



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 18, 2018 – 10:13 AM EST

PDB ID : 6GUJ  
Title : Molybdenum storage protein with two occupied ATP binding sites  
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Deposited on : 2018-06-19  
Resolution : 2.10 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : rb-20031633  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20031633

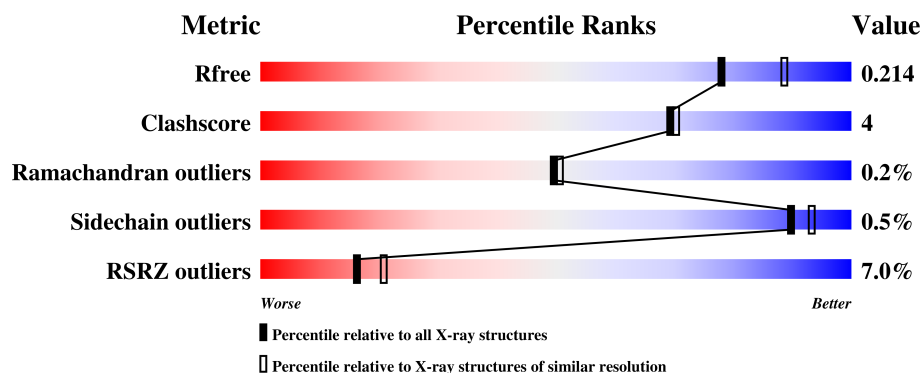
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	111664	4608 (2.10-2.10)
Clashscore	122126	5109 (2.10-2.10)
Ramachandran outliers	120053	5059 (2.10-2.10)
Sidechain outliers	120020	5060 (2.10-2.10)
RSRZ outliers	108989	4497 (2.10-2.10)

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  $\geq 3$ , 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	B	269	<div> <div>9%</div> <div>89%</div> <div>10%</div> </div>
2	A	275	<div> <div>4%</div> <div>84%</div> <div>5%</div> <div>11%</div> </div>

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4351 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.

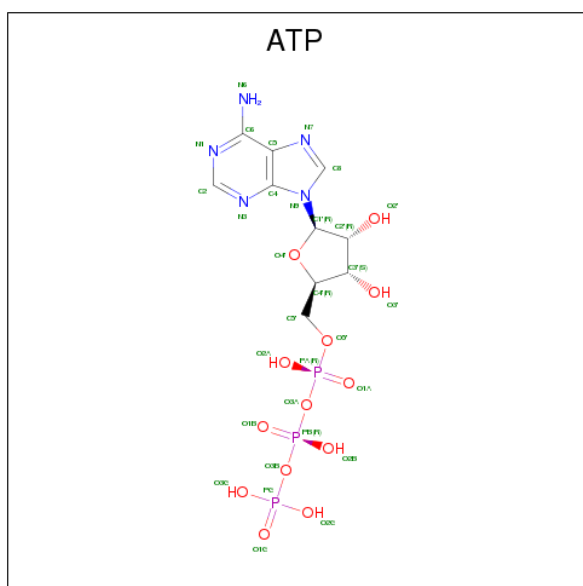
- Molecule 1 is a protein called Molybdenum storage protein subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	268	Total	C	N	O	S	0	10	0
			2023	1287	354	374	8			

- Molecule 2 is a protein called Molybdenum storage protein subunit alpha.

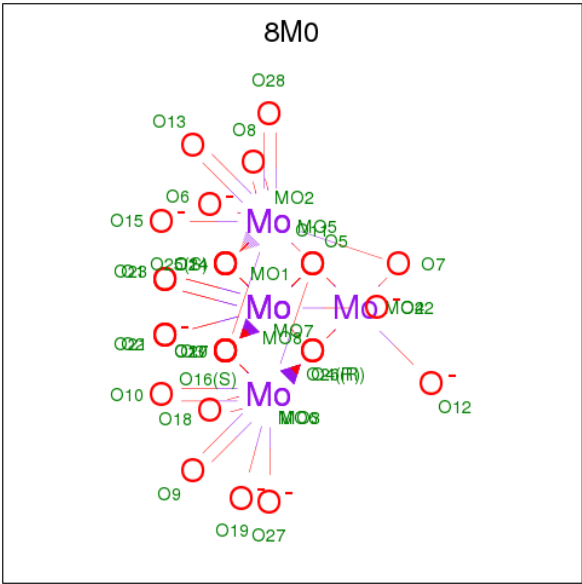
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	245	Total	C	N	O	S	0	4	0
			1847	1169	344	331	3			

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula:  $C_{10}H_{16}N_5O_{13}P_3$ ).



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

- Molecule 4 is bis(mu4-oxo)-tetrakis(mu3-oxo)-hexakis(mu2-oxo)-hexadeca-oxo-octamolybdenum (VI) (three-letter code: 8M0) (formula: Mo<sub>8</sub>O<sub>28</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	Mo	O	0	0
			36	8	28		
4	A	1	Total	Mo	O	0	0
			34	8	26		

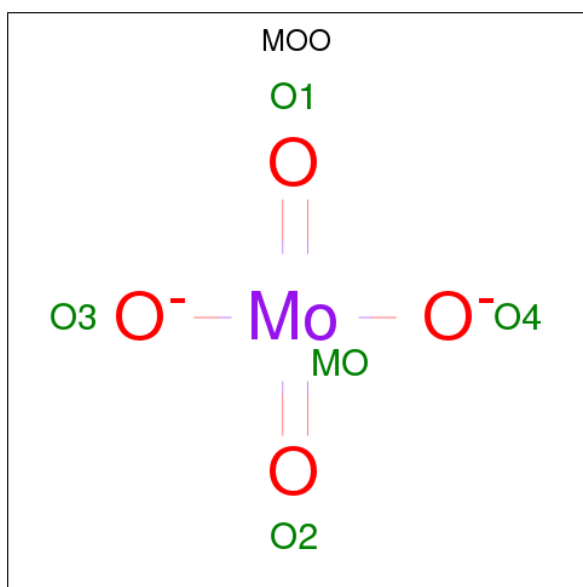
- Molecule 5 is MOLYBDENUM ATOM (three-letter code: MO) (formula: Mo).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	4	Total	Mo	0	0
			4	4		
5	A	3	Total	Mo	0	0
			3	3		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		

- Molecule 7 is MOLYBDATE ION (three-letter code: MOO) (formula: MoO<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	Mo	O	0	0
			5	1	4		

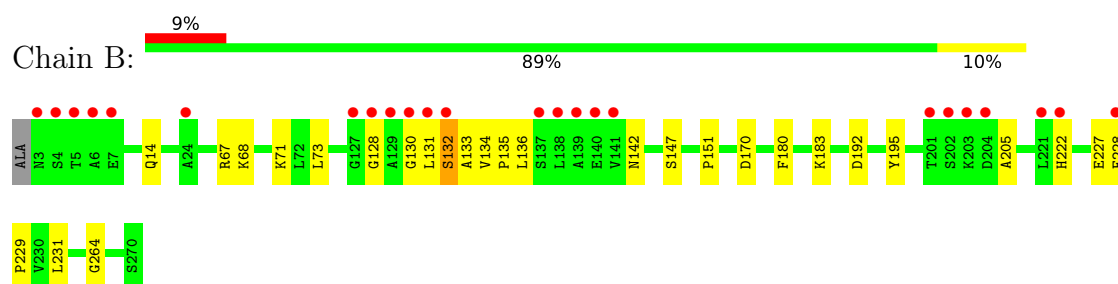
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	153	Total	O	0	0
			153	153		
8	A	183	Total	O	0	0
			183	183		

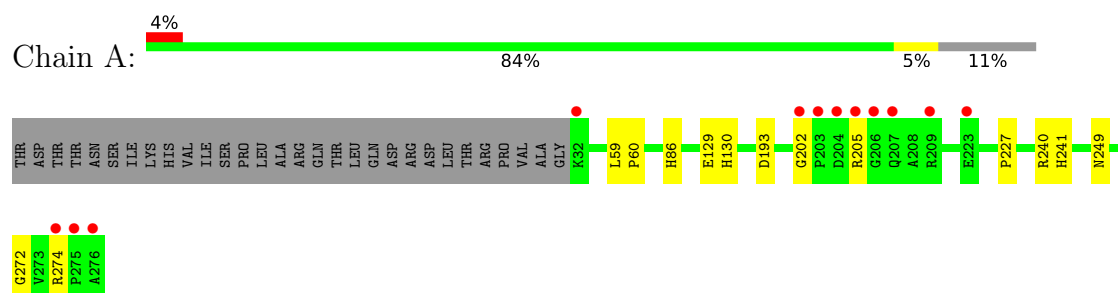
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Molybdenum storage protein subunit beta



- Molecule 2: Molybdenum storage protein subunit alpha



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	115.15Å 115.15Å 234.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.77 – 2.10 48.77 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.77-2.10) 99.2 (48.77-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.61 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, $R_{free}$	0.179 , 0.214 0.179 , 0.214	Depositor DCC
$R_{free}$ test set	2733 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.7	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 58.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4351	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 8M0, MG, MOO, MO, ATP

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	B	0.32	0/2091	0.56	0/2839
2	A	0.33	0/1899	0.52	0/2587
All	All	0.33	0/3990	0.54	0/5426

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	274	ARG	Peptide

### 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	B	2023	0	2105	19	0
2	A	1847	0	1900	11	0
3	A	31	0	12	1	0
3	B	31	0	12	0	0
4	A	34	0	0	4	0
4	B	36	0	0	2	0
5	A	3	0	0	0	0
5	B	4	0	0	0	0
6	A	1	0	0	0	0
7	A	5	0	0	0	0
8	A	183	0	0	4	0
8	B	153	0	0	3	0
All	All	4351	0	4029	32	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 4.

All (32) 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:B:73:LEU:HD21	1:B:136:LEU:HD11	1.68	0.75
1:B:183:LYS:NZ	8:B:401:HOH:O	2.25	0.69
1:B:131:LEU:O	1:B:133:ALA:N	2.31	0.63
2:A:130:HIS:HB2	4:A:303:8M0:O21	1.99	0.62
4:A:303:8M0:O6	4:A:303:8M0:MO2	1.72	0.60
4:B:302:8M0:O8	4:B:302:8M0:MO2	1.73	0.60
1:B:151:PRO:HB3	4:B:302:8M0:O4	2.04	0.57
1:B:134:VAL:HG12	1:B:135:PRO:HD3	1.86	0.57
1:B:68:LYS:NZ	8:B:403:HOH:O	2.38	0.54
1:B:71:LYS:NZ	1:B:142:ASN:OD1	2.43	0.52
1:B:228:PHE:CD2	1:B:229:PRO:HD3	2.44	0.52
1:B:222:HIS:HD2	1:B:231:LEU:HD12	1.75	0.51
1:B:130:GLY:HA3	1:B:180:PHE:CZ	2.47	0.50
2:A:272:GLY:N	8:A:405:HOH:O	2.45	0.49
1:B:132:SER:O	1:B:135:PRO:HD2	2.13	0.48
1:B:228:PHE:CG	1:B:229:PRO:HD3	2.49	0.48
1:B:128:GLY:O	1:B:132:SER:N	2.48	0.47
1:B:130:GLY:HA3	1:B:180:PHE:CE2	2.50	0.47
2:A:241:HIS:HD2	8:A:485:HOH:O	1.98	0.47
2:A:240:ARG:NH1	8:A:403:HOH:O	2.33	0.46
2:A:227:PRO:HD3	3:A:301:ATP:C6	2.51	0.45
2:A:129:GLU:OE1	4:A:303:8M0:O27	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:TYR:HB3	1:B:205:ALA:HB1	1.99	0.44
2:A:129:GLU:CD	4:A:303:8M0:O27	2.55	0.44
2:A:202:GLY:O	2:A:205:ARG:NH1	2.52	0.42
1:B:192:ASP:O	1:B:264:GLY:HA2	2.19	0.42
2:A:193:ASP:HA	2:A:249:ASN:HB2	2.01	0.41
1:B:170:ASP:HB2	1:B:227:GLU:CD	2.40	0.41
1:B:14[A]:GLN:NE2	8:B:410:HOH:O	2.53	0.41
2:A:86:HIS:HE1	8:A:565:HOH:O	2.03	0.41
1:B:222:HIS:CD2	1:B:231:LEU:HD12	2.55	0.41
2:A:59:LEU:HB2	2:A:60:PRO:HD3	2.02	0.41

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	B	276/269 (103%)	265 (96%)	10 (4%)	1 (0%)	36	34
2	A	247/275 (90%)	240 (97%)	7 (3%)	0	100	100
All	All	523/544 (96%)	505 (97%)	17 (3%)	1 (0%)	49	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	132	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	B	215/205 (105%)	213 (99%)	2 (1%)	81	85
2	A	192/215 (89%)	192 (100%)	0	100	100
All	All	407/420 (97%)	405 (100%)	2 (0%)	90	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	67	ARG
1	B	147	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	23	GLN
1	B	222	HIS

### 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 13 ligands modelled in this entry, 8 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
3	ATP	A	301	6	27,33,33	1.03	1 (3%)	27,52,52	2.03	5 (18%)
4	8M0	A	303	-	18,46,48	7.03	9 (50%)	0,140,150	0.00	-
7	MOO	A	304	-	2,4,4	9.32	1 (50%)	0,6,6	0.00	-
3	ATP	B	301	-	27,33,33	1.11	2 (7%)	27,52,52	1.81	5 (18%)
4	8M0	B	302	-	22,48,48	8.04	12 (54%)	0,150,150	0.00	-

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
3	ATP	A	301	6	-	0/18/38/38	0/3/3/3
4	8M0	A	303	-	-	0/0/290/300	0/0/13/13
7	MOO	A	304	-	-	0/0/0/0	0/0/0/0
3	ATP	B	301	-	-	0/18/38/38	0/3/3/3
4	8M0	B	302	-	-	0/0/300/300	0/0/13/13

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	ATP	PG-O3B	2.07	1.63	1.60
4	B	302	8M0	O8-MO2	2.72	1.73	1.68
4	A	303	8M0	O3-MO3	2.98	2.00	1.89
3	A	301	ATP	C5-C4	3.24	1.47	1.40
3	B	301	ATP	C5-C4	3.35	1.48	1.40
4	A	303	8M0	O8-MO2	3.87	1.75	1.68
4	B	302	8M0	O13-MO5	4.53	1.77	1.71
4	A	303	8M0	O13-MO5	4.53	1.77	1.67
4	A	303	8M0	O18-MO6	6.58	1.80	1.68
4	B	302	8M0	O25-MO8	6.80	2.14	1.88
4	B	302	8M0	O18-MO6	6.88	1.80	1.68
4	B	302	8M0	O26-MO4	7.98	2.18	1.88
4	B	302	8M0	O28-MO5	8.19	1.81	1.71
4	A	303	8M0	O1-MO1	8.94	1.84	1.68
4	B	302	8M0	O1-MO1	9.93	1.86	1.68
4	A	303	8M0	O9-MO3	10.72	1.90	1.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	302	8M0	O23-MO7	11.06	1.88	1.68
4	A	303	8M0	O23-MO7	11.27	1.88	1.68
4	B	302	8M0	O10-MO3	12.46	1.87	1.71
7	A	304	MOO	O2-MO	13.17	2.01	1.73
4	A	303	8M0	O7-MO4	15.22	1.89	1.74
4	B	302	8M0	O9-MO3	15.22	1.90	1.71
4	A	303	8M0	O20-MO8	15.53	1.89	1.74
4	B	302	8M0	O20-MO8	15.96	1.89	1.74
4	B	302	8M0	O7-MO4	17.39	1.91	1.74

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	ATP	N3-C2-N1	-6.24	123.52	128.86
3	B	301	ATP	N3-C2-N1	-6.19	123.56	128.86
3	A	301	ATP	C4'-O4'-C1'	-5.33	104.27	109.83
3	B	301	ATP	PB-O3B-PG	-3.09	122.24	132.63
3	A	301	ATP	O5'-C5'-C4'	-2.97	98.66	109.00
3	B	301	ATP	C4-C5-N7	-2.66	106.84	109.41
3	B	301	ATP	C4'-O4'-C1'	-2.64	107.08	109.83
3	B	301	ATP	PA-O3A-PB	-2.50	124.22	132.63
3	A	301	ATP	C4-C5-N7	-2.40	107.09	109.41
3	A	301	ATP	PB-O3B-PG	-2.14	125.44	132.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	ATP	1	0
4	A	303	8M0	4	0
4	B	302	8M0	2	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	268/269 (99%)	0.28	24 (8%) 9 12	17, 26, 60, 101	0
2	A	245/275 (89%)	-0.08	12 (4%) 29 36	14, 20, 49, 86	0
All	All	513/544 (94%)	0.11	36 (7%) 16 21	14, 23, 55, 101	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	128	GLY	9.5
1	B	3	ASN	9.3
1	B	130	GLY	8.6
1	B	131	LEU	7.2
2	A	276	ALA	5.8
1	B	139	ALA	5.6
2	A	32	LYS	5.3
1	B	5	THR	4.8
1	B	129	ALA	4.7
2	A	275	PRO	4.6
1	B	203	LYS	4.4
1	B	140	GLU	4.2
1	B	4	SER	3.7
1	B	127	GLY	3.6
1	B	138	LEU	3.6
1	B	6	ALA	3.5
2	A	274	ARG	3.3
2	A	203	PRO	3.3
1	B	137	SER	3.3
1	B	204	ASP	3.2
1	B	202	SER	2.9
2	A	207	GLN	2.9
2	A	209	ARG	2.7
2	A	206	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	7	GLU	2.6
1	B	132	SER	2.6
2	A	202	GLY	2.6
2	A	205	ARG	2.5
1	B	228	PHE	2.5
1	B	141	VAL	2.4
1	B	24	ALA	2.3
2	A	204	ASP	2.3
1	B	221	LEU	2.2
1	B	222	HIS	2.2
2	A	223	GLU	2.1
1	B	201	THR	2.1

## 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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
6	MG	A	302	1/1	0.91	0.52	43,43,43,43	0
5	MO	B	306	1/1	0.94	0.10	85,85,85,85	1
5	MO	A	306	1/1	0.95	0.32	59,59,59,59	1
3	ATP	B	301	31/31	0.96	0.16	26,42,50,52	19
5	MO	A	305	1/1	0.96	0.25	52,52,52,52	1
5	MO	B	304	1/1	0.96	0.10	56,56,56,56	1
4	8M0	B	302	36/36	0.97	0.16	24,31,43,52	34
5	MO	A	307	1/1	0.97	0.20	73,73,73,73	1
5	MO	B	305	1/1	0.98	0.10	51,51,51,51	1
3	ATP	A	301	31/31	0.98	0.13	17,25,34,34	0
5	MO	B	303	1/1	0.98	0.09	54,54,54,54	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
7	MOO	A	304	5/5	0.99	0.15	12,14,17,39	4
4	8M0	A	303	34/36	0.99	0.13	13,24,29,51	34

## 6.5 Other polymers [i](#)

There are no such residues in this entry.