



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:52 AM UTC

PDB ID : 5NEK / pdb_00005nek
Title : Crystal structure of the polysaccharide deacetylase Bc1974 from *Bacillus cereus* in complex with acetazolamide
Authors : Andreou, A.; Giastas, P.; Eliopoulos, E.E.
Deposited on : 2017-03-10
Resolution : 3.06 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

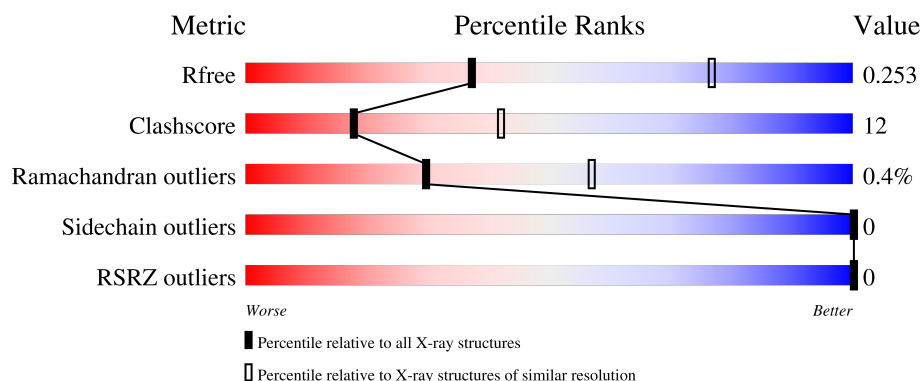
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	247	
1	B	247	
1	C	247	
1	D	247	

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
4	AZM	B	302	-	X	-	-
4	AZM	D	302	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6616 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.

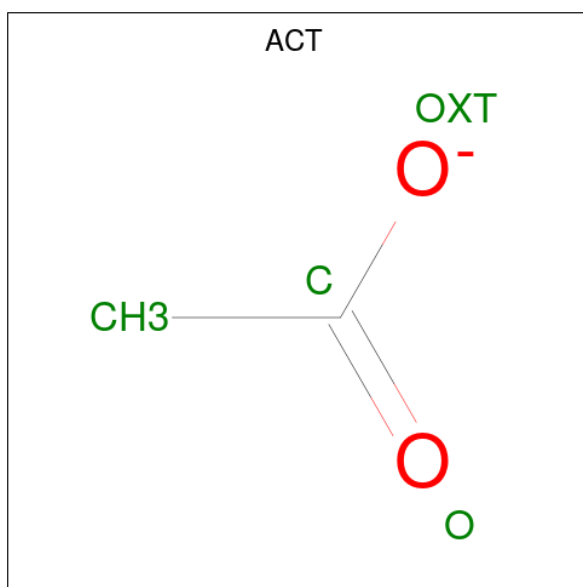
- Molecule 1 is a protein called Peptidoglycan N-acetylglucosamine deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1624	1040	277	300	7			
1	B	206	Total	C	N	O	S	0	0	0
			1645	1054	281	303	7			
1	C	205	Total	C	N	O	S	0	0	0
			1616	1034	275	300	7			
1	D	206	Total	C	N	O	S	0	1	0
			1657	1061	284	305	7			

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

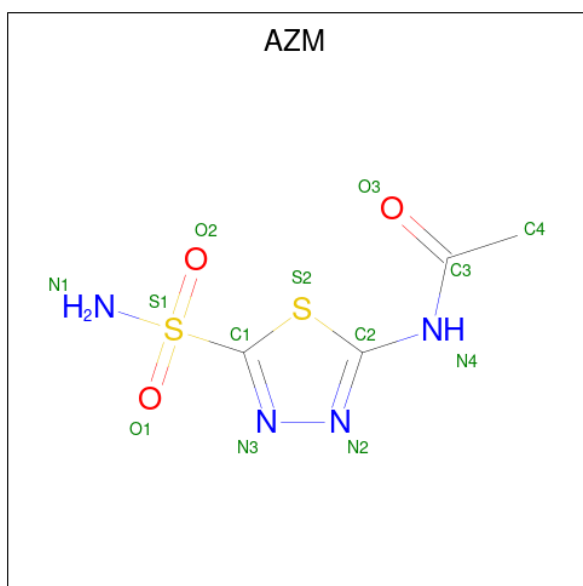
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 4 is 5-ACETAMIDO-1,3,4-THIADIAZOLE-2-SULFONAMIDE (CCD ID: AZM) (formula: $C_4H_6N_4O_3S_2$).



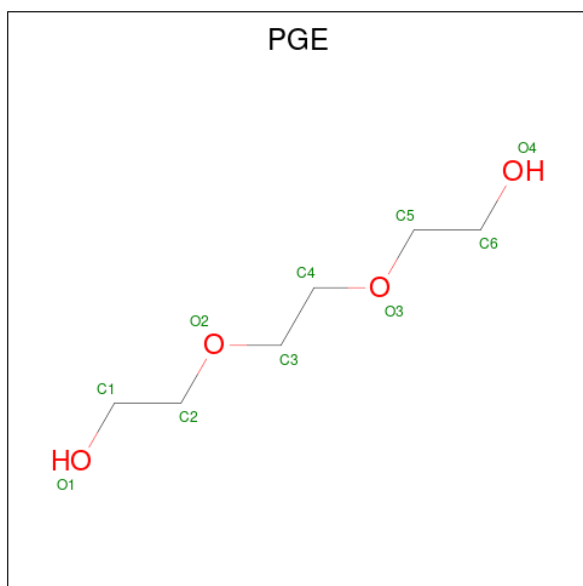
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			13	4	4	3	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	S	0	0
			13	4	4	3	2		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			10	6	4		

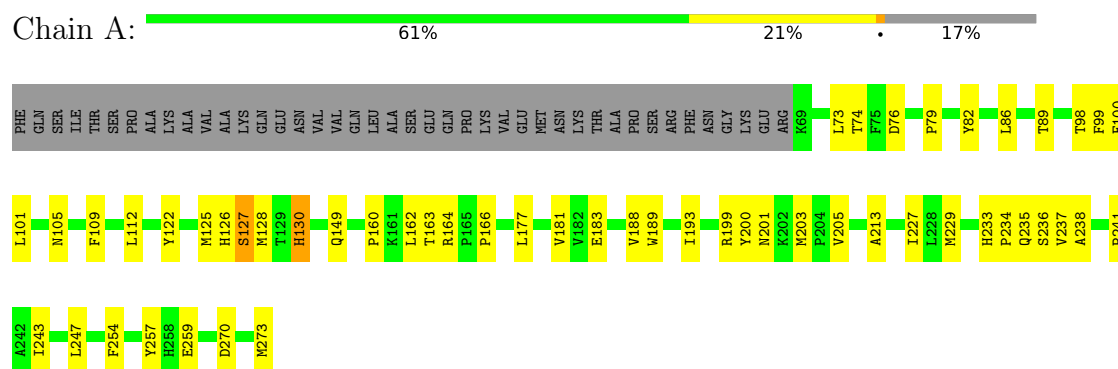
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	5	Total	O	0	0
			5	5		
6	B	9	Total	O	0	0
			9	9		
6	C	5	Total	O	0	0
			5	5		
6	D	7	Total	O	0	0
			7	7		

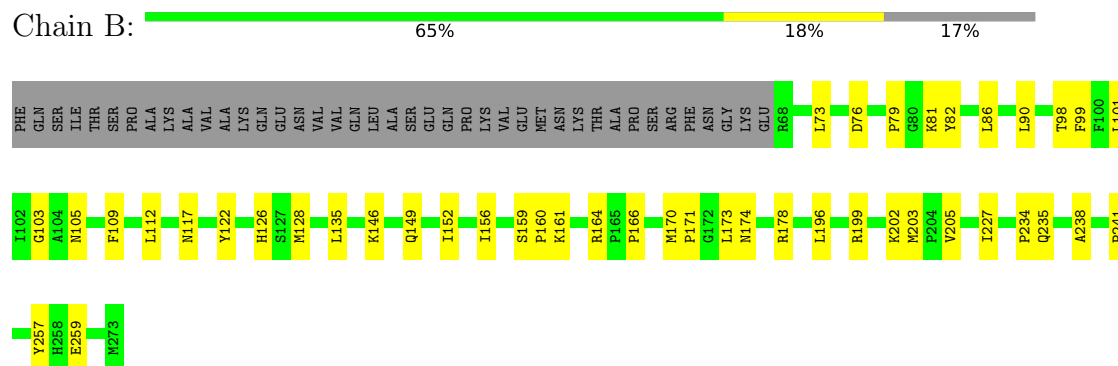
3 Residue-property plots

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.

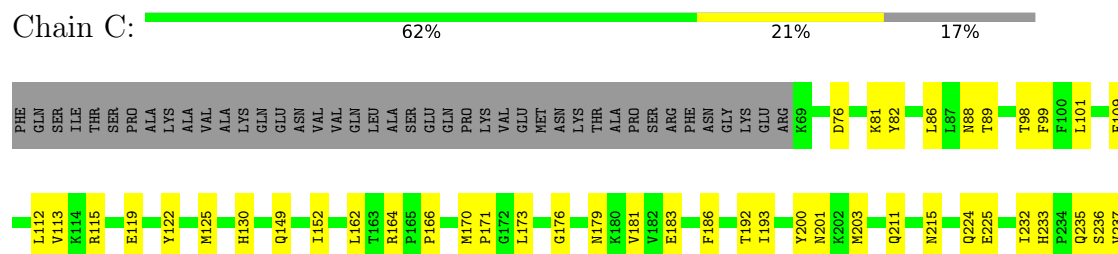
• Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase



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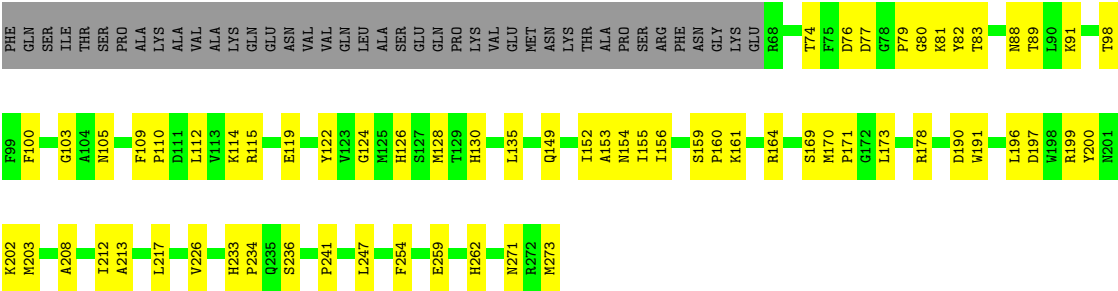


• Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase





● Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.39Å 118.21Å 98.27Å 90.00° 102.09° 90.00°	Depositor
Resolution (Å)	48.29 – 3.06 48.29 – 3.06	Depositor EDS
% Data completeness (in resolution range)	96.3 (48.29-3.06) 96.6 (48.29-3.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 3.07Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.178 , 0.253 0.181 , 0.253	Depositor DCC
R_{free} test set	1030 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.930	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 78.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6616	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.59% 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, AZM, ZN, PGE

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	0.43	0/1669	0.63	0/2264
1	B	0.42	0/1690	0.64	0/2288
1	C	0.37	0/1661	0.60	0/2256
1	D	0.43	0/1705	0.64	0/2307
All	All	0.41	0/6725	0.63	0/9115

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	1624	0	1561	38	0
1	B	1645	0	1600	34	0
1	C	1616	0	1539	38	0
1	D	1657	0	1617	55	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4	0	3	0	0
4	B	13	0	5	2	0
4	D	13	0	5	1	0
5	D	10	0	14	3	0
6	A	5	0	0	0	0
6	B	9	0	0	0	0
6	C	5	0	0	1	0
6	D	7	0	0	1	0
All	All	6616	0	6347	159	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 (159) 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:D:91:LYS:NZ	6:D:401:HOH:O	2.00	0.94
1:C:125:MET:HE3	1:C:152:ILE:HD12	1.56	0.86
1:C:201:ASN:OD1	1:C:235:GLN:NE2	2.16	0.78
1:C:270:ASP:HB3	1:C:273:MET:HE2	1.67	0.75
1:D:161:LYS:HE3	1:D:259:GLU:HB3	1.68	0.74
1:C:98:THR:HG23	1:C:122:TYR:HB3	1.69	0.73
1:D:196:LEU:HD13	5:D:303:PGE:H1	1.71	0.72
1:B:196:LEU:HD13	5:D:303:PGE:H6	1.71	0.71
1:A:109:PHE:HB3	1:A:112:LEU:HD12	1.73	0.70
1:D:98:THR:HG23	1:D:122:TYR:HB3	1.72	0.70
1:D:109:PHE:HB3	1:D:112:LEU:HD12	1.73	0.69
1:C:89:THR:HG21	1:C:241:PRO:HG3	1.75	0.68
1:A:127:SER:OG	1:A:128:MET:N	2.23	0.67
1:D:89:THR:HG21	1:D:241:PRO:HG3	1.79	0.64
1:B:149:GLN:HG3	1:B:160:PRO:HD2	1.80	0.64
1:A:270:ASP:HB3	1:A:273:MET:HE2	1.79	0.63
1:C:88:ASN:OD1	1:C:115:ARG:NH2	2.22	0.63
1:C:125:MET:HE1	1:C:149:GLN:CA	2.29	0.62
1:D:88:ASN:OD1	1:D:115:ARG:NH2	2.26	0.62
1:D:161:LYS:HG2	1:D:259:GLU:OE2	2.01	0.60
1:B:161:LYS:HD2	1:B:259:GLU:CD	2.27	0.60
1:C:125:MET:HE1	1:C:149:GLN:N	2.16	0.60
1:D:153:ALA:C	1:D:155:ILE:H	2.10	0.59
1:A:183:GLU:OE2	1:B:81:LYS:NZ	2.24	0.59
1:D:114:LYS:HD3	1:D:155:ILE:HG23	1.82	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:MET:HE2	1:A:163:THR:HG22	1.84	0.59
1:A:130:HIS:CE1	1:A:166:PRO:HB3	2.37	0.59
1:B:103:GLY:HA3	1:B:128:MET:SD	2.42	0.59
1:A:126:HIS:O	1:A:127:SER:HB3	2.03	0.58
1:D:262:HIS:NE2	1:D:273:MET:OXT	2.31	0.58
1:D:152:ILE:O	1:D:156:ILE:HG12	2.03	0.58
1:C:200:TYR:O	1:C:203:MET:HB2	2.03	0.58
1:C:125:MET:HE1	1:C:149:GLN:HA	1.86	0.58
1:D:114:LYS:CE	1:D:155:ILE:HG23	2.36	0.56
1:A:99:PHE:HB3	1:A:101:LEU:HD21	1.88	0.56
1:D:155:ILE:HG22	1:D:156:ILE:HG23	1.88	0.56
1:D:82:TYR:CZ	1:D:234:PRO:HD3	2.41	0.56
1:B:199:ARG:NH1	1:D:202:LYS:HG3	2.22	0.55
1:B:257:TYR:HE2	1:B:259:GLU:HG2	1.71	0.55
1:B:82:TYR:CZ	1:B:234:PRO:HD3	2.42	0.55
1:C:181:VAL:HG13	1:C:186:PHE:HB2	1.89	0.55
1:C:99:PHE:HB3	1:C:101:LEU:HD21	1.90	0.54
1:C:109:PHE:HB3	1:C:112:LEU:HD12	1.88	0.54
1:D:91:LYS:HE3	1:D:119:GLU:CD	2.33	0.54
1:B:73:LEU:HD23	1:B:227:ILE:HB	1.90	0.54
1:A:193:ILE:HB	1:A:227:ILE:HG12	1.89	0.53
1:B:149:GLN:NE2	1:B:159:SER:HB3	2.24	0.53
1:C:130:HIS:CE1	1:C:166:PRO:HB3	2.42	0.53
1:A:201:ASN:OD1	1:A:235:GLN:NE2	2.42	0.52
1:C:224:GLN:O	1:C:225:GLU:HG2	2.10	0.52
1:D:80:GLY:N	1:D:83:THR:OG1	2.35	0.52
1:A:200:TYR:O	1:A:203:MET:HB2	2.10	0.52
1:C:119:GLU:HB3	6:C:402:HOH:O	2.09	0.51
1:A:162:LEU:HD22	1:A:257:TYR:CD2	2.45	0.51
1:B:109:PHE:HB3	1:B:112:LEU:HD12	1.92	0.51
1:D:74:THR:HG22	1:D:98:THR:HB	1.92	0.51
1:B:98:THR:HG23	1:B:122:TYR:HB3	1.92	0.51
1:C:233:HIS:HB3	1:C:235:GLN:NE2	2.25	0.51
1:C:86:LEU:HB2	1:C:237:VAL:HG13	1.93	0.51
1:B:86:LEU:O	1:B:90:LEU:HG	2.10	0.50
1:A:238:ALA:O	1:A:241:PRO:HD2	2.12	0.49
1:A:247:LEU:HB2	1:A:254:PHE:HZ	1.77	0.49
1:D:77:ASP:N	1:D:77:ASP:OD1	2.43	0.49
1:B:99:PHE:HB3	1:B:101:LEU:HD21	1.94	0.49
1:D:203:MET:HE2	1:D:208:ALA:N	2.28	0.49
1:D:126:HIS:CE1	4:D:302:AZM:N1	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:TYR:HB2	1:C:237:VAL:HG21	1.95	0.49
1:D:173:LEU:O	1:D:178:ARG:HG3	2.13	0.49
1:A:82:TYR:CE2	1:A:234:PRO:HD3	2.48	0.48
1:A:149:GLN:NE2	1:A:160:PRO:O	2.37	0.48
1:C:238:ALA:O	1:C:241:PRO:HD2	2.14	0.48
1:A:205:VAL:HG13	1:A:235:GLN:HA	1.95	0.48
1:B:203:MET:HE3	1:B:203:MET:HB2	1.81	0.48
1:D:114:LYS:CD	1:D:155:ILE:HG23	2.44	0.47
1:D:149:GLN:NE2	1:D:159:SER:HB3	2.29	0.47
1:A:199:ARG:O	1:D:271:ASN:ND2	2.37	0.47
1:B:117:ASN:ND2	1:B:156:ILE:HB	2.28	0.47
1:A:188:VAL:O	1:A:189:TRP:HD1	1.97	0.47
1:D:100:PHE:HB3	1:D:126:HIS:HB2	1.96	0.47
1:D:200:TYR:HB3	1:D:208:ALA:HB1	1.97	0.47
1:C:183:GLU:OE2	1:D:81:LYS:NZ	2.31	0.47
1:D:100:PHE:HA	1:D:124:GLY:O	2.14	0.47
1:A:125:MET:HE3	1:A:149:GLN:HB2	1.97	0.47
1:D:233:HIS:HB2	1:D:236:SER:OG	2.15	0.47
1:B:202:LYS:HD2	1:D:199:ARG:NH1	2.30	0.46
1:B:257:TYR:CE2	1:B:259:GLU:HG2	2.50	0.46
1:C:81:LYS:HE3	1:C:82:TYR:CZ	2.50	0.46
1:C:76:ASP:OD1	1:C:164:ARG:NH1	2.49	0.46
1:C:232:ILE:HG13	1:C:233:HIS:CD2	2.50	0.46
1:D:197:ASP:HA	1:D:212:ILE:HG12	1.97	0.46
1:D:203:MET:HE3	1:D:203:MET:HB2	1.67	0.46
1:B:166:PRO:HA	4:B:302:AZM:HN11	1.80	0.46
1:B:174:ASN:O	1:B:178:ARG:HG3	2.16	0.46
1:D:153:ALA:C	1:D:155:ILE:N	2.74	0.46
1:A:200:TYR:CD1	1:A:200:TYR:N	2.83	0.46
1:C:130:HIS:ND1	1:C:166:PRO:HB3	2.31	0.46
1:D:247:LEU:HB2	1:D:254:PHE:CZ	2.51	0.45
1:D:170:MET:HA	1:D:171:PRO:HA	1.71	0.45
1:B:73:LEU:HD23	1:B:73:LEU:HA	1.76	0.45
1:C:109:PHE:O	1:C:113:VAL:HG23	2.17	0.45
1:D:173:LEU:HD12	1:D:173:LEU:HA	1.70	0.45
1:A:74:THR:OG1	1:A:227:ILE:O	2.32	0.45
1:B:202:LYS:HD2	1:D:199:ARG:HH12	1.82	0.45
1:D:152:ILE:HG21	1:D:160:PRO:HG2	2.00	0.45
1:B:170:MET:HA	1:B:171:PRO:HA	1.73	0.44
1:B:259:GLU:H	1:B:259:GLU:HG3	1.50	0.44
1:C:170:MET:HA	1:C:171:PRO:HA	1.67	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:192:THR:OG1	1:C:193:ILE:HG13	2.17	0.44
1:D:76:ASP:HB3	1:D:77:ASP:OD1	2.17	0.44
1:A:177:LEU:HD23	1:A:177:LEU:HA	1.76	0.44
1:A:98:THR:HG23	1:A:122:TYR:HB3	2.00	0.44
1:D:79:PRO:HD2	1:D:105:ASN:HB3	2.00	0.44
1:A:100:PHE:HB3	1:A:126:HIS:HB2	2.00	0.43
1:A:213:ALA:HA	1:A:243:ILE:HG12	1.99	0.43
1:C:211:GLN:NE2	1:C:215:ASN:OD1	2.49	0.43
1:D:202:LYS:N	5:D:303:PGE:O4	2.51	0.43
1:D:247:LEU:HB2	1:D:254:PHE:HZ	1.83	0.43
1:A:76:ASP:OD1	1:A:164:ARG:NH1	2.51	0.43
1:A:233:HIS:HB3	1:A:235:GLN:NE2	2.33	0.43
1:C:173:LEU:HD12	1:C:173:LEU:HA	1.90	0.43
1:A:247:LEU:HB2	1:A:254:PHE:CZ	2.53	0.43
1:B:205:VAL:HG13	1:B:235:GLN:HA	2.00	0.43
1:C:203:MET:HB3	1:C:203:MET:HE2	1.67	0.43
1:C:247:LEU:HB2	1:C:254:PHE:CZ	2.54	0.43
1:B:173:LEU:HA	1:B:173:LEU:HD12	1.58	0.43
1:A:73:LEU:HD23	1:A:73:LEU:HA	1.82	0.42
1:A:229:MET:HB3	1:A:229:MET:HE2	1.64	0.42
1:A:130:HIS:ND1	1:A:166:PRO:HB3	2.34	0.42
1:B:126:HIS:ND1	1:B:164:ARG:O	2.48	0.42
1:D:100:PHE:CG	1:D:164:ARG:HB3	2.54	0.42
1:D:114:LYS:NZ	1:D:155:ILE:HG23	2.34	0.42
1:B:152:ILE:O	1:B:156:ILE:HG12	2.18	0.42
1:C:247:LEU:HB2	1:C:254:PHE:HZ	1.83	0.42
1:D:135:LEU:HA	1:D:135:LEU:HD23	1.81	0.42
1:D:169:SER:OG	1:D:190:ASP:OD2	2.35	0.42
1:A:79:PRO:HD2	1:A:105:ASN:HB3	2.01	0.42
1:C:162:LEU:HD22	1:C:257:TYR:CD2	2.54	0.42
1:D:213:ALA:O	1:D:217:LEU:HG	2.20	0.42
1:B:146:LYS:HB2	1:B:146:LYS:HE3	1.78	0.42
1:B:166:PRO:HA	4:B:302:AZM:N1	2.34	0.42
1:C:115:ARG:O	1:C:119:GLU:HG3	2.20	0.42
1:C:233:HIS:HB2	1:C:236:SER:OG	2.20	0.42
1:B:79:PRO:HD2	1:B:105:ASN:HB3	2.02	0.42
1:B:238:ALA:O	1:B:241:PRO:HD2	2.20	0.42
1:D:103:GLY:HA3	1:D:128:MET:SD	2.60	0.41
1:B:135:LEU:HD23	1:B:135:LEU:HA	1.90	0.41
1:A:86:LEU:HB2	1:A:237:VAL:HG13	2.02	0.41
1:B:76:ASP:OD1	1:B:164:ARG:NH1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:MET:HB3	1:C:125:MET:HE2	1.68	0.41
1:A:89:THR:HG21	1:A:241:PRO:HG3	2.03	0.41
1:D:200:TYR:CD2	1:D:200:TYR:N	2.88	0.41
1:D:110:PRO:O	1:D:114:LYS:NZ	2.53	0.41
1:A:122:TYR:OH	1:A:259:GLU:OE2	2.35	0.41
1:C:176:GLY:O	1:C:179:ASN:HB2	2.20	0.41
1:A:233:HIS:HB2	1:A:236:SER:OG	2.20	0.41
1:D:115:ARG:O	1:D:119:GLU:HG3	2.21	0.41
1:D:126:HIS:CD2	1:D:130:HIS:HD2	2.39	0.40
1:D:191:TRP:HB2	1:D:226:VAL:O	2.22	0.40
1:A:181:VAL:HG11	1:A:188:VAL:HG23	2.02	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	203/247 (82%)	192 (95%)	9 (4%)	2 (1%)	12	37
1	B	204/247 (83%)	192 (94%)	12 (6%)	0	100	100
1	C	203/247 (82%)	196 (97%)	7 (3%)	0	100	100
1	D	205/247 (83%)	192 (94%)	12 (6%)	1 (0%)	24	52
All	All	815/988 (82%)	772 (95%)	40 (5%)	3 (0%)	30	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	154	ASN
1	A	130	HIS
1	A	127	SER

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	168/208 (81%)	168 (100%)	0	100	100
1	B	172/208 (83%)	172 (100%)	0	100	100
1	C	166/208 (80%)	166 (100%)	0	100	100
1	D	174/208 (84%)	174 (100%)	0	100	100
All	All	680/832 (82%)	680 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	B	269	HIS
1	C	117	ASN
1	C	224	GLN
1	D	93	HIS
1	D	138	ASN
1	D	219	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	AZM	B	302	2	13,13,13	8.13	9 (69%)	19,19,19	12.40	13 (68%)
5	PGE	D	303	-	9,9,9	0.37	0	8,8,8	0.41	0
4	AZM	D	302	1,2	13,13,13	8.26	10 (76%)	19,19,19	11.21	12 (63%)
3	ACT	C	302	2	3,3,3	0.83	0	3,3,3	1.34	0
3	ACT	A	302	2	3,3,3	0.85	0	3,3,3	1.47	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
4	AZM	B	302	2	-	2/10/10/10	0/1/1/1
5	PGE	D	303	-	-	5/7/7/7	-
4	AZM	D	302	1,2	-	6/10/10/10	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	302	AZM	C1-S2	-17.46	1.54	1.73
4	B	302	AZM	C1-N3	17.11	1.42	1.29
4	B	302	AZM	C1-S2	-16.70	1.55	1.73
4	D	302	AZM	C1-N3	15.67	1.41	1.29
4	D	302	AZM	C2-S2	-11.84	1.54	1.73
4	B	302	AZM	C2-S2	-10.78	1.55	1.73
4	B	302	AZM	N3-N2	8.99	1.58	1.39
4	D	302	AZM	N3-N2	8.44	1.57	1.39
4	D	302	AZM	C2-N2	8.09	1.43	1.31
4	B	302	AZM	C2-N2	7.52	1.42	1.31
4	D	302	AZM	O2-S1	4.36	1.50	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	302	AZM	C2-N4	3.69	1.43	1.38
4	D	302	AZM	C2-N4	3.42	1.43	1.38
4	D	302	AZM	S1-N1	3.13	1.65	1.58
4	D	302	AZM	C3-N4	3.05	1.43	1.37
4	B	302	AZM	C3-N4	2.55	1.42	1.37
4	D	302	AZM	C1-S1	-2.49	1.76	1.78
4	B	302	AZM	S1-N1	2.39	1.64	1.58
4	B	302	AZM	C1-S1	2.29	1.79	1.78

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	302	AZM	C1-S2-C2	46.65	107.61	85.47
4	D	302	AZM	C1-S2-C2	41.34	105.09	85.47
4	B	302	AZM	S2-C1-N3	-17.09	107.10	116.10
4	B	302	AZM	O2-S1-O1	-13.50	100.96	120.27
4	D	302	AZM	O2-S1-O1	-12.63	102.21	120.27
4	D	302	AZM	N4-C2-N2	12.19	131.70	120.81
4	D	302	AZM	S2-C1-N3	-9.60	111.04	116.10
4	B	302	AZM	N4-C2-N2	9.49	129.28	120.81
4	D	302	AZM	O2-S1-N1	9.28	121.31	108.39
4	D	302	AZM	C1-N3-N2	-8.70	106.09	111.48
4	B	302	AZM	O1-S1-C1	7.01	116.09	105.65
4	B	302	AZM	S2-C2-N2	-6.34	107.45	114.52
4	D	302	AZM	C1-S1-N1	-5.84	101.59	107.40
4	D	302	AZM	S2-C2-N2	-5.74	108.11	114.52
4	B	302	AZM	S2-C1-S1	5.30	128.66	121.62
4	B	302	AZM	C4-C3-N4	3.64	121.15	115.27
4	B	302	AZM	C1-N3-N2	-3.62	109.23	111.48
4	D	302	AZM	O1-S1-N1	3.59	113.39	108.39
4	D	302	AZM	S2-C1-S1	3.52	126.29	121.62
4	B	302	AZM	O2-S1-N1	3.19	112.83	108.39
4	D	302	AZM	O2-S1-C1	2.94	110.03	105.65
4	B	302	AZM	C2-N2-N3	-2.68	108.61	112.02
4	D	302	AZM	S2-C2-N4	-2.66	120.16	123.72
4	B	302	AZM	O3-C3-N4	-2.46	118.02	121.90
4	B	302	AZM	C3-N4-C2	-2.16	120.49	124.05

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	302	AZM	N2-C2-N4-C3
4	B	302	AZM	S2-C2-N4-C3
4	D	302	AZM	N2-C2-N4-C3
4	D	302	AZM	S2-C2-N4-C3
5	D	303	PGE	O3-C5-C6-O4
5	D	303	PGE	O1-C1-C2-O2
5	D	303	PGE	O2-C3-C4-O3
4	D	302	AZM	N3-C1-S1-N1
5	D	303	PGE	C3-C4-O3-C5
4	D	302	AZM	N3-C1-S1-O2
4	D	302	AZM	S2-C1-S1-N1
4	D	302	AZM	S2-C1-S1-O2
5	D	303	PGE	C4-C3-O2-C2

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	302	AZM	2	0
5	D	303	PGE	3	0
4	D	302	AZM	1	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	205/247 (82%)	-0.31	0 100 100	41, 70, 105, 124	0
1	B	206/247 (83%)	-0.36	0 100 100	43, 67, 96, 132	0
1	C	205/247 (82%)	-0.23	0 100 100	43, 72, 118, 135	0
1	D	206/247 (83%)	-0.43	0 100 100	37, 61, 86, 113	1 (0%)
All	All	822/988 (83%)	-0.33	0 100 100	37, 68, 106, 135	1 (0%)

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
3	ACT	A	302	4/4	0.82	0.26	69,84,87,87	0
4	AZM	D	302	13/13	0.83	0.17	77,94,109,111	13
3	ACT	C	302	4/4	0.84	0.27	83,84,85,87	0
4	AZM	B	302	13/13	0.90	0.16	75,89,99,99	13

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	PGE	D	303	10/10	0.93	0.19	58,65,74,79	10
2	ZN	B	301	1/1	0.99	0.06	96,96,96,96	0
2	ZN	A	301	1/1	0.99	0.04	67,67,67,67	0
2	ZN	D	301	1/1	1.00	0.04	82,82,82,82	0
2	ZN	C	301	1/1	1.00	0.02	66,66,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.