



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:56 PM GMT

PDB ID : 4B0T
Title : Structure of the Pup Ligase PafA of the Prokaryotic Ubiquitin-like Modification Pathway in Complex with ADP
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Deposited on : 2012-07-04
Resolution : 2.16 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

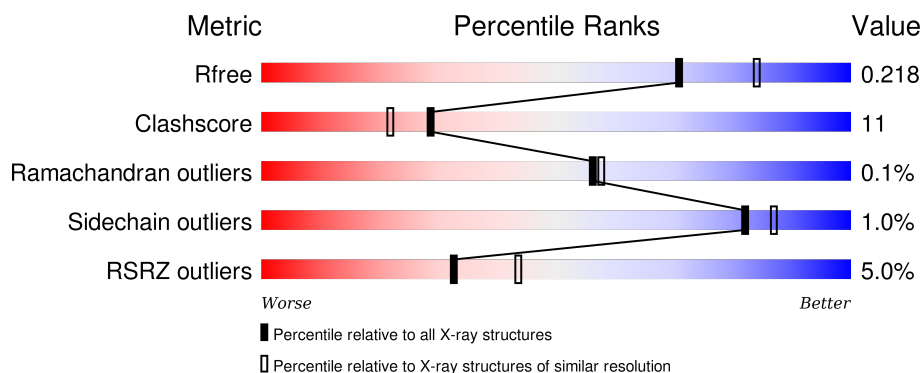
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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	493	<div> <div>4%</div> <div>77%</div> <div>17%</div> <div>• 5%</div> </div>
1	B	493	<div> <div>5%</div> <div>79%</div> <div>17%</div> <div>•</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7939 atoms, of which 34 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 PUP-PROTEIN LIGASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	1	0
			3688	2310	652	709	17			
1	B	472	Total	C	N	O	S	0	0	0
			3700	2318	653	712	17			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	483	LYS	-	EXPRESSION TAG	UNP Q8NQE1
A	484	SER	-	EXPRESSION TAG	UNP Q8NQE1
A	485	SER	-	EXPRESSION TAG	UNP Q8NQE1
A	486	VAL	-	EXPRESSION TAG	UNP Q8NQE1
A	487	GLU	-	EXPRESSION TAG	UNP Q8NQE1
A	488	HIS	-	EXPRESSION TAG	UNP Q8NQE1
A	489	HIS	-	EXPRESSION TAG	UNP Q8NQE1
A	490	HIS	-	EXPRESSION TAG	UNP Q8NQE1
A	491	HIS	-	EXPRESSION TAG	UNP Q8NQE1
A	492	HIS	-	EXPRESSION TAG	UNP Q8NQE1
A	493	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	483	LYS	-	EXPRESSION TAG	UNP Q8NQE1
B	484	SER	-	EXPRESSION TAG	UNP Q8NQE1
B	485	SER	-	EXPRESSION TAG	UNP Q8NQE1
B	486	VAL	-	EXPRESSION TAG	UNP Q8NQE1
B	487	GLU	-	EXPRESSION TAG	UNP Q8NQE1
B	488	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	489	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	490	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	491	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	492	HIS	-	EXPRESSION TAG	UNP Q8NQE1
B	493	HIS	-	EXPRESSION TAG	UNP Q8NQE1

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	
			76	20	22	10	20	4	1
2	B	1	Total	C	H	N	O	P	
			39	10	12	5	10	2	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg		
			1	1	0	0
3	A	1	Total	Mg		
			1	1	0	0

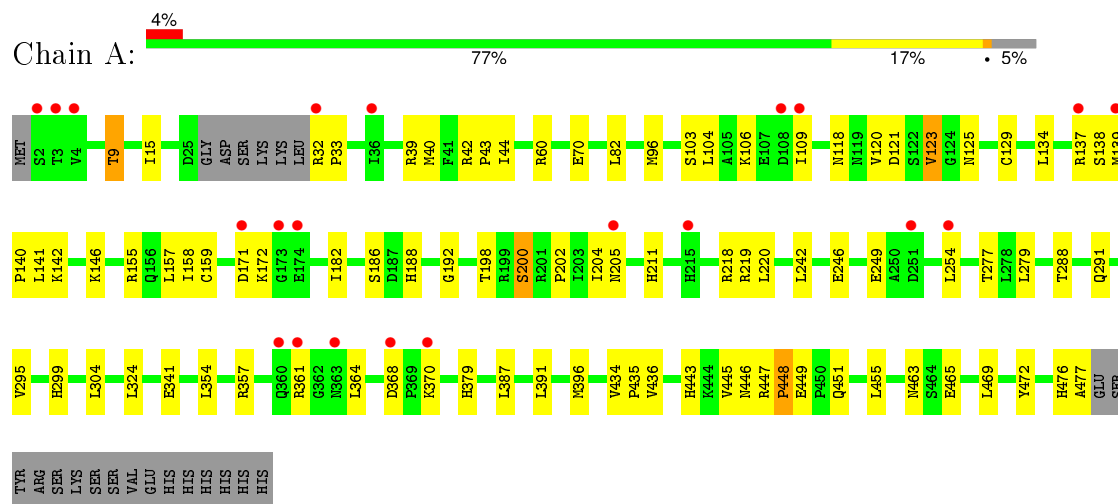
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	252	Total	O		
			252	252	0	0
4	B	182	Total	O		
			182	182	0	0

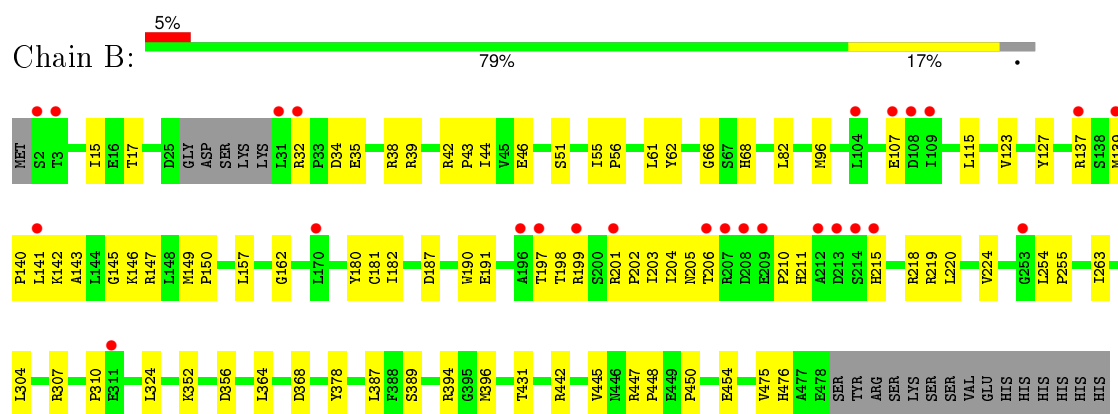
3 Residue-property plots [i](#)

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: PUP-PROTEIN LIGASE



• Molecule 1: PUP-PROTEIN LIGASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.10Å 118.81Å 163.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.42 – 2.16 49.42 – 2.16	Depositor EDS
% Data completeness (in resolution range)	98.5 (49.42-2.16) 98.5 (49.42-2.16)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.16Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.177 , 0.219 0.178 , 0.218	Depositor DCC
R_{free} test set	3274 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 64620 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7939	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% 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: MG, ADP

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	0/3761	0.58	1/5097 (0.0%)
1	B	0.38	0/3770	0.54	0/5109
All	All	0.40	0/7531	0.56	1/10206 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	CYS	N-CA-C	5.10	124.78	111.00

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	3688	0	3660	95	0
1	B	3700	0	3671	93	0
2	A	54	22	24	7	0
2	B	27	12	12	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	252	0	0	20	0
4	B	182	0	0	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7905	34	7367	170	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 11.

All (170) 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:141:LEU:HB2	1:B:206:THR:HG21	1.14	1.13
1:A:219:ARG:HG2	4:A:2102:HOH:O	1.46	1.13
1:A:9:THR:HG22	1:B:310:PRO:HB3	1.31	1.07
1:B:202:PRO:HG2	1:B:205:ASN:OD1	1.58	1.03
1:B:141:LEU:CB	1:B:206:THR:HG21	1.91	1.01
1:A:9:THR:CG2	1:B:310:PRO:HB3	1.92	0.99
1:A:445:VAL:CG2	1:A:469:LEU:HD11	1.93	0.98
4:A:2237:HOH:O	1:B:51:SER:HB2	1.65	0.95
1:B:307:ARG:HB2	4:B:2114:HOH:O	1.66	0.94
1:A:445:VAL:HG23	1:A:469:LEU:HD11	1.49	0.93
1:A:32:ARG:HB3	1:A:33:PRO:HD2	1.50	0.93
1:A:137:ARG:CD	1:A:218:ARG:HG3	2.03	0.88
1:A:211:HIS:CD2	2:A:1478[B]:ADP:O1A	2.27	0.87
1:A:435:PRO:HG2	4:A:2228:HOH:O	1.75	0.86
1:A:60:ARG:NH1	2:A:1478[B]:ADP:O1B	2.10	0.85
1:A:103:SER:HA	1:A:106:LYS:HE3	1.60	0.84
1:A:141:LEU:HD23	4:A:2105:HOH:O	1.78	0.83
1:A:104:LEU:HD23	1:B:39:ARG:NH1	1.97	0.79
1:A:60:ARG:HH12	2:A:1478[B]:ADP:PB	2.06	0.79
1:B:32:ARG:HD3	4:B:2005:HOH:O	1.81	0.79
1:B:448:PRO:HG2	1:B:476:HIS:ND1	2.00	0.77
1:A:391:LEU:HD22	1:A:396:MET:HE2	1.67	0.77
1:A:445:VAL:HG21	1:A:469:LEU:HD11	1.67	0.76
2:A:1478[A]:ADP:O3B	4:A:2251:HOH:O	2.03	0.76
1:A:137:ARG:HD2	1:A:218:ARG:HG3	1.66	0.76
1:B:203:ILE:HG23	1:B:224:VAL:HG13	1.68	0.76
1:B:141:LEU:HB2	1:B:206:THR:CG2	2.08	0.75
1:B:304:LEU:HA	4:B:2114:HOH:O	1.87	0.74
1:A:211:HIS:NE2	2:A:1478[B]:ADP:O1A	2.20	0.74
1:A:120:VAL:HB	4:A:2095:HOH:O	1.86	0.74
1:B:137:ARG:HD3	1:B:215:HIS:O	1.88	0.74
4:A:2098:HOH:O	1:B:17:THR:HG23	1.89	0.73
1:A:202:PRO:HG2	1:A:205[A]:ASN:OD1	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:ARG:O	1:A:361:ARG:HG2	1.88	0.72
1:B:141:LEU:HD12	1:B:142:LYS:N	2.05	0.72
1:A:204:ILE:HG23	1:A:220:LEU:HD11	1.71	0.72
1:B:304:LEU:HD22	1:B:324:LEU:HD12	1.73	0.71
1:A:40:MET:HE3	1:B:115:LEU:HD11	1.73	0.70
1:A:121:ASP:OD2	1:A:125:ASN:HB2	1.92	0.69
1:B:146:LYS:HE3	4:B:2056:HOH:O	1.91	0.69
1:B:201:ARG:HB2	4:B:2085:HOH:O	1.92	0.69
1:A:186:SER:HB2	4:A:2128:HOH:O	1.92	0.68
1:B:198:THR:HA	4:B:2083:HOH:O	1.92	0.68
1:A:370:LYS:NZ	1:A:370:LYS:HB2	2.09	0.68
1:B:34:ASP:HB3	4:B:2006:HOH:O	1.95	0.67
1:A:129:CYS:HB2	4:A:2098:HOH:O	1.94	0.67
1:A:198:THR:HG21	1:B:475:VAL:HG11	1.76	0.67
1:B:147:ARG:NH2	1:B:254:LEU:O	2.24	0.66
1:A:9:THR:CG2	1:B:310:PRO:CB	2.70	0.66
1:A:141:LEU:HA	4:A:2105:HOH:O	1.95	0.65
1:B:394:ARG:HB2	1:B:396:MET:HG2	1.78	0.64
1:B:352:LYS:HE2	1:B:356:ASP:OD1	1.97	0.64
1:B:210:PRO:HA	1:B:219:ARG:HG3	1.81	0.63
1:A:39:ARG:HH22	1:B:107:GLU:HG3	1.64	0.63
1:A:204:ILE:HG23	1:A:220:LEU:CD1	2.30	0.62
1:B:211:HIS:CE1	2:B:1479:ADP:O1A	2.53	0.62
1:B:61:LEU:HG	4:B:2017:HOH:O	2.00	0.61
1:A:254:LEU:HA	4:A:2149:HOH:O	2.00	0.61
1:A:304:LEU:HD22	1:A:324:LEU:HD12	1.83	0.60
4:A:2227:HOH:O	1:B:66:GLY:HA3	2.01	0.60
1:B:34:ASP:HB2	4:B:2005:HOH:O	2.02	0.60
1:B:147:ARG:HA	4:B:2055:HOH:O	2.02	0.59
1:B:396:MET:HE3	4:B:2132:HOH:O	2.02	0.59
1:A:435:PRO:HG3	1:B:62:TYR:OH	2.03	0.58
1:A:157:LEU:HA	1:A:182:ILE:HD12	1.85	0.58
1:B:123:VAL:HG12	1:B:123:VAL:O	2.03	0.58
1:B:442:ARG:HG2	1:B:454:GLU:HG2	1.86	0.57
1:A:291:GLN:O	1:A:295:VAL:HG23	2.05	0.57
1:A:32:ARG:CB	1:A:33:PRO:HD2	2.29	0.57
1:A:96:MET:SD	1:B:44:ILE:HD12	2.44	0.56
1:A:434:VAL:HG13	1:A:435:PRO:HD2	1.88	0.56
1:A:137:ARG:HD3	1:A:218:ARG:HG3	1.88	0.56
1:B:143:ALA:HB3	4:B:2051:HOH:O	2.05	0.56
1:B:204:ILE:HG23	1:B:220:LEU:HD11	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:MET:CE	1:B:115:LEU:HD11	2.37	0.55
1:A:192:GLY:O	1:A:200:SER:HA	2.06	0.55
1:A:341:GLU:HB2	4:A:2185:HOH:O	2.06	0.55
1:A:70:GLU:CD	1:B:447:ARG:HH22	2.11	0.53
1:A:60:ARG:NH1	2:A:1478[A]:ADP:O2A	2.38	0.53
1:A:446:ASN:HB3	4:A:2236:HOH:O	2.09	0.53
1:A:39:ARG:NH1	1:A:39:ARG:HG3	2.24	0.53
1:A:451:GLN:HG2	1:A:472:TYR:CE2	2.44	0.53
1:B:139:MET:HE3	4:B:2051:HOH:O	2.08	0.52
1:A:463:ASN:OD1	1:A:465:GLU:HB2	2.10	0.52
1:B:203:ILE:HG23	1:B:224:VAL:CG1	2.38	0.52
1:A:39:ARG:HG3	1:A:39:ARG:HH11	1.74	0.51
1:A:123:VAL:O	1:A:123:VAL:CG1	2.58	0.51
1:A:123:VAL:O	1:A:123:VAL:HG13	2.11	0.51
1:A:139:MET:CE	1:A:249:GLU:HG3	2.41	0.50
1:B:68:HIS:CE1	1:B:127:TYR:CE2	2.99	0.50
1:A:202:PRO:CG	1:A:205[A]:ASN:OD1	2.60	0.50
1:B:149:MET:CE	1:B:204:ILE:HD12	2.42	0.50
1:B:142:LYS:O	1:B:146:LYS:HG2	2.12	0.50
1:B:149:MET:HE2	1:B:204:ILE:HD12	1.94	0.49
1:A:364:LEU:HD22	1:A:368:ASP:OD2	2.11	0.49
1:A:391:LEU:HD22	1:A:396:MET:CE	2.39	0.49
1:B:137:ARG:HG2	1:B:137:ARG:O	2.13	0.48
1:B:139:MET:CE	1:B:140:PRO:HD2	2.44	0.48
1:B:431:THR:HG23	4:B:2158:HOH:O	2.13	0.48
1:B:35:GLU:HA	1:B:35:GLU:OE1	2.12	0.48
1:A:139:MET:O	1:A:218:ARG:NH2	2.41	0.48
1:A:299:HIS:CG	4:A:2149:HOH:O	2.67	0.48
1:A:254:LEU:CA	4:A:2149:HOH:O	2.61	0.48
1:A:42:ARG:HB3	1:A:43:PRO:HD3	1.96	0.47
1:B:42:ARG:HB3	1:B:43:PRO:HD3	1.95	0.47
1:A:9:THR:HG21	1:B:310:PRO:CB	2.43	0.47
1:B:197:THR:HB	1:B:199:ARG:NH1	2.30	0.47
1:A:446:ASN:HB2	4:A:2228:HOH:O	2.15	0.47
1:B:55:ILE:HB	1:B:56:PRO:CD	2.45	0.47
1:B:445:VAL:O	1:B:450:PRO:HA	2.14	0.47
1:A:137:ARG:HE	1:A:218:ARG:HH11	1.63	0.47
1:A:188:HIS:HD2	4:A:2094:HOH:O	1.98	0.47
1:B:364:LEU:HD22	1:B:368:ASP:OD2	2.16	0.46
1:A:104:LEU:HD23	1:B:39:ARG:CZ	2.46	0.46
1:B:180:TYR:CD2	1:B:352:LYS:HD3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ARG:HB3	1:A:33:PRO:CD	2.34	0.46
1:B:139:MET:HA	1:B:139:MET:CE	2.46	0.46
1:A:142:LYS:HE3	1:A:146:LYS:NZ	2.30	0.46
1:B:38:ARG:HD3	4:B:2007:HOH:O	2.15	0.46
1:B:139:MET:HB3	1:B:139:MET:HE3	1.89	0.46
1:B:145:GLY:HA2	1:B:204:ILE:HD13	1.97	0.46
1:A:211:HIS:HE2	2:A:1478[B]:ADP:PA	2.39	0.45
1:B:55:ILE:HB	1:B:56:PRO:HD2	1.97	0.45
1:B:157:LEU:HD23	1:B:182:ILE:CD1	2.46	0.45
1:A:476:HIS:O	1:A:477:ALA:C	2.53	0.45
1:B:149:MET:HB3	1:B:150:PRO:HD3	1.99	0.45
1:A:44:ILE:HD12	1:B:96:MET:SD	2.55	0.45
1:B:219:ARG:HD3	2:B:1479:ADP:C2	2.51	0.45
1:A:449:GLU:C	1:A:451:GLN:HE21	2.20	0.45
1:B:190:TRP:CD1	1:B:263:ILE:HD12	2.51	0.45
1:A:279:LEU:HD12	1:A:279:LEU:N	2.32	0.45
1:B:157:LEU:HA	1:B:182:ILE:HD12	1.98	0.45
1:A:15:ILE:HG12	1:B:82:LEU:HD11	1.99	0.45
1:A:202:PRO:HD2	4:A:2134:HOH:O	2.17	0.45
1:A:204:ILE:CG2	1:A:220:LEU:CD1	2.95	0.44
1:B:307:ARG:HD2	4:B:2114:HOH:O	2.17	0.44
1:A:451:GLN:CG	1:A:472:TYR:CE2	3.00	0.44
1:B:139:MET:HE1	1:B:140:PRO:HD2	1.99	0.44
1:A:370:LYS:HZ2	1:A:370:LYS:HB2	1.79	0.44
1:B:447:ARG:HA	1:B:448:PRO:C	2.39	0.43
1:A:447:ARG:HA	1:A:448:PRO:C	2.39	0.43
1:A:354:LEU:HD23	1:A:387:LEU:HD21	1.99	0.43
1:B:378:TYR:HA	1:B:387:LEU:HB2	1.99	0.43
1:B:143:ALA:O	1:B:147:ARG:HG3	2.18	0.43
1:B:55:ILE:HG12	4:B:2017:HOH:O	2.17	0.43
1:A:299:HIS:CD2	4:A:2149:HOH:O	2.71	0.43
1:A:288:THR:OG1	1:A:291:GLN:HG3	2.18	0.43
1:A:134:LEU:HD23	1:A:134:LEU:C	2.39	0.43
1:B:141:LEU:H	1:B:218:ARG:HH12	1.67	0.43
1:A:9:THR:HG21	1:B:310:PRO:HB3	1.92	0.43
1:A:242:LEU:O	1:A:246:GLU:HG3	2.19	0.43
1:A:155:ARG:O	1:A:158:ILE:HG12	2.18	0.43
1:A:118:ASN:HB3	1:A:379:HIS:CG	2.53	0.43
1:B:204:ILE:HG23	1:B:220:LEU:CD1	2.50	0.42
1:B:197:THR:HG21	1:B:199:ARG:HH22	1.83	0.42
1:B:394:ARG:CB	1:B:396:MET:HG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:THR:HG21	1:B:310:PRO:CD	2.50	0.42
1:A:445:VAL:HG21	1:A:469:LEU:CD1	2.44	0.41
1:B:368:ASP:OD1	1:B:368:ASP:C	2.58	0.41
1:B:141:LEU:N	1:B:218:ARG:HH12	2.17	0.41
1:A:82:LEU:HD11	1:B:15:ILE:HG12	2.01	0.41
1:A:443:HIS:CE1	1:A:455:LEU:HD12	2.55	0.41
1:B:254:LEU:HD12	1:B:255:PRO:HD2	2.02	0.41
1:B:162:GLY:HA2	1:B:181:CYS:O	2.21	0.41
1:B:142:LYS:HG2	1:B:146:LYS:HE2	2.03	0.41
1:A:9:THR:HG21	1:B:310:PRO:HD3	2.03	0.41
1:A:277:THR:O	1:A:288:THR:HA	2.21	0.41
1:A:139:MET:HA	1:A:140:PRO:HD3	1.80	0.40
1:A:171:ASP:OD1	1:A:172:LYS:N	2.54	0.40
1:B:187:ASP:O	1:B:191:GLU:HG3	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	467/493 (95%)	461 (99%)	5 (1%)	1 (0%)	52	51
1	B	468/493 (95%)	466 (100%)	2 (0%)	0	100	100
All	All	935/986 (95%)	927 (99%)	7 (1%)	1 (0%)	56	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	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	405/426 (95%)	399 (98%)	6 (2%)	72	78
1	B	406/426 (95%)	404 (100%)	2 (0%)	92	95
All	All	811/852 (95%)	803 (99%)	8 (1%)	82	87

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	109	ILE
1	A	123	VAL
1	A	138	SER
1	A	200	SER
1	A	436	VAL
1	B	46	GLU
1	B	389	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	A	451	GLN
1	B	443	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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	ADP	A	1478[A]	3	22,29,29	1.03	2 (9%)	27,45,45	2.04	3 (11%)
2	ADP	A	1478[B]	3	22,29,29	0.99	1 (4%)	27,45,45	2.07	6 (22%)
2	ADP	B	1479	3	22,29,29	0.98	1 (4%)	27,45,45	2.04	6 (22%)

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	ADP	A	1478[A]	3	-	0/12/32/32	0/3/3/3
2	ADP	A	1478[B]	3	-	0/12/32/32	0/3/3/3
2	ADP	B	1479	3	-	0/12/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1478[A]	ADP	O4'-C1'	2.02	1.43	1.41
2	B	1479	ADP	C5-C4	2.80	1.46	1.40
2	A	1478[A]	ADP	C5-C4	2.92	1.47	1.40
2	A	1478[B]	ADP	C5-C4	2.97	1.47	1.40

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1478[A]	ADP	N3-C2-N1	-8.01	122.76	128.89
2	A	1478[B]	ADP	N3-C2-N1	-7.95	122.81	128.89
2	B	1479	ADP	N3-C2-N1	-7.31	123.29	128.89
2	B	1479	ADP	PA-O3A-PB	-4.44	117.78	132.67
2	B	1479	ADP	C4-C5-N7	-3.28	106.47	109.48
2	A	1478[B]	ADP	C4-C5-N7	-3.23	106.51	109.48
2	A	1478[A]	ADP	C4-C5-N7	-3.20	106.53	109.48
2	A	1478[B]	ADP	PA-O3A-PB	-2.99	122.63	132.67
2	A	1478[A]	ADP	PA-O3A-PB	-2.43	124.53	132.67
2	B	1479	ADP	C1'-N9-C4	-2.24	123.56	126.94
2	A	1478[B]	ADP	C1'-N9-C4	-2.02	123.89	126.94
2	A	1478[B]	ADP	C2-N1-C6	2.01	122.35	118.77
2	B	1479	ADP	O3B-PB-O2B	2.01	115.04	107.38
2	A	1478[B]	ADP	O3B-PB-O2B	2.07	115.28	107.38
2	B	1479	ADP	O3B-PB-O1B	2.20	117.67	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1478[A]	ADP	2	0
2	A	1478[B]	ADP	5	0
2	B	1479	ADP	2	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 ⓘ

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	470/493 (95%)	-0.01	21 (4%)	37 48	27, 47, 88, 133	0
1	B	472/493 (95%)	0.12	26 (5%)	29 39	30, 49, 109, 141	0
All	All	942/986 (95%)	0.06	47 (4%)	32 43	27, 48, 100, 141	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	SER	7.2
1	B	170	LEU	6.7
1	B	212	ALA	6.7
1	B	141	LEU	6.6
1	B	139	MET	6.0
1	B	214	SER	5.5
1	B	207	ARG	4.9
1	A	4	VAL	4.7
1	B	206	THR	4.6
1	B	2	SER	4.5
1	B	213	ASP	4.5
1	B	108	ASP	4.4
1	A	173	GLY	4.2
1	B	209	GLU	4.1
1	A	32	ARG	3.7
1	B	199	ARG	3.7
1	A	361	ARG	3.6
1	A	3	THR	3.3
1	B	201	ARG	3.1
1	B	104	LEU	3.1
1	B	215	HIS	3.1
1	B	109	ILE	3.1
1	B	196	ALA	3.1
1	A	215	HIS	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	36	ILE	2.8
1	B	107	GLU	2.7
1	A	363	ASN	2.7
1	A	108	ASP	2.7
1	B	253	GLY	2.6
1	A	174	GLU	2.6
1	A	360	GLN	2.6
1	B	32	ARG	2.6
1	B	31	LEU	2.6
1	A	251	ASP	2.5
1	A	109	ILE	2.4
1	B	208	ASP	2.3
1	A	368	ASP	2.2
1	A	139	MET	2.2
1	A	137	ARG	2.2
1	B	197	THR	2.2
1	A	171	ASP	2.1
1	B	137	ARG	2.1
1	A	370	LYS	2.1
1	A	254	LEU	2.1
1	A	205[A]	ASN	2.1
1	B	3	THR	2.0
1	B	311	GLU	2.0

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	ADP	B	1479	27/27	0.97	0.11	-0.47	36,43,58,64	0
2	ADP	A	1478[B]	27/27	0.98	0.11	-0.84	33,38,45,48	38
2	ADP	A	1478[A]	27/27	0.98	0.11	-0.87	31,38,46,53	38
3	MG	B	1480	1/1	0.90	0.25	-	60,60,60,60	0
3	MG	A	1479	1/1	0.99	0.18	-	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.