



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 12:00 PM UTC

PDB ID : 7MAS / pdb_00007mas
Title : Drug Resistant HIV-1 Protease (L10F, M46I, I50V, F53L, L63P, G73S) in Complex with DRV
Authors : Lockbaum, G.J.; Schiffer, C.A.
Deposited on : 2021-03-31
Resolution : 1.50 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

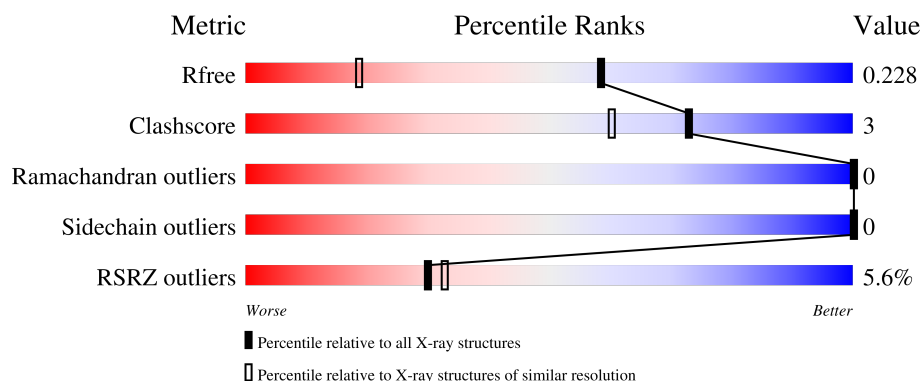
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	99	6% 98% .
1	B	99	5% 94% 6%
1	C	99	5% 92% 8%
1	D	99	6% 90% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6662 atoms, of which 3187 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 Protease.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	99	Total	C	H	N	O	S	0	0	0
			1521	482	775	126	135	3			
1	D	99	Total	C	H	N	O	S	0	0	0
			1538	485	788	127	135	3			
1	C	99	Total	C	H	N	O	S	0	0	0
			1523	482	777	126	135	3			
1	A	99	Total	C	H	N	O	S	0	0	0
			1517	480	773	128	133	3			

There are 32 discrepancies between the modelled and reference sequences:

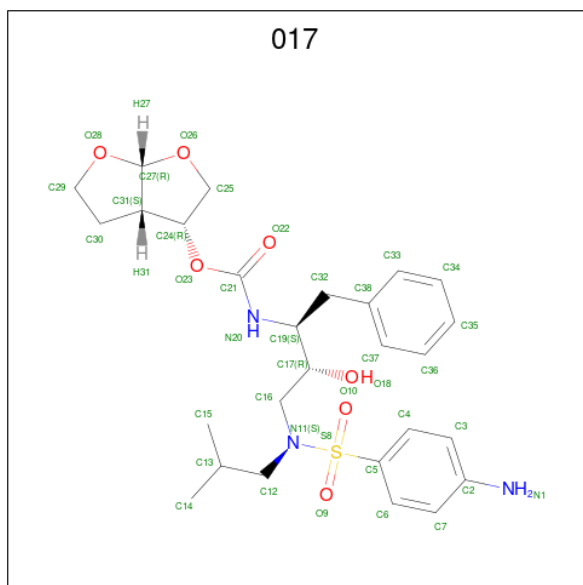
Chain	Residue	Modelled	Actual	Comment	Reference
B	7	LYS	GLN	conflict	UNP P03367
B	10	PHE	LEU	engineered mutation	UNP P03367
B	37	ASN	SER	conflict	UNP P03367
B	46	ILE	MET	engineered mutation	UNP P03367
B	50	VAL	ILE	engineered mutation	UNP P03367
B	53	LEU	PHE	engineered mutation	UNP P03367
B	63	PRO	LEU	engineered mutation	UNP P03367
B	73	SER	GLY	engineered mutation	UNP P03367
D	7	LYS	GLN	conflict	UNP P03367
D	10	PHE	LEU	engineered mutation	UNP P03367
D	37	ASN	SER	conflict	UNP P03367
D	46	ILE	MET	engineered mutation	UNP P03367
D	50	VAL	ILE	engineered mutation	UNP P03367
D	53	LEU	PHE	engineered mutation	UNP P03367
D	63	PRO	LEU	engineered mutation	UNP P03367
D	73	SER	GLY	engineered mutation	UNP P03367
C	7	LYS	GLN	conflict	UNP P03367
C	10	PHE	LEU	engineered mutation	UNP P03367
C	37	ASN	SER	conflict	UNP P03367
C	46	ILE	MET	engineered mutation	UNP P03367
C	50	VAL	ILE	engineered mutation	UNP P03367

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Chain	Residue	Modelled	Actual	Comment	Reference
C	53	LEU	PHE	engineered mutation	UNP P03367
C	63	PRO	LEU	engineered mutation	UNP P03367
C	73	SER	GLY	engineered mutation	UNP P03367
A	7	LYS	GLN	conflict	UNP P03367
A	10	PHE	LEU	engineered mutation	UNP P03367
A	37	ASN	SER	conflict	UNP P03367
A	46	ILE	MET	engineered mutation	UNP P03367
A	50	VAL	ILE	engineered mutation	UNP P03367
A	53	LEU	PHE	engineered mutation	UNP P03367
A	63	PRO	LEU	engineered mutation	UNP P03367
A	73	SER	GLY	engineered mutation	UNP P03367

- Molecule 2 is (3R,3AS,6AR)-HEXAHYDROFURO[2,3-B]FURAN-3-YL(1S,2R)-3-[(4-AMINOPHENYL)SULFONYL](ISOBUTYL)AMINO]-1-BENZYL-2-HYDROXYPROPYLCARBAMATE (CCD ID: 017) (formula: C₂₇H₃₇N₃O₇S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	C	1	Total	C	H	N	O	S	0	0
			75	27	37	3	7	1		
2	A	1	Total	C	H	N	O	S	0	0
			75	27	37	3	7	1		

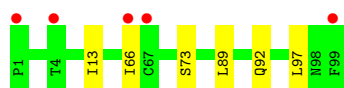
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	104	Total 104	O 104	0	0
3	D	94	Total 94	O 94	0	0
3	C	108	Total 108	O 108	0	0
3	A	107	Total 107	O 107	0	0

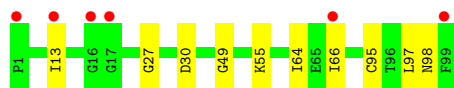
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

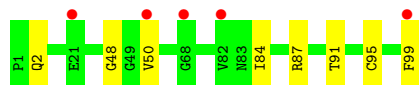
- Molecule 1: Protease



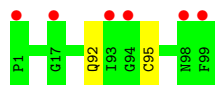
- Molecule 1: Protease



- Molecule 1: Protease



- Molecule 1: Protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.74Å 36.28Å 103.51Å 90.00° 98.66° 90.00°	Depositor
Resolution (Å)	46.34 – 1.50 46.34 – 1.50	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.34-1.50) 93.5 (46.34-1.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.83 (at 1.50Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.202 , 0.225 0.205 , 0.228	Depositor DCC
R_{free} test set	2000 reflections (3.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.304	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.6	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	0.96	EDS
Total number of atoms	6662	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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 41.64 % 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 2.2931e-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:
017

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	0.53	0/758	0.69	0/1032
1	B	0.49	0/760	0.62	0/1034
1	C	0.54	0/760	0.65	0/1034
1	D	0.44	0/764	0.59	0/1038
All	All	0.50	0/3042	0.64	0/4138

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 [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	744	773	773	2	0
1	B	746	775	777	5	0
1	C	746	777	777	8	0
1	D	750	788	788	9	0
2	A	38	37	37	0	0
2	C	38	37	37	6	0
3	A	107	0	0	1	0
3	B	104	0	0	0	0
3	C	108	0	0	0	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	94	0	0	2	2
All	All	3475	3187	3189	20	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 3.

All (20) 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:D:13:ILE:HG22	1:D:66:ILE:HG12	1.84	0.57
1:C:50:VAL:HG23	2:C:101:017:O9	2.05	0.57
1:C:84:ILE:HD11	2:C:101:017:H141	1.87	0.57
1:B:13:ILE:HG22	1:B:66:ILE:HG12	1.87	0.55
1:D:98:ASN:ND2	1:C:2:GLN:HG2	2.25	0.51
1:D:13:ILE:HD12	1:D:64:ILE:HD11	1.92	0.51
1:D:55:LYS:NZ	3:D:105:HOH:O	2.45	0.50
1:D:49:GLY:HA3	2:C:101:017:O10	2.14	0.48
1:C:87:ARG:O	1:C:91:THR:HG23	2.15	0.46
1:B:73:SER:HB2	1:B:92:GLN:HE22	1.81	0.46
1:D:95:CYS:HA	1:C:99:PHE:HB3	1.98	0.46
1:C:48:GLY:O	2:C:101:017:H302	2.16	0.45
1:D:97:LEU:HD22	1:C:95:CYS:SG	2.57	0.45
1:B:66:ILE:HD11	1:B:89:LEU:CD2	2.48	0.44
1:B:66:ILE:HD11	1:B:89:LEU:HD21	2.00	0.43
1:A:92:GLN:HG3	3:A:254:HOH:O	2.17	0.43
1:B:97:LEU:HD22	1:A:95:CYS:SG	2.59	0.42
1:D:30:ASP:HB3	3:D:133:HOH:O	2.20	0.41
1:C:50:VAL:CG2	2:C:101:017:O9	2.67	0.41
1:D:27:GLY:O	2:C:101:017:H121	2.22	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:D:119:HOH:O	3:C:271:HOH:O[2_444]	2.00	0.20
3:D:119:HOH:O	3:C:222:HOH:O[2_444]	2.08	0.12

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	97/99 (98%)	97 (100%)	0	0	100	100
1	B	97/99 (98%)	97 (100%)	0	0	100	100
1	C	97/99 (98%)	96 (99%)	1 (1%)	0	100	100
1	D	97/99 (98%)	97 (100%)	0	0	100	100
All	All	388/396 (98%)	387 (100%)	1 (0%)	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	80/84 (95%)	80 (100%)	0	100	100
1	B	81/84 (96%)	81 (100%)	0	100	100
1	C	81/84 (96%)	81 (100%)	0	100	100
1	D	82/84 (98%)	82 (100%)	0	100	100
All	All	324/336 (96%)	324 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	92	GLN
1	C	2	GLN
1	A	18	GLN
1	A	61	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	017	A	101	-	41,41,41	3.74	17 (41%)	52,58,58	1.49	11 (21%)
2	017	C	101	-	41,41,41	3.90	20 (48%)	52,58,58	2.11	11 (21%)

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	017	A	101	-	-	0/36/55/55	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	017	C	101	-	-	1/36/55/55	0/4/4/4

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	101	017	C31-C24	-10.23	1.33	1.53
2	C	101	017	C31-C24	-9.43	1.35	1.53
2	C	101	017	S8-N11	9.17	1.76	1.63
2	A	101	017	S8-N11	8.22	1.75	1.63
2	A	101	017	O26-C27	8.12	1.58	1.41
2	C	101	017	O26-C27	7.86	1.57	1.41
2	C	101	017	C25-C24	7.49	1.66	1.52
2	A	101	017	C25-C24	7.44	1.66	1.52
2	C	101	017	O28-C27	-7.25	1.27	1.41
2	C	101	017	C30-C29	-7.10	1.32	1.51
2	C	101	017	O28-C29	6.95	1.63	1.43
2	A	101	017	C30-C29	-6.76	1.33	1.51
2	A	101	017	O28-C29	6.52	1.61	1.43
2	C	101	017	C5-S8	6.02	1.84	1.76
2	A	101	017	O28-C27	-5.64	1.30	1.41
2	A	101	017	O10-S8	5.25	1.49	1.43
2	A	101	017	C5-S8	5.19	1.83	1.76
2	C	101	017	C27-C31	4.88	1.60	1.52
2	C	101	017	C21-N20	4.85	1.46	1.34
2	A	101	017	C27-C31	4.30	1.59	1.52
2	A	101	017	C21-N20	4.20	1.45	1.34
2	A	101	017	C16-N11	4.20	1.55	1.47
2	C	101	017	O26-C25	-3.56	1.36	1.43
2	C	101	017	C16-N11	3.45	1.53	1.47
2	C	101	017	C2-N1	3.16	1.49	1.38
2	A	101	017	O26-C25	-2.87	1.37	1.43
2	C	101	017	O23-C21	2.84	1.40	1.35
2	A	101	017	C16-C17	2.67	1.56	1.52
2	C	101	017	O10-S8	2.63	1.46	1.43
2	C	101	017	C16-C17	2.40	1.56	1.52
2	A	101	017	O23-C21	2.34	1.39	1.35
2	C	101	017	C6-C5	-2.21	1.35	1.38
2	C	101	017	C3-C2	-2.19	1.35	1.40
2	C	101	017	C19-N20	2.15	1.49	1.46
2	A	101	017	C2-N1	2.13	1.45	1.38
2	C	101	017	O18-C17	-2.03	1.39	1.43
2	A	101	017	C6-C5	-2.01	1.35	1.38

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	101	017	O10-S8-O9	-7.04	108.61	119.59
2	C	101	017	O10-S8-N11	7.01	113.28	106.69
2	A	101	017	O26-C27-C31	-4.18	100.97	107.02
2	C	101	017	O26-C27-C31	-4.09	101.11	107.02
2	C	101	017	C32-C19-N20	-4.03	104.36	110.08
2	A	101	017	C32-C19-N20	-4.00	104.40	110.08
2	C	101	017	O26-C25-C24	-3.78	97.87	104.42
2	A	101	017	O10-S8-O9	-3.69	113.83	119.59
2	C	101	017	C29-C30-C31	3.45	108.09	103.38
2	C	101	017	C32-C19-C17	3.31	117.04	111.66
2	A	101	017	O26-C25-C24	-3.26	98.77	104.42
2	A	101	017	C17-C19-N20	-2.77	104.77	109.97
2	C	101	017	C38-C32-C19	-2.48	109.19	113.40
2	A	101	017	O9-S8-C5	2.30	110.97	108.10
2	C	101	017	O23-C21-N20	2.24	113.12	110.26
2	A	101	017	C29-C30-C31	2.18	106.37	103.38
2	C	101	017	O9-S8-N11	2.17	108.72	106.69
2	A	101	017	O23-C24-C31	2.16	116.05	108.88
2	C	101	017	C7-C2-C3	2.12	121.34	118.16
2	A	101	017	O28-C27-C31	2.10	110.05	107.02
2	A	101	017	O28-C29-C30	2.08	110.44	106.34
2	A	101	017	C32-C19-C17	2.03	114.95	111.66

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	101	017	C25-C24-O23-C21

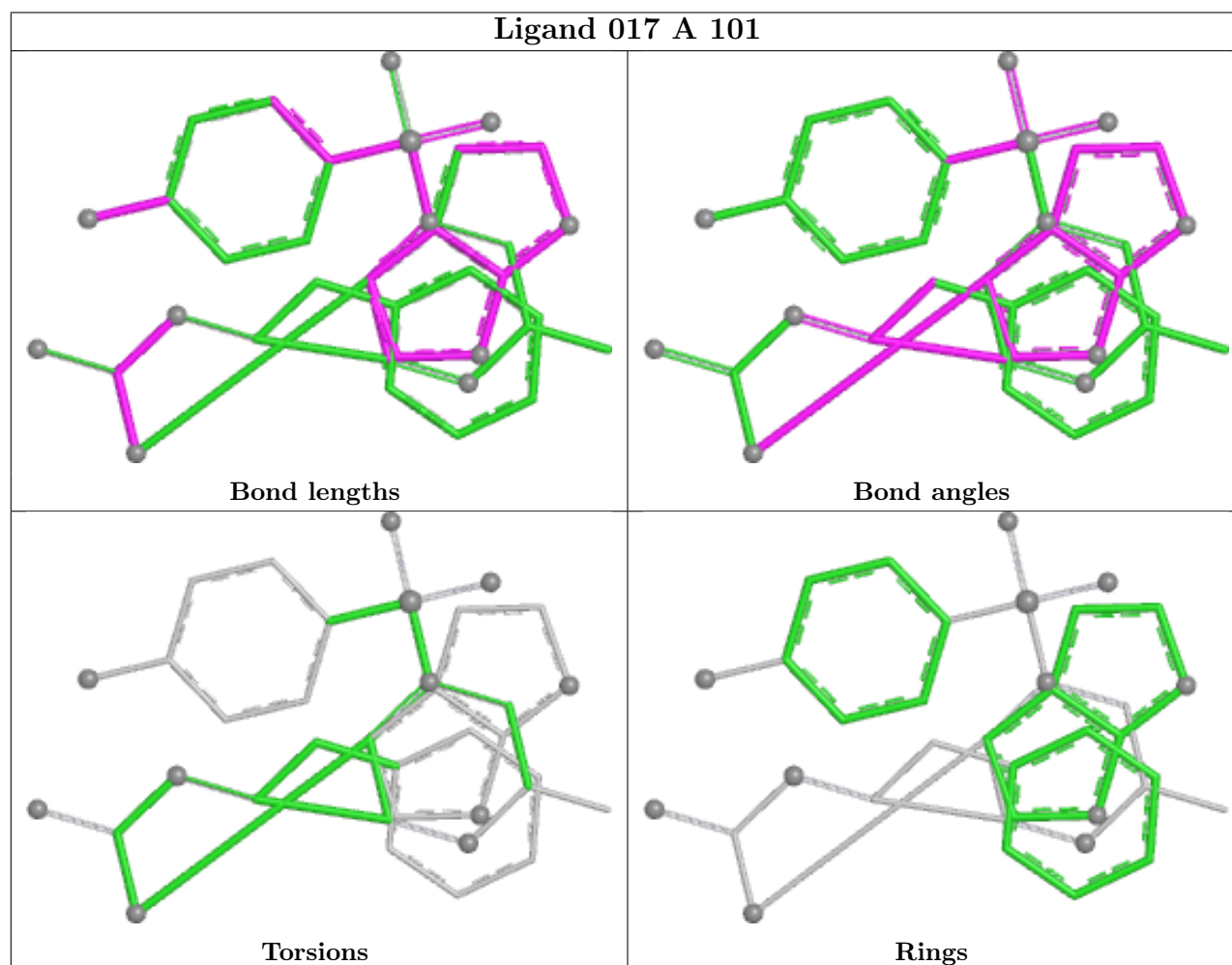
There are no ring outliers.

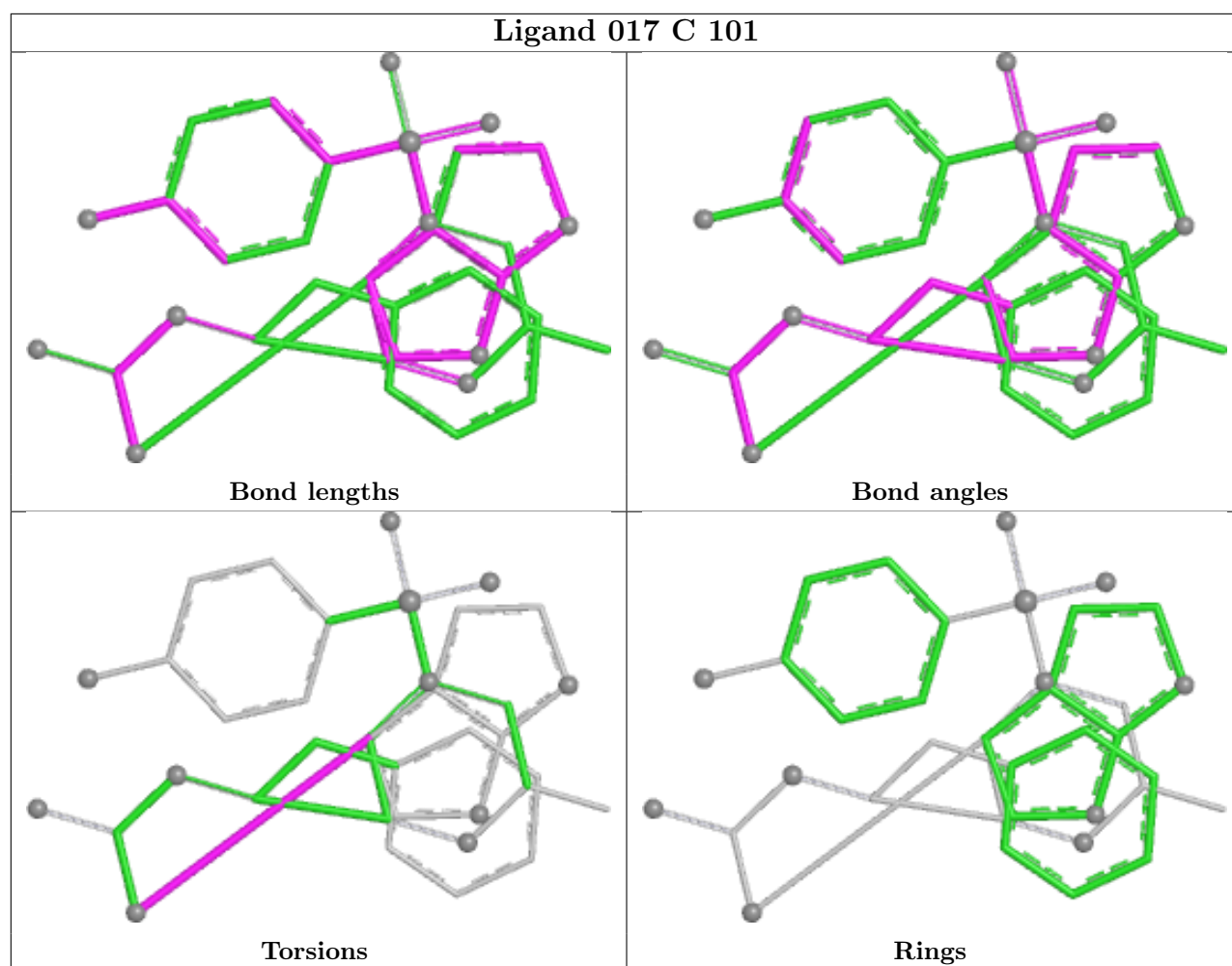
1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	101	017	6	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	99/99 (100%)	0.56	6 (6%)	27	29	17, 24, 38, 52	1 (1%)
1	B	99/99 (100%)	0.55	5 (5%)	33	36	17, 26, 42, 56	1 (1%)
1	C	99/99 (100%)	0.57	5 (5%)	33	36	18, 24, 35, 40	1 (1%)
1	D	99/99 (100%)	0.82	6 (6%)	27	29	19, 29, 45, 51	1 (1%)
All	All	396/396 (100%)	0.62	22 (5%)	30	33	17, 26, 43, 56	4 (1%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	99	PHE	5.4
1	D	1	PRO	4.3
1	A	1	PRO	4.3
1	B	1	PRO	4.1
1	D	17	GLY	3.7
1	A	99	PHE	3.7
1	C	50	VAL	3.2
1	B	4	THR	2.8
1	D	99	PHE	2.8
1	A	17	GLY	2.7
1	C	68	GLY	2.7
1	D	13	ILE	2.7
1	B	67	CYS	2.7
1	D	16	GLY	2.6
1	C	99	PHE	2.6
1	C	82	VAL	2.6
1	B	66	ILE	2.4
1	A	93	ILE	2.3
1	A	98	ASN	2.2
1	A	94	GLY	2.2
1	D	66	ILE	2.1
1	C	21	GLU	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](#)

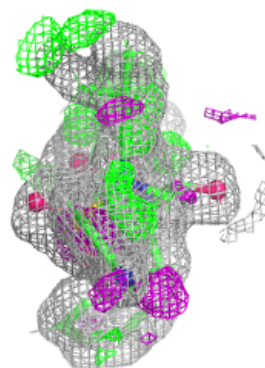
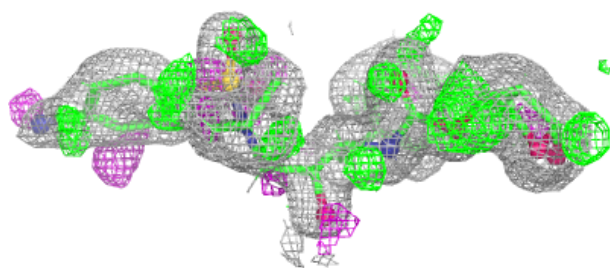
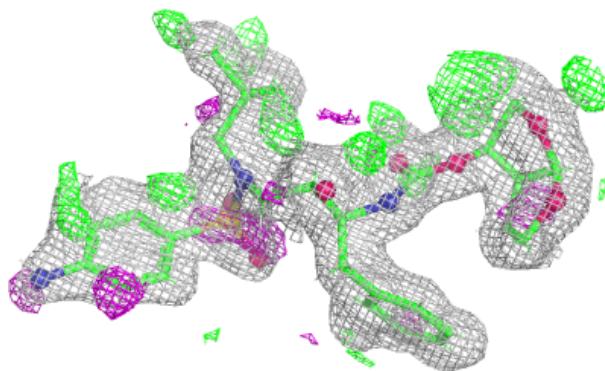
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	017	C	101	38/38	0.85	0.13	21,30,40,40	0
2	017	A	101	38/38	0.95	0.08	16,23,30,33	0

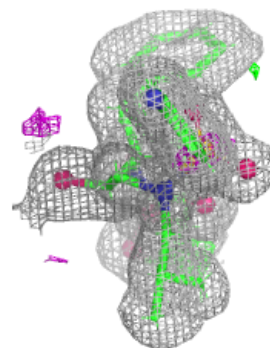
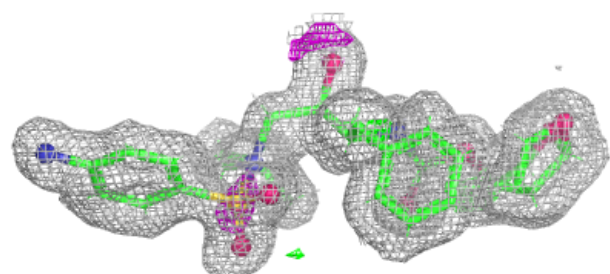
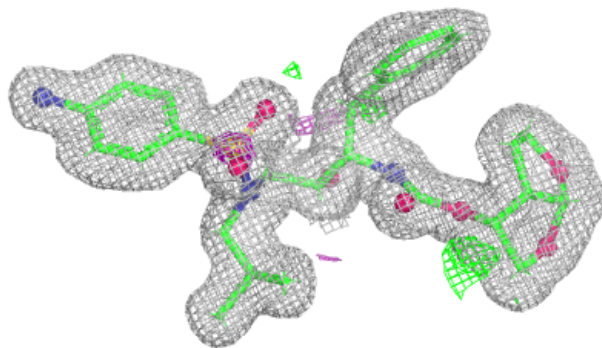
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 017 C 101:

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

**Electron density around 017 A 101:**

$2mF_o-DF_c$ (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.