



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:04 AM GMT

PDB ID : 2FGK
Title : Crystal structure of the ABC-cassette E631Q mutant of HlyB with bound ATP
Authors : Zaitseva, J.; Oswald, C.; Jumpertz, T.; Jenewein, S.; Holland, I.B.; Schmitt, L.
Deposited on : 2005-12-22
Resolution : 2.70 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
<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

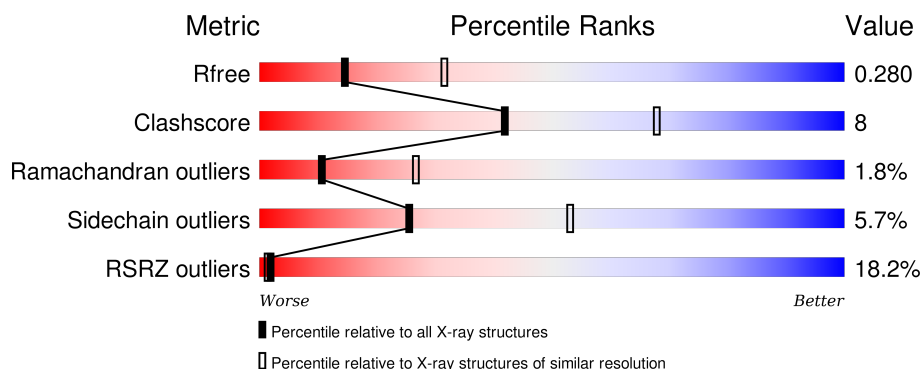
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	<div> <div>20%</div> <div>73%</div> <div>21%</div> <div>• •</div> </div>
1	B	247	<div> <div>16%</div> <div>73%</div> <div>24%</div> <div>• •</div> </div>
1	C	247	<div> <div>17%</div> <div>73%</div> <div>22%</div> <div>• •</div> </div>
1	D	247	<div> <div>17%</div> <div>74%</div> <div>23%</div> <div>• •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7848 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 Alpha-hemolysin translocation ATP-binding protein hlyB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	B	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	C	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			
1	D	241	Total	C	N	O	S	0	0	0
			1893	1194	342	352	5			

There are 28 discrepancies between the modelled and reference sequences:

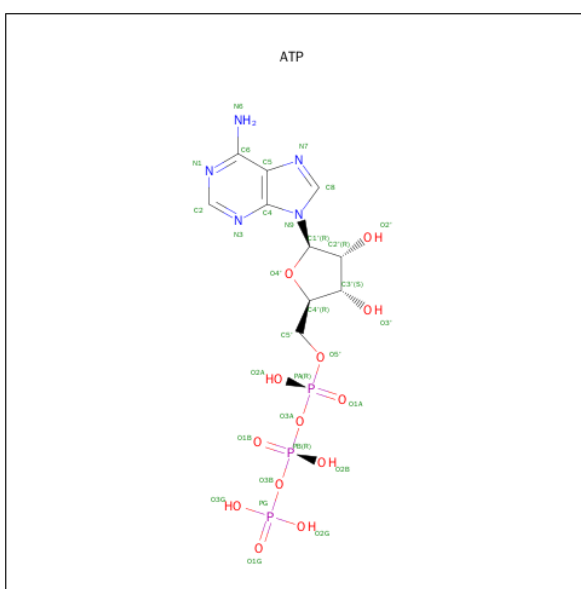
Chain	Residue	Modelled	Actual	Comment	Reference
A	461	HIS	-	EXPRESSION TAG	UNP P08716
A	462	HIS	-	EXPRESSION TAG	UNP P08716
A	463	HIS	-	EXPRESSION TAG	UNP P08716
A	464	HIS	-	EXPRESSION TAG	UNP P08716
A	465	HIS	-	EXPRESSION TAG	UNP P08716
A	466	HIS	-	EXPRESSION TAG	UNP P08716
A	631	GLN	GLU	ENGINEERED	UNP P08716
B	461	HIS	-	EXPRESSION TAG	UNP P08716
B	462	HIS	-	EXPRESSION TAG	UNP P08716
B	463	HIS	-	EXPRESSION TAG	UNP P08716
B	464	HIS	-	EXPRESSION TAG	UNP P08716
B	465	HIS	-	EXPRESSION TAG	UNP P08716
B	466	HIS	-	EXPRESSION TAG	UNP P08716
B	631	GLN	GLU	ENGINEERED	UNP P08716
C	461	HIS	-	EXPRESSION TAG	UNP P08716
C	462	HIS	-	EXPRESSION TAG	UNP P08716
C	463	HIS	-	EXPRESSION TAG	UNP P08716
C	464	HIS	-	EXPRESSION TAG	UNP P08716
C	465	HIS	-	EXPRESSION TAG	UNP P08716
C	466	HIS	-	EXPRESSION TAG	UNP P08716
C	631	GLN	GLU	ENGINEERED	UNP P08716

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Chain	Residue	Modelled	Actual	Comment	Reference
D	461	HIS	-	EXPRESSION TAG	UNP P08716
D	462	HIS	-	EXPRESSION TAG	UNP P08716
D	463	HIS	-	EXPRESSION TAG	UNP P08716
D	464	HIS	-	EXPRESSION TAG	UNP P08716
D	465	HIS	-	EXPRESSION TAG	UNP P08716
D	466	HIS	-	EXPRESSION TAG	UNP P08716
D	631	GLN	GLU	ENGINEERED	UNP P08716

- 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		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	36	Total	O	0	0
			36	36		

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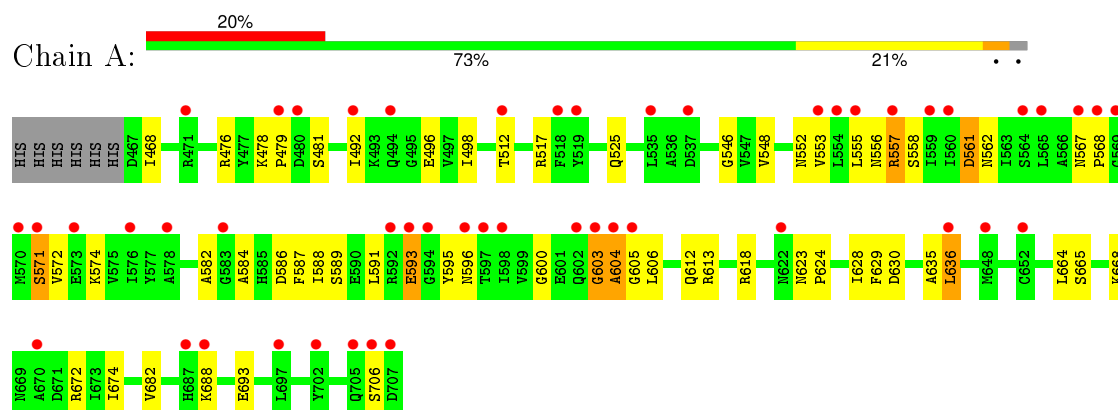
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	44	Total 44	O 44	0	0
3	C	32	Total 32	O 32	0	0
3	D	40	Total 40	O 40	0	0

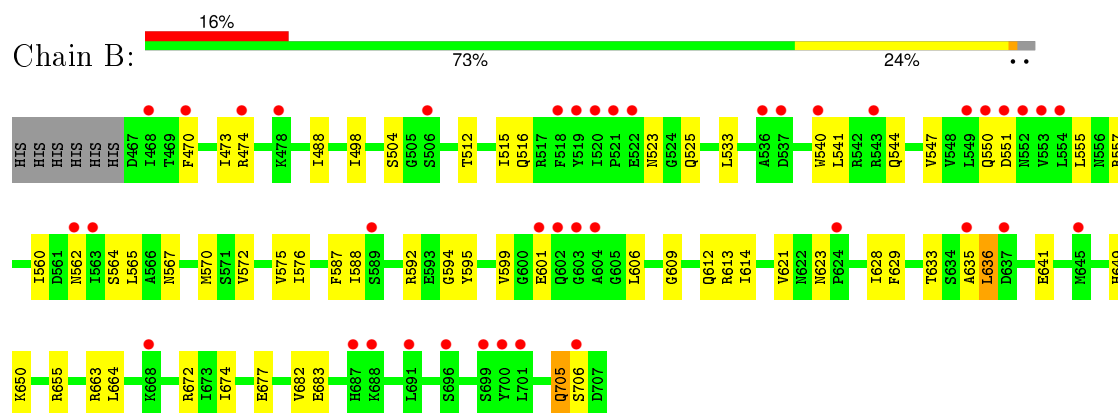
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

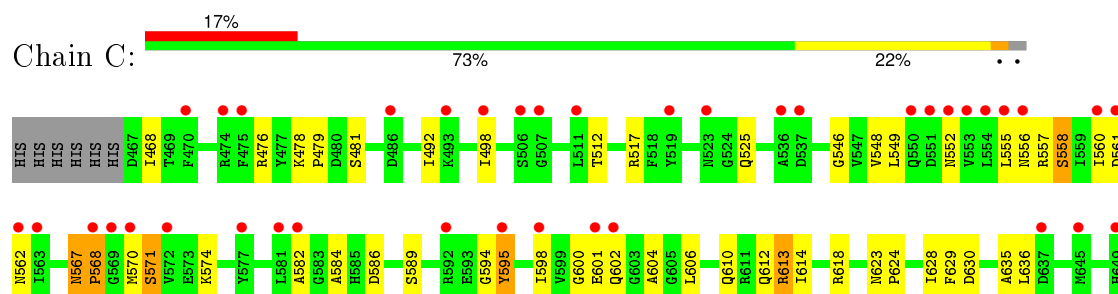
- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB

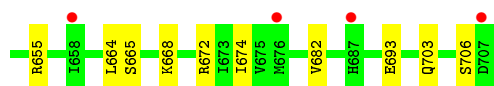


- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB

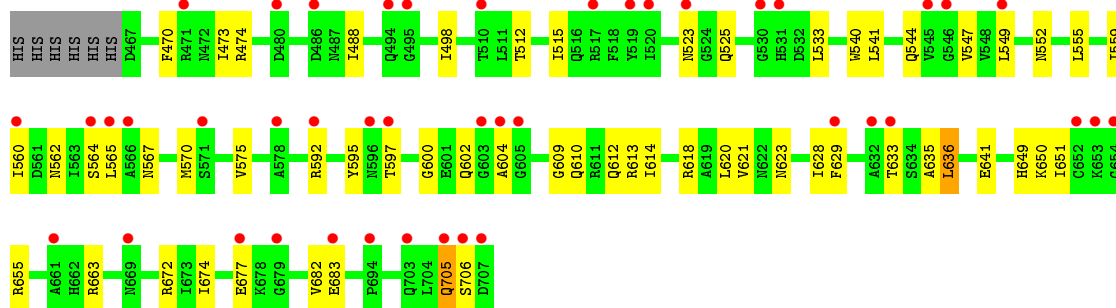
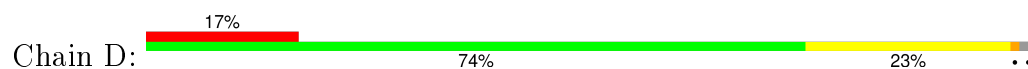


- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB





- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.14Å 189.26Å 63.48Å 90.00° 111.87° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 20.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (20.00-2.70) 98.8 (20.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.223 , 0.280 0.222 , 0.280	Depositor DCC
R_{free} test set	1406 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.9	EDS
Estimated twinning fraction	0.406 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 27971 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7848	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	A	1.27	8/1920 (0.4%)	0.68	4/2589 (0.2%)
1	B	0.41	0/1920	0.58	0/2589
1	C	0.73	8/1920 (0.4%)	0.67	4/2589 (0.2%)
1	D	0.41	0/1920	0.57	0/2589
All	All	0.79	16/7680 (0.2%)	0.63	8/10356 (0.1%)

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	593	GLU	CD-OE2	43.26	1.73	1.25
1	A	593	GLU	CD-OE1	18.30	1.45	1.25
1	C	558	SER	CB-OG	14.44	1.61	1.42
1	A	595	TYR	CE2-CZ	13.03	1.55	1.38
1	C	595	TYR	CE2-CZ	12.90	1.55	1.38

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	557	ARG	NE-CZ-NH2	-10.21	115.19	120.30
1	C	557	ARG	NE-CZ-NH2	-9.98	115.31	120.30
1	A	593	GLU	OE1-CD-OE2	9.53	134.74	123.30
1	C	595	TYR	CB-CG-CD2	-6.88	116.87	121.00
1	A	595	TYR	CB-CG-CD2	-6.80	116.92	121.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1893	0	1949	29	0
1	B	1893	0	1949	36	0
1	C	1893	0	1949	31	0
1	D	1893	0	1949	37	0
2	A	31	0	12	0	0
2	B	31	0	12	1	0
2	C	31	0	12	0	0
2	D	31	0	12	0	0
3	A	36	0	0	5	0
3	B	44	0	0	2	0
3	C	32	0	0	3	0
3	D	40	0	0	1	0
All	All	7848	0	7844	130	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 8.

The worst 5 of 130 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:593:GLU:CD	1:A:593:GLU:OE2	1.73	1.25
1:C:571:SER:HB2	1:C:574:LYS:HB2	1.36	1.06
1:B:612:GLN:HE22	1:B:635:ALA:H	1.13	0.90
1:C:560:ILE:HB	3:C:139:HOH:O	1.73	0.88
1:C:612:GLN:HE22	1:C:635:ALA:H	1.27	0.83

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	239/247 (97%)	223 (93%)	10 (4%)	6 (2%)	7	18
1	B	239/247 (97%)	226 (95%)	10 (4%)	3 (1%)	15	37
1	C	239/247 (97%)	222 (93%)	11 (5%)	6 (2%)	7	18
1	D	239/247 (97%)	224 (94%)	13 (5%)	2 (1%)	24	51
All	All	956/988 (97%)	895 (94%)	44 (5%)	17 (2%)	11	27

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	603	GLY
1	C	601	GLU
1	C	571	SER
1	D	565	LEU
1	A	604	ALA

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	207/213 (97%)	192 (93%)	15 (7%)	18	41
1	B	207/213 (97%)	196 (95%)	11 (5%)	28	57
1	C	207/213 (97%)	194 (94%)	13 (6%)	22	48
1	D	207/213 (97%)	199 (96%)	8 (4%)	39	70
All	All	828/852 (97%)	781 (94%)	47 (6%)	25	53

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	629	PHE
1	C	517	ARG
1	D	629	PHE

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Mol	Chain	Res	Type
1	B	664	LEU
1	C	525	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	705	GLN
1	C	612	GLN
1	D	623	ASN
1	C	539	ASN
1	C	623	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	ATP	A	800	-	24,33,33	1.17	2 (8%)	31,52,52	2.16	7 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ATP	B	801	-	24,33,33	1.22	2 (8%)	31,52,52	2.09	5 (16%)
2	ATP	C	802	-	24,33,33	0.97	1 (4%)	31,52,52	2.29	7 (22%)
2	ATP	D	803	-	24,33,33	1.24	2 (8%)	31,52,52	2.07	5 (16%)

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	800	-	-	0/18/38/38	0/3/3/3
2	ATP	B	801	-	-	0/18/38/38	0/3/3/3
2	ATP	C	802	-	-	0/18/38/38	0/3/3/3
2	ATP	D	803	-	-	0/18/38/38	0/3/3/3

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	802	ATP	PG-O1G	2.03	1.57	1.51
2	B	801	ATP	O4'-C1'	2.86	1.44	1.41
2	A	800	ATP	PG-O1G	2.88	1.60	1.51
2	A	800	ATP	O4'-C1'	3.04	1.45	1.41
2	D	803	ATP	PG-O1G	3.09	1.61	1.51

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	802	ATP	N3-C2-N1	-9.56	121.57	128.89
2	A	800	ATP	N3-C2-N1	-9.13	121.91	128.89
2	D	803	ATP	N3-C2-N1	-8.83	122.13	128.89
2	B	801	ATP	N3-C2-N1	-8.74	122.20	128.89
2	B	801	ATP	PA-O3A-PB	-4.04	121.39	132.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	ATP	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	241/247 (97%)	1.30	49 (20%) 1 1	39, 51, 62, 71	0
1	B	241/247 (97%)	1.14	40 (16%) 2 2	40, 51, 63, 70	0
1	C	241/247 (97%)	1.29	43 (17%) 2 1	37, 50, 61, 71	0
1	D	241/247 (97%)	1.19	43 (17%) 2 1	40, 51, 63, 71	0
All	All	964/988 (97%)	1.23	175 (18%) 2 1	37, 51, 63, 71	0

The worst 5 of 175 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	604	ALA	14.5
1	D	707	ASP	10.7
1	C	551	ASP	9.2
1	B	550	GLN	9.2
1	A	569	GLY	8.4

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 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. 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(\AA^2)	Q<0.9
2	ATP	D	803	31/31	0.92	0.24	0.09	38,46,47,50	0
2	ATP	C	802	31/31	0.92	0.17	-0.65	28,40,43,44	0
2	ATP	A	800	31/31	0.91	0.19	-0.77	32,41,44,44	0
2	ATP	B	801	31/31	0.93	0.16	-0.89	37,42,44,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.