



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:41 PM UTC

PDB ID : 6E43 / pdb_00006e43
Title : Crystal structure of human indoleamine 2,3-dioxygenase 1 (IDO1) in complex with a BMS-978587 analog
Authors : Luo, S.; Tong, L.
Deposited on : 2018-07-16
Resolution : 1.71 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

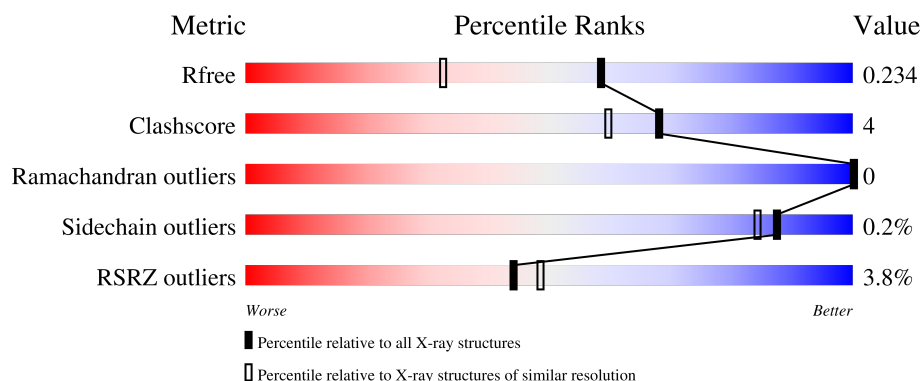
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	405	3% 82% 10% 8%
1	B	405	5% 86% 6% 8%
1	C	405	3% 87% 5% 7%
1	D	405	3% 85% 7% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13245 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 Indoleamine 2,3-dioxygenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	2	0
			2943	1893	501	530	19			
1	B	373	Total	C	N	O	S	0	3	0
			2951	1898	502	531	20			
1	C	375	Total	C	N	O	S	0	3	0
			2964	1905	504	536	19			
1	D	375	Total	C	N	O	S	0	2	0
			2962	1903	504	536	19			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P14902
A	0	GLY	-	expression tag	UNP P14902
A	1	SER	-	expression tag	UNP P14902
A	2	SER	-	expression tag	UNP P14902
A	3	HIS	-	expression tag	UNP P14902
A	4	HIS	-	expression tag	UNP P14902
A	5	HIS	-	expression tag	UNP P14902
A	6	HIS	-	expression tag	UNP P14902
A	7	HIS	-	expression tag	UNP P14902
A	8	HIS	-	expression tag	UNP P14902
A	9	SER	-	expression tag	UNP P14902
A	10	SER	-	expression tag	UNP P14902
A	11	GLY	-	expression tag	UNP P14902
A	12	SER	-	expression tag	UNP P14902
A	13	ALA	-	expression tag	UNP P14902
A	14	ALA	-	expression tag	UNP P14902
A	116	ALA	LYS	engineered mutation	UNP P14902
A	117	ALA	LYS	engineered mutation	UNP P14902
B	-1	MET	-	initiating methionine	UNP P14902
B	0	GLY	-	expression tag	UNP P14902
B	1	SER	-	expression tag	UNP P14902

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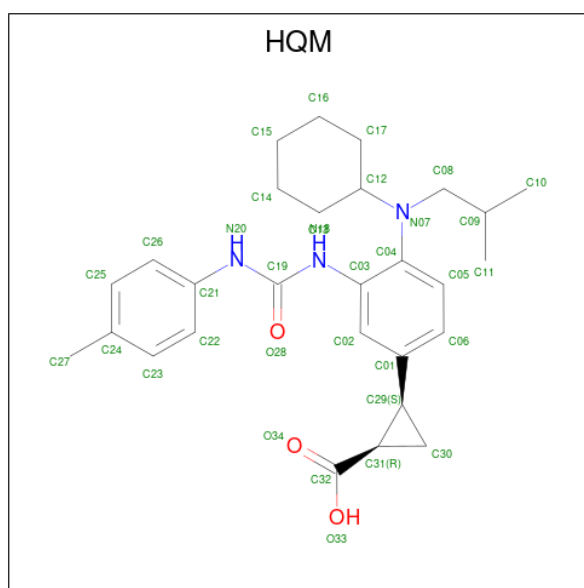
Chain	Residue	Modelled	Actual	Comment	Reference
B	2	SER	-	expression tag	UNP P14902
B	3	HIS	-	expression tag	UNP P14902
B	4	HIS	-	expression tag	UNP P14902
B	5	HIS	-	expression tag	UNP P14902
B	6	HIS	-	expression tag	UNP P14902
B	7	HIS	-	expression tag	UNP P14902
B	8	HIS	-	expression tag	UNP P14902
B	9	SER	-	expression tag	UNP P14902
B	10	SER	-	expression tag	UNP P14902
B	11	GLY	-	expression tag	UNP P14902
B	12	SER	-	expression tag	UNP P14902
B	13	ALA	-	expression tag	UNP P14902
B	14	ALA	-	expression tag	UNP P14902
B	116	ALA	LYS	engineered mutation	UNP P14902
B	117	ALA	LYS	engineered mutation	UNP P14902
C	-1	MET	-	initiating methionine	UNP P14902
C	0	GLY	-	expression tag	UNP P14902
C	1	SER	-	expression tag	UNP P14902
C	2	SER	-	expression tag	UNP P14902
C	3	HIS	-	expression tag	UNP P14902
C	4	HIS	-	expression tag	UNP P14902
C	5	HIS	-	expression tag	UNP P14902
C	6	HIS	-	expression tag	UNP P14902
C	7	HIS	-	expression tag	UNP P14902
C	8	HIS	-	expression tag	UNP P14902
C	9	SER	-	expression tag	UNP P14902
C	10	SER	-	expression tag	UNP P14902
C	11	GLY	-	expression tag	UNP P14902
C	12	SER	-	expression tag	UNP P14902
C	13	ALA	-	expression tag	UNP P14902
C	14	ALA	-	expression tag	UNP P14902
C	116	ALA	LYS	engineered mutation	UNP P14902
C	117	ALA	LYS	engineered mutation	UNP P14902
D	-1	MET	-	initiating methionine	UNP P14902
D	0	GLY	-	expression tag	UNP P14902
D	1	SER	-	expression tag	UNP P14902
D	2	SER	-	expression tag	UNP P14902
D	3	HIS	-	expression tag	UNP P14902
D	4	HIS	-	expression tag	UNP P14902
D	5	HIS	-	expression tag	UNP P14902
D	6	HIS	-	expression tag	UNP P14902
D	7	HIS	-	expression tag	UNP P14902

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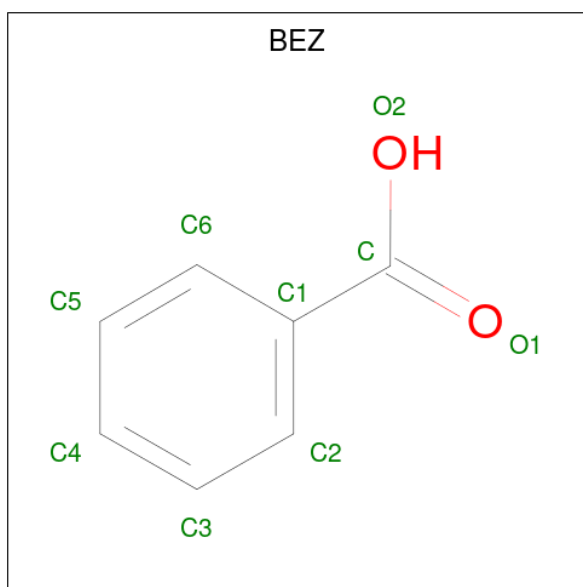
Chain	Residue	Modelled	Actual	Comment	Reference
D	8	HIS	-	expression tag	UNP P14902
D	9	SER	-	expression tag	UNP P14902
D	10	SER	-	expression tag	UNP P14902
D	11	GLY	-	expression tag	UNP P14902
D	12	SER	-	expression tag	UNP P14902
D	13	ALA	-	expression tag	UNP P14902
D	14	ALA	-	expression tag	UNP P14902
D	116	ALA	LYS	engineered mutation	UNP P14902
D	117	ALA	LYS	engineered mutation	UNP P14902

- Molecule 2 is (1R,2S)-2-(4-[cyclohexyl(2-methylpropyl)amino]-3-[(4-methylphenyl)carbamoyl]amino}phenyl)cyclopropane-1-carboxylic acid (CCD ID: HQM) (formula: $C_{28}H_{37}N_3O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			34	28	3	3		
2	B	1	Total	C	N	O	0	0
			34	28	3	3		
2	C	1	Total	C	N	O	0	0
			34	28	3	3		
2	D	1	Total	C	N	O	0	0
			34	28	3	3		

- Molecule 3 is BENZOIC ACID (CCD ID: BEZ) (formula: $C_7H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			9	7	2		
3	B	1	Total	C	O	0	0
			9	7	2		
3	C	1	Total	C	O	0	0
			9	7	2		
3	D	1	Total	C	O	0	0
			9	7	2		

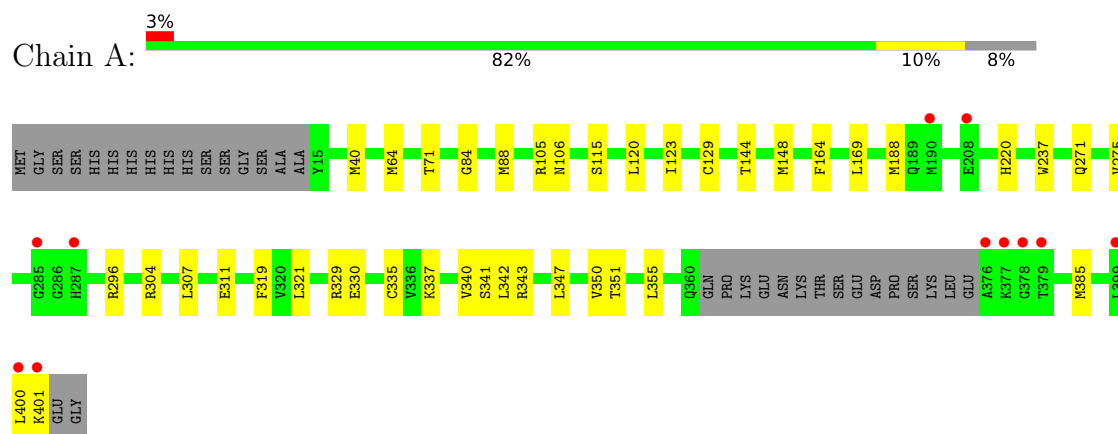
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	272	Total	O	0	0
			272	272		
4	B	271	Total	O	0	0
			271	271		
4	C	338	Total	O	0	0
			338	338		
4	D	372	Total	O	0	0
			372	372		

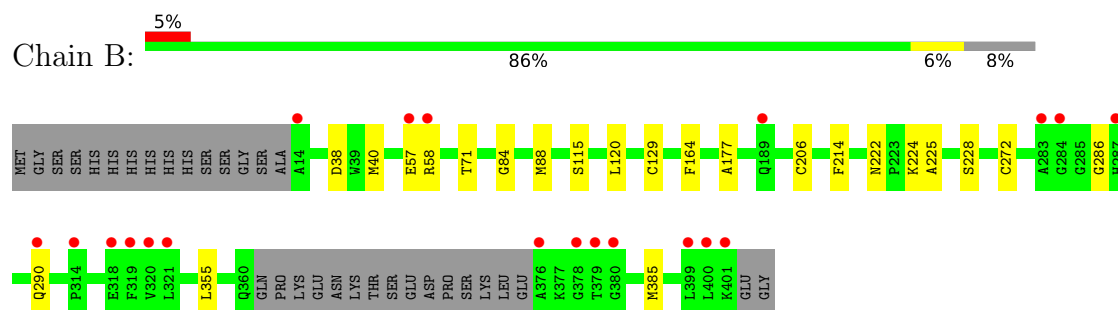
3 Residue-property plots [i](#)

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.

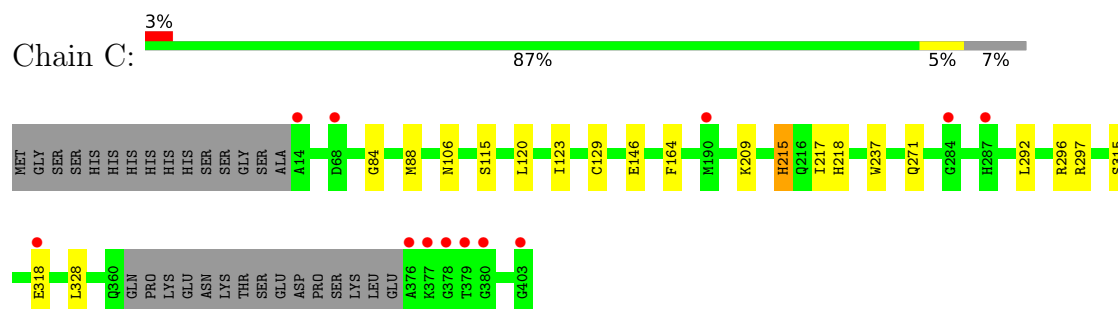
- Molecule 1: Indoleamine 2,3-dioxygenase 1



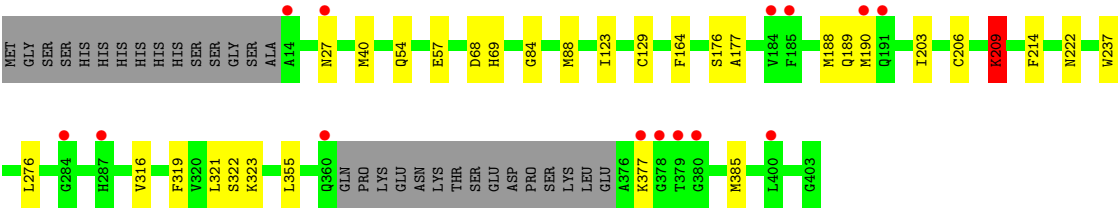
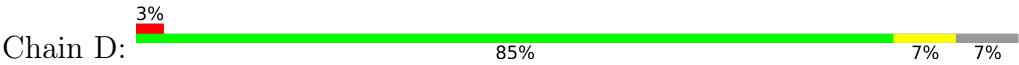
- Molecule 1: Indoleamine 2,3-dioxygenase 1



- Molecule 1: Indoleamine 2,3-dioxygenase 1



- Molecule 1: Indoleamine 2,3-dioxygenase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.48Å 195.39Å 118.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.64 – 1.71 47.64 – 1.71	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.64-1.71) 99.1 (47.64-1.71)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 1.71Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.194 , 0.231 0.200 , 0.234	Depositor DCC
R_{free} test set	3011 reflections (1.48%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13245	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 44.22 % 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 1.5745e-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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HQM, BEZ

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	0.43	6/3017 (0.2%)	0.49	0/4082
1	B	0.27	0/3028	0.47	0/4097
1	C	0.46	4/3041 (0.1%)	0.51	0/4114
1	D	0.33	1/3036 (0.0%)	0.51	0/4106
All	All	0.38	11/12122 (0.1%)	0.50	0/16399

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	217	ILE	C-O	-6.95	1.16	1.24
1	C	218	HIS	C-O	-6.39	1.16	1.24
1	C	215	HIS	C-O	-6.16	1.16	1.24
1	A	343	ARG	C-O	-6.13	1.17	1.24
1	A	337	LYS	C-O	-5.92	1.17	1.24
1	A	341	SER	C-O	-5.69	1.17	1.24
1	A	340	VAL	C-O	-5.54	1.17	1.24
1	C	292	LEU	C-O	-5.45	1.17	1.24
1	A	335	CYS	C-O	-5.42	1.17	1.24
1	A	342	LEU	C-O	-5.38	1.17	1.24
1	D	209	LYS	C-O	-5.29	1.18	1.24

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2943	0	2961	32	0
1	B	2951	0	2971	18	0
1	C	2964	0	2980	15	0
1	D	2962	0	2975	23	0
2	A	34	0	0	0	0
2	B	34	0	0	0	0
2	C	34	0	0	0	0
2	D	34	0	0	0	0
3	A	9	0	5	0	0
3	B	9	0	5	1	0
3	C	9	0	5	0	0
3	D	9	0	5	1	0
4	A	272	0	0	6	0
4	B	271	0	0	7	0
4	C	338	0	0	3	1
4	D	372	0	0	3	1
All	All	13245	0	11907	85	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 (85) 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:88[A]:MET:SD	4:B:715:HOH:O	2.35	0.83
1:B:222:ASN:HD22	1:B:225:ALA:H	1.26	0.83
1:B:58:ARG:NH1	4:B:601:HOH:O	2.15	0.79
1:D:189:GLN:NE2	1:D:190:MET:SD	2.57	0.78
1:C:297:ARG:NH1	4:C:602:HOH:O	2.25	0.69
1:A:271:GLN:HE22	1:A:296:ARG:HE	1.41	0.67
1:A:350:VAL:HB	1:A:385:MET:HE2	1.76	0.67
1:A:329:ARG:HE	1:A:401:LYS:HE2	1.62	0.65
1:A:105:ARG:NH1	4:A:605:HOH:O	2.30	0.64
1:B:57:GLU:HG3	4:B:610:HOH:O	1.98	0.63
1:A:296:ARG:O	1:A:304:ARG:HD3	1.99	0.63
1:D:355:LEU:HD11	1:D:385:MET:HG3	1.82	0.62
1:C:129[A]:CYS:SG	1:C:164:PHE:CZ	2.94	0.60
1:A:350:VAL:HB	1:A:385:MET:CE	2.32	0.60
1:B:129[B]:CYS:SG	1:B:164:PHE:CZ	2.95	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:ASN:ND2	1:B:225:ALA:H	1.97	0.60
1:D:129[B]:CYS:SG	1:D:164:PHE:CZ	2.95	0.60
1:D:68:ASP:OD2	1:D:69:HIS:ND1	2.35	0.59
1:A:355:LEU:HD11	1:A:385:MET:HG3	1.84	0.59
1:B:286:GLY:O	1:B:290:GLN:HG3	2.04	0.58
1:A:88[B]:MET:HE1	1:A:123:ILE:HG13	1.86	0.58
1:C:271:GLN:HE22	1:C:296:ARG:HE	1.53	0.57
1:A:148:MET:HE1	1:A:169:LEU:HD21	1.87	0.57
1:A:330:GLU:OE2	4:A:601:HOH:O	2.18	0.56
1:B:71:THR:OG1	4:B:602:HOH:O	2.18	0.56
1:B:355:LEU:HD11	1:B:385:MET:HG3	1.90	0.54
1:A:88[B]:MET:HE1	1:A:237:TRP:CH2	2.43	0.54
1:C:88[B]:MET:HE1	1:C:123:ILE:HG13	1.91	0.52
1:C:146:GLU:HG2	1:D:222:ASN:HB2	1.90	0.52
1:B:224:LYS:HE3	1:B:228:SER:HB3	1.91	0.52
1:A:71:THR:HB	1:D:54:GLN:HG2	1.92	0.52
1:D:88[A]:MET:HE1	1:D:237:TRP:CH2	2.45	0.52
1:A:129[B]:CYS:SG	1:A:164:PHE:CZ	3.03	0.51
1:D:88[A]:MET:HE1	1:D:123:ILE:HG13	1.93	0.51
1:C:209:LYS:HA	1:C:209:LYS:HE2	1.93	0.51
1:A:220:HIS:HE1	4:A:718:HOH:O	1.92	0.51
1:A:271:GLN:NE2	1:A:296:ARG:HE	2.07	0.50
1:D:84:GLY:O	1:D:88[B]:MET:HG2	2.11	0.50
1:B:84:GLY:O	1:B:88[A]:MET:HG2	2.12	0.50
1:A:329:ARG:HH21	1:A:401:LYS:NZ	2.10	0.49
1:A:84:GLY:O	1:A:88[A]:MET:HG2	2.14	0.48
1:A:115:SER:HB3	1:A:120:LEU:O	2.13	0.48
1:D:322:SER:OG	1:D:323:LYS:NZ	2.39	0.48
1:A:329:ARG:HE	1:A:401:LYS:CE	2.26	0.48
1:C:315:SER:HB3	1:C:318:GLU:HB3	1.96	0.48
1:B:40:MET:HG3	4:B:721:HOH:O	2.14	0.47
1:D:176:SER:HB2	1:D:206:CYS:SG	2.55	0.47
1:B:214:PHE:HZ	3:B:502:BEZ:H2	1.80	0.46
1:C:84:GLY:O	1:C:88[A]:MET:HG2	2.15	0.46
1:C:271:GLN:NE2	1:C:296:ARG:HE	2.13	0.46
1:A:307:LEU:O	1:A:311:GLU:HG3	2.16	0.46
1:C:88[B]:MET:HE1	1:C:237:TRP:CH2	2.51	0.46
1:A:106:ASN:HB2	4:A:789:HOH:O	2.15	0.46
1:C:106:ASN:HB2	4:C:760:HOH:O	2.16	0.45
1:C:209:LYS:HE3	4:C:809:HOH:O	2.16	0.45
1:A:275:VAL:HG13	1:A:311:GLU:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:THR:C	1:A:148:MET:HE2	2.42	0.45
1:A:347:LEU:O	1:A:385:MET:HE1	2.17	0.44
1:A:64:MET:HB2	1:A:106:ASN:ND2	2.32	0.44
1:D:377:LYS:NZ	4:D:604:HOH:O	2.43	0.44
1:D:214:PHE:HZ	3:D:502:BEZ:H6	1.82	0.44
1:D:40:MET:HG3	4:D:789:HOH:O	2.18	0.44
1:A:106:ASN:HB2	4:A:750:HOH:O	2.17	0.43
1:D:188:MET:SD	1:D:316:VAL:HG22	2.58	0.43
1:A:351:THR:HA	1:A:355:LEU:HB2	1.98	0.43
1:D:203:ILE:HD13	1:D:276:LEU:HD21	2.00	0.43
1:B:58:ARG:NE	4:B:606:HOH:O	2.37	0.43
1:B:115:SER:HB3	1:B:120:LEU:O	2.18	0.43
1:A:71:THR:OG1	1:D:57:GLU:OE1	2.24	0.42
1:C:115:SER:HB3	1:C:120:LEU:O	2.19	0.42
1:A:40:MET:HE2	1:A:40:MET:HB2	1.82	0.42
1:A:321:LEU:HD21	1:A:400:LEU:HD22	2.01	0.42
1:D:27:ASN:ND2	4:D:617:HOH:O	2.53	0.42
1:D:188:MET:HE3	1:D:319:PHE:HB2	2.02	0.42
1:D:209:LYS:O	1:D:209:LYS:HG3	2.20	0.42
1:A:355:LEU:O	4:A:602:HOH:O	2.22	0.41
1:A:188:MET:HE3	1:A:319:PHE:HB2	2.02	0.41
1:B:38:ASP:HA	4:B:601:HOH:O	2.21	0.41
1:C:209:LYS:HE2	1:C:209:LYS:CA	2.50	0.41
1:D:177:ALA:HB2	1:D:206:CYS:HB2	2.02	0.41
1:B:222:ASN:HD21	1:B:224:LYS:HB3	1.85	0.41
1:C:328:LEU:HD23	1:C:328:LEU:HA	1.94	0.41
1:D:88[A]:MET:HE1	1:D:237:TRP:HH2	1.86	0.40
1:B:177:ALA:HB2	1:B:206:CYS:HB2	2.02	0.40
1:D:321:LEU:HD23	1:D:321:LEU:HA	1.97	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:C:713:HOH:O	4:D:900:HOH:O[1_455]	2.15	0.05

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	370/405 (91%)	363 (98%)	7 (2%)	0	100	100
1	B	372/405 (92%)	365 (98%)	7 (2%)	0	100	100
1	C	374/405 (92%)	366 (98%)	8 (2%)	0	100	100
1	D	373/405 (92%)	368 (99%)	5 (1%)	0	100	100
All	All	1489/1620 (92%)	1462 (98%)	27 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	320/346 (92%)	320 (100%)	0	100	100
1	B	321/346 (93%)	319 (99%)	2 (1%)	78	72
1	C	322/346 (93%)	321 (100%)	1 (0%)	86	83
1	D	321/346 (93%)	320 (100%)	1 (0%)	86	83
All	All	1284/1384 (93%)	1280 (100%)	4 (0%)	87	83

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	272[A]	CYS
1	B	272[B]	CYS

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Mol	Chain	Res	Type
1	C	215	HIS
1	D	209	LYS

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

Mol	Chain	Res	Type
1	A	106	ASN
1	A	140	ASN
1	A	220	HIS
1	A	271	GLN
1	B	191	GLN
1	B	222	ASN
1	B	360	GLN
1	B	386	ASN
1	C	27	ASN
1	C	73	HIS
1	C	271	GLN
1	C	293	GLN
1	C	386	ASN
1	D	73	HIS
1	D	76	GLN
1	D	222	ASN
1	D	242	GLN
1	D	313	ASN
1	D	348	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

8 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
3	BEZ	B	502	-	9,9,9	0.77	0	11,11,11	1.06	2 (18%)
2	HQM	D	501	-	37,37,37	1.61	7 (18%)	50,52,52	2.04	11 (22%)
2	HQM	A	501	-	37,37,37	1.74	10 (27%)	50,52,52	1.86	9 (18%)
3	BEZ	A	502	-	9,9,9	0.67	0	11,11,11	1.13	1 (9%)
2	HQM	C	501	-	37,37,37	1.57	8 (21%)	50,52,52	1.95	11 (22%)
2	HQM	B	501	-	37,37,37	1.62	9 (24%)	50,52,52	1.75	9 (18%)
3	BEZ	C	502	-	9,9,9	0.79	0	11,11,11	1.06	0
3	BEZ	D	502	-	9,9,9	0.71	0	11,11,11	1.24	2 (18%)

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
3	BEZ	B	502	-	-	0/4/4/4	0/1/1/1
2	HQM	D	501	-	-	6/28/41/41	0/4/4/4
2	HQM	A	501	-	-	9/28/41/41	0/4/4/4
3	BEZ	A	502	-	-	0/4/4/4	0/1/1/1
2	HQM	C	501	-	-	9/28/41/41	0/4/4/4
2	HQM	B	501	-	-	7/28/41/41	0/4/4/4
3	BEZ	C	502	-	-	0/4/4/4	0/1/1/1
3	BEZ	D	502	-	-	0/4/4/4	0/1/1/1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HQM	C19-N18	4.69	1.47	1.37
2	D	501	HQM	C19-N18	3.90	1.46	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HQM	C19-N20	3.88	1.46	1.37
2	B	501	HQM	C19-N18	3.85	1.45	1.37
2	C	501	HQM	C19-N20	3.75	1.45	1.37
2	C	501	HQM	C19-N18	3.55	1.45	1.37
2	B	501	HQM	C19-N20	3.52	1.45	1.37
2	D	501	HQM	C19-N20	3.45	1.45	1.37
2	B	501	HQM	C04-N07	3.41	1.51	1.41
2	D	501	HQM	C04-N07	3.27	1.51	1.41
2	A	501	HQM	C04-N07	3.20	1.51	1.41
2	C	501	HQM	C04-N07	2.99	1.50	1.41
2	A	501	HQM	C21-N20	2.98	1.47	1.41
2	D	501	HQM	C01-C29	2.98	1.56	1.51
2	A	501	HQM	C03-N18	2.77	1.47	1.41
2	C	501	HQM	C12-N07	2.72	1.52	1.47
2	A	501	HQM	C12-N07	2.70	1.52	1.47
2	A	501	HQM	C01-C29	2.68	1.55	1.51
2	B	501	HQM	C08-N07	2.67	1.51	1.47
2	C	501	HQM	C01-C29	2.66	1.55	1.51
2	D	501	HQM	C12-N07	2.64	1.52	1.47
2	A	501	HQM	C31-C32	2.60	1.55	1.51
2	C	501	HQM	C08-N07	2.50	1.51	1.47
2	D	501	HQM	C08-N07	2.43	1.51	1.47
2	B	501	HQM	C12-N07	2.39	1.51	1.47
2	B	501	HQM	C01-C29	2.38	1.55	1.51
2	B	501	HQM	C21-N20	2.36	1.46	1.41
2	B	501	HQM	C31-C32	2.28	1.55	1.51
2	A	501	HQM	C08-N07	2.27	1.51	1.47
2	C	501	HQM	C21-N20	2.23	1.46	1.41
2	D	501	HQM	C03-C04	-2.23	1.38	1.40
2	B	501	HQM	C03-C04	-2.19	1.38	1.40
2	A	501	HQM	O28-C19	-2.11	1.18	1.23
2	C	501	HQM	C03-N18	2.09	1.45	1.41

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	HQM	C13-C12-N07	-8.22	101.77	112.80
2	C	501	HQM	C17-C12-N07	-7.74	102.42	112.80
2	B	501	HQM	C13-C12-N07	-7.06	103.32	112.80
2	A	501	HQM	C13-C12-N07	-6.12	104.58	112.80
2	D	501	HQM	C30-C29-C01	6.01	133.53	122.27
2	A	501	HQM	C30-C29-C01	5.95	133.41	122.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	HQM	C30-C29-C01	5.28	132.16	122.27
2	B	501	HQM	C30-C29-C01	4.98	131.59	122.27
2	D	501	HQM	C21-N20-C19	-3.99	118.47	126.61
2	C	501	HQM	C21-N20-C19	-3.91	118.64	126.61
2	A	501	HQM	C21-N20-C19	-3.53	119.41	126.61
2	B	501	HQM	C21-N20-C19	-3.39	119.68	126.61
2	D	501	HQM	C26-C21-C22	3.36	123.50	119.04
2	C	501	HQM	C01-C29-C31	3.32	127.16	120.67
2	A	501	HQM	C01-C29-C31	3.31	127.14	120.67
2	A	501	HQM	O33-C32-C31	3.22	122.28	113.93
2	D	501	HQM	C26-C25-C24	-3.13	117.77	121.37
2	D	501	HQM	C23-C22-C21	-3.11	116.72	120.30
2	C	501	HQM	C22-C23-C24	-3.05	117.85	121.37
2	B	501	HQM	C26-C21-C22	3.00	123.03	119.04
2	C	501	HQM	O33-C32-C31	2.98	121.66	113.93
2	D	501	HQM	O33-C32-C31	2.87	121.39	113.93
2	C	501	HQM	C26-C21-C22	2.71	122.64	119.04
2	B	501	HQM	O33-C32-O34	-2.66	118.04	124.08
2	A	501	HQM	C08-N07-C04	2.59	125.03	118.12
2	A	501	HQM	O33-C32-O34	-2.57	118.24	124.08
2	D	501	HQM	C25-C24-C23	2.35	122.67	118.11
2	D	501	HQM	O33-C32-O34	-2.27	118.94	124.08
2	C	501	HQM	C25-C26-C21	-2.26	117.70	120.30
3	A	502	BEZ	O2-C-C1	2.24	120.60	114.84
2	A	501	HQM	C22-C23-C24	-2.22	118.81	121.37
2	B	501	HQM	O33-C32-C31	2.20	119.63	113.93
3	B	502	BEZ	O2-C-C1	2.16	120.39	114.84
2	D	501	HQM	C05-C04-N07	-2.16	117.86	122.53
3	B	502	BEZ	O2-C-O1	-2.14	118.75	123.35
2	B	501	HQM	C23-C22-C21	-2.13	117.85	120.30
2	B	501	HQM	C26-C25-C24	-2.12	118.93	121.37
2	C	501	HQM	C05-C06-C01	-2.11	119.08	121.18
2	A	501	HQM	C26-C21-C22	2.10	121.83	119.04
3	D	502	BEZ	O2-C-C1	2.05	120.11	114.84
3	D	502	BEZ	O2-C-O1	-2.04	118.96	123.35
2	B	501	HQM	C01-C29-C31	2.03	124.64	120.67
2	C	501	HQM	O33-C32-O34	-2.03	119.48	124.08
2	D	501	HQM	C30-C31-C32	-2.02	115.81	118.94
2	C	501	HQM	C25-C24-C23	2.01	122.01	118.11

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	HQM	C09-C08-N07-C04
2	B	501	HQM	C09-C08-N07-C04
2	C	501	HQM	C09-C08-N07-C04
2	D	501	HQM	C09-C08-N07-C04
2	A	501	HQM	N07-C08-C09-C11
2	A	501	HQM	N07-C08-C09-C10
2	A	501	HQM	C03-C04-N07-C08
2	A	501	HQM	C29-C31-C32-O33
2	A	501	HQM	C29-C31-C32-O34
2	C	501	HQM	C29-C31-C32-O33
2	C	501	HQM	C29-C31-C32-O34
2	D	501	HQM	C29-C31-C32-O33
2	D	501	HQM	C29-C31-C32-O34
2	B	501	HQM	C03-C04-N07-C08
2	C	501	HQM	C03-C04-N07-C08
2	D	501	HQM	C03-C04-N07-C08
2	C	501	HQM	N07-C08-C09-C11
2	B	501	HQM	C29-C31-C32-O33
2	B	501	HQM	C29-C31-C32-O34
2	A	501	HQM	C17-C12-N07-C08
2	A	501	HQM	C05-C04-N07-C08
2	B	501	HQM	C17-C12-N07-C08
2	C	501	HQM	C13-C12-N07-C08
2	C	501	HQM	N07-C08-C09-C10
2	B	501	HQM	C05-C04-N07-C08
2	D	501	HQM	C26-C21-N20-C19
2	B	501	HQM	C26-C21-N20-C19
2	C	501	HQM	C05-C04-N07-C08
2	D	501	HQM	C05-C04-N07-C08
2	A	501	HQM	C30-C31-C32-O34
2	C	501	HQM	C30-C31-C32-O34

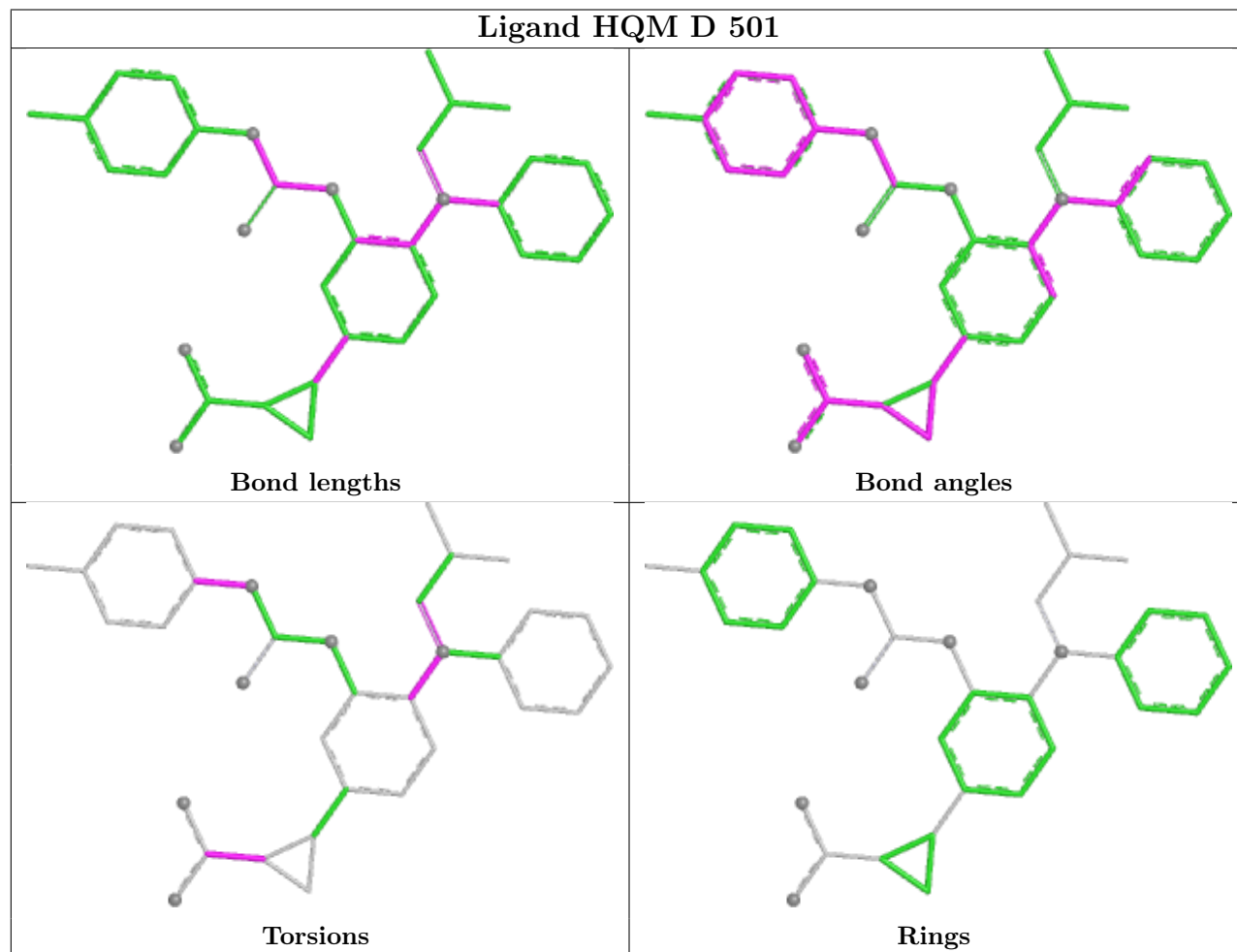
There are no ring outliers.

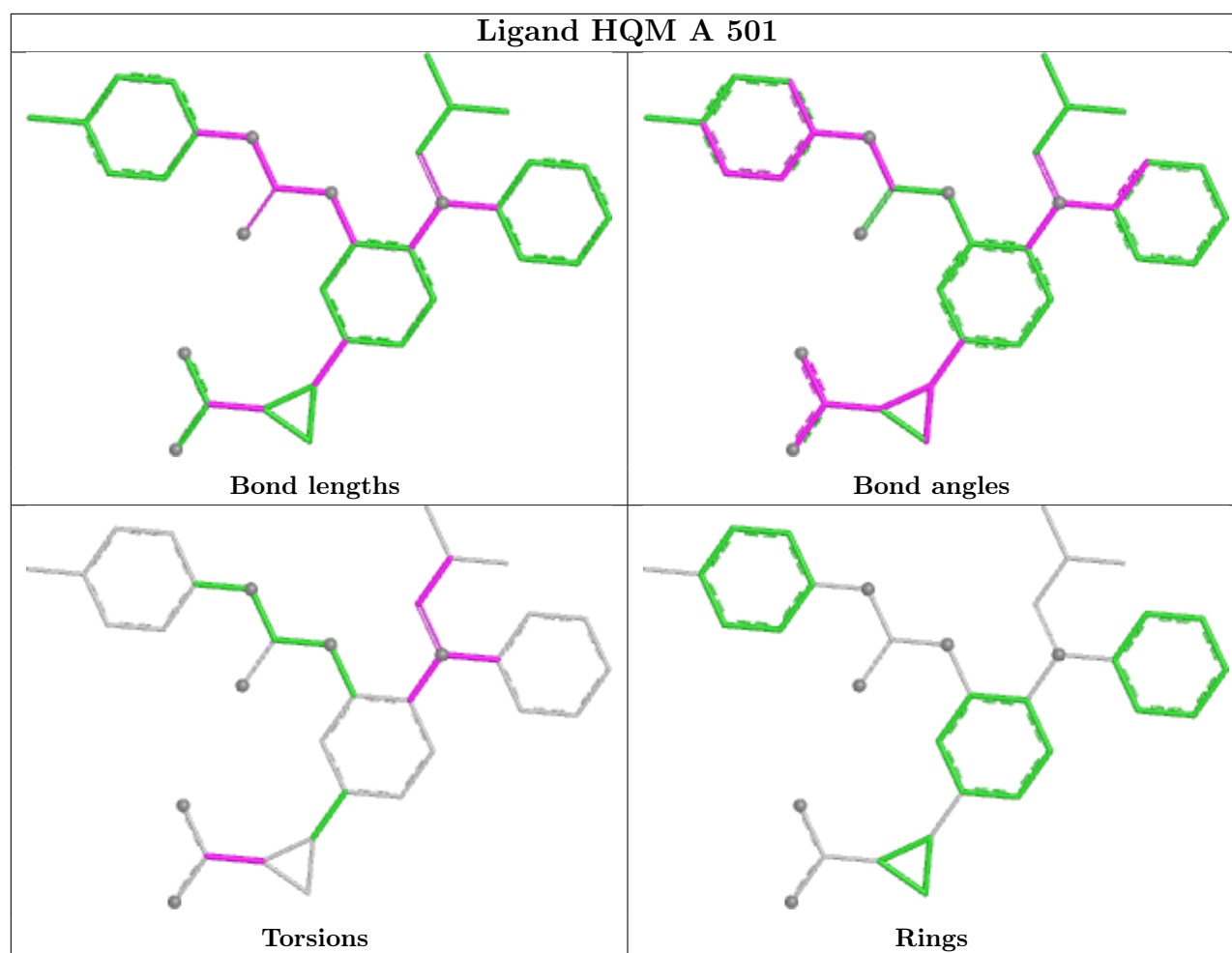
2 monomers are involved in 2 short contacts:

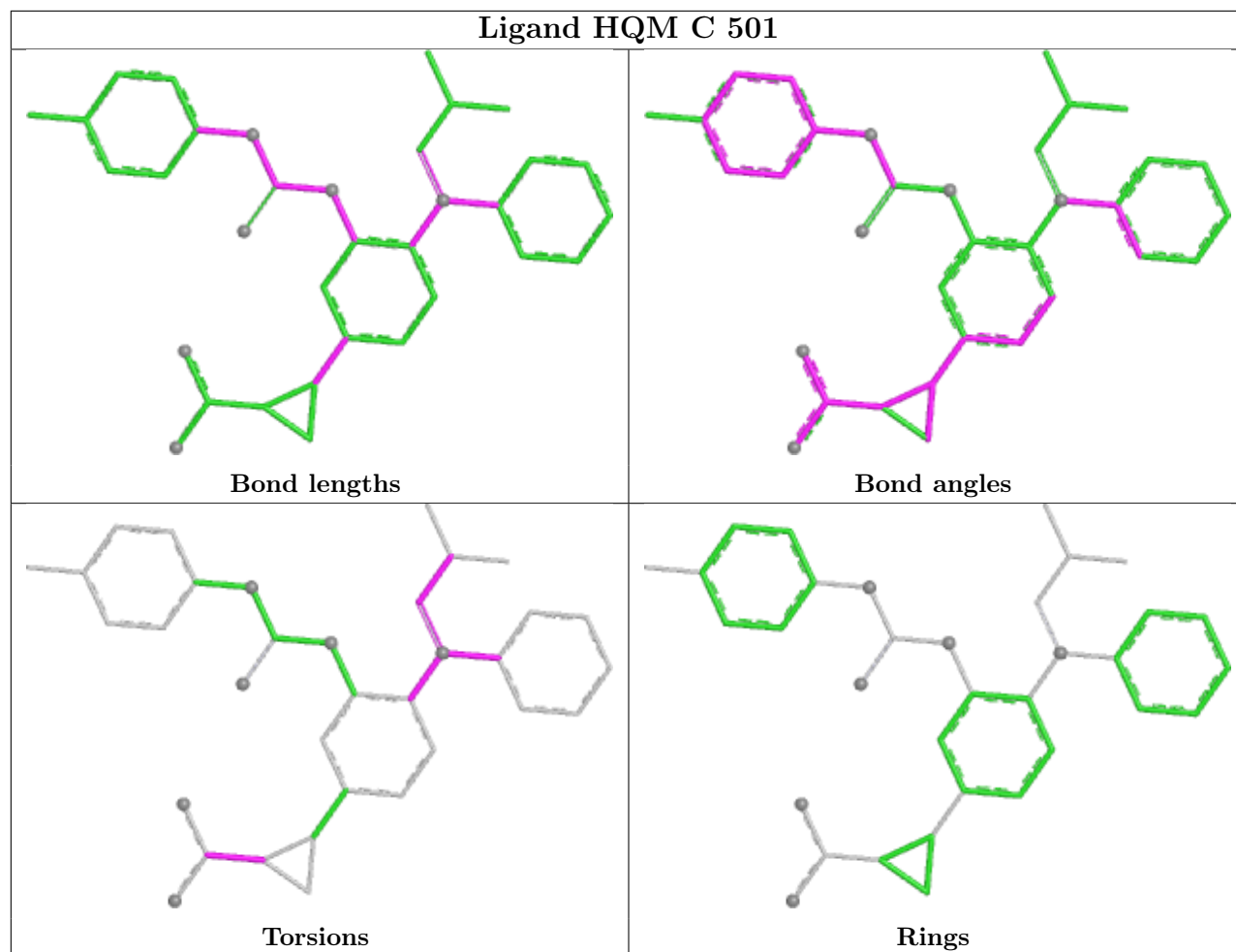
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	BEZ	1	0
3	D	502	BEZ	1	0

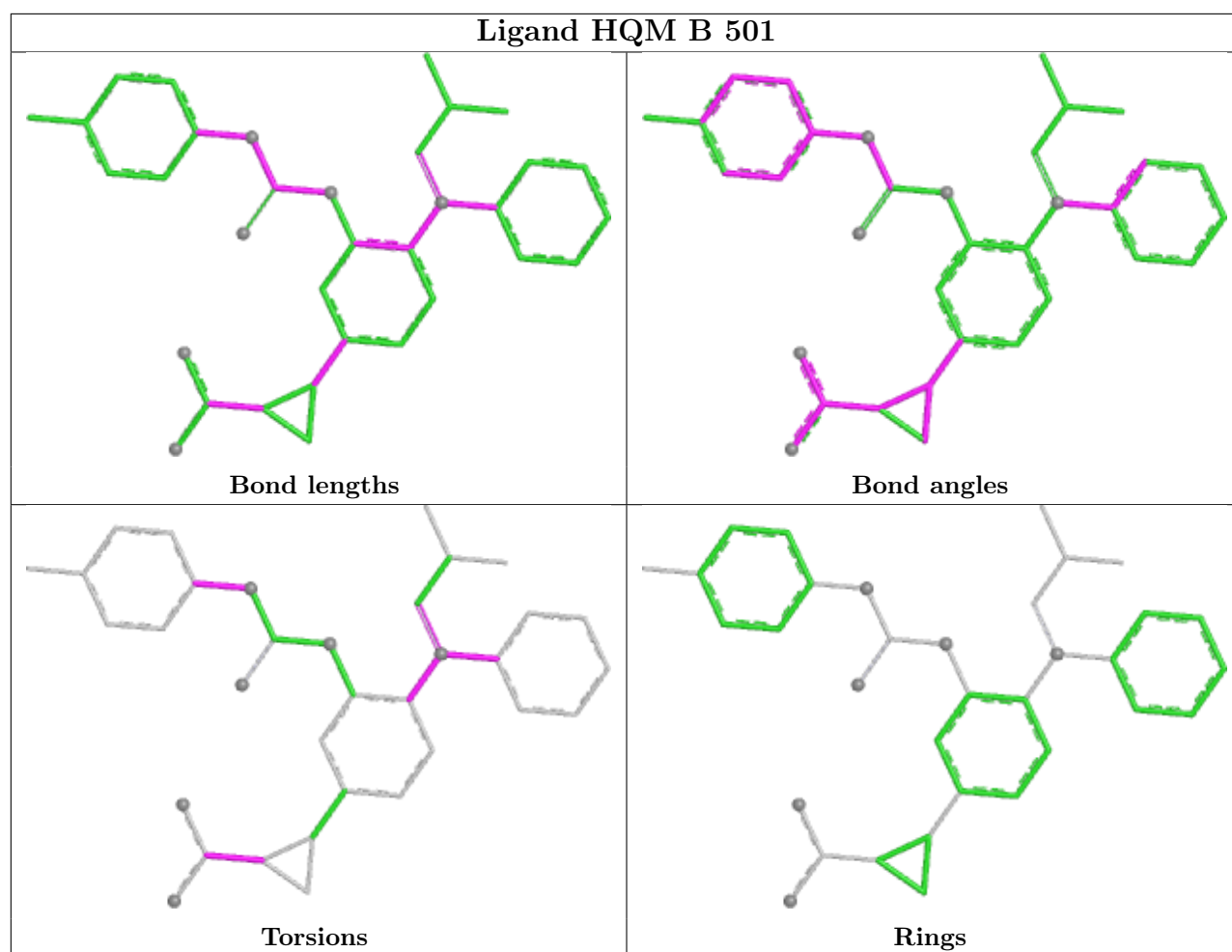
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	372/405 (91%)	0.47	11 (2%) 52 56	14, 34, 50, 60	2 (0%)
1	B	373/405 (92%)	0.50	20 (5%) 31 34	14, 33, 51, 62	3 (0%)
1	C	375/405 (92%)	0.28	12 (3%) 50 54	12, 30, 48, 65	3 (0%)
1	D	375/405 (92%)	0.20	14 (3%) 45 49	11, 28, 45, 57	2 (0%)
All	All	1495/1620 (92%)	0.36	57 (3%) 44 48	11, 32, 49, 65	10 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	14	ALA	6.5
1	C	379	THR	5.8
1	A	379	THR	5.2
1	B	14	ALA	4.6
1	B	379	THR	4.5
1	D	378	GLY	4.3
1	C	378	GLY	4.2
1	D	379	THR	4.1
1	A	378	GLY	3.8
1	C	403	GLY	3.8
1	C	380	GLY	3.7
1	B	378	GLY	3.7
1	C	14	ALA	3.4
1	C	377	LYS	3.4
1	B	376	ALA	3.3
1	A	376	ALA	3.3
1	C	376	ALA	3.3
1	D	190	MET	3.2
1	A	377	LYS	3.2
1	B	283	ALA	3.1
1	D	185	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	401	LYS	3.1
1	B	380	GLY	3.0
1	B	401	LYS	3.0
1	A	400	LEU	2.9
1	D	284	GLY	2.9
1	B	189	GLN	2.8
1	B	400	LEU	2.8
1	A	208	GLU	2.7
1	B	284	GLY	2.6
1	B	287	HIS	2.6
1	A	285	GLY	2.5
1	D	400	LEU	2.5
1	D	191	GLN	2.5
1	B	319	PHE	2.5
1	C	287	HIS	2.5
1	B	318	GLU	2.5
1	A	287	HIS	2.4
1	D	27	ASN	2.4
1	C	68	ASP	2.4
1	D	287	HIS	2.4
1	B	290	GLN	2.4
1	B	58	ARG	2.4
1	C	284	GLY	2.3
1	A	399	LEU	2.3
1	B	57	GLU	2.2
1	D	377	LYS	2.2
1	C	318	GLU	2.2
1	B	321	LEU	2.2
1	D	380	GLY	2.2
1	B	314	PRO	2.2
1	D	184	VAL	2.1
1	A	190	MET	2.1
1	C	190	MET	2.1
1	D	360	GLN	2.1
1	B	320	VAL	2.0
1	B	399	LEU	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

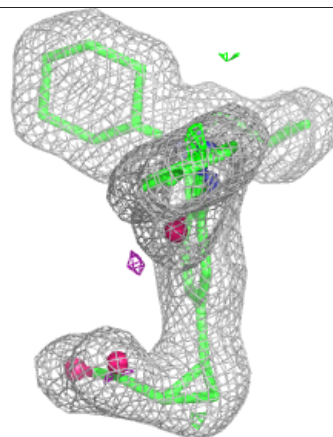
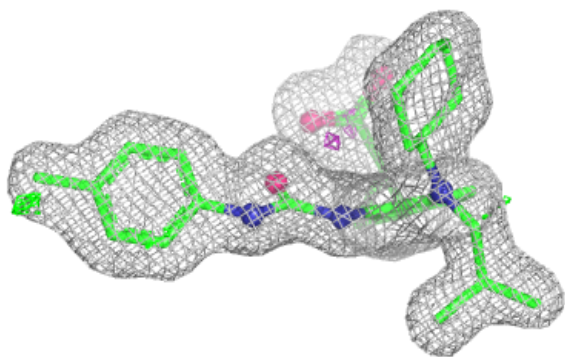
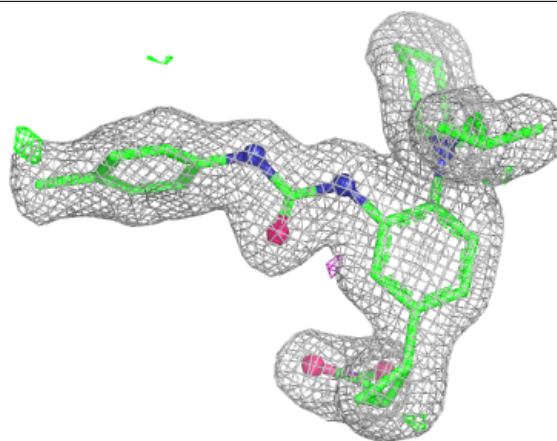
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	BEZ	B	502	9/9	0.91	0.10	31,34,35,36	0
3	BEZ	C	502	9/9	0.91	0.09	28,30,31,32	0
3	BEZ	A	502	9/9	0.92	0.09	33,34,37,38	0
3	BEZ	D	502	9/9	0.94	0.07	27,28,29,29	0
2	HQM	A	501	34/34	0.95	0.08	25,29,40,42	0
2	HQM	C	501	34/34	0.96	0.07	20,24,36,37	0
2	HQM	D	501	34/34	0.96	0.07	19,23,33,35	0
2	HQM	B	501	34/34	0.96	0.07	22,26,35,39	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.

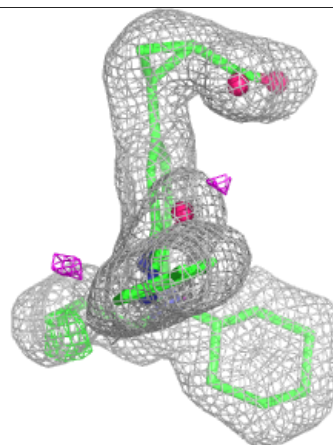
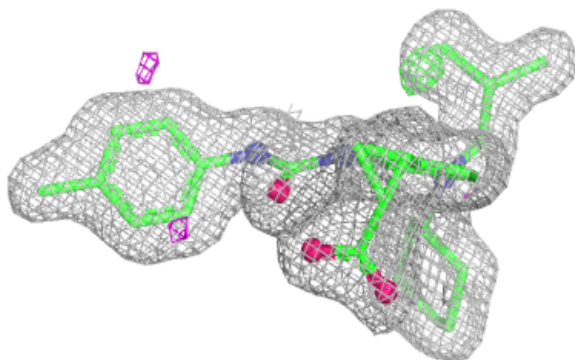
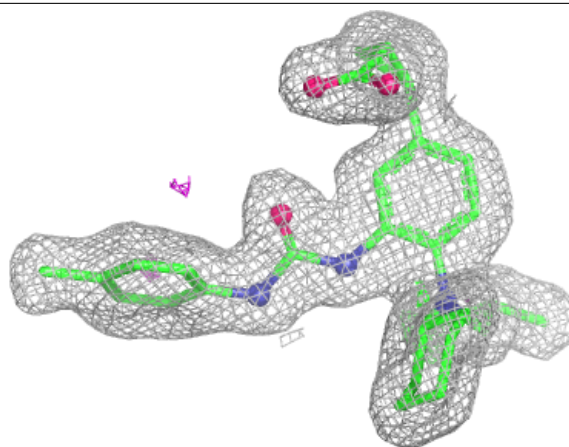
Electron density around HQM A 501:

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



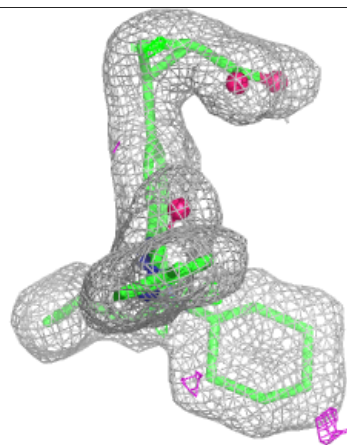
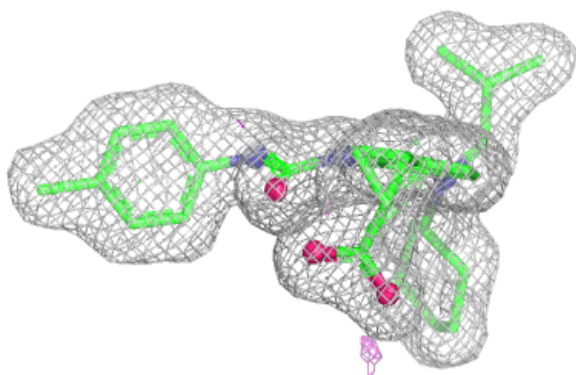
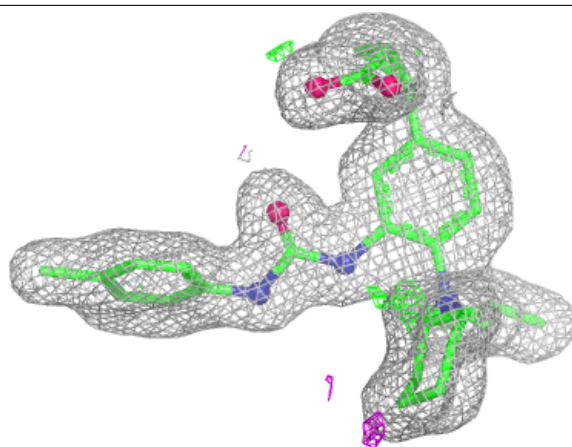
Electron density around HQM C 501:

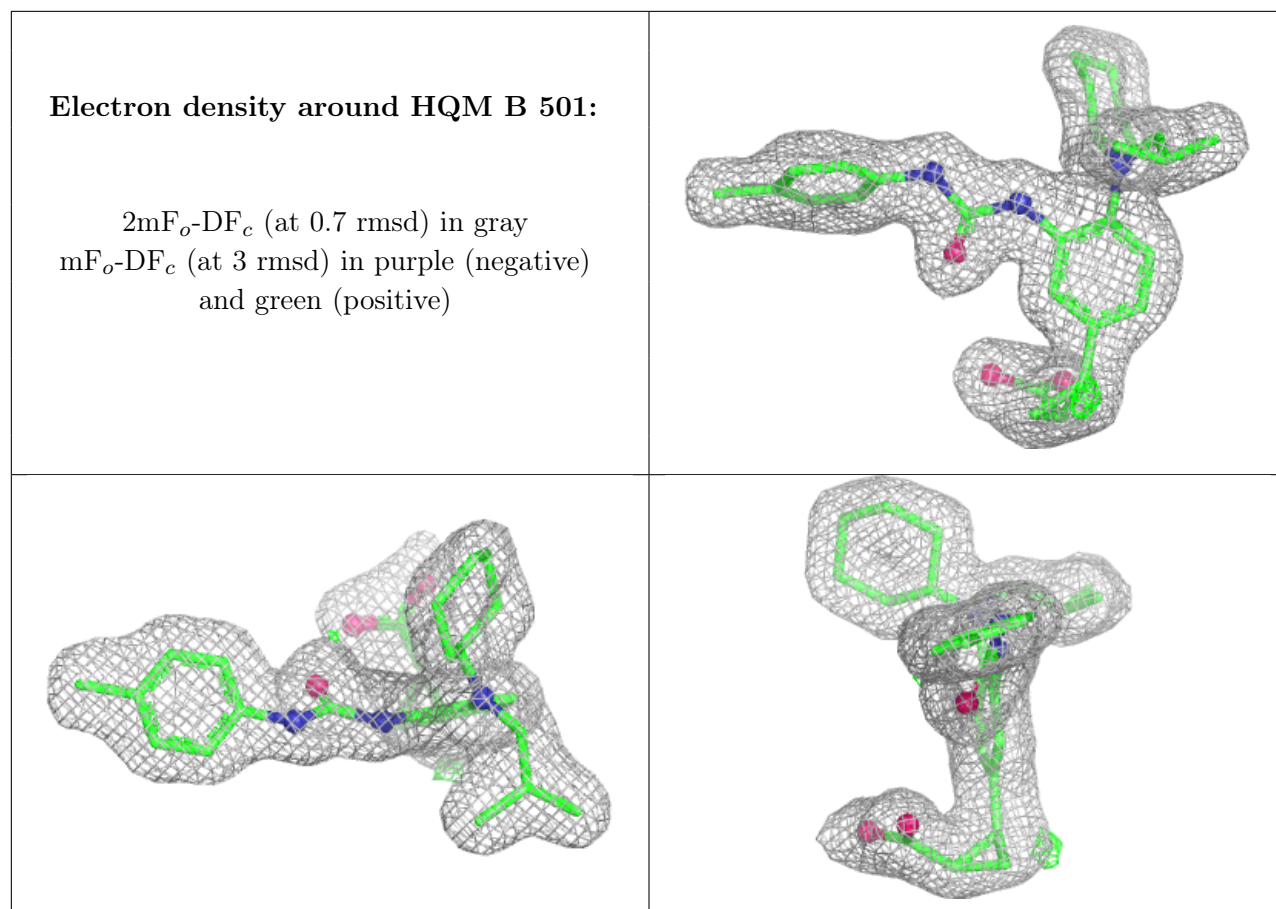
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HQM D 501:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.