



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2024 – 04:08 AM EDT

PDB ID : 4H2U  
Title : Crystal structure of Bradyrhizobium japonicum glycine:[carrier protein] ligase complexed with cognate carrier protein and ATP  
Authors : Luic, M.; Weygand-Durasevic, I.; Ivic, N.; Mocibob, M.  
Deposited on : 2012-09-13  
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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

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

MolProbity : 4.02b-467  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.20.1  
EDS : 2.37.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.37.1

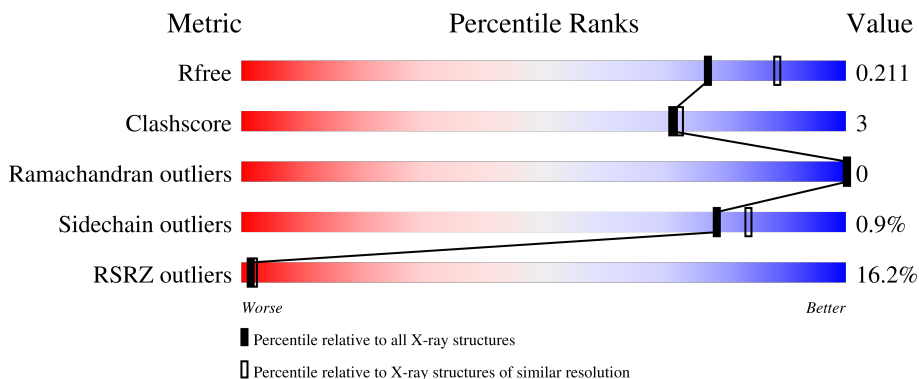
# 1 Overall quality at a glance

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))
$R_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	346	
1	B	346	
2	C	110	
2	D	110	

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 5853 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 Amino acid--[acyl-carrier-protein] ligase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	295	2317	1463	410	427	17	0	6	0
1	B	297	2311	1461	411	421	18	0	6	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q89VT8
A	-18	GLY	-	EXPRESSION TAG	UNP Q89VT8
A	-17	SER	-	EXPRESSION TAG	UNP Q89VT8
A	-16	SER	-	EXPRESSION TAG	UNP Q89VT8
A	-15	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-14	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-13	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-12	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-11	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-10	HIS	-	EXPRESSION TAG	UNP Q89VT8
A	-9	SER	-	EXPRESSION TAG	UNP Q89VT8
A	-8	SER	-	EXPRESSION TAG	UNP Q89VT8
A	-7	GLY	-	EXPRESSION TAG	UNP Q89VT8
A	-6	LEU	-	EXPRESSION TAG	UNP Q89VT8
A	-5	VAL	-	EXPRESSION TAG	UNP Q89VT8
A	-4	PRO	-	EXPRESSION TAG	UNP Q89VT8
A	-3	ARG	-	EXPRESSION TAG	UNP Q89VT8
A	-2	GLY	-	EXPRESSION TAG	UNP Q89VT8
A	-1	SER	-	EXPRESSION TAG	UNP Q89VT8
A	0	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-19	MET	-	EXPRESSION TAG	UNP Q89VT8
B	-18	GLY	-	EXPRESSION TAG	UNP Q89VT8
B	-17	SER	-	EXPRESSION TAG	UNP Q89VT8
B	-16	SER	-	EXPRESSION TAG	UNP Q89VT8
B	-15	HIS	-	EXPRESSION TAG	UNP Q89VT8

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-13	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-12	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-11	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-10	HIS	-	EXPRESSION TAG	UNP Q89VT8
B	-9	SER	-	EXPRESSION TAG	UNP Q89VT8
B	-8	SER	-	EXPRESSION TAG	UNP Q89VT8
B	-7	GLY	-	EXPRESSION TAG	UNP Q89VT8
B	-6	LEU	-	EXPRESSION TAG	UNP Q89VT8
B	-5	VAL	-	EXPRESSION TAG	UNP Q89VT8
B	-4	PRO	-	EXPRESSION TAG	UNP Q89VT8
B	-3	ARG	-	EXPRESSION TAG	UNP Q89VT8
B	-2	GLY	-	EXPRESSION TAG	UNP Q89VT8
B	-1	SER	-	EXPRESSION TAG	UNP Q89VT8
B	0	HIS	-	EXPRESSION TAG	UNP Q89VT8

- Molecule 2 is a protein called Aminoacyl carrier protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	38	Total	C	N	O	S	0	0	0
			232	144	40	45	3			
2	D	78	Total	C	N	O	S	0	0	0
			537	339	85	109	4			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	MET	-	EXPRESSION TAG	UNP Q89VT6
C	-18	GLY	-	EXPRESSION TAG	UNP Q89VT6
C	-17	SER	-	EXPRESSION TAG	UNP Q89VT6
C	-16	SER	-	EXPRESSION TAG	UNP Q89VT6
C	-15	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-14	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-13	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-12	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-11	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-10	HIS	-	EXPRESSION TAG	UNP Q89VT6
C	-9	SER	-	EXPRESSION TAG	UNP Q89VT6
C	-8	SER	-	EXPRESSION TAG	UNP Q89VT6
C	-7	GLY	-	EXPRESSION TAG	UNP Q89VT6
C	-6	LEU	-	EXPRESSION TAG	UNP Q89VT6
C	-5	VAL	-	EXPRESSION TAG	UNP Q89VT6

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	EXPRESSION TAG	UNP Q89VT6
C	-3	ARG	-	EXPRESSION TAG	UNP Q89VT6
C	-2	GLY	-	EXPRESSION TAG	UNP Q89VT6
C	-1	SER	-	EXPRESSION TAG	UNP Q89VT6
C	0	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-19	MET	-	EXPRESSION TAG	UNP Q89VT6
D	-18	GLY	-	EXPRESSION TAG	UNP Q89VT6
D	-17	SER	-	EXPRESSION TAG	UNP Q89VT6
D	-16	SER	-	EXPRESSION TAG	UNP Q89VT6
D	-15	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-14	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-13	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-12	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-11	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-10	HIS	-	EXPRESSION TAG	UNP Q89VT6
D	-9	SER	-	EXPRESSION TAG	UNP Q89VT6
D	-8	SER	-	EXPRESSION TAG	UNP Q89VT6
D	-7	GLY	-	EXPRESSION TAG	UNP Q89VT6
D	-6	LEU	-	EXPRESSION TAG	UNP Q89VT6
D	-5	VAL	-	EXPRESSION TAG	UNP Q89VT6
D	-4	PRO	-	EXPRESSION TAG	UNP Q89VT6
D	-3	ARG	-	EXPRESSION TAG	UNP Q89VT6
D	-2	GLY	-	EXPRESSION TAG	UNP Q89VT6
D	-1	SER	-	EXPRESSION TAG	UNP Q89VT6
D	0	HIS	-	EXPRESSION TAG	UNP Q89VT6

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



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

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

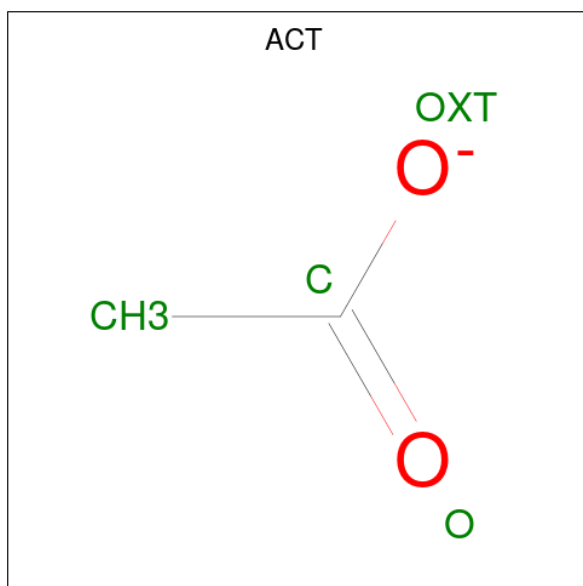
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
5	A	1	1	1	0	0
5	B	1	1	1	0	0

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



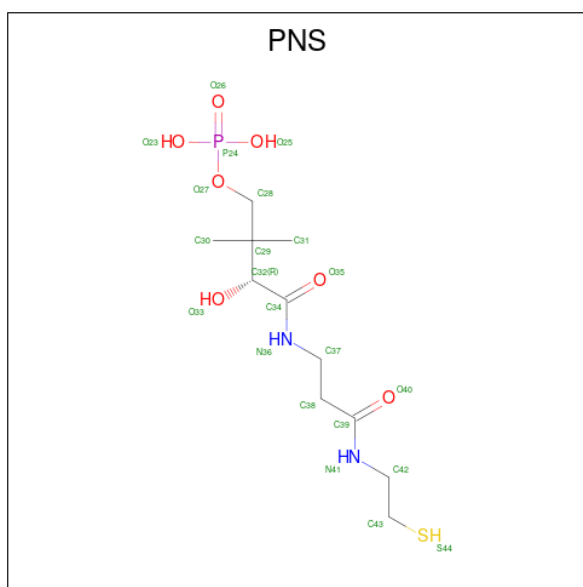
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
6	B	1	5	4	1	0	0

- Molecule 7 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
7	B	1	4	2	2	0	0

- Molecule 8 is 4'-PHOSPHOPANTETHEINE (three-letter code: PNS) (formula:  $C_{11}H_{23}N_2O_7PS$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
8	C	1	21	11	2	6	1	1	0	0
8	D	1	21	11	2	6	1	1	0	0

- Molecule 9 is water.

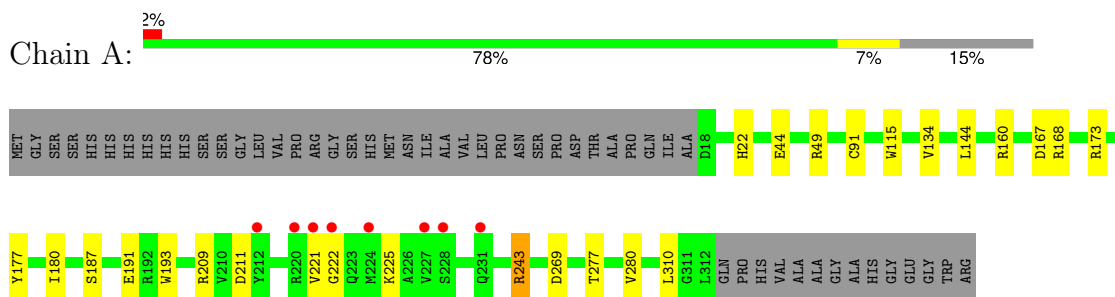
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	178	Total	O	0	0
			178	178		
9	B	157	Total	O	0	0
			157	157		
9	D	4	Total	O	0	0
			4	4		



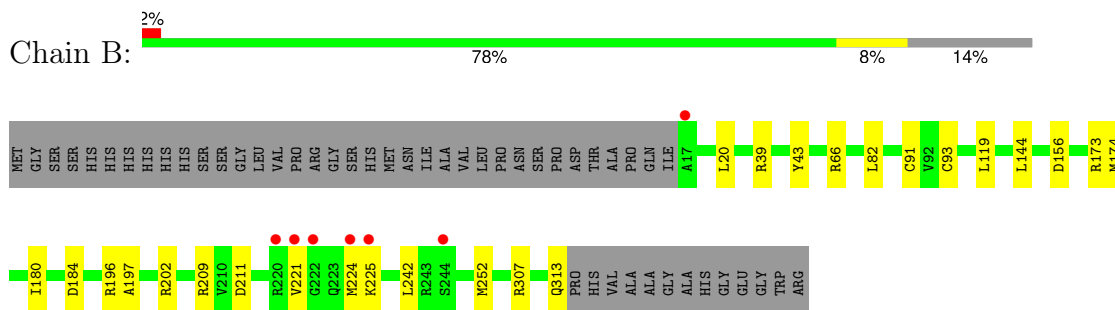
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

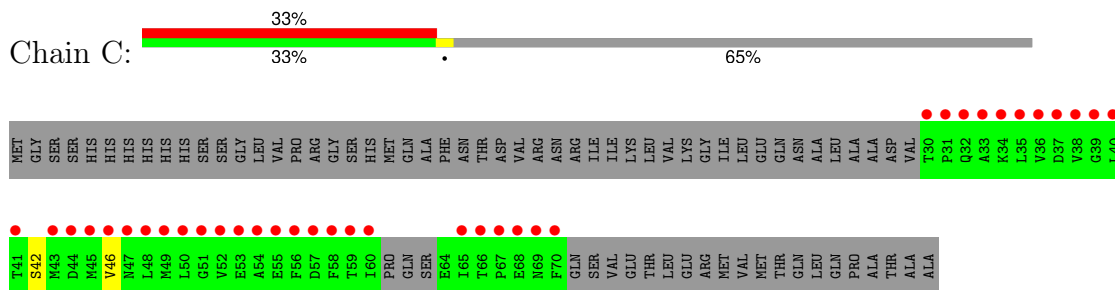
- Molecule 1: Amino acid--[acyl-carrier-protein] ligase 1



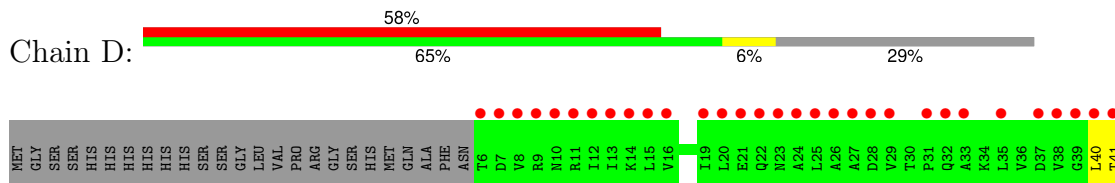
- Molecule 1: Amino acid--[acyl-carrier-protein] ligase 1

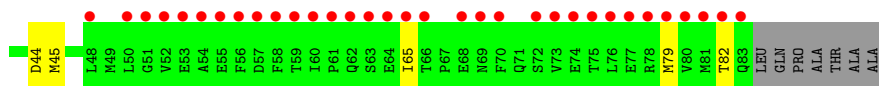


- Molecule 2: Aminoacyl carrier protein 1



- Molecule 2: Aminoacyl carrier protein 1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.71Å 100.90Å 104.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.32 – 2.10 46.32 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.32-2.10) 99.7 (46.32-2.10)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.31 (at 2.10Å)	Xtrriage
Refinement program	PHENIX dev_1116	Depositor
R, $R_{free}$	0.175 , 0.209 0.179 , 0.211	Depositor DCC
$R_{free}$ test set	2833 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtrriage
Anisotropy	0.328	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 54.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.019 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5853	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: ZN, MG, ATP, ACT, SO4, PNS

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	0.40	0/2390	0.54	0/3243
1	B	0.39	0/2383	0.53	0/3231
2	C	0.24	0/233	0.48	0/318
2	D	0.25	0/542	0.42	0/741
All	All	0.38	0/5548	0.52	0/7533

There are no bond length outliers.

There are no bond angle outliers.

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	2317	0	2241	19	0
1	B	2311	0	2240	18	0
2	C	232	0	176	1	0
2	D	537	0	480	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	31	0	12	1	0
4	B	31	0	12	2	0
5	A	1	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	1	0	0	0	0
6	B	5	0	0	0	0
7	B	4	0	3	0	0
8	C	21	0	21	0	0
8	D	21	0	21	0	0
9	A	178	0	0	2	0
9	B	157	0	0	3	0
9	D	4	0	0	0	0
All	All	5853	0	5206	37	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 3.

All (37) 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:168:ARG:NH1	4:A:402:ATP:O3G	2.24	0.66
1:B:196:ARG:NH2	9:B:542:HOH:O	2.35	0.60
1:A:44:GLU:HG2	1:B:66:ARG:HG3	1.87	0.57
1:A:243:ARG:CZ	1:A:243:ARG:HB3	2.36	0.56
1:A:222:GLY:HA2	1:A:225:LYS:HD2	1.88	0.56
1:A:187:SER:O	1:A:191:GLU:HG2	2.08	0.53
1:A:115:TRP:HH2	1:B:91[B]:CYS:SG	2.33	0.52
1:A:115:TRP:HH2	1:B:91[B]:CYS:HG	1.56	0.52
1:A:49:ARG:HD3	1:A:310:LEU:O	2.09	0.52
1:B:184:ASP:OD1	9:B:576:HOH:O	2.19	0.51
1:B:202:ARG:NH2	9:B:520:HOH:O	2.45	0.50
4:B:402:ATP:H5'1	4:B:402:ATP:H8	1.76	0.50
1:A:22:HIS:HD2	9:A:667:HOH:O	1.93	0.50
1:A:167:ASP:OD1	1:A:243:ARG:NH2	2.41	0.49
1:B:93[B]:CYS:SG	1:B:119:LEU:HD13	2.52	0.49
1:A:144:LEU:HD12	1:A:269:ASP:HA	1.95	0.49
1:A:160:ARG:HG2	9:A:666:HOH:O	2.11	0.49
1:A:115:TRP:CE2	1:B:82:LEU:HD11	2.48	0.49
1:A:177:TYR:HB2	1:A:280[A]:VAL:CG1	2.43	0.48
2:D:79:MET:O	2:D:82:THR:HG22	2.13	0.47
1:B:209:ARG:NH1	1:B:211[B]:ASP:OD2	2.42	0.47
1:A:221:VAL:HG12	1:A:225:LYS:HE3	1.96	0.47
1:B:221:VAL:HG12	1:B:225:LYS:HE3	1.96	0.46
1:B:174[B]:MET:SD	4:B:402:ATP:O4'	2.73	0.46
1:B:144:LEU:HD11	1:B:180:ILE:HG22	1.97	0.46

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:ALA:HB1	1:B:252:MET:SD	2.57	0.45
1:B:20:LEU:HD21	1:B:242:LEU:HB3	1.99	0.44
1:B:307:ARG:HH21	1:B:313:GLN:C	2.21	0.44
2:D:40:LEU:HD22	2:D:44:ASP:HB3	1.99	0.43
2:C:42:SER:O	2:C:46:VAL:HG23	2.18	0.43
1:A:180:ILE:HG23	1:A:277:THR:HG22	2.00	0.43
1:B:224:MET:HE2	2:D:65:ILE:HG23	2.00	0.42
2:D:41:THR:O	2:D:45:MET:HG2	2.19	0.42
1:A:91:CYS:HB3	1:B:93[B]:CYS:SG	2.60	0.42
1:A:193:TRP:CB	1:A:280[B]:VAL:HG21	2.50	0.41
1:A:209:ARG:NH1	1:A:211:ASP:OD2	2.53	0.41
1:B:39:ARG:HG2	1:B:43:TYR:CD2	2.56	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	299/346 (86%)	295 (99%)	4 (1%)	0	100	100
1	B	301/346 (87%)	298 (99%)	3 (1%)	0	100	100
2	C	34/110 (31%)	29 (85%)	5 (15%)	0	100	100
2	D	76/110 (69%)	72 (95%)	4 (5%)	0	100	100
All	All	710/912 (78%)	694 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	243/280 (87%)	240 (99%)	3 (1%)	71	77
1	B	239/280 (85%)	237 (99%)	2 (1%)	81	86
2	C	16/95 (17%)	16 (100%)	0	100	100
2	D	50/95 (53%)	50 (100%)	0	100	100
All	All	548/750 (73%)	543 (99%)	5 (1%)	78	84

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	VAL
1	A	173	ARG
1	A	243	ARG
1	B	156	ASP
1	B	173	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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
4	ATP	B	402	5	28,33,33	1.17	2 (7%)	34,52,52	1.68	7 (20%)
8	PNS	C	1000	2	15,20,21	0.31	0	18,26,29	0.52	0
4	ATP	A	402	5	28,33,33	1.06	3 (10%)	34,52,52	1.71	5 (14%)
6	SO4	B	403	-	4,4,4	0.27	0	6,6,6	0.09	0
8	PNS	D	1000	2	15,20,21	0.26	0	18,26,29	0.57	0
7	ACT	B	405	-	3,3,3	0.87	0	3,3,3	1.25	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
4	ATP	B	402	5	-	4/18/38/38	0/3/3/3
8	PNS	D	1000	2	-	0/24/26/27	-
8	PNS	C	1000	2	-	1/24/26/27	-
4	ATP	A	402	5	-	3/18/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	ATP	PB-O3A	4.18	1.64	1.59
4	A	402	ATP	PA-O3A	2.50	1.62	1.59
4	B	402	ATP	PA-O3A	2.30	1.62	1.59
4	A	402	ATP	PB-O3A	2.16	1.61	1.59
4	A	402	ATP	PG-O1G	2.04	1.56	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	402	ATP	O2B-PB-O3B	-5.04	93.64	107.27
4	A	402	ATP	O3A-PB-O1B	-4.91	95.95	110.70
4	B	402	ATP	O2B-PB-O3B	-4.40	95.38	107.27

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	402	ATP	O3A-PA-O1A	-4.08	98.43	110.70
4	B	402	ATP	O3A-PA-O1A	-3.53	100.08	110.70
4	B	402	ATP	O2A-PA-O5'	-3.30	92.59	107.57
4	A	402	ATP	O2B-PB-O1B	3.22	127.42	112.44
4	B	402	ATP	O3A-PB-O1B	-3.21	101.06	110.70
4	B	402	ATP	O2A-PA-O1A	3.00	126.40	112.44
4	B	402	ATP	C5-C6-N6	2.12	123.53	120.31
4	B	402	ATP	O2B-PB-O1B	2.10	122.23	112.44
4	A	402	ATP	O2A-PA-O1A	2.04	121.95	112.44

There are no chirality outliers.

All (8) torsion outliers are listed below:

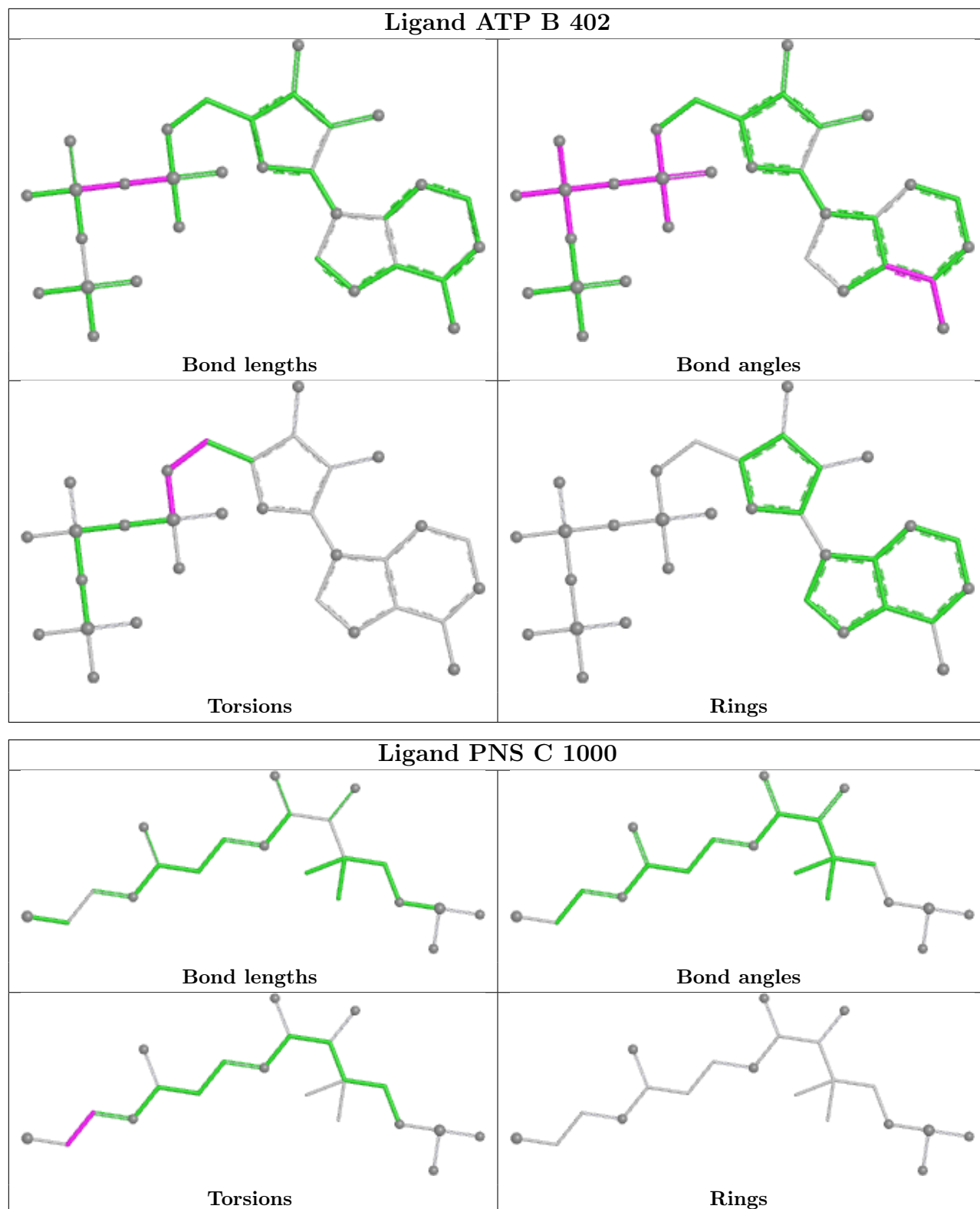
Mol	Chain	Res	Type	Atoms
4	A	402	ATP	C5'-O5'-PA-O2A
4	A	402	ATP	C5'-O5'-PA-O3A
4	B	402	ATP	C5'-O5'-PA-O1A
8	C	1000	PNS	N41-C42-C43-S44
4	B	402	ATP	C5'-O5'-PA-O2A
4	B	402	ATP	C5'-O5'-PA-O3A
4	A	402	ATP	C4'-C5'-O5'-PA
4	B	402	ATP	C4'-C5'-O5'-PA

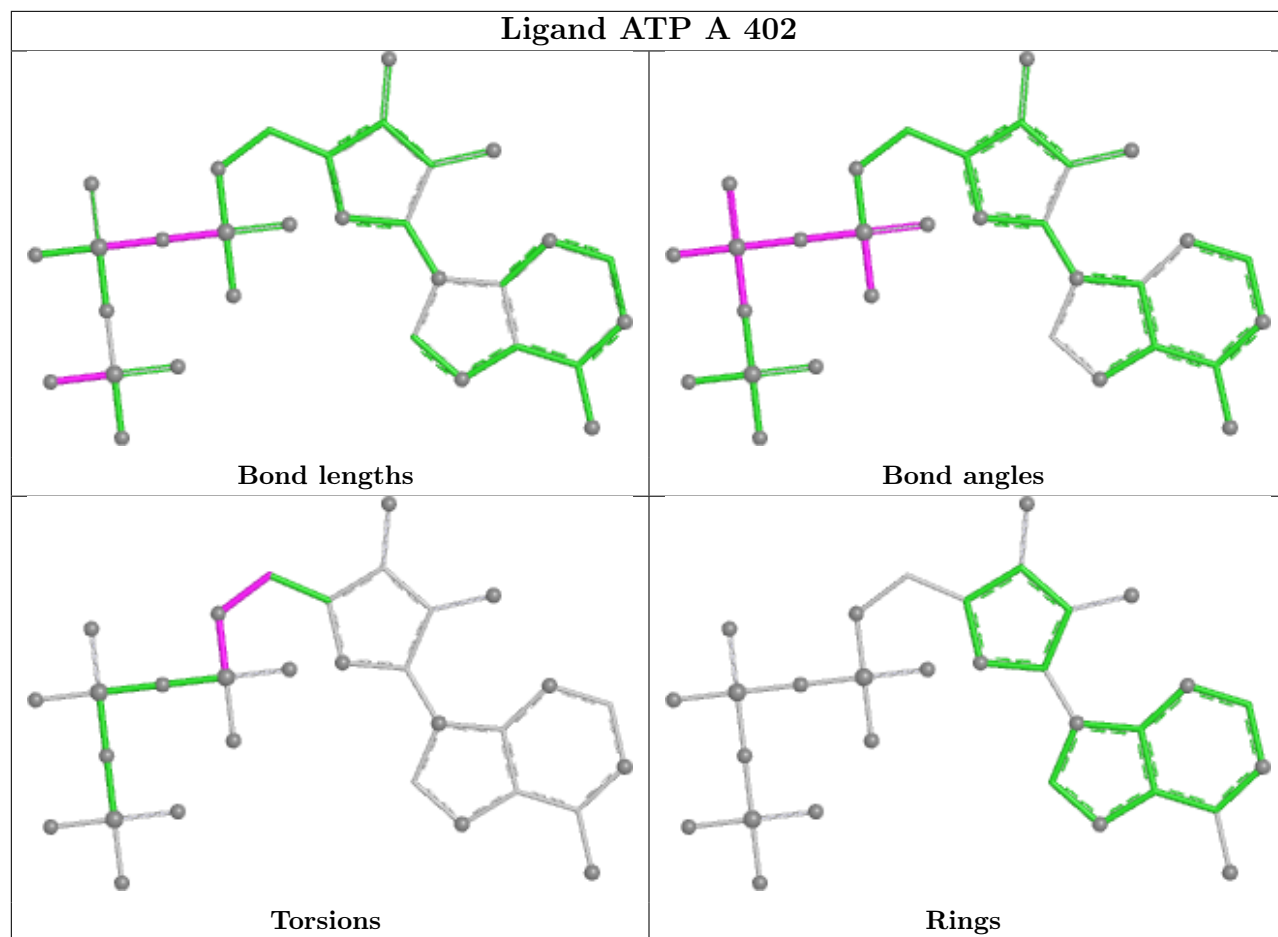
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	402	ATP	2	0
4	A	402	ATP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 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, 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	A	295/346 (85%)	0.33	8 (2%) 54 60	14, 26, 50, 72	0
1	B	297/346 (85%)	0.20	7 (2%) 59 64	13, 27, 48, 70	0
2	C	38/110 (34%)	5.84	36 (94%) 0 0	72, 95, 115, 120	0
2	D	78/110 (70%)	3.86	64 (82%) 0 0	53, 75, 101, 104	0
All	All	708/912 (77%)	0.96	115 (16%) 1 2	13, 29, 90, 120	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	50	LEU	13.4
2	C	40	LEU	11.0
2	C	51	GLY	10.8
2	D	12	ILE	9.5
2	D	54	ALA	9.1
2	D	56	PHE	9.0
2	C	52	VAL	8.9
2	C	33	ALA	8.2
2	D	58	PHE	7.9
2	C	54	ALA	7.9
2	C	48	LEU	7.8
2	C	66	THR	7.7
2	C	31	PRO	7.2
2	C	46	VAL	7.1
2	C	65	ILE	6.9
1	B	17	ALA	6.8
2	D	24	ALA	6.8
2	C	67	PRO	6.7
2	D	38	VAL	6.7
2	C	70	PHE	6.5
1	A	221	VAL	6.3

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	D	10	ASN	6.3
2	D	8	VAL	6.3
2	D	13	ILE	6.3
2	D	80	VAL	6.3
2	D	26	ALA	6.2
2	D	79	MET	6.2
2	D	59	THR	6.1
2	C	53	GLU	6.1
2	C	43	MET	6.1
2	C	47	ASN	6.1
2	D	65	ILE	6.1
2	D	81	MET	6.0
2	D	52	VAL	5.8
2	D	62	GLN	5.8
2	C	36	VAL	5.7
2	D	76	LEU	5.7
2	C	38	VAL	5.6
2	C	59	THR	5.5
2	C	57	ASP	5.5
2	D	25	LEU	5.4
2	C	44	ASP	5.3
2	C	49	MET	5.3
2	D	31	PRO	5.2
2	D	19	ILE	5.1
2	C	55	GLU	5.1
2	C	39	GLY	5.1
2	D	60	ILE	4.9
2	D	16	VAL	4.9
2	C	69	ASN	4.9
2	C	37	ASP	4.8
2	D	41	THR	4.8
2	D	57	ASP	4.7
2	C	58	PHE	4.7
2	C	45	MET	4.6
2	D	75	THR	4.4
2	D	7	ASP	4.3
2	D	64	GLU	4.2
2	D	82	THR	4.1
2	D	39	GLY	4.1
2	D	55	GLU	4.1
2	D	48	LEU	4.1
2	D	50	LEU	4.1

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	D	27	ALA	4.1
2	D	15	LEU	4.0
1	A	224	MET	4.0
2	C	30	THR	3.9
2	D	22	GLN	3.9
2	C	35	LEU	3.8
2	D	51	GLY	3.8
2	C	56	PHE	3.8
2	C	32	GLN	3.7
2	D	6	THR	3.7
2	D	23	ASN	3.7
2	C	41	THR	3.6
2	D	28	ASP	3.6
2	C	34	LYS	3.6
2	D	32	GLN	3.6
2	D	74	GLU	3.5
2	D	9	ARG	3.5
2	C	60	ILE	3.5
2	C	68	GLU	3.3
2	D	63	SER	3.3
2	D	35	LEU	3.3
2	D	14	LYS	3.2
2	D	40	LEU	3.2
2	D	73	VAL	3.2
2	D	11	ARG	3.2
2	D	77	GLU	3.2
2	D	83	GLN	3.1
1	A	227	VAL	3.1
1	A	222	GLY	3.1
1	A	231	GLN	3.1
1	B	220	ARG	3.1
2	D	33	ALA	3.0
1	B	222	GLY	3.0
2	D	61	PRO	2.9
1	A	212	TYR	2.9
2	D	53	GLU	2.8
1	B	221	VAL	2.8
1	A	220	ARG	2.7
2	D	66	THR	2.7
2	D	69	ASN	2.6
2	D	20	LEU	2.6
1	A	228	SER	2.5

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	D	72	SER	2.5
2	D	37	ASP	2.5
2	D	70	PHE	2.5
1	B	244	SER	2.4
2	D	78	ARG	2.3
1	B	225	LYS	2.3
2	D	21	GLU	2.3
1	B	224	MET	2.1
2	D	29	VAL	2.1
2	D	68	GLU	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 monosaccharides in this entry.

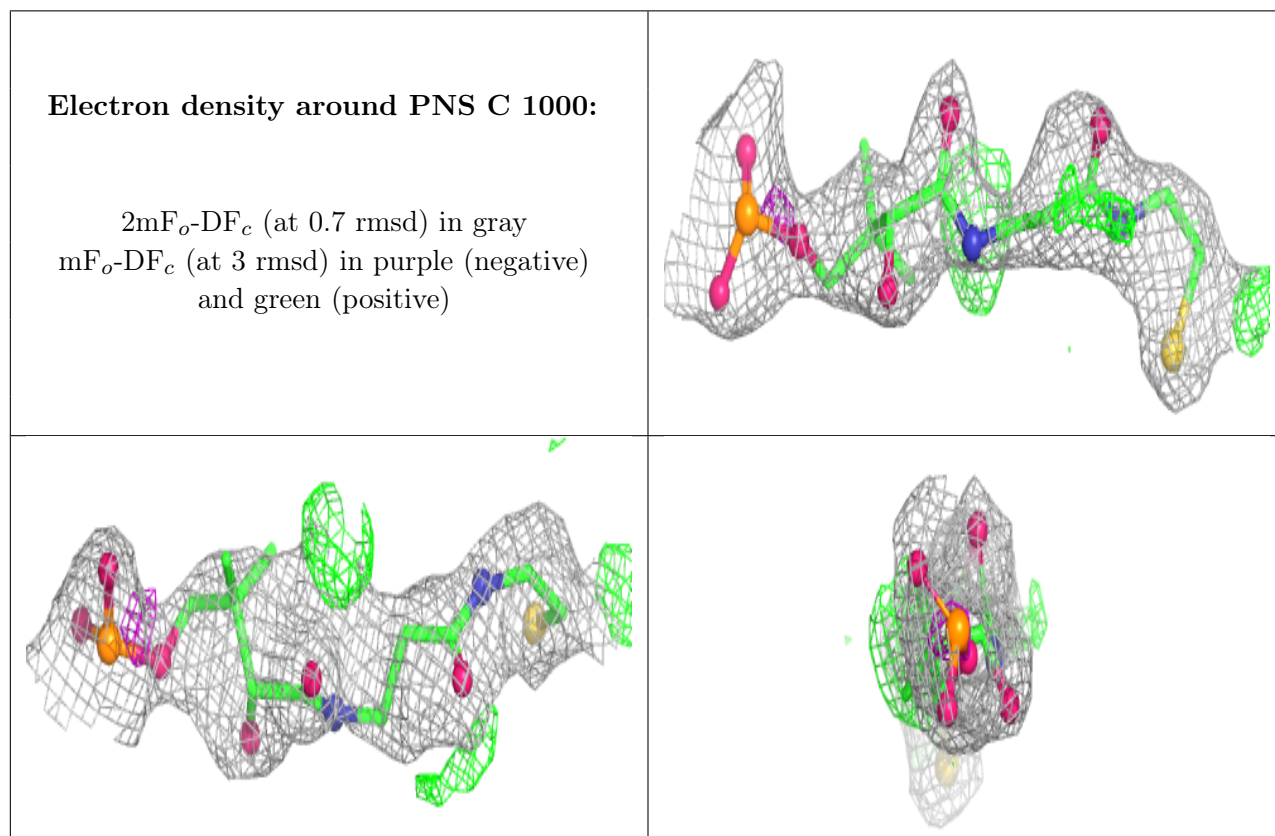
## 6.4 Ligands [i](#)

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
7	ACT	B	405	4/4	0.83	0.21	41,47,47,50	0
8	PNS	C	1000	21/22	0.89	0.24	45,59,70,80	0
5	MG	A	403	1/1	0.90	0.23	45,45,45,45	0
4	ATP	A	402	31/31	0.91	0.19	13,23,51,56	31
8	PNS	D	1000	21/22	0.91	0.20	49,58,67,77	0
5	MG	B	404	1/1	0.92	0.07	48,48,48,48	0
4	ATP	B	402	31/31	0.92	0.16	16,29,45,47	31
6	SO4	B	403	5/5	0.97	0.19	42,46,51,56	0
3	ZN	B	401	1/1	0.97	0.08	31,31,31,31	1
3	ZN	A	401	1/1	0.98	0.08	27,27,27,27	1

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

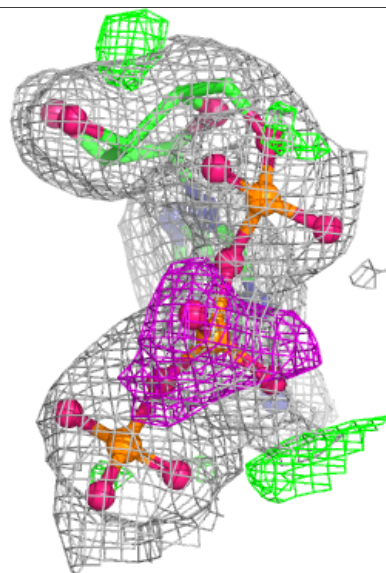
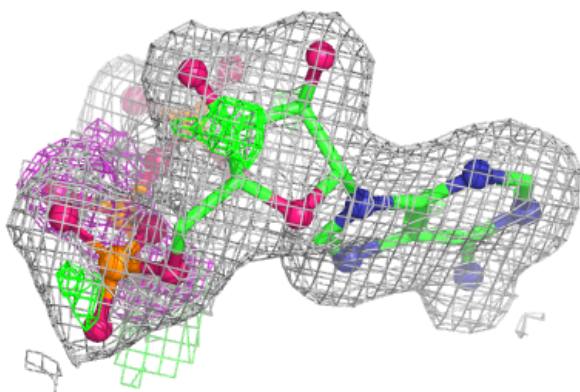
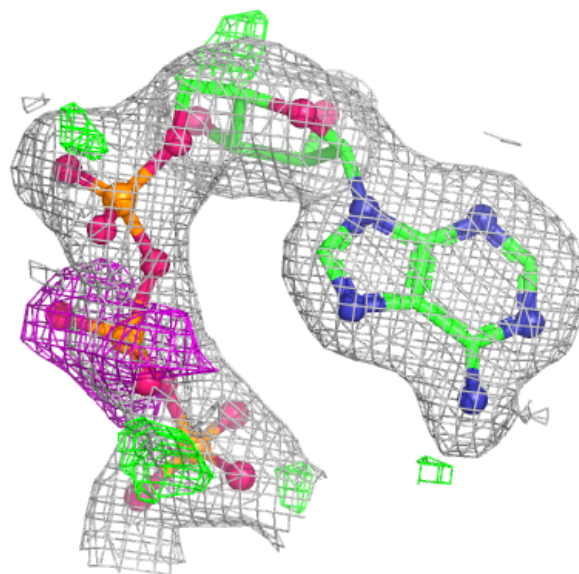
as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

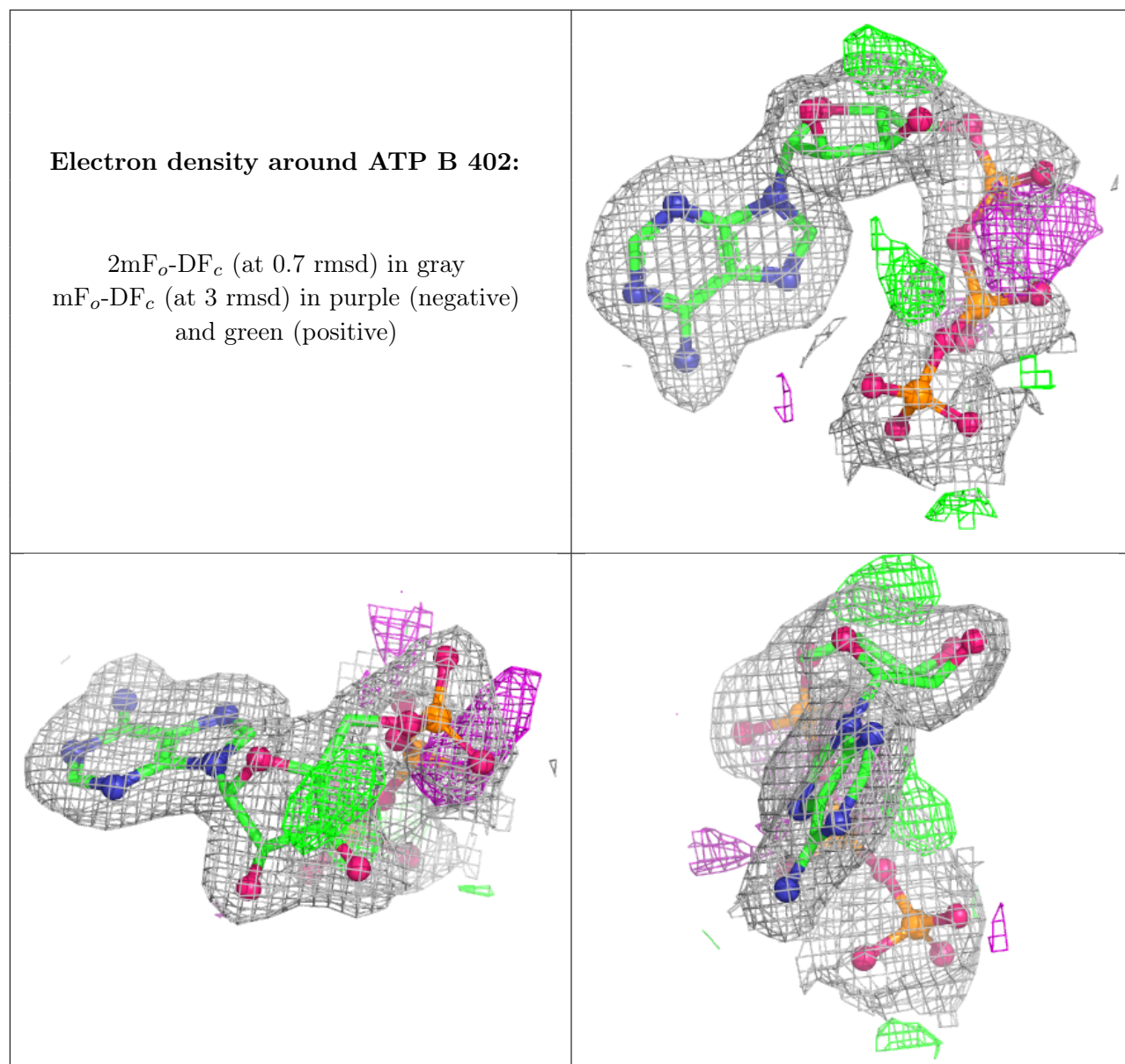




**Electron density around ATP A 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers ⓘ

There are no such residues in this entry.