



Full wwPDB X-ray Structure Validation Report

Mar 31, 2014 – 01:39 PM BST

PDB ID : 4MQO
Title : Mycobacterium tuberculosis transaminase BioA complexed with 5,6-DIHYDR
O-BENZO[H]CINNOLIN-3-YLAMINE
Authors : Finzel, B.C; Dai, R; Geders, T.W
Deposited on : 2013-09-16
Resolution : 1.70 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

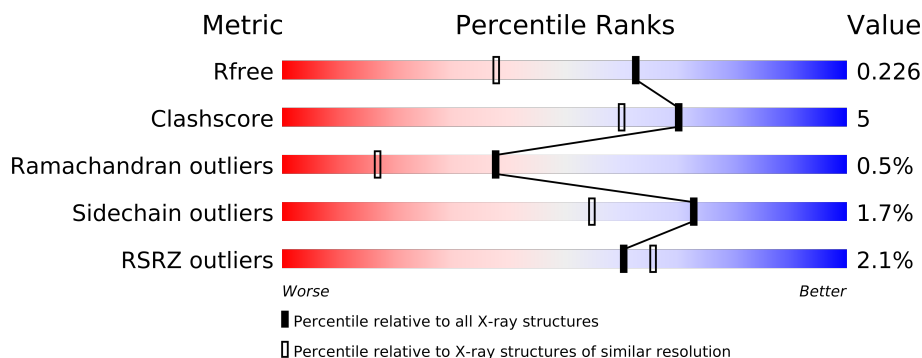
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable23004
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) : stable23004

1 Overall quality at a glance

The reported resolution of this entry is 1.70 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	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	457	
1	B	457	

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	2BG	A	501	X	X
4	SO4	A	503	-	X
5	EDO	A	504	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6756 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 Adenosylmethionine-8-amino-7-oxononanoateaminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	1	10	0
			3229	2056	570	581	22			
1	B	427	Total	C	N	O	S	4	7	0
			3239	2064	570	582	23			

There are 40 discrepancies between the modelled and reference sequences:

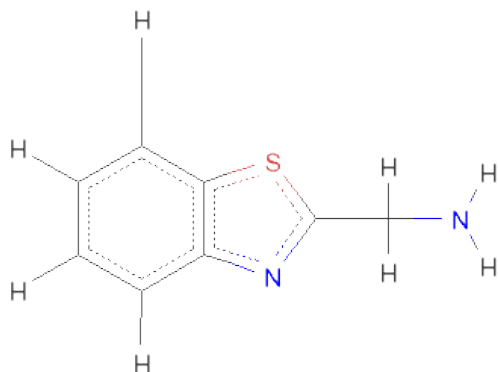
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	INITIATING METHIONINE	UNP P0A4X6
A	-18	GLY	-	EXPRESSION TAG	UNP P0A4X6
A	-17	SER	-	EXPRESSION TAG	UNP P0A4X6
A	-16	SER	-	EXPRESSION TAG	UNP P0A4X6
A	-15	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-14	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-13	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-12	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-11	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-10	HIS	-	EXPRESSION TAG	UNP P0A4X6
A	-9	SER	-	EXPRESSION TAG	UNP P0A4X6
A	-8	SER	-	EXPRESSION TAG	UNP P0A4X6
A	-7	GLY	-	EXPRESSION TAG	UNP P0A4X6
A	-6	LEU	-	EXPRESSION TAG	UNP P0A4X6
A	-5	VAL	-	EXPRESSION TAG	UNP P0A4X6
A	-4	PRO	-	EXPRESSION TAG	UNP P0A4X6
A	-3	ARG	-	EXPRESSION TAG	UNP P0A4X6
A	-2	GLY	-	EXPRESSION TAG	UNP P0A4X6
A	-1	SER	-	EXPRESSION TAG	UNP P0A4X6
A	0	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-19	MET	-	INITIATING METHIONINE	UNP P0A4X6
B	-18	GLY	-	EXPRESSION TAG	UNP P0A4X6
B	-17	SER	-	EXPRESSION TAG	UNP P0A4X6
B	-16	SER	-	EXPRESSION TAG	UNP P0A4X6
B	-15	HIS	-	EXPRESSION TAG	UNP P0A4X6

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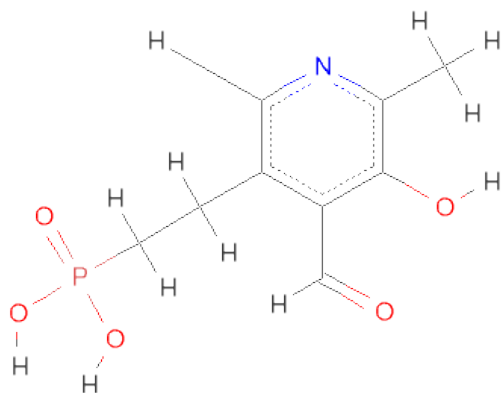
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-13	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-12	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-11	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-10	HIS	-	EXPRESSION TAG	UNP P0A4X6
B	-9	SER	-	EXPRESSION TAG	UNP P0A4X6
B	-8	SER	-	EXPRESSION TAG	UNP P0A4X6
B	-7	GLY	-	EXPRESSION TAG	UNP P0A4X6
B	-6	LEU	-	EXPRESSION TAG	UNP P0A4X6
B	-5	VAL	-	EXPRESSION TAG	UNP P0A4X6
B	-4	PRO	-	EXPRESSION TAG	UNP P0A4X6
B	-3	ARG	-	EXPRESSION TAG	UNP P0A4X6
B	-2	GLY	-	EXPRESSION TAG	UNP P0A4X6
B	-1	SER	-	EXPRESSION TAG	UNP P0A4X6
B	0	HIS	-	EXPRESSION TAG	UNP P0A4X6

- Molecule 2 is 1-(1,3-BENZOTHAZOL-2-YL)METHANAMINE (three-letter code: 2BG) (formula: C₈H₈N₂S).



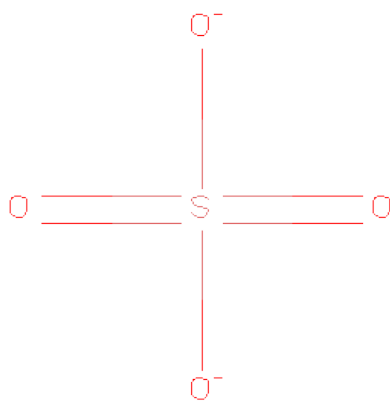
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	S	1	0
			11	8	2	1		

- Molecule 3 is [2-(4-FORMYL-5-HYDROXY-6-METHYLPYRIDIN-3-YL)ETHYL]PHOSPHONICACID (three-letter code: 2B0) (formula: C₉H₁₂NO₅P).



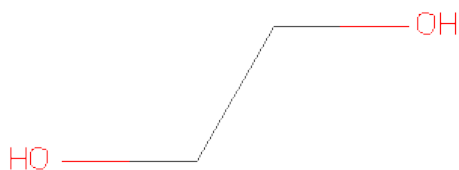
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	9	1	4	1		
3	B	1	Total	C	N	O	P	0	0
			16	9	1	5	1		

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	128	Total	O	0	0
			128	128		
6	B	104	Total	O	0	0
			104	104		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.94Å 66.08Å 201.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	97.48 – 1.70 100.95 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.5 (97.48-1.70) 97.5 (100.95-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.69 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.197 , 0.226 0.197 , 0.226	Depositor DCC
R_{free} test set	4559 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 31.9	EDS
Estimated twinning fraction	0.038 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 91186 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6756	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: 2BG, 2B0, EDO, 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	0.74	4/3339 (0.1%)	0.80	8/4559 (0.2%)
1	B	0.59	4/3337 (0.1%)	0.65	0/4562
All	All	0.67	8/6676 (0.1%)	0.73	8/9121 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	403[A]	ARG	CZ-NH2	14.35	1.51	1.33
1	A	403[B]	ARG	CZ-NH2	14.35	1.51	1.33
1	A	403[A]	ARG	NE-CZ	-13.28	1.15	1.33
1	A	403[B]	ARG	NE-CZ	-13.28	1.15	1.33
1	B	22	TRP	CD2-CE2	5.37	1.47	1.41
1	B	398	TRP	CD2-CE2	5.31	1.47	1.41
1	B	139	TRP	CD2-CE2	5.29	1.47	1.41
1	B	65	TRP	CD2-CE2	5.21	1.47	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	403[A]	ARG	NE-CZ-NH2	-15.64	112.48	120.30
1	A	403[B]	ARG	NE-CZ-NH2	-15.64	112.48	120.30
1	A	403[A]	ARG	CD-NE-CZ	11.79	140.10	123.60
1	A	403[B]	ARG	CD-NE-CZ	11.79	140.10	123.60
1	A	403[A]	ARG	NH1-CZ-NH2	9.97	130.36	119.40
1	A	403[B]	ARG	NH1-CZ-NH2	9.97	130.36	119.40
1	A	403[A]	ARG	NE-CZ-NH1	-6.99	116.80	120.30
1	A	403[B]	ARG	NE-CZ-NH1	-6.99	116.80	120.30

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	3229	0	3225	24	0
1	B	3239	0	3217	43	0
2	A	11	0	8	0	0
3	A	15	0	10	1	0
3	B	16	0	12	8	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	4	0	6	0	0
6	A	128	0	0	1	0
6	B	104	0	0	4	0
All	All	6756	0	6478	59	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 5.

All (59) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:283:LYS:HZ2	3:B:501:2B0:C1	1.40	1.31
1:B:283:LYS:NZ	3:B:501:2B0:O2	1.69	1.23
1:B:283:LYS:HZ1	3:B:501:2B0:C1	1.77	0.98
1:B:124:GLY:HA3	3:B:501:2B0:H4	1.49	0.92
1:A:314:MET:CE	1:B:25[A]:TYR:HD1	1.90	0.84
1:A:314:MET:HE3	1:B:25[A]:TYR:HD1	1.55	0.69
1:B:178:TRP:N	6:B:703:HOH:O	2.25	0.69
1:B:178:TRP:CD1	6:B:703:HOH:O	2.47	0.68
3:A:502:2B0:C1	3:A:502:2B0:H4	2.22	0.67
1:B:283:LYS:CE	3:B:501:2B0:O2	2.44	0.66
1:A:314:MET:HE3	1:B:25[A]:TYR:CD1	2.30	0.66
1:B:283:LYS:HZ1	3:B:501:2B0:H12	1.62	0.65
1:B:65:TRP:CE2	1:B:409[B]:MET:SD	2.90	0.65
1:B:65:TRP:HB2	1:B:283:LYS:HD3	1.80	0.62
1:B:25[A]:TYR:OH	1:B:160:ASP:OD1	2.17	0.62
1:A:89:HIS:HA	1:A:323:PRO:HD2	1.82	0.61
1:B:102:ARG:NH2	1:B:334:GLU:OE1	2.34	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:65:TRP:CZ2	1:B:409[B]:MET:SD	2.95	0.60
1:A:314:MET:HE1	1:B:25[A]:TYR:HD1	1.65	0.59
1:A:283:LYS:O	6:A:693:HOH:O	2.18	0.57
3:B:501:2B0:O2	3:B:501:2B0:H3	2.04	0.57
1:B:89:HIS:HA	1:B:323:PRO:HD2	1.87	0.56
1:B:124:GLY:CA	3:B:501:2B0:H4	2.31	0.56
1:A:286:THR:HG22	1:A:329[A]:SER:OG	2.06	0.55
1:B:407:TYR:CE2	1:B:409[A]:MET:SD	2.99	0.55
1:B:174:MET:HB3	1:B:177:LEU:HD13	1.89	0.54
1:A:314:MET:CE	1:B:25[A]:TYR:CD1	2.80	0.53
1:B:97:HIS:HE1	6:B:610:HOH:O	1.91	0.53
1:A:348:LEU:HD11	1:A:412:TYR:HA	1.92	0.50
1:B:395:ARG:CZ	1:B:428:GLU:HG2	2.42	0.50
1:B:389:THR:OG1	1:B:390:PRO:HD3	2.11	0.49
1:A:137:GLN:HG3	1:B:178:TRP:CH2	2.48	0.49
1:A:66:THR:HB	1:A:288:GLY:HA2	1.95	0.49
1:B:88:ASN:HD22	1:B:88:ASN:N	2.11	0.48
1:A:88:ASN:N	1:A:88:ASN:HD22	2.12	0.47
1:A:14:ILE:HD11	1:A:37:VAL:HG23	1.98	0.46
1:A:133:LYS:HG3	1:B:178:TRP:CH2	2.51	0.46
1:A:181:VAL:HG12	1:B:148:ARG:HH11	1.80	0.46
1:A:110:ILE:HG21	1:A:337:LEU:HD11	1.96	0.46
1:B:407:TYR:HE2	1:B:409[A]:MET:SD	2.39	0.46
1:A:318:THR:HG21	1:B:283:LYS:HG2	1.98	0.46
1:B:348:LEU:HD11	1:B:412:TYR:HA	1.98	0.45
1:A:163:LEU:HD13	1:A:163:LEU:C	2.37	0.45
1:B:174:MET:O	6:B:703:HOH:O	2.20	0.44
1:B:88:ASN:HD22	1:B:88:ASN:H	1.64	0.44
1:B:88:ASN:ND2	1:B:89:HIS:H	2.15	0.44
1:B:175[B]:HIS:O	1:B:175[B]:HIS:ND1	2.50	0.44
1:B:264:GLY:HA2	1:B:345:ILE:HG21	1.99	0.44
1:B:50:ARG:NH2	1:B:394:ASP:OD1	2.44	0.43
1:A:23:HIS:HB3	1:A:24:PRO:HD2	1.99	0.43
1:B:65:TRP:CD2	1:B:409[B]:MET:SD	3.11	0.43
1:A:264:GLY:HA2	1:A:345:ILE:HG21	1.99	0.42
1:A:65:TRP:CD1	1:A:65:TRP:N	2.88	0.42
1:A:129:GLU:OE1	1:A:163:LEU:HB3	2.19	0.42
1:B:65:TRP:CH2	1:B:409[B]:MET:SD	3.13	0.42
1:A:314:MET:HE1	1:B:25[A]:TYR:CD1	2.50	0.42
1:A:88:ASN:ND2	1:A:89:HIS:H	2.19	0.41
1:B:64:TRP:O	1:B:65:TRP:HB2	2.21	0.41
1:B:175[B]:HIS:C	1:B:175[B]:HIS:ND1	2.74	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	431/457 (94%)	419 (97%)	10 (2%)	2 (0%)	38	17
1	B	432/457 (94%)	422 (98%)	8 (2%)	2 (0%)	38	17
All	All	863/914 (94%)	841 (98%)	18 (2%)	4 (0%)	38	17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	283	LYS
1	A	283	LYS
1	A	316	GLY
1	B	316	GLY

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	333/346 (96%)	329 (99%)	4 (1%)	82	69
1	B	331/346 (96%)	323 (98%)	8 (2%)	61	39
All	All	664/692 (96%)	652 (98%)	12 (2%)	73	53

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ARG
1	A	34	SER

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Mol	Chain	Res	Type
1	A	88	ASN
1	A	153	TRP
1	B	30	ARG
1	B	78	GLN
1	B	88	ASN
1	B	122	ASP
1	B	153	TRP
1	B	180	ASP
1	B	409[A]	MET
1	B	409[B]	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	B	88	ASN
1	B	97	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 ⓘ

6 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	2BG	A	501	-	12,12,12	5.41	6 (50%)	14,16,16	4.22	8 (57%)
3	2B0	A	502	1	14,15,16	2.55	7 (50%)	20,22,23	2.41	6 (30%)
4	SO4	A	503	-	4,4,4	0.59	0	6,6,6	0.23	0
5	EDO	A	504	-	3,3,3	0.49	0	2,2,2	0.15	0
3	2B0	B	501	1	16,16,16	2.71	8 (50%)	23,23,23	1.75	4 (17%)
4	SO4	B	502	-	4,4,4	0.64	0	6,6,6	0.14	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	2BG	A	501	-	-	0/0/2/2	0/0/2/2
3	2B0	A	502	1	-	0/6/6/8	0/1/1/1
4	SO4	A	503	-	-	0/0/0/0	0/0/0/0
5	EDO	A	504	-	-	0/1/1/1	0/0/0/0
3	2B0	B	501	1	-	0/8/8/8	0/1/1/1
4	SO4	B	502	-	-	0/0/0/0	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	2BG	CAB-CAD	16.20	1.73	1.36
2	A	501	2BG	CAE-CAK	-5.75	1.33	1.40
3	A	502	2B0	CAI-CAN	-5.14	1.42	1.51
2	A	501	2BG	CAD-CAJ	-4.62	1.33	1.41
3	B	501	2B0	CAO-C1	-4.56	1.36	1.46
3	B	501	2B0	CAA-CAL	-4.54	1.41	1.50
3	B	501	2B0	CAI-CAN	-4.42	1.43	1.51
3	B	501	2B0	PAP-OAF	4.37	1.63	1.54
3	A	502	2B0	PAP-CAJ	-4.18	1.74	1.79
3	A	502	2B0	CAA-CAL	-4.15	1.42	1.50
2	A	501	2BG	CAI-SAH	-3.81	1.69	1.73
3	A	502	2B0	PAP-OAE	3.49	1.58	1.50
3	B	501	2B0	PAP-CAJ	-3.30	1.75	1.79
3	B	501	2B0	CAH-NAK	2.94	1.41	1.34
3	A	502	2B0	CAH-NAK	2.70	1.40	1.34
3	B	501	2B0	PAP-OAC	-2.67	1.49	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	2BG	CAK-SAH	-2.59	1.71	1.74
3	B	501	2B0	CAM-CAL	-2.41	1.39	1.40
2	A	501	2BG	CAJ-CAK	-2.30	1.34	1.42
3	A	502	2B0	CAM-CAL	-2.11	1.39	1.40
3	A	502	2B0	PAP-OAC	-2.07	1.50	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	2BG	CAC-CAB-CAD	-11.95	102.98	120.47
3	A	502	2B0	OAE-PAP-CAJ	-6.80	96.41	111.17
2	A	501	2BG	CAC-CAE-CAK	5.53	126.34	118.86
3	B	501	2B0	O2-C1-CAO	-4.55	115.50	125.32
3	A	502	2B0	CAI-CAN-CAO	4.13	125.50	121.43
3	B	501	2B0	CAM-CAO-CAN	4.00	121.32	118.15
3	A	502	2B0	OAF-PAP-CAJ	3.85	116.17	106.90
2	A	501	2BG	CAB-CAD-CAJ	3.60	125.67	120.02
2	A	501	2BG	CAB-CAC-CAE	3.54	125.65	120.47
2	A	501	2BG	CAE-CAK-SAH	3.45	131.89	125.58
2	A	501	2BG	CAJ-CAK-SAH	-3.28	107.91	112.03
2	A	501	2BG	CAK-CAJ-NAG	3.00	115.41	108.37
2	A	501	2BG	CAI-NAG-CAJ	2.99	113.95	107.69
3	B	501	2B0	OAF-PAP-CAJ	-2.98	99.73	106.90
3	A	502	2B0	CAM-CAL-NAK	-2.90	116.68	120.66
3	B	501	2B0	CAM-CAL-NAK	-2.26	117.56	120.66
3	A	502	2B0	PAP-CAJ-CAI	-2.17	109.36	113.11
3	A	502	2B0	CAO-CAM-CAL	2.14	122.13	119.67

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, 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	425/457 (92%)	-0.07	5 (1%) 75 82	9, 16, 29, 46	5 (1%)
1	B	427/457 (93%)	0.04	13 (3%) 48 53	8, 17, 36, 53	11 (2%)
All	All	852/914 (93%)	-0.02	18 (2%) 60 66	8, 17, 34, 53	16 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	32	ALA	4.9
1	B	179	THR	4.6
1	B	175[A]	HIS	4.4
1	B	177	LEU	4.1
1	B	310	ALA	3.4
1	B	180	ASP	3.1
1	A	309	ALA	3.0
1	A	30	ARG	3.0
1	A	307	ALA	2.8
1	B	174	MET	2.8
1	B	176	SER	2.7
1	B	65	TRP	2.6
1	B	7	GLY	2.5
1	B	42	HIS	2.4
1	B	178	TRP	2.3
1	B	33	VAL	2.3
1	A	34	SER	2.2
1	A	122	ASP	2.1

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	EDO	A	504	4/4	0.30	7.44	33,33,33,38	0
4	SO4	A	503	5/5	0.13	2.44	26,26,28,30	0
2	2BG	A	501	11/11	0.13	2.42	19,24,26,32	2
4	SO4	B	502	5/5	0.17	1.83	36,39,40,40	0
3	2B0	A	502	15/16	0.10	1.32	10,11,17,17	0
3	2B0	B	501	16/16	0.09	0.20	14,15,18,21	0

6.5 Other polymers

There are no such residues in this entry.