



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2020 – 10:37 AM BST

PDB ID : 2V43  
Title : Crystal structure of RseB: a sensor for periplasmic stress response in E. coli  
Authors : Wollmann, P.; Zeth, K.  
Deposited on : 2007-06-27  
Resolution : 2.37 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : **FAILED**  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.4.dev1

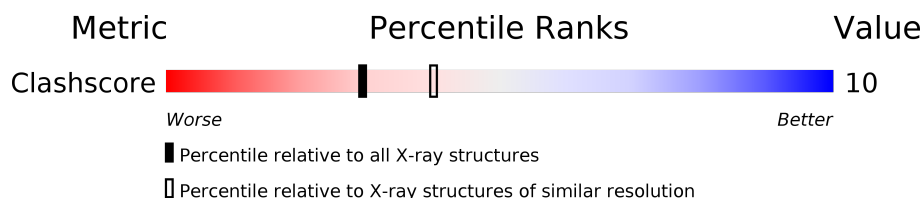
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.37 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))
Clashscore	141614	6082 (2.40-2.36)

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 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  $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	303	 70% 14% • 14%
1	B	303	 73% 16% • 9%
1	C	303	 72% 17% • 9%

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	CYS	A	1315	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6687 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.

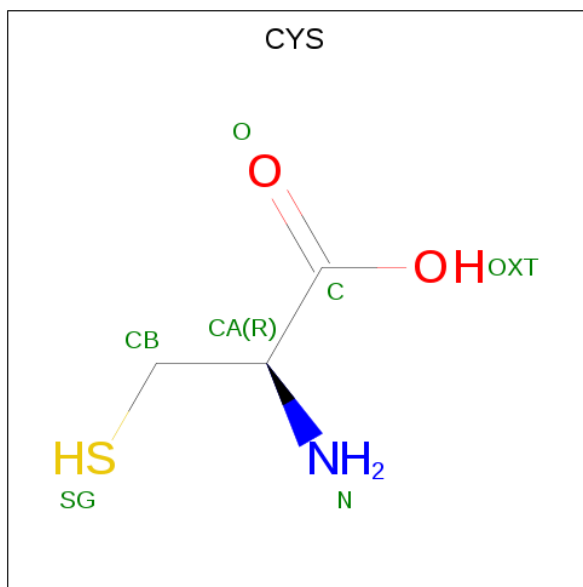
- Molecule 1 is a protein called SIGMA-E FACTOR REGULATORY PROTEIN RSEB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			2077	1307	365	398	7			
1	B	276	Total	C	N	O	S	0	0	0
			2201	1382	391	420	8			
1	C	276	Total	C	N	O	S	0	1	0
			2187	1374	391	413	9			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	MET	-	expression tag	UNP P0AFX9
A	319	HIS	-	expression tag	UNP P0AFX9
A	320	HIS	-	expression tag	UNP P0AFX9
A	321	HIS	-	expression tag	UNP P0AFX9
A	322	HIS	-	expression tag	UNP P0AFX9
A	323	HIS	-	expression tag	UNP P0AFX9
A	324	HIS	-	expression tag	UNP P0AFX9
B	22	MET	-	expression tag	UNP P0AFX9
B	319	HIS	-	expression tag	UNP P0AFX9
B	320	HIS	-	expression tag	UNP P0AFX9
B	321	HIS	-	expression tag	UNP P0AFX9
B	322	HIS	-	expression tag	UNP P0AFX9
B	323	HIS	-	expression tag	UNP P0AFX9
B	324	HIS	-	expression tag	UNP P0AFX9
C	22	MET	-	expression tag	UNP P0AFX9
C	319	HIS	-	expression tag	UNP P0AFX9
C	320	HIS	-	expression tag	UNP P0AFX9
C	321	HIS	-	expression tag	UNP P0AFX9
C	322	HIS	-	expression tag	UNP P0AFX9
C	323	HIS	-	expression tag	UNP P0AFX9
C	324	HIS	-	expression tag	UNP P0AFX9

- Molecule 2 is CYSTEINE (three-letter code: CYS) (formula: C<sub>3</sub>H<sub>7</sub>NO<sub>2</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			7	3	1	2	1		
2	A	1	Total	C	N	O	S	0	0
			7	3	1	2	1		
2	B	1	Total	C	N	O	S	0	0
			7	3	1	2	1		
2	B	1	Total	C	N	O	S	0	0
			7	3	1	2	1		

- Molecule 3 is water.

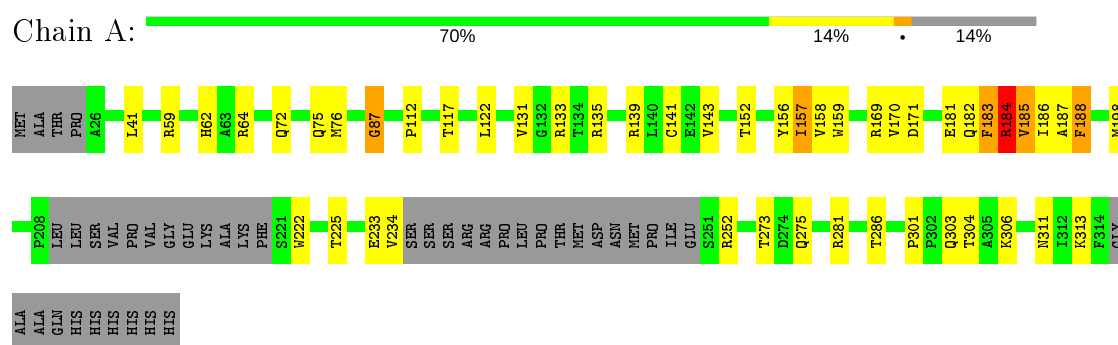
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	81	Total	O	0	0
			81	81		
3	B	76	Total	O	0	0
			76	76		
3	C	37	Total	O	0	0
			37	37		

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

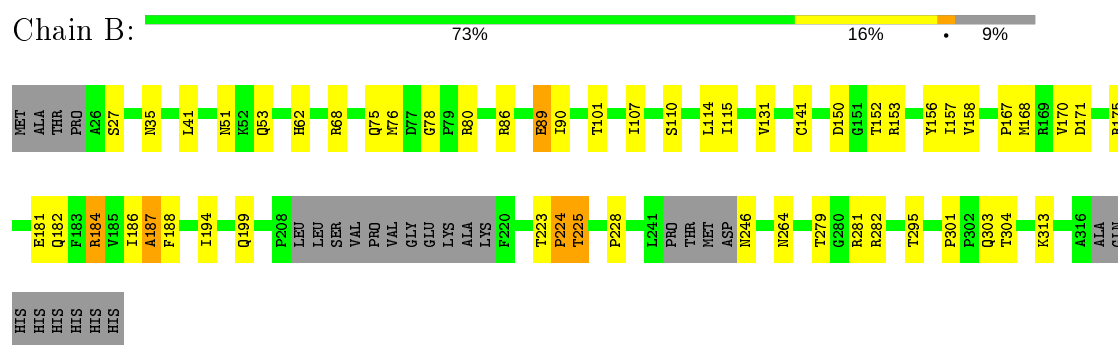
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

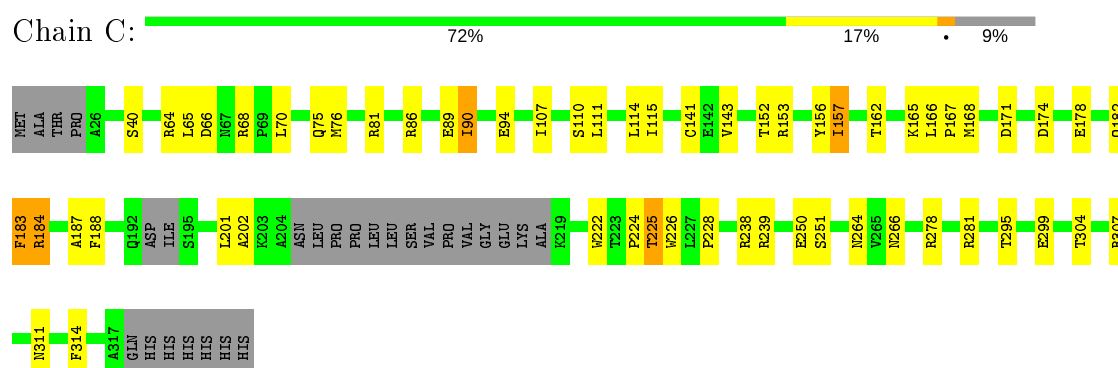
#### • Molecule 1: SIGMA-E FACTOR REGULATORY PROTEIN RSEB



#### • Molecule 1: SIGMA-E FACTOR REGULATORY PROTEIN RSEB



#### • Molecule 1: SIGMA-E FACTOR REGULATORY PROTEIN RSEB



## 4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.60Å 200.70Å 109.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.37	Depositor
% Data completeness (in resolution range)	98.5 (20.00-2.37)	Depositor
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.72 (at 2.38Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.228 , 0.278	Depositor
Wilson B-factor (Å <sup>2</sup> )	49.2	Xtriage
Anisotropy	0.235	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6687	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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

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	0.56	1/2116 (0.0%)	0.85	10/2870 (0.3%)
1	B	0.56	1/2243 (0.0%)	0.86	10/3040 (0.3%)
1	C	0.46	0/2228	0.74	6/3018 (0.2%)
All	All	0.53	2/6587 (0.0%)	0.82	26/8928 (0.3%)

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	7
1	B	0	7
1	C	0	6
All	All	0	20

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	246	ASN	CG-ND2	6.67	1.49	1.32
1	A	313	LYS	C-N	5.54	1.46	1.34

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	225	THR	N-CA-C	-8.16	88.96	111.00
1	B	224	PRO	N-CA-C	8.15	133.29	112.10
1	A	185	VAL	N-CA-C	7.95	132.46	111.00
1	B	90	ILE	N-CA-C	7.57	131.43	111.00
1	A	184	ARG	NE-CZ-NH2	-7.39	116.60	120.30
1	C	76	MET	N-CA-C	-7.12	91.76	111.00
1	B	225	THR	N-CA-C	-7.03	92.02	111.00
1	B	225	THR	N-CA-CB	6.26	122.19	110.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	114	LEU	CA-CB-CG	6.16	129.47	115.30
1	C	157	ILE	N-CA-C	6.11	127.49	111.00
1	B	76	MET	N-CA-C	-5.91	95.04	111.00
1	B	188	PHE	N-CA-CB	5.90	121.21	110.60
1	A	184	ARG	N-CA-C	-5.85	95.20	111.00
1	A	76	MET	N-CA-C	-5.83	95.26	111.00
1	A	157	ILE	N-CA-C	5.69	126.36	111.00
1	C	184	ARG	N-CA-CB	5.68	120.82	110.60
1	B	187	ALA	N-CA-CB	-5.65	102.19	110.10
1	C	90	ILE	N-CA-C	-5.57	95.95	111.00
1	A	186	ILE	N-CA-C	-5.56	95.99	111.00
1	A	87	GLY	N-CA-C	-5.38	99.65	113.10
1	C	114	LEU	CA-CB-CG	5.34	127.59	115.30
1	A	184	ARG	N-CA-CB	5.32	120.17	110.60
1	A	184	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	B	184	ARG	N-CA-CB	-5.12	101.38	110.60
1	B	157	ILE	N-CA-C	5.10	124.78	111.00
1	A	188	PHE	N-CA-CB	5.10	119.78	110.60

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	TYR	Peptide
1	A	183	PHE	Peptide
1	A	184	ARG	Peptide
1	A	185	VAL	Peptide
1	A	187	ALA	Peptide
1	A	75	GLN	Peptide
1	A	87	GLY	Peptide
1	B	156	TYR	Peptide
1	B	186	ILE	Peptide
1	B	187	ALA	Peptide
1	B	223	THR	Peptide
1	B	224	PRO	Peptide
1	B	75	GLN	Peptide
1	B	89	GLU	Peptide
1	C	156	TYR	Peptide
1	C	183	PHE	Peptide
1	C	187	ALA	Peptide
1	C	224	PRO	Peptide
1	C	75	GLN	Peptide

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Group
1	C	89	GLU	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	2077	0	2049	45	0
1	B	2201	0	2179	38	0
1	C	2187	0	2159	43	0
2	A	14	0	8	6	0
2	B	14	0	8	3	0
3	A	81	0	0	1	0
3	B	76	0	0	1	0
3	C	37	0	0	3	0
All	All	6687	0	6403	123	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 10.

All (123) 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:C:225:THR:HB	1:C:311:ASN:O	1.54	1.06
1:C:152:THR:HG21	1:C:228:PRO:HB3	1.40	1.00
1:A:183:PHE:HB2	2:A:1315:CYS:HB2	1.41	0.99
1:B:301:PRO:HB3	2:B:1317:CYS:HB3	1.47	0.96
1:C:182:GLN:NE2	1:C:184:ARG:HD3	1.81	0.95
1:C:152:THR:CG2	1:C:153:ARG:HD2	1.98	0.94
1:B:152:THR:HG22	1:B:153:ARG:HD2	1.50	0.93
1:A:183:PHE:CB	2:A:1315:CYS:HB2	2.00	0.91
1:A:182:GLN:HE22	1:A:184:ARG:HH11	1.18	0.90
1:A:64:ARG:HG2	1:A:64:ARG:HH11	1.34	0.90
1:A:182:GLN:NE2	1:A:184:ARG:HD3	1.89	0.87
1:B:182:GLN:HE22	1:B:184:ARG:HH11	1.20	0.86
1:A:152:THR:O	1:A:304:THR:HG21	1.79	0.83
1:B:152:THR:O	1:B:304:THR:HG21	1.78	0.82

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:ASP:OD1	1:C:182:GLN:HG3	1.81	0.81
1:B:152:THR:HG21	1:B:228:PRO:HB3	1.63	0.80
1:B:107:ILE:O	1:B:110:SER:HB2	1.81	0.79
1:C:182:GLN:HE22	1:C:184:ARG:HD3	1.46	0.78
1:A:112:PRO:HG3	2:A:1315:CYS:HB3	1.67	0.77
1:C:152:THR:HG23	1:C:153:ARG:HD2	1.68	0.76
1:A:225:THR:HB	1:A:311:ASN:O	1.85	0.76
1:C:152:THR:O	1:C:304:THR:HG21	1.86	0.75
1:A:59:ARG:HH21	1:A:72:GLN:HE22	1.34	0.74
1:A:64:ARG:HG2	1:A:64:ARG:NH1	1.97	0.74
1:B:295:THR:HG22	3:B:2072:HOH:O	1.88	0.74
1:A:182:GLN:HE21	1:A:184:ARG:HD3	1.51	0.73
1:C:152:THR:HG22	1:C:153:ARG:HD2	1.73	0.69
1:C:115:ILE:HG12	1:C:167:PRO:HG3	1.75	0.69
1:A:234:VAL:HG13	1:C:66:ASP:HB3	1.74	0.69
1:B:35:ASN:C	1:B:35:ASN:HD22	1.96	0.69
1:B:27:SER:H	2:B:1318:CYS:HB2	1.57	0.68
1:B:181:GLU:OE1	1:B:281:ARG:NH2	2.25	0.67
1:B:182:GLN:NE2	1:B:184:ARG:HD3	2.10	0.67
1:C:143:VAL:HG12	1:C:157:ILE:HD11	1.76	0.65
1:B:150:ASP:OD1	1:B:152:THR:HB	1.97	0.65
1:A:41:LEU:O	1:A:62:HIS:HD2	1.81	0.64
1:C:281:ARG:HG2	1:C:299:GLU:HG3	1.80	0.64
1:A:182:GLN:HE22	1:A:184:ARG:NH1	1.94	0.63
1:C:110:SER:OG	1:C:111:LEU:N	2.32	0.63
1:C:182:GLN:HE22	1:C:184:ARG:HH11	1.46	0.63
1:C:70:LEU:HD21	1:C:202:ALA:HB2	1.80	0.62
1:A:143:VAL:CG1	1:A:157:ILE:HD11	2.29	0.62
1:A:275:GLN:HG2	1:A:286:THR:HB	1.80	0.61
1:C:81:ARG:HB3	1:C:81:ARG:HH11	1.65	0.61
1:B:182:GLN:HE21	1:B:184:ARG:HD3	1.65	0.61
1:B:264:ASN:HB2	1:B:295:THR:HG23	1.83	0.60
1:C:40:SER:HA	1:C:64:ARG:HD2	1.84	0.60
1:C:143:VAL:CG1	1:C:157:ILE:HD11	2.32	0.59
1:B:41:LEU:O	1:B:62:HIS:HD2	1.86	0.58
1:A:273:THR:O	1:A:286:THR:HG22	2.02	0.58
1:A:198:MET:CE	3:A:2021:HOH:O	2.53	0.57
1:C:281:ARG:HG2	1:C:299:GLU:CG	2.35	0.56
1:A:135:ARG:HG3	1:B:78:GLY:O	2.06	0.56
1:A:233:GLU:OE2	1:A:252:ARG:HG2	2.06	0.56
1:B:35:ASN:C	1:B:35:ASN:ND2	2.59	0.56

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:THR:HG22	1:C:153:ARG:CD	2.36	0.55
1:C:251:SER:OG	1:C:264:ASN:ND2	2.36	0.55
1:B:152:THR:HG22	1:B:153:ARG:CD	2.30	0.55
1:C:182:GLN:HE21	1:C:184:ARG:HD3	1.68	0.54
1:B:194:ILE:HG12	1:B:199:GLN:HG2	1.88	0.54
1:A:133:ARG:CB	1:A:141:CYS:O	2.54	0.54
1:B:301:PRO:CB	2:B:1317:CYS:HB3	2.30	0.54
1:C:162:THR:O	1:C:165:LYS:HE3	2.07	0.54
1:A:181:GLU:OE2	1:A:281:ARG:HG3	2.07	0.54
1:A:59:ARG:HH21	1:A:72:GLN:NE2	2.05	0.54
1:C:182:GLN:HE21	1:C:184:ARG:HB2	1.73	0.54
1:C:183:PHE:C	1:C:183:PHE:CD2	2.81	0.53
1:B:152:THR:CG2	1:B:153:ARG:HD2	2.33	0.53
1:C:107:ILE:O	1:C:110:SER:HB3	2.09	0.52
1:B:89:GLU:HB3	1:B:101:THR:CG2	2.40	0.52
1:C:81:ARG:HG2	1:C:94:GLU:HG3	1.92	0.51
1:A:183:PHE:HB3	2:A:1315:CYS:HB2	1.86	0.51
1:A:301:PRO:HB2	1:A:304:THR:HG23	1.93	0.50
1:B:115:ILE:HG12	1:B:167:PRO:HG3	1.94	0.49
1:A:143:VAL:HG12	1:A:157:ILE:HD11	1.94	0.48
1:A:139:ARG:CZ	1:A:188:PHE:HE1	2.25	0.48
1:A:222:TRP:CH2	1:A:252:ARG:HD3	2.49	0.48
1:A:133:ARG:HA	1:A:143:VAL:HG23	1.96	0.47
1:C:174:ASP:HB2	1:C:178:GLU:O	2.14	0.47
1:A:112:PRO:CG	2:A:1315:CYS:HB3	2.41	0.47
1:A:64:ARG:HH11	1:A:64:ARG:CG	2.14	0.47
1:C:295:THR:HG22	3:C:2027:HOH:O	2.15	0.47
1:A:158:VAL:HG13	1:A:170:VAL:HG22	1.96	0.46
1:C:174:ASP:OD2	1:C:278:ARG:NH2	2.48	0.46
1:A:112:PRO:HG3	2:A:1315:CYS:CB	2.41	0.46
1:C:86:ARG:HD3	1:C:202:ALA:HA	1.97	0.46
1:B:89:GLU:HB3	1:B:101:THR:HG21	1.96	0.46
1:A:133:ARG:HB3	1:A:141:CYS:O	2.15	0.46
1:B:175:ARG:HH12	1:B:303:GLN:NE2	2.14	0.45
1:B:152:THR:HG21	1:B:228:PRO:CB	2.41	0.45
1:C:250:GLU:O	1:C:264:ASN:HA	2.17	0.45
1:C:86:ARG:HG3	1:C:201:LEU:HD22	1.98	0.45
1:A:171:ASP:OD1	1:A:182:GLN:HG3	2.16	0.45
1:A:303:GLN:HA	1:A:306:LYS:HE2	1.99	0.45
1:B:225:THR:HG21	1:B:313:LYS:CG	2.47	0.44
1:B:141:CYS:SG	1:B:168:MET:HG3	2.58	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:239:ARG:HG2	3:C:2025:HOH:O	2.18	0.44
1:C:266:ASN:ND2	3:C:2027:HOH:O	2.50	0.44
1:A:117:THR:HG21	1:A:122:LEU:HD22	1.98	0.44
1:C:238:ARG:HD3	1:C:250:GLU:HG2	2.00	0.44
1:B:131:VAL:HG22	1:B:131:VAL:O	2.18	0.43
1:B:158:VAL:HG13	1:B:170:VAL:HG22	2.00	0.43
1:A:182:GLN:HE21	1:A:184:ARG:HB2	1.83	0.43
1:C:226:TRP:CE3	1:C:307:ARG:HG2	2.54	0.43
1:B:51:ASN:OD1	1:B:53:GLN:HB2	2.19	0.42
1:C:65:LEU:O	1:C:68:ARG:HB3	2.19	0.42
1:A:135:ARG:CZ	1:B:80:ARG:HG3	2.49	0.42
1:C:141:CYS:SG	1:C:168:MET:HG3	2.60	0.42
1:C:90:ILE:HD13	1:C:90:ILE:HA	1.95	0.42
1:C:166:LEU:HD12	1:C:188:PHE:HE2	1.85	0.42
1:B:279:THR:OG1	1:B:282:ARG:HB3	2.20	0.41
1:A:59:ARG:NH2	1:A:72:GLN:HE22	2.10	0.41
1:A:169:ARG:HA	1:A:183:PHE:O	2.21	0.41
1:B:131:VAL:O	1:B:131:VAL:CG2	2.69	0.41
1:C:222:TRP:HB3	1:C:314:PHE:CD2	2.56	0.41
1:B:264:ASN:HB2	1:B:295:THR:CG2	2.48	0.41
1:B:175:ARG:HD3	1:B:301:PRO:CG	2.51	0.40
1:A:182:GLN:HE22	1:A:184:ARG:HD3	1.79	0.40
1:B:171:ASP:OD1	1:B:182:GLN:HG3	2.21	0.40
1:A:157:ILE:HG12	1:A:159:TRP:CZ3	2.56	0.40
1:A:117:THR:HG21	1:A:122:LEU:CD2	2.51	0.40
1:A:131:VAL:HG12	1:A:143:VAL:HB	2.04	0.40
1:B:68:ARG:NH1	1:B:86:ARG:HG2	2.36	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

### 5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

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	$\# Z  > 2$
2	CYS	B	1318	-	3,6,6	1.43	0	1,7,7	1.10	0
2	CYS	A	1315	-	3,6,6	1.13	0	1,7,7	3.15	1 (100%)
2	CYS	B	1317	-	3,6,6	1.01	0	1,7,7	0.12	0
2	CYS	A	1316	-	3,6,6	0.80	0	1,7,7	0.26	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	CYS	B	1318	-	-	2/2/6/6	-
2	CYS	A	1315	-	-	0/2/6/6	-
2	CYS	B	1317	-	-	2/2/6/6	-
2	CYS	A	1316	-	-	1/2/6/6	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	A	1315	CYS	CA-CB-SG	3.15	121.22	114.44

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1318	CYS	N-CA-CB-SG
2	B	1318	CYS	C-CA-CB-SG
2	B	1317	CYS	N-CA-CB-SG
2	B	1317	CYS	C-CA-CB-SG
2	A	1316	CYS	N-CA-CB-SG

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1318	CYS	1	0
2	A	1315	CYS	6	0
2	B	1317	CYS	2	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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

EDS failed to run properly - this section is therefore empty.

### 6.4 Ligands ⓘ

EDS failed to run properly - this section is therefore empty.

### 6.5 Other polymers ⓘ

EDS failed to run properly - this section is therefore empty.