wwPDB X-ray Structure Validation Summary Report (i)

Oct 26, 2026 – 11:18 PM JST

PDB ID	:	9XDS / pdb_00009xds
Title	:	Crystal Structure of Fructose-1,6-bisphosphatase Complexed with a Covalent Inhibitor
Authors	:	Cao, H.; Huang, Z.; Hu, X.; Li, W.; Chang, M.; Zhang, Y.; Chu, Y.
Deposited on	:	2025-10-27
Resolution	:	2.08 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 (i) 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 (i)) were used in the production of this report:

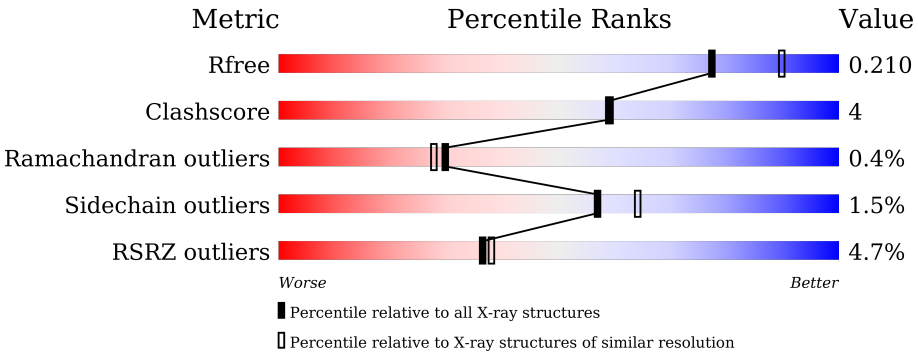
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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 <=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	3% 87% 5% • 8%
1	B	346	3% 84% 9% • 6%
1	C	346	4% 81% 10% • 8%
1	D	346	7% 84% 9% • 6%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	A1EZR	A	409	-	X	-	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 10726 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 Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2435	1553	406	459	17			
1	B	325	Total	C	N	O	S	0	0	0
			2481	1577	416	471	17			
1	C	319	Total	C	N	O	S	0	0	0
			2441	1556	409	459	17			
1	D	325	Total	C	N	O	S	0	0	0
			2486	1581	417	471	17			

There are 44 discrepancies between the modelled and reference sequences:

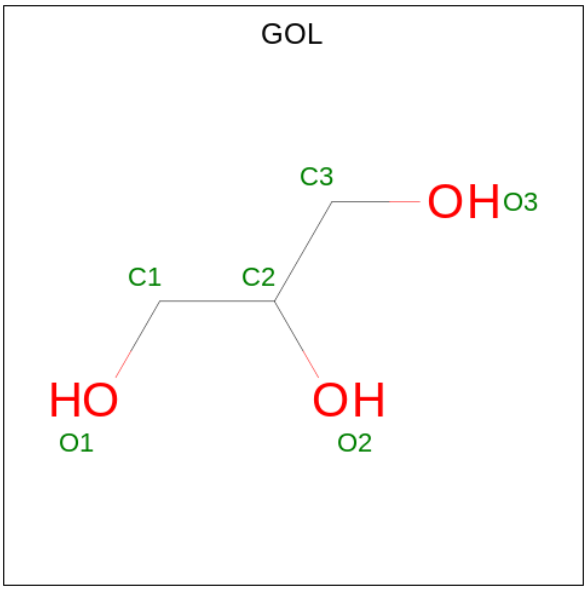
Chain	Residue	Modelled	Actual	Comment	Reference
A	217	LYS	ARG	variant	UNP P09467
A	336	LEU	-	expression tag	UNP P09467
A	337	GLU	-	expression tag	UNP P09467
A	338	HIS	-	expression tag	UNP P09467
A	339	HIS	-	expression tag	UNP P09467
A	340	HIS	-	expression tag	UNP P09467
A	341	HIS	-	expression tag	UNP P09467
A	342	HIS	-	expression tag	UNP P09467
A	343	HIS	-	expression tag	UNP P09467
A	344	HIS	-	expression tag	UNP P09467
A	345	HIS	-	expression tag	UNP P09467
B	217	LYS	ARG	variant	UNP P09467
B	336	LEU	-	expression tag	UNP P09467
B	337	GLU	-	expression tag	UNP P09467
B	338	HIS	-	expression tag	UNP P09467
B	339	HIS	-	expression tag	UNP P09467
B	340	HIS	-	expression tag	UNP P09467
B	341	HIS	-	expression tag	UNP P09467
B	342	HIS	-	expression tag	UNP P09467
B	343	HIS	-	expression tag	UNP P09467
B	344	HIS	-	expression tag	UNP P09467

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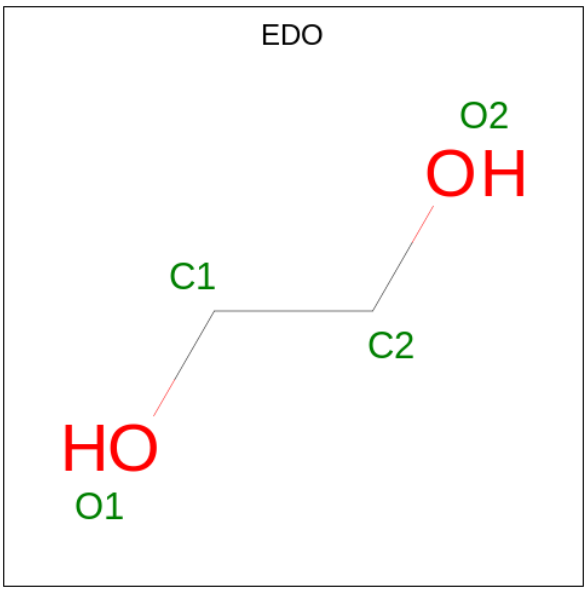
Chain	Residue	Modelled	Actual	Comment	Reference
B	345	HIS	-	expression tag	UNP P09467
C	217	LYS	ARG	variant	UNP P09467
C	336	LEU	-	expression tag	UNP P09467
C	337	GLU	-	expression tag	UNP P09467
C	338	HIS	-	expression tag	UNP P09467
C	339	HIS	-	expression tag	UNP P09467
C	340	HIS	-	expression tag	UNP P09467
C	341	HIS	-	expression tag	UNP P09467
C	342	HIS	-	expression tag	UNP P09467
C	343	HIS	-	expression tag	UNP P09467
C	344	HIS	-	expression tag	UNP P09467
C	345	HIS	-	expression tag	UNP P09467
D	217	LYS	ARG	variant	UNP P09467
D	336	LEU	-	expression tag	UNP P09467
D	337	GLU	-	expression tag	UNP P09467
D	338	HIS	-	expression tag	UNP P09467
D	339	HIS	-	expression tag	UNP P09467
D	340	HIS	-	expression tag	UNP P09467
D	341	HIS	-	expression tag	UNP P09467
D	342	HIS	-	expression tag	UNP P09467
D	343	HIS	-	expression tag	UNP P09467
D	344	HIS	-	expression tag	UNP P09467
D	345	HIS	-	expression tag	UNP P09467

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



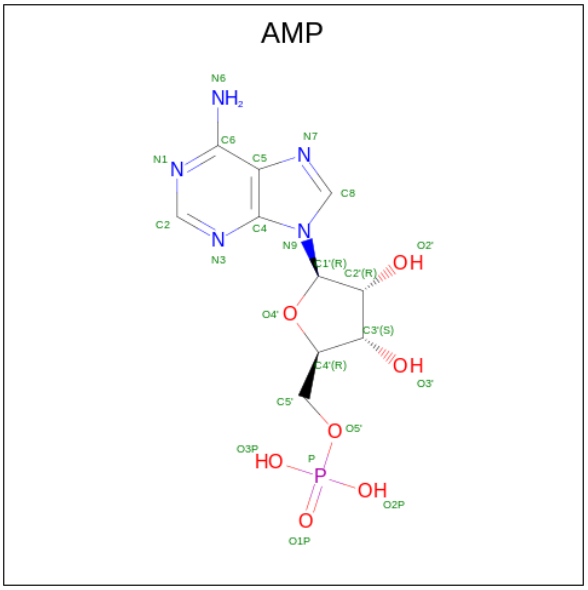
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0

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

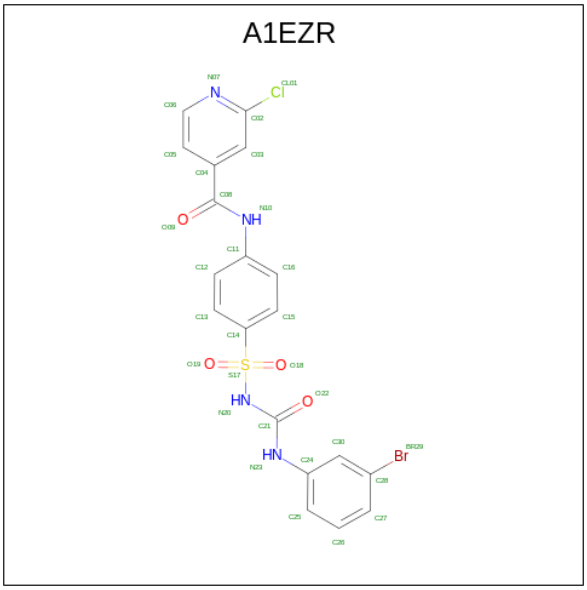
- Molecule 4 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O P 23 10 5 7 1	0	0
4	B	1	Total C N O P 23 10 5 7 1	0	0
4	C	1	Total C N O P 23 10 5 7 1	0	0
4	D	1	Total C N O P 23 10 5 7 1	0	0

- Molecule 5 is {N}-[4-[(3-bromophenyl)carbamoylsulfamoyl]phenyl]-2-chloranyl-pyridine-4-c

arboxamide (CCD ID: A1EZR) (formula: $C_{19}H_{14}BrClN_4O_4S$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	A	1	Total	Br	C	Cl	N	O	S	0	0
			30	1	19	1	4	4	1		
5	B	1	Total	Br	C	Cl	N	O	S	0	0
			30	1	19	1	4	4	1		
5	C	1	Total	Br	C	Cl	N	O	S	0	0
			30	1	19	1	4	4	1		
5	D	1	Total	Br	C	Cl	N	O	S	0	0
			30	1	19	1	4	4	1		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 7 is water.

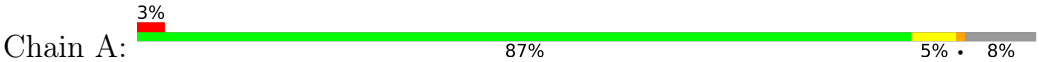
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	172	Total	O	0	0
			172	172		
7	B	151	Total	O	0	0
			151	151		
7	C	136	Total	O	0	0
			136	136		
7	D	119	Total	O	0	0
			119	119		

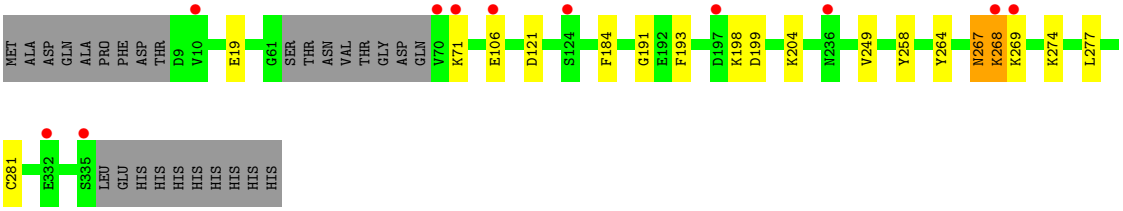
3 Residue-property plots (i)

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: Fructose-1,6-bisphosphatase 1

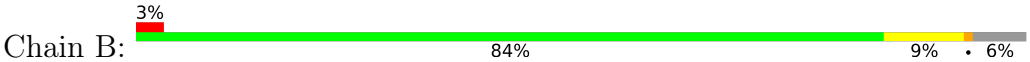
Chain A:

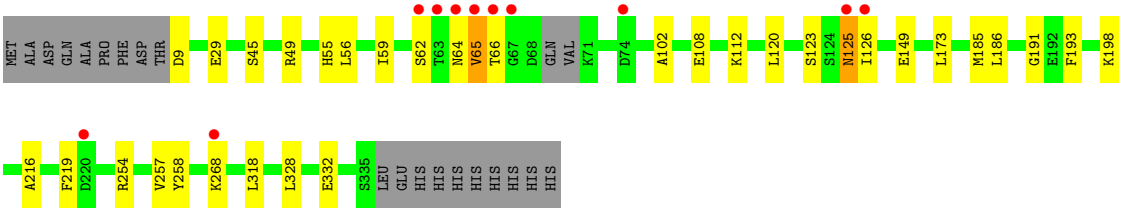




• Molecule 1: Fructose-1,6-bisphosphatase 1

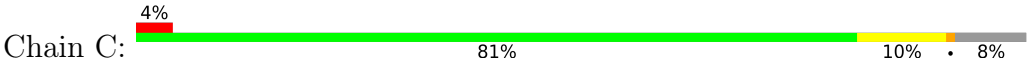
Chain B:

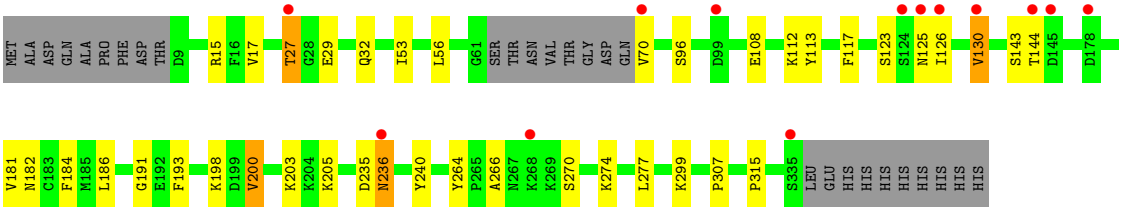




• Molecule 1: Fructose-1,6-bisphosphatase 1

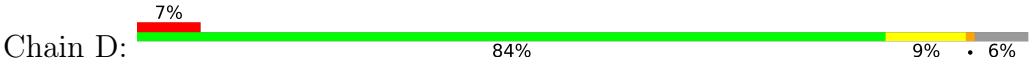
Chain C:

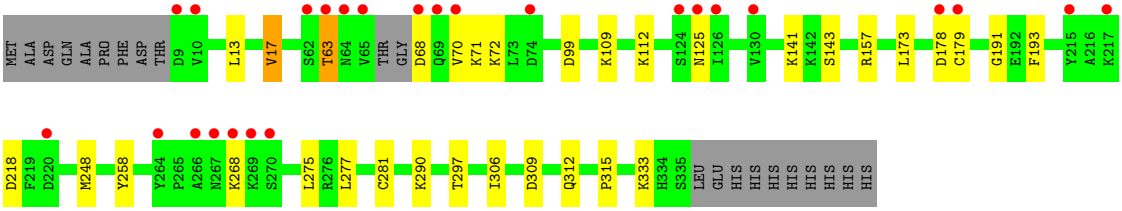




• Molecule 1: Fructose-1,6-bisphosphatase 1

Chain D:




MET
ALA
ASP
GLN
ALA
PRO
PHE
ASP
THR
D9
V10
L13
V17
S62
T63
N64
V65
THR
GLY
D68
Q69
V70
K71
K72
L73
D74
D99
K109
K112
S124
N125
I126
V130
K141
K142
S143
R157
L173
D178
C179
G191
E192
F193
Y215
A216
K217
D218
F219
D220
M248
Y258
Y264
P265
A266
N267
K268
K269
S270
L275
R276
L277
C281
K290
T297
I306
D309
Q312
P315
K333
H334
S335
LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	67.67Å 83.33Å 279.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.64 – 2.08 71.64 – 2.08	Depositor EDS
% Data completeness (in resolution range)	98.6 (71.64-2.08) 98.6 (71.64-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.37 (at 2.08Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.176 , 0.212 0.176 , 0.210	Depositor DCC
R_{free} test set	4692 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 40.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10726	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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, 0.333 respectively for untwinned datasets, and 0.375, 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, EDO, GOL, A1EZR, AMP

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	0.34	0/2478	0.52	0/3347
1	B	0.32	0/2524	0.52	0/3409
1	C	0.32	0/2484	0.55	0/3354
1	D	0.29	0/2529	0.50	0/3416
All	All	0.32	0/10015	0.52	0/13526

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

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	2435	0	2480	16	0
1	B	2481	0	2523	25	0
1	C	2441	0	2491	27	0
1	D	2486	0	2530	26	0
2	A	30	0	40	1	0
2	D	30	0	40	4	0
3	A	12	0	18	0	0
3	C	16	0	24	1	0
3	D	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	23	0	12	0	0
4	B	23	0	12	0	0
4	C	23	0	12	2	0
4	D	23	0	12	0	0
5	A	30	0	0	1	0
5	B	30	0	0	1	0
5	C	30	0	0	0	0
5	D	30	0	0	0	0
6	A	1	0	0	0	0
7	A	172	0	0	2	0
7	B	151	0	0	4	0
7	C	136	0	0	4	0
7	D	119	0	0	4	0
All	All	10726	0	10200	89	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 4.

The worst 5 of 89 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:179:CYS:SG	7:D:613:HOH:O	2.29	0.90
1:B:123:SER:HA	1:B:126:ILE:HG23	1.56	0.88
1:C:182:ASN:OD1	1:C:198:LYS:HD2	1.87	0.74
1:B:268:LYS:NZ	7:B:501:HOH:O	2.21	0.73
1:C:125:ASN:ND2	7:C:501:HOH:O	2.20	0.73

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	315/346 (91%)	308 (98%)	5 (2%)	2 (1%)	21	17
1	B	321/346 (93%)	311 (97%)	8 (2%)	2 (1%)	21	17
1	C	315/346 (91%)	307 (98%)	7 (2%)	1 (0%)	36	36
1	D	321/346 (93%)	312 (97%)	9 (3%)	0	100	100
All	All	1272/1384 (92%)	1238 (97%)	29 (2%)	5 (0%)	30	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	LYS
1	B	62	SER
1	C	236	ASN
1	B	65	VAL
1	A	267	ASN

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	265/290 (91%)	265 (100%)	0	100	100
1	B	271/290 (93%)	267 (98%)	4 (2%)	57	64
1	C	266/290 (92%)	259 (97%)	7 (3%)	40	45
1	D	272/290 (94%)	267 (98%)	5 (2%)	51	58
All	All	1074/1160 (93%)	1058 (98%)	16 (2%)	57	64

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	125	ASN
1	D	70	VAL
1	C	143	SER
1	D	63	THR
1	C	130	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	236	ASN
1	B	101	HIS
1	B	125	ASN
1	D	64	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 27 ligands modelled in this entry, 1 is monoatomic - leaving 26 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	GOL	A	402	-	5,5,5	0.95	0	5,5,5	1.34	1 (20%)
2	GOL	D	405	-	5,5,5	0.93	0	5,5,5	0.98	0
4	AMP	D	408	-	25,25,25	1.38	3 (12%)	38,38,38	1.85	6 (15%)
2	GOL	A	401	-	5,5,5	1.25	1 (20%)	5,5,5	0.88	0
3	EDO	C	404	-	3,3,3	0.52	0	2,2,2	0.38	0
2	GOL	A	410	-	5,5,5	1.01	0	5,5,5	0.88	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	GOL	D	404	-	5,5,5	0.88	0	5,5,5	1.01	0
3	EDO	C	401	-	3,3,3	0.45	0	2,2,2	0.59	0
3	EDO	C	403	-	3,3,3	0.51	0	2,2,2	0.47	0
2	GOL	A	403	-	5,5,5	1.09	0	5,5,5	0.96	0
4	AMP	A	408	-	25,25,25	1.29	3 (12%)	38,38,38	1.90	8 (21%)
3	EDO	D	406	-	3,3,3	0.48	0	2,2,2	0.35	0
4	AMP	C	406	-	25,25,25	1.46	4 (16%)	38,38,38	1.86	6 (15%)
3	EDO	A	406	-	3,3,3	0.47	0	2,2,2	0.55	0
2	GOL	D	401	-	5,5,5	1.05	0	5,5,5	0.93	0
3	EDO	A	407	-	3,3,3	0.47	0	2,2,2	0.45	0
3	EDO	C	402	-	3,3,3	0.46	0	2,2,2	0.48	0
2	GOL	D	402	-	5,5,5	0.85	0	5,5,5	1.06	0
4	AMP	B	402	-	25,25,25	1.36	3 (12%)	38,38,38	1.79	8 (21%)
3	EDO	A	405	-	3,3,3	0.45	0	2,2,2	0.46	0
2	GOL	D	403	-	5,5,5	0.69	0	5,5,5	1.04	0
2	GOL	A	404	-	5,5,5	0.84	0	5,5,5	0.89	0
5	A1EZR	C	405	-	32,32,32	1.99	13 (40%)	44,45,45	2.58	11 (25%)
5	A1EZR	B	401	-	32,32,32	1.94	11 (34%)	44,45,45	2.61	9 (20%)
5	A1EZR	D	407	-	32,32,32	1.88	9 (28%)	44,45,45	2.88	16 (36%)
5	A1EZR	A	409	-	32,32,32	2.15	13 (40%)	44,45,45	3.67	18 (40%)

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	GOL	A	402	-	-	2/4/4/4	-
2	GOL	D	405	-	-	4/4/4/4	-
4	AMP	D	408	-	-	0/10/26/26	0/3/3/3
2	GOL	A	401	-	-	2/4/4/4	-
3	EDO	C	404	-	-	0/1/1/1	-
2	GOL	A	410	-	-	2/4/4/4	-
2	GOL	D	404	-	-	0/4/4/4	-
3	EDO	C	401	-	-	0/1/1/1	-
3	EDO	C	403	-	-	0/1/1/1	-
2	GOL	A	403	-	-	0/4/4/4	-
4	AMP	A	408	-	-	0/10/26/26	0/3/3/3
3	EDO	D	406	-	-	1/1/1/1	-
4	AMP	C	406	-	-	0/10/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	406	-	-	1/1/1/1	-
2	GOL	D	401	-	-	1/4/4/4	-
3	EDO	A	407	-	-	1/1/1/1	-
3	EDO	C	402	-	-	0/1/1/1	-
2	GOL	D	402	-	-	0/4/4/4	-
4	AMP	B	402	-	-	0/10/26/26	0/3/3/3
3	EDO	A	405	-	-	1/1/1/1	-
2	GOL	D	403	-	-	2/4/4/4	-
2	GOL	A	404	-	-	2/4/4/4	-
5	A1EZR	C	405	-	-	4/23/23/23	0/3/3/3
5	A1EZR	B	401	-	-	2/23/23/23	0/3/3/3
5	A1EZR	D	407	-	-	8/23/23/23	0/3/3/3
5	A1EZR	A	409	-	-	13/23/23/23	0/3/3/3

The worst 5 of 60 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	406	AMP	C5-C4	4.90	1.48	1.39
4	D	408	AMP	C5-C4	4.72	1.47	1.39
5	D	407	A1EZR	C21-N23	4.61	1.47	1.37
5	A	409	A1EZR	C21-N23	4.49	1.46	1.37
4	B	402	AMP	C5-C4	4.48	1.47	1.39

The worst 5 of 83 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	409	A1EZR	O19-S17-O18	-16.84	98.84	119.55
5	D	407	A1EZR	O19-S17-O18	-13.46	103.01	119.55
5	B	401	A1EZR	O19-S17-O18	-12.81	103.80	119.55
5	C	405	A1EZR	O19-S17-O18	-12.69	103.95	119.55
4	C	406	AMP	C5-C4-N3	-6.51	118.26	126.75

There are no chirality outliers.

5 of 46 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	GOL	O2-C2-C3-O3
2	D	405	GOL	C1-C2-C3-O3
5	A	409	A1EZR	C04-C08-N10-C11
5	A	409	A1EZR	O09-C08-N10-C11

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Mol	Chain	Res	Type	Atoms
5	D	407	A1EZR	C16-C11-N10-C08

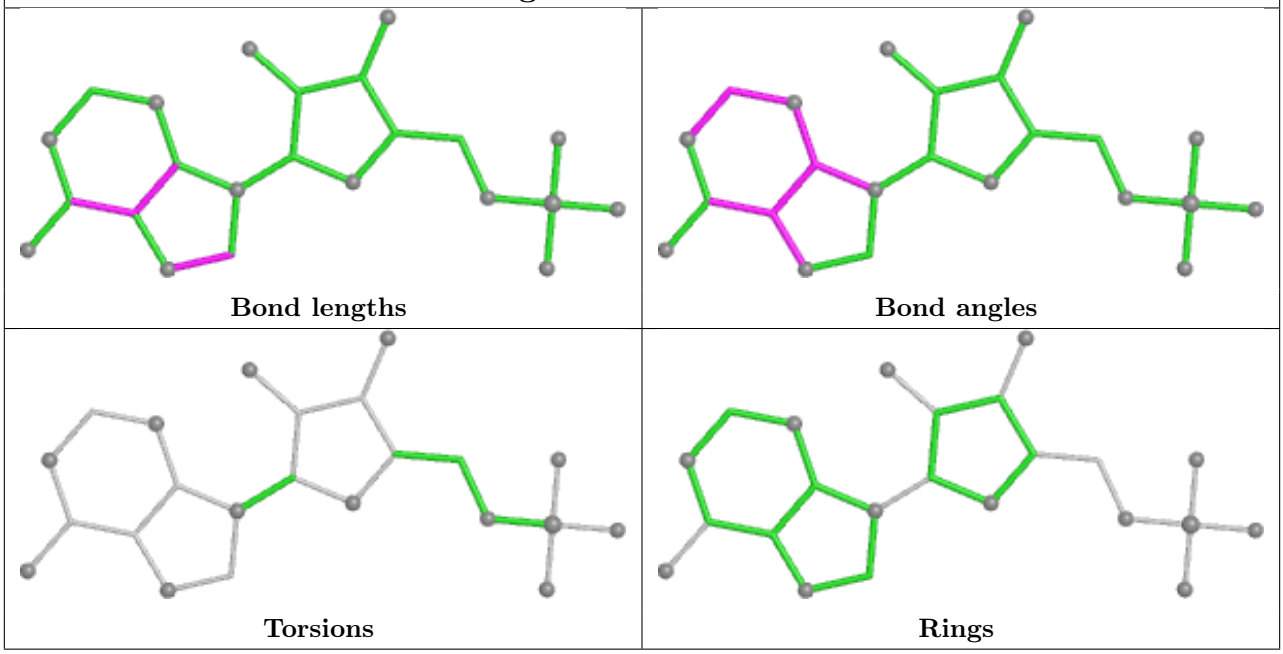
There are no ring outliers.

9 monomers are involved in 10 short contacts:

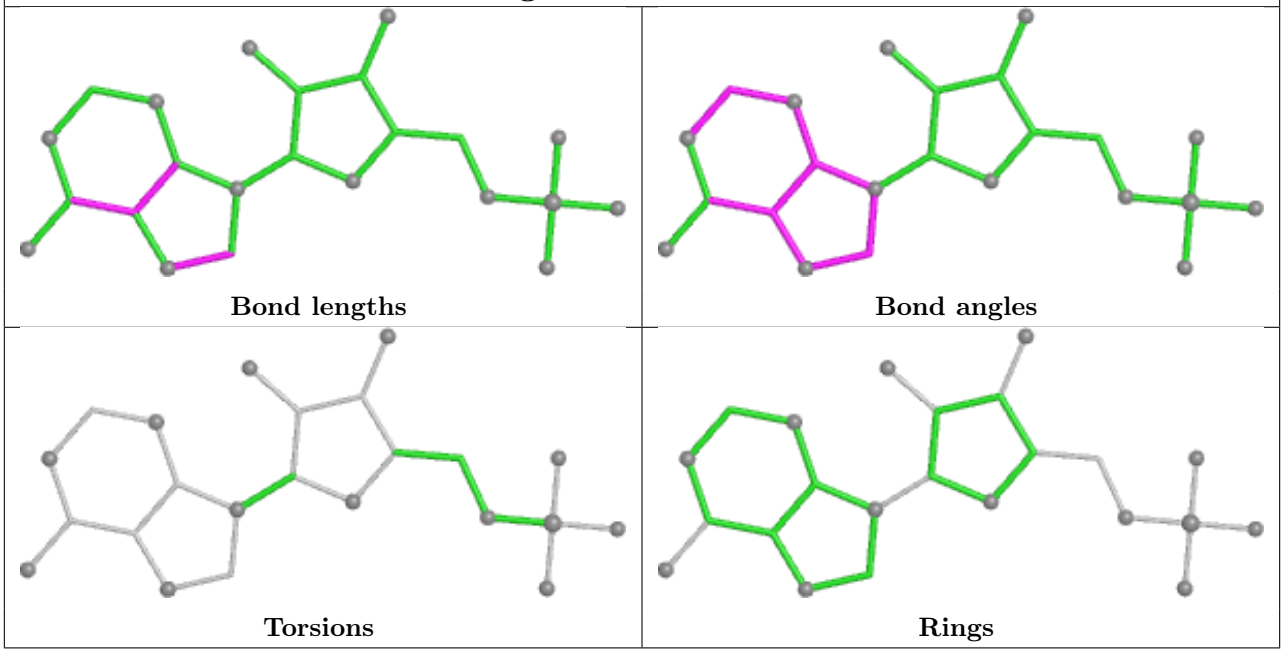
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	GOL	1	0
2	D	405	GOL	2	0
3	C	404	EDO	1	0
2	D	404	GOL	1	0
4	C	406	AMP	2	0
2	D	403	GOL	1	0
2	A	404	GOL	1	0
5	B	401	A1EZR	1	0
5	A	409	A1EZR	1	0

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 then 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.

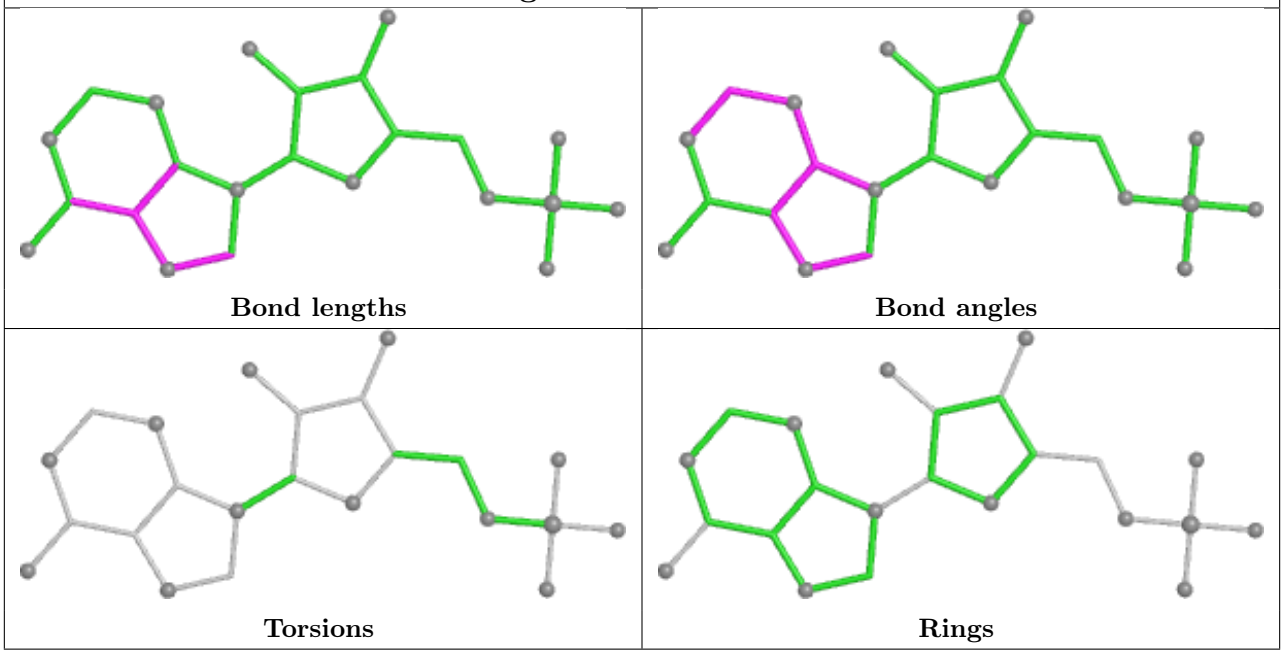
Ligand AMP D 408



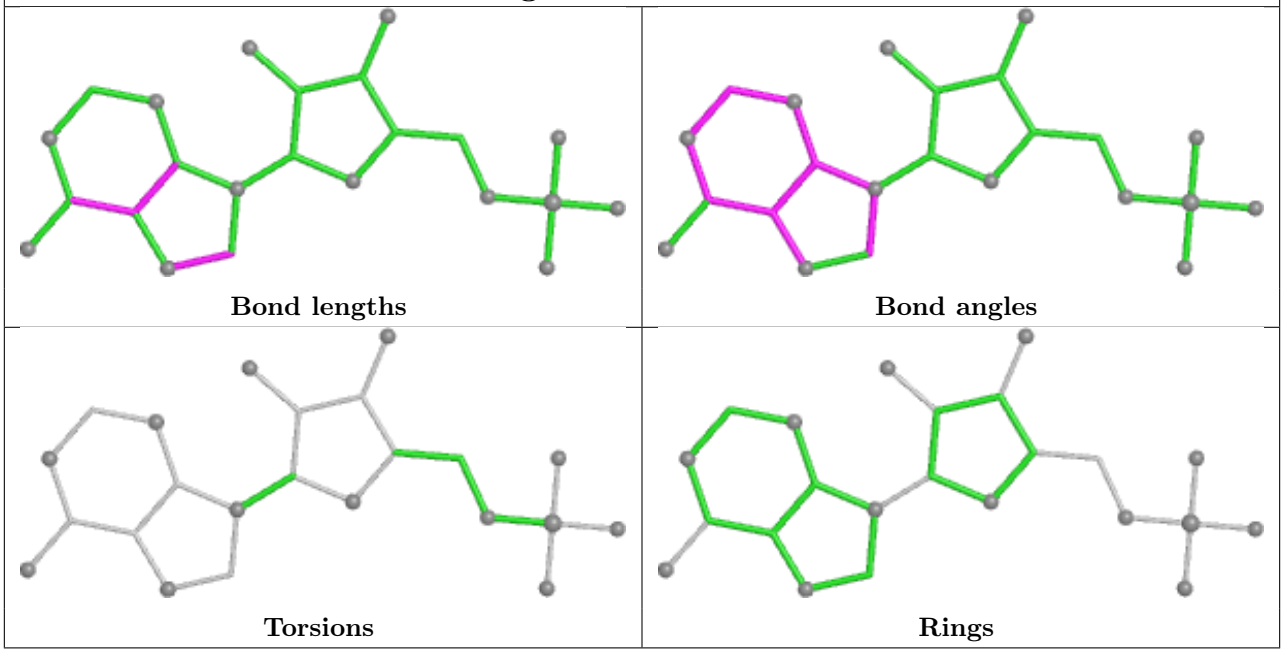
Ligand AMP A 408



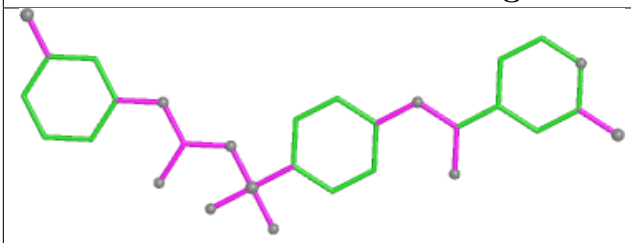
Ligand AMP C 406



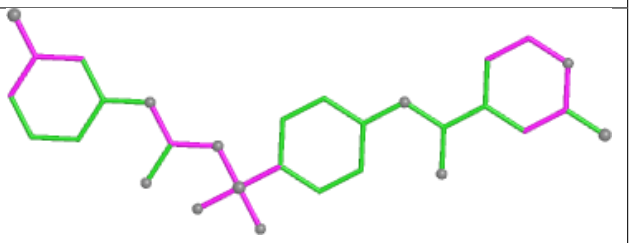
Ligand AMP B 402



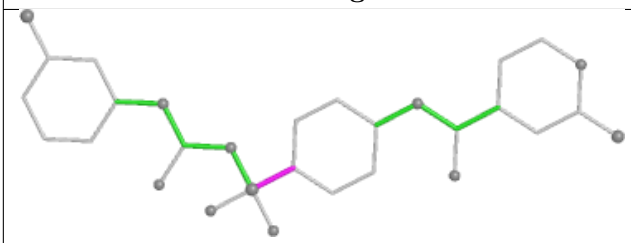
Ligand A1EZR C 405



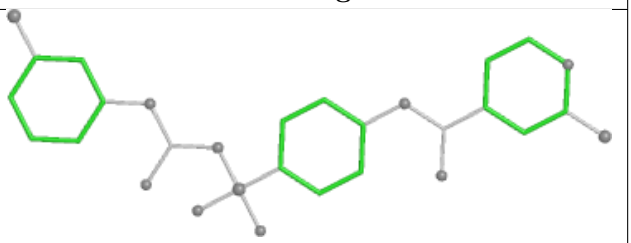
Bond lengths



Bond angles

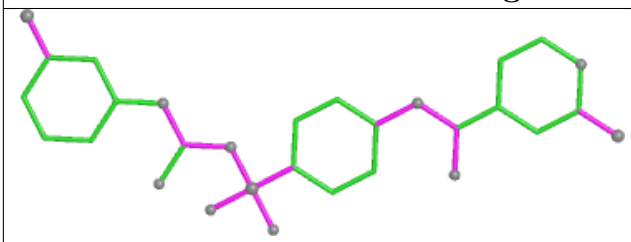


Torsions

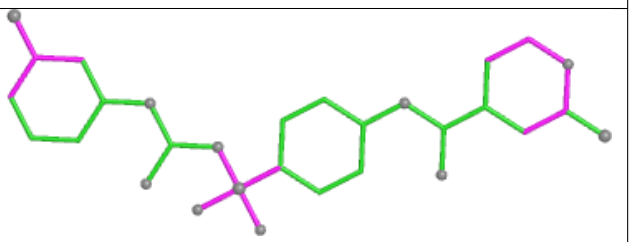


Rings

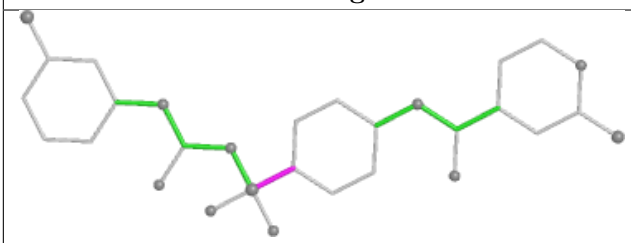
Ligand A1EZR B 401



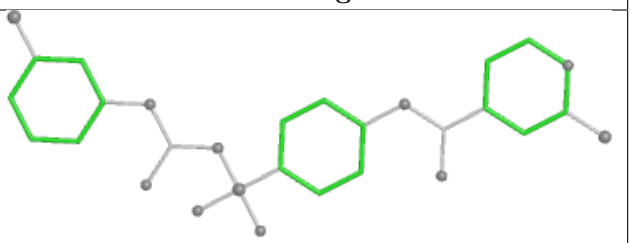
Bond lengths



Bond angles

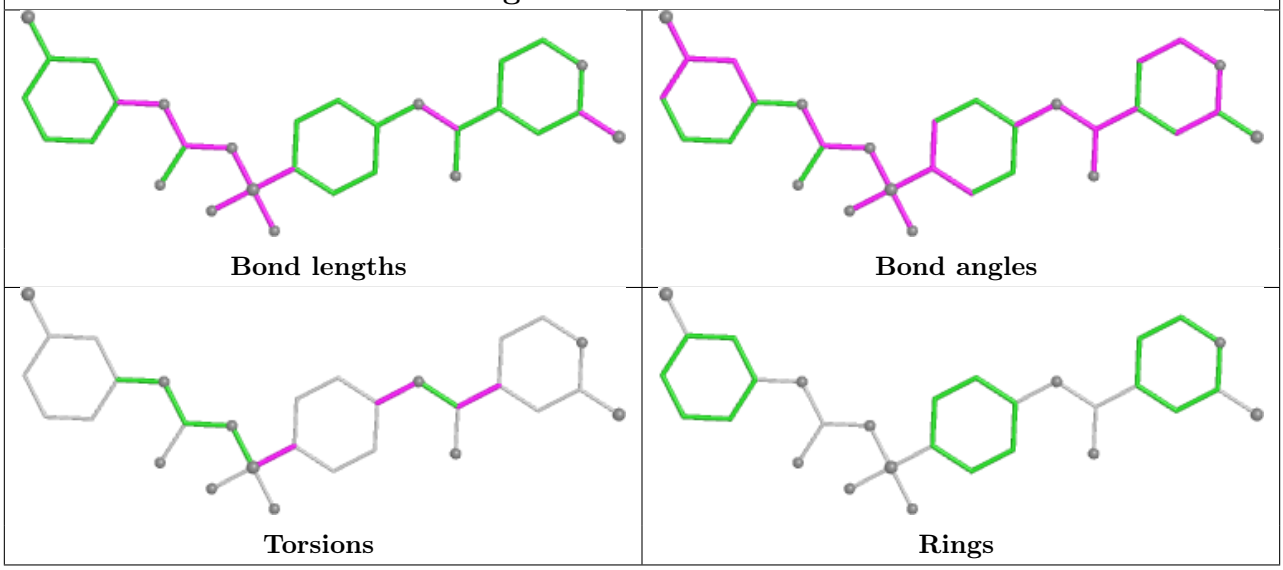


Torsions

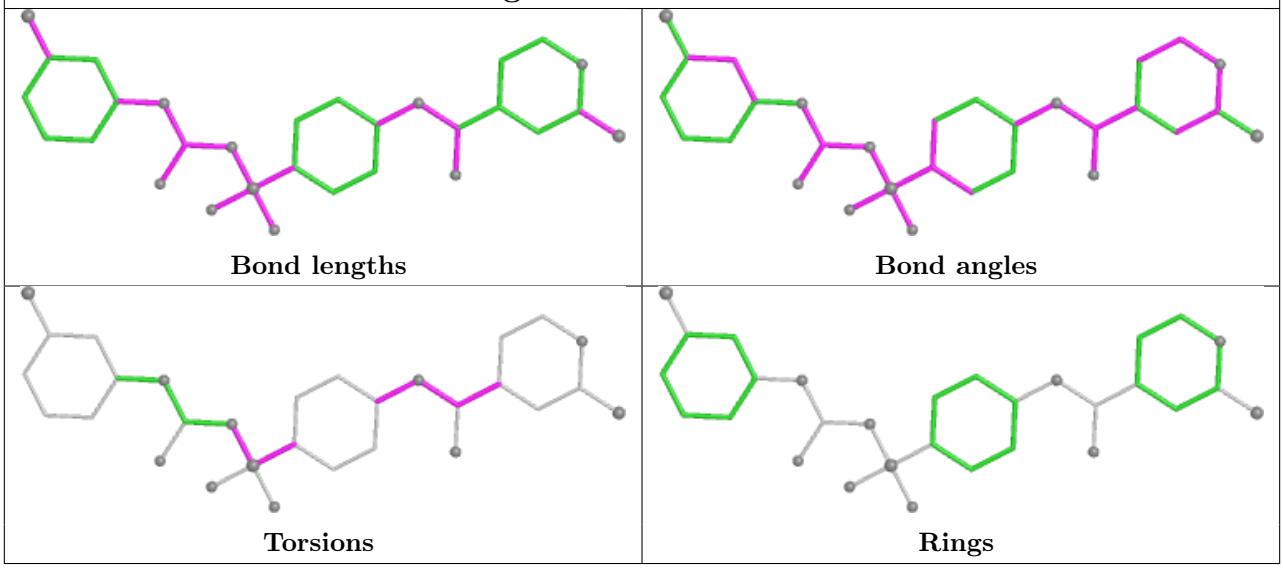


Rings

Ligand A1EZR D 407



Ligand A1EZR A 409



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, 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	319/346 (92%)	-0.14	11 (3%)	48	50	20, 28, 46, 72	0
1	B	325/346 (93%)	0.08	11 (3%)	48	50	21, 30, 52, 98	0
1	C	319/346 (92%)	0.07	13 (4%)	41	43	21, 31, 50, 68	0
1	D	325/346 (93%)	0.32	25 (7%)	19	20	23, 35, 64, 81	0
All	All	1288/1384 (93%)	0.09	60 (4%)	36	38	20, 31, 54, 98	0

The worst 5 of 60 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	63	THR	6.3
1	B	66	THR	6.1
1	D	269	LYS	5.8
1	B	64	ASN	5.7
1	D	63	THR	5.5

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

There are no oligosaccharides in this entry.

6.4 Ligands

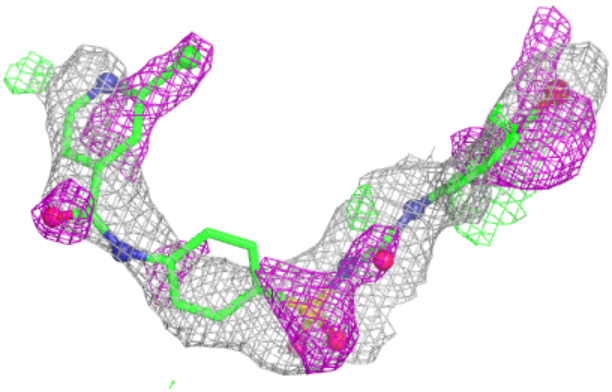
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

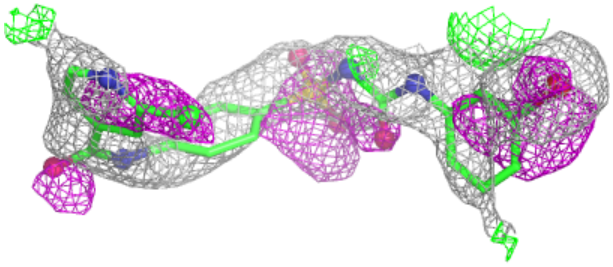
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	A1EZR	A	409	30/30	0.64	0.19	56,66,86,99	0
3	EDO	C	402	4/4	0.73	0.20	52,54,55,59	0
3	EDO	C	404	4/4	0.73	0.19	45,50,53,61	0
2	GOL	A	410	6/6	0.73	0.20	47,52,56,62	0
2	GOL	D	403	6/6	0.74	0.15	48,51,57,59	0
2	GOL	D	405	6/6	0.75	0.21	64,69,74,75	0
2	GOL	A	404	6/6	0.75	0.15	53,54,56,69	0
2	GOL	A	403	6/6	0.76	0.16	46,49,56,67	0
3	EDO	D	406	4/4	0.76	0.18	47,48,51,51	0
3	EDO	C	403	4/4	0.76	0.15	40,44,45,53	0
3	EDO	C	401	4/4	0.78	0.18	49,53,56,56	0
2	GOL	A	402	6/6	0.78	0.15	37,44,51,53	0
5	A1EZR	D	407	30/30	0.79	0.17	53,62,73,93	0
2	GOL	D	402	6/6	0.82	0.15	45,47,49,53	0
2	GOL	D	401	6/6	0.82	0.15	39,46,51,51	0
5	A1EZR	C	405	30/30	0.83	0.13	33,47,58,103	0
3	EDO	A	405	4/4	0.83	0.15	52,59,61,65	0
3	EDO	A	406	4/4	0.84	0.14	46,50,54,60	0
2	GOL	A	401	6/6	0.85	0.14	33,46,51,53	0
6	MG	A	411	1/1	0.85	0.17	51,51,51,51	0
5	A1EZR	B	401	30/30	0.86	0.12	34,51,61,89	0
2	GOL	D	404	6/6	0.86	0.15	53,54,58,65	0
3	EDO	A	407	4/4	0.87	0.12	51,55,59,60	0
4	AMP	A	408	23/23	0.93	0.08	29,33,38,39	0
4	AMP	D	408	23/23	0.94	0.08	34,38,40,42	0
4	AMP	B	402	23/23	0.97	0.06	27,35,37,38	0
4	AMP	C	406	23/23	0.97	0.06	27,31,33,35	0

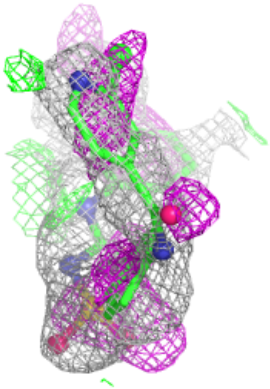
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around A1EZR A 409:

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

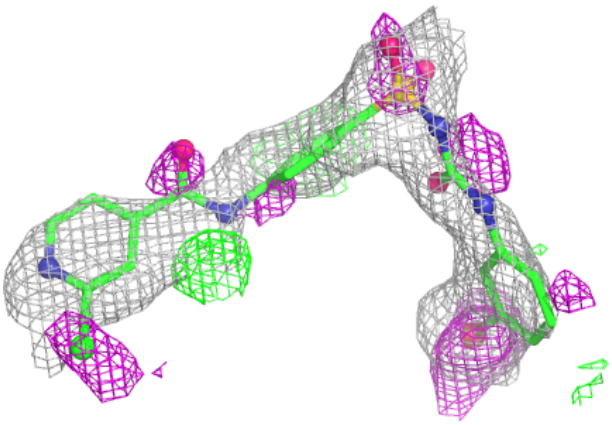


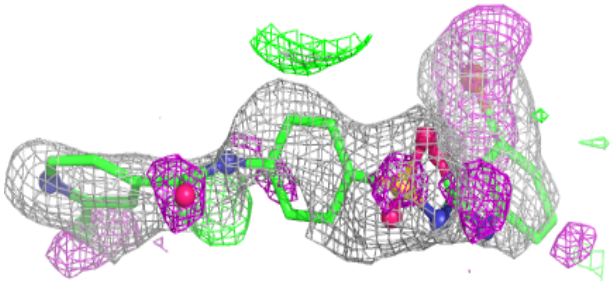


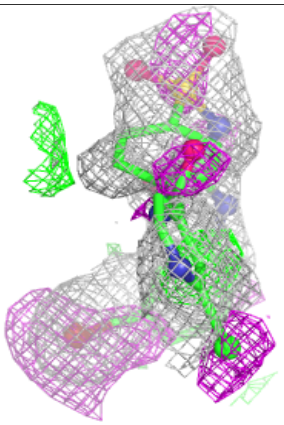


Electron density around A1EZR D 407:

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

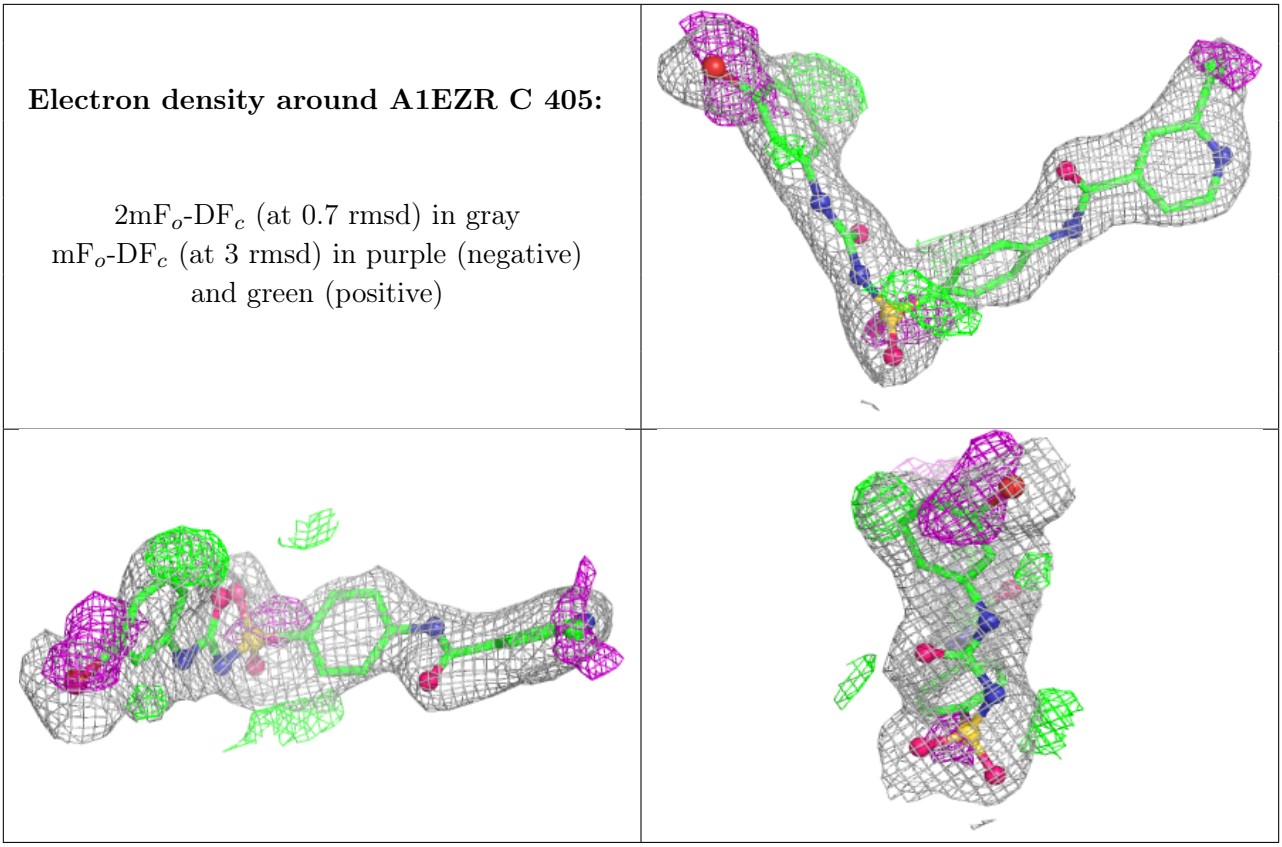






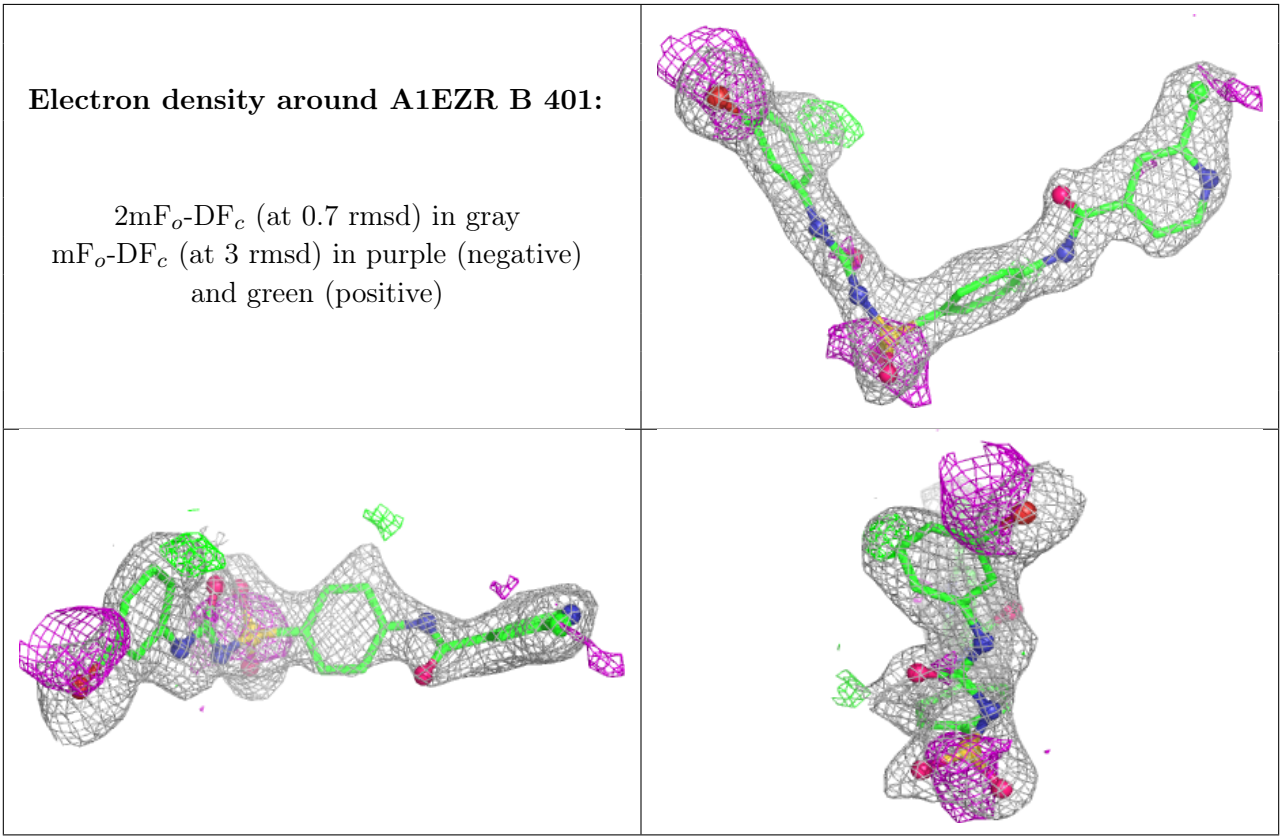
Electron density around A1EZR C 405:

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



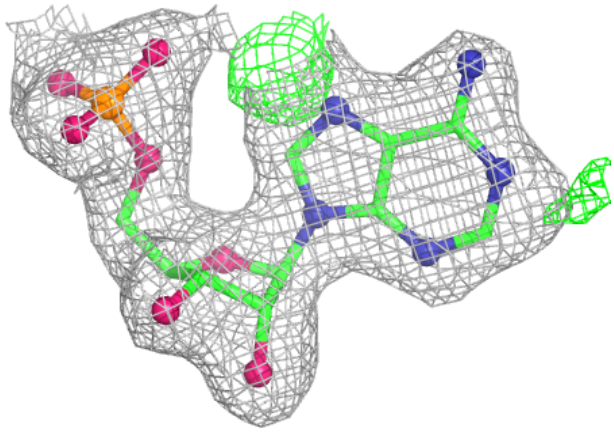
Electron density around A1EZR B 401:

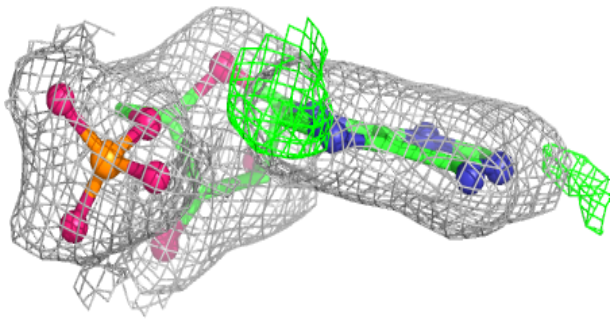
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

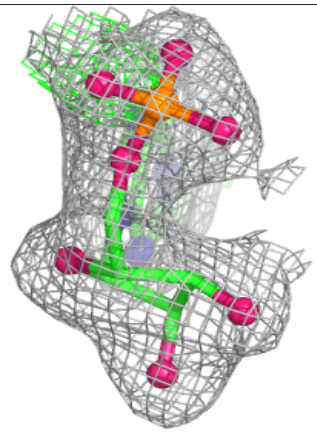


Electron density around AMP A 408:

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

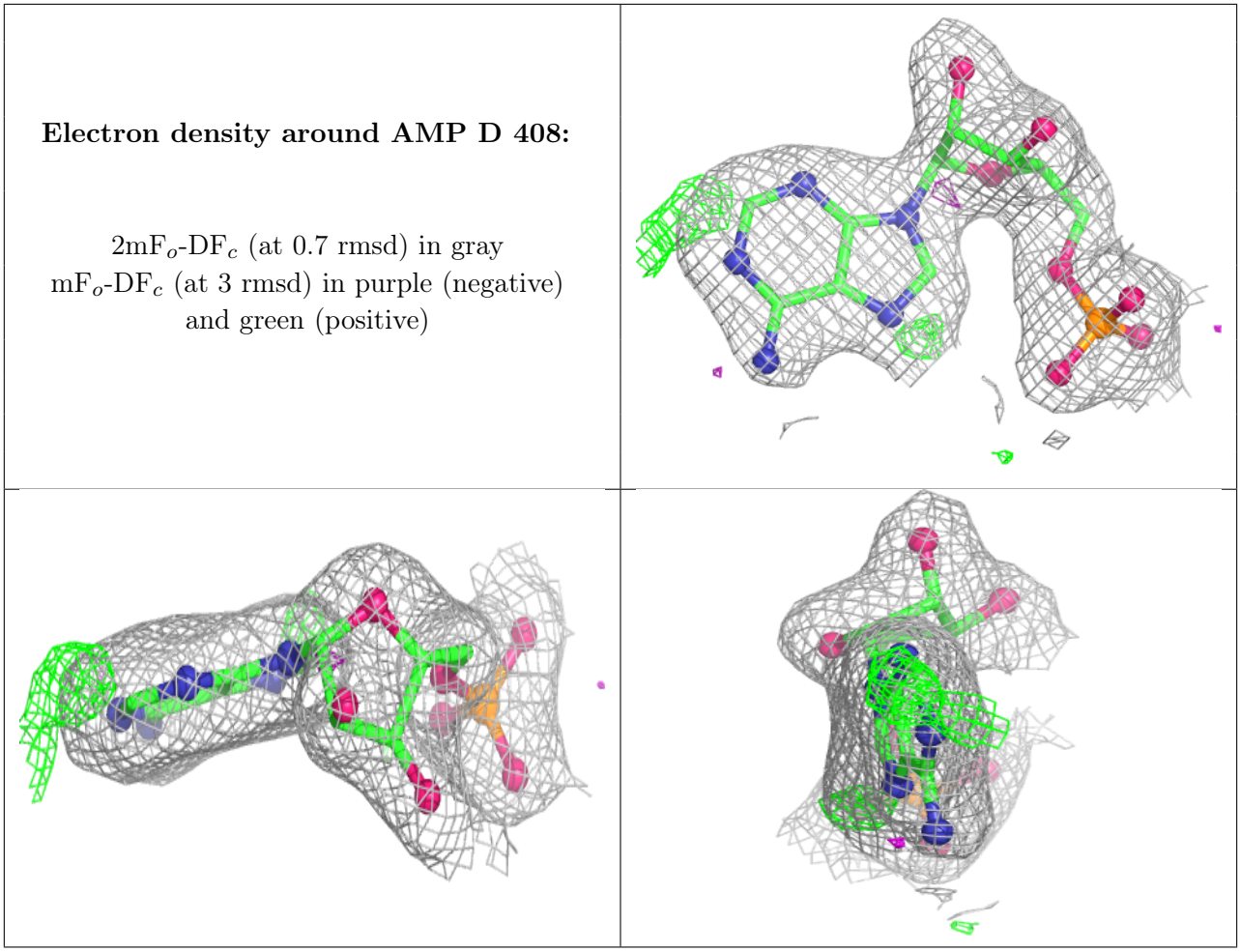






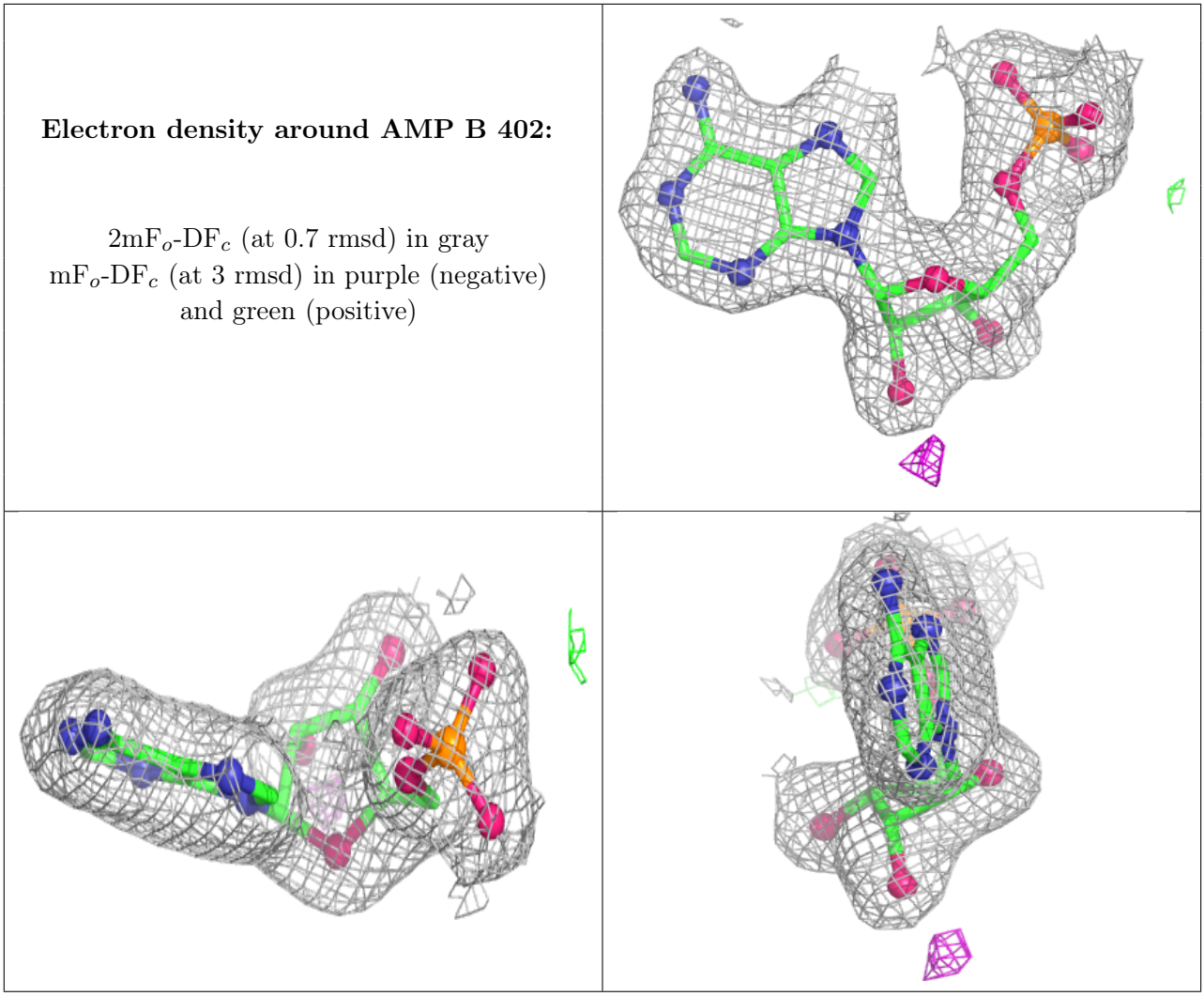
Electron density around AMP D 408:

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



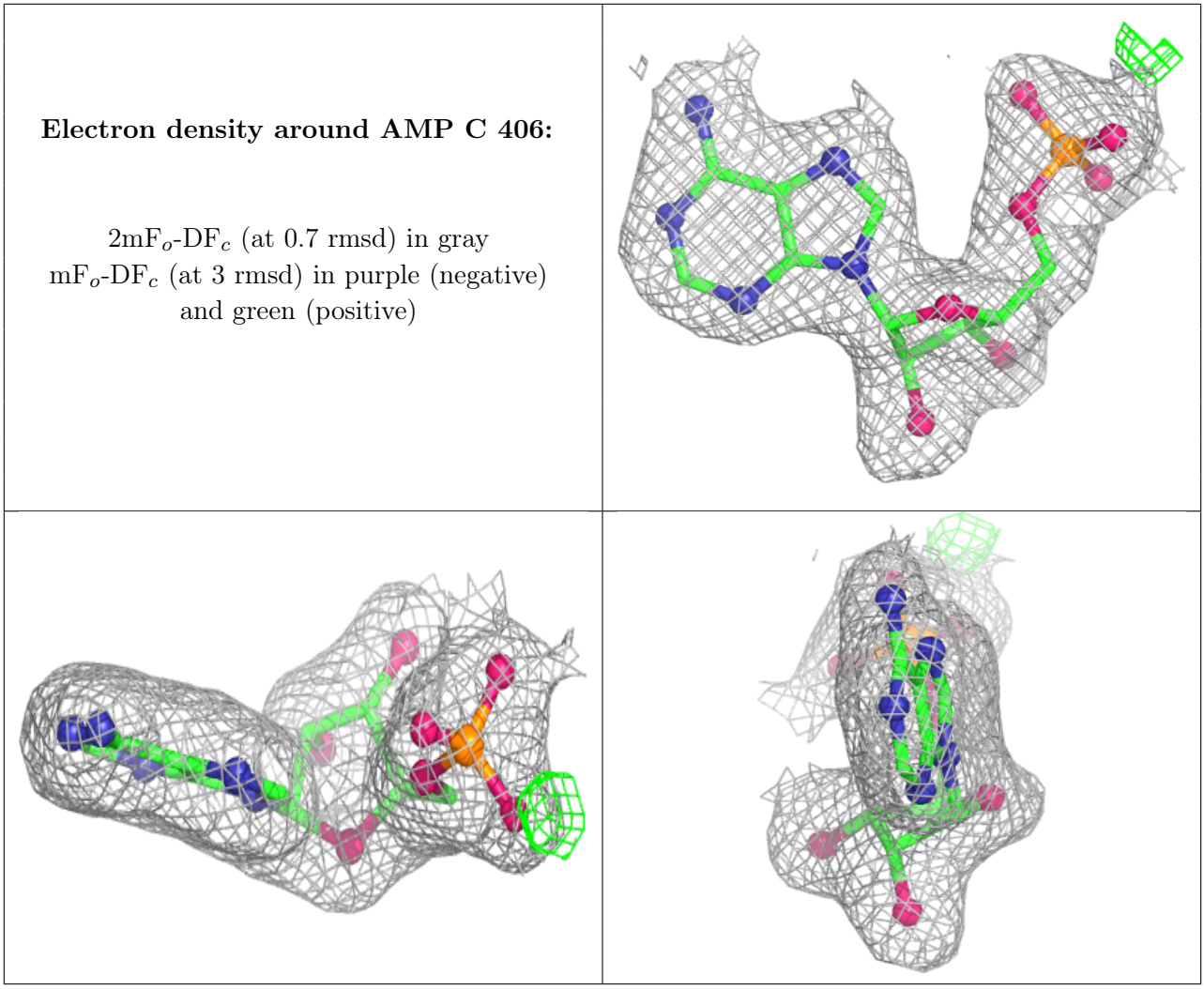
Electron density around AMP B 402:

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



Electron density around AMP C 406:

$2mF_o$-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.