



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:05 AM UTC

PDB ID : 4JH8 / pdb_00004jh8
Title : Crystal Structure of FosB from Bacillus cereus with Zinc and L-Cysteine-Fosfomycin Ternary Complex
Authors : Thompson, M.K.; Harp, J.; Keithly, M.E.; Jagessar, K.; Cook, P.D.; Armstrong, R.N.
Deposited on : 2013-03-04
Resolution : 1.41 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

<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>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

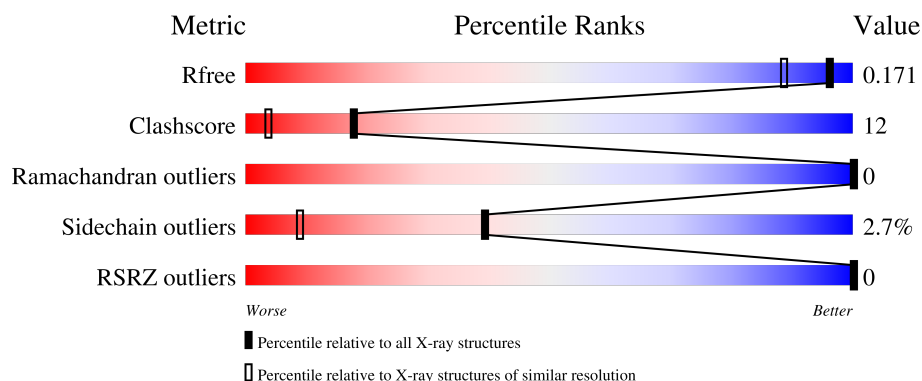
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.41 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}	180053	4041 (1.44-1.40)
Clashscore	190562	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)
RSRZ outliers	180081	4039 (1.44-1.40)

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 ≥ 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	138	 81% 17% ..
1	B	138	 80% 18% .

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
5	FCN	A	204	-	-	X	-
5	FCN	B	203	-	-	X	-
6	CYS	A	205	-	X	X	-
6	CYS	B	204	-	X	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2751 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 Metallothiol transferase FosB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	138	Total	C	N	O	S	0	8	0
			1216	775	203	232	6			
1	B	138	Total	C	N	O	S	0	11	0
			1228	789	204	229	6			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

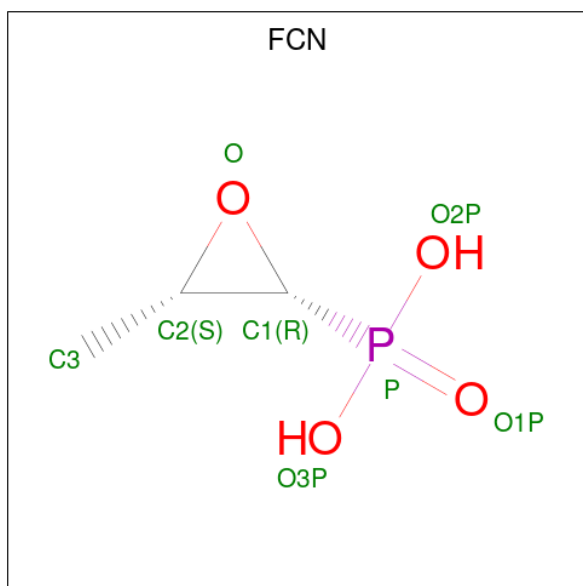
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



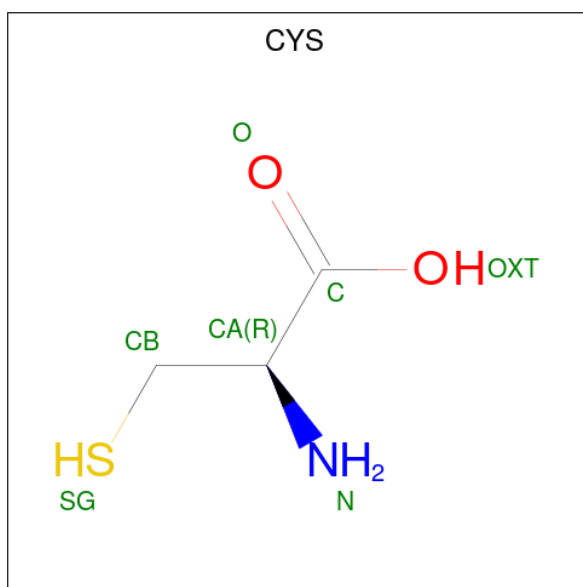
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is FOSFOMYCIN (CCD ID: FCN) (formula: $C_3H_7O_4P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	P	0	0
			8	3	4	1		
5	B	1	Total	C	O	P	0	0
			8	3	4	1		

- Molecule 6 is CYSTEINE (CCD ID: CYS) (formula: $C_3H_7NO_2S$).



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

- Molecule 7 is water.

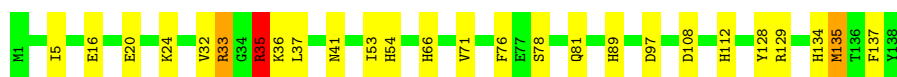
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	118	Total	O	0	2
			120	120		
7	B	132	Total	O	0	1
			133	133		

3 Residue-property plots

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: Metallothiol transferase FosB

Chain A:  81% 17% ..



- Molecule 1: Metallothiol transferase FosB

Chain B:  80% 18% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.31Å 68.53Å 69.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.92 – 1.41 48.92 – 1.41	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.92-1.41) 99.1 (48.92-1.41)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 1.41Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.137 , 0.174 0.134 , 0.171	Depositor DCC
R_{free} test set	3018 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.383	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,l,k	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2751	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.73% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, MG, FCN

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	1.46	10/1262 (0.8%)	1.18	4/1699 (0.2%)
1	B	1.38	4/1284 (0.3%)	1.19	4/1731 (0.2%)
All	All	1.42	14/2546 (0.5%)	1.18	8/3430 (0.2%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	112	HIS	ND1-CE1	7.29	1.39	1.32
1	B	134	HIS	CE1-NE2	7.20	1.39	1.32
1	A	54	HIS	N-CA	6.88	1.55	1.46
1	A	66	HIS	CE1-NE2	6.60	1.39	1.32
1	A	66	HIS	ND1-CE1	6.12	1.38	1.32
1	A	35	ARG	NE-CZ	6.02	1.39	1.33
1	A	54	HIS	CG-CD2	5.81	1.42	1.35
1	A	89	HIS	CE1-NE2	5.70	1.38	1.32
1	B	66	HIS	CE1-NE2	5.65	1.38	1.32
1	B	54	HIS	ND1-CE1	5.58	1.38	1.32
1	A	41	ASN	CG-OD1	-5.52	1.13	1.23
1	A	89	HIS	ND1-CE1	5.48	1.38	1.32
1	A	66	HIS	CA-C	-5.37	1.46	1.52
1	A	97	ASP	CG-OD1	-5.22	1.15	1.25

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	35	ARG	NE-CZ-NH2	8.37	126.74	119.20
1	B	129	ARG	N-CA-C	-7.89	102.67	111.82
1	B	92	LYS	CA-C-N	-6.85	107.98	121.41
1	B	92	LYS	C-N-CA	-6.85	107.98	121.41
1	A	53[A]	ILE	N-CA-C	6.61	118.21	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53[B]	ILE	N-CA-C	6.61	118.21	111.00
1	B	32	VAL	CB-CA-C	-5.42	101.28	111.30
1	A	32	VAL	CB-CA-C	-5.30	101.49	111.30

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	1216	0	1164	21	0
1	B	1228	0	1195	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	6	0	8	2	0
5	A	8	0	5	8	0
5	B	8	0	5	10	0
6	A	14	0	7	10	0
6	B	14	0	7	12	0
7	A	120	0	0	1	0
7	B	133	0	0	3	0
All	All	2751	0	2391	60	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 12.

All (60) 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)
5:B:203:FCN:O3P	6:B:204:CYS:HB2	1.35	1.22
5:B:203:FCN:P	6:B:204:CYS:HB2	1.82	1.18
1:A:16[A]:GLU:HG2	7:A:362:HOH:O	1.46	1.16
1:A:5:ILE:O	4:A:203:GOL:H12	1.43	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16[A]:GLU:OE1	1:A:33[A]:ARG:NH2	1.79	1.12
5:A:204:FCN:O2P	6:A:205:CYS:HB2	1.49	1.13
1:A:35:ARG:HH11	1:A:35:ARG:HG2	1.17	1.06
1:B:71[B]:VAL:HG11	1:B:79:LEU:CD1	1.90	1.01
1:B:41[A]:ASN:OD1	7:B:395:HOH:O	1.91	0.87
1:B:71[B]:VAL:HG11	1:B:79:LEU:HD11	1.56	0.87
1:B:20:GLU:OE2	1:B:24:LYS:HD2	1.76	0.85
5:B:203:FCN:P	6:B:204:CYS:CB	2.66	0.82
5:B:203:FCN:O3P	6:B:204:CYS:CB	2.26	0.82
5:A:204:FCN:P	6:A:205:CYS:HB2	2.19	0.81
5:A:204:FCN:C1	6:A:205:CYS:SG	2.70	0.79
6:B:204:CYS:N	7:B:425:HOH:O	2.14	0.79
5:A:204:FCN:C2	6:A:205:CYS:SG	2.71	0.78
1:B:71[B]:VAL:CG1	1:B:79:LEU:HD11	2.16	0.75
1:A:35:ARG:HD3	1:A:36:LYS:HG2	1.69	0.74
1:A:16[A]:GLU:CD	1:A:33[A]:ARG:HH22	1.93	0.71
6:A:205:CYS:HA	1:B:50:ASN:HD21	1.56	0.69
6:A:205:CYS:HA	1:B:50:ASN:ND2	2.08	0.68
1:A:35:ARG:HH11	1:A:35:ARG:CG	1.98	0.67
1:A:35:ARG:HG2	1:A:35:ARG:NH1	1.97	0.66
1:A:134:HIS:O	1:A:135[B]:MET:HG2	1.96	0.65
5:B:203:FCN:C2	6:B:204:CYS:SG	2.84	0.65
1:B:83:LEU:HB3	1:B:90[B]:ILE:HD11	1.79	0.65
1:B:16:GLU:OE2	1:B:33[B]:ARG:NH2	2.31	0.64
5:B:203:FCN:C1	6:B:204:CYS:SG	2.88	0.62
5:B:203:FCN:H1	6:B:204:CYS:HB2	1.86	0.57
1:A:134:HIS:C	1:A:135[B]:MET:HG2	2.30	0.56
5:A:204:FCN:P	6:A:205:CYS:CB	2.91	0.56
5:B:203:FCN:C1	6:B:204:CYS:HB2	2.36	0.56
1:A:128:TYR:HD2	1:A:135[B]:MET:HE1	1.71	0.55
1:A:78:SER:HA	1:A:81:GLN:HE21	1.71	0.55
1:A:128:TYR:CD2	1:A:135[B]:MET:HE1	2.42	0.54
1:A:20[B]:GLU:OE2	1:A:24:LYS:HD2	2.08	0.52
1:A:71:VAL:HG23	1:A:76:PHE:CE1	2.45	0.52
1:A:71:VAL:CG2	1:A:76:PHE:CE1	2.93	0.52
5:A:204:FCN:P	6:A:205:CYS:SG	3.08	0.51
1:B:20:GLU:CD	1:B:24:LYS:HD2	2.36	0.51
1:B:37:LEU:C	1:B:37:LEU:HD12	2.36	0.50
5:A:204:FCN:H2	6:A:205:CYS:SG	2.51	0.50
1:B:108:ASP:OD2	1:B:112:HIS:HD2	1.95	0.49
1:B:71[B]:VAL:CG1	1:B:79:LEU:CD1	2.75	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:ASP:OD2	1:A:112:HIS:HD2	1.99	0.46
1:A:35:ARG:CD	1:A:36:LYS:HG2	2.41	0.46
4:A:203:GOL:H32	1:B:70:SER:OG	2.16	0.46
1:B:129:ARG:HA	1:B:137:PHE:CZ	2.51	0.46
1:A:129:ARG:HA	1:A:137:PHE:CZ	2.53	0.43
1:B:24:LYS:HD3	1:B:86:ASN:HD21	1.83	0.43
1:B:71[B]:VAL:HG11	1:B:79:LEU:HD12	1.88	0.43
1:B:107[B]:VAL:CG1	1:B:111:GLY:HA2	2.49	0.42
1:A:135[B]:MET:HE1	1:B:32:VAL:HG11	2.00	0.42
5:B:203:FCN:C1	6:B:204:CYS:CB	2.97	0.42
5:B:203:FCN:H1	6:B:204:CYS:CB	2.49	0.42
5:A:204:FCN:O2P	6:A:205:CYS:CB	2.42	0.41
6:B:204:CYS:O	7:B:431:HOH:O	2.22	0.41
1:A:71:VAL:HG23	1:A:76:PHE:HE1	1.84	0.41
1:B:132:LYS:HD3	1:B:135[A]:MET:HE3	2.03	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	144/138 (104%)	141 (98%)	3 (2%)	0	100	100
1	B	147/138 (106%)	145 (99%)	2 (1%)	0	100	100
All	All	291/276 (105%)	286 (98%)	5 (2%)	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	136/128 (106%)	130 (96%)	6 (4%)	25 3
1	B	139/128 (109%)	137 (99%)	2 (1%)	59 27
All	All	275/256 (107%)	267 (97%)	8 (3%)	39 8

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33[A]	ARG
1	A	33[B]	ARG
1	A	35	ARG
1	A	37	LEU
1	A	135[A]	MET
1	A	135[B]	MET
1	B	43	CYS
1	B	58	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	112	HIS
1	B	50	ASN
1	B	86	ASN
1	B	89	HIS
1	B	112	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 for Mogul analysis.

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
4	GOL	A	203	-	5,5,5	0.97	0	5,5,5	1.60	1 (20%)
5	FCN	B	203	2	5,8,8	0.43	0	5,13,13	2.75	1 (20%)
6	CYS	A	206	1	5,6,6	1.27	1 (20%)	3,7,7	3.42	2 (66%)
6	CYS	B	204	-	5,6,6	1.79	2 (40%)	3,7,7	3.46	2 (66%)
5	FCN	A	204	2	5,8,8	2.00	3 (60%)	5,13,13	4.27	3 (60%)
6	CYS	B	202	1	5,6,6	1.59	1 (20%)	3,7,7	1.09	0
6	CYS	A	205	-	5,6,6	1.37	1 (20%)	3,7,7	2.99	3 (100%)

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
4	GOL	A	203	-	-	4/4/4/4	-
5	FCN	B	203	2	-	0/0/11/11	0/1/1/1
6	CYS	A	206	1	-	3/6/6/6	-
6	CYS	B	204	-	-	5/6/6/6	-
5	FCN	A	204	2	-	0/0/11/11	0/1/1/1
6	CYS	B	202	1	-	2/6/6/6	-
6	CYS	A	205	-	-	5/6/6/6	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	204	FCN	P-O2P	3.10	1.59	1.54
6	B	204	CYS	CB-CA	-2.87	1.49	1.53
6	A	205	CYS	CB-CA	-2.73	1.50	1.53
6	B	204	CYS	OXT-C	-2.56	1.22	1.30
6	B	202	CYS	OXT-C	-2.56	1.22	1.30
5	A	204	FCN	P-O3P	2.40	1.58	1.54
6	A	206	CYS	O-C	2.14	1.28	1.22
5	A	204	FCN	P-O1P	2.12	1.53	1.49

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	204	FCN	O1P-P-C1	-7.66	96.14	113.37
5	B	203	FCN	O-C2-C3	-5.75	105.97	116.61
5	A	204	FCN	O-C2-C3	-4.93	107.48	116.61
6	B	204	CYS	CB-CA-C	-4.66	105.25	109.89
6	A	206	CYS	OXT-C-O	-4.59	113.66	124.08
6	A	205	CYS	CB-CA-C	-4.31	105.60	109.89
6	A	206	CYS	CB-CA-C	3.69	113.56	109.89
6	B	204	CYS	OXT-C-O	-3.45	116.26	124.08
4	A	203	GOL	O1-C1-C2	2.81	123.01	110.38
5	A	204	FCN	O3P-P-O1P	-2.13	108.12	113.45
6	A	205	CYS	CA-CB-SG	2.04	118.75	114.40
6	A	205	CYS	OXT-C-O	-2.03	119.48	124.08

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	203	GOL	C1-C2-C3-O3
6	A	205	CYS	OXT-C-CA-CB
6	A	205	CYS	N-CA-CB-SG
6	A	206	CYS	N-CA-CB-SG
6	A	206	CYS	C-CA-CB-SG
6	B	202	CYS	C-CA-CB-SG
6	B	204	CYS	O-C-CA-CB
6	B	204	CYS	OXT-C-CA-CB
6	B	204	CYS	N-CA-CB-SG
6	B	204	CYS	C-CA-CB-SG
6	A	205	CYS	OXT-C-CA-N
4	A	203	GOL	O1-C1-C2-C3
4	A	203	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
6	A	205	CYS	O-C-CA-CB
6	A	205	CYS	O-C-CA-N
4	A	203	GOL	O1-C1-C2-O2
6	B	204	CYS	O-C-CA-N
6	B	202	CYS	N-CA-CB-SG
6	A	206	CYS	O-C-CA-N

There are no ring outliers.

5 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	203	GOL	2	0
5	B	203	FCN	10	0
6	B	204	CYS	12	0
5	A	204	FCN	8	0
6	A	205	CYS	10	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	138/138 (100%)	-0.35	0 100 100	9, 17, 38, 52	8 (5%)
1	B	138/138 (100%)	-0.61	0 100 100	6, 14, 27, 35	11 (7%)
All	All	276/276 (100%)	-0.48	0 100 100	6, 15, 34, 52	19 (6%)

There are no RSRZ outliers to report.

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 oligosaccharides 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, 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	B-factors(Å ²)	Q<0.9
6	CYS	A	206	7/7	0.88	0.15	43,62,75,76	0
6	CYS	A	205	7/7	0.89	0.14	26,41,60,62	0
4	GOL	A	203	6/6	0.93	0.10	21,31,45,58	0
6	CYS	B	202	7/7	0.96	0.11	28,31,41,43	0
6	CYS	B	204	7/7	0.96	0.10	27,40,47,60	0
3	MG	B	205	1/1	0.98	0.13	39,39,39,39	0
5	FCN	A	204	8/8	0.98	0.06	16,20,24,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	202	1/1	0.99	0.15	25,25,25,25	0
5	FCN	B	203	8/8	0.99	0.05	11,15,21,22	0
2	ZN	B	201	1/1	1.00	0.01	10,10,10,10	0
2	ZN	A	201	1/1	1.00	0.01	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.