



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:53 PM UTC

PDB ID : 6IUD / pdb_00006iud
Title : Structure of Helicobacter pylori Soj-ADP complex bound to DNA
Authors : Yen, C.Y.; Chu, C.H.; Sun, Y.J.
Deposited on : 2018-11-28
Resolution : 2.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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

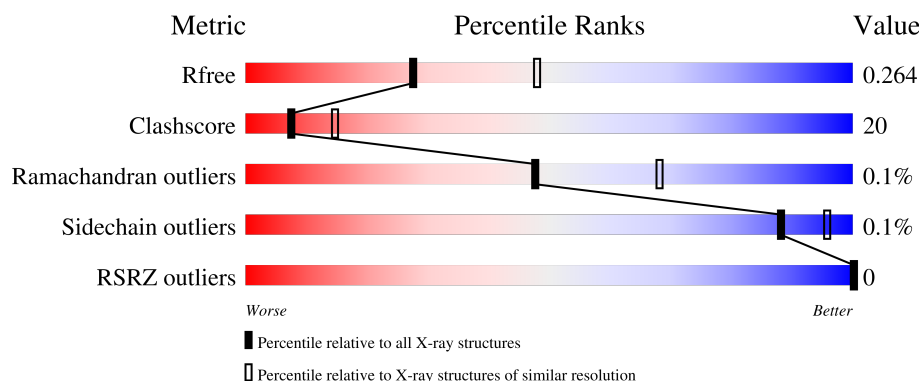
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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.



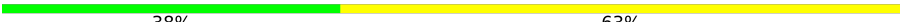
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	276	
1	B	276	
1	C	276	
1	D	276	
2	E	24	

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Mol	Chain	Length	Quality of chain
3	F	24	 38% 63%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9308 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 SpoOJ regulator (Soj).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2053	1328	337	380	8			
1	B	264	Total	C	N	O	S	0	0	0
			2053	1328	337	380	8			
1	C	264	Total	C	N	O	S	0	0	0
			2053	1328	337	380	8			
1	D	264	Total	C	N	O	S	0	0	0
			2053	1328	337	380	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP O25759
A	-10	ARG	-	expression tag	UNP O25759
A	-9	GLY	-	expression tag	UNP O25759
A	-8	SER	-	expression tag	UNP O25759
A	-7	HIS	-	expression tag	UNP O25759
A	-6	HIS	-	expression tag	UNP O25759
A	-5	HIS	-	expression tag	UNP O25759
A	-4	HIS	-	expression tag	UNP O25759
A	-3	HIS	-	expression tag	UNP O25759
A	-2	HIS	-	expression tag	UNP O25759
A	-1	GLY	-	expression tag	UNP O25759
A	0	SER	-	expression tag	UNP O25759
B	-11	MET	-	initiating methionine	UNP O25759
B	-10	ARG	-	expression tag	UNP O25759
B	-9	GLY	-	expression tag	UNP O25759
B	-8	SER	-	expression tag	UNP O25759
B	-7	HIS	-	expression tag	UNP O25759
B	-6	HIS	-	expression tag	UNP O25759
B	-5	HIS	-	expression tag	UNP O25759
B	-4	HIS	-	expression tag	UNP O25759
B	-3	HIS	-	expression tag	UNP O25759

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP O25759
B	-1	GLY	-	expression tag	UNP O25759
B	0	SER	-	expression tag	UNP O25759
C	-11	MET	-	initiating methionine	UNP O25759
C	-10	ARG	-	expression tag	UNP O25759
C	-9	GLY	-	expression tag	UNP O25759
C	-8	SER	-	expression tag	UNP O25759
C	-7	HIS	-	expression tag	UNP O25759
C	-6	HIS	-	expression tag	UNP O25759
C	-5	HIS	-	expression tag	UNP O25759
C	-4	HIS	-	expression tag	UNP O25759
C	-3	HIS	-	expression tag	UNP O25759
C	-2	HIS	-	expression tag	UNP O25759
C	-1	GLY	-	expression tag	UNP O25759
C	0	SER	-	expression tag	UNP O25759
D	-11	MET	-	initiating methionine	UNP O25759
D	-10	ARG	-	expression tag	UNP O25759
D	-9	GLY	-	expression tag	UNP O25759
D	-8	SER	-	expression tag	UNP O25759
D	-7	HIS	-	expression tag	UNP O25759
D	-6	HIS	-	expression tag	UNP O25759
D	-5	HIS	-	expression tag	UNP O25759
D	-4	HIS	-	expression tag	UNP O25759
D	-3	HIS	-	expression tag	UNP O25759
D	-2	HIS	-	expression tag	UNP O25759
D	-1	GLY	-	expression tag	UNP O25759
D	0	SER	-	expression tag	UNP O25759

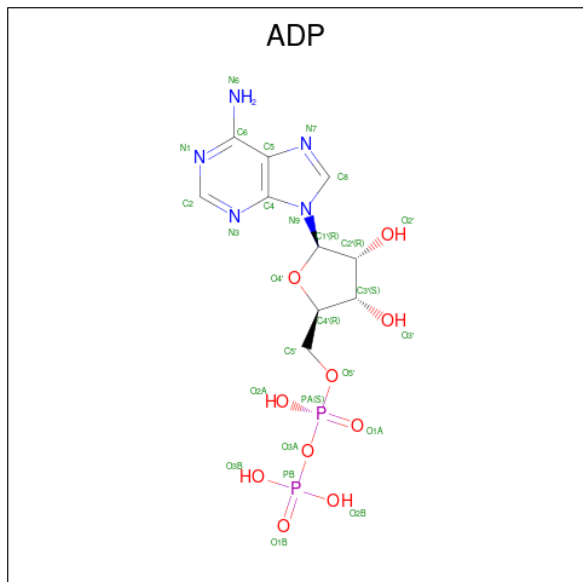
- Molecule 2 is a DNA chain called DNA (5'-D(P*TP*CP*CP*CP*TP*GP*TP*TP*TP*CP*AP*CP*GP*TP*GP*GP*AP*AP*CP*AP*CP*CP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	24	Total	C	N	O	P	0	0	0
			483	231	81	147	24			

- Molecule 3 is a DNA chain called DNA (5'-D(P*AP*GP*GP*GP*TP*GP*TP*TP*CP*CP*AP*CP*GP*TP*GP*AP*AP*AP*CP*AP*GP*GP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	24	Total	C	N	O	P	0	0	0
			501	236	100	141	24			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

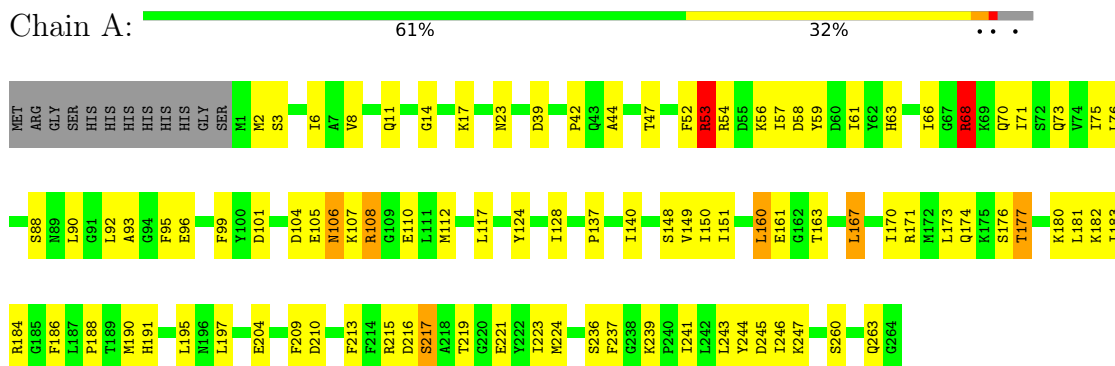
- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

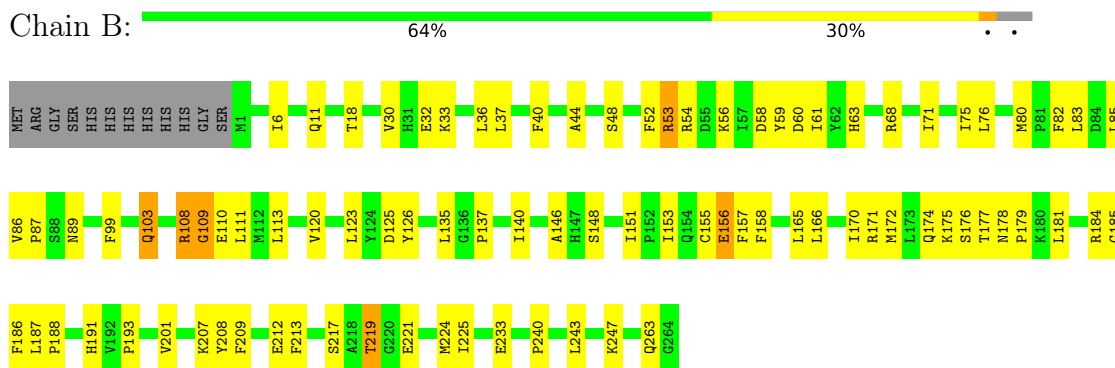
3 Residue-property plots

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.

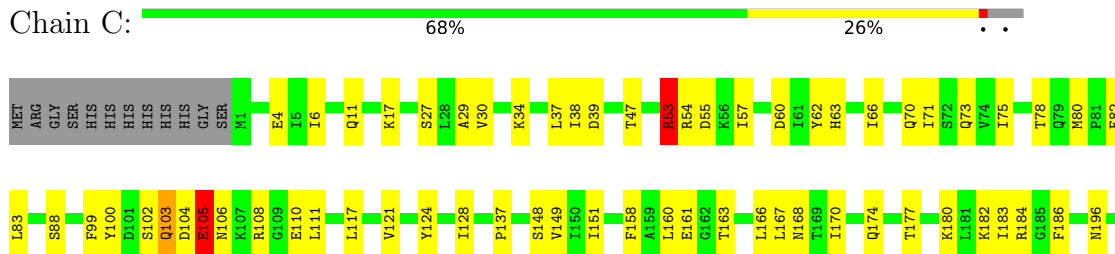
• Molecule 1: SpoOJ regulator (Soj)



• Molecule 1: SpoOJ regulator (Soj)



• Molecule 1: SpoOJ regulator (Soj)





- Molecule 1: SpoOJ regulator (Soj)

Chain D: 61% 32%



- Molecule 2: DNA (5'-D(P*TP*CP*CP*CP*TP*GP*TP*TP*TP*CP*AP*CP*GP*TP*GP*GP*AP*AP*CP*AP*CP*CP*CP*T)-3')

Chain E: 54% 46%



- Molecule 3: DNA (5'-D(P*AP*GP*GP*GP*TP*GP*TP*TP*CP*CP*AP*CP*GP*TP*GP*A*P*AP*AP*CP*AP*GP*GP*GP*A)-3')

Chain F: 38% 63%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	74.77Å 74.69Å 80.75Å 71.54° 71.59° 67.73°	Depositor
Resolution (Å)	29.95 – 2.51 29.95 – 2.51	Depositor EDS
% Data completeness (in resolution range)	95.7 (29.95-2.51) 91.5 (29.95-2.51)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.205 , 0.258 0.210 , 0.264	Depositor DCC
R_{free} test set	2443 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.457 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9308	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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, ADP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.69	2/2091 (0.1%)	1.12	12/2826 (0.4%)
1	B	0.86	5/2091 (0.2%)	1.10	11/2826 (0.4%)
1	C	0.67	2/2091 (0.1%)	1.04	10/2826 (0.4%)
1	D	0.77	3/2091 (0.1%)	1.10	14/2826 (0.5%)
2	E	0.55	0/538	0.76	0/826
3	F	0.60	0/564	0.80	0/870
All	All	0.73	12/9466 (0.1%)	1.06	47/13000 (0.4%)

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	A	0	3
1	B	0	1
1	C	0	2
1	D	0	1
All	All	0	7

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	108	ARG	CG-CD	14.55	1.96	1.52
1	B	108	ARG	NE-CZ	12.06	1.46	1.33
1	B	108	ARG	N-CA	-8.23	1.34	1.46
1	B	108	ARG	CZ-NH1	7.93	1.43	1.32
1	A	177	THR	CB-CG2	-7.41	1.28	1.52
1	D	199	LYS	CG-CD	6.93	1.73	1.52
1	D	54	ARG	CG-CD	-6.41	1.33	1.52
1	C	103	GLN	CG-CD	-6.27	1.36	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	68	ARG	CG-CD	-5.39	1.36	1.52
1	B	109	GLY	CA-C	5.20	1.60	1.51
1	D	59	TYR	C-N	5.02	1.40	1.33
1	C	103	GLN	CB-CG	-5.02	1.37	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	53	ARG	CB-CG-CD	-15.15	76.44	111.30
1	A	108	ARG	CG-CD-NE	-13.04	83.31	112.00
1	C	180	LYS	CB-CG-CD	11.80	138.44	111.30
1	A	180	LYS	CB-CG-CD	-11.67	84.47	111.30
1	A	68	ARG	CB-CG-CD	11.08	136.78	111.30
1	A	68	ARG	CA-CB-CG	10.89	135.89	114.10
1	C	180	LYS	CG-CD-CE	-10.19	87.87	111.30
1	C	53	ARG	CB-CG-CD	9.77	133.76	111.30
1	D	54	ARG	CB-CG-CD	-9.61	89.19	111.30
1	D	68	ARG	N-CA-CB	-9.08	96.24	110.22
1	C	53	ARG	NE-CZ-NH2	-8.81	111.27	119.20
1	D	171	ARG	CG-CD-NE	-8.67	92.92	112.00
1	C	103	GLN	N-CA-CB	-8.56	97.04	110.22
1	B	108	ARG	NE-CZ-NH1	8.53	130.03	121.50
1	C	53	ARG	CD-NE-CZ	-8.27	112.82	124.40
1	B	108	ARG	CG-CD-NE	-8.11	94.16	112.00
1	A	53	ARG	CA-CB-CG	-8.09	97.91	114.10
1	C	105	GLU	CB-CG-CD	6.70	123.99	112.60
1	D	53	ARG	CA-C-N	-6.53	110.56	122.38
1	D	53	ARG	C-N-CA	-6.53	110.56	122.38
1	A	68	ARG	CG-CD-NE	-6.46	97.78	112.00
1	A	180	LYS	CA-CB-CG	6.24	126.58	114.10
1	D	54	ARG	NE-CZ-NH2	-6.08	113.73	119.20
1	D	60	ASP	CB-CA-C	-6.08	98.41	111.78
1	D	107	LYS	CA-C-N	5.91	128.19	120.28
1	D	107	LYS	C-N-CA	5.91	128.19	120.28
1	A	53	ARG	CD-NE-CZ	5.90	132.66	124.40
1	C	53	ARG	NE-CZ-NH1	5.82	127.32	121.50
1	A	167	LEU	CA-CB-CG	-5.80	96.00	116.30
1	A	108	ARG	N-CA-C	5.75	120.43	113.41
1	B	108	ARG	O-C-N	-5.66	115.01	122.42
1	B	52	PHE	CA-C-N	-5.64	112.87	120.87
1	B	52	PHE	C-N-CA	-5.64	112.87	120.87
1	C	53	ARG	CB-CA-C	-5.63	98.56	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	53	ARG	N-CA-CB	5.47	118.76	110.45
1	A	106	ASN	N-CA-C	5.45	122.30	114.39
1	A	217	SER	N-CA-CB	5.37	118.02	110.12
1	B	108	ARG	CA-C-N	-5.36	108.90	121.34
1	B	108	ARG	C-N-CA	-5.36	108.90	121.34
1	B	219	THR	CA-CB-OG1	-5.30	101.66	109.60
1	B	103	GLN	CA-CB-CG	-5.25	103.60	114.10
1	D	170	ILE	CA-C-N	-5.20	112.50	121.66
1	D	170	ILE	C-N-CA	-5.20	112.50	121.66
1	D	68	ARG	NE-CZ-NH2	-5.16	114.56	119.20
1	D	59	TYR	CB-CA-C	5.06	118.56	110.16
1	B	18	THR	OG1-CB-CG2	-5.06	99.19	109.30
1	D	118	GLU	CA-CB-CG	-5.02	104.06	114.10

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	GLU	Peptide
1	A	53	ARG	Sidechain
1	A	68	ARG	Sidechain
1	B	156	GLU	Sidechain
1	C	105	GLU	Peptide
1	C	53	ARG	Mainchain
1	D	60	ASP	Sidechain

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	2053	0	2133	87	2
1	B	2053	0	2133	86	1
1	C	2053	0	2133	86	1
1	D	2053	0	2133	96	2
2	E	483	0	272	15	0
3	F	501	0	269	24	0
4	A	27	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	27	0	12	0	0
4	C	27	0	12	0	0
4	D	27	0	12	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
All	All	9308	0	9121	370	3

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 20.

All (370) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ARG:CD	1:B:108:ARG:CG	1.96	1.41
1:C:237:PHE:HE2	1:C:244:TYR:CE2	1.45	1.33
1:C:237:PHE:CE2	1:C:244:TYR:CE2	2.25	1.24
1:D:60:ASP:OD2	1:D:62:TYR:HB3	1.46	1.14
1:B:60:ASP:OD2	1:B:89:ASN:HB3	1.52	1.08
2:E:23:DC:N3	3:F:2:DG:N1	2.16	0.94
1:C:78:THR:HG22	1:C:80:MET:H	1.33	0.94
1:C:161:GLU:OE2	1:D:54:ARG:NH1	2.01	0.92
1:C:161:GLU:OE2	1:D:54:ARG:NH2	2.04	0.90
2:E:24:DT:H3	3:F:1:DA:H61	1.17	0.90
2:E:23:DC:N4	3:F:2:DG:O6	2.06	0.89
1:C:161:GLU:OE2	1:D:54:ARG:CZ	2.22	0.87
1:B:54:ARG:HD3	1:B:89:ASN:HA	1.56	0.86
1:A:217:SER:OG	1:A:263:GLN:NE2	2.08	0.86
1:A:110:GLU:OE1	1:A:177:THR:HG21	1.77	0.84
1:D:23:ASN:HD21	1:D:241:ILE:H	1.23	0.83
2:E:24:DT:O4	3:F:1:DA:N6	2.12	0.82
1:B:151:ILE:HD12	1:B:166:LEU:HD21	1.60	0.82
1:D:105:GLU:HB3	1:D:106:ASN:OD1	1.79	0.82
1:C:237:PHE:CD2	1:C:244:TYR:CD2	2.67	0.82
1:A:217:SER:OG	1:A:263:GLN:CD	2.24	0.81
1:A:6:ILE:HB	1:A:128:ILE:HD13	1.63	0.80
1:D:59:TYR:HA	1:D:63:HIS:CE1	2.16	0.80
1:D:53:ARG:HB2	1:D:55:ASP:OD2	1.81	0.79
1:A:173:LEU:HA	1:A:176:SER:HB3	1.61	0.79
1:C:237:PHE:HE2	1:C:244:TYR:HE2	1.29	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:PHE:HD2	1:C:244:TYR:CD2	2.03	0.76
1:D:36:LEU:HB3	1:D:127:ILE:HD13	1.66	0.76
1:A:176:SER:OG	1:A:177:THR:HG23	1.85	0.75
1:A:219:THR:HG23	1:A:221:GLU:H	1.52	0.75
1:D:3:SER:HB2	1:D:124:TYR:O	1.87	0.74
1:A:161:GLU:OE2	1:B:54:ARG:NH2	2.20	0.73
1:C:219:THR:HB	1:C:221:GLU:HG3	1.69	0.73
1:A:167:LEU:HD22	1:A:171:ARG:HH12	1.54	0.73
1:B:174:GLN:HB2	1:B:175:LYS:HE2	1.71	0.73
1:D:59:TYR:CD2	1:D:63:HIS:CE1	2.76	0.73
1:C:99:PHE:CE1	1:C:103:GLN:NE2	2.57	0.72
1:B:108:ARG:CG	1:B:108:ARG:NE	2.51	0.72
1:B:171:ARG:HD2	1:B:175:LYS:HE3	1.72	0.72
1:C:237:PHE:CD2	1:C:244:TYR:CE2	2.77	0.72
1:A:106:ASN:HB3	1:A:108:ARG:HE	1.55	0.72
1:B:103:GLN:OE1	1:B:109:GLY:N	2.24	0.71
1:A:70:GLN:H	1:A:73:GLN:NE2	1.89	0.71
1:A:54:ARG:HA	1:A:57:ILE:HD12	1.73	0.70
1:B:170:ILE:O	1:B:174:GLN:HG3	1.91	0.70
1:A:6:ILE:HB	1:A:128:ILE:CD1	2.20	0.70
1:C:106:ASN:HB2	1:C:108:ARG:HG3	1.74	0.70
1:C:237:PHE:HE2	1:C:244:TYR:CZ	2.07	0.69
1:C:53:ARG:HB2	1:C:55:ASP:OD2	1.92	0.69
1:C:233:GLU:O	1:C:236:SER:HB3	1.92	0.69
1:D:111:LEU:HD22	1:D:114:LYS:HD3	1.74	0.69
1:D:102:SER:HB3	1:D:108:ARG:HD2	1.73	0.68
1:A:6:ILE:HG12	1:A:148:SER:OG	1.93	0.68
1:C:237:PHE:CE2	1:C:244:TYR:CZ	2.80	0.68
1:C:100:TYR:O	1:C:103:GLN:HB2	1.94	0.68
1:C:70:GLN:OE1	1:C:73:GLN:NE2	2.27	0.68
1:D:102:SER:CB	1:D:108:ARG:HD2	2.24	0.67
1:D:110:GLU:OE2	1:D:177:THR:HG21	1.95	0.66
1:D:207:LYS:HE2	1:D:208:TYR:CZ	2.30	0.66
1:A:54:ARG:HA	1:A:57:ILE:CD1	2.25	0.66
1:A:161:GLU:CD	1:B:54:ARG:HH22	2.03	0.66
1:C:11:GLN:OE1	1:C:166:LEU:HB2	1.95	0.66
1:D:59:TYR:HD2	1:D:63:HIS:CE1	2.11	0.66
1:C:54:ARG:HA	1:C:57:ILE:HD13	1.77	0.66
1:B:148:SER:HB2	1:B:184:ARG:HB2	1.76	0.65
1:C:106:ASN:CB	1:C:108:ARG:HG3	2.26	0.65
1:B:108:ARG:NH1	1:B:111:LEU:HD12	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:4:DG:H2''	3:F:5:DT:H5'	1.78	0.65
1:A:260:SER:HA	1:A:263:GLN:NE2	2.13	0.64
1:D:53:ARG:HD3	1:D:53:ARG:N	2.11	0.64
1:C:105:GLU:HG3	1:C:106:ASN:ND2	2.13	0.64
1:A:39:ASP:OD2	1:A:44:ALA:HA	1.98	0.64
1:D:148:SER:HB2	1:D:184:ARG:HB2	1.80	0.64
2:E:24:DT:C4	3:F:1:DA:N6	2.66	0.63
2:E:20:DA:H2''	2:E:21:DC:H5''	1.80	0.63
1:D:36:LEU:HD11	1:D:86:VAL:HG23	1.82	0.62
1:A:70:GLN:HB2	1:A:73:GLN:HE21	1.63	0.62
1:C:70:GLN:OE1	1:C:70:GLN:N	2.32	0.62
1:C:6:ILE:HG12	1:C:148:SER:OG	2.00	0.62
1:A:210:ASP:OD1	1:A:215:ARG:NH2	2.32	0.62
1:B:54:ARG:HG2	1:B:89:ASN:OD1	2.00	0.61
1:B:33:LYS:HD2	1:B:126:TYR:HE1	1.65	0.61
3:F:10:DC:H2'	3:F:11:DA:C8	2.35	0.61
1:A:70:GLN:H	1:A:73:GLN:HE21	1.48	0.61
1:A:170:ILE:HG23	1:A:181:LEU:HD23	1.80	0.61
1:B:156:GLU:OE1	1:B:157:PHE:N	2.34	0.61
1:D:37:LEU:HD21	1:D:46:ALA:HB1	1.80	0.61
1:D:59:TYR:HA	1:D:63:HIS:HE1	1.66	0.60
1:D:171:ARG:HG2	1:D:174:GLN:OE1	2.00	0.60
1:B:191:HIS:HE1	1:B:193:PRO:HB3	1.67	0.60
1:A:58:ASP:OD2	1:A:59:TYR:CD2	2.55	0.60
1:B:36:LEU:HD11	1:B:86:VAL:HG23	1.84	0.59
1:B:120:VAL:HA	1:B:123:LEU:HD12	1.83	0.59
1:D:23:ASN:ND2	1:D:241:ILE:H	1.98	0.59
1:C:4:GLU:OE2	1:C:184:ARG:NH1	2.35	0.59
1:C:99:PHE:CG	1:C:137:PRO:HB3	2.37	0.59
1:B:53:ARG:O	1:B:56:LYS:HG2	2.02	0.59
1:C:39:ASP:HB3	1:C:88:SER:HB2	1.84	0.59
1:D:23:ASN:ND2	1:D:241:ILE:HG22	2.18	0.59
1:D:54:ARG:HA	1:D:57:ILE:HG12	1.84	0.59
1:D:51:GLY:O	1:D:53:ARG:NE	2.36	0.58
1:C:196:ASN:OD1	1:C:199:LYS:HE3	2.03	0.58
1:A:71:ILE:O	1:A:75:ILE:HG13	2.03	0.58
1:C:111:LEU:HD21	1:C:177:THR:CG2	2.33	0.58
3:F:5:DT:H2''	3:F:6:DG:C8	2.39	0.58
2:E:24:DT:H3	3:F:1:DA:N6	1.96	0.58
1:A:216:ASP:CG	1:A:219:THR:HG22	2.29	0.58
1:D:11:GLN:HE21	1:D:135:LEU:HG	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:MET:HB2	1:D:83:LEU:HB3	1.85	0.58
1:C:106:ASN:CB	1:C:108:ARG:HE	2.17	0.58
3:F:20:DA:OP2	3:F:20:DA:H2'	2.04	0.58
1:B:171:ARG:CD	1:B:175:LYS:HE3	2.34	0.57
1:B:188:PRO:HG2	1:B:224:MET:HE2	1.86	0.57
1:C:78:THR:HB	1:C:83:LEU:O	2.04	0.57
1:C:54:ARG:NH1	1:D:161:GLU:CD	2.62	0.57
1:C:110:GLU:OE1	1:C:177:THR:OG1	2.21	0.57
1:C:62:TYR:CZ	1:C:66:ILE:HD11	2.38	0.57
1:C:245:ASP:OD2	1:C:245:ASP:C	2.45	0.57
3:F:22:DG:H1'	3:F:23:DG:H5'	1.87	0.57
1:D:63:HIS:O	1:D:68:ARG:HB2	2.06	0.56
1:D:99:PHE:CG	1:D:137:PRO:HB3	2.40	0.56
1:A:188:PRO:HG2	1:A:224:MET:SD	2.45	0.56
1:D:151:ILE:HD12	1:D:166:LEU:CD2	2.36	0.56
1:A:223:ILE:C	1:A:224:MET:HG2	2.31	0.56
1:A:246:ILE:HG23	1:A:247:LYS:HG2	1.87	0.56
1:B:240:PRO:HD2	1:B:243:LEU:HD12	1.88	0.56
1:C:237:PHE:CE2	1:C:244:TYR:CD2	2.85	0.56
1:D:16:GLY:O	1:D:20:THR:HG23	2.07	0.55
2:E:21:DC:H2''	2:E:22:DC:H5''	1.88	0.55
3:F:21:DG:H2''	3:F:22:DG:H8	1.70	0.55
1:C:106:ASN:CG	1:C:108:ARG:HE	2.14	0.55
1:D:11:GLN:NE2	1:D:135:LEU:HG	2.21	0.55
1:C:170:ILE:O	1:C:174:GLN:HG3	2.07	0.55
1:C:27:SER:HA	1:C:242:LEU:HD11	1.87	0.55
3:F:4:DG:H2'	3:F:5:DT:C6	2.42	0.55
1:D:184:ARG:NH2	1:D:261:ILE:HG23	2.22	0.54
1:A:66:ILE:HG22	1:A:68:ARG:HB2	1.88	0.54
1:B:99:PHE:CG	1:B:137:PRO:HB3	2.43	0.54
1:A:161:GLU:CD	1:B:54:ARG:NH2	2.64	0.54
1:B:217:SER:OG	1:B:263:GLN:NE2	2.40	0.54
2:E:22:DC:H2'	2:E:23:DC:O4'	2.08	0.54
1:A:66:ILE:CG2	1:A:68:ARG:HB2	2.39	0.53
1:A:191:HIS:CD2	1:A:224:MET:HE3	2.42	0.53
1:D:61:ILE:HG23	1:D:92:LEU:HD13	1.89	0.53
1:B:33:LYS:HD2	1:B:126:TYR:CE1	2.44	0.53
1:A:42:PRO:O	1:A:90:LEU:HD13	2.09	0.53
1:A:151:ILE:HD12	1:A:186:PHE:HE1	1.72	0.53
1:A:171:ARG:HG2	1:A:174:GLN:OE1	2.09	0.53
1:B:59:TYR:HA	1:B:63:HIS:CE1	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:TYR:HD1	1:A:63:HIS:ND1	2.06	0.53
1:A:101:ASP:O	1:A:104:ASP:HB2	2.09	0.53
1:D:20:THR:HB	1:D:254:TYR:HE1	1.74	0.53
1:C:71:ILE:O	1:C:75:ILE:HG13	2.09	0.52
1:A:245:ASP:OD2	1:A:245:ASP:C	2.52	0.52
1:B:207:LYS:HE3	1:B:208:TYR:CE1	2.44	0.52
1:D:150:ILE:HD11	1:D:261:ILE:HD11	1.91	0.52
1:B:30:VAL:C	1:B:32:GLU:H	2.16	0.52
1:B:146:ALA:O	1:B:178:ASN:ND2	2.42	0.52
1:D:186:PHE:HE1	1:D:209:PHE:CG	2.27	0.52
1:A:137:PRO:HA	1:A:140:ILE:HG22	1.91	0.52
1:B:188:PRO:HG3	1:B:213:PHE:HE1	1.75	0.52
1:A:8:VAL:HG22	1:A:150:ILE:HB	1.93	0.51
1:A:11:GLN:OE1	1:A:11:GLN:N	2.42	0.51
1:A:148:SER:HB2	1:A:184:ARG:HB2	1.91	0.51
1:A:239:LYS:HB3	1:A:243:LEU:HD12	1.93	0.51
1:A:99:PHE:CG	1:A:137:PRO:HB3	2.46	0.51
1:D:60:ASP:OD2	1:D:62:TYR:CB	2.38	0.51
1:D:149:VAL:HG23	1:D:181:LEU:HD11	1.91	0.51
1:D:154:GLN:HG2	1:D:156:GLU:HG3	1.93	0.51
1:A:246:ILE:O	1:A:246:ILE:HG13	2.11	0.51
1:D:151:ILE:HD12	1:D:166:LEU:HD21	1.92	0.51
1:B:44:ALA:HB3	1:B:54:ARG:NE	2.26	0.50
1:A:216:ASP:OD2	1:A:219:THR:HG22	2.11	0.50
1:C:184:ARG:NH2	1:C:261:ILE:O	2.45	0.50
1:A:23:ASN:HD21	1:A:241:ILE:H	1.59	0.50
1:A:95:PHE:HE1	1:A:112:MET:SD	2.34	0.50
1:C:54:ARG:HH12	1:D:161:GLU:CD	2.19	0.50
1:C:99:PHE:CD1	1:C:137:PRO:HG3	2.47	0.50
1:C:151:ILE:HD12	1:C:186:PHE:HE2	1.77	0.50
1:D:60:ASP:H	1:D:63:HIS:CE1	2.29	0.50
1:D:71:ILE:O	1:D:74:VAL:HG22	2.12	0.50
1:D:58:ASP:O	1:D:59:TYR:CD2	2.65	0.50
3:F:1:DA:H2''	3:F:2:DG:H8	1.76	0.50
1:A:163:THR:HG22	1:A:167:LEU:HD12	1.94	0.49
1:A:219:THR:HG21	1:A:221:GLU:OE1	2.11	0.49
1:D:120:VAL:HA	1:D:123:LEU:HD12	1.94	0.49
1:A:95:PHE:CE1	1:A:112:MET:SD	3.05	0.49
1:A:110:GLU:OE1	1:A:177:THR:CG2	2.56	0.49
1:D:58:ASP:O	1:D:59:TYR:HD2	1.95	0.49
1:D:102:SER:OG	1:D:108:ARG:HD2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:GLU:OE1	1:B:54:ARG:NH2	2.44	0.49
1:A:151:ILE:HD12	1:A:186:PHE:CE1	2.48	0.49
1:B:108:ARG:HH21	1:B:110:GLU:N	2.11	0.49
1:B:108:ARG:HH21	1:B:110:GLU:H	1.61	0.49
1:C:148:SER:HB2	1:C:184:ARG:HB2	1.93	0.49
1:D:229:VAL:O	1:D:233:GLU:HG2	2.13	0.49
1:A:186:PHE:HE2	1:A:209:PHE:CG	2.31	0.48
2:E:14:DT:H2''	2:E:15:DG:C8	2.48	0.48
3:F:21:DG:H2''	3:F:22:DG:C8	2.47	0.48
1:A:14:GLY:O	1:A:190:MET:HE1	2.13	0.48
1:A:59:TYR:HA	1:A:63:HIS:ND1	2.28	0.48
1:A:106:ASN:O	1:A:107:LYS:HG2	2.13	0.48
1:A:236:SER:HA	1:B:157:PHE:CE1	2.48	0.48
1:D:170:ILE:O	1:D:174:GLN:HG3	2.13	0.48
1:A:61:ILE:HG23	1:A:92:LEU:HD13	1.95	0.48
1:A:151:ILE:HB	1:A:186:PHE:HD1	1.78	0.48
1:B:108:ARG:NE	1:B:108:ARG:O	2.43	0.48
1:C:149:VAL:HG13	1:C:183:ILE:HA	1.95	0.48
1:C:105:GLU:HG3	1:C:106:ASN:CG	2.38	0.48
1:B:151:ILE:HD12	1:B:166:LEU:CD2	2.38	0.48
1:A:2:MET:HE3	1:A:2:MET:O	2.14	0.48
1:A:197:LEU:HB2	1:B:233:GLU:HG2	1.96	0.48
1:D:163:THR:O	1:D:167:LEU:HG	2.14	0.48
1:B:110:GLU:HG2	1:B:140:ILE:HG21	1.97	0.47
1:C:38:ILE:HD11	1:C:117:LEU:HD21	1.96	0.47
1:A:149:VAL:HG12	1:A:182:LYS:O	2.13	0.47
1:B:80:MET:HB2	1:B:83:LEU:HB3	1.95	0.47
1:D:36:LEU:HB3	1:D:127:ILE:CD1	2.41	0.47
1:B:186:PHE:HE2	1:B:209:PHE:CG	2.33	0.47
1:D:60:ASP:C	1:D:60:ASP:OD1	2.54	0.47
1:C:151:ILE:HD12	1:C:186:PHE:CE2	2.49	0.47
1:B:76:LEU:HD11	1:B:87:PRO:HG3	1.96	0.47
3:F:2:DG:H2'	3:F:3:DG:C8	2.50	0.47
1:C:99:PHE:O	1:C:102:SER:HB3	2.15	0.47
2:E:21:DC:H2''	2:E:22:DC:O4'	2.15	0.47
3:F:6:DG:OP2	3:F:6:DG:H3'	2.15	0.46
1:B:108:ARG:NE	1:B:108:ARG:HA	2.30	0.46
1:D:149:VAL:CG1	1:D:151:ILE:HD11	2.45	0.46
1:A:75:ILE:O	1:A:76:LEU:HD12	2.15	0.46
1:B:58:ASP:C	1:B:59:TYR:HD2	2.23	0.46
1:B:247:LYS:HE3	1:B:247:LYS:HB2	1.73	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:ARG:NH1	1:D:161:GLU:OE1	2.48	0.46
1:D:10:ASN:HA	1:D:166:LEU:HD22	1.97	0.46
1:D:103:GLN:C	1:D:105:GLU:H	2.24	0.46
1:D:184:ARG:NH1	1:D:264:GLY:C	2.74	0.46
1:A:23:ASN:ND2	1:A:241:ILE:H	2.14	0.46
1:A:191:HIS:CE1	1:A:224:MET:HE1	2.51	0.46
1:C:100:TYR:HA	1:C:103:GLN:CD	2.41	0.46
1:A:137:PRO:HA	1:A:140:ILE:CG2	2.45	0.46
1:D:45:ASN:O	1:D:49:SER:OG	2.28	0.46
1:D:59:TYR:HA	1:D:63:HIS:ND1	2.30	0.46
1:B:44:ALA:HB3	1:B:54:ARG:CZ	2.46	0.46
1:C:158:PHE:CZ	1:D:45:ASN:HB3	2.51	0.46
1:D:216:ASP:O	1:D:220:GLY:N	2.45	0.46
2:E:24:DT:N3	3:F:1:DA:N6	2.57	0.46
1:A:188:PRO:HG3	1:A:213:PHE:HE2	1.81	0.45
1:B:71:ILE:O	1:B:75:ILE:HG13	2.17	0.45
1:D:108:ARG:HD3	1:D:112:MET:SD	2.56	0.45
1:C:4:GLU:CD	1:C:184:ARG:HH12	2.23	0.45
1:D:20:THR:HG22	1:D:254:TYR:OH	2.16	0.45
1:D:170:ILE:HG22	1:D:171:ARG:HG3	1.98	0.45
1:A:54:ARG:HH11	1:A:90:LEU:HD21	1.80	0.45
1:A:167:LEU:CD2	1:A:183:ILE:HD11	2.47	0.45
1:A:59:TYR:HD1	1:A:63:HIS:CE1	2.35	0.45
1:B:40:PHE:CZ	1:B:113:LEU:HD22	2.52	0.45
1:B:156:GLU:CD	1:B:158:PHE:H	2.23	0.45
1:D:60:ASP:OD1	1:D:60:ASP:N	2.48	0.45
3:F:1:DA:C2'	3:F:2:DG:H8	2.30	0.45
1:D:95:PHE:CE2	1:D:112:MET:HE2	2.52	0.45
1:C:53:ARG:HH21	1:C:53:ARG:HD3	1.11	0.45
1:C:34:LYS:HE3	1:C:124:TYR:CE1	2.52	0.45
1:C:99:PHE:CD1	1:C:103:GLN:NE2	2.85	0.45
1:D:57:ILE:HD12	1:D:57:ILE:HG23	1.74	0.45
1:A:167:LEU:HD23	1:A:167:LEU:HA	1.42	0.44
1:B:53:ARG:HD2	1:B:53:ARG:HA	1.42	0.44
1:C:60:ASP:OD1	1:C:63:HIS:ND1	2.49	0.44
1:D:44:ALA:HB3	1:D:54:ARG:NH1	2.32	0.44
1:C:30:VAL:HG21	1:C:242:LEU:HD13	1.98	0.44
1:B:32:GLU:HA	1:B:82:PHE:CD1	2.53	0.44
1:B:191:HIS:CE1	1:B:193:PRO:HB3	2.50	0.44
1:C:117:LEU:O	1:C:121:VAL:HG23	2.17	0.44
1:B:61:ILE:HB	1:B:86:VAL:CG1	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:236:SER:HA	1:D:157:PHE:CE1	2.53	0.44
1:D:117:LEU:HD23	1:D:117:LEU:HA	1.62	0.44
1:B:108:ARG:O	1:B:108:ARG:HG3	2.17	0.44
1:A:160:LEU:HD13	1:A:204:GLU:HG2	1.98	0.44
1:D:184:ARG:NH1	1:D:264:GLY:O	2.48	0.44
3:F:9:DC:H2''	3:F:10:DC:C6	2.52	0.44
2:E:5:DT:H2'	2:E:6:DG:C8	2.53	0.44
1:A:167:LEU:HD22	1:A:171:ARG:NH1	2.28	0.43
1:A:182:LYS:HB2	1:A:182:LYS:HE3	1.94	0.43
1:B:156:GLU:OE2	1:B:158:PHE:HB2	2.18	0.43
1:D:231:LEU:HD11	1:D:250:GLY:HA3	2.00	0.43
1:A:53:ARG:O	1:A:56:LYS:HB3	2.18	0.43
1:A:96:GLU:CD	1:A:96:GLU:H	2.26	0.43
1:C:53:ARG:O	1:C:57:ILE:HD12	2.18	0.43
1:C:106:ASN:HB3	1:C:108:ARG:HG3	2.00	0.43
2:E:22:DC:H1'	3:F:4:DG:N2	2.33	0.43
1:B:187:LEU:HD11	1:B:225:ILE:HD11	2.01	0.43
3:F:23:DG:H2''	3:F:24:DA:H5''	1.99	0.43
1:B:11:GLN:HE21	1:B:135:LEU:HG	1.84	0.43
1:C:160:LEU:CD1	1:C:204:GLU:HG2	2.48	0.43
1:D:55:ASP:OD2	1:D:56:LYS:N	2.51	0.43
1:D:102:SER:O	1:D:105:GLU:HB2	2.18	0.43
1:A:170:ILE:O	1:A:174:GLN:HG3	2.18	0.43
1:B:174:GLN:HG2	1:B:181:LEU:HB3	2.00	0.43
1:C:70:GLN:N	1:C:70:GLN:CD	2.77	0.43
1:C:151:ILE:HB	1:C:186:PHE:HD2	1.83	0.43
1:B:59:TYR:CG	1:B:63:HIS:ND1	2.87	0.43
1:C:38:ILE:CD1	1:C:117:LEU:HD21	2.49	0.43
1:C:11:GLN:HG3	1:D:43:GLN:OE1	2.19	0.43
1:C:106:ASN:HB3	1:C:108:ARG:HE	1.83	0.43
1:D:110:GLU:OE2	1:D:177:THR:CG2	2.66	0.43
1:A:117:LEU:HD23	1:A:117:LEU:HA	1.68	0.43
1:C:242:LEU:HD23	1:C:242:LEU:HA	1.83	0.43
1:B:37:LEU:O	1:B:85:LEU:HD12	2.18	0.43
1:B:59:TYR:CG	1:B:63:HIS:CE1	3.06	0.43
1:C:37:LEU:HD12	1:C:128:ILE:O	2.19	0.43
1:C:71:ILE:HD13	1:C:117:LEU:HD23	2.00	0.43
1:D:32:GLU:HG2	1:D:82:PHE:CE2	2.53	0.43
1:B:32:GLU:HB3	1:B:82:PHE:CE1	2.53	0.42
1:B:110:GLU:OE1	1:B:177:THR:HG21	2.19	0.42
1:B:148:SER:HB2	1:B:184:ARG:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:LYS:HG2	1:D:13:GLY:N	2.34	0.42
1:C:39:ASP:OD1	1:C:47:THR:OG1	2.26	0.42
1:D:103:GLN:NE2	1:D:107:LYS:HA	2.34	0.42
1:B:188:PRO:HG3	1:B:213:PHE:CE1	2.55	0.42
1:C:163:THR:O	1:C:167:LEU:HD22	2.19	0.42
1:C:168:ASN:N	1:C:168:ASN:HD22	2.16	0.42
1:D:184:ARG:HH12	1:D:264:GLY:C	2.28	0.42
1:A:47:THR:HG22	1:A:52:PHE:HB2	2.00	0.42
1:B:174:GLN:O	1:B:179:PRO:HA	2.19	0.42
1:D:71:ILE:O	1:D:75:ILE:HG13	2.19	0.42
1:B:11:GLN:NE2	1:B:135:LEU:HG	2.35	0.42
1:D:150:ILE:HD11	1:D:261:ILE:CD1	2.50	0.42
1:A:39:ASP:HB3	1:A:88:SER:HB2	2.02	0.42
1:B:40:PHE:CE2	1:B:113:LEU:HD22	2.54	0.42
1:B:153:ILE:O	1:B:188:PRO:HA	2.20	0.42
1:B:185:GLY:HA2	1:B:212:GLU:O	2.19	0.42
1:B:207:LYS:HE3	1:B:208:TYR:CZ	2.54	0.42
1:B:172:MET:O	1:B:176:SER:OG	2.29	0.42
1:B:48:SER:HB2	1:B:54:ARG:NH1	2.35	0.41
3:F:3:DG:H8	3:F:3:DG:OP2	2.03	0.41
1:B:6:ILE:HA	1:B:148:SER:O	2.20	0.41
1:B:110:GLU:HG2	1:B:140:ILE:CG2	2.50	0.41
1:C:263:GLN:OE1	1:C:263:GLN:HA	2.20	0.41
1:D:151:ILE:HB	1:D:186:PHE:HD2	1.85	0.41
1:C:149:VAL:CG1	1:C:183:ILE:HA	2.51	0.41
1:D:171:ARG:CG	1:D:174:GLN:OE1	2.67	0.41
1:D:103:GLN:C	1:D:105:GLU:N	2.76	0.41
1:D:159:ALA:O	1:D:163:THR:HG23	2.20	0.41
1:B:176:SER:HB2	1:B:177:THR:HG23	2.03	0.41
1:A:3:SER:HB2	1:A:124:TYR:O	2.21	0.41
1:A:93:ALA:O	1:A:96:GLU:OE1	2.39	0.41
1:D:99:PHE:O	1:D:102:SER:HB2	2.20	0.41
1:A:17:LYS:HE2	1:A:17:LYS:HB2	1.80	0.41
1:B:11:GLN:HG2	1:B:165:LEU:HD23	2.02	0.41
1:D:188:PRO:HG3	1:D:213:PHE:HE1	1.86	0.41
1:D:226:PRO:HG3	1:D:249:ASN:HB3	2.03	0.41
2:E:2:DC:H2"	2:E:3:DC:C6	2.56	0.41
1:D:99:PHE:CD1	1:D:137:PRO:HB3	2.56	0.41
1:D:207:LYS:HE2	1:D:208:TYR:CE2	2.55	0.41
1:B:207:LYS:HE2	1:B:207:LYS:HB3	1.75	0.40
1:C:54:ARG:HA	1:C:57:ILE:CD1	2.47	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:GLY:HA3	1:D:214:PHE:CE1	2.56	0.40
1:A:195:LEU:CD2	3:F:17:DA:H4'	2.52	0.40
1:B:80:MET:HE2	1:B:80:MET:HB3	1.89	0.40
1:B:219:THR:OG1	1:B:221:GLU:HG2	2.21	0.40
1:C:80:MET:HE3	1:C:243:LEU:HD21	2.02	0.40
1:B:155:CYS:HA	1:B:201:VAL:HG11	2.04	0.40
1:C:17:LYS:HB2	1:C:17:LYS:HE2	1.65	0.40
1:B:33:LYS:HD3	1:B:125:ASP:HB3	2.02	0.40
1:B:36:LEU:HD11	1:B:86:VAL:CG2	2.49	0.40
1:C:4:GLU:CD	1:C:184:ARG:NH1	2.80	0.40
1:C:117:LEU:HD23	1:C:117:LEU:HA	1.76	0.40
1:A:237:PHE:HE1	1:A:244:TYR:CE2	2.40	0.40
1:B:148:SER:CB	1:B:184:ARG:HD3	2.51	0.40
1:C:29:ALA:HB1	1:C:82:PHE:HB2	2.03	0.40
1:C:149:VAL:HG12	1:C:182:LYS:O	2.20	0.40
1:C:184:ARG:HH11	1:C:184:ARG:HD3	1.65	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ASP:OD2	1:D:68:ARG:NH2[1_446]	1.61	0.59
1:A:104:ASP:CG	1:D:68:ARG:NH2[1_446]	2.17	0.03
1:B:68:ARG:NH1	1:C:104:ASP:OD2[1_446]	2.18	0.02

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	262/276 (95%)	252 (96%)	9 (3%)	1 (0%)	30	49
1	B	262/276 (95%)	250 (95%)	12 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	262/276 (95%)	248 (95%)	14 (5%)	0	100	100
1	D	262/276 (95%)	249 (95%)	13 (5%)	0	100	100
All	All	1048/1104 (95%)	999 (95%)	48 (5%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	LEU

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	228/238 (96%)	228 (100%)	0	100	100
1	B	228/238 (96%)	228 (100%)	0	100	100
1	C	228/238 (96%)	228 (100%)	0	100	100
1	D	228/238 (96%)	227 (100%)	1 (0%)	84	93
All	All	912/952 (96%)	911 (100%)	1 (0%)	88	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	20	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	70	GLN
1	A	73	GLN
1	A	79	GLN
1	A	178	ASN

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Mol	Chain	Res	Type
1	A	255	GLN
1	A	263	GLN
1	B	11	GLN
1	B	45	ASN
1	B	191	HIS
1	B	259	GLN
1	C	79	GLN
1	C	168	ASN
1	C	174	GLN
1	C	259	GLN
1	D	11	GLN
1	D	23	ASN
1	D	103	GLN
1	D	259	GLN

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	ADP	D	500	5	28,29,29	1.63	4 (14%)	43,45,45	1.89	11 (25%)
4	ADP	C	500	5	28,29,29	1.51	4 (14%)	43,45,45	2.38	14 (32%)
4	ADP	A	500	5	28,29,29	1.49	3 (10%)	43,45,45	2.07	11 (25%)
4	ADP	B	500	5	28,29,29	1.60	5 (17%)	43,45,45	1.97	13 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ADP	D	500	5	-	1/16/32/32	0/3/3/3
4	ADP	C	500	5	-	0/16/32/32	0/3/3/3
4	ADP	A	500	5	-	0/16/32/32	0/3/3/3
4	ADP	B	500	5	-	1/16/32/32	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	500	ADP	C5-C4	5.12	1.48	1.39
4	B	500	ADP	C5-C4	4.81	1.47	1.39
4	A	500	ADP	C5-C4	4.70	1.47	1.39
4	D	500	ADP	PA-O3A	4.44	1.64	1.59
4	B	500	ADP	PA-O3A	4.40	1.64	1.59
4	C	500	ADP	C5-C4	4.33	1.46	1.39
4	C	500	ADP	C5-C6	3.70	1.51	1.41
4	A	500	ADP	C5-C6	3.57	1.50	1.41
4	A	500	ADP	C8-N7	2.70	1.36	1.31
4	D	500	ADP	C5-C6	2.58	1.48	1.41
4	C	500	ADP	PA-O3A	2.47	1.62	1.59
4	B	500	ADP	C5-C6	2.41	1.47	1.41
4	C	500	ADP	C8-N7	2.24	1.36	1.31
4	D	500	ADP	C8-N7	2.23	1.36	1.31
4	B	500	ADP	C5-N7	-2.23	1.35	1.39
4	B	500	ADP	C8-N7	2.07	1.35	1.31

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	500	ADP	C5-C4-N3	-6.84	117.30	126.72
4	A	500	ADP	C5-C4-N3	-5.95	118.53	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	500	ADP	C5-C4-N3	-5.68	118.89	126.72
4	D	500	ADP	C5-C4-N3	-5.52	119.11	126.72
4	C	500	ADP	N3-C4-N9	5.22	136.05	127.17
4	C	500	ADP	C4-C5-N7	-5.20	104.64	110.58
4	B	500	ADP	N3-C4-N9	4.85	135.42	127.17
4	D	500	ADP	N3-C4-N9	4.84	135.39	127.17
4	A	500	ADP	N3-C4-N9	4.71	135.19	127.17
4	C	500	ADP	C2-N3-C4	4.48	122.77	111.83
4	C	500	ADP	C4-N9-C8	4.19	110.14	105.74
4	C	500	ADP	C5-N7-C8	4.19	110.03	103.45
4	A	500	ADP	C2-N3-C4	4.16	121.99	111.83
4	C	500	ADP	N9-C8-N7	-3.76	108.60	113.94
4	B	500	ADP	O3'-C3'-C4'	-3.73	100.37	111.08
4	C	500	ADP	O3'-C3'-C4'	-3.67	100.53	111.08
4	A	500	ADP	C4-C5-N7	-3.65	106.41	110.58
4	A	500	ADP	C4-N9-C8	3.41	109.31	105.74
4	D	500	ADP	C2-N3-C4	3.38	120.08	111.83
4	B	500	ADP	C2-N3-C4	3.36	120.03	111.83
4	A	500	ADP	O3'-C3'-C4'	-3.32	101.56	111.08
4	B	500	ADP	N3-C2-N1	-3.07	123.93	128.58
4	D	500	ADP	N3-C2-N1	-3.02	124.01	128.58
4	A	500	ADP	N3-C2-N1	-3.02	124.02	128.58
4	B	500	ADP	C4-C5-N7	-3.00	107.15	110.58
4	C	500	ADP	C6-C5-N7	2.93	137.75	132.09
4	A	500	ADP	O2B-PB-O3A	-2.87	95.02	104.64
4	C	500	ADP	N3-C2-N1	-2.86	124.26	128.58
4	D	500	ADP	C4-N9-C8	2.85	108.73	105.74
4	C	500	ADP	O2A-PA-O1A	2.80	125.45	112.44
4	D	500	ADP	C4-C5-N7	-2.77	107.41	110.58
4	B	500	ADP	C4-N9-C8	2.77	108.65	105.74
4	A	500	ADP	C5-N7-C8	2.75	107.77	103.45
4	B	500	ADP	C5-N7-C8	2.70	107.69	103.45
4	D	500	ADP	O2A-PA-O3A	2.65	114.44	107.27
4	D	500	ADP	C2'-C3'-C4'	2.63	107.70	102.61
4	A	500	ADP	N9-C8-N7	-2.59	110.27	113.94
4	A	500	ADP	C6-C5-N7	2.49	136.89	132.09
4	D	500	ADP	O3'-C3'-C4'	-2.45	104.05	111.08
4	D	500	ADP	C5-N7-C8	2.19	106.90	103.45
4	B	500	ADP	O2A-PA-O3A	2.19	113.19	107.27
4	B	500	ADP	N9-C8-N7	-2.19	110.84	113.94
4	B	500	ADP	C2'-C3'-C4'	2.15	106.76	102.61
4	B	500	ADP	O3B-PB-O1B	2.13	119.14	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	500	ADP	C4'-O4'-C1'	2.13	114.17	109.47
4	C	500	ADP	O2'-C2'-C3'	2.12	118.60	111.82
4	C	500	ADP	O3A-PB-O1B	-2.06	100.18	111.04
4	C	500	ADP	O4'-C1'-C2'	-2.06	102.21	106.62
4	B	500	ADP	C2-N1-C6	2.04	122.07	118.73

There are no chirality outliers.

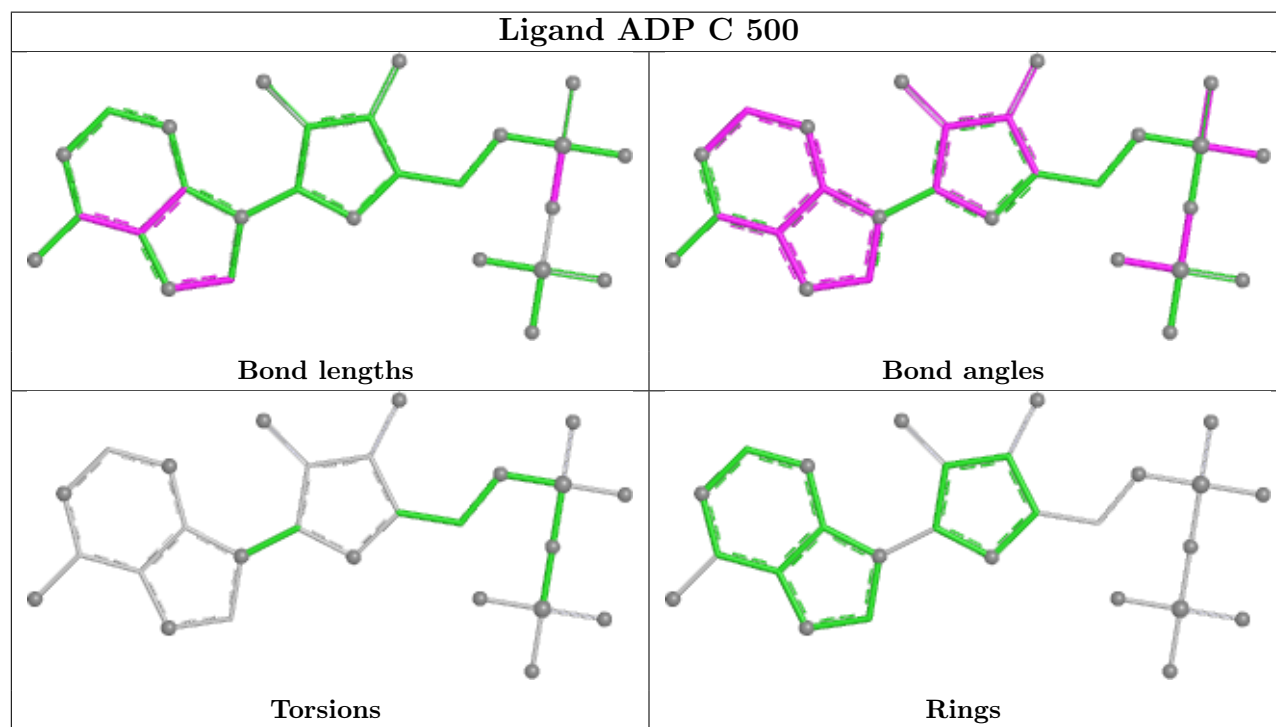
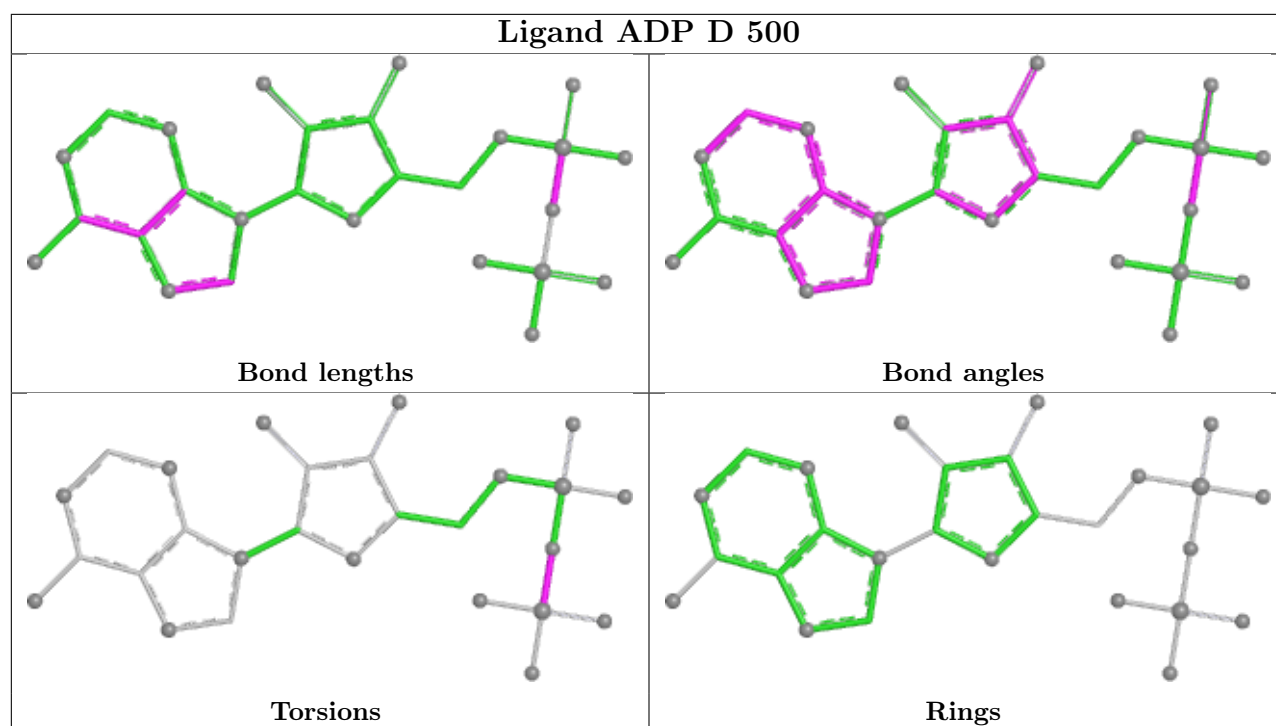
All (2) torsion outliers are listed below:

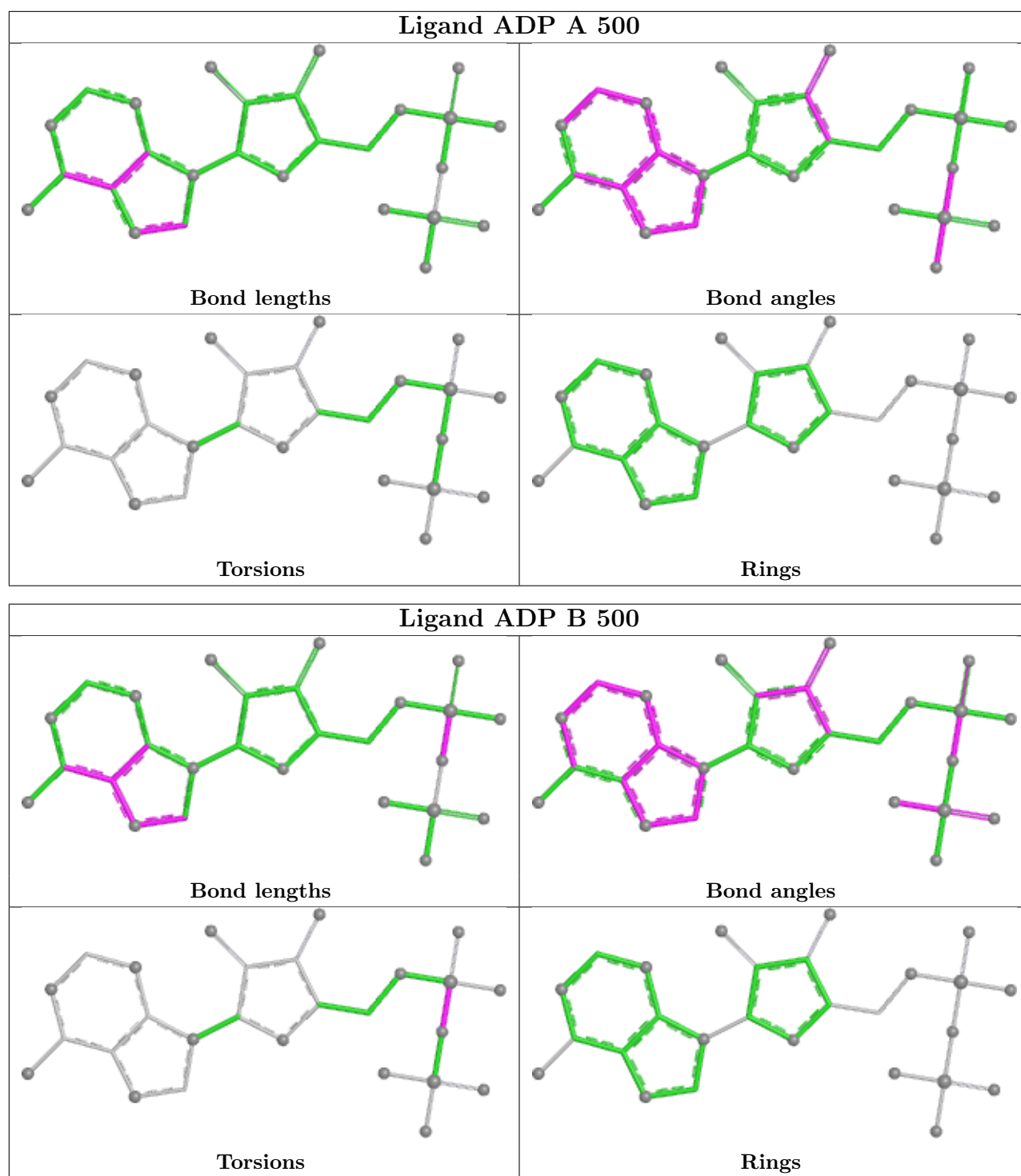
Mol	Chain	Res	Type	Atoms
4	B	500	ADP	PB-O3A-PA-O2A
4	D	500	ADP	PA-O3A-PB-O2B

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	264/276 (95%)	-1.18	0 100 100	32, 52, 86, 118	0
1	B	264/276 (95%)	-1.21	0 100 100	31, 52, 95, 124	0
1	C	264/276 (95%)	-1.18	0 100 100	30, 53, 85, 133	0
1	D	264/276 (95%)	-1.18	0 100 100	34, 53, 100, 124	0
2	E	24/24 (100%)	-0.68	0 100 100	42, 72, 170, 174	0
3	F	24/24 (100%)	-0.88	0 100 100	43, 76, 159, 169	0
All	All	1104/1152 (95%)	-1.17	0 100 100	30, 53, 101, 174	0

There are no RSRZ outliers to report.

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
5	MG	A	501	1/1	0.97	0.11	56,56,56,56	0

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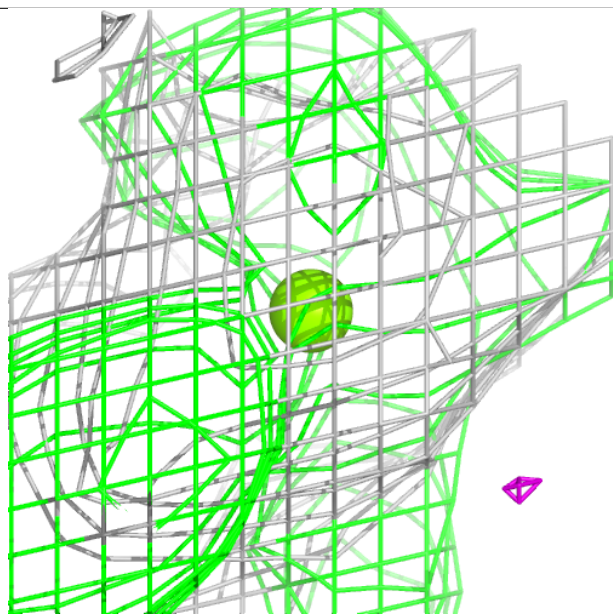
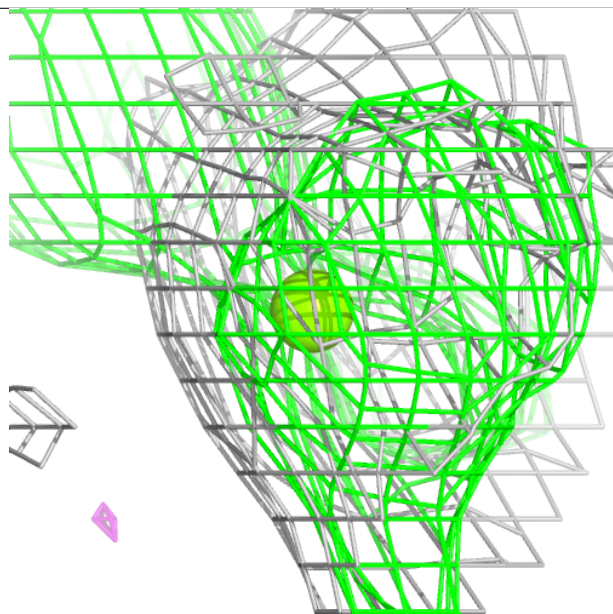
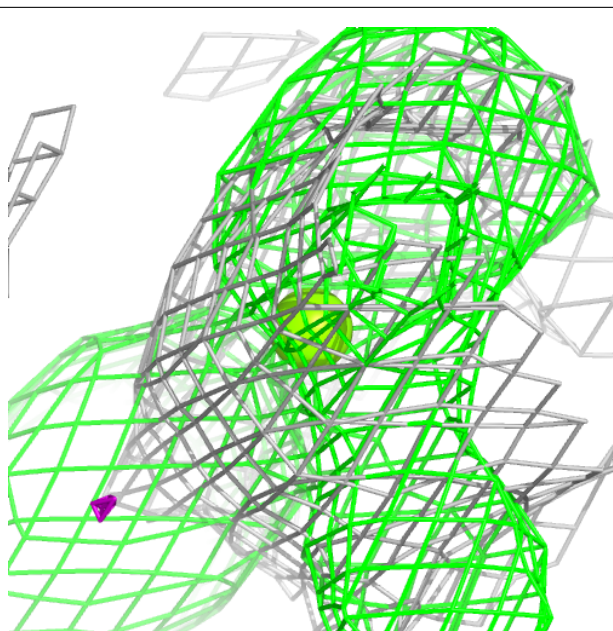
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	B	501	1/1	0.98	0.11	64,64,64,64	0
4	ADP	C	500	27/27	0.99	0.04	35,37,42,43	0
4	ADP	D	500	27/27	0.99	0.03	37,39,42,43	0
4	ADP	A	500	27/27	0.99	0.03	36,39,40,41	0
4	ADP	B	500	27/27	0.99	0.03	40,42,51,55	0
5	MG	C	501	1/1	0.99	0.07	50,50,50,50	0
5	MG	D	501	1/1	0.99	0.12	59,59,59,59	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.

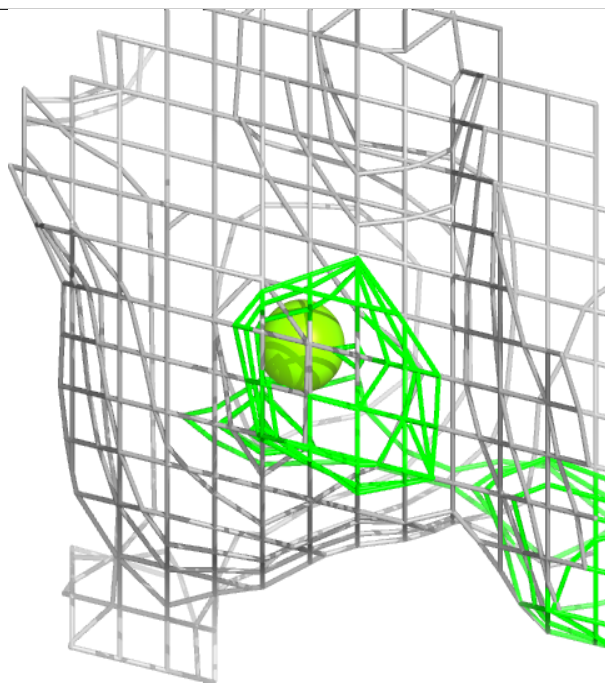
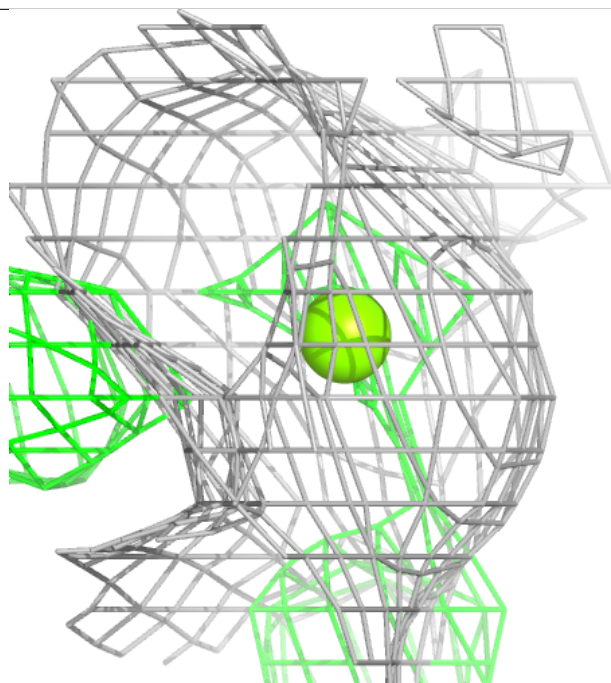
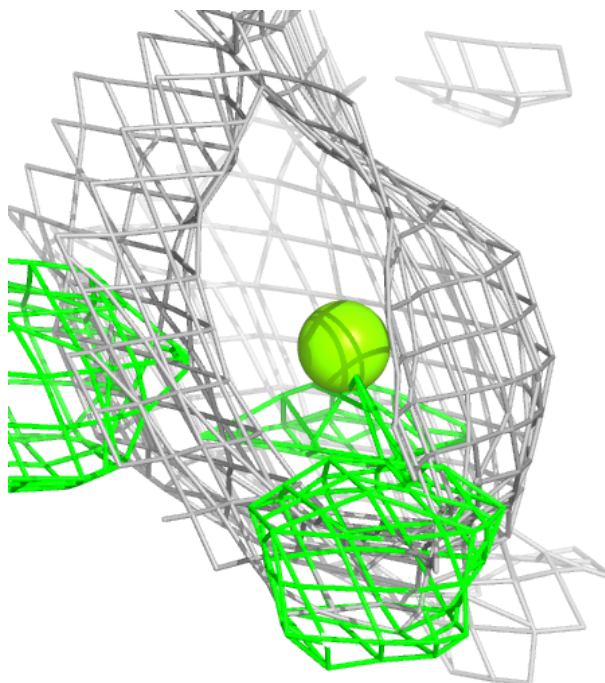
Electron density around MG A 501:

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



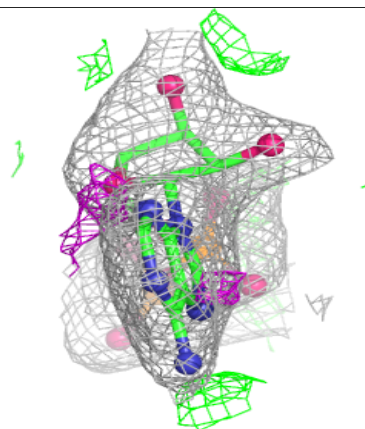
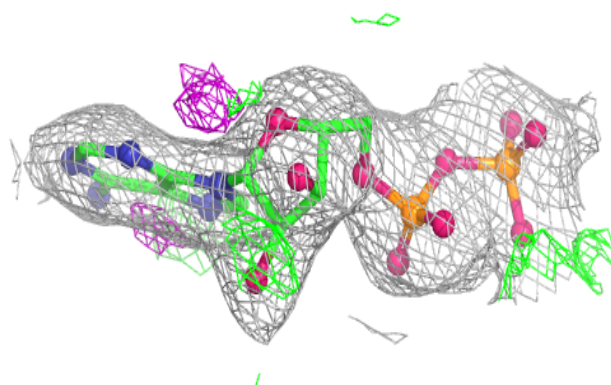
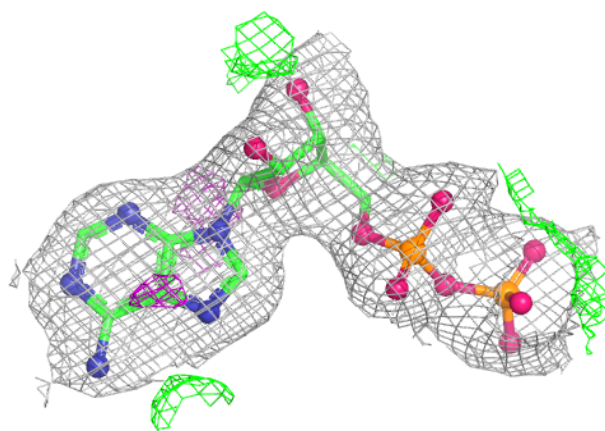
Electron density around MG B 501:

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



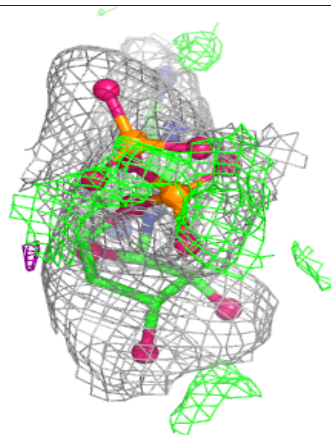
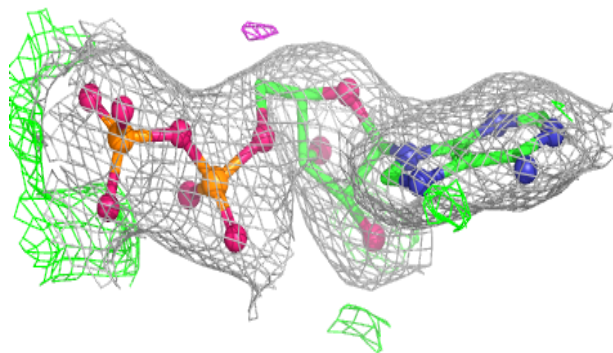
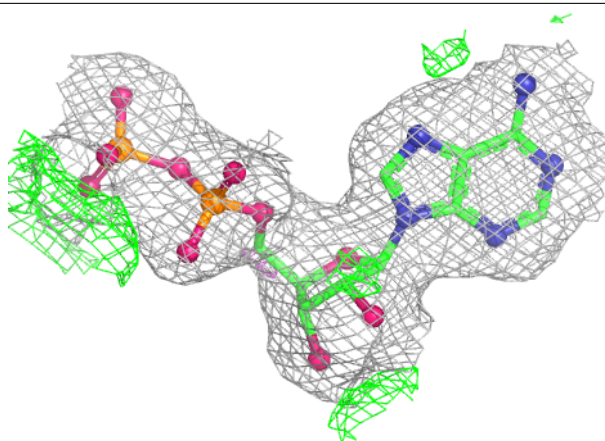
Electron density around ADP C 500:

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



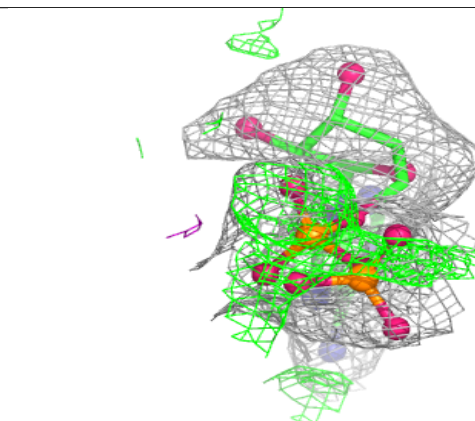
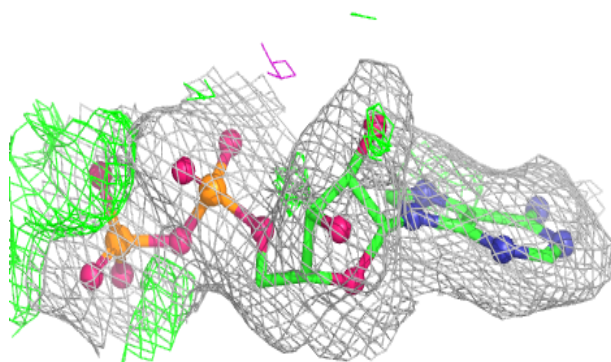
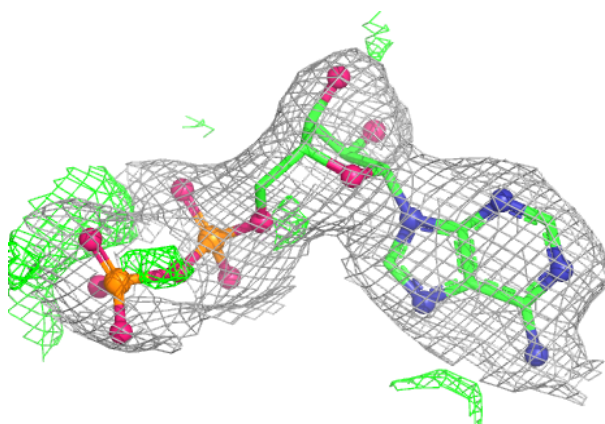
Electron density around ADP D 500:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

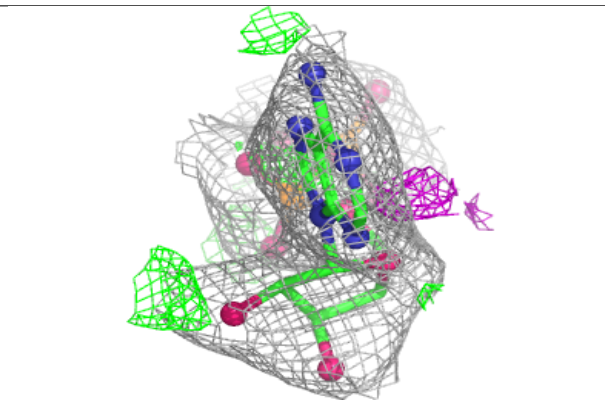
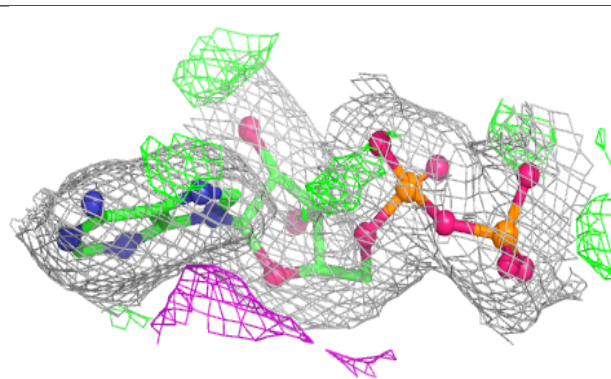
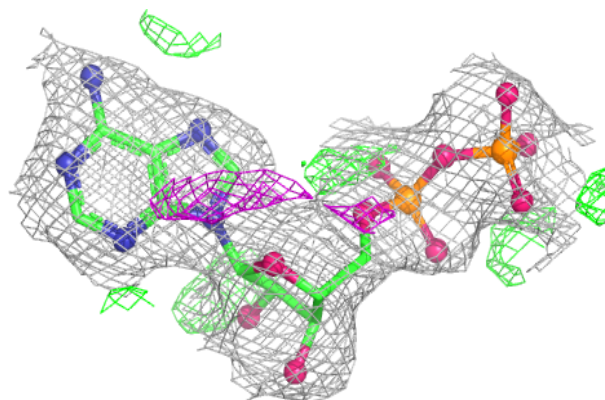


Electron density around ADP A 500:

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

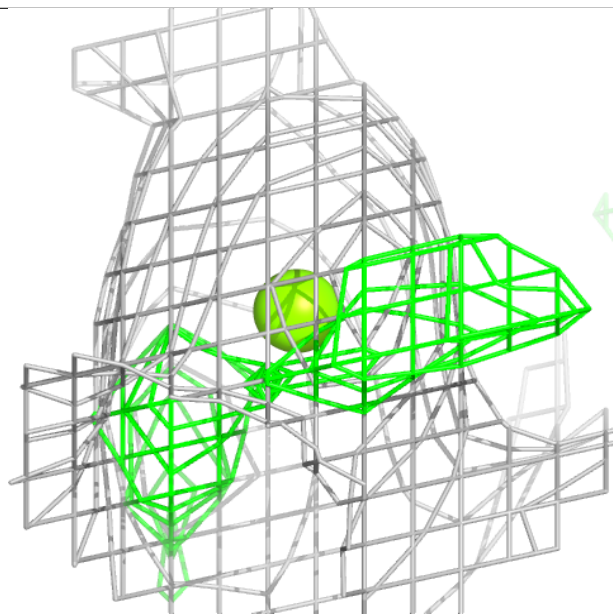
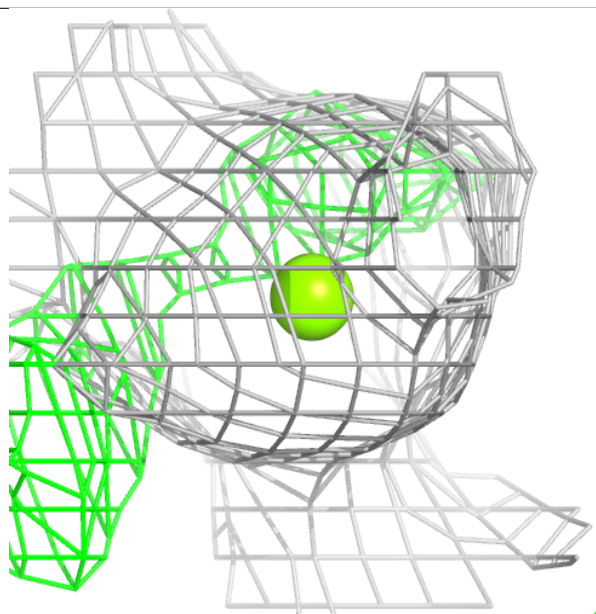
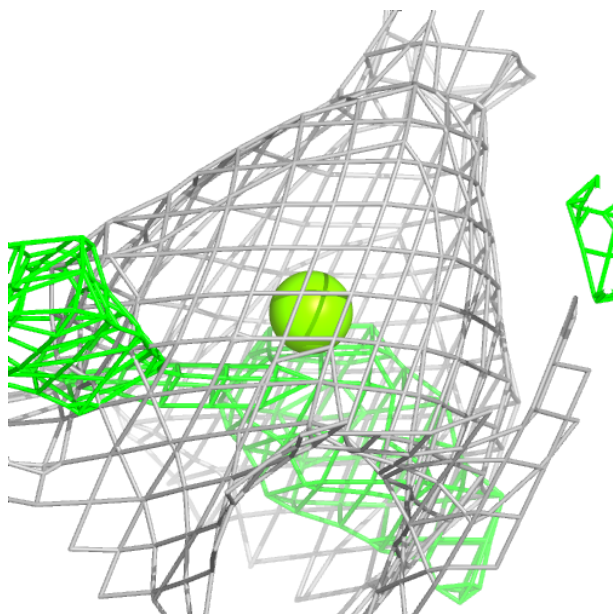
**Electron density around ADP B 500:**

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



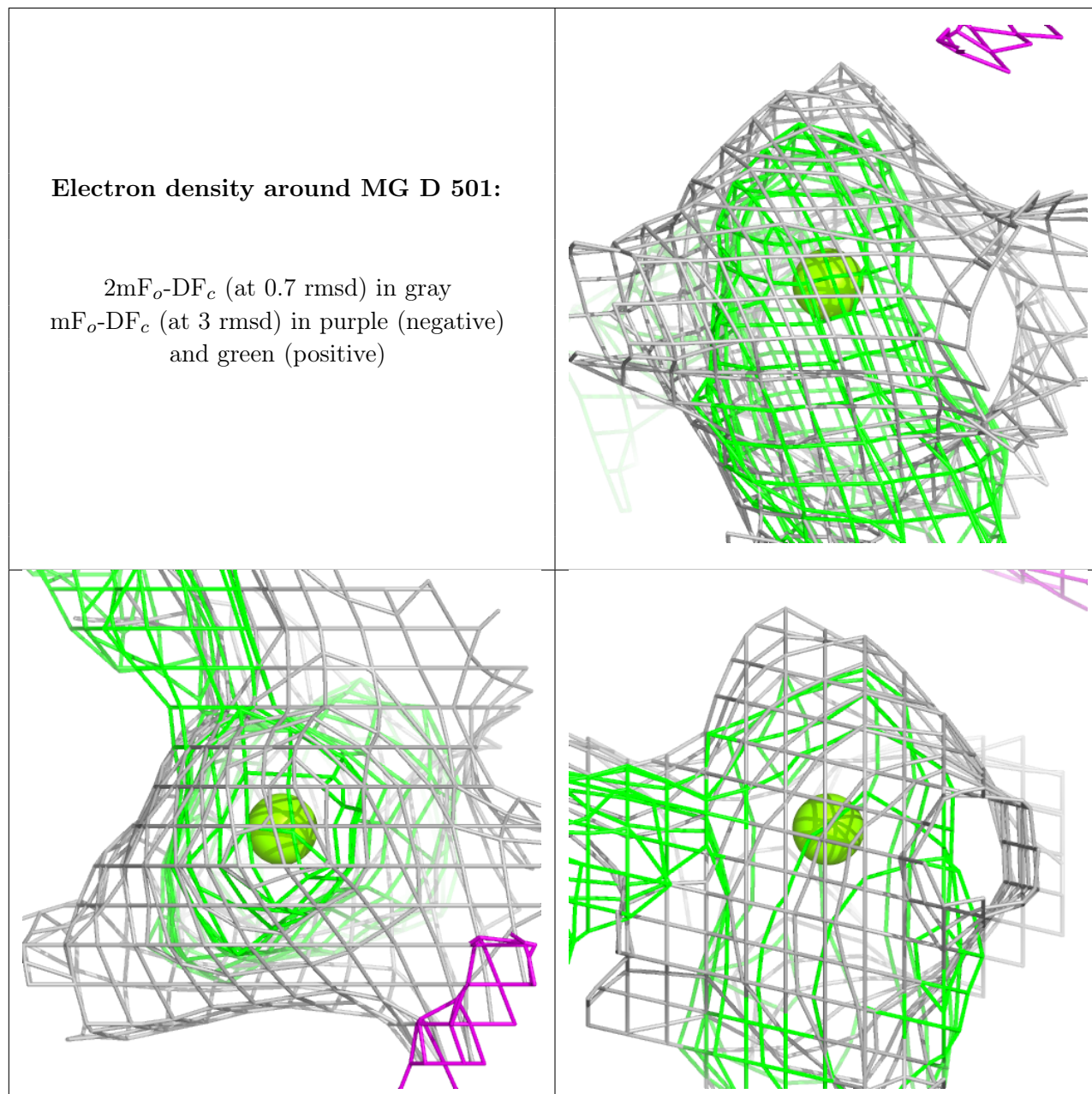
Electron density around MG C 501:

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



Electron density around MG D 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.