



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 16, 2023 – 04:17 PM JST

PDB ID : 8J5M
Title : Structure of GH1 Br2 beta-glucosidase E350G mutant from bovine rumen metagenome
Authors : Kaenyng, W.; Kongsaree, P.T.; Tagami, T.
Deposited on : 2023-04-23
Resolution : 1.62 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

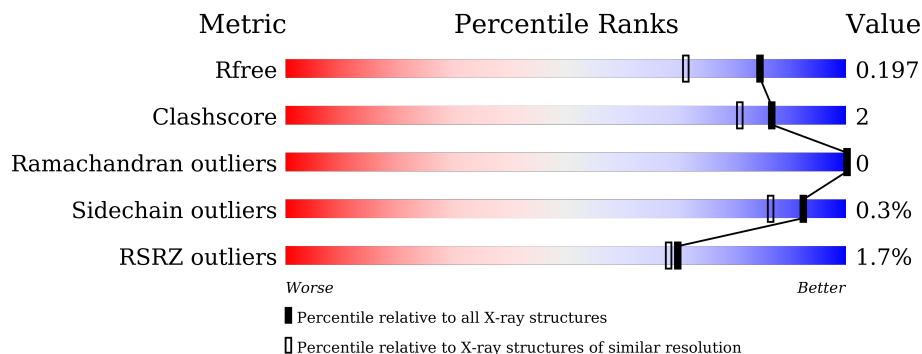
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.62 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}	130704	4693 (1.64-1.60)
Clashscore	141614	5002 (1.64-1.60)
Ramachandran outliers	138981	4888 (1.64-1.60)
Sidechain outliers	138945	4887 (1.64-1.60)
RSRZ outliers	127900	4609 (1.64-1.60)

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	465	
1	B	465	
1	C	465	
1	D	465	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	501	-	-	X	-
2	ACT	B	501	-	-	X	-
2	ACT	C	501	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16812 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 Beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	452	3652	2355	611	670	16	0	1	0
1	B	452	3658	2357	612	673	16	0	2	0
1	D	444	3594	2319	599	661	15	0	1	0
1	C	444	3600	2322	600	663	15	0	2	0

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A1S5SJM8
A	-18	GLY	-	expression tag	UNP A0A1S5SJM8
A	-17	SER	-	expression tag	UNP A0A1S5SJM8
A	-16	SER	-	expression tag	UNP A0A1S5SJM8
A	-15	HIS	-	expression tag	UNP A0A1S5SJM8
A	-14	HIS	-	expression tag	UNP A0A1S5SJM8
A	-13	HIS	-	expression tag	UNP A0A1S5SJM8
A	-12	HIS	-	expression tag	UNP A0A1S5SJM8
A	-11	HIS	-	expression tag	UNP A0A1S5SJM8
A	-10	HIS	-	expression tag	UNP A0A1S5SJM8
A	-9	SER	-	expression tag	UNP A0A1S5SJM8
A	-8	SER	-	expression tag	UNP A0A1S5SJM8
A	-7	GLY	-	expression tag	UNP A0A1S5SJM8
A	-6	LEU	-	expression tag	UNP A0A1S5SJM8
A	-5	VAL	-	expression tag	UNP A0A1S5SJM8
A	-4	PRO	-	expression tag	UNP A0A1S5SJM8
A	-3	ARG	-	expression tag	UNP A0A1S5SJM8
A	-2	GLY	-	expression tag	UNP A0A1S5SJM8
A	-1	SER	-	expression tag	UNP A0A1S5SJM8
A	0	HIS	-	expression tag	UNP A0A1S5SJM8
A	350	GLY	GLU	engineered mutation	UNP A0A1S5SJM8

Continued on next page...

Continued from previous page...

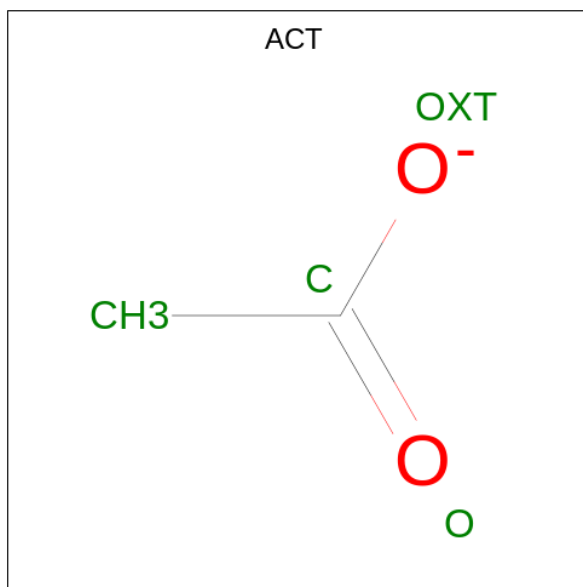
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP A0A1S5SJM8
B	-18	GLY	-	expression tag	UNP A0A1S5SJM8
B	-17	SER	-	expression tag	UNP A0A1S5SJM8
B	-16	SER	-	expression tag	UNP A0A1S5SJM8
B	-15	HIS	-	expression tag	UNP A0A1S5SJM8
B	-14	HIS	-	expression tag	UNP A0A1S5SJM8
B	-13	HIS	-	expression tag	UNP A0A1S5SJM8
B	-12	HIS	-	expression tag	UNP A0A1S5SJM8
B	-11	HIS	-	expression tag	UNP A0A1S5SJM8
B	-10	HIS	-	expression tag	UNP A0A1S5SJM8
B	-9	SER	-	expression tag	UNP A0A1S5SJM8
B	-8	SER	-	expression tag	UNP A0A1S5SJM8
B	-7	GLY	-	expression tag	UNP A0A1S5SJM8
B	-6	LEU	-	expression tag	UNP A0A1S5SJM8
B	-5	VAL	-	expression tag	UNP A0A1S5SJM8
B	-4	PRO	-	expression tag	UNP A0A1S5SJM8
B	-3	ARG	-	expression tag	UNP A0A1S5SJM8
B	-2	GLY	-	expression tag	UNP A0A1S5SJM8
B	-1	SER	-	expression tag	UNP A0A1S5SJM8
B	0	HIS	-	expression tag	UNP A0A1S5SJM8
B	350	GLY	GLU	engineered mutation	UNP A0A1S5SJM8
D	-19	MET	-	initiating methionine	UNP A0A1S5SJM8
D	-18	GLY	-	expression tag	UNP A0A1S5SJM8
D	-17	SER	-	expression tag	UNP A0A1S5SJM8
D	-16	SER	-	expression tag	UNP A0A1S5SJM8
D	-15	HIS	-	expression tag	UNP A0A1S5SJM8
D	-14	HIS	-	expression tag	UNP A0A1S5SJM8
D	-13	HIS	-	expression tag	UNP A0A1S5SJM8
D	-12	HIS	-	expression tag	UNP A0A1S5SJM8
D	-11	HIS	-	expression tag	UNP A0A1S5SJM8
D	-10	HIS	-	expression tag	UNP A0A1S5SJM8
D	-9	SER	-	expression tag	UNP A0A1S5SJM8
D	-8	SER	-	expression tag	UNP A0A1S5SJM8
D	-7	GLY	-	expression tag	UNP A0A1S5SJM8
D	-6	LEU	-	expression tag	UNP A0A1S5SJM8
D	-5	VAL	-	expression tag	UNP A0A1S5SJM8
D	-4	PRO	-	expression tag	UNP A0A1S5SJM8
D	-3	ARG	-	expression tag	UNP A0A1S5SJM8
D	-2	GLY	-	expression tag	UNP A0A1S5SJM8
D	-1	SER	-	expression tag	UNP A0A1S5SJM8
D	0	HIS	-	expression tag	UNP A0A1S5SJM8
D	350	GLY	GLU	engineered mutation	UNP A0A1S5SJM8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	MET	-	initiating methionine	UNP A0A1S5SJM8
C	-18	GLY	-	expression tag	UNP A0A1S5SJM8
C	-17	SER	-	expression tag	UNP A0A1S5SJM8
C	-16	SER	-	expression tag	UNP A0A1S5SJM8
C	-15	HIS	-	expression tag	UNP A0A1S5SJM8
C	-14	HIS	-	expression tag	UNP A0A1S5SJM8
C	-13	HIS	-	expression tag	UNP A0A1S5SJM8
C	-12	HIS	-	expression tag	UNP A0A1S5SJM8
C	-11	HIS	-	expression tag	UNP A0A1S5SJM8
C	-10	HIS	-	expression tag	UNP A0A1S5SJM8
C	-9	SER	-	expression tag	UNP A0A1S5SJM8
C	-8	SER	-	expression tag	UNP A0A1S5SJM8
C	-7	GLY	-	expression tag	UNP A0A1S5SJM8
C	-6	LEU	-	expression tag	UNP A0A1S5SJM8
C	-5	VAL	-	expression tag	UNP A0A1S5SJM8
C	-4	PRO	-	expression tag	UNP A0A1S5SJM8
C	-3	ARG	-	expression tag	UNP A0A1S5SJM8
C	-2	GLY	-	expression tag	UNP A0A1S5SJM8
C	-1	SER	-	expression tag	UNP A0A1S5SJM8
C	0	HIS	-	expression tag	UNP A0A1S5SJM8
C	350	GLY	GLU	engineered mutation	UNP A0A1S5SJM8

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



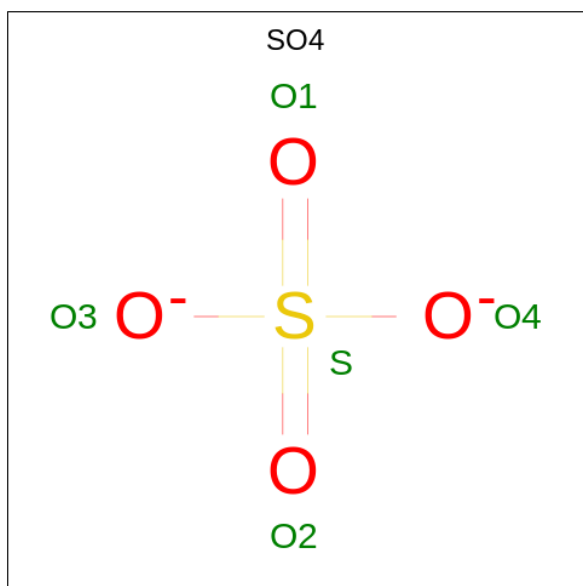
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

Continued on next page...

Continued from previous page...

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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

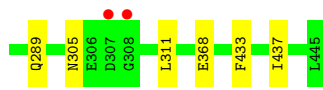
Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	532	Total O 532 532	0	0
4	B	540	Total O 540 540	0	0
4	D	510	Total O 510 510	0	0
4	C	525	Total O 525 525	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.30Å 113.61Å 180.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.24 – 1.62 42.24 – 1.62	Depositor EDS
% Data completeness (in resolution range)	99.2 (42.24-1.62) 99.2 (42.24-1.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 1.62Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.161 , 0.188 0.172 , 0.197	Depositor DCC
R_{free} test set	2001 reflections (0.74%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtrriage
Anisotropy	0.272	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	16812	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.20% 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: ACT, 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	0.63	0/3760	0.74	0/5098
1	B	0.61	0/3766	0.74	0/5107
1	C	0.61	0/3706	0.73	0/5025
1	D	0.63	0/3700	0.74	0/5017
All	All	0.62	0/14932	0.74	0/20247

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3652	0	3508	16	0
1	B	3658	0	3504	13	0
1	C	3600	0	3448	18	0
1	D	3594	0	3444	8	0
2	A	8	0	6	3	0
2	B	16	0	12	4	0
2	C	8	0	6	3	0
2	D	4	0	3	0	0
3	A	40	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	35	0	0	0	0
3	C	35	0	0	0	0
3	D	55	0	0	0	0
4	A	532	0	0	9	0
4	B	540	0	0	5	0
4	C	525	0	0	6	0
4	D	510	0	0	2	0
All	All	16812	0	13931	61	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 2.

The worst 5 of 61 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:267:TRP:CZ3	1:C:271:ILE:HD11	2.05	0.92
1:C:267:TRP:HZ3	1:C:271:ILE:HD11	1.36	0.90
2:C:501:ACT:H3	4:C:604:HOH:O	1.73	0.88
2:B:501:ACT:H1	4:B:622:HOH:O	1.83	0.77
4:A:788:HOH:O	1:C:368:GLU:HG3	1.85	0.75

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	451/465 (97%)	440 (98%)	11 (2%)	0	100	100
1	B	452/465 (97%)	441 (98%)	11 (2%)	0	100	100
1	C	444/465 (96%)	432 (97%)	12 (3%)	0	100	100
1	D	443/465 (95%)	433 (98%)	10 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1790/1860 (96%)	1746 (98%)	44 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	385/395 (98%)	384 (100%)	1 (0%)	92	86
1	B	385/395 (98%)	385 (100%)	0	100	100
1	C	379/395 (96%)	378 (100%)	1 (0%)	92	86
1	D	378/395 (96%)	375 (99%)	3 (1%)	81	69
All	All	1527/1580 (97%)	1522 (100%)	5 (0%)	92	86

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	347	MET
1	D	5	LYS
1	D	225	PRO
1	D	306	GLU
1	C	225	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	238	ASN
1	D	435	GLN
1	C	407	ASN
1	A	118	HIS
1	A	17	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

42 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	C	503	-	4,4,4	0.45	0	6,6,6	0.05	0
3	SO4	C	507	-	4,4,4	0.24	0	6,6,6	0.14	0
2	ACT	B	509	-	3,3,3	0.92	0	3,3,3	1.00	0
3	SO4	B	508	-	4,4,4	0.42	0	6,6,6	0.13	0
3	SO4	C	504	-	4,4,4	0.36	0	6,6,6	0.13	0
3	SO4	A	509	-	4,4,4	0.37	0	6,6,6	0.07	0
2	ACT	B	511	-	3,3,3	0.98	0	3,3,3	0.75	0
3	SO4	B	505	-	4,4,4	0.34	0	6,6,6	0.09	0
3	SO4	D	508	-	4,4,4	0.37	0	6,6,6	0.08	0
3	SO4	B	507	-	4,4,4	0.23	0	6,6,6	0.12	0
3	SO4	A	503	-	4,4,4	0.35	0	6,6,6	0.07	0
2	ACT	B	501	-	3,3,3	0.66	0	3,3,3	0.86	0
3	SO4	D	505	-	4,4,4	0.38	0	6,6,6	0.10	0
2	ACT	A	501	-	3,3,3	0.86	0	3,3,3	0.81	0
3	SO4	B	506	-	4,4,4	0.20	0	6,6,6	0.14	0
3	SO4	D	504	-	4,4,4	0.38	0	6,6,6	0.08	0
3	SO4	D	506	-	4,4,4	0.36	0	6,6,6	0.07	0
3	SO4	C	502	-	4,4,4	0.34	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	504	-	4,4,4	0.35	0	6,6,6	0.05	0
3	SO4	A	506	-	4,4,4	0.34	0	6,6,6	0.08	0
2	ACT	A	510	-	3,3,3	1.04	0	3,3,3	0.76	0
3	SO4	A	505	-	4,4,4	0.40	0	6,6,6	0.12	0
3	SO4	D	511	-	4,4,4	0.24	0	6,6,6	0.09	0
2	ACT	D	501	-	3,3,3	0.85	0	3,3,3	0.90	0
3	SO4	A	507	-	4,4,4	0.36	0	6,6,6	0.11	0
3	SO4	D	510	-	4,4,4	0.37	0	6,6,6	0.08	0
2	ACT	C	509	-	3,3,3	0.99	0	3,3,3	0.98	0
3	SO4	A	508	-	4,4,4	0.18	0	6,6,6	0.13	0
3	SO4	D	512	-	4,4,4	0.37	0	6,6,6	0.09	0
3	SO4	D	503	-	4,4,4	0.37	0	6,6,6	0.10	0
2	ACT	B	510	-	3,3,3	0.96	0	3,3,3	0.78	0
3	SO4	D	509	-	4,4,4	0.37	0	6,6,6	0.12	0
3	SO4	D	502	-	4,4,4	0.39	0	6,6,6	0.07	0
3	SO4	B	502	-	4,4,4	0.36	0	6,6,6	0.12	0
3	SO4	A	502	-	4,4,4	0.34	0	6,6,6	0.10	0
3	SO4	C	508	-	4,4,4	0.22	0	6,6,6	0.10	0
3	SO4	D	507	-	4,4,4	0.35	0	6,6,6	0.09	0
3	SO4	B	503	-	4,4,4	0.41	0	6,6,6	0.15	0
3	SO4	B	504	-	4,4,4	0.41	0	6,6,6	0.09	0
3	SO4	C	506	-	4,4,4	0.35	0	6,6,6	0.17	0
2	ACT	C	501	-	3,3,3	1.00	0	3,3,3	0.93	0
3	SO4	C	505	-	4,4,4	0.36	0	6,6,6	0.10	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	ACT	4	0
2	A	501	ACT	3	0
2	C	501	ACT	3	0

5.7 Other polymers [\(i\)](#)

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 [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	452/465 (97%)	-0.14	7 (1%) 73 72	15, 20, 36, 56	0
1	B	452/465 (97%)	-0.12	10 (2%) 62 60	15, 21, 36, 61	0
1	C	444/465 (95%)	-0.22	7 (1%) 72 71	15, 21, 37, 58	0
1	D	444/465 (95%)	-0.17	6 (1%) 75 74	15, 23, 38, 53	0
All	All	1792/1860 (96%)	-0.16	30 (1%) 70 68	15, 21, 37, 61	0

The worst 5 of 30 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	270	TYR	4.3
1	B	310	ILE	4.2
1	A	310	ILE	3.8
1	A	312	GLN	3.7
1	C	308	GLY	3.4

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 monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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(\AA^2)	Q<0.9
3	SO4	A	506	5/5	0.56	0.25	73,76,84,85	0
3	SO4	B	505	5/5	0.69	0.21	69,72,81,81	0
3	SO4	D	507	5/5	0.76	0.26	69,78,85,87	0
3	SO4	D	506	5/5	0.81	0.22	82,84,91,92	0
3	SO4	C	508	5/5	0.84	0.21	41,42,45,53	0
3	SO4	C	505	5/5	0.87	0.17	65,71,78,79	0
2	ACT	B	510	4/4	0.87	0.14	59,59,60,60	0
3	SO4	A	504	5/5	0.88	0.20	81,87,88,91	0
3	SO4	D	511	5/5	0.88	0.23	44,50,52,59	0
3	SO4	D	510	5/5	0.89	0.34	72,80,85,85	0
3	SO4	A	502	5/5	0.89	0.24	52,61,68,69	0
3	SO4	A	508	5/5	0.90	0.18	39,41,46,51	0
3	SO4	A	507	5/5	0.90	0.20	63,63,66,66	0
2	ACT	B	509	4/4	0.91	0.20	47,52,52,53	0
3	SO4	D	503	5/5	0.91	0.30	74,75,78,78	0
3	SO4	C	504	5/5	0.93	0.17	55,56,60,60	0
3	SO4	B	506	5/5	0.93	0.19	39,45,48,53	0
3	SO4	D	508	5/5	0.93	0.20	54,55,58,65	0
2	ACT	A	510	4/4	0.94	0.15	34,38,41,44	0
3	SO4	B	502	5/5	0.94	0.24	53,57,62,69	0
3	SO4	B	507	5/5	0.94	0.19	42,48,49,53	0
2	ACT	B	511	4/4	0.95	0.12	43,44,45,48	0
2	ACT	C	509	4/4	0.95	0.10	29,32,37,39	0
3	SO4	A	505	5/5	0.96	0.16	50,53,56,62	0
3	SO4	C	502	5/5	0.96	0.28	48,49,55,56	0
2	ACT	D	501	4/4	0.96	0.10	28,29,32,38	0
2	ACT	A	501	4/4	0.96	0.13	26,28,29,38	0
3	SO4	D	504	5/5	0.96	0.17	54,54,57,59	0
3	SO4	A	503	5/5	0.97	0.11	59,59,60,62	0
3	SO4	D	509	5/5	0.97	0.11	41,45,47,47	0
3	SO4	B	508	5/5	0.97	0.18	55,55,60,63	0
3	SO4	D	502	5/5	0.97	0.19	60,63,64,67	0
3	SO4	D	512	5/5	0.97	0.16	36,38,42,43	0
3	SO4	B	503	5/5	0.97	0.16	45,45,52,54	0
3	SO4	B	504	5/5	0.97	0.17	61,62,65,66	0
2	ACT	B	501	4/4	0.97	0.13	26,29,31,35	0
2	ACT	C	501	4/4	0.97	0.11	26,32,37,42	0
3	SO4	A	509	5/5	0.98	0.15	51,53,60,61	0
3	SO4	D	505	5/5	0.98	0.16	43,43,49,50	0
3	SO4	C	507	5/5	0.98	0.12	33,34,35,36	0
3	SO4	C	503	5/5	0.98	0.12	55,60,61,63	0
3	SO4	C	506	5/5	0.99	0.09	38,43,45,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.