



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2023 – 04:55 PM JST

PDB ID : 5WSC
Title : Crystal of pyruvate kinase (PYK) from Mycobacterium tuberculosis in complex with Oxalate, soaked with allosteric activators AMP and Glucose 6-Phosphate
Authors : Zhong, W.; Cai, Q.; El Sahili, A.; Lescar, J.; Dedon, P.C.
Deposited on : 2016-12-06
Resolution : 2.40 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

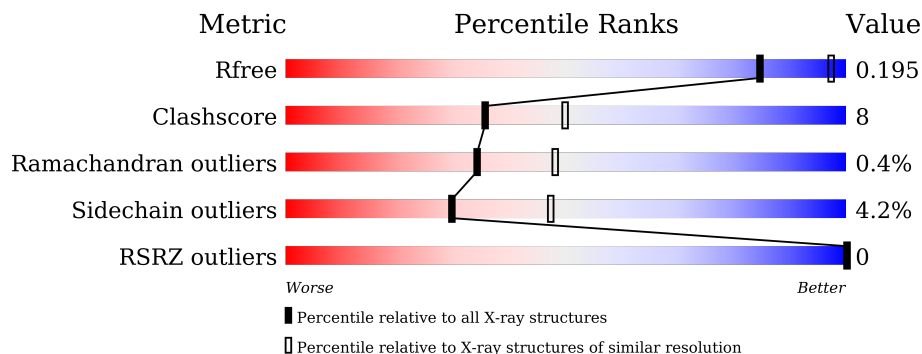
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	475	 82% 16% ..
1	B	475	 82% 15% ..
1	C	475	 81% 16% ..
1	D	475	 81% 16% ..

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
7	PO4	C	505	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 14702 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.

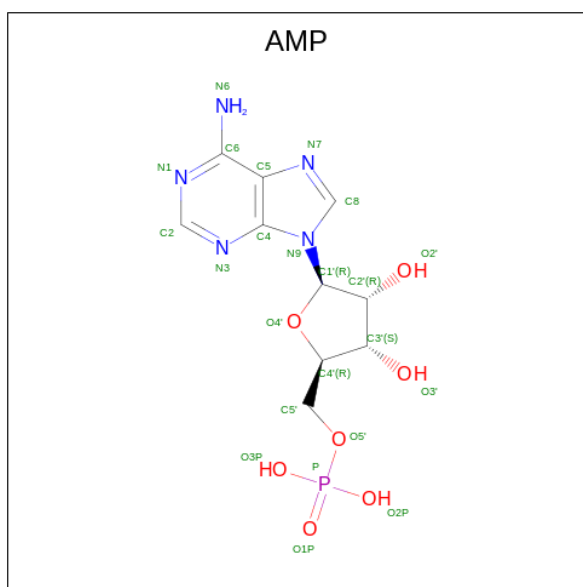
- Molecule 1 is a protein called Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	471	3543	2206	640	679	18	0	0	0
1	B	471	3543	2206	640	679	18	0	0	0
1	C	471	3543	2206	640	679	18	0	0	0
1	D	471	3543	2206	640	679	18	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

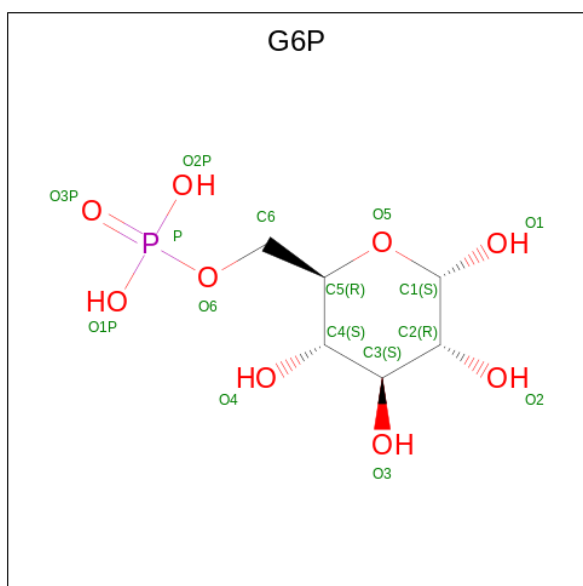
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P9WKE5
A	-1	GLY	-	expression tag	UNP P9WKE5
A	0	HIS	-	expression tag	UNP P9WKE5
B	-2	GLY	-	expression tag	UNP P9WKE5
B	-1	GLY	-	expression tag	UNP P9WKE5
B	0	HIS	-	expression tag	UNP P9WKE5
C	-2	GLY	-	expression tag	UNP P9WKE5
C	-1	GLY	-	expression tag	UNP P9WKE5
C	0	HIS	-	expression tag	UNP P9WKE5
D	-2	GLY	-	expression tag	UNP P9WKE5
D	-1	GLY	-	expression tag	UNP P9WKE5
D	0	HIS	-	expression tag	UNP P9WKE5

- Molecule 2 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C₁₀H₁₄N₅O₇P).



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

- Molecule 3 is 6-O-phosphono-alpha-D-glucopyranose (three-letter code: G6P) (formula: $C_6H_{13}O_9P$).

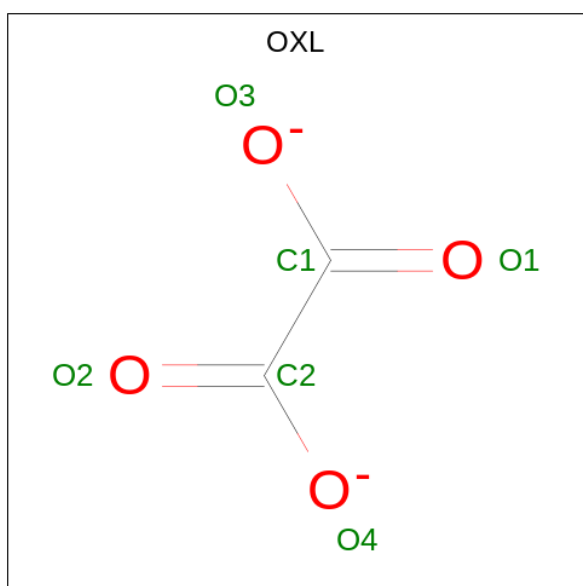


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		
3	C	1	Total	C	O	P	0	0
			16	6	9	1		
3	D	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is OXALATE ION (three-letter code: OXL) (formula: C₂O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	2	4		
5	B	1	Total	C	O	0	0
			6	2	4		

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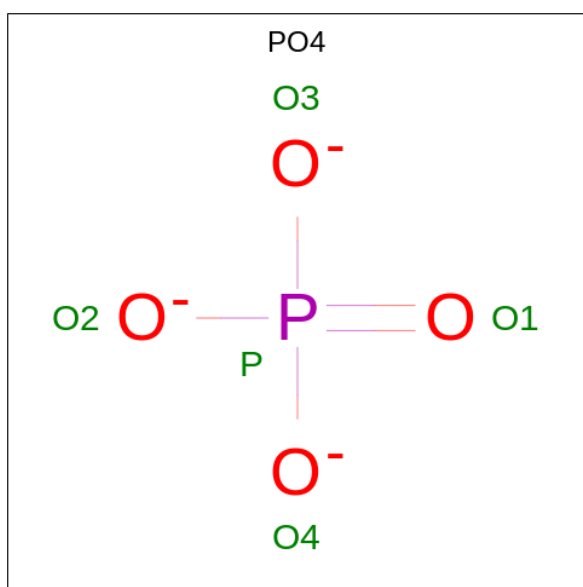
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total C O 6 2 4	0	0
5	D	1	Total C O 6 2 4	0	0

- Molecule 6 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total K 1 1	0	0

- Molecule 7 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	1	Total O P 5 4 1	0	0
7	D	1	Total O P 5 4 1	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	83	Total O 83 83	0	0
8	B	107	Total O 107 107	0	0

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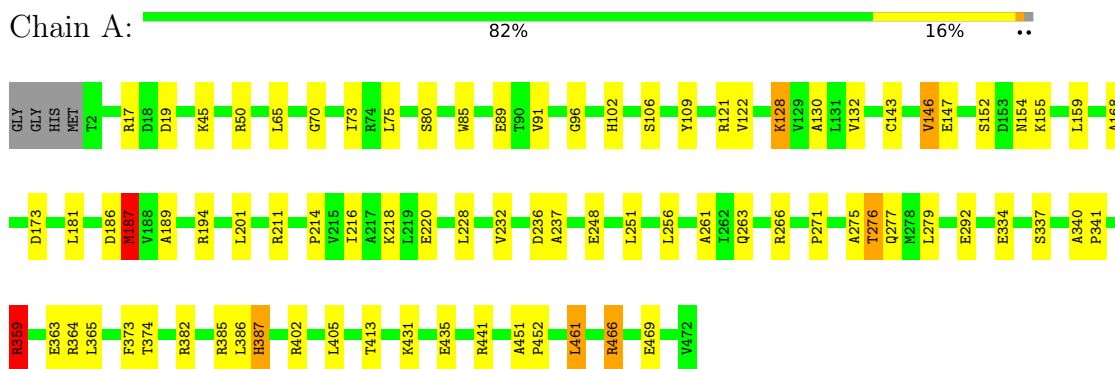
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	C	73	Total O 73 73	0	0
8	D	72	Total O 72 72	0	0

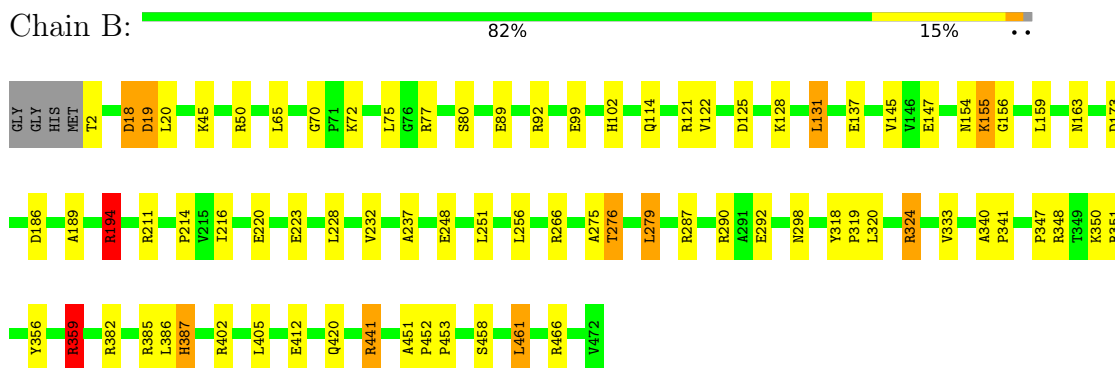
3 Residue-property plots [i](#)

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.

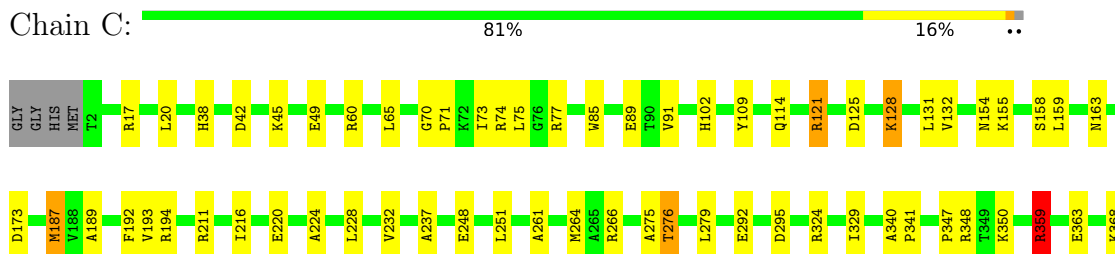
- Molecule 1: Pyruvate kinase



- Molecule 1: Pyruvate kinase



- Molecule 1: Pyruvate kinase





- Molecule 1: Pyruvate kinase

Chain D: 81% 16% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	124.37Å 124.37Å 144.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	53.85 – 2.40 47.09 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (53.85-2.40) 100.0 (47.09-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.179 , 0.227 0.148 , 0.195	Depositor DCC
R_{free} test set	4788 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtrriage
Anisotropy	0.596	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 7.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	0.469 for -h,-k,l 0.477 for h,-h-k,-l 0.470 for -k,-h,-l	Xtrriage
Reported twinning fraction	0.245 for H, K, L 0.248 for -K, -H, -L 0.254 for -h,-k,l 0.253 for K, H, -L	Depositor
Outliers	0 of 97489 reflections	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14702	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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: K, MG, G6P, PO4, OXL, 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	0.71	0/3593	0.79	3/4881 (0.1%)
1	B	0.71	0/3593	0.81	6/4881 (0.1%)
1	C	0.69	0/3593	0.79	4/4881 (0.1%)
1	D	0.70	0/3593	0.80	3/4881 (0.1%)
All	All	0.70	0/14372	0.80	16/19524 (0.1%)

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	B	0	1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	194	ARG	NE-CZ-NH2	7.61	124.10	120.30
1	A	359	ARG	NE-CZ-NH2	-7.01	116.79	120.30
1	C	211	ARG	NE-CZ-NH1	6.65	123.63	120.30
1	C	211	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	B	359	ARG	NE-CZ-NH1	-6.18	117.21	120.30
1	D	359	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	C	60	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	B	211	ARG	NE-CZ-NH2	-5.69	117.45	120.30
1	A	187	MET	CG-SD-CE	5.68	109.29	100.20
1	D	60	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	A	211	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	B	324	ARG	NE-CZ-NH2	5.45	123.03	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	359	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	D	211	ARG	NE-CZ-NH1	5.37	122.99	120.30
1	B	211	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	B	77	ARG	NE-CZ-NH2	-5.23	117.69	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	2	THR	Peptide

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	3543	0	3608	56	1
1	B	3543	0	3608	62	3
1	C	3543	0	3608	64	2
1	D	3543	0	3608	58	0
2	A	23	0	12	2	0
2	B	23	0	12	2	0
2	C	23	0	12	1	0
2	D	23	0	12	1	0
3	A	16	0	11	1	0
3	B	16	0	11	2	0
3	C	16	0	11	3	0
3	D	16	0	11	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	6	0	0	0	0
5	B	6	0	0	0	0
5	C	6	0	0	0	0
5	D	6	0	0	1	0
6	B	1	0	0	0	0
7	C	5	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	5	0	0	0	0
8	A	83	0	0	8	1
8	B	107	0	0	8	0
8	C	73	0	0	4	0
8	D	72	0	0	8	1
All	All	14702	0	14524	234	4

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 8.

All (234) 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:C:74:ARG:HH12	7:C:505:PO4:P	1.82	1.02
1:C:163:ASN:OD1	1:C:163:ASN:O	1.81	0.99
1:C:74:ARG:NH1	7:C:505:PO4:O1	1.96	0.97
1:C:38:HIS:NE2	7:C:505:PO4:O4	1.98	0.96
1:D:155:LYS:HE3	1:D:155:LYS:HA	1.50	0.94
1:A:466:ARG:HB3	1:A:469:GLU:HG3	1.50	0.90
1:B:298:ASN:HB3	8:B:647:HOH:O	1.73	0.87
1:A:334:GLU:OE1	8:A:601:HOH:O	1.94	0.82
1:D:331:CYS:O	1:D:335:GLU:HG3	1.83	0.78
1:D:72:LYS:HG3	1:D:74:ARG:HD2	1.67	0.77
1:A:122:VAL:HG22	1:A:159:LEU:HD11	1.66	0.77
1:A:431:LYS:HD3	8:A:605:HOH:O	1.85	0.76
1:C:74:ARG:NH1	7:C:505:PO4:P	2.59	0.73
1:C:382:ARG:HH12	3:C:502:G6P:H62	1.55	0.72
1:C:121:ARG:HD2	1:C:132:VAL:HG22	1.72	0.70
1:B:155:LYS:HA	1:B:155:LYS:NZ	2.06	0.69
1:C:125:ASP:HB2	1:C:155:LYS:HG2	1.74	0.69
1:D:432:SER:OG	8:D:601:HOH:O	2.07	0.69
1:B:194:ARG:HH21	1:B:194:ARG:HG3	1.57	0.68
1:D:122:VAL:HG22	1:D:159:LEU:HD11	1.78	0.66
1:B:122:VAL:HG22	1:B:159:LEU:HD11	1.78	0.66
5:D:504:OXL:O4	8:D:602:HOH:O	2.13	0.65
1:C:387:HIS:NE2	8:C:603:HOH:O	2.28	0.65
1:C:74:ARG:NH1	7:C:505:PO4:O4	2.30	0.65
1:D:382:ARG:HH22	3:D:502:G6P:H1	1.62	0.64
1:C:121:ARG:HG3	1:C:121:ARG:NH1	2.13	0.63
1:B:382:ARG:HH22	3:B:502:G6P:H1	1.63	0.63
1:A:122:VAL:HG22	1:A:159:LEU:CD1	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:50:ARG:NH2	8:D:604:HOH:O	2.19	0.62
1:B:348:ARG:HH11	1:B:348:ARG:HA	1.63	0.61
1:A:461:LEU:C	1:A:461:LEU:HD23	2.20	0.61
1:B:92:ARG:NH1	1:B:137:GLU:OE2	2.32	0.61
1:D:431:LYS:O	1:D:435:GLU:HG3	2.01	0.60
1:B:320:LEU:HD22	1:B:324:ARG:HH11	1.66	0.60
1:C:382:ARG:NH1	3:C:502:G6P:H62	2.17	0.60
1:B:279:LEU:HD12	1:B:292:GLU:HB3	1.82	0.59
1:B:276:THR:N	8:B:609:HOH:O	2.34	0.59
1:D:374:THR:HA	2:D:501:AMP:H5'2	1.85	0.59
1:A:431:LYS:O	1:A:435:GLU:HG3	2.03	0.59
1:A:128:LYS:HD3	1:A:128:LYS:N	2.18	0.59
1:C:125:ASP:OD2	1:C:155:LYS:HE2	2.03	0.58
1:A:85:TRP:CZ2	1:A:91:VAL:HG11	2.38	0.58
1:C:74:ARG:NE	1:C:154:ASN:O	2.29	0.58
1:B:155:LYS:HA	1:B:155:LYS:HZ2	1.67	0.58
1:D:15:THR:HA	1:D:20:LEU:HD23	1.84	0.57
1:C:85:TRP:CZ2	1:C:91:VAL:HG11	2.40	0.57
1:D:228:LEU:HD21	1:D:261:ALA:HA	1.85	0.57
1:A:70:GLY:HA2	1:A:173:ASP:OD2	2.05	0.56
1:B:461:LEU:HD12	1:B:461:LEU:C	2.26	0.56
1:C:65:LEU:HD13	1:C:216:ILE:HD12	1.87	0.56
1:D:186:ASP:C	1:D:187:MET:HG3	2.26	0.56
1:D:73:ILE:HG23	1:D:109:TYR:HB2	1.89	0.55
1:C:17:ARG:HG2	1:C:20:LEU:HG	1.89	0.55
1:A:218:LYS:NZ	8:A:604:HOH:O	2.36	0.55
1:B:70:GLY:HA2	1:B:173:ASP:OD2	2.07	0.54
1:A:128:LYS:HE3	8:A:661:HOH:O	2.06	0.54
1:B:223:GLU:OE2	8:B:601:HOH:O	2.18	0.54
1:B:125:ASP:HB2	1:B:155:LYS:HG2	1.88	0.54
1:C:386:LEU:O	1:C:387:HIS:O	2.25	0.54
1:C:128:LYS:HD2	1:C:128:LYS:N	2.21	0.54
1:D:461:LEU:C	1:D:461:LEU:HD12	2.28	0.54
1:A:228:LEU:HD21	1:A:261:ALA:HA	1.90	0.53
1:C:374:THR:HA	2:C:501:AMP:H5'2	1.90	0.53
1:D:186:ASP:O	1:D:187:MET:HG3	2.09	0.53
1:B:359:ARG:HD2	1:B:359:ARG:C	2.28	0.53
1:A:130:ALA:HB3	1:A:147:GLU:HG3	1.91	0.53
1:A:256:LEU:HD11	1:B:333:VAL:HA	1.90	0.52
1:C:38:HIS:CD2	7:C:505:PO4:O4	2.60	0.52
1:C:70:GLY:HA2	1:C:173:ASP:OD2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:LEU:HD13	1:D:216:ILE:HD12	1.90	0.52
1:C:228:LEU:HD21	1:C:261:ALA:HA	1.91	0.52
1:B:461:LEU:HD12	1:B:461:LEU:O	2.10	0.52
1:D:386:LEU:O	1:D:387:HIS:O	2.27	0.52
1:C:232:VAL:HG21	1:C:264:MET:CG	2.40	0.51
1:A:277:GLN:OE1	1:B:290:ARG:HG2	2.10	0.51
1:A:374:THR:HA	2:A:501:AMP:H5'2	1.91	0.51
1:D:70:GLY:HA2	1:D:173:ASP:OD2	2.10	0.51
1:C:73:ILE:HG23	1:C:109:TYR:HB2	1.92	0.51
1:A:386:LEU:O	1:A:387:HIS:O	2.29	0.51
1:B:50:ARG:NH1	8:B:619:HOH:O	2.44	0.51
1:D:348:ARG:HB3	8:D:607:HOH:O	2.10	0.51
1:B:451:ALA:HA	1:B:452:PRO:C	2.31	0.51
1:C:125:ASP:CB	1:C:155:LYS:HG2	2.39	0.51
1:C:419:MET:CE	8:C:654:HOH:O	2.59	0.51
1:D:85:TRP:CZ2	1:D:91:VAL:HG11	2.46	0.51
1:D:444:LEU:HD21	1:D:464:VAL:HG13	1.93	0.51
1:B:122:VAL:HG22	1:B:159:LEU:CD1	2.41	0.50
1:D:173:ASP:OD1	8:D:605:HOH:O	2.19	0.50
1:B:320:LEU:HD22	1:B:324:ARG:NH1	2.27	0.50
1:C:461:LEU:HD12	1:C:461:LEU:O	2.10	0.50
1:B:458:SER:OG	2:B:501:AMP:O3'	2.28	0.50
1:A:65:LEU:HD13	1:A:216:ILE:HD12	1.93	0.50
1:D:72:LYS:HG3	1:D:74:ARG:CD	2.39	0.50
1:C:216:ILE:HG12	1:C:237:ALA:HB3	1.93	0.49
1:D:17:ARG:HD2	1:D:18:ASP:H	1.77	0.49
1:B:402:ARG:HD3	8:B:620:HOH:O	2.12	0.49
1:D:220:GLU:O	1:D:248:GLU:HG3	2.12	0.49
1:B:386:LEU:O	1:B:387:HIS:O	2.30	0.49
1:C:121:ARG:CD	1:C:132:VAL:HG22	2.42	0.49
1:A:216:ILE:HG12	1:A:237:ALA:HB3	1.94	0.49
1:A:402:ARG:HG3	1:A:413:THR:HB	1.95	0.49
1:D:122:VAL:HG22	1:D:159:LEU:CD1	2.42	0.48
1:C:275:ALA:O	1:C:276:THR:HB	2.13	0.48
1:D:121:ARG:O	1:D:159:LEU:HD13	2.14	0.48
1:D:444:LEU:HD21	1:D:464:VAL:CG1	2.44	0.48
1:A:168:ALA:HB3	1:A:201:LEU:HD12	1.96	0.48
1:A:266:ARG:O	1:A:385:ARG:HD2	2.14	0.48
1:D:275:ALA:O	1:D:276:THR:HB	2.14	0.48
1:C:232:VAL:HG21	1:C:264:MET:HG3	1.96	0.47
1:C:158:SER:OG	1:C:248:GLU:OE1	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:451:ALA:HA	1:A:452:PRO:C	2.33	0.47
1:B:155:LYS:HA	1:B:155:LYS:HZ3	1.77	0.47
1:B:216:ILE:HG12	1:B:237:ALA:HB3	1.97	0.47
1:C:461:LEU:HD12	1:C:461:LEU:C	2.35	0.47
1:C:38:HIS:HE2	7:C:505:PO4:P	2.35	0.47
1:A:121:ARG:O	1:A:159:LEU:HD13	2.15	0.47
1:D:14:ALA:O	1:D:20:LEU:HD23	2.14	0.47
1:D:266:ARG:O	1:D:385:ARG:HD2	2.15	0.47
1:D:359:ARG:HD3	1:D:363:GLU:OE2	2.14	0.47
1:B:65:LEU:HD13	1:B:216:ILE:HD12	1.96	0.47
1:A:73:ILE:HG23	1:A:109:TYR:HB2	1.96	0.46
1:C:451:ALA:HA	1:C:452:PRO:C	2.35	0.46
1:B:359:ARG:NH2	8:B:622:HOH:O	2.47	0.46
1:B:65:LEU:HD11	1:B:189:ALA:HB2	1.97	0.46
1:B:382:ARG:HH12	3:B:502:G6P:H1	1.81	0.46
1:C:266:ARG:O	1:C:385:ARG:HD2	2.16	0.46
1:A:263:GLN:HA	8:A:627:HOH:O	2.15	0.46
1:A:461:LEU:HD23	1:A:461:LEU:O	2.15	0.46
1:C:432:SER:O	1:C:436:LEU:HD13	2.16	0.45
1:C:85:TRP:CH2	1:C:91:VAL:HG11	2.51	0.45
1:C:431:LYS:O	1:C:435:GLU:HG3	2.17	0.45
1:D:11:LEU:HD13	1:D:47:ALA:HB1	1.97	0.45
1:C:65:LEU:HD11	1:C:189:ALA:HB2	1.98	0.45
1:A:80:SER:HB2	8:A:655:HOH:O	2.16	0.45
1:C:220:GLU:O	1:C:248:GLU:HG3	2.17	0.45
1:A:365:LEU:HD21	1:C:350:LYS:HA	1.98	0.45
1:B:121:ARG:O	1:B:159:LEU:HD13	2.17	0.45
1:D:452:PRO:HA	1:D:453:PRO:HD3	1.85	0.45
1:A:382:ARG:HH22	3:A:502:G6P:H1	1.82	0.45
1:D:279:LEU:HD23	1:D:292:GLU:HB3	1.99	0.45
1:B:45:LYS:NZ	8:B:628:HOH:O	2.49	0.44
1:D:216:ILE:HG12	1:D:237:ALA:HB3	1.99	0.44
1:D:461:LEU:HD12	1:D:461:LEU:O	2.17	0.44
1:B:340:ALA:O	1:B:341:PRO:C	2.55	0.44
1:B:347:PRO:O	1:B:348:ARG:NH1	2.50	0.44
1:D:340:ALA:O	1:D:341:PRO:C	2.56	0.44
1:A:50:ARG:HD2	8:A:615:HOH:O	2.18	0.44
1:C:386:LEU:C	1:C:387:HIS:O	2.56	0.44
1:A:147:GLU:OE1	1:B:287:ARG:NH2	2.50	0.44
8:A:628:HOH:O	1:B:147:GLU:HB3	2.16	0.44
1:D:451:ALA:HA	1:D:452:PRO:C	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:LEU:HB2	1:C:187:MET:SD	2.57	0.44
1:B:72:LYS:HE3	1:B:156:GLY:HA3	1.99	0.44
1:B:266:ARG:O	1:B:385:ARG:HD2	2.18	0.44
1:A:96:GLY:O	1:A:106:SER:OG	2.34	0.44
1:C:279:LEU:HD23	1:C:292:GLU:HB3	1.99	0.44
1:A:279:LEU:HD23	1:A:292:GLU:HB3	1.99	0.44
1:A:65:LEU:HD11	1:A:189:ALA:HB2	1.99	0.43
1:A:75:LEU:O	1:A:154:ASN:HA	2.18	0.43
1:D:65:LEU:HD11	1:D:189:ALA:HB2	1.99	0.43
1:A:186:ASP:C	1:A:187:MET:HG3	2.39	0.43
1:B:75:LEU:O	1:B:154:ASN:HA	2.17	0.43
1:B:256:LEU:HD23	1:B:256:LEU:HA	1.79	0.43
1:A:152:SER:OG	1:A:155:LYS:HE2	2.18	0.43
1:B:186:ASP:O	1:B:214:PRO:HD2	2.19	0.43
1:B:220:GLU:O	1:B:248:GLU:HG3	2.18	0.43
1:D:72:LYS:CG	1:D:74:ARG:HD2	2.42	0.43
1:A:85:TRP:CH2	1:A:91:VAL:HG11	2.53	0.43
1:B:350:LYS:HA	1:D:365:LEU:HD21	2.00	0.43
1:D:386:LEU:C	1:D:387:HIS:O	2.57	0.43
1:B:89:GLU:OE1	1:B:102:HIS:NE2	2.48	0.43
1:B:131:LEU:HD22	1:B:145:VAL:HG22	2.01	0.43
1:D:267:GLU:OE2	1:D:345:HIS:ND1	2.52	0.43
1:D:72:LYS:HB2	1:D:192:PHE:HE2	1.83	0.43
1:D:85:TRP:CH2	1:D:91:VAL:HG11	2.54	0.43
1:D:155:LYS:HA	1:D:155:LYS:CE	2.34	0.43
1:A:340:ALA:O	1:A:341:PRO:C	2.56	0.43
1:B:194:ARG:HH21	1:B:194:ARG:CG	2.30	0.43
1:C:251:LEU:HD12	1:C:251:LEU:C	2.39	0.43
1:D:75:LEU:O	1:D:154:ASN:HA	2.18	0.43
1:B:356:TYR:O	1:B:359:ARG:HG3	2.19	0.42
1:C:121:ARG:HG3	1:C:121:ARG:HH11	1.83	0.42
1:B:125:ASP:OD2	1:B:128:LYS:CD	2.67	0.42
1:C:45:LYS:HE3	1:C:49:GLU:OE1	2.19	0.42
1:D:186:ASP:HB2	1:D:187:MET:CE	2.49	0.42
1:A:275:ALA:O	1:A:276:THR:HB	2.19	0.42
1:A:186:ASP:O	1:A:214:PRO:HD2	2.20	0.42
1:C:193:VAL:O	1:C:224:ALA:HB2	2.20	0.42
1:B:441:ARG:NH2	1:B:441:ARG:H	2.16	0.42
1:C:340:ALA:O	1:C:341:PRO:C	2.55	0.42
1:D:89:GLU:OE1	1:D:102:HIS:NE2	2.52	0.42
1:C:71:PRO:HD2	1:C:192:PHE:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:GLU:OE1	1:C:102:HIS:NE2	2.53	0.42
1:C:329:ILE:HA	1:D:252:GLU:HB2	2.02	0.42
1:D:218:LYS:NZ	8:D:602:HOH:O	2.40	0.42
1:A:466:ARG:HB3	1:A:469:GLU:CG	2.36	0.41
1:D:134:ASP:OD2	1:D:144:THR:OG1	2.29	0.41
1:D:236:ASP:O	1:D:271:PRO:HD2	2.20	0.41
1:B:318:TYR:N	1:B:319:PRO:HD3	2.35	0.41
1:B:461:LEU:C	1:B:461:LEU:CD1	2.88	0.41
1:C:461:LEU:C	1:C:461:LEU:CD1	2.88	0.41
1:A:132:VAL:O	1:A:143:CYS:HB3	2.21	0.41
1:A:364:ARG:HG2	1:C:348:ARG:HH12	1.86	0.41
1:B:251:LEU:C	1:B:251:LEU:HD12	2.40	0.41
1:C:324:ARG:NH1	8:C:622:HOH:O	2.54	0.41
1:C:359:ARG:HD3	1:C:363:GLU:OE2	2.21	0.41
1:A:359:ARG:HD3	1:A:363:GLU:OE2	2.21	0.41
1:B:228:LEU:O	1:B:232:VAL:HG23	2.21	0.41
1:C:75:LEU:O	1:C:154:ASN:HA	2.20	0.41
1:C:295:ASP:HA	8:C:619:HOH:O	2.20	0.41
1:A:236:ASP:O	1:A:271:PRO:HD2	2.21	0.41
1:B:348:ARG:HA	1:B:348:ARG:NH1	2.33	0.41
1:B:275:ALA:O	1:B:276:THR:HB	2.20	0.41
1:A:181:LEU:HD23	1:A:181:LEU:HA	1.96	0.41
1:A:220:GLU:O	1:A:248:GLU:HG3	2.21	0.41
1:A:251:LEU:HD12	1:A:251:LEU:C	2.41	0.41
1:A:373:PHE:HD2	2:A:501:AMP:N7	2.18	0.41
1:D:72:LYS:NZ	8:D:603:HOH:O	2.19	0.41
1:D:324:ARG:NH1	8:D:609:HOH:O	2.34	0.41
1:A:146:VAL:HG12	1:A:147:GLU:HG3	2.03	0.41
1:C:382:ARG:HH12	3:C:502:G6P:C6	2.30	0.41
1:B:131:LEU:CD2	1:B:145:VAL:HG22	2.51	0.40
1:C:347:PRO:O	1:C:348:ARG:NH1	2.53	0.40
1:D:193:VAL:O	1:D:224:ALA:HB2	2.21	0.40
1:A:89:GLU:OE1	1:A:102:HIS:NE2	2.55	0.40
1:B:351:ARG:NH1	2:B:501:AMP:O3P	2.53	0.40
1:B:386:LEU:C	1:B:387:HIS:O	2.59	0.40
1:A:386:LEU:C	1:A:387:HIS:O	2.59	0.40
1:C:402:ARG:HG3	1:C:413:THR:HB	2.03	0.40
1:C:452:PRO:HA	1:C:453:PRO:HD3	1.78	0.40
1:A:228:LEU:O	1:A:232:VAL:HG23	2.21	0.40
1:B:276:THR:HA	8:B:609:HOH:O	2.21	0.40
1:B:452:PRO:HA	1:B:453:PRO:HD3	1.80	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:LEU:O	1:D:160:PRO:C	2.60	0.40

All (4) 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)
8:A:655:HOH:O	8:D:618:HOH:O[1_545]	1.56	0.64
1:A:45:LYS:NZ	1:B:99:GLU:OE2[2_445]	2.13	0.07
1:B:18:ASP:O	1:C:17:ARG:NH1[3_444]	2.15	0.05
1:B:19:ASP:OD2	1:C:17:ARG:NH1[3_444]	2.16	0.04

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	469/475 (99%)	461 (98%)	6 (1%)	2 (0%)	34	48
1	B	469/475 (99%)	459 (98%)	8 (2%)	2 (0%)	34	48
1	C	469/475 (99%)	459 (98%)	8 (2%)	2 (0%)	34	48
1	D	469/475 (99%)	455 (97%)	12 (3%)	2 (0%)	34	48
All	All	1876/1900 (99%)	1834 (98%)	34 (2%)	8 (0%)	34	48

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	387	HIS
1	B	387	HIS
1	C	387	HIS
1	D	387	HIS
1	A	276	THR
1	C	276	THR

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Mol	Chain	Res	Type
1	D	276	THR
1	B	276	THR

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	379/381 (100%)	367 (97%)	12 (3%)	39 59
1	B	379/381 (100%)	362 (96%)	17 (4%)	27 44
1	C	379/381 (100%)	362 (96%)	17 (4%)	27 44
1	D	379/381 (100%)	362 (96%)	17 (4%)	27 44
All	All	1516/1524 (100%)	1453 (96%)	63 (4%)	30 47

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	19	ASP
1	A	128	LYS
1	A	146	VAL
1	A	187	MET
1	A	194	ARG
1	A	337	SER
1	A	359	ARG
1	A	405	LEU
1	A	441	ARG
1	A	461	LEU
1	A	466	ARG
1	B	18	ASP
1	B	19	ASP
1	B	20	LEU
1	B	80	SER
1	B	114	GLN
1	B	131	LEU
1	B	155	LYS

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Mol	Chain	Res	Type
1	B	163	ASN
1	B	194	ARG
1	B	279	LEU
1	B	359	ARG
1	B	405	LEU
1	B	412	GLU
1	B	420	GLN
1	B	441	ARG
1	B	461	LEU
1	B	466	ARG
1	C	42	ASP
1	C	77	ARG
1	C	114	GLN
1	C	121	ARG
1	C	128	LYS
1	C	131	LEU
1	C	159	LEU
1	C	187	MET
1	C	194	ARG
1	C	359	ARG
1	C	368	LYS
1	C	400	GLU
1	C	405	LEU
1	C	412	GLU
1	C	438	ARG
1	C	461	LEU
1	C	466	ARG
1	D	18	ASP
1	D	45	LYS
1	D	74	ARG
1	D	106	SER
1	D	131	LEU
1	D	155	LYS
1	D	163	ASN
1	D	204	GLU
1	D	319	PRO
1	D	348	ARG
1	D	359	ARG
1	D	405	LEU
1	D	438	ARG
1	D	441	ARG
1	D	444	LEU

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Mol	Chain	Res	Type
1	D	461	LEU
1	D	466	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	404	GLN
1	B	227	ASN
1	B	404	GLN
1	C	163	ASN
1	C	227	ASN
1	C	404	GLN
1	D	227	ASN
1	D	404	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 5 are monoatomic - leaving 14 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	AMP	A	501	-	22,25,25	0.98	1 (4%)	25,38,38	1.47	4 (16%)
2	AMP	B	501	-	22,25,25	0.85	0	25,38,38	1.38	4 (16%)
3	G6P	A	502	-	16,16,16	0.73	0	24,24,24	1.87	4 (16%)
5	OXL	A	504	4	5,5,5	2.05	1 (20%)	6,6,6	1.31	0
3	G6P	D	502	-	16,16,16	0.73	0	24,24,24	1.68	5 (20%)
2	AMP	C	501	-	22,25,25	0.80	0	25,38,38	1.41	4 (16%)
5	OXL	C	504	4	5,5,5	1.57	1 (20%)	6,6,6	1.47	2 (33%)
7	PO4	D	505	-	4,4,4	0.64	0	6,6,6	0.67	0
2	AMP	D	501	-	22,25,25	0.91	1 (4%)	25,38,38	1.63	6 (24%)
5	OXL	B	504	4	5,5,5	2.12	1 (20%)	6,6,6	1.05	0
5	OXL	D	504	4	5,5,5	1.71	1 (20%)	6,6,6	1.37	1 (16%)
3	G6P	B	502	-	16,16,16	0.85	0	24,24,24	1.67	4 (16%)
7	PO4	C	505	-	4,4,4	0.86	0	6,6,6	0.66	0
3	G6P	C	502	-	16,16,16	0.59	0	24,24,24	1.09	1 (4%)

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	AMP	A	501	-	-	0/6/26/26	0/3/3/3
2	AMP	B	501	-	-	0/6/26/26	0/3/3/3
3	G6P	A	502	-	-	1/6/26/26	0/1/1/1
5	OXL	A	504	4	-	0/4/4/4	-
3	G6P	D	502	-	-	0/6/26/26	0/1/1/1
2	AMP	C	501	-	-	0/6/26/26	0/3/3/3
5	OXL	C	504	4	-	1/4/4/4	-
2	AMP	D	501	-	-	0/6/26/26	0/3/3/3
5	OXL	B	504	4	-	1/4/4/4	-
5	OXL	D	504	4	-	0/4/4/4	-
3	G6P	B	502	-	-	0/6/26/26	0/1/1/1
3	G6P	C	502	-	-	2/6/26/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	504	OXL	C2-C1	-3.73	1.43	1.54
5	B	504	OXL	C2-C1	-3.61	1.44	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	504	OXL	C2-C1	-2.67	1.46	1.54
5	C	504	OXL	C2-C1	-2.31	1.47	1.54
2	A	501	AMP	O4'-C1'	2.26	1.44	1.41
2	D	501	AMP	C5-N7	-2.10	1.32	1.39

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	G6P	O5-C1-C2	4.99	119.19	110.28
3	D	502	G6P	O5-C1-C2	4.75	118.76	110.28
2	D	501	AMP	N3-C2-N1	-4.65	121.41	128.68
3	A	502	G6P	C1-O5-C5	4.56	122.27	113.66
3	B	502	G6P	O5-C1-C2	4.26	117.89	110.28
2	B	501	AMP	N3-C2-N1	-4.09	122.28	128.68
2	C	501	AMP	N3-C2-N1	-4.09	122.29	128.68
2	A	501	AMP	N3-C2-N1	-4.05	122.36	128.68
3	D	502	G6P	C1-O5-C5	3.96	121.13	113.66
3	B	502	G6P	C1-O5-C5	3.75	120.74	113.66
2	A	501	AMP	C1'-N9-C4	-3.68	120.18	126.64
2	C	501	AMP	C3'-C2'-C1'	3.19	105.78	100.98
3	B	502	G6P	O3-C3-C2	-2.98	103.47	110.35
3	A	502	G6P	O5-C5-C6	2.74	112.19	106.67
2	B	501	AMP	N6-C6-N1	2.67	124.12	118.57
2	D	501	AMP	C3'-C2'-C1'	2.66	104.98	100.98
2	D	501	AMP	N6-C6-N1	2.63	124.03	118.57
2	B	501	AMP	C1'-N9-C4	-2.58	122.12	126.64
2	D	501	AMP	O2'-C2'-C3'	-2.53	103.64	111.82
2	A	501	AMP	N6-C6-N1	2.47	123.70	118.57
5	D	504	OXL	O3-C1-C2	2.47	120.49	113.16
3	D	502	G6P	C3-C4-C5	-2.46	105.86	110.24
2	C	501	AMP	C2-N1-C6	2.45	122.95	118.75
2	A	501	AMP	C3'-C2'-C1'	2.39	104.58	100.98
3	B	502	G6P	C1-C2-C3	2.34	115.17	110.31
3	C	502	G6P	O1-C1-C2	2.27	115.43	109.03
3	D	502	G6P	O2-C2-C1	2.22	114.30	109.16
5	C	504	OXL	O4-C2-C1	2.18	119.64	113.16
5	C	504	OXL	O3-C1-C2	2.17	119.61	113.16
2	D	501	AMP	O3P-P-O2P	2.17	115.91	107.64
2	C	501	AMP	C1'-N9-C4	-2.14	122.88	126.64
3	D	502	G6P	C4-C3-C2	-2.04	107.26	110.82
2	B	501	AMP	C5-C6-N6	-2.02	117.29	120.35
2	D	501	AMP	C5-C6-N6	-2.01	117.29	120.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	G6P	O5-C5-C4	2.00	113.33	109.69

There are no chirality outliers.

All (5) torsion outliers are listed below:

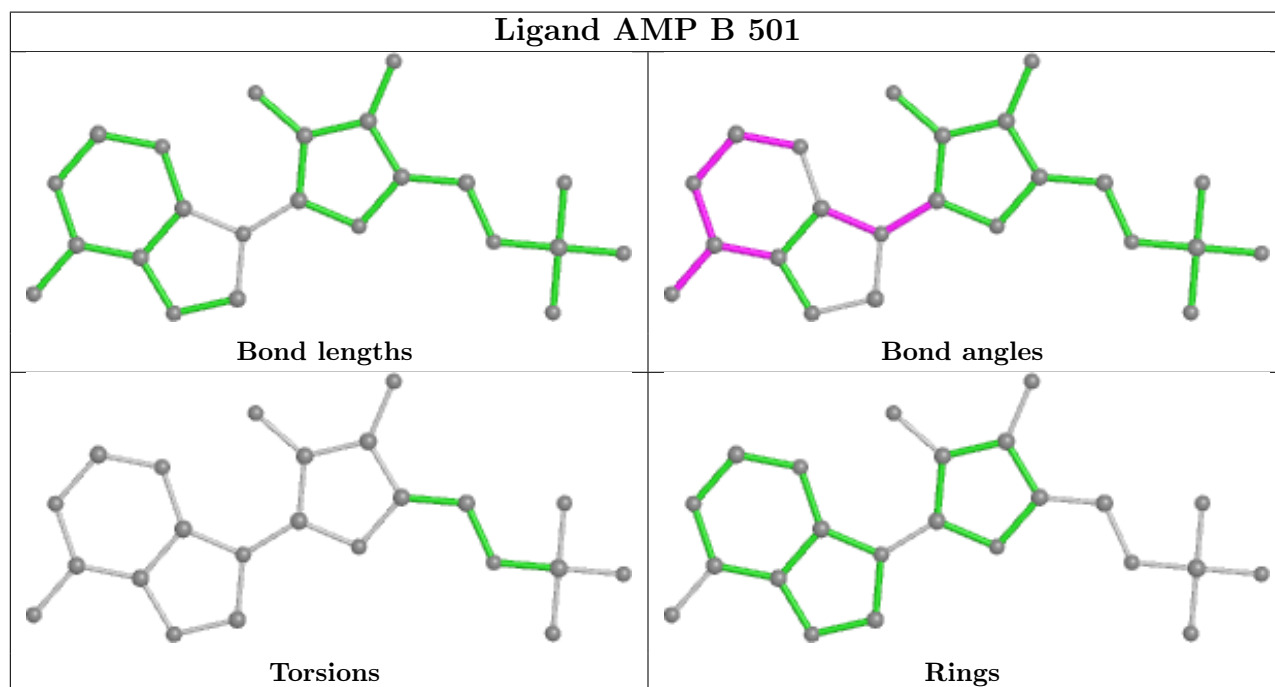
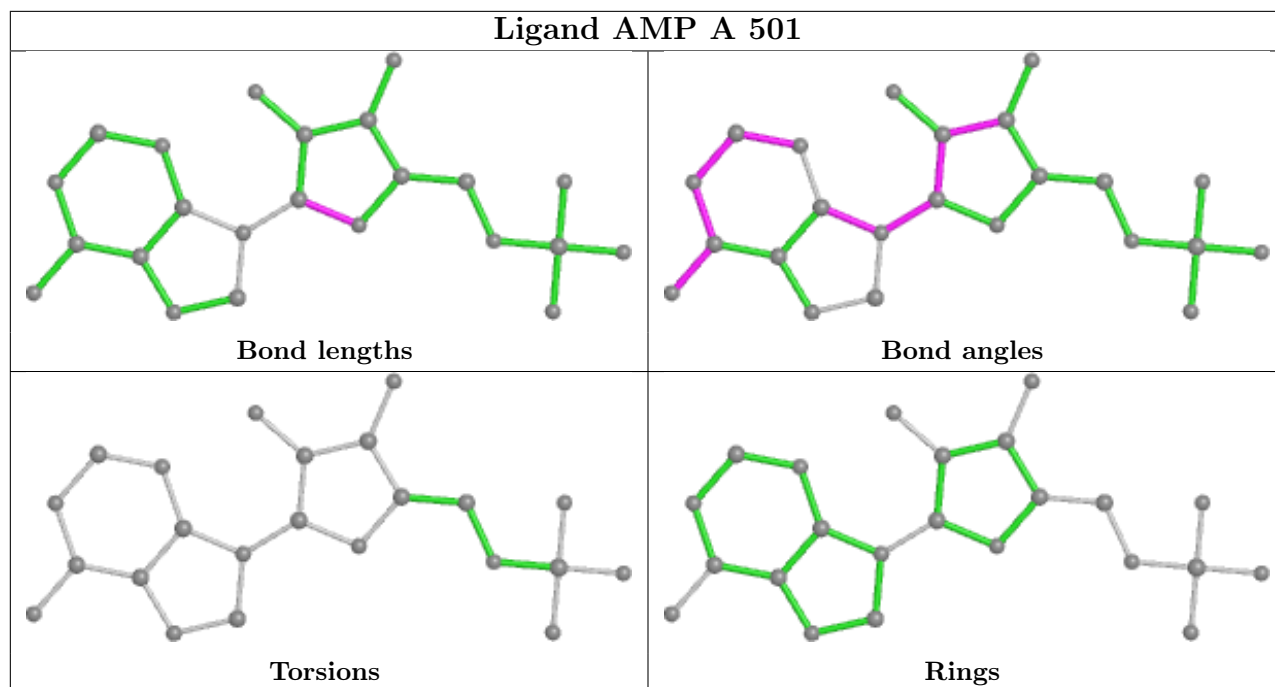
Mol	Chain	Res	Type	Atoms
3	A	502	G6P	C6-O6-P-O1P
3	C	502	G6P	C6-O6-P-O3P
3	C	502	G6P	C6-O6-P-O2P
5	B	504	OXL	O3-C1-C2-O4
5	C	504	OXL	O3-C1-C2-O4

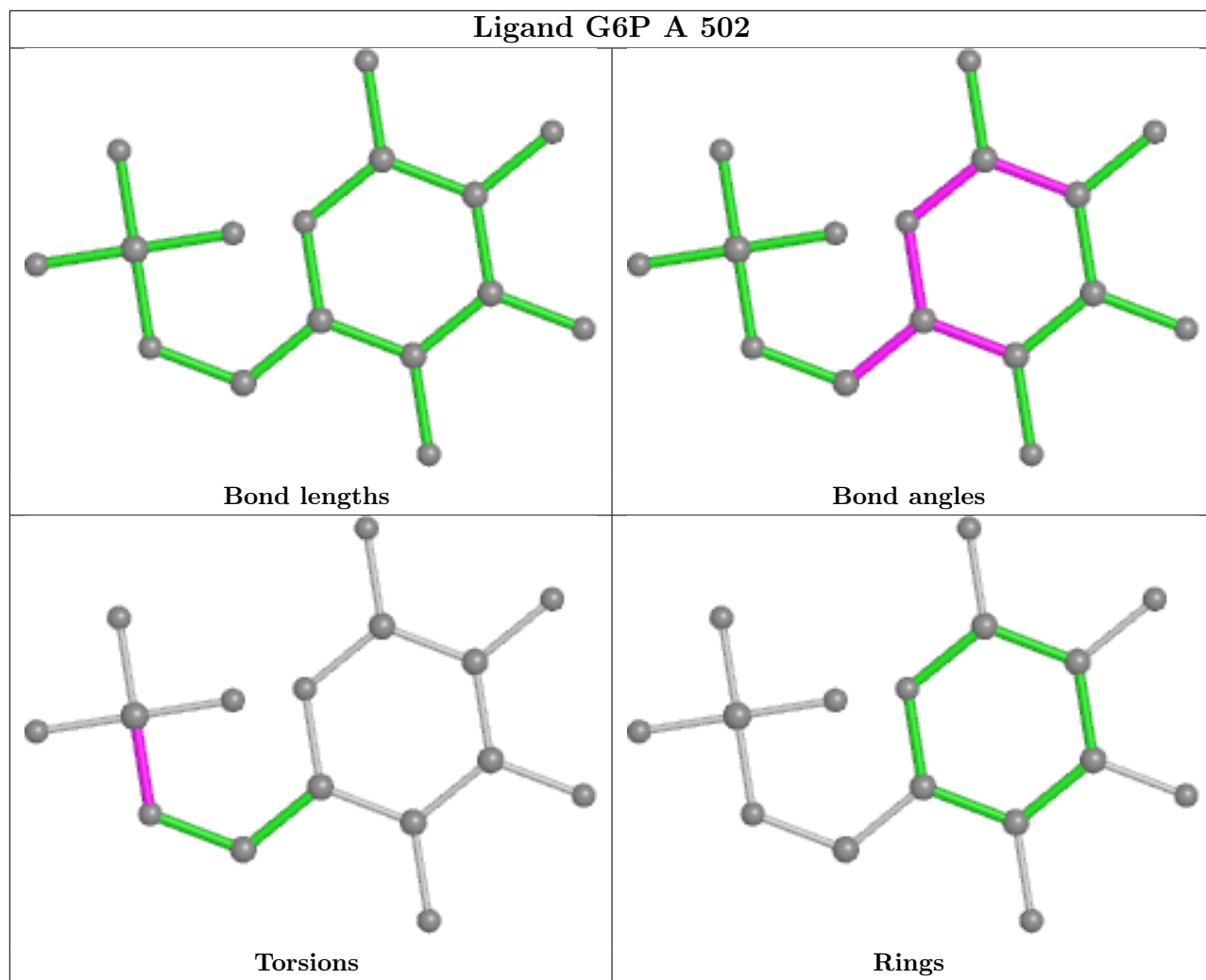
There are no ring outliers.

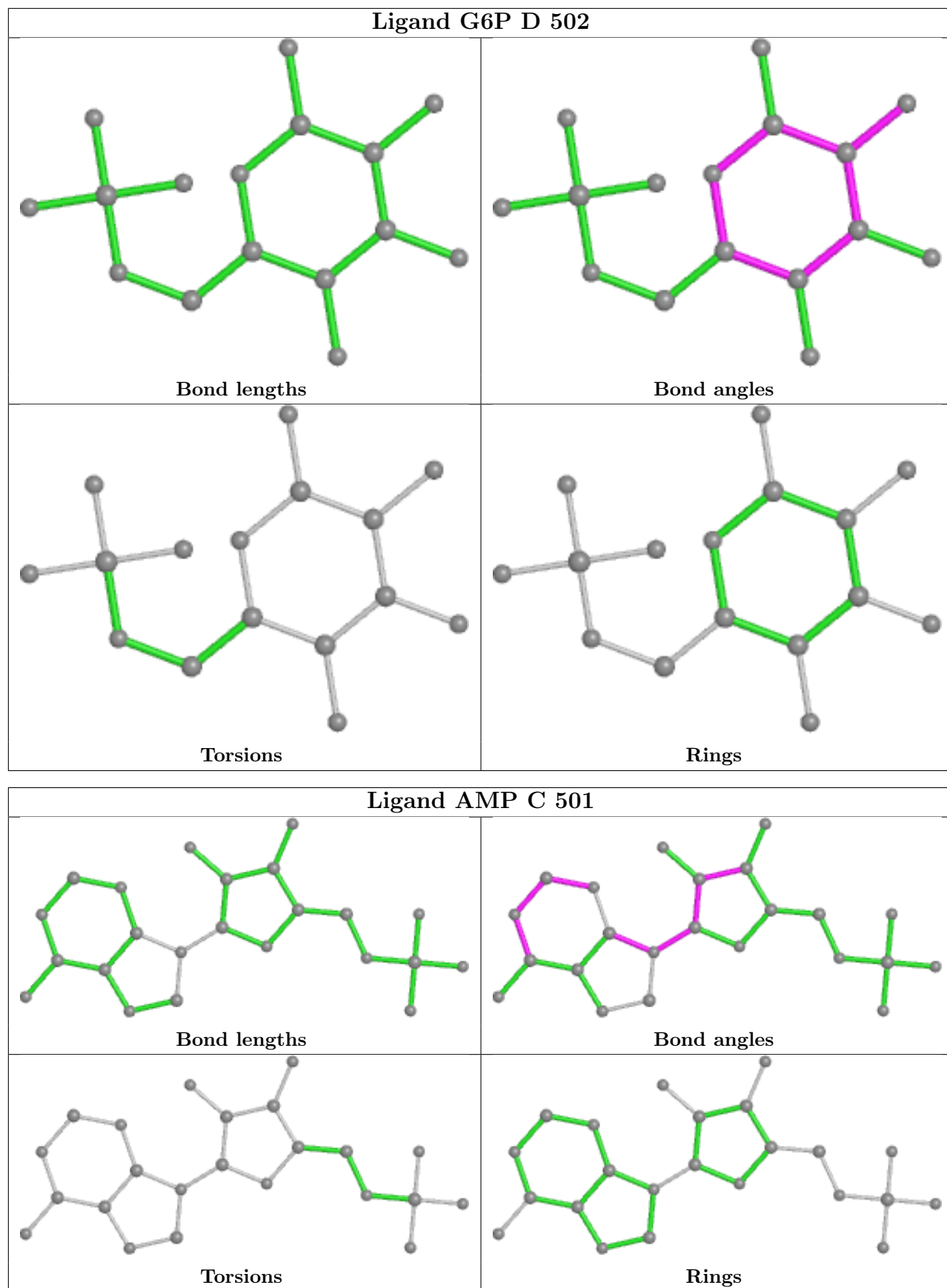
10 monomers are involved in 21 short contacts:

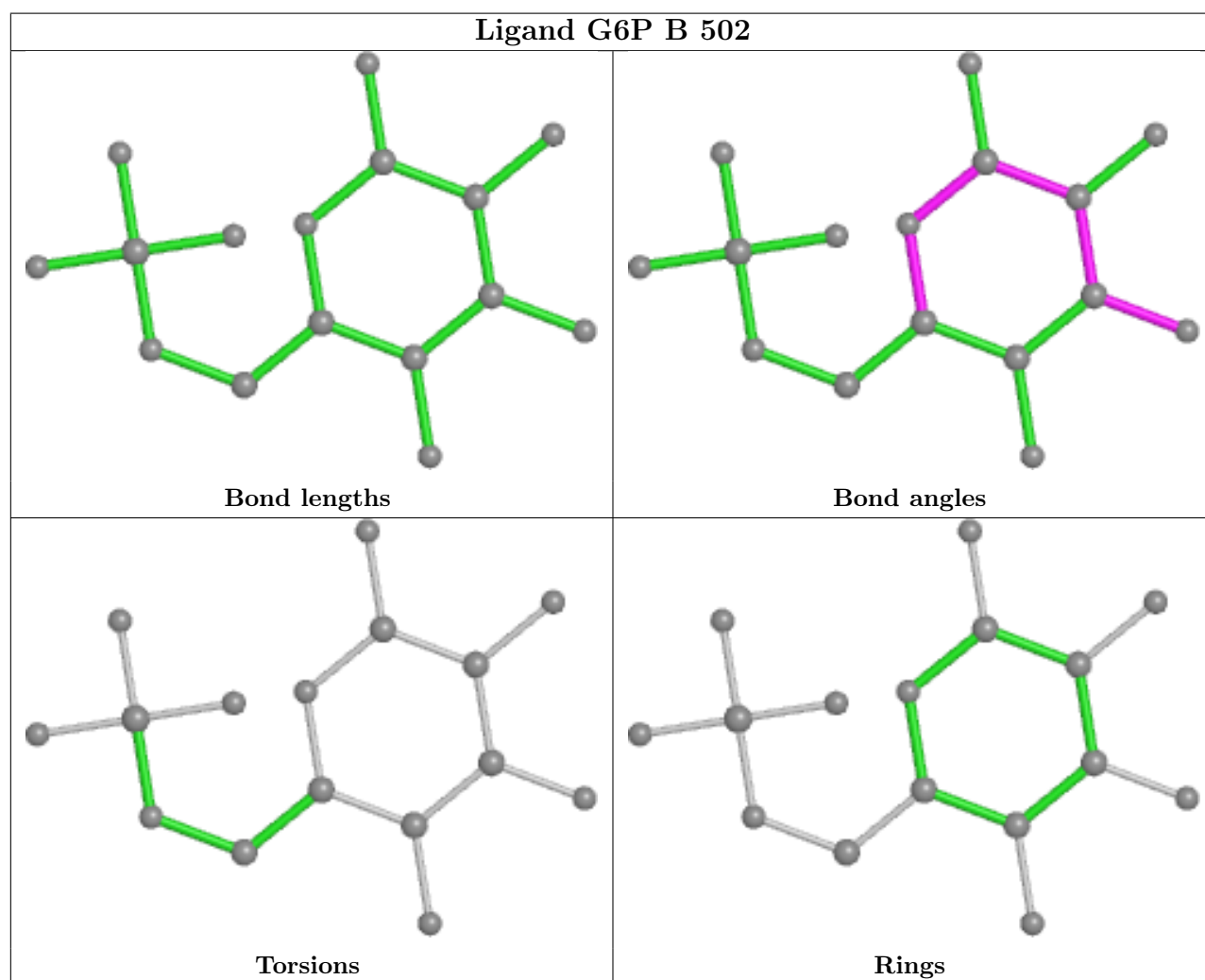
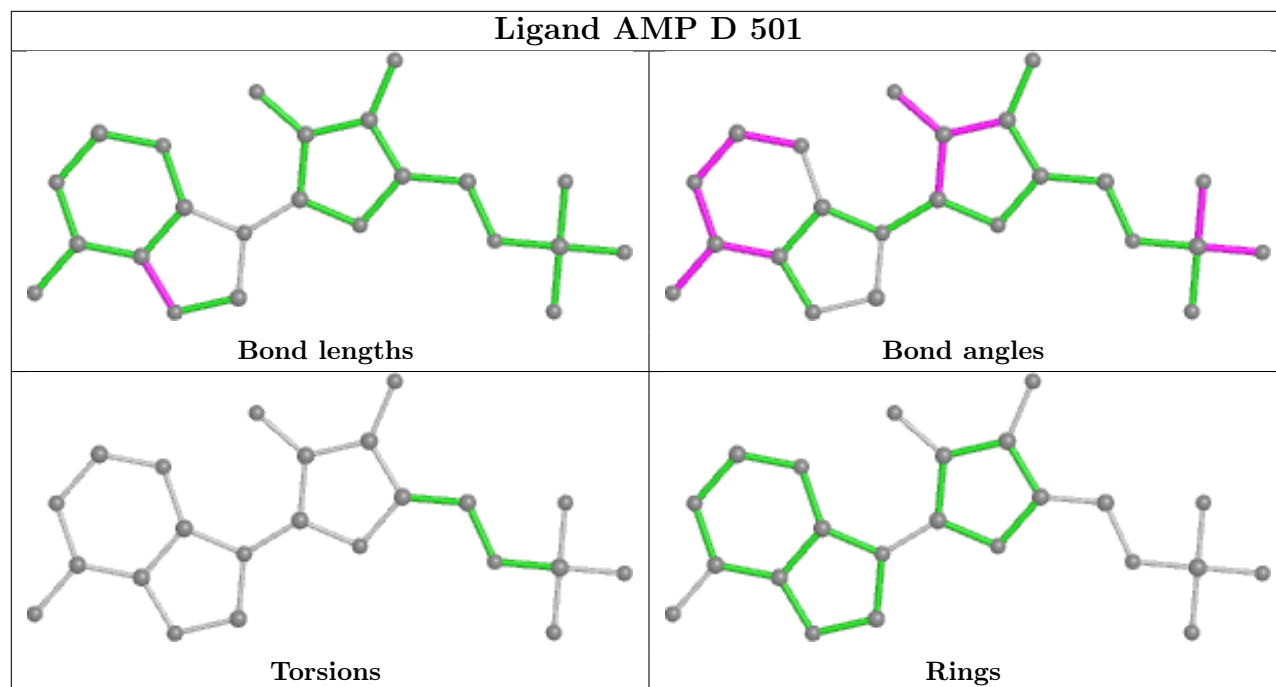
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	AMP	2	0
2	B	501	AMP	2	0
3	A	502	G6P	1	0
3	D	502	G6P	1	0
2	C	501	AMP	1	0
2	D	501	AMP	1	0
5	D	504	OXL	1	0
3	B	502	G6P	2	0
7	C	505	PO4	7	0
3	C	502	G6P	3	0

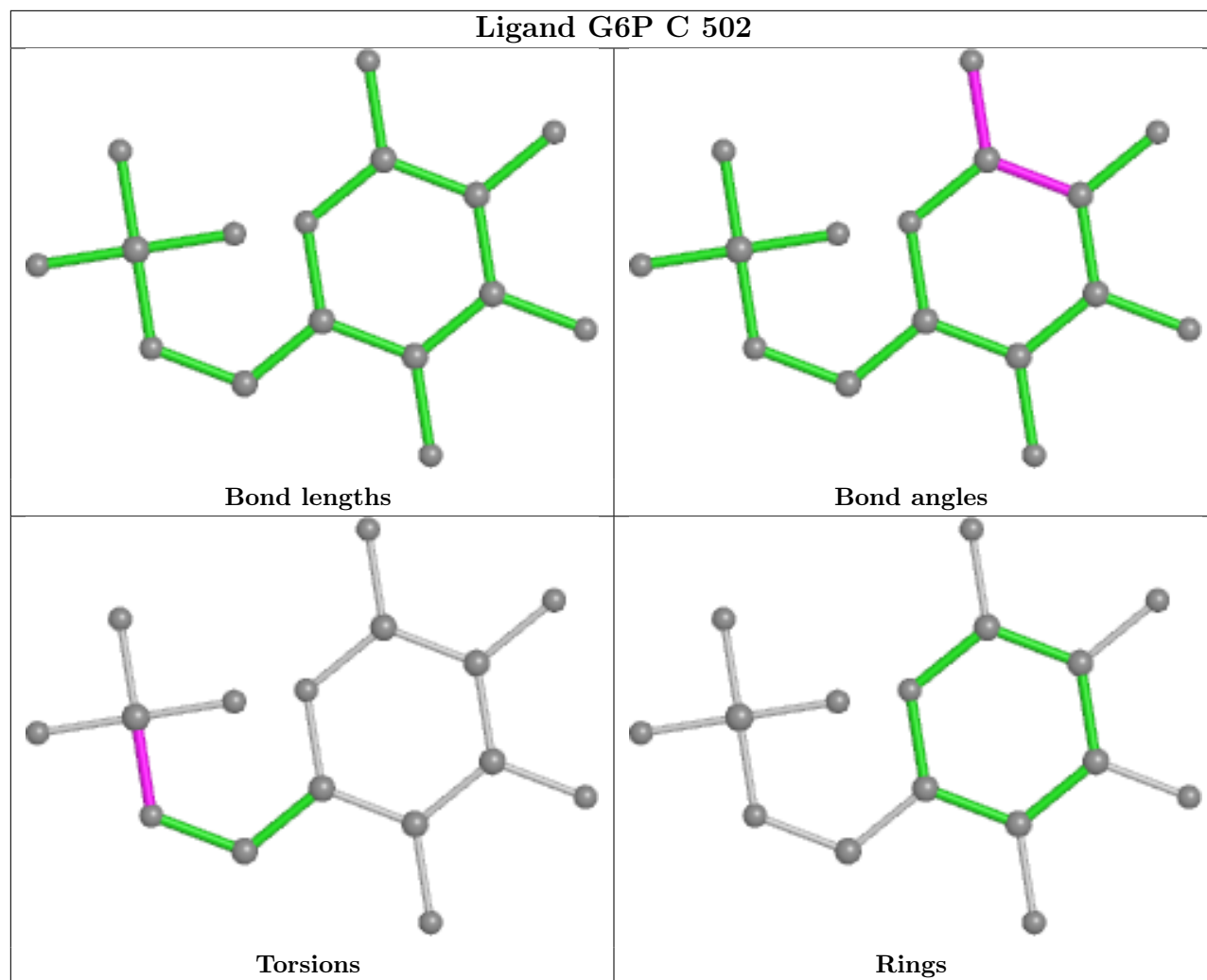
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	471/475 (99%)	-0.70	0 100 100	21, 27, 34, 42	0
1	B	471/475 (99%)	-0.70	0 100 100	21, 27, 34, 40	0
1	C	471/475 (99%)	-0.67	0 100 100	22, 28, 36, 42	0
1	D	471/475 (99%)	-0.68	0 100 100	21, 28, 36, 44	0
All	All	1884/1900 (99%)	-0.69	0 100 100	21, 27, 35, 44	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 monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

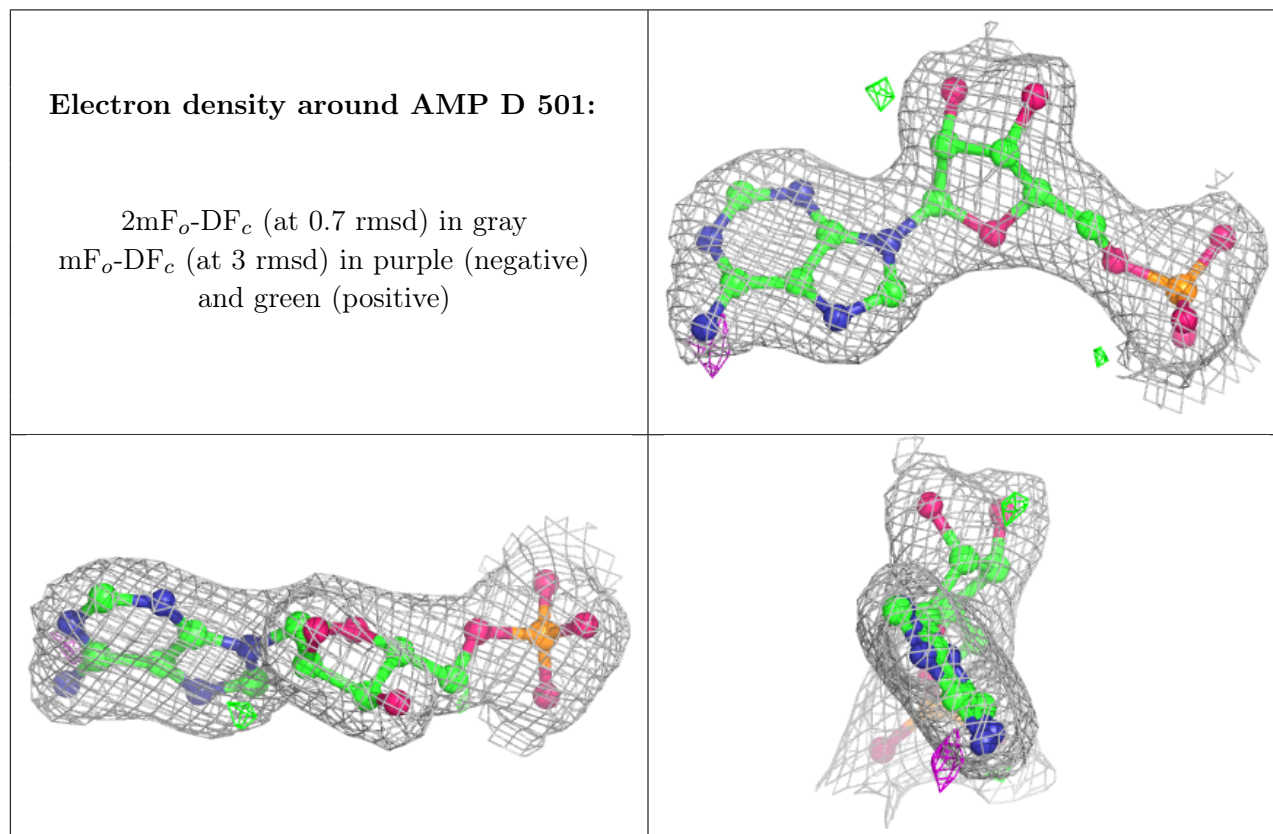
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	B	503	1/1	0.98	0.07	23,23,23,23	0
4	MG	D	503	1/1	0.98	0.05	21,21,21,21	0
7	PO4	C	505	5/5	0.98	0.06	24,24,29,31	0
2	AMP	D	501	23/23	0.99	0.10	21,24,27,28	0

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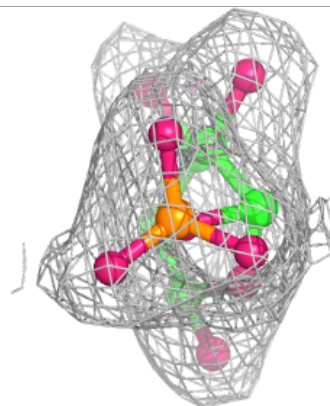
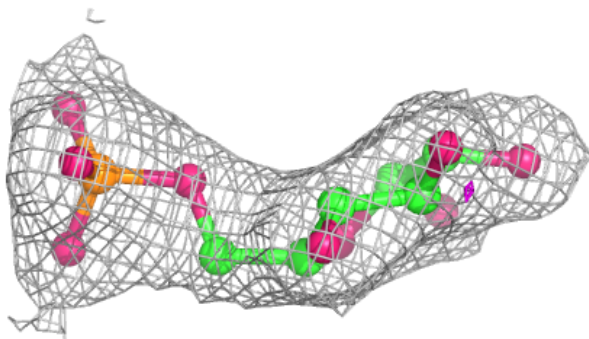
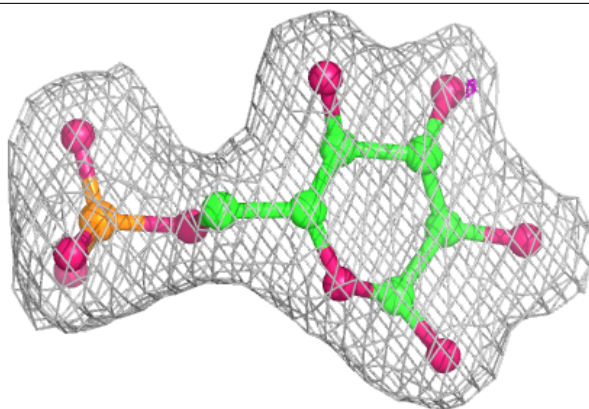
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	G6P	A	502	16/16	0.99	0.12	20,22,24,24	0
3	G6P	B	502	16/16	0.99	0.11	19,21,25,25	0
3	G6P	C	502	16/16	0.99	0.10	18,21,23,24	0
3	G6P	D	502	16/16	0.99	0.10	19,22,24,24	0
4	MG	A	503	1/1	0.99	0.06	22,22,22,22	0
2	AMP	A	501	23/23	0.99	0.10	18,23,25,27	0
4	MG	C	503	1/1	0.99	0.06	19,19,19,19	0
2	AMP	B	501	23/23	0.99	0.10	18,23,26,26	0
5	OXL	A	504	6/6	0.99	0.12	18,20,21,23	0
5	OXL	B	504	6/6	0.99	0.11	18,20,21,21	0
5	OXL	C	504	6/6	0.99	0.10	19,20,23,24	0
5	OXL	D	504	6/6	0.99	0.09	22,22,23,23	0
2	AMP	C	501	23/23	0.99	0.10	21,25,27,32	0
7	PO4	D	505	5/5	0.99	0.06	23,27,28,29	0
6	K	B	505	1/1	1.00	0.10	23,23,23,23	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.

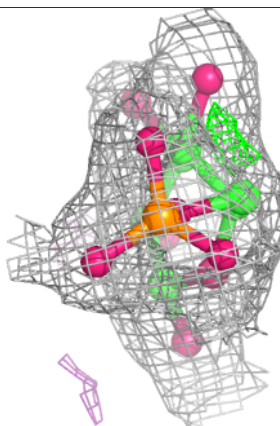
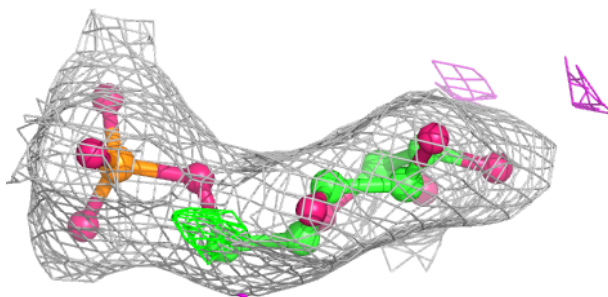
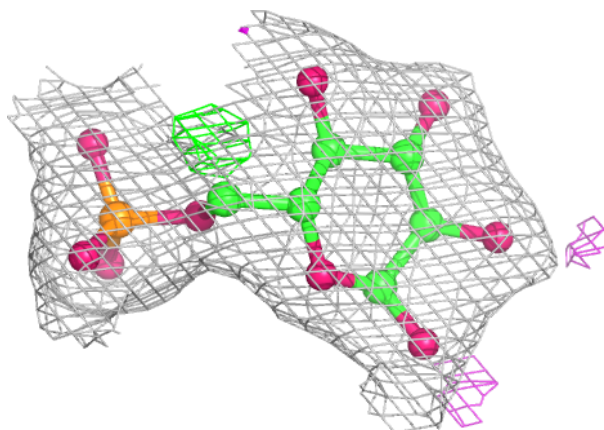


Electron density around G6P A 502:

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

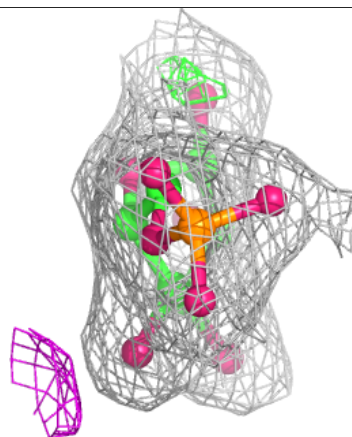
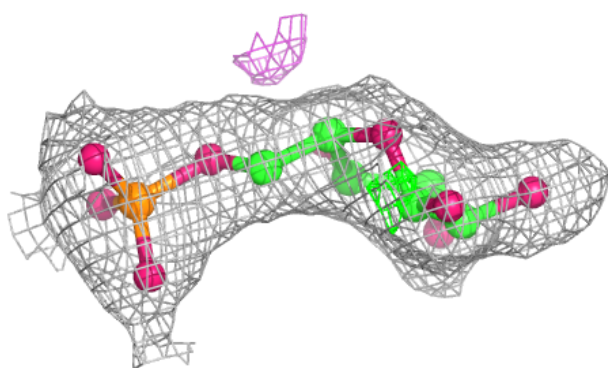
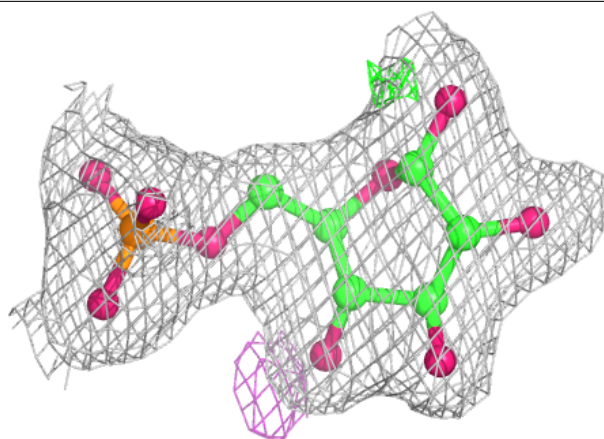
**Electron density around G6P B 502:**

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



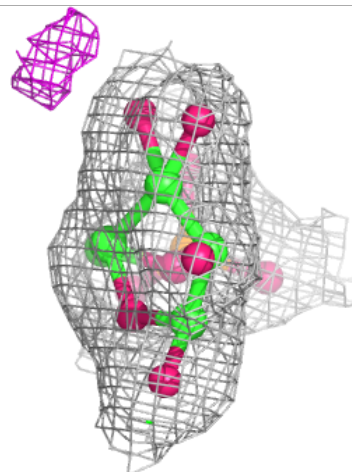
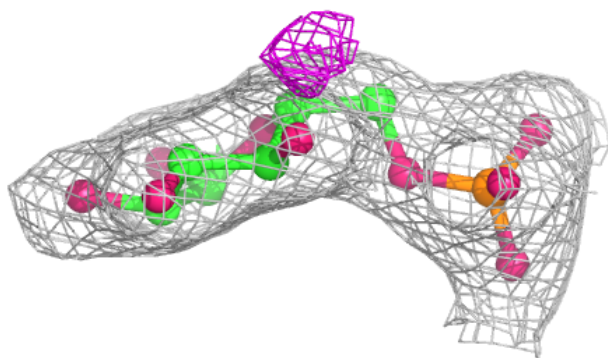
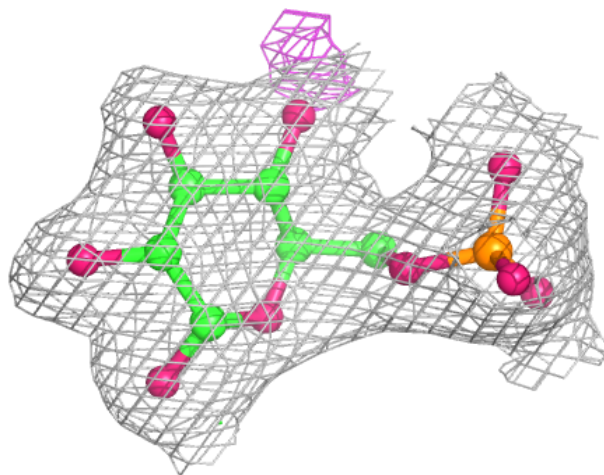
Electron density around G6P C 502:

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



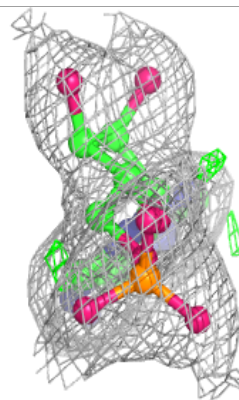
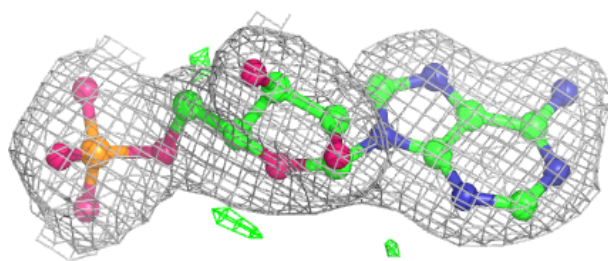
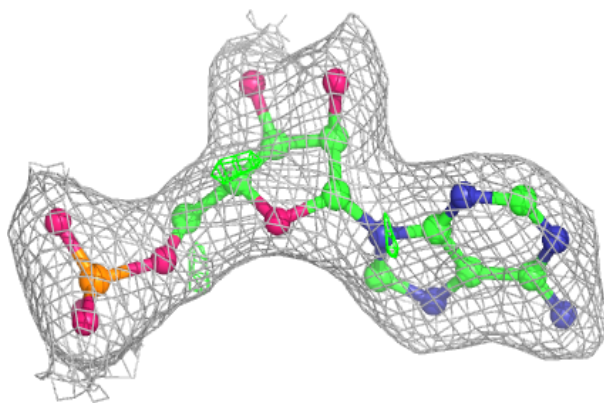
Electron density around G6P D 502:

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

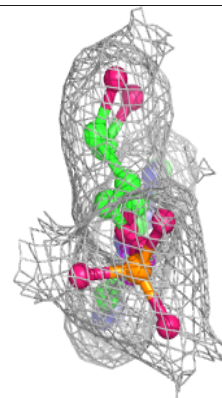
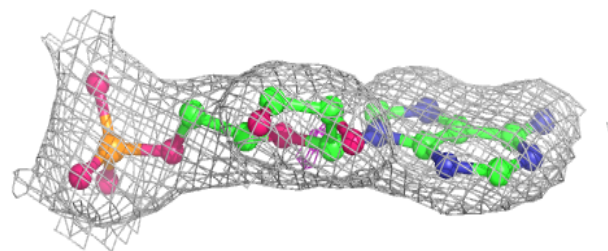
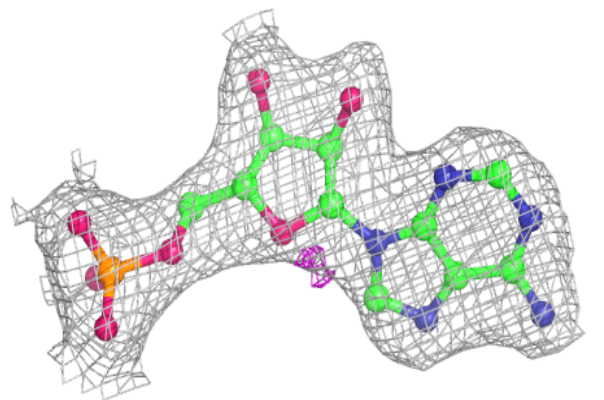


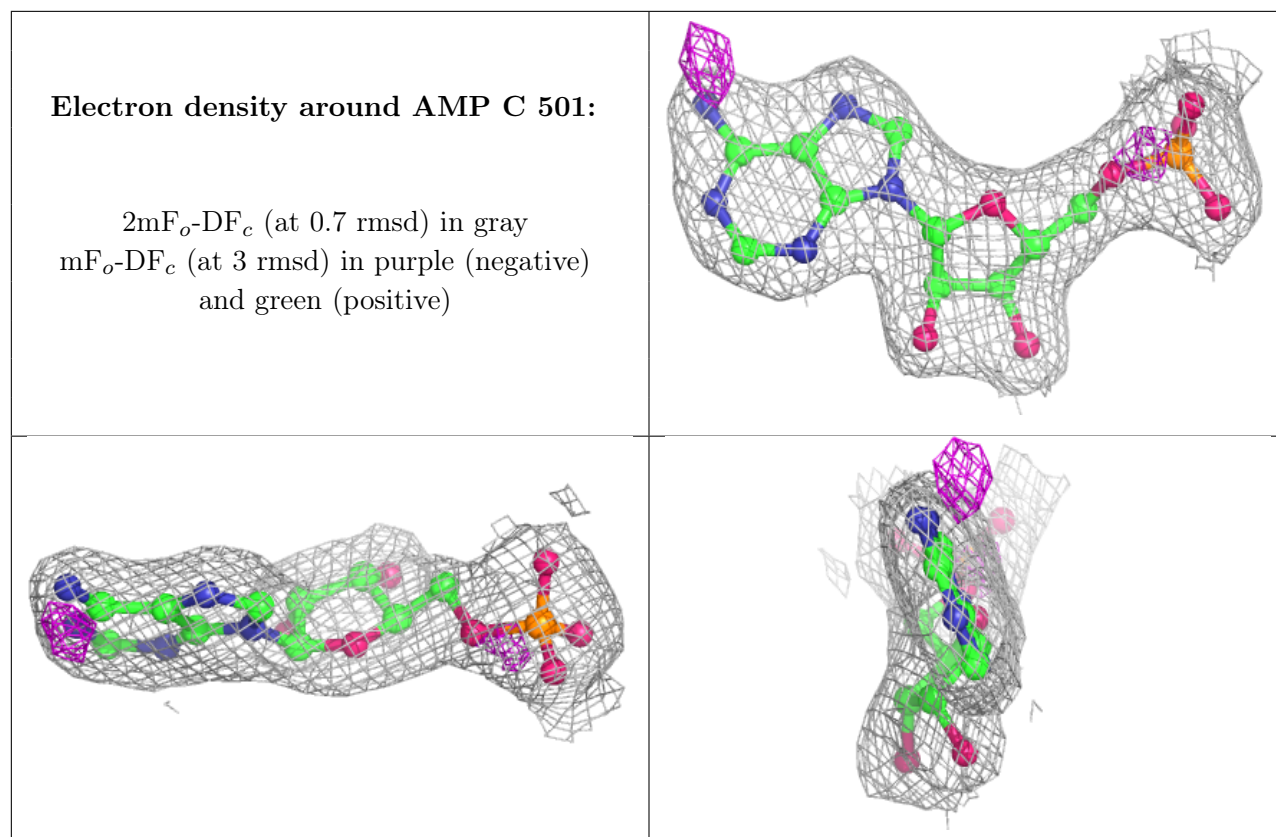
Electron density around AMP A 501:

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

**Electron density around AMP 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)





6.5 Other polymers [i](#)

There are no such residues in this entry.