1:C:87:PHE:HE2	1.84	0.42
1:C:33:ILE:CG1	1:C:33:ILE:O	2.46	0.42
1:B:103:HIS:HA	1:B:106:ILE:HG22	2.01	0.42
1:C:127:PHE:CD2	1:C:130:LEU:HB3	2.54	0.42
1:C:139:LEU:HD13	1:C:190:ARG:CA	2.48	0.42
1:B:57:THR:HG23	1:B:58:LEU:N	2.34	0.42
1:B:86:ARG:HA	1:B:181:PHE:CE1	2.54	0.42
1:D:184:MET:CE	1:D:188:LEU:HD12	2.50	0.42
1:B:76:ARG:O	1:B:77:SER:CB	2.67	0.42
1:B:169:ASP:OD2	1:B:172:ARG:NH1	2.51	0.42
1:B:135:HIS:HA	1:B:143:MET:CE	2.49	0.42
1:D:59:TYR:HE1	1:D:106:ILE:HD11	1.84	0.42
1:D:173:ALA:O	1:D:177:ILE:HG13	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:34:ALA:CB	1:D:41:GLU:HA	2.50	0.41
1:A:65:ASP:OD2	1:A:102:ARG:NH1	2.48	0.41
1:B:152:ASP:HB3	2:B:301:BER:H192	2.01	0.41
1:B:35:ARG:HG3	1:B:36:ASN:OD1	2.19	0.41
1:B:43:THR:O	1:B:46:ARG:CG	2.62	0.41
1:D:140:MET:HE3	1:D:190:ARG:NH1	2.35	0.41
1:C:84:MET:HG3	1:C:133:LEU:HD21	2.03	0.41
1:A:126:MET:C	1:A:128:PRO:HD3	2.41	0.41
1:A:127:PHE:CE2	1:A:130:LEU:HB2	2.56	0.41
1:A:127:PHE:CD2	1:A:130:LEU:HD23	2.56	0.41
1:C:85:THR:HG21	2:C:301:BER:H172	2.03	0.41
1:B:127:PHE:C	1:B:129:GLU:H	2.24	0.41
1:C:15:GLU:O	1:C:19:GLN:HG2	2.20	0.41
1:C:131:ARG:HG3	1:C:135:HIS:CE1	2.56	0.41
1:A:92:TYR:CE2	1:A:163:MET:HE1	2.56	0.40
1:C:139:LEU:HD22	1:C:188:LEU:C	2.42	0.40
1:B:26:ILE:HD12	1:B:114:LYS:HZ2	1.87	0.40
1:A:185:TRP:CE3	1:A:189:THR:HG21	2.56	0.40
1:B:54:LEU:CA	1:B:57:THR:HG22	2.51	0.40
1:D:112:SER:O	1:D:113:GLU:HB2	2.21	0.40
1:D:60:LEU:HD23	1:D:119:THR:HG23	1.95	0.40
1:B:60:LEU:HD11	1:B:119:THR:CG2	2.33	0.40
1:C:128:PRO:HB3	1:C:131:ARG:HH21	1.83	0.40
1:D:138:VAL:HB	1:D:143:MET:CE	2.51	0.40
1:C:169:ASP:N	1:C:170:PRO:HD3	2.37	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	182/194 (94%)	178 (98%)	4 (2%)	0	100	100
1	B	182/194 (94%)	172 (94%)	8 (4%)	2 (1%)	21	20
1	C	183/194 (94%)	173 (94%)	8 (4%)	2 (1%)	21	20

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	181/194 (93%)	172 (95%)	9 (5%)	0	100	100
All	All	728/776 (94%)	695 (96%)	29 (4%)	4 (0%)	38	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	77	SER
1	C	72	MET
1	C	135	HIS
1	B	36	ASN

### 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	149/158 (94%)	137 (92%)	12 (8%)	17	17
1	B	149/158 (94%)	142 (95%)	7 (5%)	36	46
1	C	149/158 (94%)	138 (93%)	11 (7%)	20	21
1	D	148/158 (94%)	137 (93%)	11 (7%)	20	21
All	All	595/632 (94%)	554 (93%)	41 (7%)	22	24

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	26	ILE
1	A	60	LEU
1	A	84	MET
1	A	97	LEU
1	A	109	LEU
1	A	127	PHE
1	A	135	HIS
1	A	139	LEU
1	A	158	LEU
1	A	179	LEU

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Mol	Chain	Res	Type
1	A	188	LEU
1	B	9	LYS
1	B	11	GLN
1	B	46	ARG
1	B	124	ASP
1	B	139	LEU
1	B	148	ARG
1	B	179	LEU
1	C	8	ASP
1	C	9	LYS
1	C	26	ILE
1	C	33	ILE
1	C	60	LEU
1	C	107	ARG
1	C	117	LYS
1	C	158	LEU
1	C	168	ARG
1	C	170	PRO
1	C	188	LEU
1	D	14	LEU
1	D	26	ILE
1	D	86	ARG
1	D	106	ILE
1	D	109	LEU
1	D	115	LEU
1	D	154	LEU
1	D	156	LEU
1	D	168	ARG
1	D	179	LEU
1	D	184	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	90	ASN
1	A	98	ASN
1	A	103	HIS
1	A	108	GLN
1	A	121	GLN
1	B	11	GLN
1	B	61	HIS

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Mol	Chain	Res	Type
1	B	68	GLN
1	B	90	ASN
1	B	99	HIS
1	B	103	HIS
1	B	108	GLN
1	B	121	GLN
1	C	64	GLN
1	C	103	HIS
1	D	90	ASN
1	D	103	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

8 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	BER	A	301	-	29,29,29	3.95	18 (62%)	43,43,43	4.58	23 (53%)
3	SO4	A	302	-	4,4,4	0.74	0	6,6,6	0.44	0
2	BER	B	301	-	29,29,29	3.91	18 (62%)	43,43,43	4.66	21 (48%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	B	302	-	4,4,4	0.41	0	6,6,6	0.43	0
2	BER	C	301	-	29,29,29	4.21	15 (51%)	43,43,43	3.82	23 (53%)
3	SO4	C	302	-	4,4,4	0.33	0	6,6,6	0.65	0
2	BER	D	301	-	29,29,29	3.91	18 (62%)	43,43,43	4.29	19 (44%)
3	SO4	D	302	-	4,4,4	0.52	0	6,6,6	0.31	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	BER	A	301	-	-	0/4/19/19	0/0/5/5
3	SO4	A	302	-	-	0/0/0/0	0/0/0/0
2	BER	B	301	-	-	0/4/19/19	0/0/5/5
3	SO4	B	302	-	-	0/0/0/0	0/0/0/0
2	BER	C	301	-	-	0/4/19/19	0/0/5/5
3	SO4	C	302	-	-	0/0/0/0	0/0/0/0
2	BER	D	301	-	-	0/4/19/19	0/0/5/5
3	SO4	D	302	-	-	0/0/0/0	0/0/0/0

All (69) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	BER	C7-N1	-11.44	1.33	1.48
2	A	301	BER	C7-N1	-11.38	1.33	1.48
2	B	301	BER	C7-N1	-11.33	1.33	1.48
2	C	301	BER	C7-N1	-9.31	1.36	1.48
2	C	301	BER	C15-C12	9.14	1.60	1.43
2	A	301	BER	C7-C10	-7.82	1.37	1.50
2	B	301	BER	C10-C4	-7.71	1.36	1.51
2	C	301	BER	C2-C4	7.59	1.50	1.40
2	D	301	BER	C10-C4	-7.53	1.37	1.51
2	D	301	BER	C7-C10	-7.33	1.38	1.50
2	B	301	BER	C7-C10	-7.29	1.38	1.50
2	A	301	BER	C10-C4	-7.22	1.37	1.51
2	C	301	BER	C3-C1	7.03	1.51	1.36
2	C	301	BER	O3-C15	6.84	1.48	1.38
2	C	301	BER	C10-C4	-6.15	1.39	1.51
2	C	301	BER	C18-C15	5.57	1.51	1.38
2	B	301	BER	C3-C1	5.44	1.48	1.36
2	D	301	BER	C3-C1	5.22	1.47	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	BER	C3-C1	5.04	1.47	1.36
2	D	301	BER	C2-C4	4.92	1.47	1.40
2	C	301	BER	C6-C12	4.90	1.51	1.41
2	B	301	BER	C13-C8	-4.78	1.29	1.41
2	B	301	BER	C2-C4	4.73	1.46	1.40
2	C	301	BER	C13-C16	4.65	1.46	1.36
2	B	301	BER	C16-C18	-4.60	1.29	1.39
2	D	301	BER	C16-C18	-4.25	1.30	1.39
2	A	301	BER	O1-C11	-4.19	1.32	1.38
2	D	301	BER	C13-C8	-4.19	1.31	1.41
2	A	301	BER	C2-C4	4.15	1.46	1.40
2	A	301	BER	C18-C15	4.09	1.48	1.38
2	A	301	BER	C13-C8	-4.09	1.31	1.41
2	C	301	BER	C3-C8	4.06	1.51	1.42
2	D	301	BER	C6-C12	4.03	1.49	1.41
2	A	301	BER	C13-C16	3.98	1.45	1.36
2	C	301	BER	C12-C8	3.93	1.50	1.42
2	A	301	BER	C6-C12	3.87	1.49	1.41
2	A	301	BER	O4-C18	3.69	1.43	1.37
2	C	301	BER	O4-C18	3.62	1.42	1.37
2	A	301	BER	C15-C12	3.52	1.49	1.43
2	B	301	BER	O4-C18	3.45	1.42	1.37
2	B	301	BER	C6-C12	3.32	1.47	1.41
2	B	301	BER	C13-C16	3.32	1.43	1.36
2	D	301	BER	C15-C12	3.30	1.49	1.43
2	D	301	BER	O4-C18	3.30	1.42	1.37
2	D	301	BER	C18-C15	3.26	1.46	1.38
2	B	301	BER	O1-C11	-3.13	1.33	1.38
2	D	301	BER	C13-C16	3.12	1.43	1.36
2	B	301	BER	O2-C14	-3.09	1.33	1.38
2	A	301	BER	O2-C14	-3.06	1.33	1.38
2	D	301	BER	O1-C11	-3.05	1.33	1.38
2	A	301	BER	C16-C18	-3.03	1.33	1.39
2	B	301	BER	C15-C12	3.03	1.48	1.43
2	A	301	BER	C12-C8	2.96	1.48	1.42
2	B	301	BER	C3-C8	2.78	1.48	1.42
2	D	301	BER	O2-C14	-2.76	1.34	1.38
2	D	301	BER	C1-N1	-2.59	1.31	1.36
2	D	301	BER	C3-C8	2.53	1.47	1.42
2	C	301	BER	C14-C11	2.48	1.46	1.39
2	D	301	BER	C12-C8	2.47	1.47	1.42
2	B	301	BER	C1-N1	-2.45	1.31	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	BER	C12-C8	2.43	1.47	1.42
2	B	301	BER	C18-C15	2.43	1.44	1.38
2	D	301	BER	C5-C2	-2.25	1.36	1.39
2	C	301	BER	C6-N1	2.21	1.35	1.33
2	A	301	BER	C3-C8	2.16	1.46	1.42
2	A	301	BER	C2-C1	-2.16	1.43	1.47
2	B	301	BER	C5-C2	-2.08	1.36	1.39
2	A	301	BER	C1-N1	-2.07	1.32	1.36
2	C	301	BER	C13-C8	-2.04	1.36	1.41

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	BER	C6-C12-C15	-17.15	106.44	121.77
2	A	301	BER	C6-C12-C15	-15.13	108.25	121.77
2	D	301	BER	C6-C12-C15	-12.12	110.94	121.77
2	C	301	BER	C6-C12-C15	-11.29	111.68	121.77
2	B	301	BER	C18-C15-C12	-10.97	103.34	120.19
2	D	301	BER	C18-C15-C12	-10.46	104.13	120.19
2	A	301	BER	C3-C1-N1	10.42	127.46	117.22
2	A	301	BER	C3-C1-C2	-9.77	112.58	124.56
2	C	301	BER	C7-N1-C6	-9.02	112.57	118.05
2	A	301	BER	C18-C15-C12	-8.88	106.55	120.19
2	D	301	BER	C3-C1-N1	8.76	125.83	117.22
2	B	301	BER	C10-C7-N1	8.35	124.54	110.08
2	D	301	BER	C10-C7-N1	8.32	124.47	110.08
2	C	301	BER	C3-C1-N1	8.17	125.25	117.22
2	B	301	BER	C1-C3-C8	-8.07	110.17	121.26
2	D	301	BER	C1-C3-C8	-8.00	110.27	121.26
2	D	301	BER	C3-C1-C2	-7.76	115.03	124.56
2	A	301	BER	C1-C3-C8	-7.76	110.61	121.26
2	C	301	BER	C18-C15-C12	-7.56	108.58	120.19
2	A	301	BER	O3-C15-C12	-7.45	103.97	117.23
2	B	301	BER	C3-C1-N1	7.37	124.46	117.22
2	A	301	BER	C7-C10-C4	7.37	119.91	109.87
2	C	301	BER	C3-C1-C2	-6.83	116.18	124.56
2	A	301	BER	C10-C7-N1	6.82	121.89	110.08
2	D	301	BER	C15-C12-C8	-6.73	106.56	118.66
2	B	301	BER	O3-C15-C18	-6.48	103.76	120.35
2	B	301	BER	C7-C10-C4	6.45	118.67	109.87
2	D	301	BER	C7-C10-C4	6.42	118.61	109.87
2	B	301	BER	C3-C1-C2	-6.39	116.71	124.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	BER	C10-C7-N1	6.12	120.67	110.08
2	B	301	BER	C7-N1-C6	-6.04	114.38	118.05
2	C	301	BER	C7-C10-C4	5.91	117.92	109.87
2	D	301	BER	O3-C15-C18	-5.50	106.26	120.35
2	C	301	BER	C1-C3-C8	-5.40	113.84	121.26
2	B	301	BER	O3-C15-C12	-5.38	107.66	117.23
2	B	301	BER	C6-C12-C8	-4.96	111.83	117.58
2	A	301	BER	O3-C15-C18	-4.78	108.11	120.35
2	D	301	BER	O3-C15-C12	-4.46	109.29	117.23
2	C	301	BER	C12-C6-N1	-4.38	119.01	121.85
2	D	301	BER	C6-C12-C8	-4.33	112.56	117.58
2	B	301	BER	C15-C12-C8	-4.11	111.27	118.66
2	C	301	BER	C7-N1-C1	3.93	124.87	118.81
2	A	301	BER	C7-N1-C6	-3.80	115.74	118.05
2	A	301	BER	C5-C2-C4	3.79	124.00	119.19
2	A	301	BER	O4-C18-C16	-3.68	118.30	124.37
2	D	301	BER	C5-C2-C4	3.66	123.84	119.19
2	B	301	BER	C3-C8-C12	3.65	124.12	119.18
2	C	301	BER	C16-C18-C15	3.64	123.38	119.81
2	A	301	BER	C15-C12-C8	-3.63	112.14	118.66
2	D	301	BER	C14-C9-C4	-3.61	114.32	120.98
2	B	301	BER	C14-C9-C4	-3.52	114.48	120.98
2	A	301	BER	C2-C5-C11	-3.49	113.67	120.62
2	A	301	BER	C1-C2-C4	-3.39	116.92	120.22
2	D	301	BER	C2-C5-C11	-3.37	113.92	120.62
2	B	301	BER	C6-N1-C1	3.29	124.16	122.46
2	B	301	BER	C2-C5-C11	-3.17	114.31	120.62
2	D	301	BER	C17-O2-C14	3.15	109.89	105.36
2	D	301	BER	C7-N1-C6	-3.06	116.19	118.05
2	B	301	BER	C13-C8-C3	-2.92	117.08	122.07
2	A	301	BER	C20-O4-C18	-2.87	113.33	117.59
2	C	301	BER	O2-C17-O1	-2.82	103.20	108.09
2	A	301	BER	C14-C9-C4	-2.78	115.86	120.98
2	C	301	BER	C1-C2-C4	-2.69	117.60	120.22
2	C	301	BER	C14-C9-C4	-2.67	116.06	120.98
2	D	301	BER	C16-C13-C8	-2.62	116.71	120.83
2	A	301	BER	C17-O1-C11	2.55	109.03	105.36
2	C	301	BER	C2-C5-C11	-2.54	115.57	120.62
2	A	301	BER	C7-N1-C1	2.49	122.65	118.81
2	D	301	BER	C17-O1-C11	2.47	108.91	105.36
2	B	301	BER	C5-C2-C4	2.42	122.27	119.19
2	A	301	BER	C6-C12-C8	-2.42	114.78	117.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	BER	O3-C15-C12	-2.39	112.97	117.23
2	C	301	BER	C20-O4-C18	-2.29	114.19	117.59
2	C	301	BER	C17-O2-C14	2.29	108.65	105.36
2	B	301	BER	C17-O2-C14	2.28	108.64	105.36
2	C	301	BER	C19-O3-C15	2.28	120.53	115.11
2	B	301	BER	C17-O1-C11	2.26	108.61	105.36
2	C	301	BER	C13-C8-C12	2.24	122.20	119.09
2	B	301	BER	C9-C4-C2	2.21	122.43	119.54
2	A	301	BER	C16-C18-C15	2.10	121.88	119.81
2	A	301	BER	C2-C1-N1	2.10	119.94	118.21
2	C	301	BER	O2-C14-C11	-2.09	107.12	109.78
2	A	301	BER	C13-C8-C3	-2.08	118.50	122.07
2	C	301	BER	C17-O1-C11	2.03	108.28	105.36
2	C	301	BER	C5-C2-C4	2.02	121.76	119.19
2	D	301	BER	C3-C8-C12	2.00	121.89	119.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	184/194 (94%)	0.48	10 (5%)	25 35	21, 41, 69, 84	12 (6%)
1	B	184/194 (94%)	1.50	50 (27%)	1 1	23, 66, 101, 111	3 (1%)
1	C	185/194 (95%)	0.64	18 (9%)	8 13	23, 47, 78, 85	5 (2%)
1	D	183/194 (94%)	0.83	22 (12%)	5 8	24, 51, 86, 93	4 (2%)
All	All	736/776 (94%)	0.86	100 (13%)	4 6	21, 49, 90, 111	24 (3%)

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	123	ALA	7.9
1	B	39	VAL	7.1
1	B	33	ILE	6.2
1	B	9	LYS	5.6
1	B	34	ALA	5.5
1	B	13	LEU	5.3
1	B	124	ASP	4.8
1	B	44	LEU	4.6
1	C	36	ASN	4.5
1	B	42	GLY	4.4
1	B	31	ALA	4.3
1	B	8	ASP	4.3
1	B	38	GLY	4.3
1	D	34	ALA	4.2
1	B	12	ALA	4.2
1	B	50	THR	4.1
1	D	11	GLN	4.1
1	B	78	ILE	4.1
1	B	17	ALA	4.1
1	B	27	ALA	3.9
1	B	46	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	77	SER	3.9
1	B	37	ALA	3.9
1	D	39	VAL	3.5
1	D	132	ASP	3.5
1	B	75	ASP	3.4
1	C	35	ARG	3.4
1	B	133	LEU	3.4
1	B	10	LYS	3.4
1	B	79	THR	3.2
1	B	48	PHE	3.2
1	D	78	ILE	3.1
1	D	130	LEU	3.1
1	B	36	ASN	3.1
1	B	11	GLN	3.1
1	C	127	PHE	3.0
1	C	121	GLN	3.0
1	D	13	LEU	2.9
1	B	74	LEU	2.8
1	C	7	GLU	2.8
1	D	131	ARG	2.8
1	B	191	GLU	2.7
1	A	87	PHE	2.7
1	B	57	THR	2.7
1	D	76	ARG	2.7
1	B	125	ASP	2.7
1	C	126	MET	2.7
1	B	131	ARG	2.7
1	A	123	ALA	2.6
1	B	156	LEU	2.6
1	B	45	PHE	2.6
1	C	8	ASP	2.6
1	D	133	LEU	2.6
1	B	155	PHE	2.5
1	C	44	LEU	2.5
1	C	117	LYS	2.5
1	A	135	HIS	2.5
1	C	118	GLU	2.4
1	B	158	LEU	2.4
1	A	134	CYS	2.4
1	B	132	ASP	2.4
1	B	51	LYS	2.4
1	B	14	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	135	HIS	2.4
1	B	159	ALA	2.3
1	C	32	VAL	2.3
1	D	9	LYS	2.3
1	B	29	SER	2.3
1	D	79	THR	2.3
1	B	129	GLU	2.3
1	B	24	SER	2.3
1	A	125	ASP	2.3
1	B	76	ARG	2.2
1	B	122	ARG	2.2
1	A	119	THR	2.2
1	C	13	LEU	2.2
1	C	140	MET	2.2
1	C	38	GLY	2.2
1	D	38	GLY	2.2
1	D	68	GLN	2.2
1	D	31	ALA	2.2
1	C	47	TYR	2.2
1	B	84	MET	2.2
1	D	140	MET	2.2
1	B	77	SER	2.2
1	C	31	ALA	2.1
1	A	127	PHE	2.1
1	B	163	MET	2.1
1	D	36	ASN	2.1
1	D	74	LEU	2.1
1	A	10	LYS	2.1
1	D	8	ASP	2.1
1	A	45	PHE	2.1
1	B	41	GLU	2.1
1	B	64	GLN	2.0
1	C	163	MET	2.0
1	B	127	PHE	2.0
1	A	158	LEU	2.0
1	D	154	LEU	2.0
1	C	11	GLN	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates

There are no carbohydrates 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. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	BER	D	301	25/25	0.23	1.14	57,62,71,73	0
2	BER	B	301	25/25	0.21	1.06	60,64,72,76	0
2	BER	C	301	25/25	0.17	0.32	45,49,54,58	0
2	BER	A	301	25/25	0.17	0.09	47,53,61,67	0
3	SO4	D	302	5/5	0.13	-0.56	59,60,61,64	0
3	SO4	C	302	5/5	0.12	-0.89	62,65,67,67	0
3	SO4	B	302	5/5	0.08	-1.57	79,80,81,82	0
3	SO4	A	302	5/5	0.11	-1.67	46,50,54,55	0

### 6.5 Other polymers

There are no such residues in this entry.