



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2026 – 02:35 PM UTC

PDB ID : 6NYA / pdb_00006nya
Title : Crystal Structure of ubiquitin E1 (Uba1) in complex with Ubc3 (Cdc34) and ubiquitin
Authors : Olsen, S.K.; Williams, K.M.; Atkison, J.H.
Deposited on : 2019-02-11
Resolution : 2.06 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

<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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

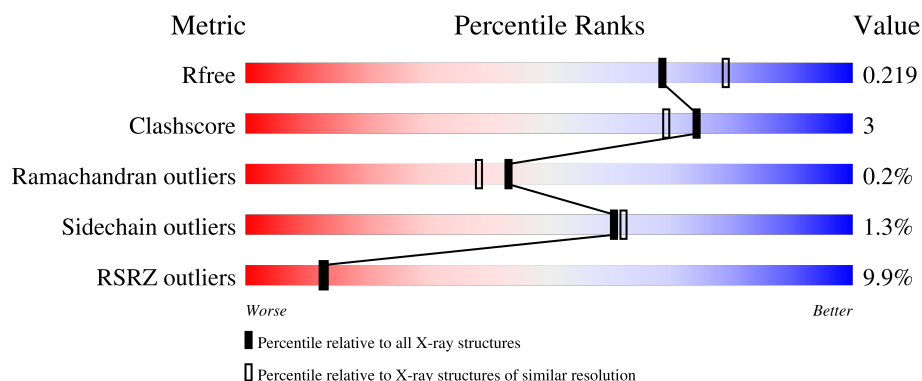
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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}	180053	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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 ≥ 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	1017	4% 89% 8% .
1	D	1017	10% 89% 8% .
2	B	85	5% 84% 7% . 8%
2	E	85	15% 79% 12% 9%
3	C	197	20% 76% 11% 14%

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Mol	Chain	Length	Quality of chain
3	F	197	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	SO4	D	1204	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 20705 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 Ubiquitin-activating enzyme E1 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	987	Total	C	N	O	S	0	0	0
			7800	4979	1288	1510	23			
1	D	987	Total	C	N	O	S	0	0	0
			7796	4977	1288	1508	23			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	GLY	-	expression tag	UNP P22515
A	9	ALA	-	expression tag	UNP P22515
A	10	MET	-	expression tag	UNP P22515
D	8	GLY	-	expression tag	UNP P22515
D	9	ALA	-	expression tag	UNP P22515
D	10	MET	-	expression tag	UNP P22515

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	78	Total	C	N	O	S	0	0	0
			622	380	121	120	1			
2	E	77	Total	C	N	O	S	0	0	0
			618	378	120	119	1			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	MET	-	initiating methionine	UNP A0A1D6RLC6
B	-7	HIS	-	expression tag	UNP A0A1D6RLC6
B	-6	HIS	-	expression tag	UNP A0A1D6RLC6
B	-5	HIS	-	expression tag	UNP A0A1D6RLC6
B	-4	HIS	-	expression tag	UNP A0A1D6RLC6
B	-3	HIS	-	expression tag	UNP A0A1D6RLC6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP A0A1D6RLC6
B	-1	GLY	-	expression tag	UNP A0A1D6RLC6
B	0	ALA	-	expression tag	UNP A0A1D6RLC6
B	6	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	11	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	27	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	29	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	33	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	48	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
B	63	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	-8	MET	-	initiating methionine	UNP A0A1D6RLC6
E	-7	HIS	-	expression tag	UNP A0A1D6RLC6
E	-6	HIS	-	expression tag	UNP A0A1D6RLC6
E	-5	HIS	-	expression tag	UNP A0A1D6RLC6
E	-4	HIS	-	expression tag	UNP A0A1D6RLC6
E	-3	HIS	-	expression tag	UNP A0A1D6RLC6
E	-2	HIS	-	expression tag	UNP A0A1D6RLC6
E	-1	GLY	-	expression tag	UNP A0A1D6RLC6
E	0	ALA	-	expression tag	UNP A0A1D6RLC6
E	6	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	11	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	27	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	29	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	33	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	48	ARG	LYS	engineered mutation	UNP A0A1D6RLC6
E	63	ARG	LYS	engineered mutation	UNP A0A1D6RLC6

- Molecule 3 is a protein called Ubiquitin-conjugating enzyme E2-34 kDa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	170	Total	C	N	O	S	0	0	0
			1381	881	235	260	5			
3	F	167	Total	C	N	O	S	0	0	0
			1356	868	229	254	5			

There are 8 discrepancies between the modelled and reference sequences:

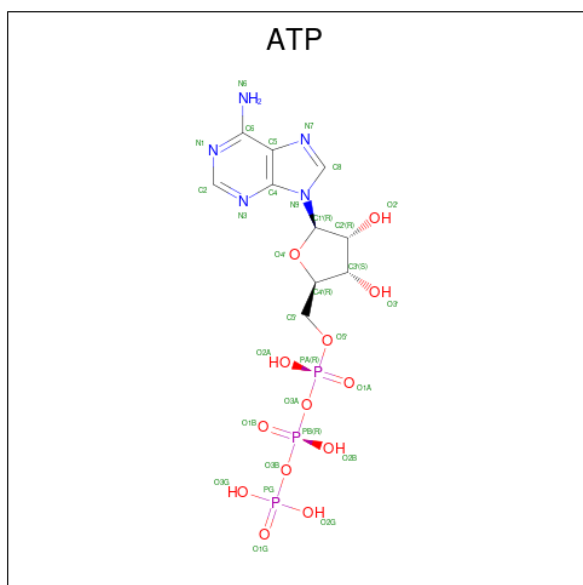
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP P14682
C	0	ALA	-	expression tag	UNP P14682
C	1	MET	-	expression tag	UNP P14682
C	2	ALA	-	expression tag	UNP P14682

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	GLY	-	expression tag	UNP P14682
F	0	ALA	-	expression tag	UNP P14682
F	1	MET	-	expression tag	UNP P14682
F	2	ALA	-	expression tag	UNP P14682

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



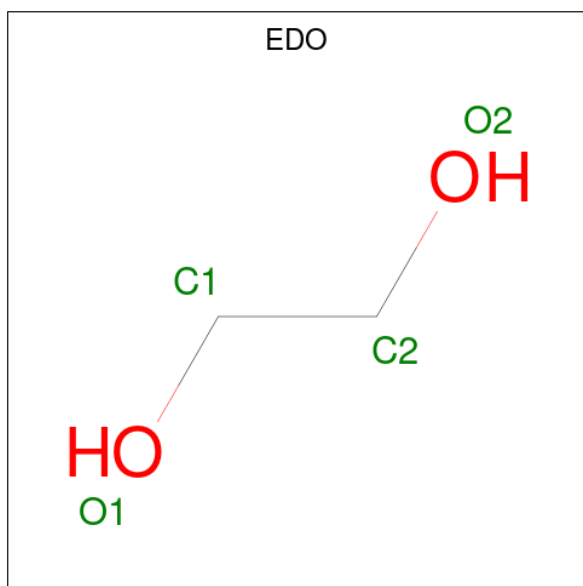
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total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

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

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

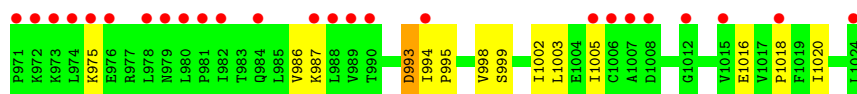
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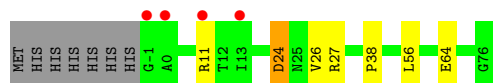
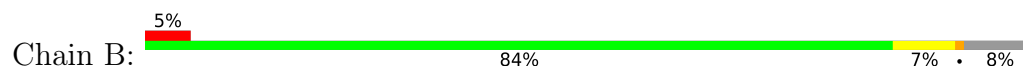
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 4 2 2	0	0
7	C	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0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

- Molecule 8 is water.

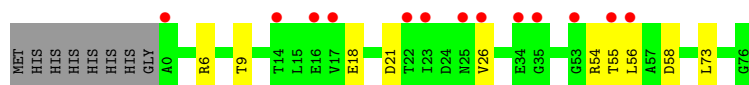
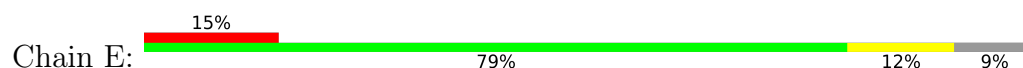
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	471	Total O 471 471	0	0
8	B	29	Total O 29 29	0	0
8	C	9	Total O 9 9	0	0
8	D	373	Total O 373 373	0	0
8	E	21	Total O 21 21	0	0
8	F	6	Total O 6 6	0	0



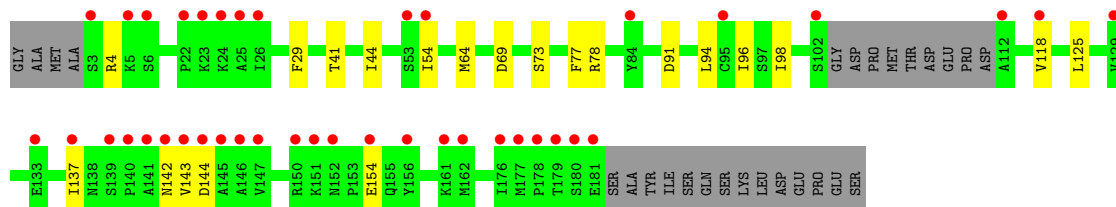
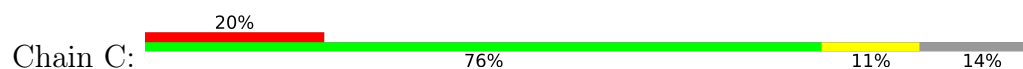
• Molecule 2: Ubiquitin



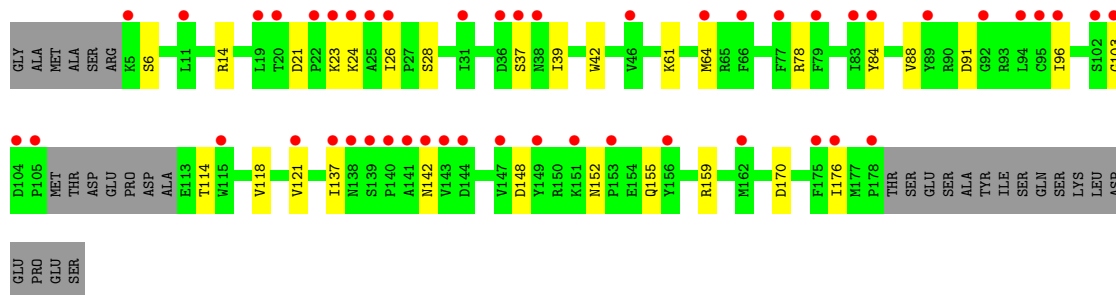
• Molecule 2: Ubiquitin



• Molecule 3: Ubiquitin-conjugating enzyme E2-34 kDa



• Molecule 3: Ubiquitin-conjugating enzyme E2-34 kDa



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	126.44Å 68.54Å 171.72Å 90.00° 110.72° 90.00°	Depositor
Resolution (Å)	117.03 – 2.06 117.03 – 2.06	Depositor EDS
% Data completeness (in resolution range)	99.0 (117.03-2.06) 89.2 (117.03-2.06)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.88 (at 2.07Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.187 , 0.217 0.188 , 0.219	Depositor DCC
R_{free} test set	2005 reflections (1.20%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20705	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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, 0.333 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: EDO, MG, ATP, SO4

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	0.16	0/7962	0.34	0/10770
1	D	0.16	0/7958	0.32	0/10766
2	B	0.14	0/627	0.36	0/843
2	E	0.14	0/623	0.35	0/838
3	C	0.14	0/1417	0.30	0/1922
3	F	0.15	0/1393	0.31	0/1891
All	All	0.16	0/19980	0.33	0/27030

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	7800	0	7721	46	0
1	D	7796	0	7721	50	0
2	B	622	0	633	4	0
2	E	618	0	630	7	0
3	C	1381	0	1351	10	0
3	F	1356	0	1325	14	0
4	A	31	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	31	0	12	0	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	40	0	0	1	0
6	D	35	0	0	2	0
7	A	44	0	66	2	0
7	C	4	0	6	1	0
7	D	36	0	54	3	0
8	A	471	0	0	7	0
8	B	29	0	0	0	0
8	C	9	0	0	0	0
8	D	373	0	0	3	0
8	E	21	0	0	2	0
8	F	6	0	0	0	0
All	All	20705	0	19531	120	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.

The worst 5 of 120 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:D:930:ARG:NH1	1:D:932:ASP:OD2	2.07	0.88
1:D:1016:GLU:OE2	3:F:14:ARG:NH2	2.21	0.73
3:F:84:TYR:OH	3:F:148:ASP:OD2	2.05	0.73
3:C:143:VAL:HB	3:C:144:ASP:HA	1.72	0.71
1:D:44:LEU:HD21	1:D:65:VAL:HB	1.76	0.66

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	981/1017 (96%)	951 (97%)	29 (3%)	1 (0%)	48	44
1	D	981/1017 (96%)	948 (97%)	31 (3%)	2 (0%)	43	38
2	B	76/85 (89%)	75 (99%)	1 (1%)	0	100	100
2	E	75/85 (88%)	73 (97%)	2 (3%)	0	100	100
3	C	166/197 (84%)	161 (97%)	5 (3%)	0	100	100
3	F	163/197 (83%)	155 (95%)	7 (4%)	1 (1%)	21	13
All	All	2442/2598 (94%)	2363 (97%)	75 (3%)	4 (0%)	43	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	ASP
1	D	220	ASP
3	F	37	SER
1	D	649	GLY

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	870/893 (97%)	864 (99%)	6 (1%)	76	79
1	D	870/893 (97%)	859 (99%)	11 (1%)	61	62
2	B	67/74 (90%)	65 (97%)	2 (3%)	36	32
2	E	67/74 (90%)	67 (100%)	0	100	100
3	C	157/179 (88%)	151 (96%)	6 (4%)	29	23
3	F	154/179 (86%)	150 (97%)	4 (3%)	40	37
All	All	2185/2292 (95%)	2156 (99%)	29 (1%)	61	62

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	79	GLN

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Mol	Chain	Res	Type
3	F	137	ILE
1	D	369	LEU
1	D	993	ASP
1	D	222	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	725	ASN
1	D	844	ASN
3	F	30	HIS
1	D	997	HIS
2	B	68	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 2 are monoatomic - leaving 38 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
6	SO4	D	1203	-	4,4,4	0.24	0	6,6,6	0.08	0
7	EDO	D	1215	-	3,3,3	0.44	0	2,2,2	0.29	0
7	EDO	A	1119	-	3,3,3	0.42	0	2,2,2	0.47	0
7	EDO	D	1216	-	3,3,3	0.45	0	2,2,2	0.40	0
7	EDO	C	201	-	3,3,3	0.45	0	2,2,2	0.22	0
7	EDO	A	1112	-	3,3,3	0.38	0	2,2,2	0.44	0
6	SO4	A	1106	-	4,4,4	0.24	0	6,6,6	0.10	0
7	EDO	A	1111	-	3,3,3	0.37	0	2,2,2	0.51	0
7	EDO	D	1210	-	3,3,3	0.41	0	2,2,2	0.49	0
6	SO4	A	1110	-	4,4,4	0.24	0	6,6,6	0.06	0
7	EDO	D	1211	-	3,3,3	0.38	0	2,2,2	0.51	0
7	EDO	D	1213	-	3,3,3	0.35	0	2,2,2	0.42	0
7	EDO	D	1217	-	3,3,3	0.46	0	2,2,2	0.28	0
6	SO4	A	1105	-	4,4,4	0.23	0	6,6,6	0.07	0
7	EDO	A	1121	-	3,3,3	0.44	0	2,2,2	0.37	0
6	SO4	A	1107	-	4,4,4	0.24	0	6,6,6	0.08	0
6	SO4	D	1205	-	4,4,4	0.23	0	6,6,6	0.08	0
7	EDO	D	1212	-	3,3,3	0.42	0	2,2,2	0.39	0
4	ATP	D	1202	5	32,33,33	1.47	7 (21%)	48,52,52	1.76	9 (18%)
7	EDO	A	1115	-	3,3,3	0.39	0	2,2,2	0.60	0
6	SO4	A	1104	-	4,4,4	0.24	0	6,6,6	0.06	0
7	EDO	D	1214	-	3,3,3	0.39	0	2,2,2	0.52	0
6	SO4	D	1206	-	4,4,4	0.23	0	6,6,6	0.10	0
6	SO4	D	1207	-	4,4,4	0.24	0	6,6,6	0.08	0
7	EDO	A	1120	-	3,3,3	0.49	0	2,2,2	0.24	0
7	EDO	D	1218	-	3,3,3	0.43	0	2,2,2	0.40	0
6	SO4	A	1108	-	4,4,4	0.25	0	6,6,6	0.08	0
6	SO4	A	1103	-	4,4,4	0.23	0	6,6,6	0.09	0
6	SO4	D	1209	-	4,4,4	0.24	0	6,6,6	0.07	0
7	EDO	A	1117	-	3,3,3	0.44	0	2,2,2	0.24	0
7	EDO	A	1113	-	3,3,3	0.40	0	2,2,2	0.36	0
6	SO4	A	1109	-	4,4,4	0.24	0	6,6,6	0.12	0
7	EDO	A	1118	-	3,3,3	0.43	0	2,2,2	0.43	0
4	ATP	A	1101	5	32,33,33	1.47	7 (21%)	48,52,52	1.74	9 (18%)
7	EDO	A	1114	-	3,3,3	0.41	0	2,2,2	0.42	0
6	SO4	D	1204	-	4,4,4	0.25	0	6,6,6	0.06	0
6	SO4	D	1208	-	4,4,4	0.25	0	6,6,6	0.06	0
7	EDO	A	1116	-	3,3,3	0.38	0	2,2,2	0.39	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
7	EDO	D	1215	-	-	0/1/1/1	-
7	EDO	A	1119	-	-	0/1/1/1	-
7	EDO	D	1216	-	-	0/1/1/1	-
7	EDO	C	201	-	-	0/1/1/1	-
7	EDO	A	1112	-	-	0/1/1/1	-
7	EDO	A	1111	-	-	0/1/1/1	-
7	EDO	D	1210	-	-	0/1/1/1	-
7	EDO	D	1211	-	-	0/1/1/1	-
7	EDO	D	1213	-	-	0/1/1/1	-
7	EDO	D	1217	-	-	0/1/1/1	-
7	EDO	A	1121	-	-	0/1/1/1	-
7	EDO	D	1212	-	-	0/1/1/1	-
4	ATP	D	1202	5	-	4/22/38/38	0/3/3/3
7	EDO	A	1115	-	-	0/1/1/1	-
7	EDO	D	1214	-	-	0/1/1/1	-
7	EDO	A	1120	-	-	0/1/1/1	-
7	EDO	D	1218	-	-	0/1/1/1	-
7	EDO	A	1117	-	-	0/1/1/1	-
7	EDO	A	1113	-	-	0/1/1/1	-
7	EDO	A	1118	-	-	0/1/1/1	-
4	ATP	A	1101	5	-	5/22/38/38	0/3/3/3
7	EDO	A	1114	-	-	0/1/1/1	-
7	EDO	A	1116	-	-	0/1/1/1	-

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1202	ATP	C5-C4	4.69	1.47	1.39
4	A	1101	ATP	C5-C4	4.46	1.47	1.39
4	A	1101	ATP	PA-O3A	2.83	1.62	1.59
4	D	1202	ATP	C5-C6	2.81	1.48	1.41
4	A	1101	ATP	C5-C6	2.70	1.48	1.41

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1202	ATP	C5-C4-N3	-5.75	118.80	126.72
4	A	1101	ATP	C5-C4-N3	-5.63	118.97	126.72
4	D	1202	ATP	N3-C4-N9	4.62	135.02	127.17
4	A	1101	ATP	N3-C4-N9	4.58	134.96	127.17
4	A	1101	ATP	C2-N3-C4	3.69	120.85	111.83

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

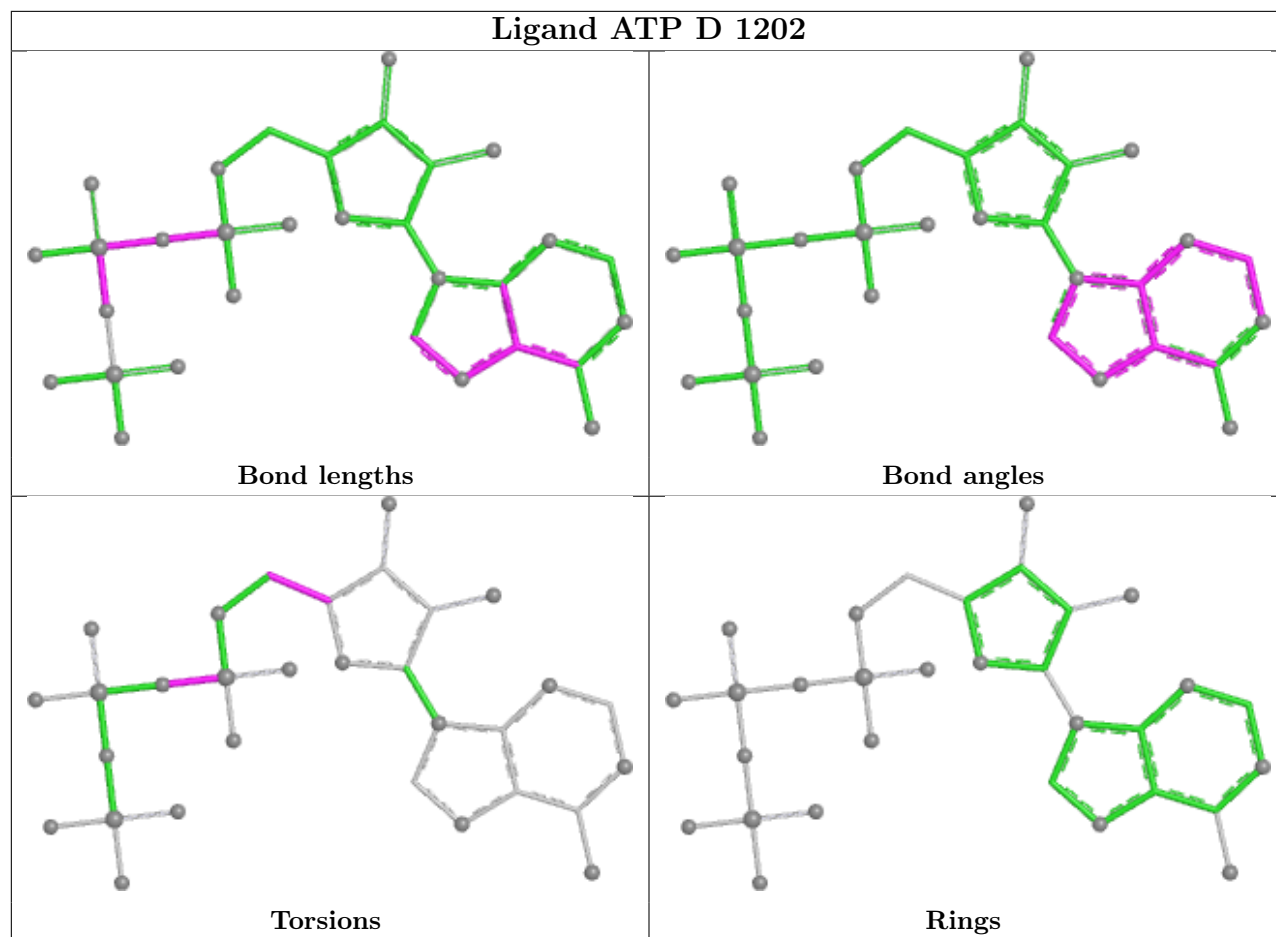
Mol	Chain	Res	Type	Atoms
4	D	1202	ATP	O4'-C4'-C5'-O5'
4	D	1202	ATP	C3'-C4'-C5'-O5'
4	A	1101	ATP	O4'-C4'-C5'-O5'
4	A	1101	ATP	C3'-C4'-C5'-O5'
4	D	1202	ATP	PB-O3A-PA-O2A

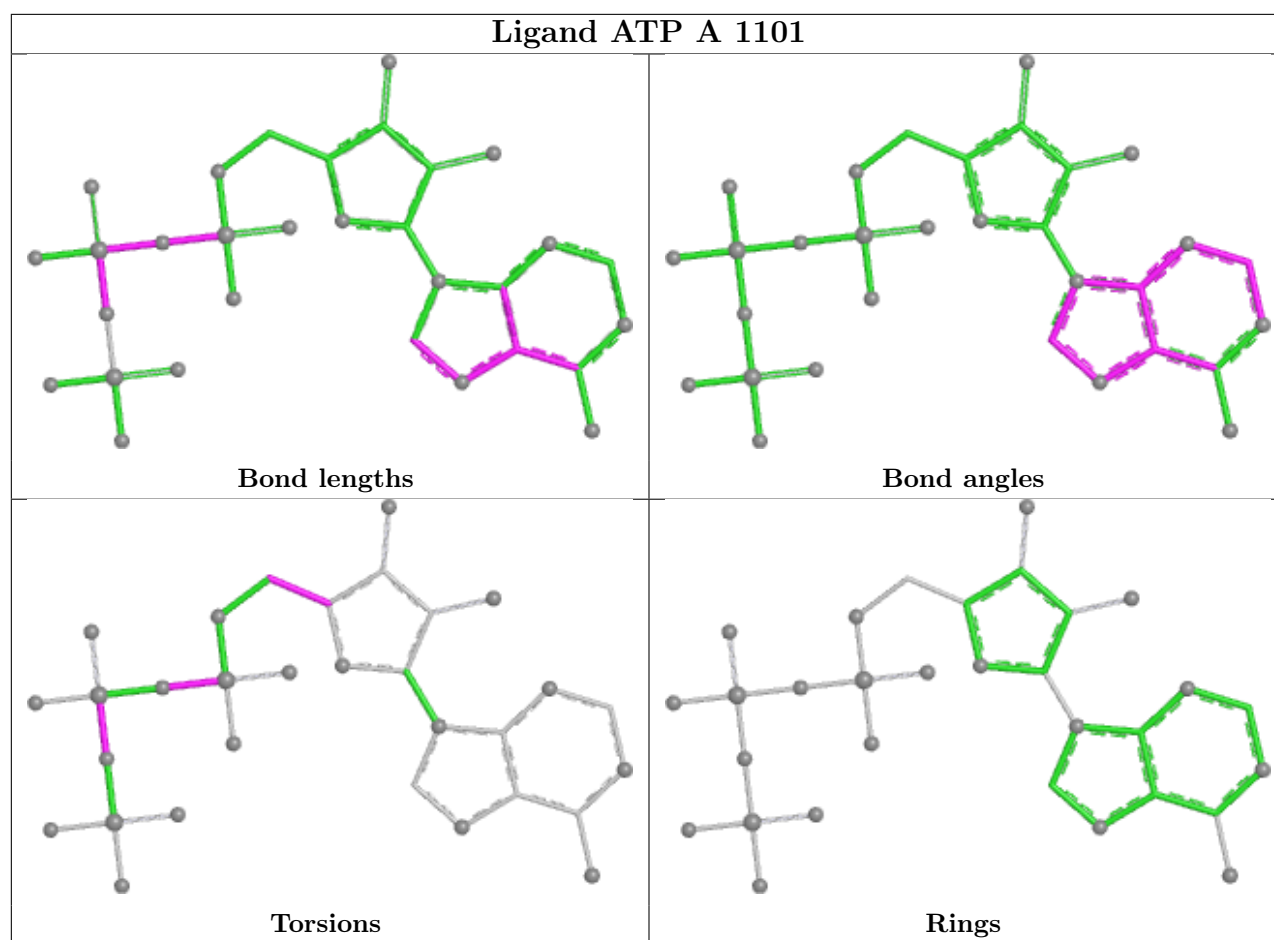
There are no ring outliers.

8 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	201	EDO	1	0
7	A	1112	EDO	1	0
7	D	1213	EDO	1	0
6	A	1107	SO4	1	0
7	D	1212	EDO	1	0
7	D	1214	EDO	1	0
7	A	1114	EDO	1	0
6	D	1204	SO4	2	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, 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	987/1017 (97%)	-0.03	37 (3%)	45 46	19, 33, 77, 152	0
1	D	987/1017 (97%)	0.32	103 (10%)	11 11	19, 39, 100, 137	0
2	B	78/85 (91%)	0.63	4 (5%)	33 33	23, 53, 81, 95	0
2	E	77/85 (90%)	1.03	13 (16%)	4 4	24, 59, 95, 106	0
3	C	170/197 (86%)	1.45	40 (23%)	2 1	43, 66, 104, 139	0
3	F	167/197 (84%)	1.69	48 (28%)	1 1	52, 79, 111, 156	0
All	All	2466/2598 (94%)	0.38	245 (9%)	13 13	19, 40, 98, 156	0

The worst 5 of 245 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	141	ALA	6.4
3	F	105	PRO	6.1
3	C	143	VAL	5.7
1	A	798	ILE	5.4
1	D	978	LEU	5.0

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

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, 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	B-factors(Å ²)	Q<0.9
6	SO4	A	1105	5/5	0.63	0.19	138,140,140,140	0
6	SO4	A	1103	5/5	0.66	0.17	105,106,107,108	0
6	SO4	D	1209	5/5	0.69	0.13	99,101,102,103	0
6	SO4	A	1107	5/5	0.71	0.21	122,124,125,127	0
6	SO4	A	1104	5/5	0.74	0.20	137,138,139,140	0
6	SO4	D	1208	5/5	0.75	0.16	68,72,76,76	0
6	SO4	D	1207	5/5	0.75	0.13	82,85,86,87	0
6	SO4	D	1206	5/5	0.77	0.13	91,93,96,98	0
7	EDO	A	1112	4/4	0.80	0.18	42,47,49,57	0
7	EDO	A	1121	4/4	0.80	0.16	64,66,68,69	0
7	EDO	D	1218	4/4	0.80	0.21	57,59,60,60	0
6	SO4	D	1203	5/5	0.81	0.24	126,127,129,130	0
6	SO4	A	1110	5/5	0.81	0.13	82,85,86,89	0
7	EDO	A	1119	4/4	0.82	0.17	51,51,56,60	0
6	SO4	D	1205	5/5	0.82	0.14	105,105,106,107	0
6	SO4	A	1108	5/5	0.82	0.14	61,72,78,78	0
7	EDO	A	1116	4/4	0.83	0.18	47,48,49,55	0
6	SO4	A	1109	5/5	0.83	0.14	86,88,92,93	0
7	EDO	A	1117	4/4	0.84	0.18	53,56,56,60	0
7	EDO	A	1115	4/4	0.84	0.18	46,51,51,52	0
7	EDO	D	1211	4/4	0.85	0.18	49,49,52,57	0
7	EDO	D	1213	4/4	0.86	0.16	42,45,51,51	0
7	EDO	D	1214	4/4	0.86	0.17	58,58,61,65	0
6	SO4	A	1106	5/5	0.86	0.22	110,111,111,113	0
6	SO4	D	1204	5/5	0.87	0.18	120,121,121,122	0
7	EDO	D	1217	4/4	0.87	0.15	47,51,52,52	0
7	EDO	D	1210	4/4	0.87	0.14	36,37,40,43	0
7	EDO	A	1118	4/4	0.88	0.15	33,38,49,60	0
7	EDO	D	1212	4/4	0.88	0.15	51,54,57,61	0
7	EDO	A	1111	4/4	0.91	0.10	27,28,34,36	0
7	EDO	D	1215	4/4	0.91	0.13	52,56,63,68	0
7	EDO	A	1113	4/4	0.92	0.13	32,38,42,45	0
7	EDO	A	1114	4/4	0.92	0.14	46,47,48,55	0
7	EDO	A	1120	4/4	0.93	0.10	34,42,43,45	0
7	EDO	D	1216	4/4	0.93	0.11	30,39,42,43	0
7	EDO	C	201	4/4	0.96	0.07	44,44,44,45	0
4	ATP	A	1101	31/31	0.98	0.04	18,24,28,31	0

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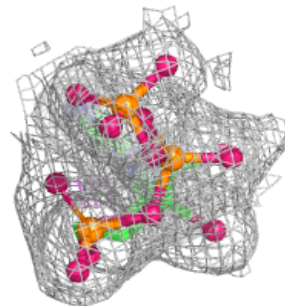
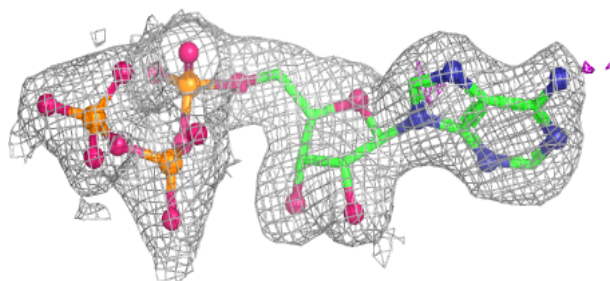
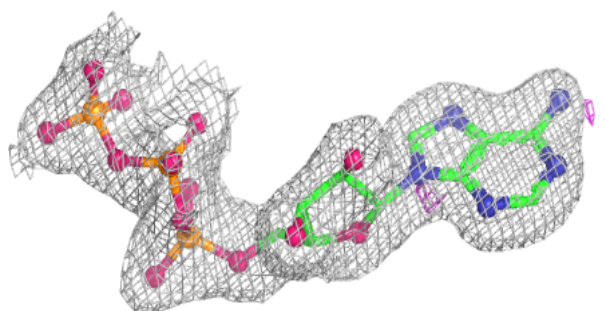
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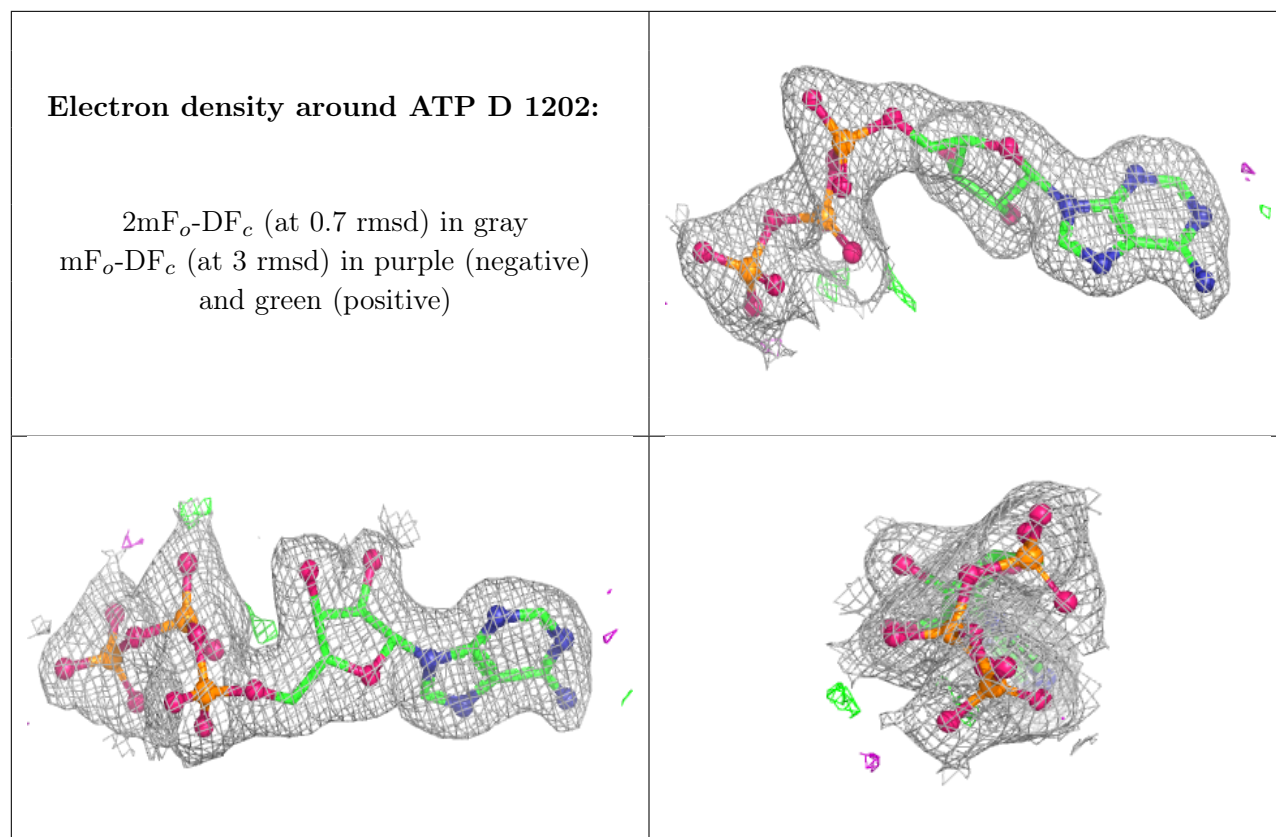
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ATP	D	1202	31/31	0.99	0.05	18,28,36,38	0
5	MG	A	1102	1/1	0.99	0.04	24,24,24,24	0
5	MG	D	1201	1/1	0.99	0.02	31,31,31,31	0

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 1101:

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 [i](#)

There are no such residues in this entry.