



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 12:38 AM BST

PDB ID : 5XSS
Title : XylFII molecule
Authors : Li, J.X.; Wang, C.Y.; Zhang, P.
Deposited on : 2017-06-15
Resolution : 2.09 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 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 : 2.13.1
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) : 2.13.1

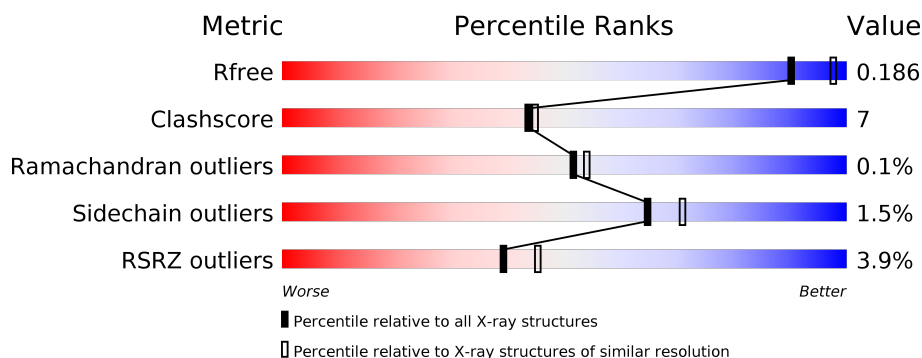
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on 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	316	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	316	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>•</div> <div>14%</div> </div> </div>
1	C	316	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div></div> <div>15%</div> </div> </div>
1	X	316	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>7%</div> <div>•</div> <div>14%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9151 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 Periplasmic binding protein/LacI transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	271	Total	C	N	O	S	0	0	0
			2068	1289	357	413	9			
1	A	274	Total	C	N	O	S	0	0	0
			2090	1305	360	416	9			
1	B	272	Total	C	N	O	S	0	0	0
			2075	1296	358	412	9			
1	C	270	Total	C	N	O	S	0	0	0
			2062	1286	356	411	9			

There are 56 discrepancies between the modelled and reference sequences:

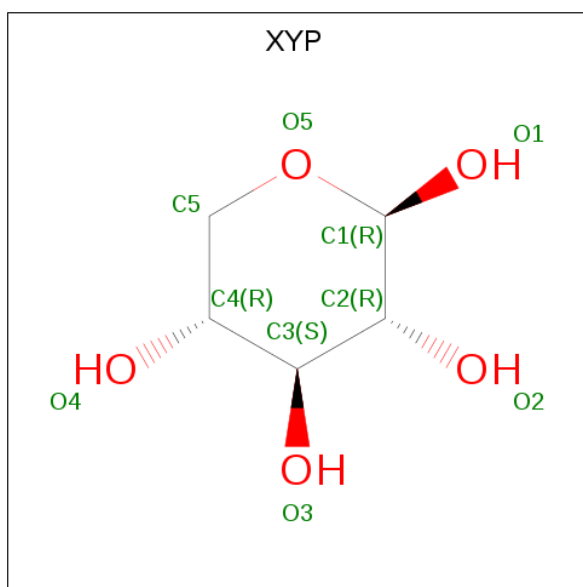
Chain	Residue	Modelled	Actual	Comment	Reference
X	-13	MET	-	expression tag	UNP A6LW07
X	-12	GLY	-	expression tag	UNP A6LW07
X	-11	SER	-	expression tag	UNP A6LW07
X	-10	SER	-	expression tag	UNP A6LW07
X	-9	HIS	-	expression tag	UNP A6LW07
X	-8	HIS	-	expression tag	UNP A6LW07
X	-7	HIS	-	expression tag	UNP A6LW07
X	-6	HIS	-	expression tag	UNP A6LW07
X	-5	HIS	-	expression tag	UNP A6LW07
X	-4	HIS	-	expression tag	UNP A6LW07
X	-3	SER	-	expression tag	UNP A6LW07
X	-2	GLN	-	expression tag	UNP A6LW07
X	-1	GLY	-	expression tag	UNP A6LW07
X	0	SER	-	expression tag	UNP A6LW07
A	-13	MET	-	expression tag	UNP A6LW07
A	-12	GLY	-	expression tag	UNP A6LW07
A	-11	SER	-	expression tag	UNP A6LW07
A	-10	SER	-	expression tag	UNP A6LW07
A	-9	HIS	-	expression tag	UNP A6LW07
A	-8	HIS	-	expression tag	UNP A6LW07
A	-7	HIS	-	expression tag	UNP A6LW07

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	HIS	-	expression tag	UNP A6LW07
A	-5	HIS	-	expression tag	UNP A6LW07
A	-4	HIS	-	expression tag	UNP A6LW07
A	-3	SER	-	expression tag	UNP A6LW07
A	-2	GLN	-	expression tag	UNP A6LW07
A	-1	GLY	-	expression tag	UNP A6LW07
A	0	SER	-	expression tag	UNP A6LW07
B	-13	MET	-	expression tag	UNP A6LW07
B	-12	GLY	-	expression tag	UNP A6LW07
B	-11	SER	-	expression tag	UNP A6LW07
B	-10	SER	-	expression tag	UNP A6LW07
B	-9	HIS	-	expression tag	UNP A6LW07
B	-8	HIS	-	expression tag	UNP A6LW07
B	-7	HIS	-	expression tag	UNP A6LW07
B	-6	HIS	-	expression tag	UNP A6LW07
B	-5	HIS	-	expression tag	UNP A6LW07
B	-4	HIS	-	expression tag	UNP A6LW07
B	-3	SER	-	expression tag	UNP A6LW07
B	-2	GLN	-	expression tag	UNP A6LW07
B	-1	GLY	-	expression tag	UNP A6LW07
B	0	SER	-	expression tag	UNP A6LW07
C	-13	MET	-	expression tag	UNP A6LW07
C	-12	GLY	-	expression tag	UNP A6LW07
C	-11	SER	-	expression tag	UNP A6LW07
C	-10	SER	-	expression tag	UNP A6LW07
C	-9	HIS	-	expression tag	UNP A6LW07
C	-8	HIS	-	expression tag	UNP A6LW07
C	-7	HIS	-	expression tag	UNP A6LW07
C	-6	HIS	-	expression tag	UNP A6LW07
C	-5	HIS	-	expression tag	UNP A6LW07
C	-4	HIS	-	expression tag	UNP A6LW07
C	-3	SER	-	expression tag	UNP A6LW07
C	-2	GLN	-	expression tag	UNP A6LW07
C	-1	GLY	-	expression tag	UNP A6LW07
C	0	SER	-	expression tag	UNP A6LW07

- Molecule 2 is beta-D-xylopyranose (three-letter code: XYP) (formula: C₅H₁₀O₅) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	X	1	Total	C	O	0	0
			10	5	5		
2	A	1	Total	C	O	0	0
			10	5	5		
2	B	1	Total	C	O	0	0
			10	5	5		
2	C	1	Total	C	O	0	0
			10	5	5		

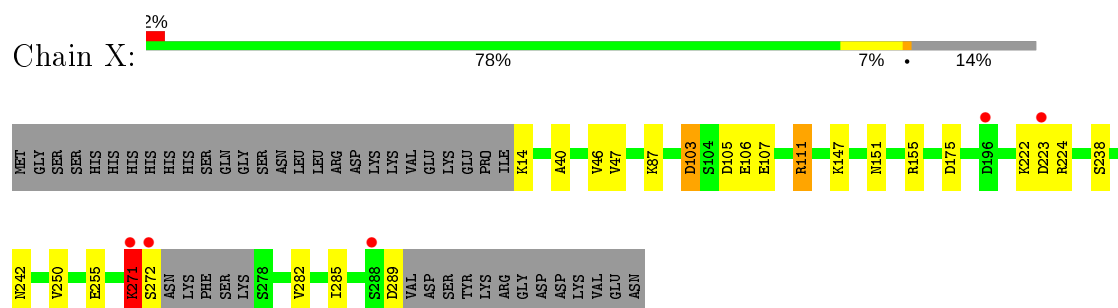
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	205	Total	O	0	0
			205	205		
3	A	196	Total	O	0	0
			196	196		
3	B	205	Total	O	0	0
			205	205		
3	C	210	Total	O	0	0
			210	210		

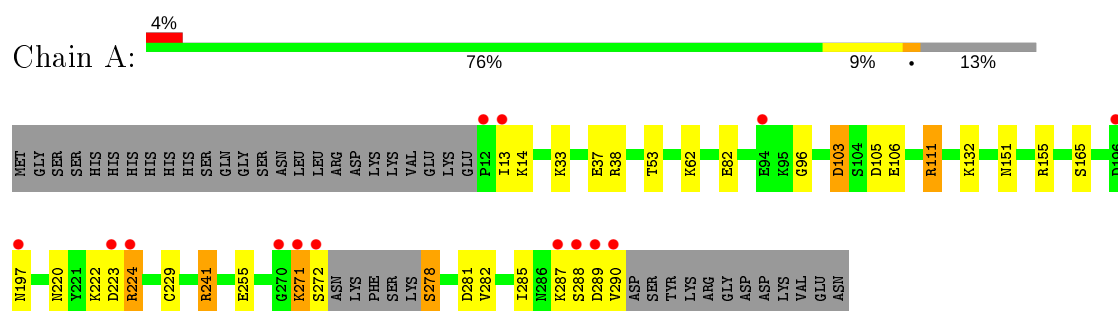
3 Residue-property plots [i](#)

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.

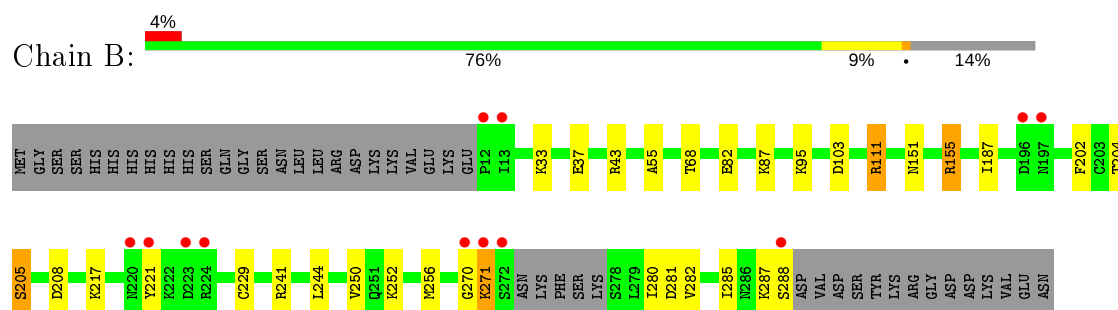
- Molecule 1: Periplasmic binding protein/LacI transcriptional regulator



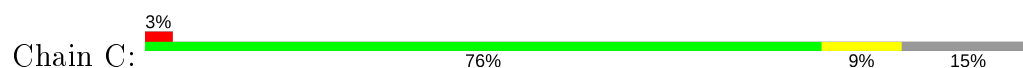
- Molecule 1: Periplasmic binding protein/LacI transcriptional regulator

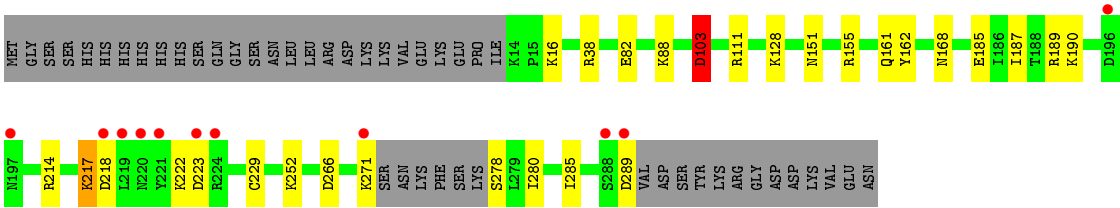


- Molecule 1: Periplasmic binding protein/LacI transcriptional regulator



- Molecule 1: Periplasmic binding protein/LacI transcriptional regulator





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	67.54Å 68.06Å 76.68Å 89.89° 65.15° 87.33°	Depositor
Resolution (Å)	30.78 – 2.09 30.78 – 2.09	Depositor EDS
% Data completeness (in resolution range)	96.8 (30.78-2.09) 96.8 (30.78-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.180 , 0.214 0.183 , 0.186	Depositor DCC
R_{free} test set	3549 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	17.2	Xtriage
Anisotropy	0.604	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for h,-k,h-l 0.066 for -h,k,-l 0.002 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9151	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% 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: XYP

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	0.48	0/2108	0.72	5/2838 (0.2%)
1	B	0.57	1/2093 (0.0%)	0.74	5/2817 (0.2%)
1	C	0.56	0/2079	0.66	1/2798 (0.0%)
1	X	0.55	1/2085 (0.0%)	0.71	5/2806 (0.2%)
All	All	0.54	2/8365 (0.0%)	0.71	16/11259 (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	X	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	205	SER	CB-OG	-6.54	1.33	1.42
1	X	238	SER	CB-OG	-5.38	1.35	1.42

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	38	ARG	NE-CZ-NH2	-14.00	113.30	120.30
1	X	271	LYS	CB-CA-C	-12.88	84.64	110.40
1	B	155	ARG	NE-CZ-NH1	12.34	126.47	120.30
1	A	38	ARG	NE-CZ-NH1	10.45	125.53	120.30
1	B	43	ARG	NE-CZ-NH2	-10.10	115.25	120.30
1	B	155	ARG	NE-CZ-NH2	-9.93	115.33	120.30
1	X	271	LYS	N-CA-C	9.58	136.87	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	CB-CA-C	-7.51	95.38	110.40
1	B	43	ARG	NE-CZ-NH1	6.92	123.76	120.30
1	A	271	LYS	CB-CA-C	-6.46	97.48	110.40
1	X	103	ASP	CB-CA-C	-6.40	97.60	110.40
1	X	223	ASP	CB-CG-OD1	5.93	123.63	118.30
1	C	103	ASP	CB-CA-C	-5.89	98.63	110.40
1	X	224	ARG	N-CA-C	-5.75	95.49	111.00
1	B	271	LYS	CB-CA-C	-5.52	99.35	110.40
1	A	241	ARG	NE-CZ-NH2	-5.48	117.56	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	271	LYS	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2090	0	2148	36	1
1	B	2075	0	2135	27	0
1	C	2062	0	2115	30	0
1	X	2068	0	2120	21	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	X	10	0	0	0	0
3	A	196	0	0	20	0
3	B	205	0	0	11	0
3	C	210	0	0	13	2
3	X	205	0	0	2	1
All	All	9151	0	8518	112	2

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 7.

All (112) 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:271:LYS:O	3:B:501:HOH:O	1.59	1.16
1:A:278:SER:CA	3:A:611:HOH:O	1.92	1.15
1:X:271:LYS:HG3	1:X:271:LYS:O	1.37	1.07
1:A:278:SER:HA	3:A:611:HOH:O	1.54	0.98
1:C:217:LYS:HE3	1:C:218:ASP:HB2	1.47	0.95
1:A:220:ASN:OD1	3:A:601:HOH:O	1.85	0.94
1:A:281:ASP:OD2	3:A:602:HOH:O	1.86	0.91
1:A:271:LYS:O	3:A:603:HOH:O	1.92	0.88
1:B:241:ARG:NH2	3:B:505:HOH:O	2.09	0.86
1:C:217:LYS:CE	1:C:218:ASP:HB2	2.07	0.84
1:C:82:GLU:OE1	3:C:601:HOH:O	1.95	0.83
1:A:278:SER:C	3:A:611:HOH:O	2.11	0.83
1:C:103:ASP:O	1:C:103:ASP:OD1	1.98	0.80
1:X:255:GLU:OE1	3:X:501:HOH:O	2.01	0.78
1:A:96:GLY:O	3:A:604:HOH:O	2.00	0.78
1:X:105:ASP:OD1	1:X:111:ARG:NH1	2.17	0.78
1:A:111:ARG:NH2	3:A:608:HOH:O	2.16	0.77
1:C:214:ARG:O	1:C:217:LYS:HB3	1.85	0.77
1:X:271:LYS:CG	1:X:271:LYS:O	2.24	0.77
1:A:62:LYS:NZ	3:A:607:HOH:O	2.16	0.76
1:C:103:ASP:O	1:C:103:ASP:CG	2.17	0.76
1:X:106:GLU:C	1:X:111:ARG:HH12	1.89	0.76
1:C:217:LYS:HE2	1:C:218:ASP:H	1.51	0.75
1:A:111:ARG:CZ	3:A:608:HOH:O	2.36	0.74
1:C:16:LYS:NZ	3:C:606:HOH:O	2.21	0.73
1:A:103:ASP:O	1:A:103:ASP:CG	2.27	0.72
1:A:289:ASP:OD2	3:A:606:HOH:O	2.07	0.72
1:C:168:ASN:ND2	3:C:605:HOH:O	2.21	0.72
1:C:82:GLU:OE2	3:C:603:HOH:O	2.07	0.72
1:C:185:GLU:OE1	3:C:602:HOH:O	2.09	0.71
1:B:82:GLU:OE2	3:B:504:HOH:O	2.08	0.69
1:C:189:ARG:NH2	3:C:610:HOH:O	2.26	0.69
1:A:13:ILE:HG13	1:A:14:LYS:H	1.57	0.68
1:C:285:ILE:HG23	1:C:289:ASP:HB2	1.75	0.68
1:B:217:LYS:HG3	1:B:244:LEU:HD11	1.77	0.67
1:B:281:ASP:OD1	3:B:506:HOH:O	2.13	0.66
1:C:266:ASP:OD2	3:C:604:HOH:O	2.14	0.65
1:B:151:ASN:O	1:B:155:ARG:HD3	1.96	0.65
1:X:103:ASP:CG	1:X:103:ASP:O	2.36	0.64
1:B:205:SER:HB3	1:B:208:ASP:OD1	1.98	0.62
1:C:217:LYS:HE2	1:C:218:ASP:N	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:CYS:HB2	3:B:554:HOH:O	2.01	0.61
1:X:14:LYS:HE3	1:X:46:VAL:HG22	1.82	0.61
1:A:272:SER:OG	1:A:272:SER:O	2.19	0.59
1:B:95:LYS:NZ	3:B:508:HOH:O	2.30	0.59
1:X:106:GLU:O	1:X:111:ARG:NH1	2.34	0.59
1:A:278:SER:OG	3:A:609:HOH:O	2.17	0.59
1:A:105:ASP:OD1	1:A:111:ARG:NH1	2.36	0.58
1:C:38:ARG:NH1	3:C:609:HOH:O	2.25	0.58
1:X:151:ASN:O	1:X:155:ARG:HD3	2.05	0.57
1:A:285:ILE:HG23	1:A:289:ASP:HB2	1.85	0.57
1:B:221:TYR:OH	3:B:503:HOH:O	1.95	0.56
1:X:87:LYS:NZ	1:A:82:GLU:HG2	2.20	0.56
1:B:103:ASP:CG	1:B:103:ASP:O	2.42	0.56
1:A:132:LYS:NZ	3:A:615:HOH:O	2.37	0.55
1:X:271:LYS:O	1:X:272:SER:OG	2.20	0.55
1:A:224:ARG:NE	3:A:616:HOH:O	2.39	0.55
1:A:241:ARG:O	1:A:287:LYS:NZ	2.37	0.55
1:C:217:LYS:CE	1:C:218:ASP:CB	2.85	0.55
1:A:229:CYS:HB2	3:A:706:HOH:O	2.07	0.54
1:A:111:ARG:NE	3:A:608:HOH:O	2.41	0.54
1:X:222:LYS:NZ	1:X:242:ASN:O	2.41	0.53
1:C:214:ARG:NH1	3:C:602:HOH:O	2.04	0.53
1:A:151:ASN:O	1:A:155:ARG:HD3	2.09	0.52
1:B:87:LYS:NZ	3:B:502:HOH:O	1.91	0.52
1:C:223:ASP:N	1:C:223:ASP:OD1	2.41	0.52
1:B:205:SER:CB	1:B:208:ASP:OD1	2.58	0.52
1:B:252:LYS:HD2	1:B:280:ILE:CG2	2.40	0.51
1:C:229:CYS:HB2	3:C:659:HOH:O	2.10	0.51
1:B:111:ARG:HD2	1:B:111:ARG:O	2.11	0.51
1:A:103:ASP:O	1:A:103:ASP:OD1	2.28	0.51
1:A:278:SER:N	3:A:611:HOH:O	2.24	0.50
1:A:197:ASN:ND2	3:A:616:HOH:O	2.42	0.49
1:C:151:ASN:O	1:C:155:ARG:HD3	2.13	0.49
1:B:68:THR:HG21	3:B:508:HOH:O	2.13	0.48
1:A:222:LYS:HA	1:A:223:ASP:HA	1.70	0.47
1:B:271:LYS:N	1:B:271:LYS:HD3	2.30	0.47
1:X:250:VAL:HG21	1:X:285:ILE:HD12	1.95	0.47
1:C:222:LYS:HA	1:C:223:ASP:HA	1.54	0.47
1:C:161:GLN:HG2	3:C:699:HOH:O	2.15	0.47
1:X:285:ILE:HG23	1:X:289:ASP:HB2	1.97	0.47
1:X:242:ASN:ND2	3:X:517:HOH:O	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LYS:HG3	3:A:696:HOH:O	2.15	0.46
1:C:217:LYS:HE3	1:C:218:ASP:CB	2.30	0.46
1:X:40:ALA:HB2	1:X:47:VAL:HG23	1.96	0.46
1:C:252:LYS:HD2	1:C:280:ILE:CG2	2.46	0.46
1:B:270:GLY:C	1:B:271:LYS:HD3	2.37	0.45
1:A:241:ARG:HE	1:A:290:VAL:HG12	1.82	0.45
1:C:217:LYS:CE	1:C:218:ASP:N	2.79	0.45
1:A:165:SER:OG	3:A:610:HOH:O	2.19	0.45
1:C:88:LYS:HG3	3:C:692:HOH:O	2.17	0.44
1:B:252:LYS:O	1:B:256:MET:HG3	2.18	0.44
1:X:107:GLU:C	1:X:111:ARG:HH22	2.21	0.44
1:B:187:ILE:HA	1:B:187:ILE:HD12	1.85	0.43
1:X:147:LYS:HZ3	1:X:175:ASP:HA	1.83	0.43
1:C:190:LYS:NZ	3:C:617:HOH:O	2.47	0.43
1:B:250:VAL:HG21	1:B:285:ILE:HD12	2.01	0.43
1:B:202:PHE:CZ	1:B:204:THR:HG22	2.54	0.42
1:B:287:LYS:O	1:B:288:SER:CB	2.66	0.42
1:X:87:LYS:HZ1	1:A:82:GLU:HG2	1.84	0.42
1:B:111:ARG:HH11	1:B:111:ARG:HD2	1.68	0.42
1:B:252:LYS:NZ	3:B:515:HOH:O	2.46	0.42
1:C:128:LYS:HA	1:C:162:TYR:CE2	2.55	0.42
1:B:55:ALA:O	3:B:507:HOH:O	2.22	0.41
1:X:106:GLU:O	1:X:111:ARG:NH2	2.53	0.41
1:B:33:LYS:O	1:B:37:GLU:HG2	2.20	0.41
1:X:14:LYS:HE3	1:X:46:VAL:CG2	2.49	0.41
1:A:105:ASP:OD1	1:A:106:GLU:N	2.53	0.41
1:A:13:ILE:HG13	1:A:14:LYS:N	2.31	0.41
1:A:241:ARG:C	1:A:287:LYS:NZ	2.73	0.41
1:C:187:ILE:HD12	1:C:187:ILE:HA	1.91	0.41
1:A:33:LYS:NZ	1:A:37:GLU:OE2	2.43	0.40

All (2) 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)
3:X:514:HOH:O	3:C:748:HOH:O[1_545]	1.72	0.48
1:A:255:GLU:OE1	3:C:609:HOH:O[1_456]	2.10	0.10

5.3 Torsion angles

5.3.1 Protein backbone

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	270/316 (85%)	259 (96%)	11 (4%)	0	100	100
1	B	268/316 (85%)	261 (97%)	7 (3%)	0	100	100
1	C	266/316 (84%)	256 (96%)	9 (3%)	1 (0%)	34	32
1	X	267/316 (84%)	261 (98%)	6 (2%)	0	100	100
All	All	1071/1264 (85%)	1037 (97%)	33 (3%)	1 (0%)	51	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	103	ASP

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	232/271 (86%)	226 (97%)	6 (3%)	46	50
1	B	230/271 (85%)	228 (99%)	2 (1%)	78	84
1	C	228/271 (84%)	224 (98%)	4 (2%)	59	65
1	X	229/271 (84%)	227 (99%)	2 (1%)	78	84
All	All	919/1084 (85%)	905 (98%)	14 (2%)	65	71

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	111	ARG
1	X	282	VAL
1	A	53	THR
1	A	111	ARG
1	A	224	ARG
1	A	278	SER
1	A	282	VAL
1	A	288	SER
1	B	111	ARG
1	B	282	VAL
1	C	111	ARG
1	C	217	LYS
1	C	271	LYS
1	C	278	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	ASN
1	C	168	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	XYP	C	500	-	10,10,10	0.47	0	14,14,14	0.45	0
2	XYP	B	401	-	10,10,10	0.47	0	14,14,14	0.45	0
2	XYP	A	501	-	10,10,10	1.74	2 (20%)	14,14,14	1.68	4 (28%)
2	XYP	X	401	-	10,10,10	0.47	0	14,14,14	0.45	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XYP	C	500	-	-	-	0/1/1/1
2	XYP	B	401	-	-	-	0/1/1/1
2	XYP	A	501	-	-	-	0/1/1/1
2	XYP	X	401	-	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	XYP	O5-C5	-3.88	1.37	1.43
2	A	501	XYP	O2-C2	-2.15	1.37	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	XYP	O3-C3-C2	2.72	116.65	110.35
2	A	501	XYP	O5-C1-C2	2.72	113.47	109.43
2	A	501	XYP	O3-C3-C4	-2.18	105.83	109.99
2	A	501	XYP	O2-C2-C3	2.14	115.31	110.35

There are no chirality outliers.

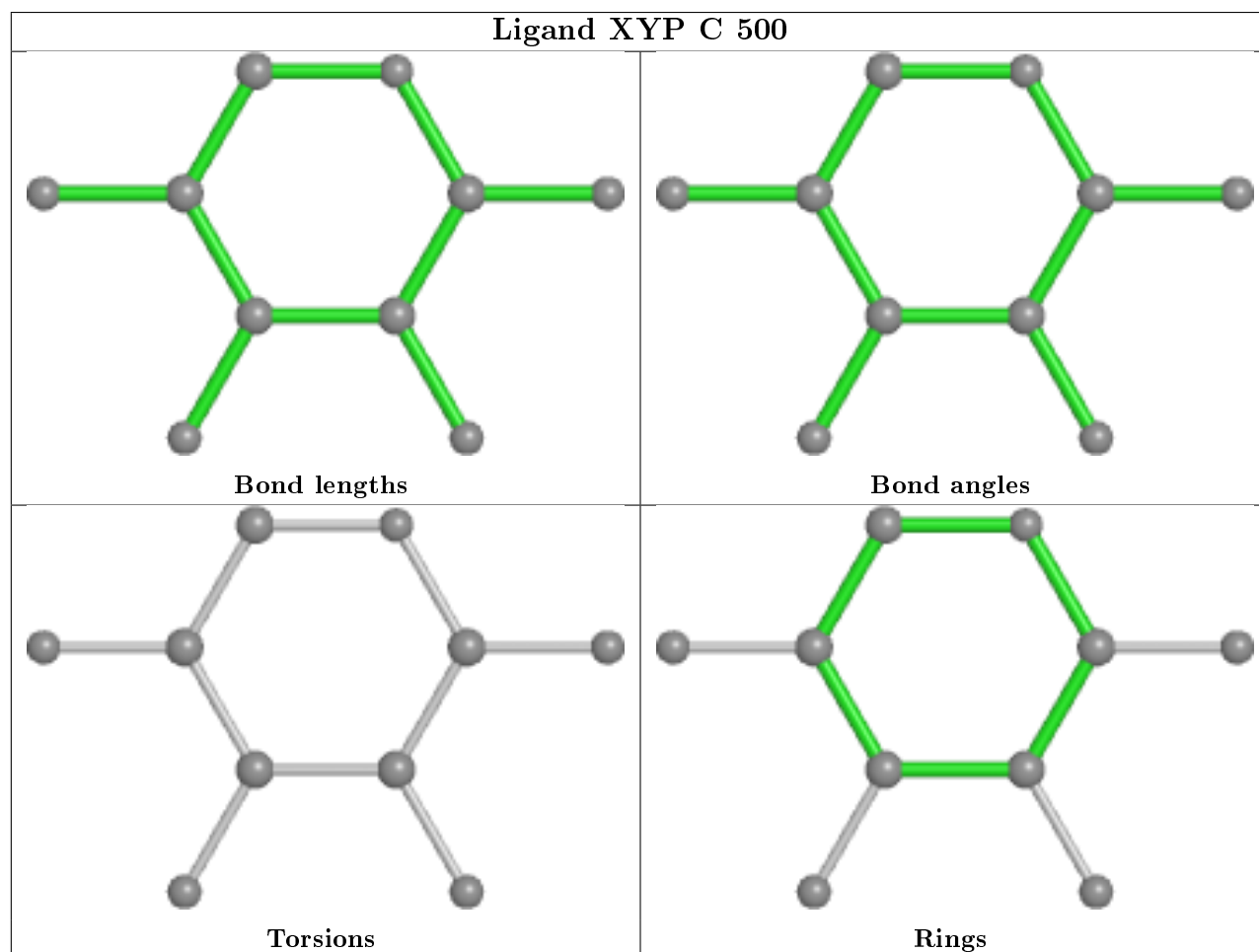
There are no torsion outliers.

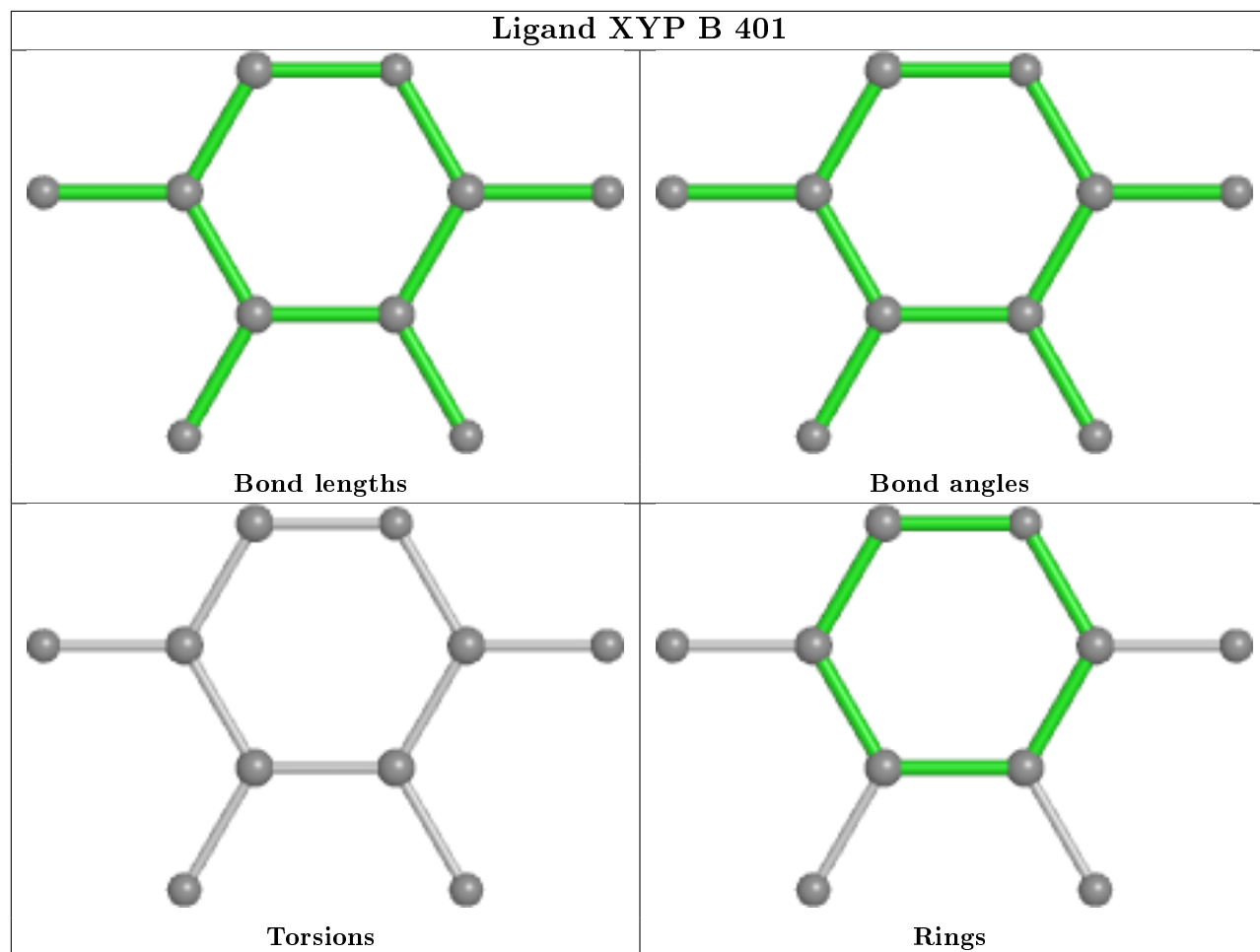
There are no ring outliers.

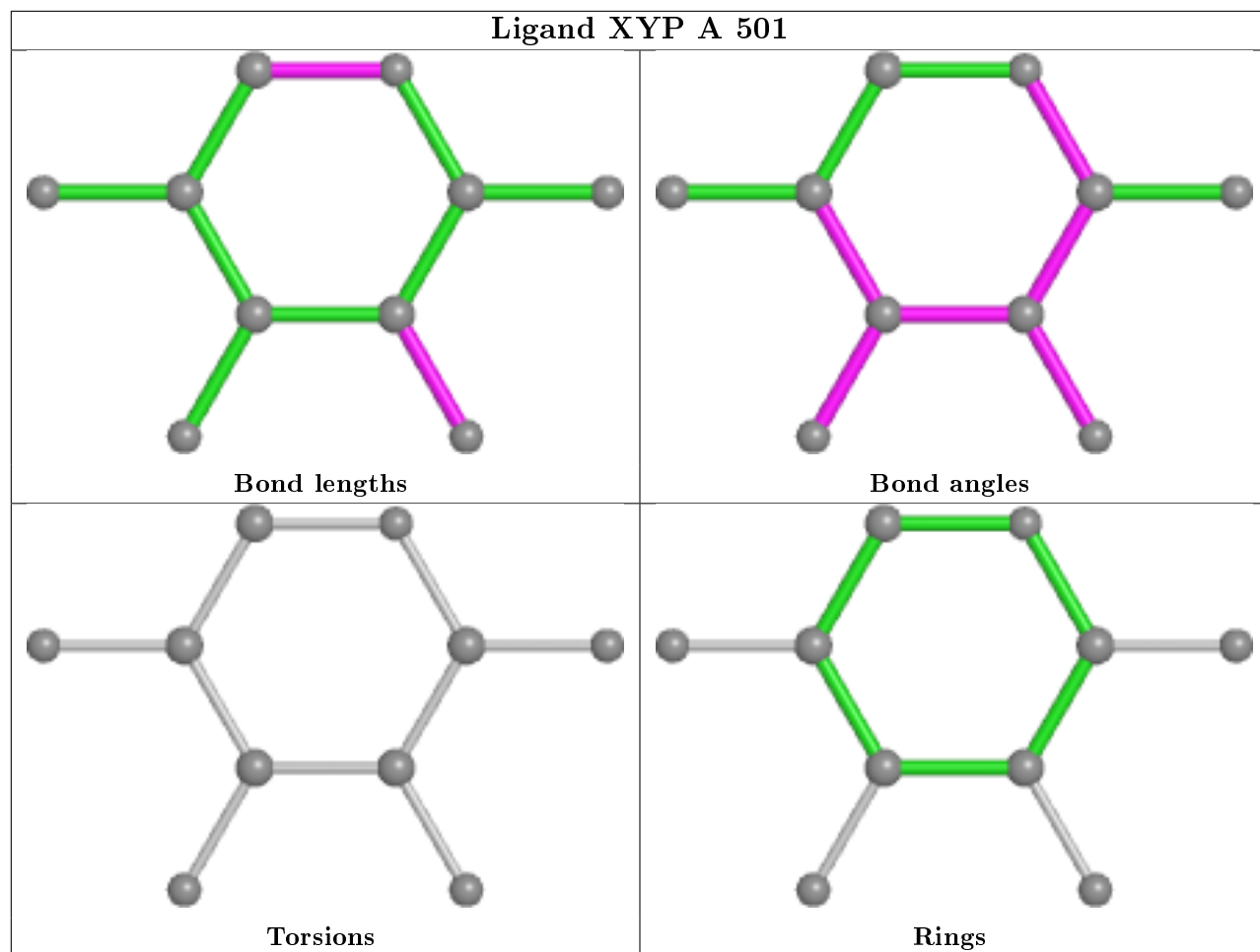
No monomer is involved in short contacts.

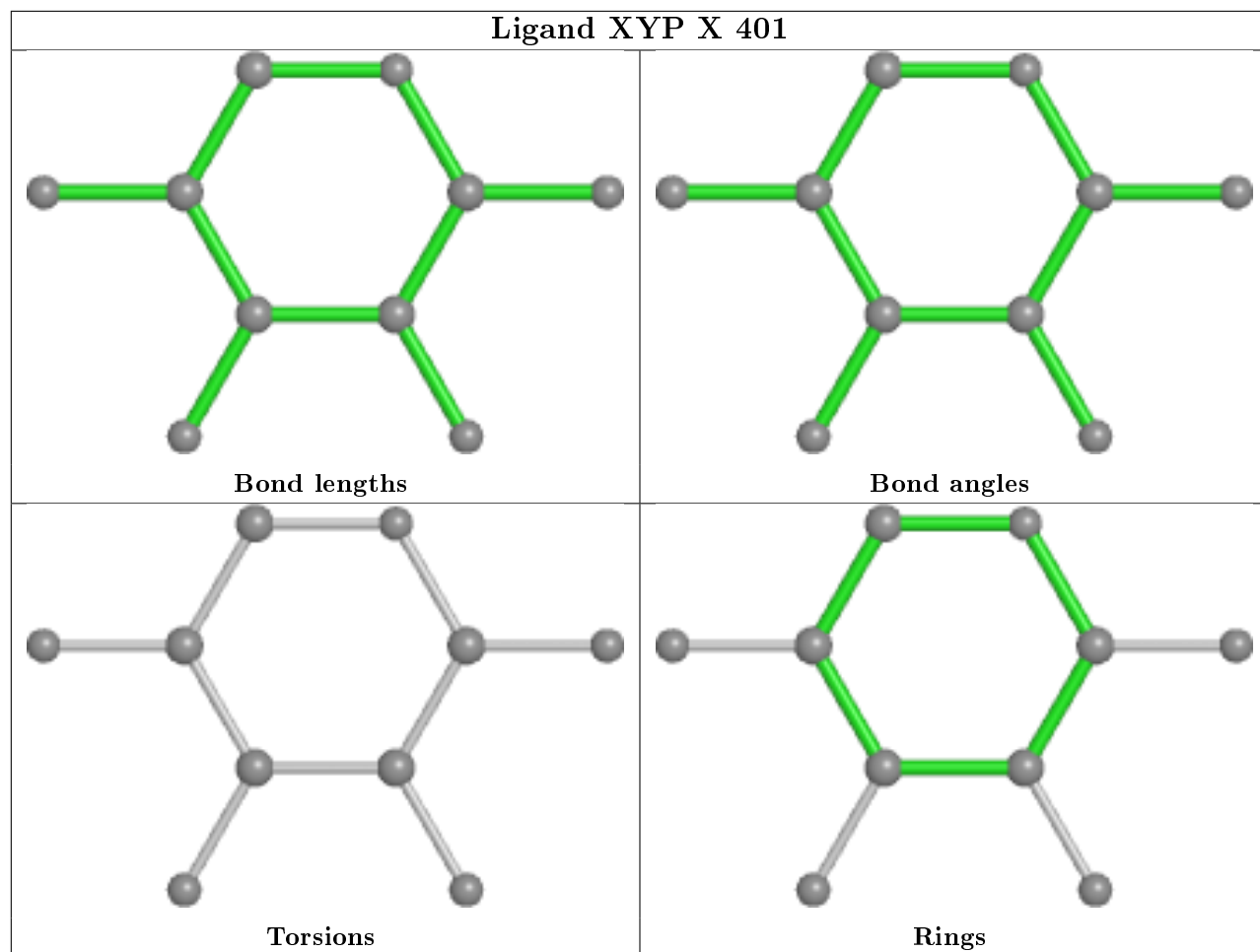
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	274/316 (86%)	0.01	14 (5%) 28 33	11, 21, 41, 70	0
1	B	272/316 (86%)	-0.12	12 (4%) 34 40	11, 19, 37, 65	0
1	C	270/316 (85%)	-0.12	11 (4%) 37 43	11, 21, 40, 52	0
1	X	271/316 (85%)	-0.17	5 (1%) 68 72	11, 19, 36, 56	0
All	All	1087/1264 (85%)	-0.10	42 (3%) 39 45	11, 20, 39, 70	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	13	ILE	9.5
1	A	12	PRO	9.2
1	B	12	PRO	9.0
1	B	13	ILE	8.6
1	B	288	SER	6.0
1	B	272	SER	5.6
1	A	223	ASP	5.3
1	C	289	ASP	4.9
1	B	223	ASP	4.9
1	A	290	VAL	4.8
1	C	223	ASP	4.4
1	C	288	SER	3.9
1	A	288	SER	3.8
1	C	224	ARG	3.8
1	B	270	GLY	3.7
1	X	272	SER	3.5
1	A	196	ASP	3.5
1	A	287	LYS	3.3
1	B	224	ARG	3.2
1	C	271	LYS	3.2
1	X	223	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	197	ASN	3.2
1	A	270	GLY	3.1
1	A	272	SER	3.0
1	A	271	LYS	3.0
1	A	289	ASP	2.9
1	C	196	ASP	2.9
1	B	271	LYS	2.9
1	C	220	ASN	2.8
1	A	94	GLU	2.7
1	C	197	ASN	2.7
1	C	221	TYR	2.5
1	B	221	TYR	2.4
1	A	224	ARG	2.4
1	X	288	SER	2.2
1	X	196	ASP	2.2
1	X	271	LYS	2.2
1	B	196	ASP	2.2
1	C	218	ASP	2.1
1	C	219	LEU	2.1
1	B	197	ASN	2.1
1	B	220	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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

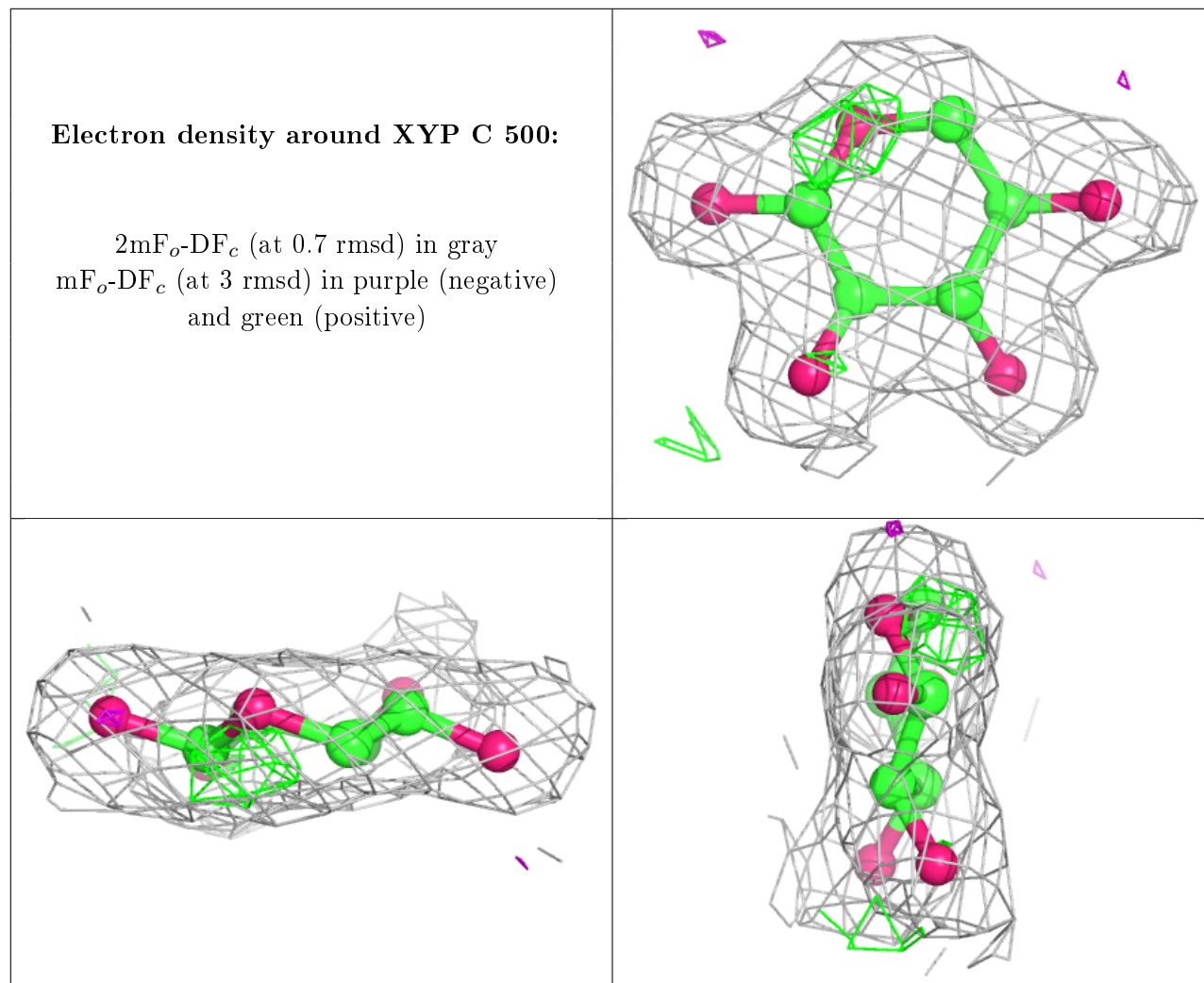
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	XYP	C	500	10/10	0.95	0.19	12,14,19,21	0
2	XYP	X	401	10/10	0.95	0.13	9,12,14,16	0

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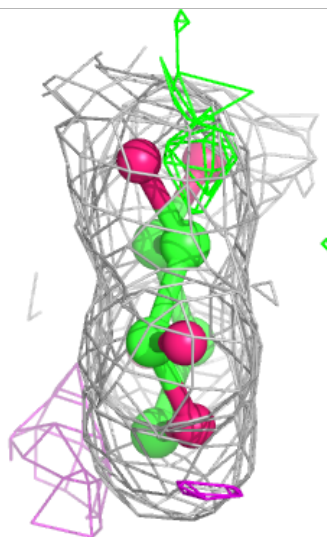
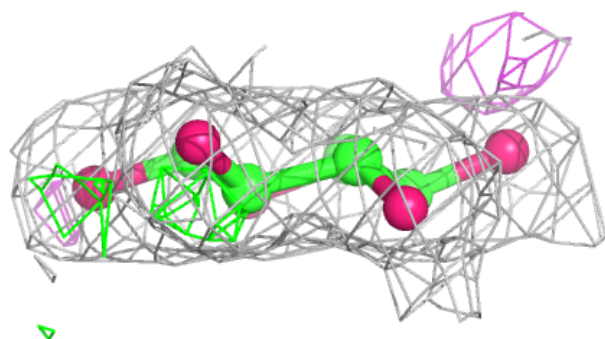
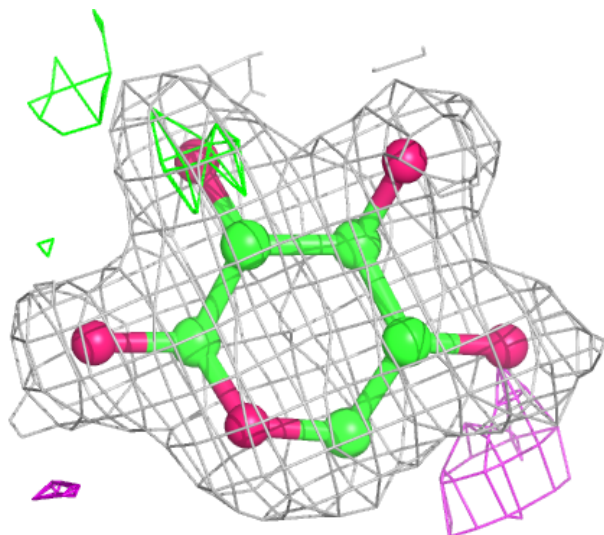
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	XYP	B	401	10/10	0.96	0.14	11,12,17,19	0
2	XYP	A	501	10/10	0.97	0.18	12,14,19,19	0

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.



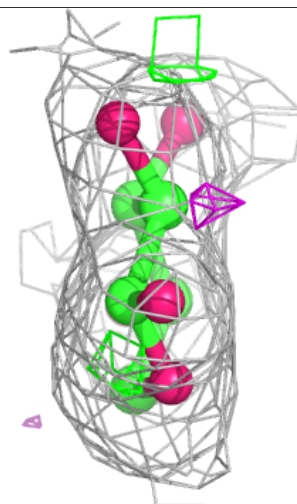
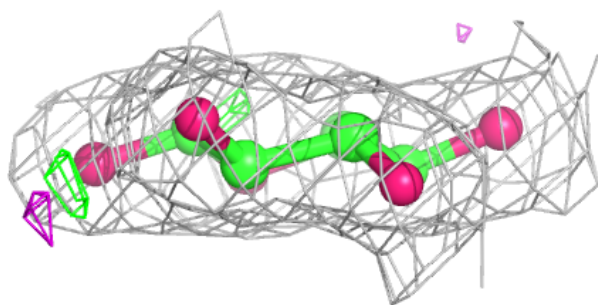
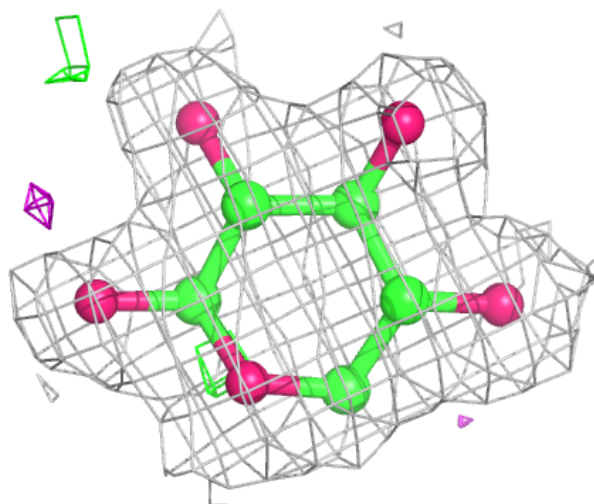
Electron density around XYP X 401:

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



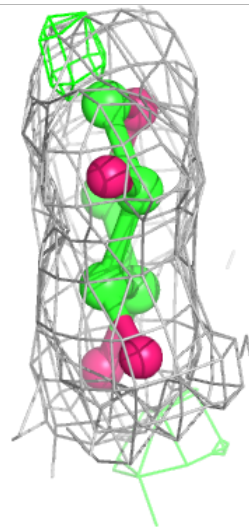
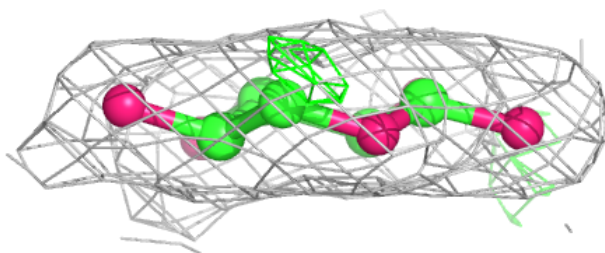
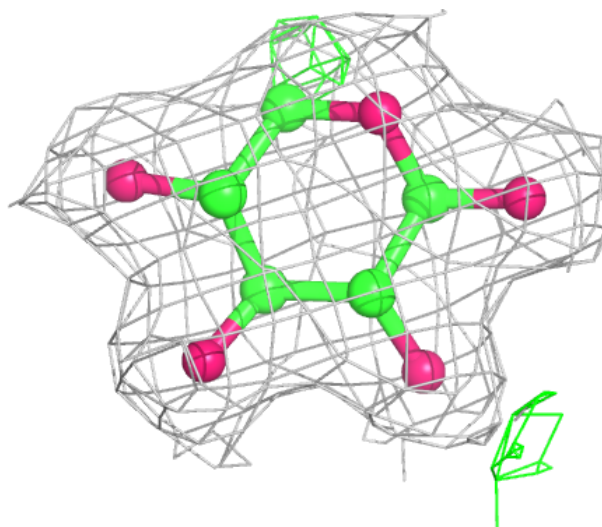
Electron density around XYP B 401:

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



Electron density around XYP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.