



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:54 PM GMT

PDB ID : 4QDI
Title : Crystal structure II of MurF from *Acinetobacter baumannii*
Authors : An, Y.J.; Jeong, C.S.; Cha, S.S.
Deposited on : 2014-05-13
Resolution : 1.80 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

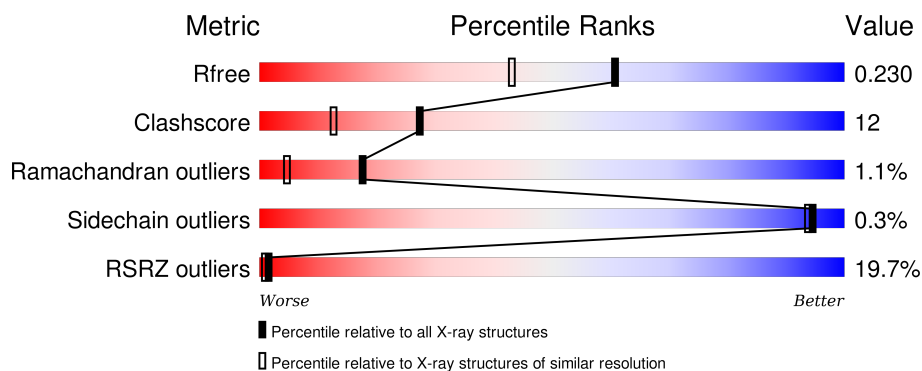
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.80 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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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. 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	472	

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	MG	A	503	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	A	505	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3774 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.

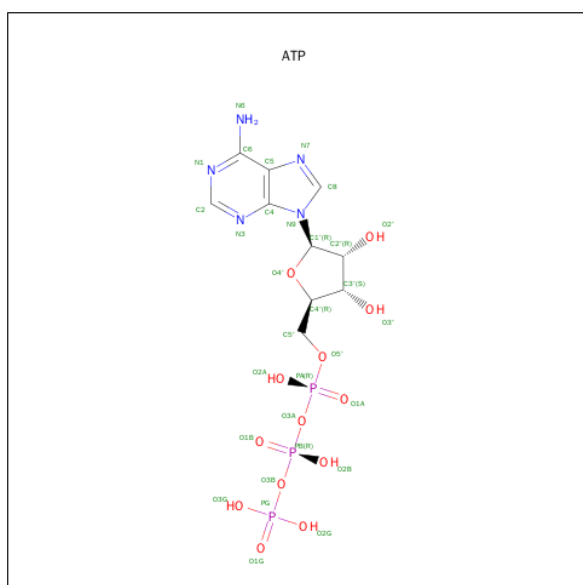
- Molecule 1 is a protein called UDP-N-acetylmuramoyl-tripeptide--D-alanyl-D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	457	3444	2164	619	647	14	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

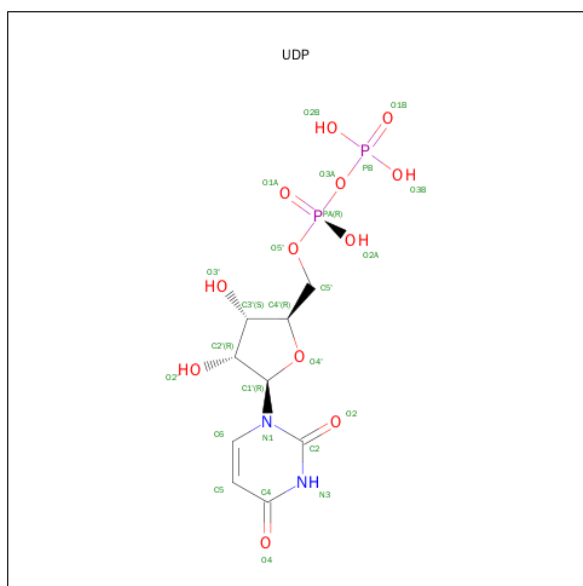
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP B7GVN5
A	-4	HIS	-	EXPRESSION TAG	UNP B7GVN5
A	-3	HIS	-	EXPRESSION TAG	UNP B7GVN5
A	-2	HIS	-	EXPRESSION TAG	UNP B7GVN5
A	-1	HIS	-	EXPRESSION TAG	UNP B7GVN5
A	0	HIS	-	EXPRESSION TAG	UNP B7GVN5

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

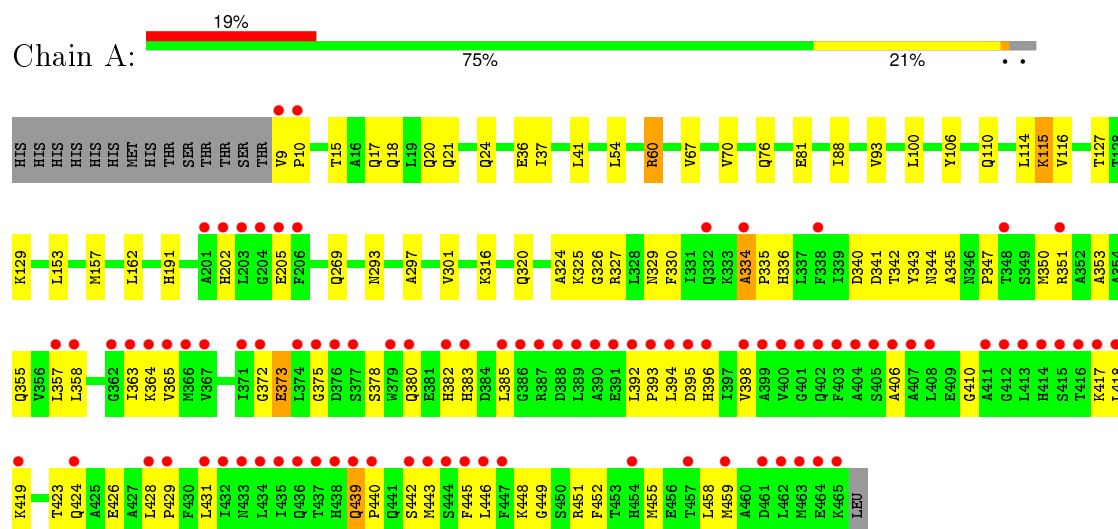
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	265	Total	O	0	0
			265	265		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: UDP-N-acetylmuramoyl-tripeptide--D-alanyl-D-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.78Å 85.78Å 130.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.48 – 1.80 37.47 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.2 (37.48-1.80) 97.2 (37.47-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.46 (at 1.81Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.227 , 0.253 0.228 , 0.230	Depositor DCC
R_{free} test set	2583 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.593	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.4	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 50545 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3774	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: UDP, MG, ATP, EDO

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	0.30	0/3508	0.57	2/4756 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	347	PRO	N-CA-CB	5.44	109.83	103.30
1	A	115	LYS	N-CA-C	-5.18	97.01	111.00

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	3444	0	3357	82	0
2	A	31	0	12	0	0
3	A	25	0	11	0	0
4	A	1	0	0	0	0
5	A	8	0	8	0	0
6	A	265	0	0	1	0
All	All	3774	0	3388	82	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 (82) 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:A:334:ALA:HB1	1:A:335:PRO:HD2	1.46	0.95
1:A:116:VAL:H	1:A:191:HIS:HD2	1.25	0.85
1:A:365:VAL:HG12	1:A:396:HIS:HB2	1.62	0.81
1:A:17:GLN:HG2	1:A:21:GLN:HE21	1.48	0.79
1:A:341:ASP:O	1:A:448:LYS:HD2	1.85	0.75
1:A:355:GLN:NE2	1:A:358:LEU:HD12	2.03	0.72
1:A:334:ALA:HB1	1:A:335:PRO:CD	2.19	0.71
1:A:350:MET:HG2	1:A:385:LEU:HD21	1.73	0.70
1:A:363:ILE:HB	1:A:443:MET:HG2	1.74	0.69
1:A:36:GLU:HG2	6:A:813:HOH:O	1.93	0.69
1:A:428:LEU:HB3	1:A:429:PRO:HD3	1.78	0.65
1:A:448:LYS:HG3	1:A:449:GLY:H	1.61	0.64
1:A:358:LEU:HD23	1:A:364:LYS:HD3	1.79	0.64
1:A:334:ALA:CB	1:A:335:PRO:HD2	2.26	0.64
1:A:202:HIS:HB2	1:A:205:GLU:HB2	1.79	0.63
1:A:326:GLY:H	1:A:329:ASN:ND2	1.98	0.61
1:A:116:VAL:H	1:A:191:HIS:CD2	2.12	0.60
1:A:269:GLN:HG2	1:A:330:PHE:CD2	2.35	0.60
1:A:353:ALA:HB1	1:A:446:LEU:HD21	1.83	0.60
1:A:372:GLY:HA3	1:A:452:PHE:HE2	1.69	0.58
1:A:17:GLN:CG	1:A:21:GLN:HE21	2.15	0.58
1:A:153:LEU:O	1:A:157:MET:HG3	2.05	0.57
1:A:439:GLN:CB	1:A:440:PRO:HD2	2.34	0.57
1:A:424:GLN:HB3	1:A:458:LEU:HD13	1.87	0.57
1:A:343:TYR:O	1:A:451:ARG:HD2	2.05	0.55
1:A:395:ASP:O	1:A:417:LYS:HG3	2.06	0.55
1:A:357:LEU:HB2	1:A:446:LEU:HD22	1.90	0.53
1:A:448:LYS:O	1:A:455:MET:HG3	2.08	0.53
1:A:340:ASP:HB2	1:A:459:MET:CE	2.39	0.53
1:A:344:ASN:O	1:A:448:LYS:HE2	2.10	0.51
1:A:372:GLY:O	1:A:373:GLU:HB2	2.11	0.51
1:A:380:GLN:HA	1:A:380:GLN:HE21	1.76	0.50
1:A:365:VAL:HG23	1:A:365:VAL:O	2.11	0.50
1:A:37:ILE:N	1:A:37:ILE:HD12	2.27	0.50
1:A:325:LYS:HA	1:A:329:ASN:HD21	1.76	0.49
1:A:418:LEU:HD23	1:A:419:LYS:N	2.27	0.49
1:A:106:TYR:O	1:A:110:GLN:HG2	2.13	0.49
1:A:357:LEU:HD13	1:A:446:LEU:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:GLN:HE21	1:A:355:GLN:HA	1.77	0.48
1:A:41:LEU:H	1:A:41:LEU:HD23	1.78	0.48
1:A:383:HIS:O	1:A:410:GLY:HA3	2.13	0.48
1:A:431:LEU:HD12	1:A:445:PHE:HE1	1.79	0.48
1:A:9:VAL:HA	1:A:10:PRO:O	2.14	0.48
1:A:60:ARG:HH11	1:A:60:ARG:HG2	1.78	0.48
1:A:334:ALA:CB	1:A:335:PRO:CD	2.89	0.48
1:A:378:SER:O	1:A:382:HIS:ND1	2.47	0.47
1:A:269:GLN:HG2	1:A:330:PHE:CE2	2.48	0.47
1:A:334:ALA:C	1:A:336:HIS:H	2.18	0.47
1:A:326:GLY:H	1:A:329:ASN:HD21	1.61	0.47
1:A:335:PRO:O	1:A:442:SER:HA	2.15	0.47
1:A:392:LEU:C	1:A:394:LEU:H	2.18	0.47
1:A:76:GLN:C	1:A:88:ILE:HG13	2.35	0.47
1:A:355:GLN:NE2	1:A:355:GLN:HA	2.30	0.46
1:A:67:VAL:O	1:A:70:VAL:HG22	2.16	0.46
1:A:423:THR:HG23	1:A:426:GLU:H	1.80	0.46
1:A:380:GLN:NE2	1:A:380:GLN:HA	2.31	0.46
1:A:81:GLU:HA	1:A:93:VAL:O	2.16	0.46
1:A:355:GLN:HE22	1:A:358:LEU:HD12	1.78	0.45
1:A:351:ARG:NH1	1:A:385:LEU:HD13	2.32	0.45
1:A:15:THR:OG1	1:A:18:GLN:HG3	2.15	0.45
1:A:54:LEU:HD21	1:A:100:LEU:HB2	1.99	0.44
1:A:316:LYS:O	1:A:320:GLN:HG2	2.17	0.44
1:A:297:ALA:O	1:A:301:VAL:HG23	2.18	0.44
1:A:324:ALA:HB3	1:A:327:ARG:HD3	1.99	0.44
1:A:365:VAL:HA	1:A:396:HIS:O	2.17	0.43
1:A:114:LEU:HD11	1:A:162:LEU:HD22	2.00	0.43
1:A:383:HIS:HA	1:A:406:ALA:O	2.17	0.43
1:A:9:VAL:HA	1:A:10:PRO:C	2.39	0.43
1:A:365:VAL:CG1	1:A:396:HIS:HB2	2.41	0.43
1:A:345:ALA:HB2	1:A:448:LYS:HG3	2.01	0.43
1:A:351:ARG:NH2	1:A:385:LEU:HA	2.34	0.43
1:A:342:THR:O	1:A:343:TYR:C	2.57	0.42
1:A:439:GLN:CB	1:A:440:PRO:CD	2.97	0.42
1:A:115:LYS:HA	1:A:115:LYS:HD3	1.89	0.42
1:A:418:LEU:C	1:A:418:LEU:HD23	2.40	0.41
1:A:398:VAL:HG11	1:A:431:LEU:HD21	2.00	0.41
1:A:372:GLY:O	1:A:373:GLU:CB	2.68	0.41
1:A:398:VAL:HG21	1:A:431:LEU:HD11	2.02	0.41
1:A:448:LYS:HG3	1:A:449:GLY:N	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LYS:HB3	1:A:129:LYS:HE3	1.81	0.41
1:A:20:GLN:O	1:A:24:GLN:N	2.53	0.41
1:A:127:THR:HG21	1:A:293:ASN:OD1	2.21	0.40

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	455/472 (96%)	426 (94%)	24 (5%)	5 (1%)	17 5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	373	GLU
1	A	334	ALA
1	A	375	GLY
1	A	439	GLN
1	A	393	PRO

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	347/382 (91%)	346 (100%)	1 (0%)	94 94

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	65	ASN
1	A	73	ASN
1	A	191	HIS
1	A	272	ASN
1	A	329	ASN
1	A	355	GLN
1	A	361	ASN
1	A	380	GLN
1	A	424	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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
2	ATP	A	501	4	24,33,33	1.42	3 (12%)	31,52,52	2.90	8 (25%)
3	UDP	A	502	-	18,26,26	1.14	2 (11%)	26,40,40	1.80	2 (7%)
5	EDO	A	504	-	3,3,3	1.71	1 (33%)	2,2,2	1.36	0
5	EDO	A	505	-	3,3,3	1.69	1 (33%)	2,2,2	1.38	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	ATP	A	501	4	-	0/18/38/38	0/3/3/3
3	UDP	A	502	-	-	0/12/32/32	0/2/2/2
5	EDO	A	504	-	-	0/1/1/1	0/0/0/0
5	EDO	A	505	-	-	0/1/1/1	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	504	EDO	O2-C2	-2.33	1.29	1.42
5	A	505	EDO	O2-C2	-2.24	1.29	1.42
2	A	501	ATP	C8-N7	2.03	1.38	1.34
3	A	502	UDP	C6-N1	2.65	1.39	1.35
3	A	502	UDP	C4-N3	2.78	1.38	1.33
2	A	501	ATP	C5-C4	3.67	1.48	1.40
2	A	501	ATP	O4'-C1'	3.74	1.45	1.41

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ATP	N3-C2-N1	-11.57	120.04	128.89
2	A	501	ATP	C1'-N9-C4	-6.24	117.53	126.94
2	A	501	ATP	PB-O3B-PG	-3.69	120.29	132.67
2	A	501	ATP	O4'-C1'-N9	-2.68	102.48	108.10
3	A	502	UDP	PA-O3A-PB	-2.53	124.18	132.67
2	A	501	ATP	O2B-PB-O1B	2.08	123.80	112.53
2	A	501	ATP	O2B-PB-O3B	2.20	115.06	105.09
2	A	501	ATP	C4'-O4'-C1'	2.68	112.66	109.72
2	A	501	ATP	C2-N1-C6	5.24	128.14	118.77
3	A	502	UDP	C4-N3-C2	8.28	122.34	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	457/472 (96%)	0.86	90 (19%) 1 1	14, 28, 100, 109	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	399	ALA	10.8
1	A	413	LEU	10.7
1	A	374	LEU	10.2
1	A	431	LEU	9.9
1	A	412	GLY	8.2
1	A	392	LEU	8.1
1	A	379	TRP	7.6
1	A	414	HIS	7.3
1	A	203	LEU	7.1
1	A	415	SER	7.0
1	A	376	ASP	7.0
1	A	375	GLY	6.8
1	A	411	ALA	6.7
1	A	416	THR	5.9
1	A	418	LEU	5.8
1	A	402	GLN	5.8
1	A	385	LEU	5.7
1	A	390	ALA	5.6
1	A	383	HIS	5.6
1	A	365	VAL	5.4
1	A	445	PHE	5.2
1	A	435	ILE	5.2
1	A	438	HIS	5.0
1	A	419	LYS	4.9
1	A	406	ALA	4.9
1	A	367	VAL	4.9
1	A	403	PHE	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	387	ARG	4.5
1	A	463	MET	4.5
1	A	204	GLY	4.4
1	A	434	LEU	4.3
1	A	461	ASP	4.2
1	A	358	LEU	4.2
1	A	396	HIS	4.2
1	A	394	LEU	4.2
1	A	404	ALA	4.1
1	A	202	HIS	4.1
1	A	442	SER	4.1
1	A	429	PRO	4.0
1	A	206	PHE	4.0
1	A	440	PRO	3.9
1	A	401	GLY	3.9
1	A	334	ALA	3.8
1	A	338	PHE	3.8
1	A	364	LYS	3.8
1	A	348	THR	3.8
1	A	351	ARG	3.8
1	A	398	VAL	3.7
1	A	372	GLY	3.6
1	A	395	ASP	3.6
1	A	393	PRO	3.6
1	A	388	ASP	3.5
1	A	433	ASN	3.4
1	A	432	ILE	3.4
1	A	386	GLY	3.4
1	A	428	LEU	3.3
1	A	443	MET	3.3
1	A	464	GLU	3.2
1	A	457	THR	3.2
1	A	437	THR	3.1
1	A	447	PHE	3.1
1	A	371	ILE	3.1
1	A	400	VAL	3.0
1	A	446	LEU	3.0
1	A	9	VAL	3.0
1	A	405	SER	3.0
1	A	408	LEU	3.0
1	A	465	LYS	2.8
1	A	424	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	439	GLN	2.8
1	A	362	GLY	2.8
1	A	10	PRO	2.8
1	A	366	MET	2.7
1	A	391	GLU	2.7
1	A	363	ILE	2.7
1	A	459	MET	2.7
1	A	332	GLN	2.6
1	A	205	GLU	2.5
1	A	462	LEU	2.5
1	A	201	ALA	2.4
1	A	382	HIS	2.4
1	A	454	HIS	2.4
1	A	417	LYS	2.4
1	A	444	SER	2.4
1	A	389	LEU	2.4
1	A	407	ALA	2.3
1	A	380	GLN	2.3
1	A	377	SER	2.2
1	A	357	LEU	2.2
1	A	436	GLN	2.2

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	MG	A	503	1/1	0.72	0.14	2.83	51,51,51,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	A	505	4/4	0.91	0.14	2.24	41,42,42,42	0
3	UDP	A	502	25/25	0.87	0.14	1.52	39,48,66,66	0
5	EDO	A	504	4/4	0.96	0.13	1.15	24,26,26,27	0
2	ATP	A	501	31/31	0.94	0.11	-0.12	28,36,40,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.