Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:58 AM GMT

PDB ID	:	2JK8
Title	:	TYPE IV SECRETION SYSTEM EFFECTOR PROTEIN BEPA COMPLEXED WITH A PYROPHOSPHATE MOIETY
Authors	:	Palanivelu, D.V.; Schirmer, T.
Deposited on	:	2008-08-22
Resolution	:	2.80 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

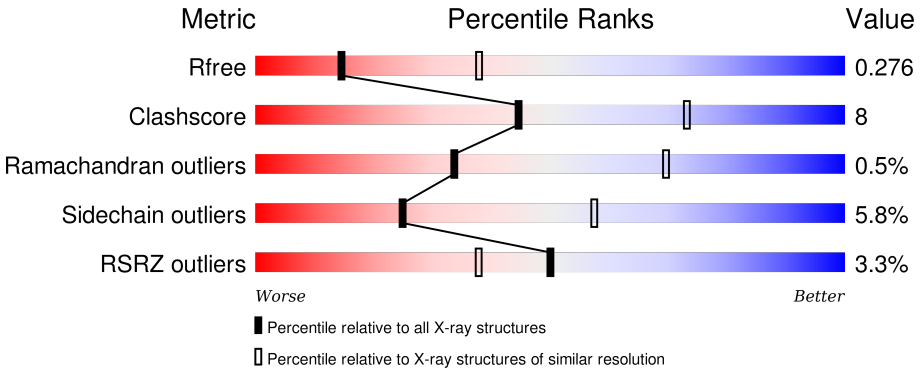
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 <=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	302	3% 76% 17%
1	B	302	3% 73% 20%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4690 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.

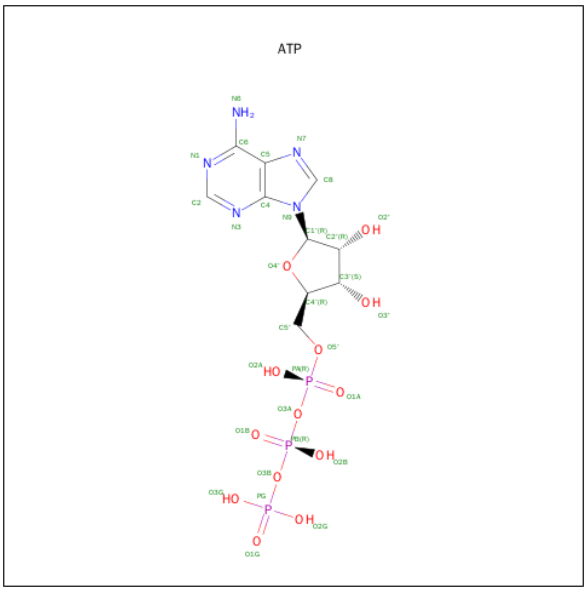
- Molecule 1 is a protein called PUTATIVE CELL FILAMENTATION PROTEIN (BEPA PROTEIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	112	0	0
			2334	1482	407	434	11			
1	B	291	Total	C	N	O	S	120	0	0
			2334	1482	407	434	11			

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Ni	0	0
			2	2		
2	A	2	Total	Ni	0	0
			2	2		

- 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	A	1	Total O P 9 7 2	0	0

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

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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Cl 1 1	0	0

- Molecule 6 is water.

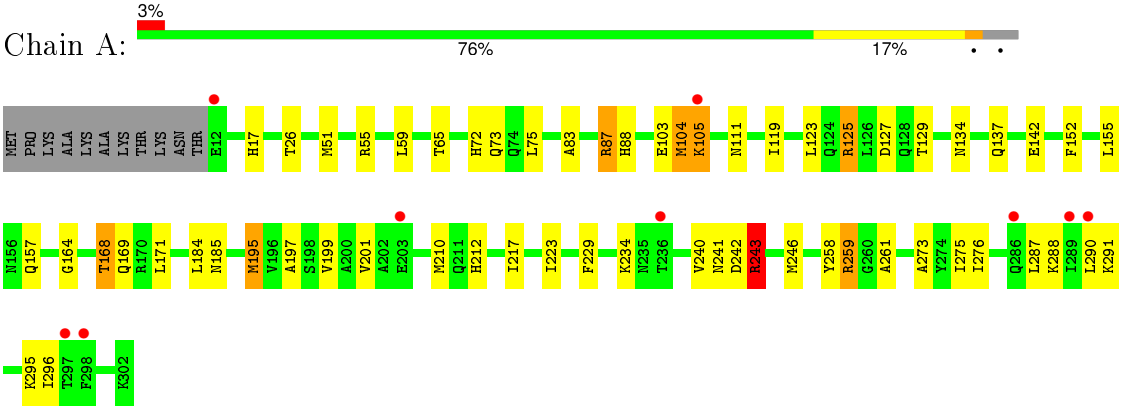
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	4	Total O 4 4	0	0
6	B	3	Total O 3 3	0	0

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.

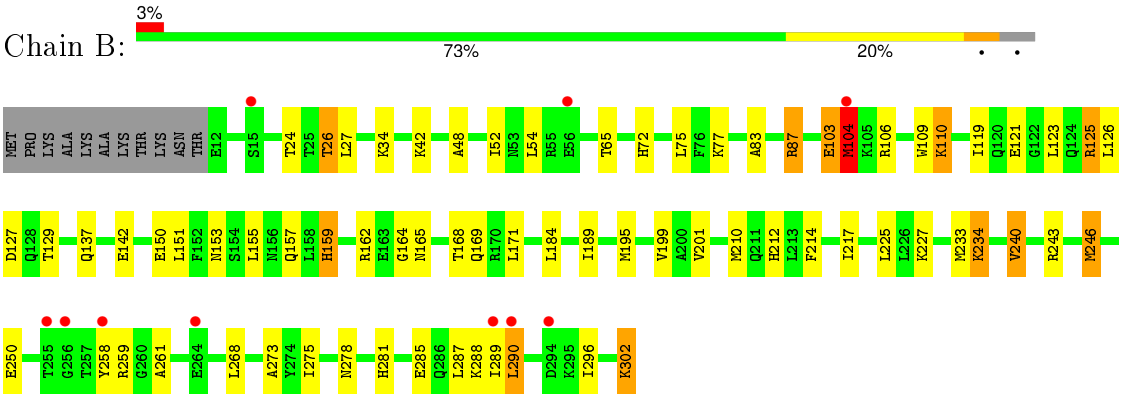
- Molecule 1: PUTATIVE CELL FILAMENTATION PROTEIN (BEPA PROTEIN)

Chain A:



- Molecule 1: PUTATIVE CELL FILAMENTATION PROTEIN (BEPA PROTEIN)

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	72.95Å 82.88Å 126.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 29.59 – 2.80	Depositor EDS
% Data completeness (in resolution range)	91.7 (30.00-2.80) 91.7 (29.59-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.10 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.216 , 0.275 0.224 , 0.276	Depositor DCC
R_{free} test set	890 reflections (5.24%)	DCC
Wilson B-factor (Å^2)	67.0	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 51.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 17866 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4690	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NI, MG, ATP, CL

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	0.80	8/2387 (0.3%)	0.76	8/3223 (0.2%)
1	B	1.04	13/2387 (0.5%)	0.86	10/3223 (0.3%)
All	All	0.93	21/4774 (0.4%)	0.81	18/6446 (0.3%)

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
1	A	0	1
1	B	0	1
All	All	0	2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	125	ARG	CD-NE	25.27	1.89	1.46
1	B	109	TRP	C-N	18.08	1.75	1.34
1	A	243	ARG	CG-CD	-16.52	1.10	1.51
1	B	159	HIS	CB-CG	-16.51	1.20	1.50
1	A	125	ARG	CD-NE	-14.53	1.21	1.46
1	B	104	MET	C-N	-12.14	1.06	1.34
1	B	234	LYS	CD-CE	11.31	1.79	1.51
1	B	103	GLU	CB-CG	10.32	1.71	1.52
1	A	111	ASN	C-N	-10.13	1.10	1.34
1	A	291	LYS	CB-CG	9.74	1.78	1.52
1	B	121	GLU	CG-CD	7.67	1.63	1.51
1	B	150	GLU	CB-CG	-7.65	1.37	1.52
1	A	240	VAL	CB-CG2	6.42	1.66	1.52
1	B	285	GLU	CB-CG	-6.19	1.40	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	105	LYS	C-N	-6.12	1.20	1.34
1	A	142	GLU	CG-CD	5.96	1.60	1.51
1	A	288	LYS	CB-CG	-5.81	1.36	1.52
1	B	289	ILE	CB-CG1	-5.57	1.38	1.54
1	B	42	LYS	CG-CD	-5.36	1.34	1.52
1	B	142	GLU	CG-CD	-5.28	1.44	1.51
1	B	110	LYS	CB-CG	-5.13	1.38	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	159	HIS	CA-CB-CG	16.59	141.80	113.60
1	B	302	LYS	CB-CG-CD	13.56	146.87	111.60
1	B	109	TRP	C-N-CA	-13.11	88.93	121.70
1	B	104	MET	O-C-N	-11.54	104.24	122.70
1	A	125	ARG	CG-CD-NE	10.30	133.44	111.80
1	A	243	ARG	CB-CG-CD	7.67	131.55	111.60
1	B	104	MET	CA-C-N	7.34	133.34	117.20
1	A	295	LYS	CB-CG-CD	-7.17	92.96	111.60
1	B	109	TRP	O-C-N	6.65	133.33	122.70
1	A	243	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	B	288	LYS	CB-CG-CD	-6.17	95.57	111.60
1	B	125	ARG	CG-CD-NE	-6.00	99.21	111.80
1	A	105	LYS	O-C-N	-5.73	113.53	122.70
1	A	291	LYS	CA-CB-CG	-5.46	101.39	113.40
1	A	142	GLU	CB-CG-CD	-5.31	99.86	114.20
1	A	259	ARG	CB-CG-CD	-5.30	97.81	111.60
1	B	42	LYS	CB-CG-CD	5.16	125.03	111.60
1	B	227	LYS	CB-CG-CD	5.07	124.78	111.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	LYS	Mainchain
1	B	104	MET	Mainchain

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	2334	0	2296	31	1
1	B	2334	0	2295	38	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	0	0	0
4	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	4	0	0	0	0
6	B	3	0	0	0	0
All	All	4690	0	4591	69	1

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.

All (69) 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:75:LEU:HD12	1:B:168:THR:HG22	1.53	0.91
1:A:164:GLY:O	1:A:168:THR:HG23	1.77	0.83
1:B:65:THR:HG23	1:B:127:ASP:OD1	1.83	0.78
1:A:103:GLU:O	1:A:104:MET:HB2	1.90	0.70
1:A:125:ARG:O	1:A:129:THR:HG23	1.92	0.70
1:B:290:LEU:HD21	1:B:296:ILE:HG21	1.74	0.68
1:B:225:LEU:HD21	1:B:278:ASN:HD22	1.61	0.65
1:A:258:TYR:OH	1:A:261:ALA:HB2	1.97	0.64
1:B:72:HIS:ND1	1:B:83:ALA:O	2.23	0.64
1:B:48:ALA:O	1:B:52:ILE:HD13	1.97	0.64
1:B:153:ASN:ND2	1:B:201:VAL:HG11	2.12	0.64
1:A:184:LEU:HA	1:A:217:ILE:HG22	1.81	0.62
1:A:65:THR:HG21	1:A:123:LEU:HB3	1.81	0.60
1:A:72:HIS:ND1	1:A:83:ALA:O	2.33	0.59
1:B:195:MET:O	1:B:199:VAL:HG23	2.04	0.58
1:B:126:LEU:CD1	1:B:151:LEU:HD22	2.34	0.58
1:A:155:LEU:HD23	1:A:169:GLN:HG3	1.85	0.57
1:B:87:ARG:NH1	1:B:119:ILE:HD11	2.19	0.57
1:A:134:ASN:O	1:A:137:GLN:HB2	2.05	0.56
1:B:258:TYR:OH	1:B:261:ALA:HB2	2.06	0.55
1:A:197:ALA:O	1:A:201:VAL:HG23	2.07	0.54
1:B:165:ASN:O	1:B:169:GLN:HB2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:THR:HG21	1:B:123:LEU:HB3	1.88	0.53
1:A:229:PHE:CG	1:A:276:ILE:HD11	2.43	0.53
1:B:126:LEU:HD12	1:B:151:LEU:HD22	1.89	0.53
1:B:26:THR:HG21	1:B:34:LYS:O	2.08	0.53
1:B:126:LEU:HD22	1:B:155:LEU:HD13	1.91	0.53
1:A:51:MET:CE	1:A:55:ARG:HG3	2.40	0.52
1:B:87:ARG:HD3	1:B:159:HIS:O	2.09	0.51
1:B:268:LEU:HD21	1:B:296:ILE:HD11	1.93	0.51
1:A:217:ILE:O	1:A:223:ILE:HD11	2.11	0.51
1:B:125:ARG:O	1:B:129:THR:HG23	2.10	0.50
1:B:210:MET:HE1	1:B:214:PHE:HE2	1.77	0.50
1:A:185:ASN:O	1:A:217:ILE:HD12	2.12	0.49
1:A:65:THR:HG23	1:A:127:ASP:OD1	2.13	0.48
1:B:240:VAL:HA	1:B:243:ARG:HG3	1.95	0.48
1:A:195:MET:O	1:A:199:VAL:HG23	2.14	0.48
1:A:246:MET:HE3	1:A:275:ILE:HD13	1.94	0.48
1:B:246:MET:HE3	1:B:275:ILE:HD13	1.96	0.47
1:B:184:LEU:HA	1:B:217:ILE:HG22	1.95	0.47
1:B:164:GLY:O	1:B:168:THR:HG23	2.13	0.47
1:B:65:THR:HG23	1:B:127:ASP:CG	2.35	0.47
1:A:290:LEU:HD21	1:A:296:ILE:HG21	1.96	0.47
1:B:246:MET:HE1	1:B:273:ALA:HB1	1.96	0.46
1:A:152:PHE:CD2	1:A:210:MET:HE1	2.51	0.46
1:B:250:GLU:HB3	1:B:302:LYS:HG2	1.98	0.46
1:B:233:MET:HE2	1:B:240:VAL:CG2	2.46	0.46
1:A:65:THR:HG23	1:A:127:ASP:CG	2.37	0.46
1:B:250:GLU:HB3	1:B:302:LYS:CG	2.46	0.45
1:A:197:ALA:HB1	1:A:210:MET:HG2	2.00	0.44
1:A:73:GLN:HA	1:A:83:ALA:HB1	1.99	0.44
1:A:87:ARG:NH1	1:A:119:ILE:HD11	2.32	0.44
1:B:233:MET:CE	1:B:240:VAL:CG2	2.97	0.43
1:A:75:LEU:HD12	1:A:168:THR:HG22	2.00	0.43
1:B:54:LEU:HD23	1:B:54:LEU:HA	1.88	0.43
1:B:278:ASN:HB3	1:B:281:HIS:CD2	2.54	0.43
1:A:88:HIS:HB3	1:A:119:ILE:HD12	2.00	0.43
1:A:290:LEU:HD11	1:A:296:ILE:CG2	2.49	0.42
1:A:246:MET:HE1	1:A:273:ALA:HB1	2.02	0.42
1:A:51:MET:HE3	1:A:55:ARG:HG3	2.01	0.42
1:B:233:MET:HE1	1:B:240:VAL:HG23	2.02	0.42
1:A:65:THR:HG23	1:A:127:ASP:OD2	2.20	0.42
1:B:189:ILE:HD11	1:B:217:ILE:HG12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:MET:CE	1:B:214:PHE:HE2	2.33	0.41
1:A:103:GLU:O	1:A:104:MET:CB	2.61	0.41
1:B:27:LEU:HD21	1:B:162:ARG:CZ	2.50	0.41
1:A:290:LEU:HD11	1:A:296:ILE:HG21	2.03	0.41
1:B:24:THR:HB	1:B:26:THR:HG23	2.03	0.41
1:B:189:ILE:HD11	1:B:217:ILE:CG1	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:LEU:O	1:A:243:ARG:NH2[4_466]	2.16	0.04

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	289/302 (96%)	279 (96%)	9 (3%)	1 (0%)	46 79
1	B	289/302 (96%)	279 (96%)	8 (3%)	2 (1%)	26 62
All	All	578/604 (96%)	558 (96%)	17 (3%)	3 (0%)	34 69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	MET
1	B	106	ARG
1	B	104	MET

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	248/257 (96%)	234 (94%)	14 (6%)	26 59
1	B	248/257 (96%)	233 (94%)	15 (6%)	24 56
All	All	496/514 (96%)	467 (94%)	29 (6%)	25 57

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	26	THR
1	A	87	ARG
1	A	157	GLN
1	A	168	THR
1	A	171	LEU
1	A	195	MET
1	A	212	HIS
1	A	234	LYS
1	A	241	ASN
1	A	242	ASP
1	A	243	ARG
1	A	259	ARG
1	A	287	LEU
1	B	26	THR
1	B	77	LYS
1	B	87	ARG
1	B	103	GLU
1	B	110	LYS
1	B	137	GLN
1	B	157	GLN
1	B	171	LEU
1	B	212	HIS
1	B	234	LYS
1	B	240	VAL
1	B	246	MET
1	B	259	ARG

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Mol	Chain	Res	Type
1	B	287	LEU
1	B	290	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	157	GLN
1	A	281	HIS
1	B	157	GLN
1	B	278	ASN
1	B	281	HIS

5.3.3 RNA

There are no RNA molecules 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

Of 7 ligands modelled in this entry, 6 are monoatomic - leaving 1 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	701	4	6,8,33	1.57	1 (16%)	11,13,52	1.26	1 (9%)

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	701	4	-	0/6/6/38	0/0/0/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	ATP	PG-O1G	3.59	1.63	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	701	ATP	O3G-PG-O2G	3.02	118.87	107.38

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, 95^{th} 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	285/302 (94%)	0.08	9 (3%) 51 39	54, 59, 67, 88	17 (5%)
1	B	286/302 (94%)	0.08	10 (3%) 48 35	46, 59, 65, 98	23 (8%)
All	All	571/604 (94%)	0.08	19 (3%) 50 38	46, 59, 65, 98	40 (7%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	104	MET	6.6
1	A	289	ILE	4.8
1	A	236	THR	3.6
1	B	258	TYR	3.5
1	B	289	ILE	3.2
1	A	203	GLU	3.1
1	B	56	GLU	3.1
1	B	264	GLU	3.1
1	A	286	GLN	3.0
1	B	290	LEU	3.0
1	A	105	LYS	2.8
1	A	298	PHE	2.8
1	A	297	THR	2.6
1	B	294	ASP	2.6
1	B	255	THR	2.2
1	A	290	LEU	2.1
1	A	12	GLU	2.1
1	B	15	SER	2.1
1	B	256	GLY	2.0

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 (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, 95^{th} 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
5	CL	B	601	1/1	0.96	0.24	0.07	71,71,71,71	0
3	ATP	A	701	9/31	0.85	0.18	-0.26	97,99,105,105	0
2	NI	A	502	1/1	0.97	0.10	-	59,59,59,59	0
2	NI	A	501	1/1	0.95	0.15	-	68,68,68,68	0
2	NI	B	502	1/1	0.96	0.13	-	64,64,64,64	0
2	NI	B	501	1/1	0.78	0.07	-	75,75,75,75	0
4	MG	A	702	1/1	0.83	0.20	-	97,97,97,97	0

6.5 Other polymers (i)

There are no such residues in this entry.