



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 04:07 AM UTC

PDB ID : 3OPY / pdb_00003opy
Title : Crystal structure of Pichia pastoris phosphofructokinase in the T-state
Authors : Strater, N.; Marek, S.; Kuettner, E.B.; Kloos, M.; Keim, A.; Bruser, A.; Kirchberger, J.; Schoneberg, T.
Deposited on : 2010-09-02
Resolution : 3.05 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)
Xtrriage (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

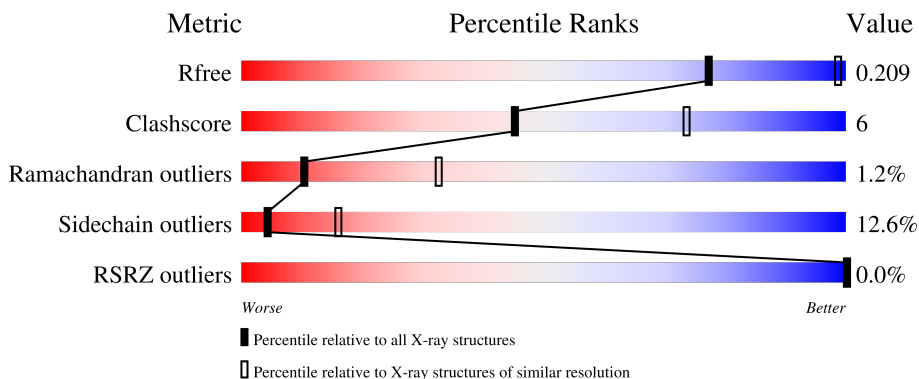
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	989	71% 21% • 5%
1	C	989	71% 20% • 5%
1	E	989	71% 20% • 5%
1	G	989	71% 20% • 5%
2	B	941	72% 18% • 6%

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Mol	Chain	Length	Quality of chain			
2	D	941				
2	F	941				
2	H	941				
3	I	351				
3	J	351				
3	K	351				
3	L	351				

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
4	SO4	C	991	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 65025 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 6-phosphofructo-1-kinase alpha-subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	936	7008	4410	1200	1373	25	0	0	0
1	C	936	7008	4410	1200	1373	25	0	0	0
1	E	936	7008	4410	1200	1373	25	0	0	0
1	G	936	7008	4410	1200	1373	25	0	0	0

- Molecule 2 is a protein called 6-phosphofructo-1-kinase beta-subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	884	6546	4093	1165	1254	34	0	0	0
2	D	884	6546	4093	1165	1254	34	0	0	0
2	F	884	6546	4093	1165	1254	34	0	0	0
2	H	884	6546	4093	1165	1254	34	0	0	0

- Molecule 3 is a protein called 6-phosphofructo-1-kinase gamma-subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	I	323	2590	1650	439	492	9	0	0	0
3	J	323	2590	1650	439	492	9	0	0	0
3	K	323	2590	1650	439	492	9	0	0	0
3	L	323	2590	1650	439	492	9	0	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		

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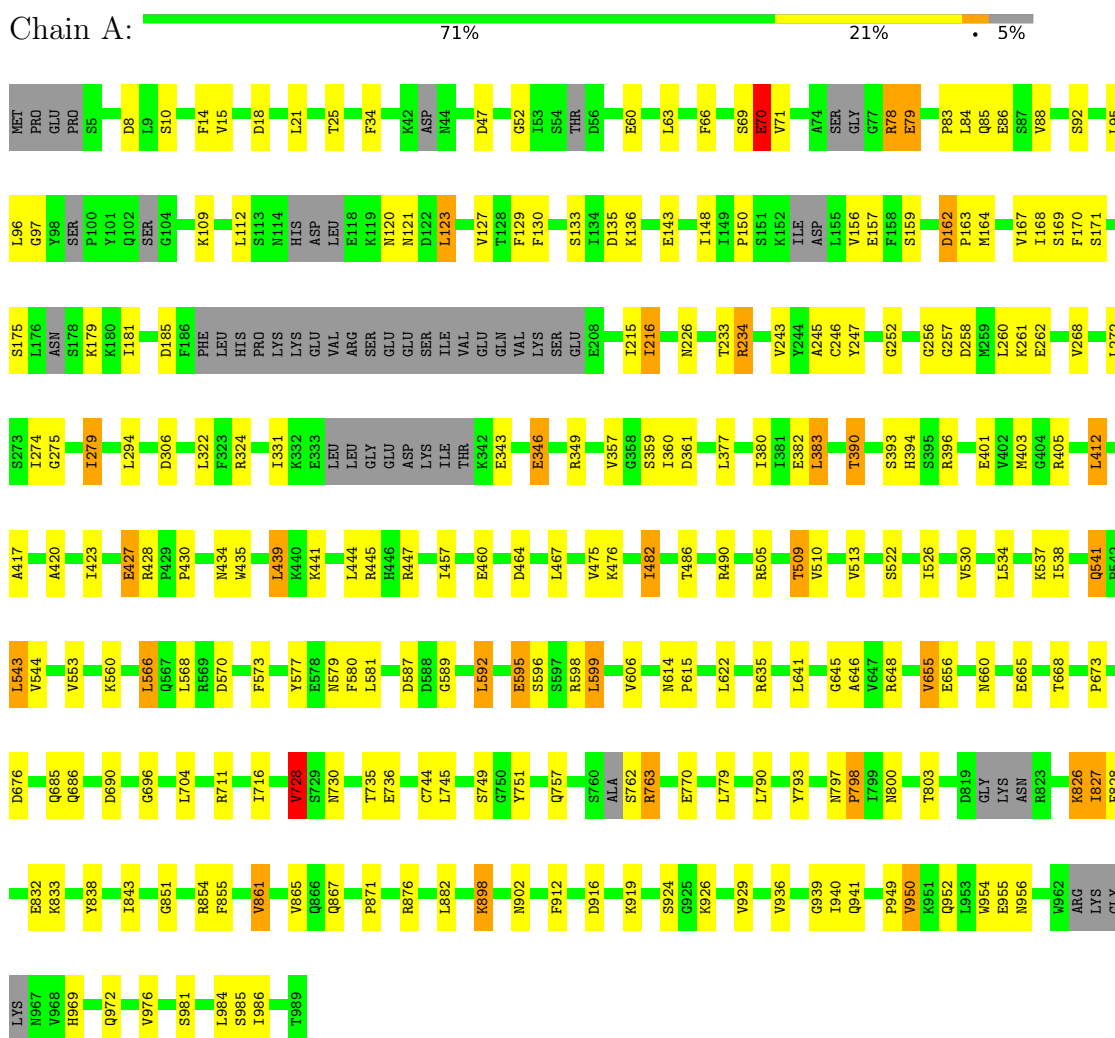
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
5	F	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
5	H	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

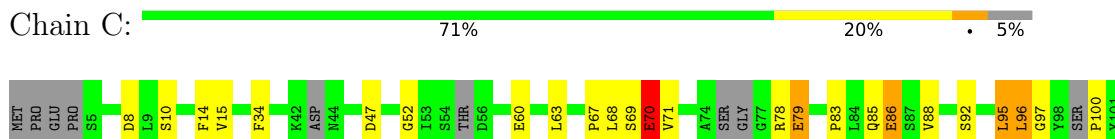
3 Residue-property plots [i](#)

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: 6-phosphofructo-1-kinase alpha-subunit



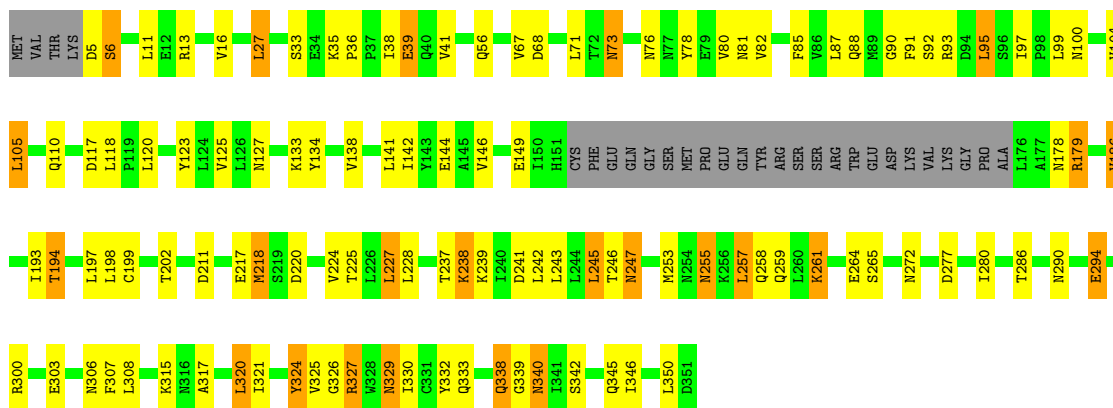
- Molecule 1: 6-phosphofructo-1-kinase alpha-subunit





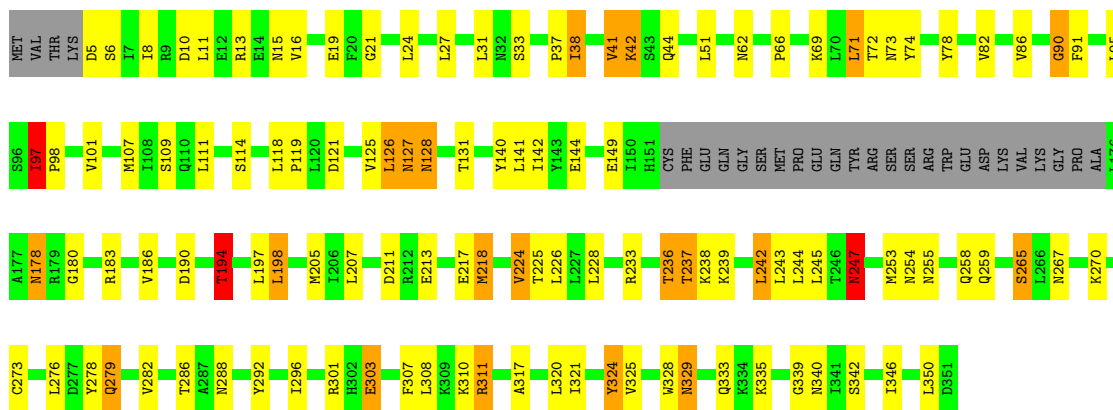
- Molecule 3: 6-phosphofructo-1-kinase gamma-subunit

Chain I: 60% 26% 7% 8%



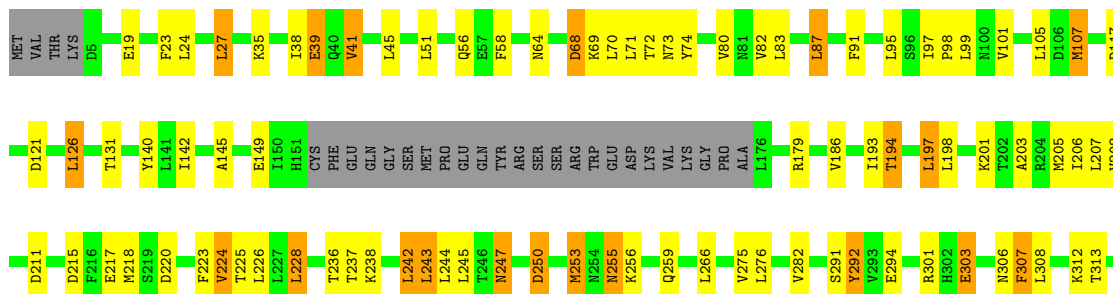
- Molecule 3: 6-phosphofructo-1-kinase gamma-subunit

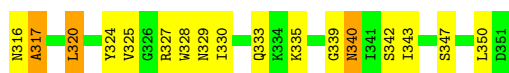
Chain J: 58% 27% 6% 8%



- Molecule 3: 6-phosphofructo-1-kinase gamma-subunit

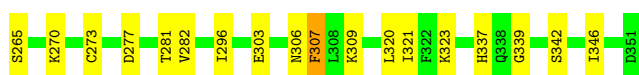
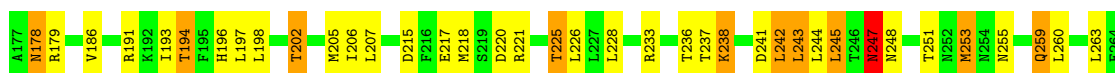
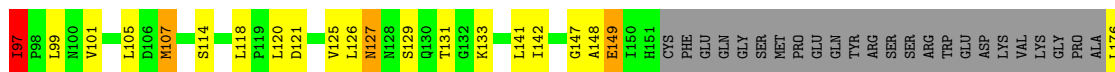
Chain K: 62% 24% 7% 8%





- Molecule 3: 6-phosphofructo-1-kinase gamma-subunit

Chain L: 60% 26% 5% 8%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	161.66Å 188.30Å 231.56Å 90.00° 92.85° 90.00°	Depositor
Resolution (Å)	29.80 – 3.05 29.80 – 3.05	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.80-3.05) 84.0 (29.80-3.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 3.06Å)	Xtrriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.200 , 0.232 0.216 , 0.209	Depositor DCC
R_{free} test set	5145 reflections (1.96%)	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtrriage
Anisotropy	0.250	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtrriage
Estimated twinning fraction	0.079 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	65025	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.64% 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: 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	0.86	0/7125	1.45	49/9664 (0.5%)
1	C	0.86	0/7125	1.44	54/9664 (0.6%)
1	E	0.81	0/7125	1.44	47/9664 (0.5%)
1	G	0.83	0/7125	1.44	46/9664 (0.5%)
2	B	0.88	0/6651	1.45	43/9002 (0.5%)
2	D	0.89	1/6651 (0.0%)	1.46	44/9002 (0.5%)
2	F	0.83	0/6651	1.45	33/9002 (0.4%)
2	H	0.84	0/6651	1.44	33/9002 (0.4%)
3	I	1.02	5/2640 (0.2%)	1.62	36/3578 (1.0%)
3	J	0.97	1/2640 (0.0%)	1.57	37/3578 (1.0%)
3	K	0.90	0/2640	1.53	25/3578 (0.7%)
3	L	0.92	3/2640 (0.1%)	1.51	26/3578 (0.7%)
All	All	0.87	10/65664 (0.0%)	1.46	473/88976 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1
2	B	0	1
2	D	0	1
2	F	0	1
2	H	0	1
All	All	0	5

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	218	MET	SD-CE	-6.57	1.63	1.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	218	MET	SD-CE	-5.72	1.65	1.79
2	D	593	MET	SD-CE	-5.55	1.65	1.79
3	I	261	LYS	CA-C	5.49	1.60	1.52
3	L	97	ILE	CA-CB	5.27	1.56	1.54

The worst 5 of 473 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	67	VAL	CA-C-N	10.74	135.04	120.54
3	I	67	VAL	C-N-CA	10.74	135.04	120.54
3	L	194	THR	CB-CA-C	9.55	125.00	111.82
3	J	127	ASN	N-CA-C	9.45	121.18	111.07
1	A	394	HIS	N-CA-C	9.07	121.17	111.28

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	915	VAL	Mainchain
2	D	915	VAL	Mainchain
1	E	123	LEU	Mainchain
2	F	915	VAL	Mainchain
2	H	915	VAL	Mainchain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7008	0	6692	86	0
1	C	7008	0	6692	85	0
1	E	7008	0	6692	86	0
1	G	7008	0	6692	81	0
2	B	6546	0	6253	73	0
2	D	6546	0	6253	75	0
2	F	6546	0	6253	72	0
2	H	6546	0	6253	73	0
3	I	2590	0	2509	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	J	2590	0	2509	54	0
3	K	2590	0	2509	40	0
3	L	2590	0	2509	46	0
4	A	65	0	0	1	0
4	B	10	0	0	0	0
4	C	50	0	0	3	0
4	D	15	0	0	0	0
4	E	40	0	0	2	0
4	F	15	0	0	0	0
4	G	40	0	0	0	0
4	H	20	0	0	0	0
4	I	25	0	0	0	0
4	J	15	0	0	0	0
4	K	15	0	0	0	0
4	L	15	0	0	0	0
5	B	31	0	12	0	0
5	D	31	0	12	0	0
5	F	31	0	12	1	0
5	H	31	0	12	1	0
All	All	65025	0	61864	750	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 6.

The worst 5 of 750 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)
2:B:650:ASP:CB	2:D:148:ASN:HB2	2.01	0.90
2:F:148:ASN:HB2	2:H:650:ASP:CB	2.06	0.85
1:C:622:LEU:HD21	1:C:655:VAL:HG11	1.59	0.84
1:E:216:ILE:HG22	1:E:246:CYS:HB3	1.60	0.84
3:I:255:ASN:O	3:I:259:GLN:HG3	1.78	0.84

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	908/989 (92%)	840 (92%)	58 (6%)	10 (1%)	11	34
1	C	908/989 (92%)	841 (93%)	57 (6%)	10 (1%)	11	34
1	E	908/989 (92%)	841 (93%)	57 (6%)	10 (1%)	11	34
1	G	908/989 (92%)	843 (93%)	55 (6%)	10 (1%)	11	34
2	B	859/941 (91%)	800 (93%)	46 (5%)	13 (2%)	8	28
2	D	859/941 (91%)	803 (94%)	45 (5%)	11 (1%)	9	31
2	F	859/941 (91%)	803 (94%)	45 (5%)	11 (1%)	9	31
2	H	859/941 (91%)	800 (93%)	46 (5%)	13 (2%)	8	28
3	I	319/351 (91%)	296 (93%)	21 (7%)	2 (1%)	21	48
3	J	319/351 (91%)	285 (89%)	29 (9%)	5 (2%)	7	26
3	K	319/351 (91%)	286 (90%)	31 (10%)	2 (1%)	21	48
3	L	319/351 (91%)	283 (89%)	33 (10%)	3 (1%)	14	39
All	All	8344/9124 (92%)	7721 (92%)	523 (6%)	100 (1%)	10	32

5 of 100 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	646	ALA
2	B	127	SER
2	B	150	PHE
2	B	303	ILE
1	C	646	ALA

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	725/835 (87%)	637 (88%)	88 (12%)	5	17
1	C	725/835 (87%)	638 (88%)	87 (12%)	5	17
1	E	725/835 (87%)	643 (89%)	82 (11%)	5	20
1	G	725/835 (87%)	642 (89%)	83 (11%)	5	20
2	B	670/799 (84%)	588 (88%)	82 (12%)	5	17
2	D	670/799 (84%)	588 (88%)	82 (12%)	5	17
2	F	670/799 (84%)	587 (88%)	83 (12%)	4	16
2	H	670/799 (84%)	587 (88%)	83 (12%)	4	16
3	I	282/322 (88%)	238 (84%)	44 (16%)	2	10
3	J	282/322 (88%)	239 (85%)	43 (15%)	3	10
3	K	282/322 (88%)	237 (84%)	45 (16%)	2	10
3	L	282/322 (88%)	238 (84%)	44 (16%)	2	10
All	All	6708/7824 (86%)	5862 (87%)	846 (13%)	4	16

5 of 846 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	550	MET
1	G	833	LYS
3	K	340	ASN
2	F	686	ILE
2	F	547	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 201 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	344	GLN
2	H	720	ASN
3	L	338	GLN
1	G	644	HIS
2	H	81	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](#)

69 ligands are modelled in this entry.

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	SO4	I	352	-	4,4,4	0.65	0	6,6,6	0.51	0
4	SO4	G	996	-	4,4,4	0.28	0	6,6,6	0.09	0
5	ATP	B	942	-	32,33,33	1.31	5 (15%)	48,52,52	1.15	4 (8%)
4	SO4	A	990	-	4,4,4	0.71	0	6,6,6	0.33	0
4	SO4	E	995	-	4,4,4	0.35	0	6,6,6	0.21	0
4	SO4	C	998	-	4,4,4	0.30	0	6,6,6	0.21	0
4	SO4	F	944	-	4,4,4	0.45	0	6,6,6	0.27	0
4	SO4	F	945	-	4,4,4	0.30	0	6,6,6	0.24	0
4	SO4	G	991	-	4,4,4	0.53	0	6,6,6	0.49	0
4	SO4	K	353	-	4,4,4	0.34	0	6,6,6	0.15	0
4	SO4	A	999	-	4,4,4	0.81	0	6,6,6	0.19	0
4	SO4	A	1002	-	4,4,4	0.35	0	6,6,6	0.16	0
4	SO4	C	992	-	4,4,4	1.21	0	6,6,6	0.23	0
4	SO4	A	994	-	4,4,4	0.22	0	6,6,6	0.38	0
4	SO4	C	995	-	4,4,4	0.37	0	6,6,6	0.18	0
4	SO4	A	995	-	4,4,4	0.60	0	6,6,6	0.26	0
4	SO4	C	991	-	4,4,4	0.62	0	6,6,6	0.25	0
4	SO4	A	992	-	4,4,4	0.39	0	6,6,6	0.18	0
4	SO4	G	995	-	4,4,4	0.19	0	6,6,6	0.26	0
4	SO4	J	352	-	4,4,4	0.32	0	6,6,6	0.22	0
4	SO4	A	1001	-	4,4,4	0.59	0	6,6,6	0.23	0
4	SO4	D	944	-	4,4,4	0.39	0	6,6,6	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	F	943	-	4,4,4	0.49	0	6,6,6	0.23	0
4	SO4	E	990	-	4,4,4	0.37	0	6,6,6	0.14	0
5	ATP	F	942	-	32,33,33	1.16	3 (9%)	48,52,52	1.19	2 (4%)
4	SO4	D	945	-	4,4,4	0.22	0	6,6,6	0.38	0
4	SO4	C	993	-	4,4,4	0.24	0	6,6,6	0.41	0
4	SO4	A	996	-	4,4,4	0.57	0	6,6,6	0.29	0
4	SO4	E	996	-	4,4,4	0.25	0	6,6,6	0.12	0
4	SO4	E	991	-	4,4,4	0.83	0	6,6,6	0.27	0
4	SO4	C	997	-	4,4,4	0.36	0	6,6,6	0.21	0
4	SO4	E	993	-	4,4,4	0.28	0	6,6,6	0.41	0
4	SO4	G	992	-	4,4,4	0.57	0	6,6,6	0.13	0
4	SO4	A	991	-	4,4,4	0.96	0	6,6,6	0.28	0
4	SO4	L	352	-	4,4,4	0.20	0	6,6,6	0.12	0
4	SO4	A	998	-	4,4,4	0.68	0	6,6,6	0.37	0
4	SO4	E	997	-	4,4,4	0.19	0	6,6,6	0.19	0
5	ATP	H	942	-	32,33,33	1.44	7 (21%)	48,52,52	1.35	6 (12%)
4	SO4	B	943	-	4,4,4	0.63	0	6,6,6	0.30	0
4	SO4	K	354	-	4,4,4	0.28	0	6,6,6	0.19	0
4	SO4	G	993	-	4,4,4	0.38	0	6,6,6	0.26	0
4	SO4	D	943	-	4,4,4	0.92	0	6,6,6	0.42	0
4	SO4	A	1000	-	4,4,4	0.32	0	6,6,6	0.21	0
4	SO4	I	355	-	4,4,4	0.42	0	6,6,6	0.16	0
4	SO4	J	353	-	4,4,4	0.39	0	6,6,6	0.16	0
4	SO4	C	994	-	4,4,4	0.50	0	6,6,6	0.30	0
4	SO4	A	997	-	4,4,4	0.32	0	6,6,6	0.24	0
5	ATP	D	942	-	32,33,33	1.41	5 (15%)	48,52,52	1.40	8 (16%)
4	SO4	H	943	-	4,4,4	0.26	0	6,6,6	0.34	0
4	SO4	H	944	-	4,4,4	0.20	0	6,6,6	0.33	0
4	SO4	B	944	-	4,4,4	0.62	0	6,6,6	0.57	0
4	SO4	E	992	-	4,4,4	0.28	0	6,6,6	0.26	0
4	SO4	E	994	-	4,4,4	0.41	0	6,6,6	0.12	0
4	SO4	G	990	-	4,4,4	0.44	0	6,6,6	0.41	0
4	SO4	I	354	-	4,4,4	0.26	0	6,6,6	0.05	0
4	SO4	H	945	-	4,4,4	0.37	0	6,6,6	0.38	0
4	SO4	C	990	-	4,4,4	0.86	0	6,6,6	0.39	0
4	SO4	G	994	-	4,4,4	0.39	0	6,6,6	0.26	0
4	SO4	J	354	-	4,4,4	0.39	0	6,6,6	0.14	0
4	SO4	L	353	-	4,4,4	0.34	0	6,6,6	0.15	0
4	SO4	C	996	-	4,4,4	0.63	0	6,6,6	0.44	0
4	SO4	G	997	-	4,4,4	0.19	0	6,6,6	0.21	0
4	SO4	I	353	-	4,4,4	0.34	0	6,6,6	0.32	0
4	SO4	C	999	-	4,4,4	0.32	0	6,6,6	0.08	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	H	946	-	4,4,4	0.32	0	6,6,6	0.26	0
4	SO4	L	354	-	4,4,4	0.27	0	6,6,6	0.14	0
4	SO4	A	993	-	4,4,4	0.33	0	6,6,6	0.19	0
4	SO4	K	352	-	4,4,4	0.34	0	6,6,6	0.35	0
4	SO4	I	356	-	4,4,4	0.36	0	6,6,6	0.13	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
5	ATP	D	942	-	-	1/22/38/38	0/3/3/3
5	ATP	F	942	-	-	4/22/38/38	0/3/3/3
5	ATP	H	942	-	-	1/22/38/38	0/3/3/3
5	ATP	B	942	-	-	0/22/38/38	0/3/3/3

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	942	ATP	O2'-C2'	4.04	1.53	1.43
5	H	942	ATP	C4-N3	3.10	1.40	1.34
5	F	942	ATP	C2-N1	2.94	1.39	1.33
5	H	942	ATP	C2-N1	2.85	1.39	1.33
5	H	942	ATP	C6-N6	2.73	1.41	1.34

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	942	ATP	O2'-C2'-C1'	3.88	123.48	110.10
5	F	942	ATP	O2G-PG-O3B	3.71	117.06	104.64
5	F	942	ATP	C2'-C1'-N9	3.69	122.47	113.30
5	D	942	ATP	O3A-PA-O1A	3.38	120.89	110.70
5	H	942	ATP	O2G-PG-O3B	3.14	115.18	104.64

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	F	942	ATP	PB-O3A-PA-O5'
5	H	942	ATP	PB-O3A-PA-O5'

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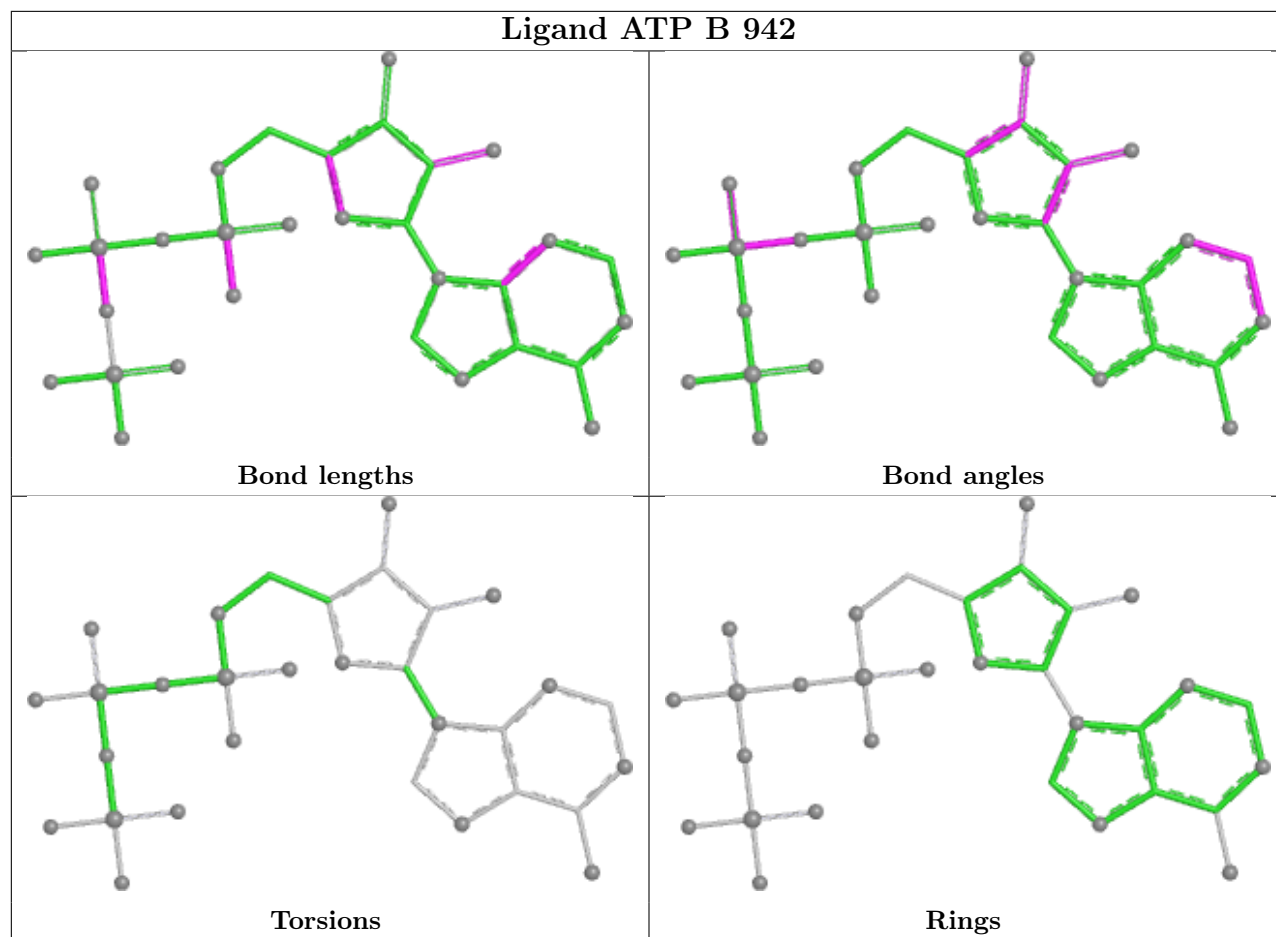
Mol	Chain	Res	Type	Atoms
5	F	942	ATP	O4'-C4'-C5'-O5'
5	F	942	ATP	C5'-O5'-PA-O1A
5	D	942	ATP	PG-O3B-PB-O2B

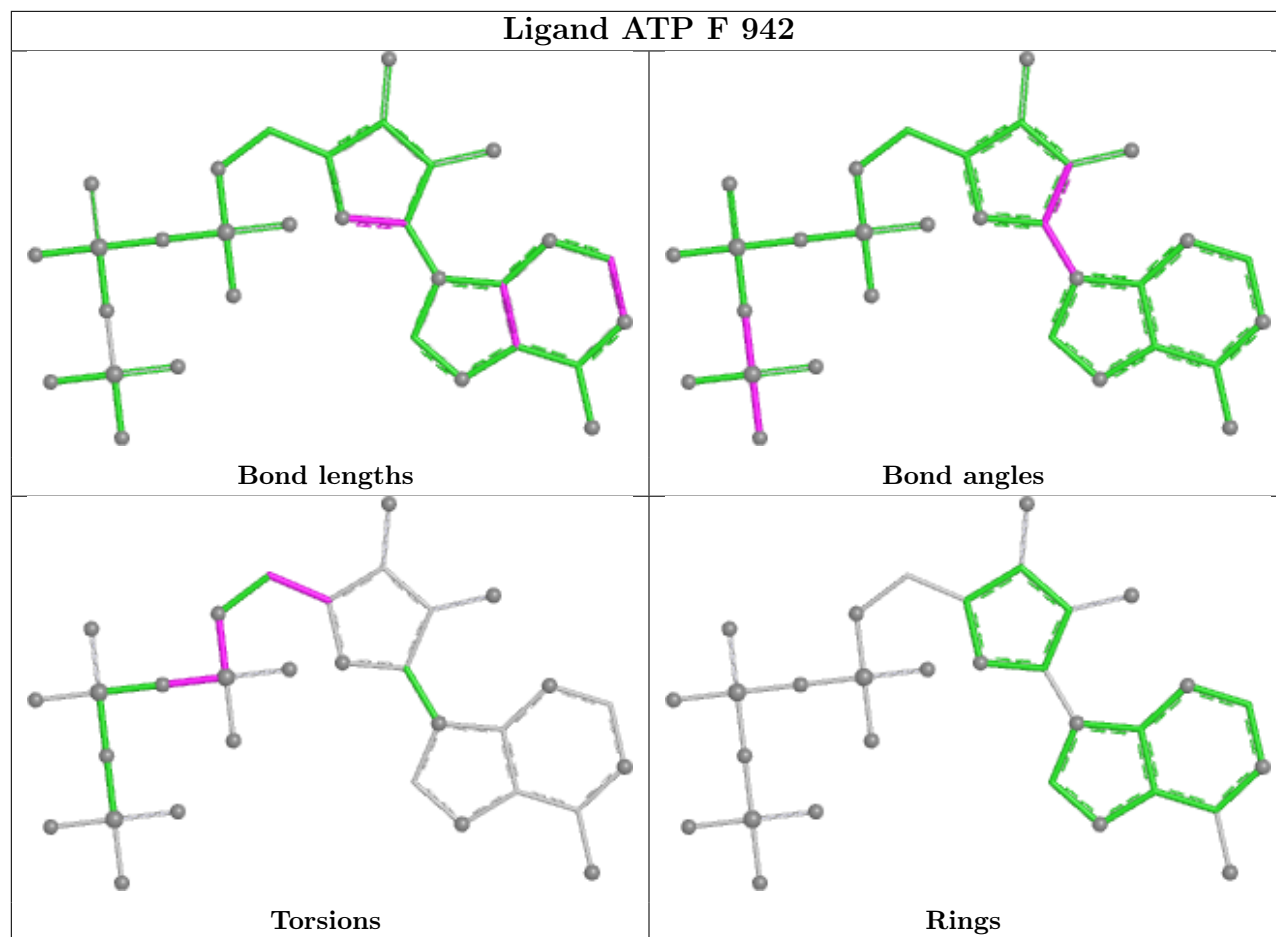
There are no ring outliers.

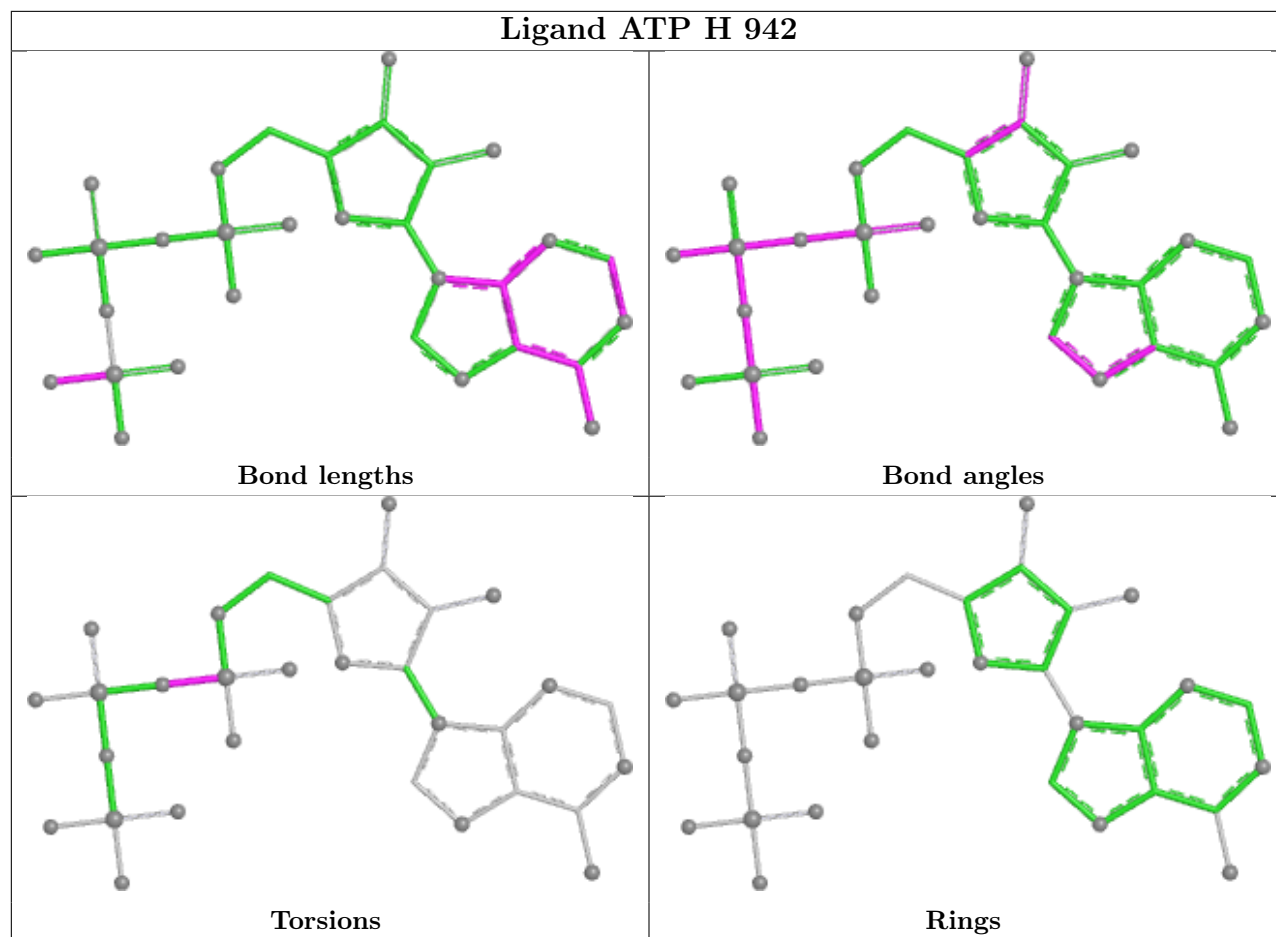
7 monomers are involved in 8 short contacts:

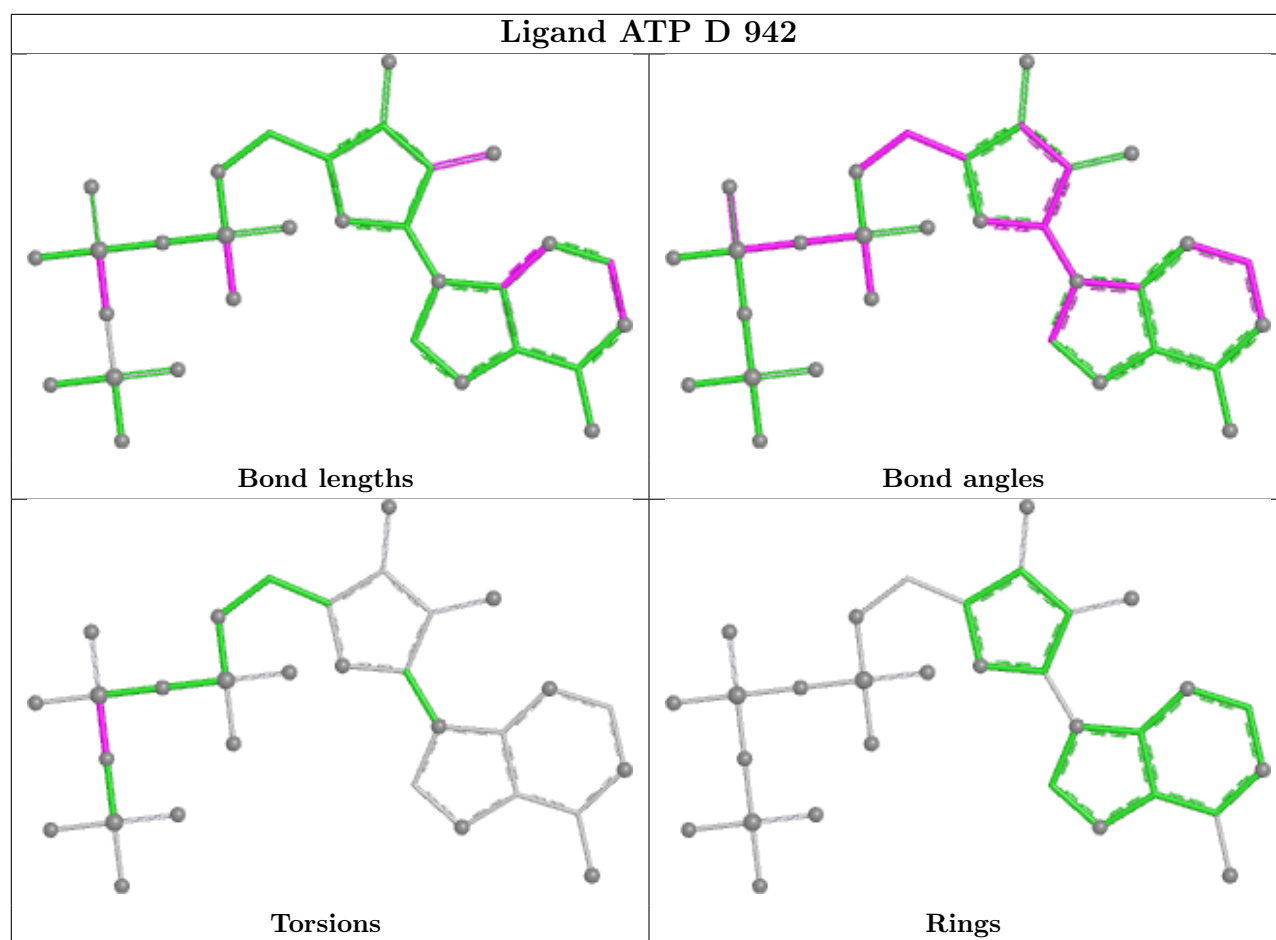
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	991	SO4	2	0
4	E	990	SO4	1	0
5	F	942	ATP	1	0
4	E	991	SO4	1	0
5	H	942	ATP	1	0
4	C	996	SO4	1	0
4	A	993	SO4	1	0

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, 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	936/989 (94%)	-1.26	0 100 100	19, 52, 102, 146	0
1	C	936/989 (94%)	-1.27	0 100 100	21, 53, 107, 163	0
1	E	936/989 (94%)	-1.06	0 100 100	31, 75, 120, 163	0
1	G	936/989 (94%)	-1.10	1 (0%) 92 86	31, 70, 114, 160	0
2	B	884/941 (93%)	-1.22	0 100 100	20, 52, 109, 162	0
2	D	884/941 (93%)	-1.21	1 (0%) 92 86	20, 51, 113, 186	0
2	F	884/941 (93%)	-0.84	1 (0%) 92 86	43, 83, 126, 191	0
2	H	884/941 (93%)	-0.98	0 100 100	38, 72, 133, 184	0
3	I	323/351 (92%)	-1.42	0 100 100	22, 47, 84, 106	0
3	J	323/351 (92%)	-1.43	0 100 100	17, 47, 84, 113	0
3	K	323/351 (92%)	-1.26	0 100 100	38, 66, 96, 121	0
3	L	323/351 (92%)	-1.11	0 100 100	47, 72, 104, 125	0
All	All	8572/9124 (93%)	-1.15	3 (0%) 100 100	17, 64, 116, 191	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	19	ASN	2.3
1	G	173	TYR	2.2
2	F	528	LYS	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 oligosaccharides 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, 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
4	SO4	E	996	5/5	0.95	0.08	123,127,128,129	0
4	SO4	I	355	5/5	0.95	0.08	131,135,136,137	0
4	SO4	J	354	5/5	0.95	0.06	140,144,145,145	0
4	SO4	L	353	5/5	0.95	0.07	165,169,170,170	0
4	SO4	K	354	5/5	0.96	0.07	136,141,141,142	0
4	SO4	I	354	5/5	0.96	0.06	160,164,165,166	0
4	SO4	G	996	5/5	0.97	0.06	133,137,138,139	0
4	SO4	C	999	5/5	0.97	0.07	147,152,152,153	0
4	SO4	L	352	5/5	0.97	0.04	122,127,128,128	0
4	SO4	A	1002	5/5	0.97	0.06	131,136,136,137	0
4	SO4	G	997	5/5	0.98	0.05	135,139,139,140	0
4	SO4	H	944	5/5	0.98	0.05	94,99,99,100	0
4	SO4	H	946	5/5	0.98	0.07	116,120,121,121	0
4	SO4	I	353	5/5	0.98	0.05	91,95,97,97	0
4	SO4	A	992	5/5	0.98	0.04	76,80,80,82	0
4	SO4	A	997	5/5	0.98	0.05	102,106,107,108	0
4	SO4	I	356	5/5	0.98	0.05	125,129,130,130	0
4	SO4	J	352	5/5	0.98	0.07	104,108,109,109	0
4	SO4	E	992	5/5	0.98	0.05	94,99,100,100	0
4	SO4	K	353	5/5	0.98	0.05	95,100,100,100	0
4	SO4	A	1000	5/5	0.98	0.05	105,109,110,110	0
4	SO4	E	997	5/5	0.98	0.05	124,128,129,129	0
4	SO4	A	1001	5/5	0.98	0.05	90,94,95,96	0
4	SO4	L	354	5/5	0.98	0.05	129,133,134,135	0
5	ATP	F	942	31/31	0.98	0.05	83,87,104,105	0
4	SO4	A	993	5/5	0.99	0.03	82,87,88,88	0
4	SO4	A	994	5/5	0.99	0.04	61,65,66,68	0
4	SO4	F	943	5/5	0.99	0.04	68,73,74,74	0
4	SO4	F	944	5/5	0.99	0.03	79,84,84,85	0
4	SO4	F	945	5/5	0.99	0.04	113,118,118,119	0
4	SO4	G	991	5/5	0.99	0.03	55,58,60,61	0
4	SO4	G	992	5/5	0.99	0.03	79,83,84,85	0

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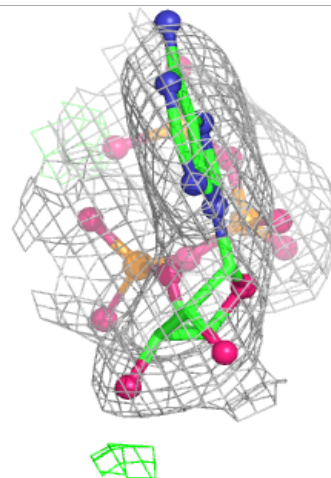
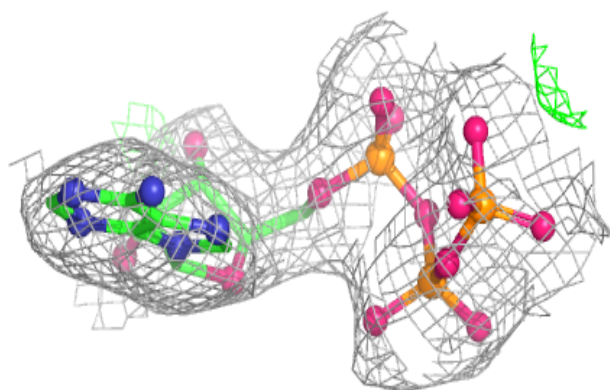
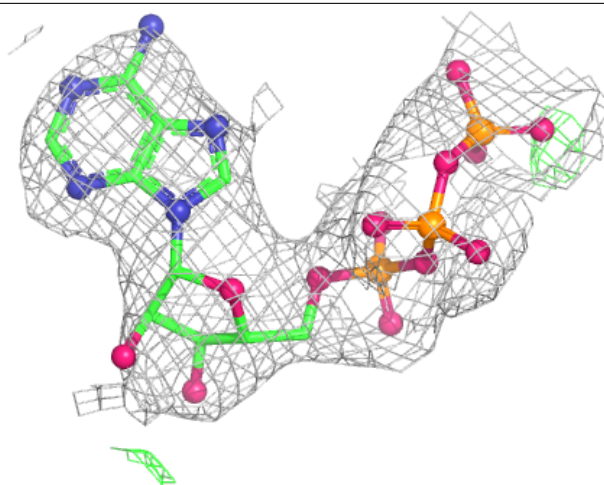
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	G	993	5/5	0.99	0.03	97,101,102,103	0
4	SO4	G	995	5/5	0.99	0.05	96,101,101,102	0
4	SO4	A	995	5/5	0.99	0.06	116,120,121,122	0
4	SO4	B	943	5/5	0.99	0.05	45,49,50,50	0
4	SO4	H	943	5/5	0.99	0.04	69,73,74,75	0
4	SO4	C	990	5/5	0.99	0.03	43,45,48,50	0
4	SO4	C	992	5/5	0.99	0.03	61,65,65,66	0
4	SO4	I	352	5/5	0.99	0.04	66,69,71,71	0
4	SO4	C	994	5/5	0.99	0.04	87,91,92,92	0
4	SO4	C	995	5/5	0.99	0.04	85,89,90,90	0
4	SO4	C	997	5/5	0.99	0.04	113,117,118,118	0
4	SO4	C	998	5/5	0.99	0.06	119,123,124,125	0
4	SO4	A	990	5/5	0.99	0.03	70,74,75,76	0
4	SO4	J	353	5/5	0.99	0.06	83,87,88,88	0
4	SO4	D	944	5/5	0.99	0.03	79,83,83,85	0
4	SO4	K	352	5/5	0.99	0.06	88,92,93,94	0
4	SO4	E	990	5/5	0.99	0.05	110,114,115,115	0
4	SO4	E	991	5/5	0.99	0.04	66,70,72,72	0
4	SO4	A	998	5/5	0.99	0.04	84,87,89,90	0
4	SO4	E	993	5/5	0.99	0.04	93,97,99,99	0
4	SO4	E	994	5/5	0.99	0.04	118,122,123,124	0
5	ATP	D	942	31/31	0.99	0.03	38,42,53,56	0
4	SO4	E	995	5/5	0.99	0.04	95,99,100,100	0
5	ATP	H	942	31/31	0.99	0.04	69,70,102,104	0
4	SO4	A	999	5/5	1.00	0.03	46,50,51,52	0
4	SO4	D	943	5/5	1.00	0.02	33,36,38,40	0
4	SO4	C	993	5/5	1.00	0.02	57,60,61,64	0
4	SO4	D	945	5/5	1.00	0.02	49,53,55,55	0
4	SO4	A	991	5/5	1.00	0.02	57,61,62,64	0
4	SO4	H	945	5/5	1.00	0.03	59,64,65,66	0
4	SO4	B	944	5/5	1.00	0.02	36,40,41,42	0
4	SO4	G	990	5/5	1.00	0.03	63,66,69,70	0
4	SO4	C	996	5/5	1.00	0.02	46,49,51,53	0
5	ATP	B	942	31/31	1.00	0.03	34,36,55,57	0
4	SO4	A	996	5/5	1.00	0.03	45,48,50,53	0
4	SO4	C	991	5/5	1.00	0.02	58,61,63,65	0
4	SO4	G	994	5/5	1.00	0.04	86,89,92,92	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.

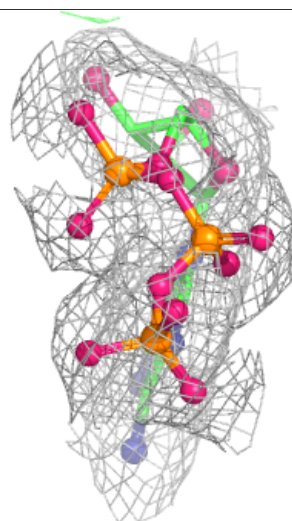
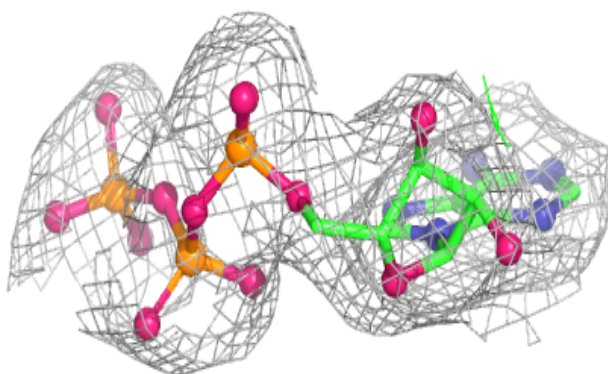
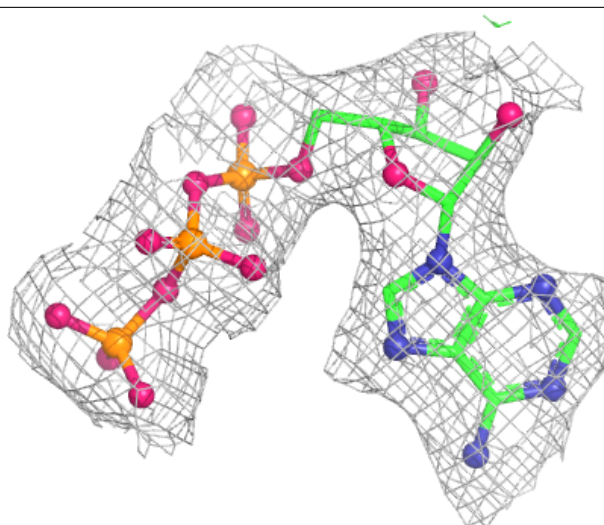
Electron density around ATP F 942:

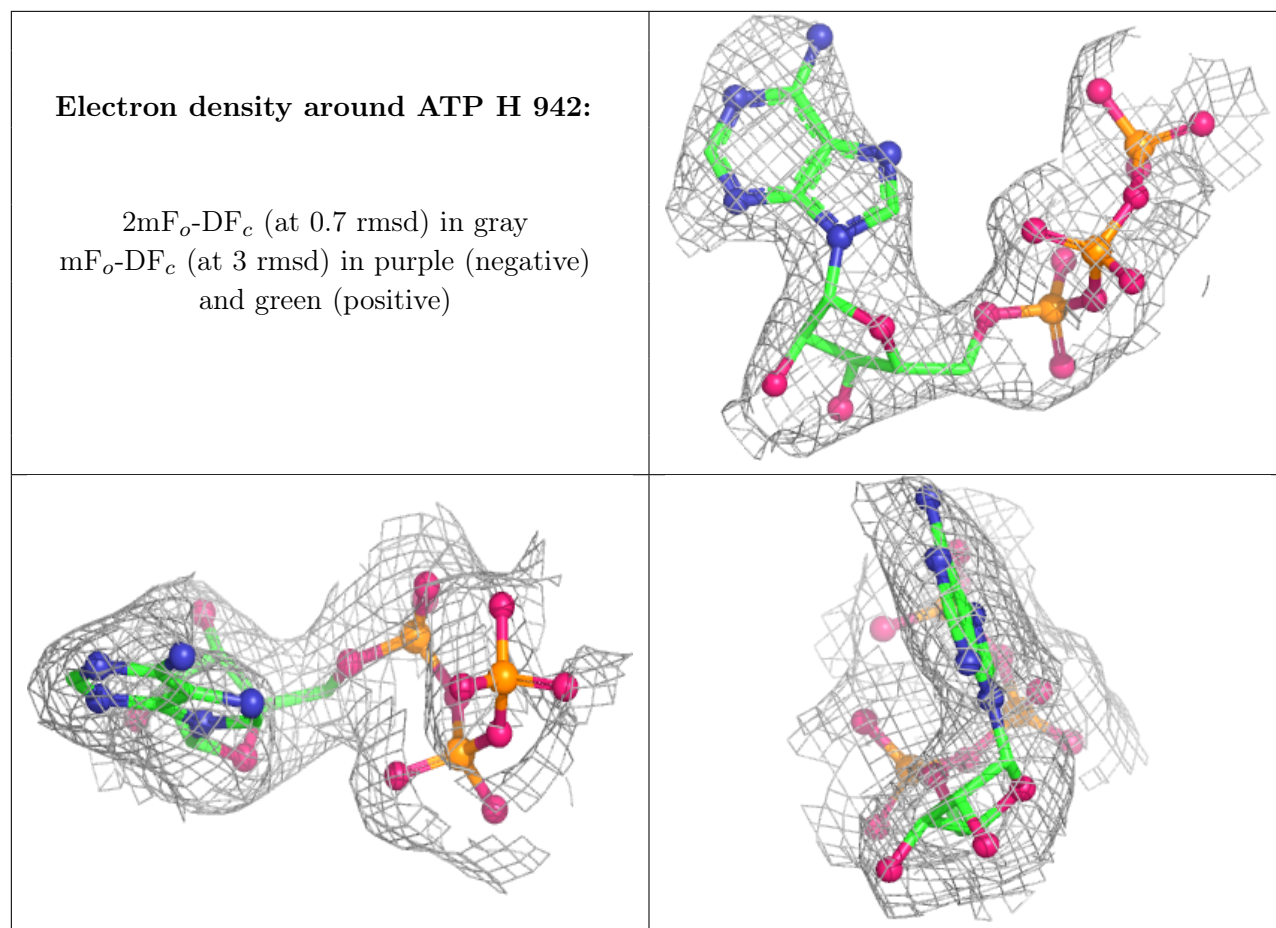
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

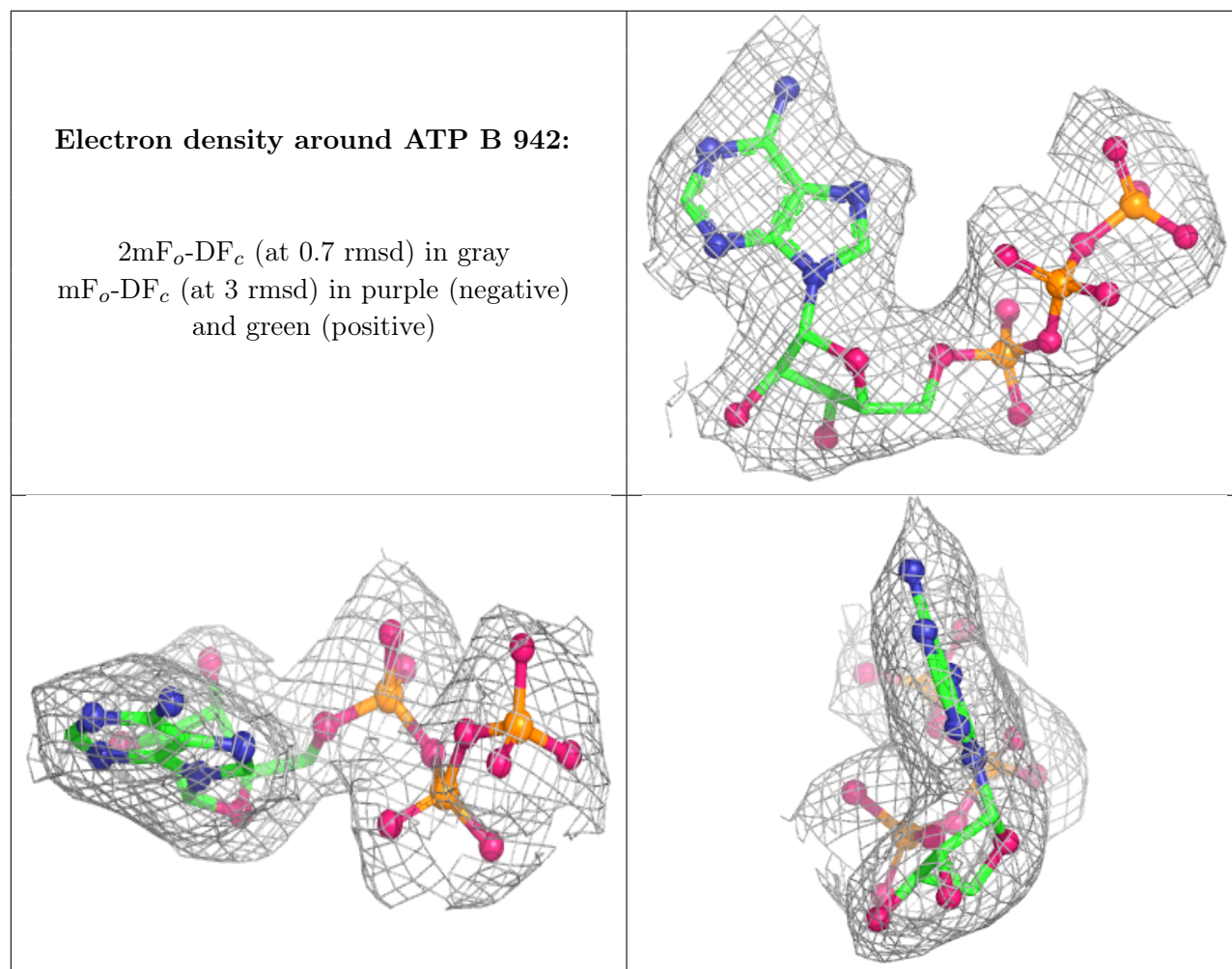


Electron density around ATP D 942:

$2mF_o-DF_c$ (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.