



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:02 PM UTC

PDB ID : 8EGM / pdb_00008egm
Title : Crystal Structure of UDP-N-acetylmuramate-L-alanine ligase (UDP-N-acetylmuramoyl-L-alanine synthetase, MurC) *Pseudomonas aeruginosa* in complex with compound AZ13644908
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2022-09-12
Resolution : 2.20 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

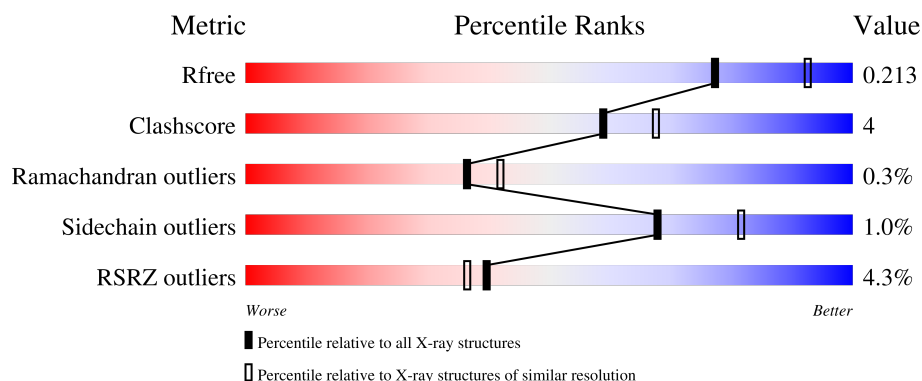
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	316	3% 84% 11% ..
1	B	316	4% 88% 8% .
1	C	316	4% 85% 10% .
1	D	316	4% 85% 10% ..

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Mol	Chain	Length	Quality of chain
1	E	316	4% 84% 10%
1	F	316	4% 85% 11%
1	G	316	4% 84% 12%
1	H	316	4% 84% 11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19366 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 UDP-N-acetylmuramate--L-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	6	0
			2264	1433	396	426	9			
1	B	303	Total	C	N	O	S	0	7	0
			2278	1441	396	432	9			
1	C	303	Total	C	N	O	S	0	3	0
			2256	1425	394	427	10			
1	D	303	Total	C	N	O	S	0	8	0
			2272	1438	396	429	9			
1	E	303	Total	C	N	O	S	0	5	0
			2262	1431	396	426	9			
1	F	303	Total	C	N	O	S	0	8	0
			2283	1446	396	432	9			
1	G	303	Total	C	N	O	S	0	6	0
			2264	1433	396	426	9			
1	H	303	Total	C	N	O	S	0	6	0
			2278	1439	398	432	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	MET	-	initiating methionine	UNP Q9HW02
A	8	ALA	-	expression tag	UNP Q9HW02
A	9	HIS	-	expression tag	UNP Q9HW02
A	10	HIS	-	expression tag	UNP Q9HW02
A	11	HIS	-	expression tag	UNP Q9HW02
A	12	HIS	-	expression tag	UNP Q9HW02
A	13	HIS	-	expression tag	UNP Q9HW02
A	14	HIS	-	expression tag	UNP Q9HW02
A	15	MET	-	expression tag	UNP Q9HW02
B	7	MET	-	initiating methionine	UNP Q9HW02
B	8	ALA	-	expression tag	UNP Q9HW02
B	9	HIS	-	expression tag	UNP Q9HW02
B	10	HIS	-	expression tag	UNP Q9HW02

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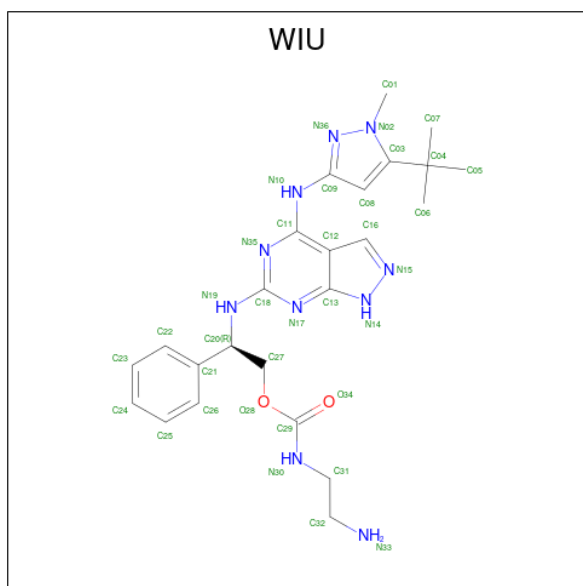
Chain	Residue	Modelled	Actual	Comment	Reference
B	11	HIS	-	expression tag	UNP Q9HW02
B	12	HIS	-	expression tag	UNP Q9HW02
B	13	HIS	-	expression tag	UNP Q9HW02
B	14	HIS	-	expression tag	UNP Q9HW02
B	15	MET	-	expression tag	UNP Q9HW02
C	7	MET	-	initiating methionine	UNP Q9HW02
C	8	ALA	-	expression tag	UNP Q9HW02
C	9	HIS	-	expression tag	UNP Q9HW02
C	10	HIS	-	expression tag	UNP Q9HW02
C	11	HIS	-	expression tag	UNP Q9HW02
C	12	HIS	-	expression tag	UNP Q9HW02
C	13	HIS	-	expression tag	UNP Q9HW02
C	14	HIS	-	expression tag	UNP Q9HW02
C	15	MET	-	expression tag	UNP Q9HW02
D	7	MET	-	initiating methionine	UNP Q9HW02
D	8	ALA	-	expression tag	UNP Q9HW02
D	9	HIS	-	expression tag	UNP Q9HW02
D	10	HIS	-	expression tag	UNP Q9HW02
D	11	HIS	-	expression tag	UNP Q9HW02
D	12	HIS	-	expression tag	UNP Q9HW02
D	13	HIS	-	expression tag	UNP Q9HW02
D	14	HIS	-	expression tag	UNP Q9HW02
D	15	MET	-	expression tag	UNP Q9HW02
E	7	MET	-	initiating methionine	UNP Q9HW02
E	8	ALA	-	expression tag	UNP Q9HW02
E	9	HIS	-	expression tag	UNP Q9HW02
E	10	HIS	-	expression tag	UNP Q9HW02
E	11	HIS	-	expression tag	UNP Q9HW02
E	12	HIS	-	expression tag	UNP Q9HW02
E	13	HIS	-	expression tag	UNP Q9HW02
E	14	HIS	-	expression tag	UNP Q9HW02
E	15	MET	-	expression tag	UNP Q9HW02
F	7	MET	-	initiating methionine	UNP Q9HW02
F	8	ALA	-	expression tag	UNP Q9HW02
F	9	HIS	-	expression tag	UNP Q9HW02
F	10	HIS	-	expression tag	UNP Q9HW02
F	11	HIS	-	expression tag	UNP Q9HW02
F	12	HIS	-	expression tag	UNP Q9HW02
F	13	HIS	-	expression tag	UNP Q9HW02
F	14	HIS	-	expression tag	UNP Q9HW02
F	15	MET	-	expression tag	UNP Q9HW02
G	7	MET	-	initiating methionine	UNP Q9HW02

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Chain	Residue	Modelled	Actual	Comment	Reference
G	8	ALA	-	expression tag	UNP Q9HW02
G	9	HIS	-	expression tag	UNP Q9HW02
G	10	HIS	-	expression tag	UNP Q9HW02
G	11	HIS	-	expression tag	UNP Q9HW02
G	12	HIS	-	expression tag	UNP Q9HW02
G	13	HIS	-	expression tag	UNP Q9HW02
G	14	HIS	-	expression tag	UNP Q9HW02
G	15	MET	-	expression tag	UNP Q9HW02
H	7	MET	-	initiating methionine	UNP Q9HW02
H	8	ALA	-	expression tag	UNP Q9HW02
H	9	HIS	-	expression tag	UNP Q9HW02
H	10	HIS	-	expression tag	UNP Q9HW02
H	11	HIS	-	expression tag	UNP Q9HW02
H	12	HIS	-	expression tag	UNP Q9HW02
H	13	HIS	-	expression tag	UNP Q9HW02
H	14	HIS	-	expression tag	UNP Q9HW02
H	15	MET	-	expression tag	UNP Q9HW02

- Molecule 2 is (2R)-2-({4-[(5-tert-butyl-1-methyl-1H-pyrazol-3-yl)amino]-1H-pyrazolo[3,4-d]pyrimidin-6-yl}amino)-2-phenylethyl (2-aminoethyl)carbamate (CCD ID: WIU) (formula: C₂₄H₃₂N₁₀O₂) (labeled as "Ligand of Interest" by depositor).



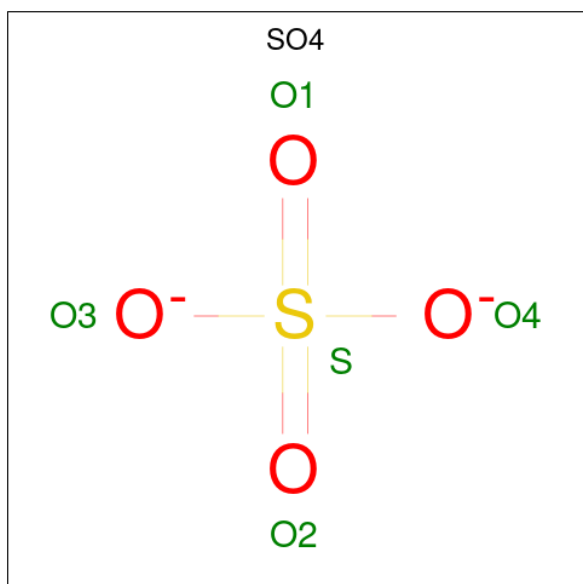
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			36	24	10	2		
2	B	1	Total	C	N	O	0	0
			36	24	10	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			36	24	10	2		
2	D	1	Total	C	N	O	0	0
			36	24	10	2		
2	E	1	Total	C	N	O	0	0
			36	24	10	2		
2	F	1	Total	C	N	O	0	0
			36	24	10	2		
2	G	1	Total	C	N	O	0	0
			36	24	10	2		
2	H	1	Total	C	N	O	0	0
			36	24	10	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

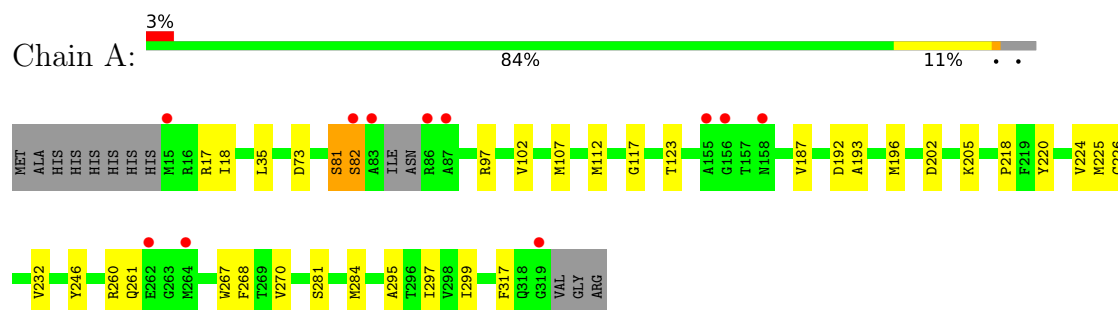
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	108	Total	O	0	0
			108	108		
4	B	123	Total	O	0	0
			123	123		
4	C	112	Total	O	0	0
			112	112		
4	D	106	Total	O	0	0
			106	106		
4	E	99	Total	O	0	0
			99	99		
4	F	110	Total	O	0	0
			110	110		
4	G	103	Total	O	0	0
			103	103		
4	H	120	Total	O	0	0
			120	120		

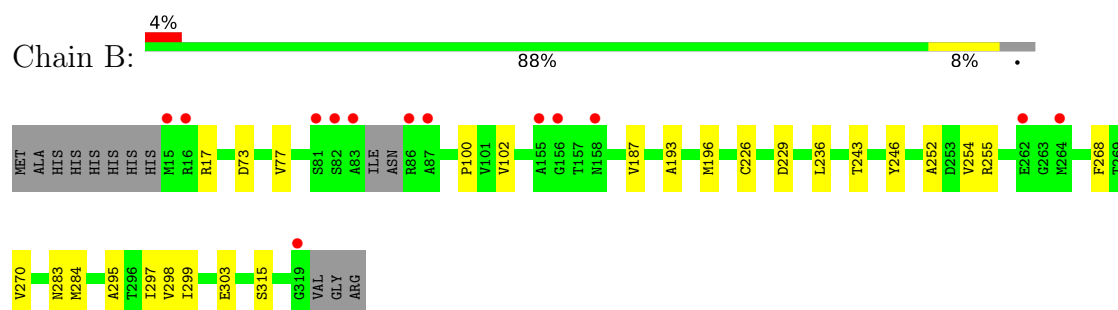
3 Residue-property plots [i](#)

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.

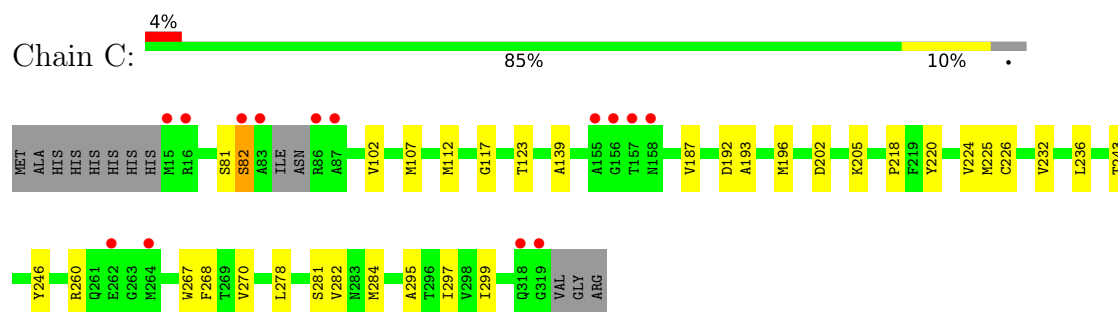
- Molecule 1: UDP-N-acetylmuramate--L-alanine ligase



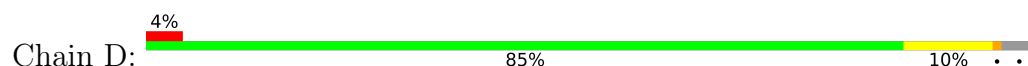
- Molecule 1: UDP-N-acetylmuramate--L-alanine ligase



- Molecule 1: UDP-N-acetylmuramate--L-alanine ligase



- Molecule 1: UDP-N-acetylmuramate--L-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	285.26Å 109.20Å 108.71Å 90.00° 112.40° 90.00°	Depositor
Resolution (Å)	42.66 – 2.20 42.66 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.66-2.20) 99.9 (42.66-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.20.1 dev 4694	Depositor
R, R_{free}	0.182 , 0.212 0.182 , 0.213	Depositor DCC
R_{free} test set	1915 reflections (1.23%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.57$, $\langle L^2 \rangle = 0.42$	Xtriage
Estimated twinning fraction	0.021 for -h-k-l,l,k 0.017 for -h+k-l,-l,-k 0.046 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19366	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.85 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0795e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: WIU, SO4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.32	0/2320	0.49	0/3158
1	B	0.33	0/2337	0.50	0/3179
1	C	0.31	0/2303	0.49	0/3133
1	D	0.30	0/2334	0.47	0/3179
1	E	0.33	0/2315	0.51	0/3150
1	F	0.33	0/2345	0.51	0/3191
1	G	0.31	0/2320	0.49	0/3157
1	H	0.32	0/2337	0.50	0/3179
All	All	0.32	0/18611	0.50	0/25326

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2264	0	2250	20	0
1	B	2278	0	2264	15	0
1	C	2256	0	2237	19	0
1	D	2272	0	2256	19	0
1	E	2262	0	2248	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2283	0	2275	21	0
1	G	2264	0	2250	23	0
1	H	2278	0	2261	23	0
2	A	36	0	0	0	0
2	B	36	0	0	1	0
2	C	36	0	0	0	0
2	D	36	0	0	0	0
2	E	36	0	0	1	0
2	F	36	0	0	0	0
2	G	36	0	0	0	0
2	H	36	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
3	G	5	0	0	0	0
3	H	5	0	0	0	0
4	A	108	0	0	1	0
4	B	123	0	0	1	0
4	C	112	0	0	0	0
4	D	106	0	0	1	0
4	E	99	0	0	2	1
4	F	110	0	0	4	1
4	G	103	0	0	2	0
4	H	120	0	0	1	0
All	All	19366	0	18041	160	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (160) 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:77[A]:VAL:HG22	1:B:100:PRO:HG2	1.74	0.70
1:A:18:ILE:HD13	1:A:35:LEU:HD13	1.71	0.70
1:C:193:ALA:HB1	1:C:196:MET:HE2	1.74	0.69
1:H:283[B]:ASN:ND2	1:H:315:SER:O	2.29	0.65
1:D:17[B]:ARG:NH2	1:D:73:ASP:O	2.30	0.64
1:G:187:VAL:CG2	1:G:299[B]:ILE:HD11	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:PHE:O	4:A:501:HOH:O	2.16	0.60
1:H:17[A]:ARG:NH2	1:H:73:ASP:O	2.35	0.60
1:C:202:ASP:HB3	1:C:205:LYS:HD2	1.85	0.59
1:D:303:GLU:OE1	4:D:501:HOH:O	2.17	0.57
1:B:283:ASN:ND2	1:B:315:SER:O	2.37	0.57
1:B:17[A]:ARG:NH2	1:B:73:ASP:O	2.37	0.57
1:E:187:VAL:CG2	1:E:299[B]:ILE:HD11	2.34	0.57
1:F:216:ASN:ND2	4:F:501:HOH:O	2.14	0.57
1:A:202:ASP:HB3	1:A:205:LYS:HD2	1.87	0.57
1:F:187:VAL:CG2	1:F:299[B]:ILE:HD11	2.35	0.57
1:F:255:ARG:NH1	4:F:502:HOH:O	2.27	0.56
1:E:226:CYS:HB2	1:E:246:TYR:CZ	2.41	0.55
1:D:77[A]:VAL:HG22	1:D:100:PRO:HG2	1.87	0.55
1:G:81:SER:O	1:G:83:ALA:N	2.39	0.55
1:H:216:ASN:ND2	4:H:501:HOH:O	2.10	0.55
1:D:207:LYS:HD3	1:D:231:VAL:HG13	1.88	0.55
1:G:16:ARG:NH1	4:G:504:HOH:O	2.38	0.55
1:A:187:VAL:CG2	1:A:299[A]:ILE:HD11	2.37	0.54
1:G:18:ILE:HD13	1:G:35:LEU:HD13	1.89	0.54
1:B:270[B]:VAL:HG21	1:B:297:ILE:HD13	1.90	0.54
1:D:187:VAL:CG2	1:D:299[A]:ILE:HD11	2.38	0.53
1:H:270[A]:VAL:HG21	1:H:297:ILE:HD13	1.89	0.53
1:B:226:CYS:HB2	1:B:246:TYR:CZ	2.44	0.53
1:C:268:PHE:CE2	1:C:270[B]:VAL:HG22	2.44	0.53
1:A:295:ALA:O	1:A:299[B]:ILE:HG12	2.08	0.53
1:E:236:LEU:HD11	1:E:243:THR:HG21	1.90	0.53
1:H:226:CYS:HB2	1:H:246:TYR:CZ	2.44	0.53
1:H:187:VAL:CG2	1:H:299[B]:ILE:HD11	2.39	0.52
1:H:193:ALA:HB1	1:H:196:MET:HE2	1.91	0.52
1:C:187:VAL:CG2	1:C:299[A]:ILE:HD11	2.40	0.52
1:A:193:ALA:HB1	1:A:196:MET:HE2	1.91	0.52
1:C:236:LEU:HD11	1:C:243:THR:HG21	1.92	0.52
1:A:225:MET:HE3	1:A:232:VAL:HG13	1.92	0.52
1:B:303:GLU:OE1	4:B:501:HOH:O	2.19	0.52
1:F:270[B]:VAL:HG21	1:F:297:ILE:HD13	1.93	0.51
1:H:270[B]:VAL:HG11	1:H:297:ILE:HD13	1.92	0.51
1:F:193:ALA:HB1	1:F:196:MET:HE2	1.93	0.51
1:E:207:LYS:HD3	1:E:231:VAL:HG13	1.92	0.51
1:B:187:VAL:CG2	1:B:299[B]:ILE:HD11	2.40	0.51
1:G:226:CYS:HB2	1:G:246:TYR:CZ	2.46	0.51
1:F:261:GLN:NE2	4:F:505:HOH:O	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:98:ARG:NE	4:E:501:HOH:O	2.26	0.51
1:D:81:SER:O	1:D:83:ALA:N	2.44	0.51
1:G:193:ALA:HB1	1:G:196:MET:HE2	1.94	0.50
1:G:102:VAL:HG21	1:G:107:MET:HE2	1.94	0.50
1:D:112:MET:HG3	1:D:117:GLY:HA3	1.93	0.50
1:F:112:MET:HG2	1:F:117:GLY:HA3	1.93	0.49
1:H:187:VAL:HG22	1:H:299[B]:ILE:HD11	1.94	0.49
1:E:77[B]:VAL:HG22	1:E:100:PRO:HG2	1.93	0.49
1:D:102[B]:VAL:HG21	1:D:107:MET:HE2	1.94	0.49
1:E:18:ILE:HG23	1:E:77[A]:VAL:HG13	1.94	0.49
1:E:81:SER:HB3	4:E:506:HOH:O	2.13	0.49
1:D:260:ARG:HG2	1:D:267:TRP:HB2	1.95	0.49
1:E:268:PHE:CE2	1:E:270[B]:VAL:HG22	2.47	0.49
1:H:265:ARG:NH2	1:H:283[B]:ASN:OD1	2.46	0.49
1:H:236:LEU:HD11	1:H:243:THR:HG21	1.95	0.48
1:E:267:TRP:CE3	1:E:281:SER:HB3	2.48	0.48
1:G:202:ASP:HB3	1:G:205:LYS:HD2	1.95	0.48
1:F:195:HIS:ND1	4:F:503:HOH:O	2.35	0.48
1:A:268:PHE:CE2	1:A:270[B]:VAL:HG22	2.48	0.48
1:F:252:ALA:HB3	1:F:255:ARG:HG2	1.95	0.48
1:D:283[B]:ASN:ND2	1:D:315:SER:O	2.47	0.48
1:E:17[B]:ARG:NH2	1:E:73:ASP:O	2.47	0.48
1:E:218:PRO:HB2	1:E:220:TYR:CE1	2.48	0.47
1:E:224:VAL:HG21	1:E:299[A]:ILE:HD13	1.95	0.47
1:G:270[B]:VAL:HG21	1:G:297:ILE:HD13	1.94	0.47
1:A:226:CYS:HB2	1:A:246:TYR:CZ	2.50	0.47
1:B:229:ASP:OD1	2:B:401:WIU:N33	2.47	0.47
1:E:295:ALA:O	1:E:299[A]:ILE:HG12	2.13	0.47
1:F:187:VAL:HG22	1:F:299[B]:ILE:HD11	1.97	0.47
1:H:102:VAL:HG21	1:H:107:MET:HE2	1.97	0.47
1:A:224:VAL:HG21	1:A:299[B]:ILE:HD13	1.96	0.47
1:A:284:MET:HE3	1:A:284:MET:HB2	1.74	0.47
1:F:284:MET:HE3	1:F:284:MET:HB2	1.78	0.47
1:C:102:VAL:HG21	1:C:107:MET:HE2	1.97	0.46
1:D:224:VAL:HG21	1:D:299[B]:ILE:HD13	1.96	0.46
1:F:236:LEU:HD11	1:F:243:THR:HG21	1.96	0.46
1:D:193:ALA:HB1	1:D:196:MET:HE2	1.96	0.46
1:E:46:ASP:OD1	1:E:47:LEU:N	2.48	0.46
1:B:252:ALA:HB3	1:B:255:ARG:HG2	1.98	0.46
1:B:295:ALA:O	1:B:299[A]:ILE:HG12	2.16	0.46
1:D:270[B]:VAL:HG21	1:D:297:ILE:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:295:ALA:O	1:D:299[B]:ILE:HG12	2.15	0.46
1:H:218:PRO:HB2	1:H:220:TYR:CE1	2.50	0.46
1:A:270[B]:VAL:HG21	1:A:297:ILE:HD13	1.97	0.46
1:D:284:MET:HE3	1:D:284:MET:HB2	1.70	0.46
1:F:224:VAL:HG21	1:F:299[A]:ILE:HD13	1.98	0.46
1:E:229:ASP:OD1	2:E:401:WIU:N33	2.49	0.46
1:G:236:LEU:HD11	1:G:243:THR:HG21	1.98	0.46
1:E:68:GLN:HE21	1:E:70:GLU:HB2	1.81	0.46
1:A:112:MET:HG2	1:A:117:GLY:HA3	1.98	0.45
1:C:270[B]:VAL:HG21	1:C:297:ILE:HD13	1.96	0.45
1:C:226:CYS:HB2	1:C:246:TYR:CZ	2.51	0.45
1:G:112:MET:HG2	1:G:117:GLY:HA3	1.99	0.45
1:A:267:TRP:CE3	1:A:281:SER:HB3	2.51	0.45
1:C:295:ALA:O	1:C:299[B]:ILE:HG12	2.16	0.45
1:F:81:SER:O	1:F:83:ALA:N	2.49	0.45
1:F:17[B]:ARG:NH2	1:F:73:ASP:O	2.50	0.45
1:H:81:SER:O	1:H:82:SER:C	2.59	0.45
1:C:112:MET:HG2	1:C:117:GLY:HA3	1.98	0.45
1:F:106[B]:GLU:OE2	1:H:241:ARG:NH2	2.44	0.45
1:F:226:CYS:HB2	1:F:246:TYR:CZ	2.52	0.45
1:F:295:ALA:O	1:F:299[A]:ILE:HG12	2.16	0.45
1:H:267:TRP:CE3	1:H:281:SER:HB3	2.51	0.45
1:C:225:MET:HE3	1:C:232:VAL:HG13	1.99	0.45
1:C:123:THR:HG21	1:C:192:ASP:HB3	1.99	0.44
1:H:123:THR:HG21	1:H:192:ASP:HB3	2.00	0.44
1:E:270[B]:VAL:HG21	1:E:297:ILE:HD13	1.99	0.44
1:G:295:ALA:O	1:G:299[A]:ILE:HG12	2.17	0.44
1:B:254:VAL:HG21	1:B:298:VAL:HG22	1.99	0.44
1:C:224:VAL:HG21	1:C:299[B]:ILE:HD13	2.00	0.43
1:D:226:CYS:HB2	1:D:246:TYR:CZ	2.54	0.43
1:E:112:MET:HG2	1:E:117:GLY:HA3	1.99	0.43
1:D:130:THR:HG23	1:D:168:VAL:HG12	2.00	0.43
1:G:224:VAL:HG21	1:G:299[A]:ILE:HD13	1.99	0.43
1:H:18:ILE:HD13	1:H:35:LEU:HD13	2.01	0.43
1:F:283:ASN:ND2	1:F:315:SER:O	2.51	0.43
1:D:236:LEU:HD11	1:D:243:THR:HG21	2.01	0.43
1:G:225:MET:HE3	1:G:232:VAL:HG13	2.01	0.43
1:C:270[A]:VAL:HG12	1:C:278:LEU:O	2.19	0.42
1:A:218:PRO:HB2	1:A:220:TYR:CE1	2.54	0.42
1:C:282:VAL:HG22	1:C:284:MET:HE2	2.01	0.42
1:G:77[B]:VAL:HG22	1:G:100:PRO:HG2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:308[A]:GLU:H	1:G:308[A]:GLU:CD	2.26	0.42
1:E:187:VAL:HG22	1:E:299[B]:ILE:HD11	2.01	0.42
1:H:284:MET:HE3	1:H:284:MET:HB2	1.90	0.42
1:A:102[B]:VAL:HG21	1:A:107:MET:HE2	2.00	0.42
1:G:254:VAL:HG21	1:G:298:VAL:HG22	2.02	0.42
1:H:225:MET:HE3	1:H:232:VAL:HG13	2.01	0.42
1:H:207:LYS:HD3	1:H:231:VAL:HG13	2.00	0.42
1:B:284:MET:HE3	1:B:284:MET:HB2	1.85	0.42
1:E:45:SER:HA	1:E:64:PHE:O	2.19	0.42
1:G:218:PRO:HB2	1:G:220:TYR:CE1	2.55	0.42
1:A:123:THR:HG21	1:A:192:ASP:HB3	2.01	0.41
1:B:268:PHE:CE2	1:B:270[B]:VAL:HG22	2.55	0.41
1:F:218:PRO:HB2	1:F:220:TYR:CE1	2.55	0.41
1:A:81:SER:O	1:A:82:SER:C	2.63	0.41
1:B:236:LEU:HD11	1:B:243:THR:HG21	2.02	0.41
1:G:187:VAL:HG22	1:G:299[B]:ILE:HD11	2.01	0.41
1:A:73:ASP:OD1	1:A:97:ARG:NH1	2.50	0.41
1:G:268:PHE:CE2	1:G:270[B]:VAL:HG22	2.56	0.41
1:G:284:MET:HE3	1:G:284:MET:HB2	1.93	0.41
1:B:193:ALA:HB1	1:B:196:MET:HE2	2.02	0.41
1:C:139:ALA:O	1:E:58:LYS:HD2	2.21	0.41
1:A:17[A]:ARG:NH2	1:A:73:ASP:O	2.54	0.41
1:H:265:ARG:NH1	1:H:281:SER:OG	2.52	0.41
1:F:77[A]:VAL:HG22	1:F:100:PRO:HG2	2.03	0.41
1:H:214:LEU:HA	1:H:214:LEU:HD23	1.90	0.41
1:G:315:SER:HB3	4:G:531:HOH:O	2.21	0.41
1:D:261:GLN:O	1:D:262:GLU:HG3	2.21	0.40
1:C:267:TRP:CE3	1:C:281:SER:HB3	2.56	0.40
1:G:270[A]:VAL:HG12	1:G:278:LEU:O	2.21	0.40
1:C:218:PRO:HB2	1:C:220:TYR:CE1	2.56	0.40
1:C:81:SER:O	1:C:82:SER:C	2.64	0.40

All (1) 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)
4:E:597:HOH:O	4:F:610:HOH:O[2_545]	2.12	0.08

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	305/316 (96%)	296 (97%)	8 (3%)	1 (0%)	36	42
1	B	306/316 (97%)	299 (98%)	7 (2%)	0	100	100
1	C	302/316 (96%)	292 (97%)	9 (3%)	1 (0%)	36	42
1	D	307/316 (97%)	298 (97%)	8 (3%)	1 (0%)	36	42
1	E	304/316 (96%)	294 (97%)	9 (3%)	1 (0%)	36	42
1	F	307/316 (97%)	299 (97%)	7 (2%)	1 (0%)	36	42
1	G	305/316 (96%)	298 (98%)	6 (2%)	1 (0%)	36	42
1	H	306/316 (97%)	297 (97%)	8 (3%)	1 (0%)	36	42
All	All	2442/2528 (97%)	2373 (97%)	62 (2%)	7 (0%)	36	42

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	82	SER
1	F	82	SER
1	G	82	SER
1	C	82	SER
1	A	82	SER
1	E	82	SER
1	H	82	SER

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	233/249 (94%)	230 (99%)	3 (1%)	61	76
1	B	236/249 (95%)	235 (100%)	1 (0%)	84	92
1	C	233/249 (94%)	232 (100%)	1 (0%)	84	92
1	D	235/249 (94%)	233 (99%)	2 (1%)	70	84
1	E	233/249 (94%)	229 (98%)	4 (2%)	53	69
1	F	237/249 (95%)	234 (99%)	3 (1%)	61	76
1	G	233/249 (94%)	231 (99%)	2 (1%)	70	84
1	H	236/249 (95%)	234 (99%)	2 (1%)	73	85
All	All	1876/1992 (94%)	1858 (99%)	18 (1%)	68	81

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	SER
1	A	260	ARG
1	A	261	GLN
1	B	102	VAL
1	C	260	ARG
1	D	81	SER
1	D	260	ARG
1	E	58	LYS
1	E	68	GLN
1	E	81	SER
1	E	102	VAL
1	F	58	LYS
1	F	102	VAL
1	F	260	ARG
1	G	81	SER
1	G	260	ARG
1	H	58	LYS
1	H	81	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	261	GLN
1	B	37	ASN
1	B	68	GLN

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Mol	Chain	Res	Type
1	B	180	HIS
1	C	37	ASN
1	C	180	HIS
1	C	312	GLN
1	D	37	ASN
1	E	37	ASN
1	E	68	GLN
1	F	37	ASN
1	F	195	HIS
1	G	37	ASN
1	G	180	HIS
1	G	195	HIS
1	G	312	GLN
1	H	37	ASN
1	H	180	HIS
1	H	258	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 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	WIU	B	401	-	39,39,39	1.42	3 (7%)	45,55,55	1.37	5 (11%)
2	WIU	D	401	-	39,39,39	1.67	6 (15%)	45,55,55	1.35	5 (11%)
3	SO4	B	402	-	4,4,4	0.79	0	6,6,6	0.59	0
2	WIU	H	401	-	39,39,39	1.41	2 (5%)	45,55,55	1.36	5 (11%)
2	WIU	G	401	-	39,39,39	1.31	2 (5%)	45,55,55	1.45	8 (17%)
3	SO4	A	402	-	4,4,4	0.85	0	6,6,6	0.21	0
3	SO4	C	402	-	4,4,4	0.84	0	6,6,6	0.41	0
3	SO4	D	402	-	4,4,4	0.74	0	6,6,6	0.41	0
3	SO4	G	402	-	4,4,4	0.71	0	6,6,6	0.49	0
2	WIU	F	401	-	39,39,39	1.64	4 (10%)	45,55,55	1.38	5 (11%)
2	WIU	A	401	-	39,39,39	1.53	5 (12%)	45,55,55	1.24	5 (11%)
2	WIU	E	401	-	39,39,39	1.66	4 (10%)	45,55,55	1.45	6 (13%)
3	SO4	H	402	-	4,4,4	0.77	0	6,6,6	0.23	0
3	SO4	E	402	-	4,4,4	0.73	0	6,6,6	0.19	0
2	WIU	C	401	-	39,39,39	1.38	3 (7%)	45,55,55	1.36	5 (11%)
3	SO4	F	402	-	4,4,4	0.81	0	6,6,6	0.54	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	WIU	B	401	-	-	2/27/27/27	0/4/4/4
2	WIU	D	401	-	-	4/27/27/27	0/4/4/4
2	WIU	H	401	-	-	4/27/27/27	0/4/4/4
2	WIU	G	401	-	-	3/27/27/27	0/4/4/4
2	WIU	F	401	-	-	2/27/27/27	0/4/4/4
2	WIU	A	401	-	-	3/27/27/27	0/4/4/4
2	WIU	E	401	-	-	2/27/27/27	0/4/4/4
2	WIU	C	401	-	-	2/27/27/27	0/4/4/4

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	WIU	N14-N15	8.44	1.43	1.36
2	F	401	WIU	N14-N15	8.15	1.43	1.36
2	D	401	WIU	N14-N15	8.12	1.43	1.36
2	A	401	WIU	N14-N15	7.23	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	401	WIU	N14-N15	6.59	1.42	1.36
2	C	401	WIU	N14-N15	6.42	1.41	1.36
2	B	401	WIU	N14-N15	6.40	1.41	1.36
2	G	401	WIU	N14-N15	5.87	1.41	1.36
2	B	401	WIU	C16-N15	-3.03	1.30	1.32
2	D	401	WIU	C16-N15	-3.03	1.30	1.32
2	H	401	WIU	C16-N15	-3.00	1.30	1.32
2	A	401	WIU	C16-N15	-2.93	1.30	1.32
2	C	401	WIU	C16-N15	-2.64	1.30	1.32
2	A	401	WIU	C09-N36	-2.53	1.30	1.33
2	F	401	WIU	C16-N15	-2.44	1.30	1.32
2	D	401	WIU	C12-C16	2.40	1.44	1.42
2	F	401	WIU	C09-N36	-2.37	1.30	1.33
2	C	401	WIU	C12-C16	2.36	1.44	1.42
2	G	401	WIU	C09-N36	-2.33	1.30	1.33
2	A	401	WIU	C12-C16	2.27	1.44	1.42
2	D	401	WIU	C09-N36	-2.27	1.30	1.33
2	E	401	WIU	C16-N15	-2.24	1.30	1.32
2	B	401	WIU	C09-N36	-2.20	1.30	1.33
2	E	401	WIU	C12-C16	2.18	1.44	1.42
2	D	401	WIU	C12-C11	-2.14	1.39	1.42
2	F	401	WIU	C12-C11	-2.04	1.40	1.42
2	E	401	WIU	C12-C11	-2.02	1.40	1.42
2	A	401	WIU	C12-C11	-2.00	1.40	1.42
2	D	401	WIU	C18-N19	2.00	1.37	1.34

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	WIU	C01-N02-C03	4.88	136.14	128.80
2	B	401	WIU	C01-N02-C03	4.82	136.04	128.80
2	H	401	WIU	C01-N02-C03	4.64	135.78	128.80
2	E	401	WIU	C01-N02-C03	4.59	135.70	128.80
2	C	401	WIU	C01-N02-C03	4.58	135.68	128.80
2	D	401	WIU	C01-N02-C03	4.43	135.46	128.80
2	G	401	WIU	C01-N02-C03	4.39	135.41	128.80
2	A	401	WIU	C01-N02-C03	4.29	135.26	128.80
2	F	401	WIU	C01-N02-N36	-3.76	113.66	119.29
2	C	401	WIU	C01-N02-N36	-3.52	114.03	119.29
2	E	401	WIU	C01-N02-N36	-3.51	114.04	119.29
2	G	401	WIU	C13-N14-N15	-3.29	110.49	111.26
2	E	401	WIU	C21-C20-N19	-3.25	109.25	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	WIU	C01-N02-N36	-3.25	114.43	119.29
2	H	401	WIU	C01-N02-N36	-3.23	114.46	119.29
2	G	401	WIU	C21-C20-N19	-3.19	109.32	113.50
2	G	401	WIU	C01-N02-N36	-3.18	114.53	119.29
2	A	401	WIU	C01-N02-N36	-3.11	114.63	119.29
2	D	401	WIU	C01-N02-N36	-3.11	114.64	119.29
2	C	401	WIU	C21-C20-N19	-2.87	109.74	113.50
2	E	401	WIU	C08-C09-N36	-2.86	110.29	112.34
2	B	401	WIU	C03-N02-N36	-2.77	109.53	112.48
2	E	401	WIU	C13-N14-N15	-2.71	110.63	111.26
2	H	401	WIU	C08-C09-N36	-2.69	110.41	112.34
2	D	401	WIU	C13-N14-N15	-2.62	110.65	111.26
2	H	401	WIU	C21-C20-N19	-2.55	110.16	113.50
2	H	401	WIU	C03-N02-N36	-2.55	109.76	112.48
2	G	401	WIU	C12-C13-N14	2.53	108.31	107.21
2	F	401	WIU	C21-C20-N19	-2.45	110.29	113.50
2	D	401	WIU	C08-C09-N36	-2.45	110.59	112.34
2	D	401	WIU	C03-N02-N36	-2.42	109.89	112.48
2	C	401	WIU	C08-C09-N36	-2.30	110.69	112.34
2	A	401	WIU	C21-C20-N19	-2.28	110.51	113.50
2	G	401	WIU	C03-N02-N36	-2.27	110.06	112.48
2	A	401	WIU	C08-C09-N36	-2.24	110.74	112.34
2	A	401	WIU	C03-N02-N36	-2.23	110.10	112.48
2	B	401	WIU	C08-C09-N36	-2.19	110.77	112.34
2	G	401	WIU	C08-C09-N36	-2.16	110.80	112.34
2	G	401	WIU	C31-N30-C29	-2.15	118.33	121.98
2	F	401	WIU	C03-N02-N36	-2.14	110.20	112.48
2	E	401	WIU	C03-N02-N36	-2.09	110.25	112.48
2	C	401	WIU	C03-N02-N36	-2.05	110.29	112.48
2	F	401	WIU	C08-C09-N36	-2.04	110.88	112.34
2	B	401	WIU	C13-N14-N15	-2.03	110.79	111.26

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	401	WIU	O28-C29-N30-C31
2	D	401	WIU	O34-C29-N30-C31
2	B	401	WIU	N30-C29-O28-C27
2	H	401	WIU	N30-C29-O28-C27
2	B	401	WIU	O34-C29-O28-C27
2	E	401	WIU	N30-C29-O28-C27

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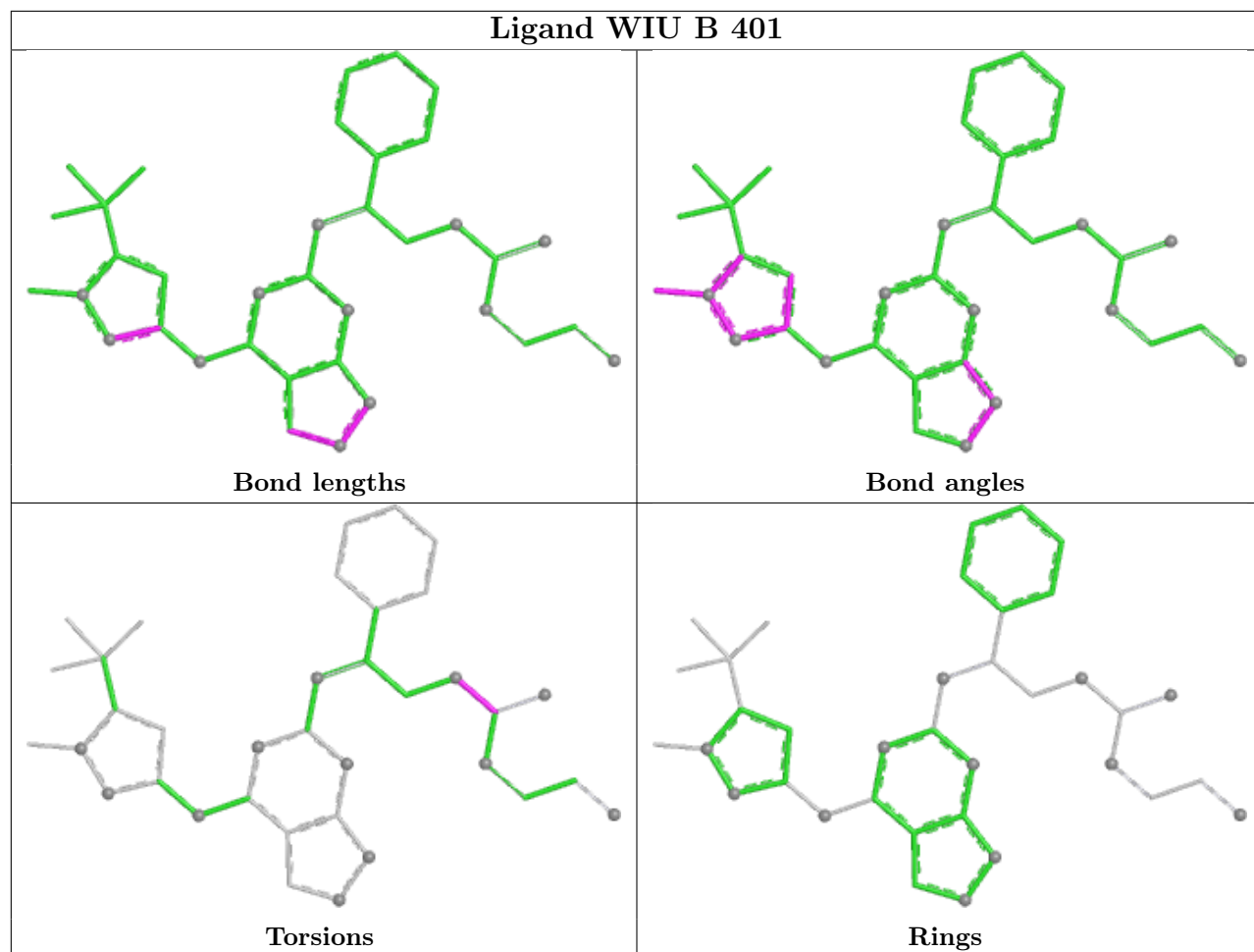
Mol	Chain	Res	Type	Atoms
2	H	401	WIU	O34-C29-O28-C27
2	F	401	WIU	N30-C29-O28-C27
2	G	401	WIU	O28-C29-N30-C31
2	E	401	WIU	O34-C29-O28-C27
2	A	401	WIU	N30-C29-O28-C27
2	G	401	WIU	N30-C29-O28-C27
2	H	401	WIU	N19-C20-C21-C26
2	G	401	WIU	O34-C29-O28-C27
2	A	401	WIU	O34-C29-O28-C27
2	F	401	WIU	O34-C29-O28-C27
2	D	401	WIU	N30-C29-O28-C27
2	C	401	WIU	N30-C29-O28-C27
2	H	401	WIU	N19-C20-C21-C22
2	A	401	WIU	O28-C29-N30-C31
2	C	401	WIU	O34-C29-O28-C27
2	D	401	WIU	O34-C29-O28-C27

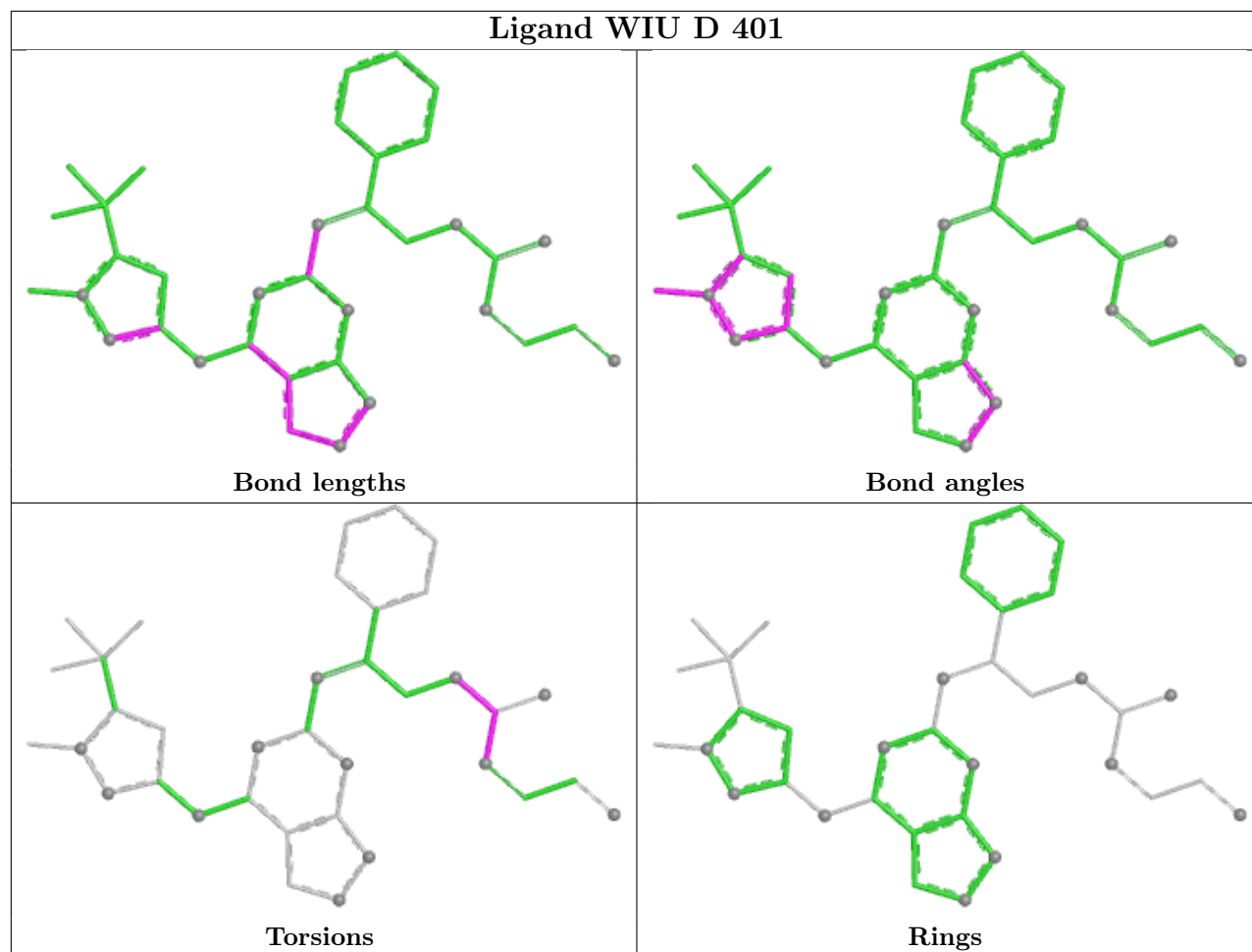
There are no ring outliers.

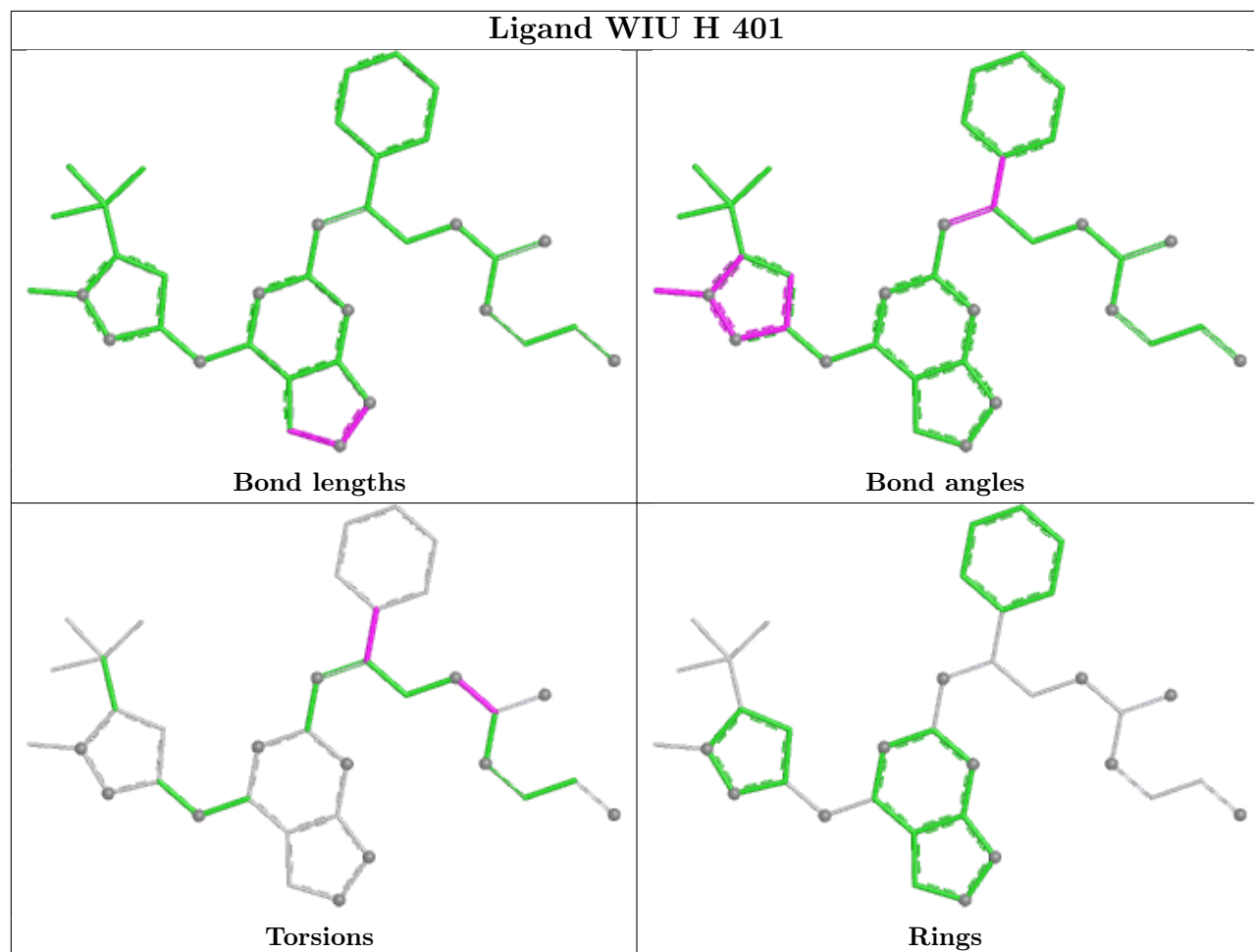
2 monomers are involved in 2 short contacts:

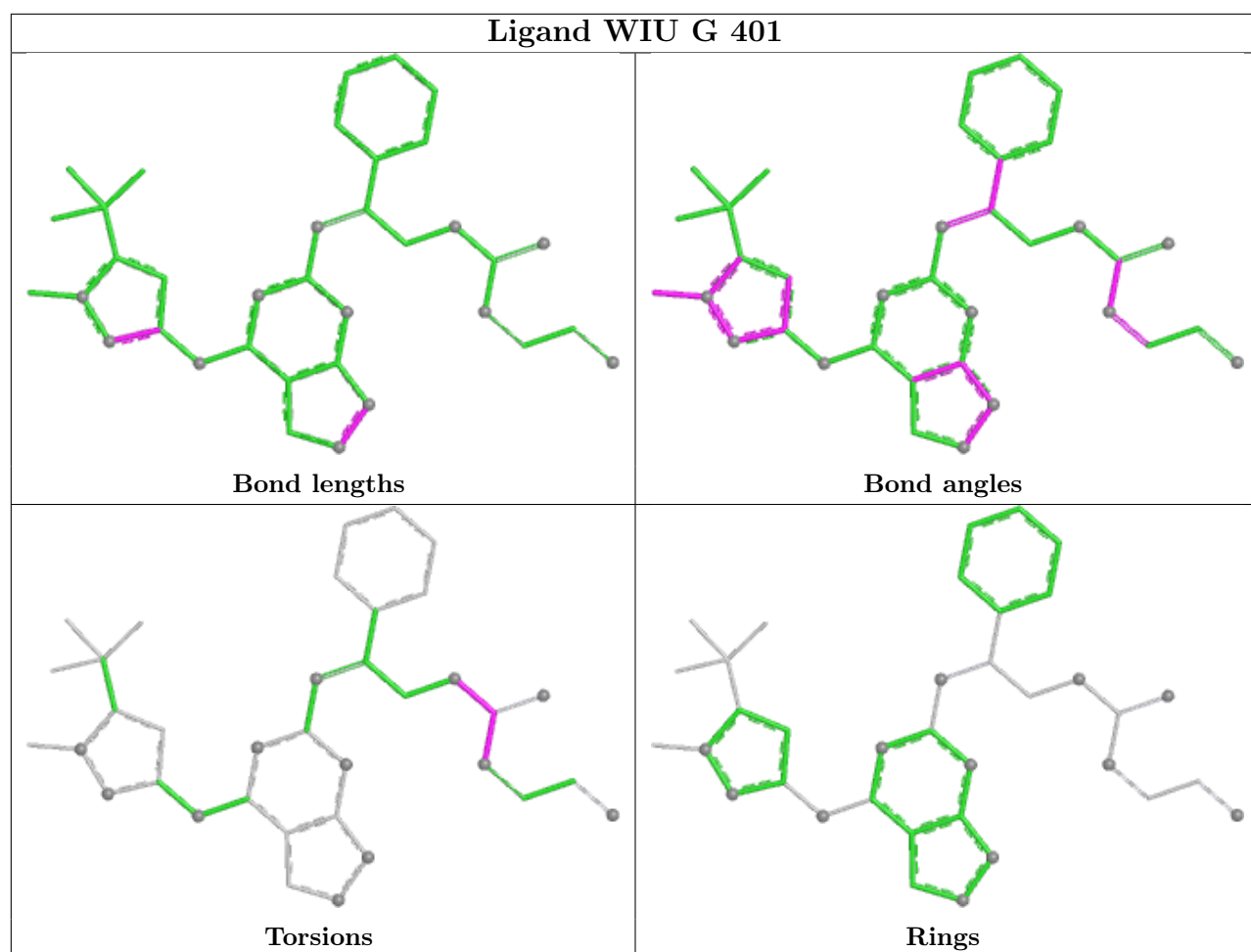
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	WIU	1	0
2	E	401	WIU	1	0

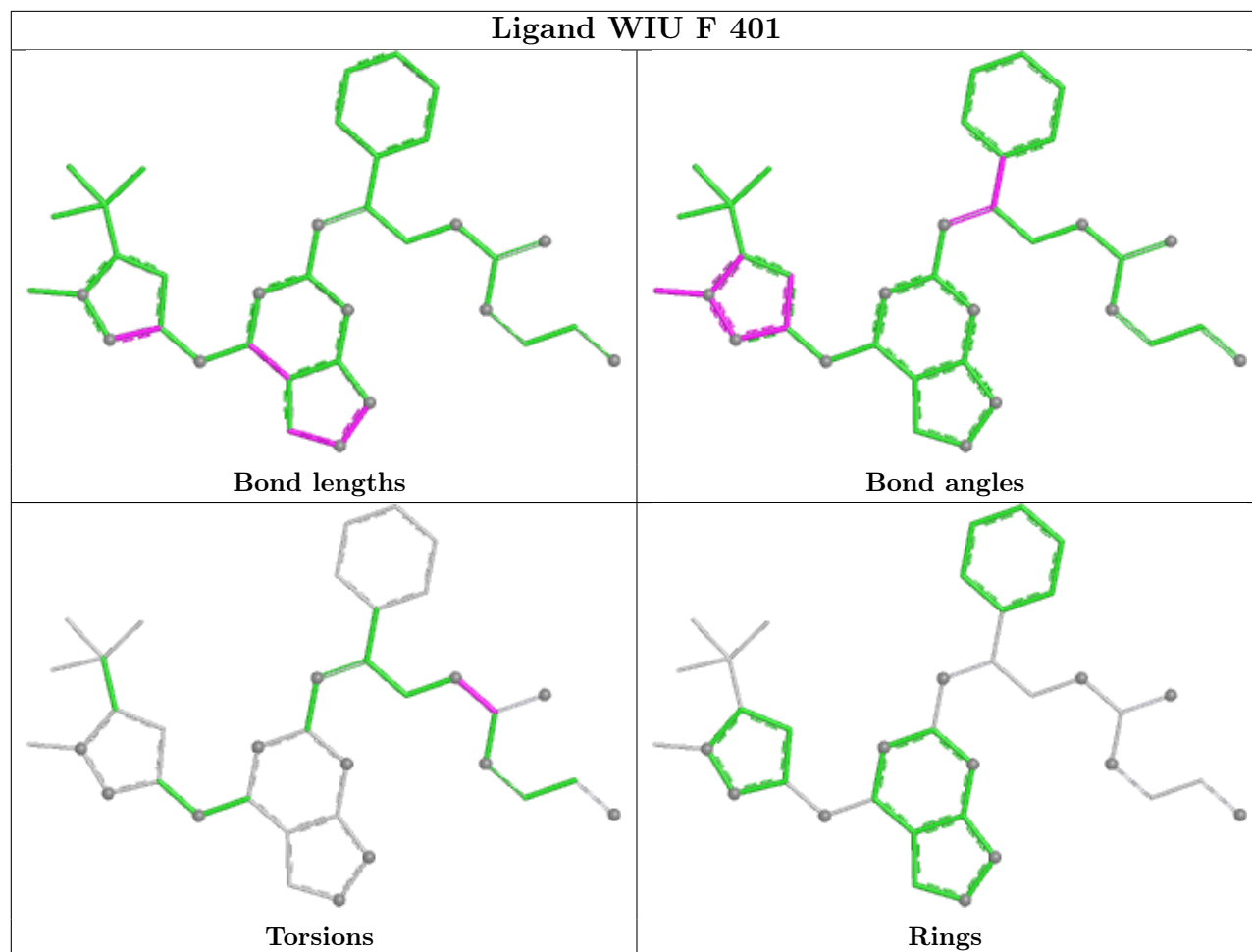
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

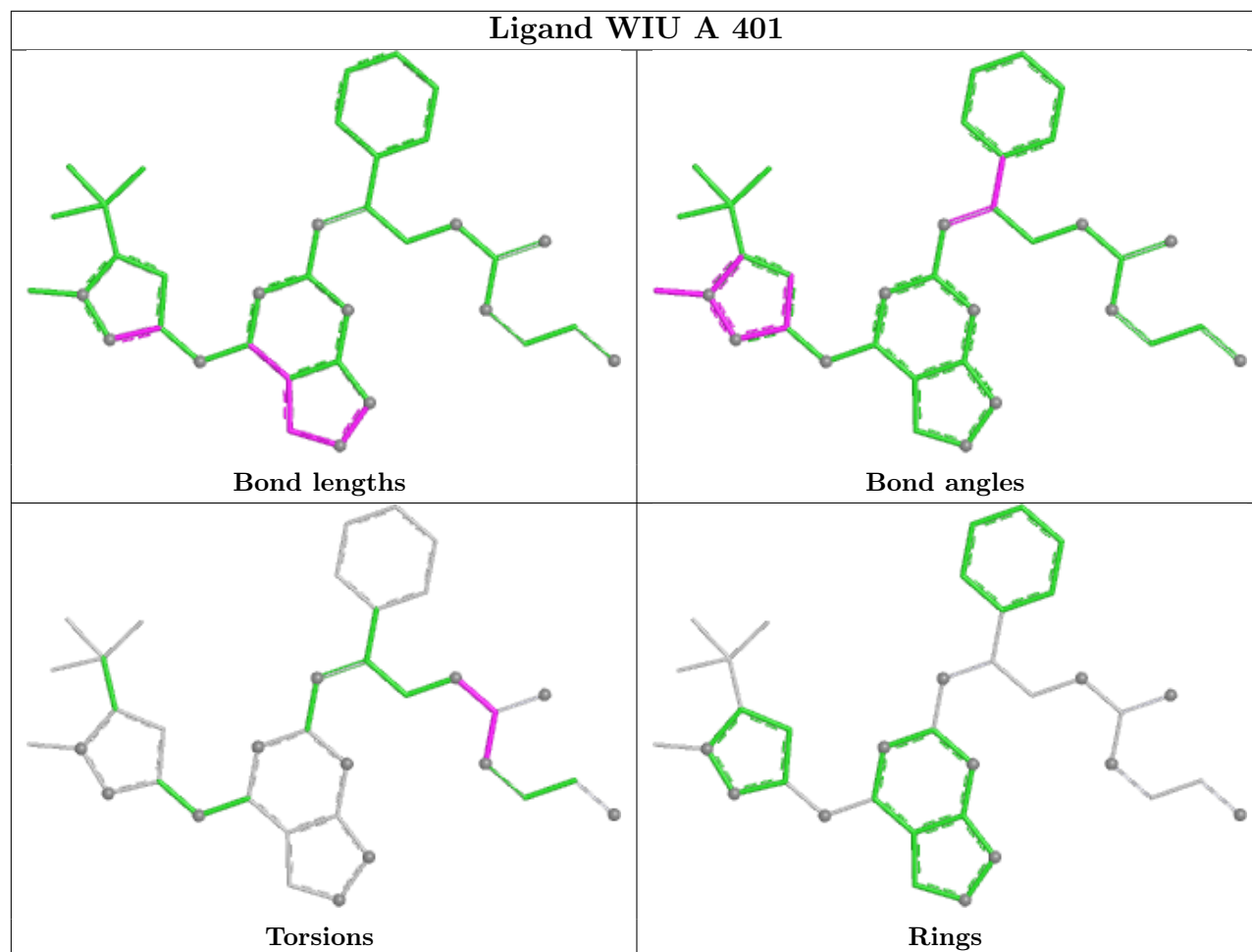




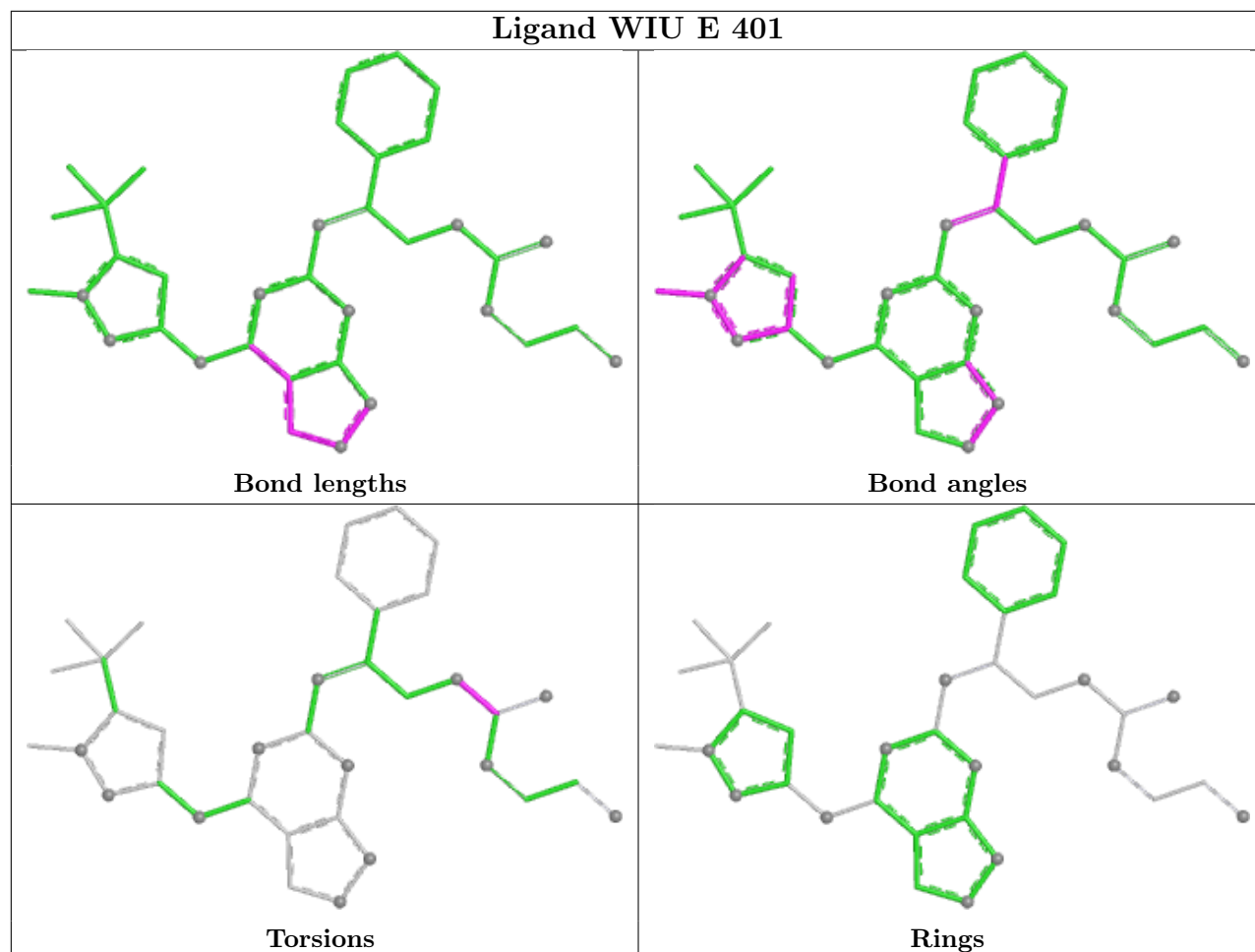


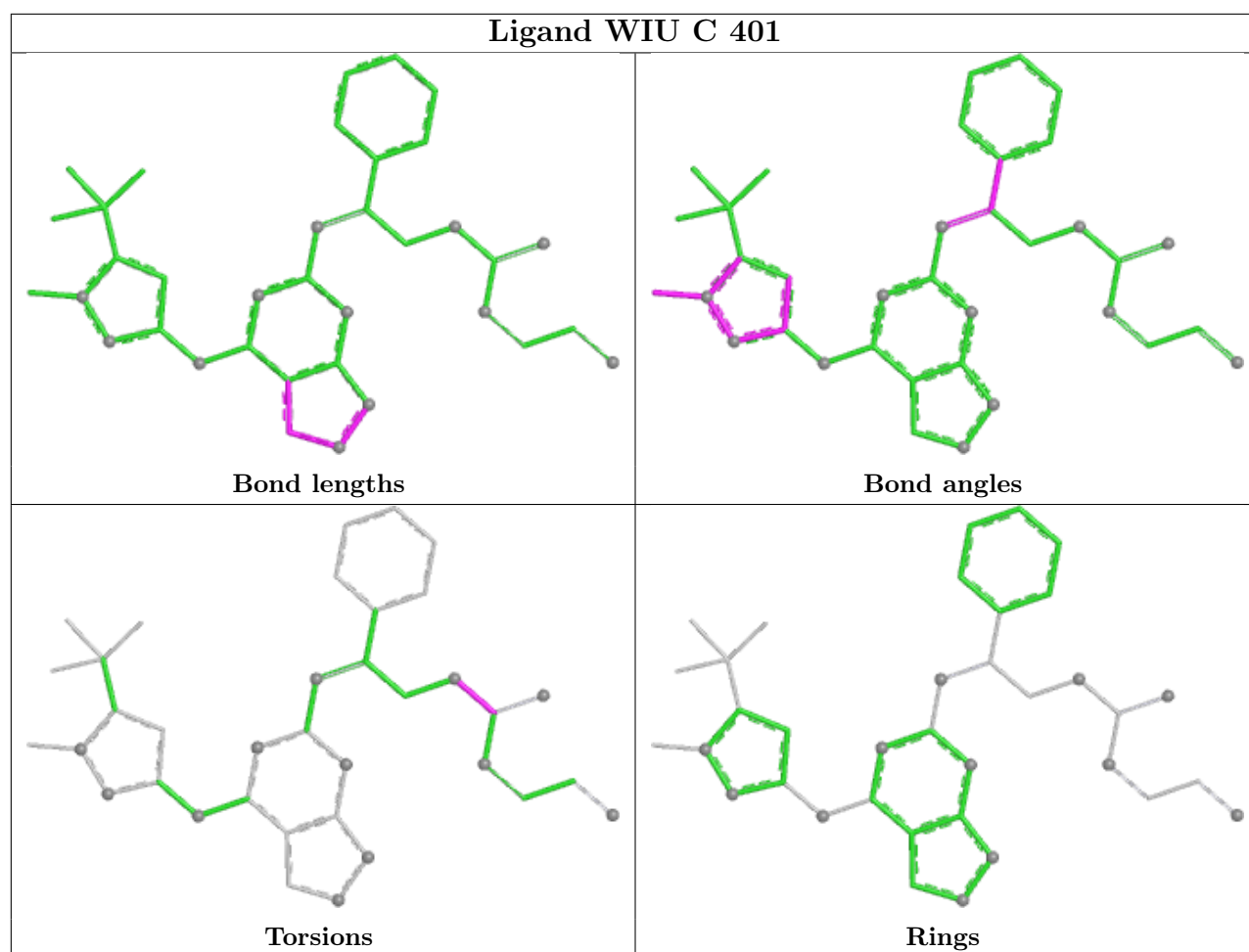






Ligand WIU E 401





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	303/316 (95%)	-0.08	11 (3%)	46	43	21, 42, 72, 99	6 (1%)
1	B	303/316 (95%)	-0.15	13 (4%)	40	36	19, 40, 71, 95	7 (2%)
1	C	303/316 (95%)	-0.10	14 (4%)	37	34	20, 42, 71, 101	3 (0%)
1	D	303/316 (95%)	-0.10	14 (4%)	37	34	21, 42, 76, 104	8 (2%)
1	E	303/316 (95%)	-0.09	14 (4%)	37	34	20, 41, 70, 106	5 (1%)
1	F	303/316 (95%)	-0.15	12 (3%)	42	39	19, 39, 65, 98	8 (2%)
1	G	303/316 (95%)	-0.09	12 (3%)	42	39	21, 42, 73, 102	6 (1%)
1	H	303/316 (95%)	-0.12	14 (4%)	37	34	18, 40, 73, 97	6 (1%)
All	All	2424/2528 (95%)	-0.11	104 (4%)	40	36	18, 41, 72, 106	49 (2%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	83	ALA	6.4
1	C	83	ALA	6.2
1	H	83	ALA	5.7
1	B	83	ALA	5.6
1	A	86	ARG	5.5
1	G	155	ALA	5.4
1	F	83	ALA	5.3
1	E	155	ALA	5.1
1	E	158	ASN	5.1
1	A	83	ALA	5.0
1	F	155	ALA	4.9
1	A	155	ALA	4.9
1	H	155	ALA	4.8
1	E	83	ALA	4.7
1	A	319	GLY	4.7
1	G	83	ALA	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	319	GLY	4.6
1	D	319	GLY	4.5
1	E	319	GLY	4.4
1	D	155	ALA	4.4
1	C	86	ARG	4.3
1	F	319	GLY	4.3
1	H	158	ASN	4.3
1	B	158	ASN	4.2
1	D	156	GLY	4.2
1	B	155	ALA	4.2
1	E	156	GLY	4.1
1	F	15	MET	4.1
1	G	264	MET	4.1
1	H	319	GLY	4.1
1	G	319	GLY	4.1
1	C	155	ALA	4.0
1	H	156	GLY	4.0
1	A	158	ASN	4.0
1	F	158	ASN	3.9
1	F	26	ALA	3.9
1	G	158	ASN	3.8
1	G	15	MET	3.6
1	B	82	SER	3.6
1	C	156	GLY	3.6
1	D	86	ARG	3.5
1	A	87	ALA	3.5
1	E	86	ARG	3.5
1	B	319	GLY	3.5
1	F	82	SER	3.4
1	B	156	GLY	3.3
1	H	15	MET	3.3
1	B	15	MET	3.3
1	E	15	MET	3.2
1	H	86	ARG	3.2
1	A	156	GLY	3.2
1	G	156	GLY	3.2
1	G	86	ARG	3.1
1	G	87	ALA	3.1
1	F	156	GLY	3.0
1	B	87	ALA	3.0
1	H	87	ALA	3.0
1	D	15	MET	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	262	GLU	3.0
1	C	157	THR	3.0
1	B	86	ARG	3.0
1	E	157	THR	3.0
1	B	81	SER	3.0
1	A	15	MET	2.9
1	A	82	SER	2.9
1	H	264	MET	2.9
1	E	82	SER	2.8
1	E	318	GLN	2.8
1	C	158	ASN	2.8
1	C	82	SER	2.8
1	F	264	MET	2.7
1	D	81	SER	2.7
1	H	82	SER	2.7
1	D	264	MET	2.6
1	H	318	GLN	2.6
1	A	264	MET	2.6
1	C	16	ARG	2.6
1	G	82	SER	2.6
1	A	262	GLU	2.5
1	H	262	GLU	2.5
1	C	15	MET	2.5
1	D	157	THR	2.5
1	D	158	ASN	2.4
1	C	87	ALA	2.4
1	F	86	ARG	2.4
1	H	81	SER	2.3
1	E	27	GLY	2.3
1	B	262	GLU	2.3
1	C	264	MET	2.3
1	D	26	ALA	2.3
1	F	87	ALA	2.3
1	B	264	MET	2.2
1	C	262	GLU	2.2
1	C	318	GLN	2.2
1	H	157	THR	2.2
1	E	264	MET	2.2
1	G	16	ARG	2.2
1	E	16	ARG	2.2
1	B	16	ARG	2.1
1	F	16	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	318	GLN	2.1
1	D	82	SER	2.1
1	G	157	THR	2.1
1	D	87	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

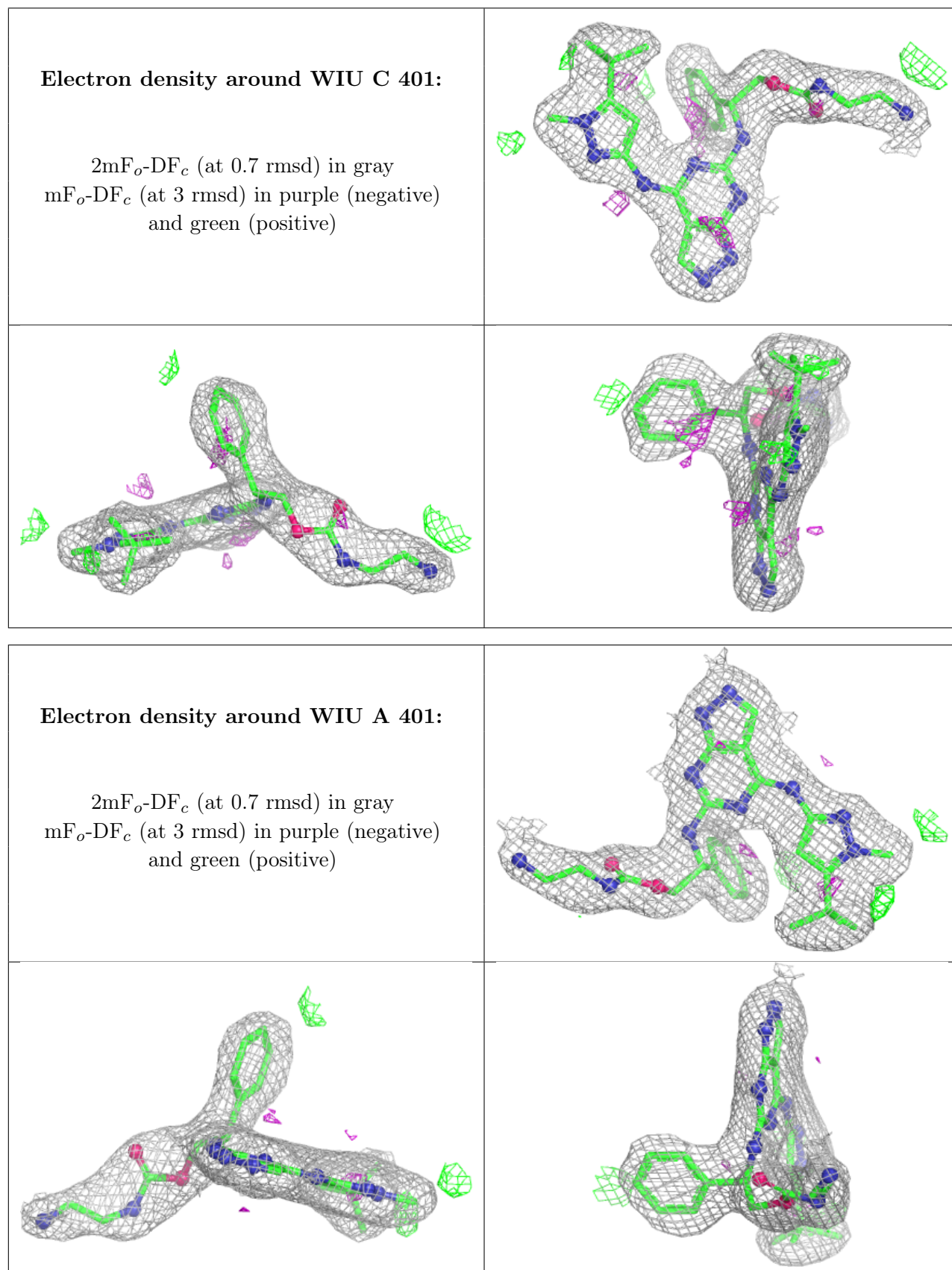
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	WIU	C	401	36/36	0.92	0.10	33,41,49,57	0
2	WIU	A	401	36/36	0.93	0.10	35,42,50,56	0
2	WIU	D	401	36/36	0.94	0.09	33,40,49,54	0
2	WIU	E	401	36/36	0.94	0.09	31,40,46,54	0
2	WIU	F	401	36/36	0.94	0.09	31,39,48,57	0
2	WIU	H	401	36/36	0.94	0.09	31,37,43,54	0
2	WIU	G	401	36/36	0.95	0.08	31,40,48,53	0
2	WIU	B	401	36/36	0.95	0.08	29,36,44,49	0
3	SO4	C	402	5/5	0.96	0.08	44,44,54,54	0
3	SO4	F	402	5/5	0.96	0.08	34,39,50,50	0
3	SO4	A	402	5/5	0.97	0.09	42,44,50,50	0
3	SO4	G	402	5/5	0.97	0.09	37,38,45,50	0
3	SO4	E	402	5/5	0.98	0.07	38,42,48,50	0
3	SO4	B	402	5/5	0.98	0.06	32,34,44,45	0
3	SO4	D	402	5/5	0.98	0.06	37,38,47,50	0
3	SO4	H	402	5/5	0.98	0.07	35,40,44,48	0

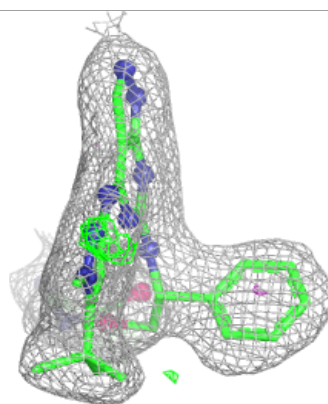
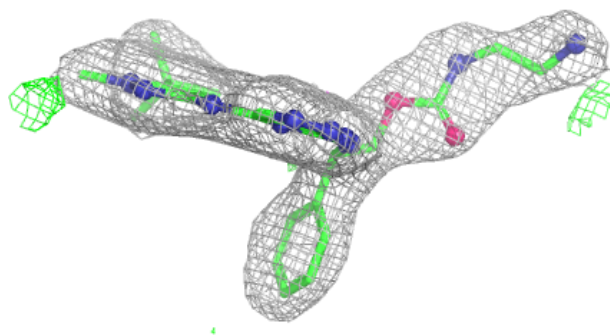
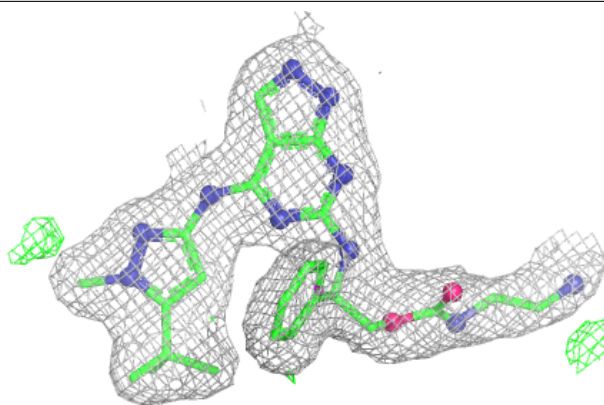
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

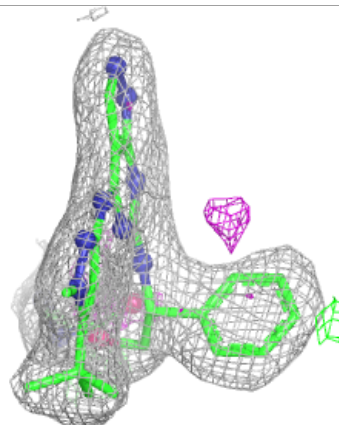
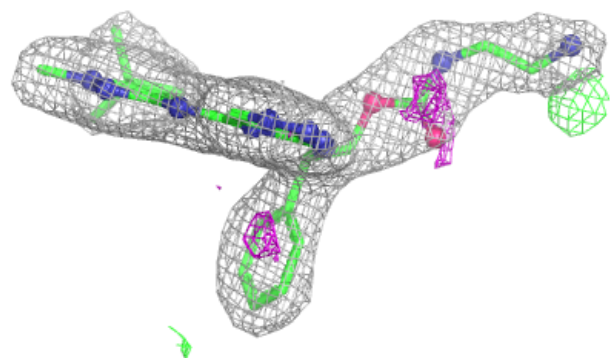
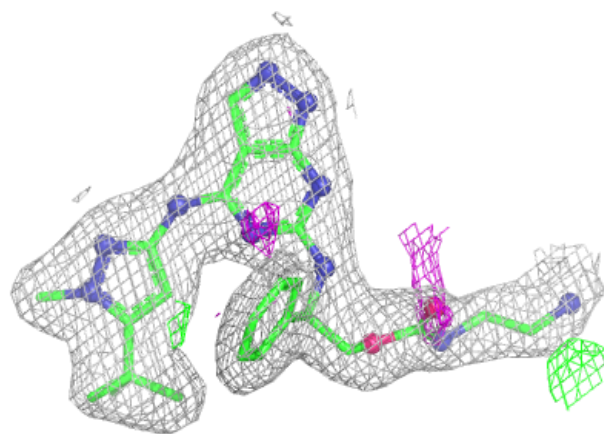


Electron density around WIU D 401:

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

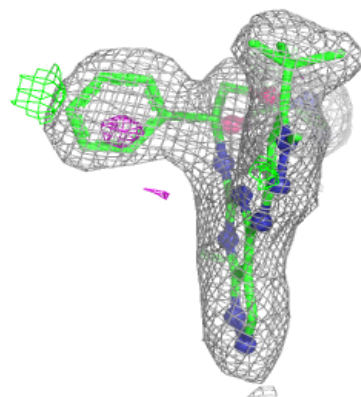
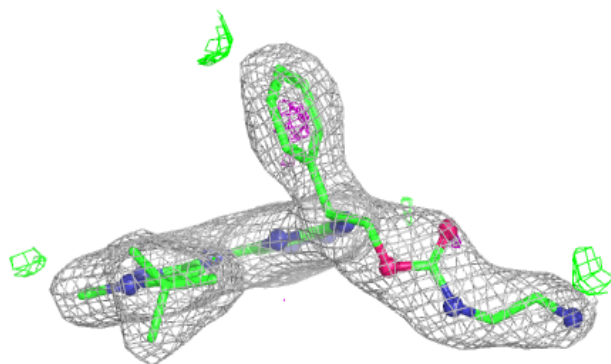
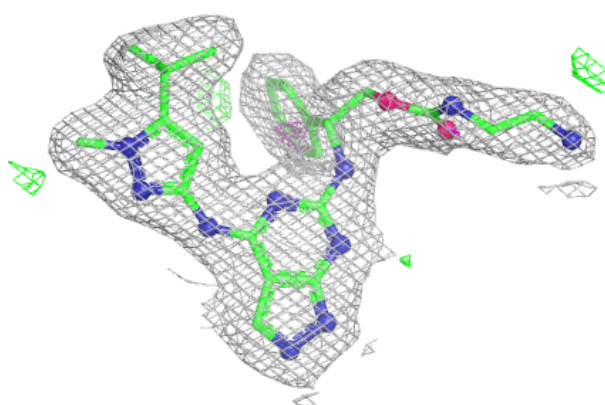
**Electron density around WIU E 401:**

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

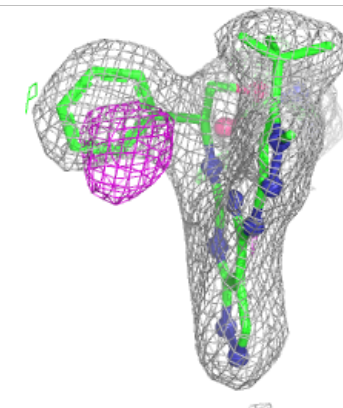
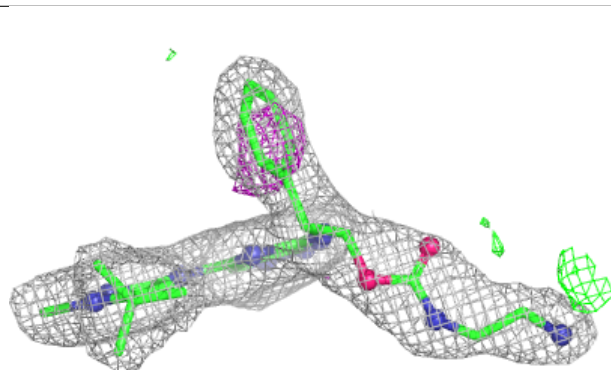
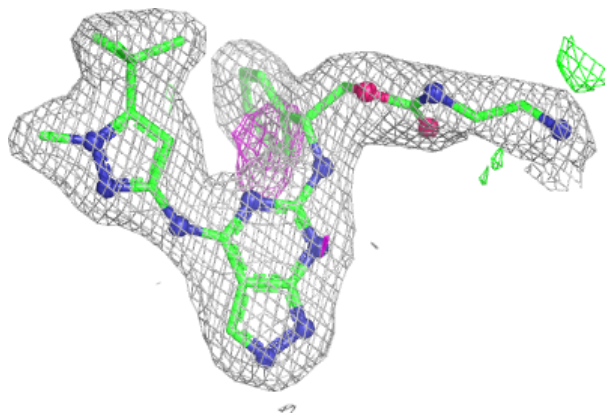


Electron density around WIU F 401:

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

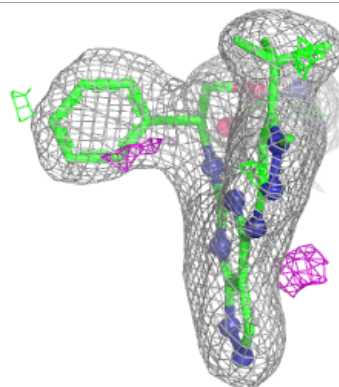
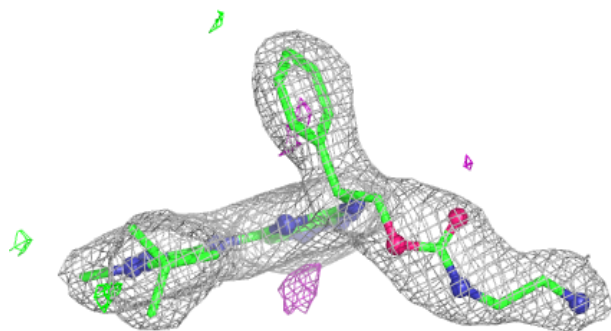
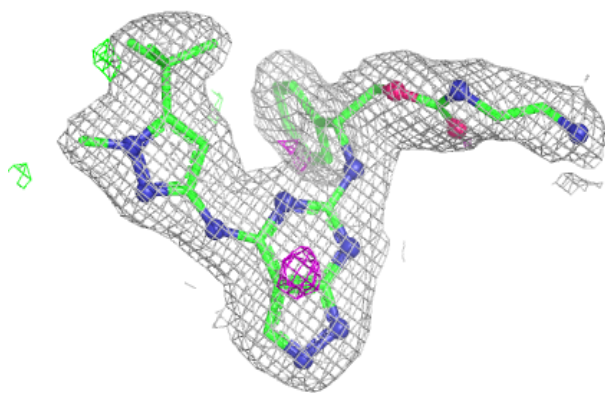
**Electron density around WIU H 401:**

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

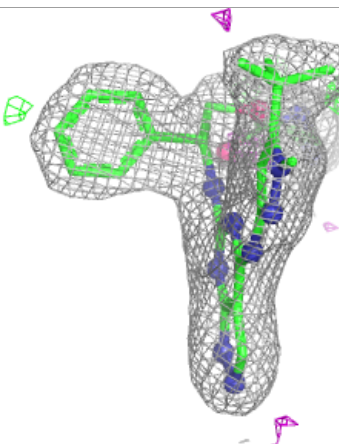
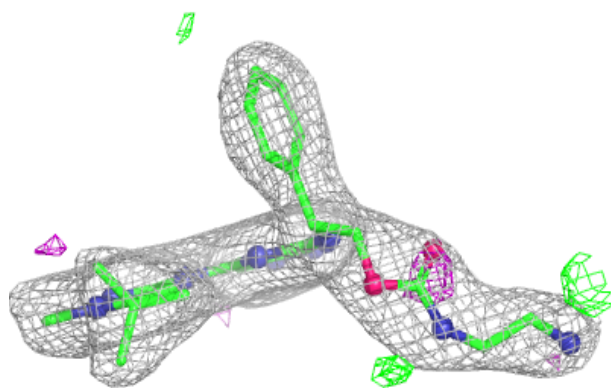
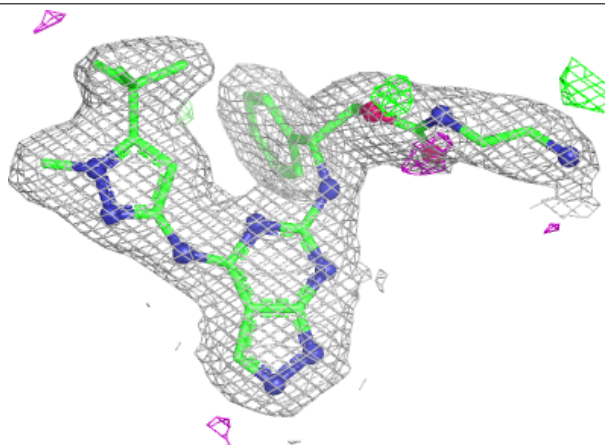


Electron density around WIU G 401:

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

**Electron density around WIU 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.