



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:00 AM UTC

PDB ID : 8GKF / pdb_00008gkf
Title : Phosphopantetheinyl transferase PptT from Mycobacterium tuberculosis in complex with Ralnitrexed.
Authors : Krieger, I.V.; Singh, A.; Sacchettini, J.C.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2023-03-18
Resolution : 2.45 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

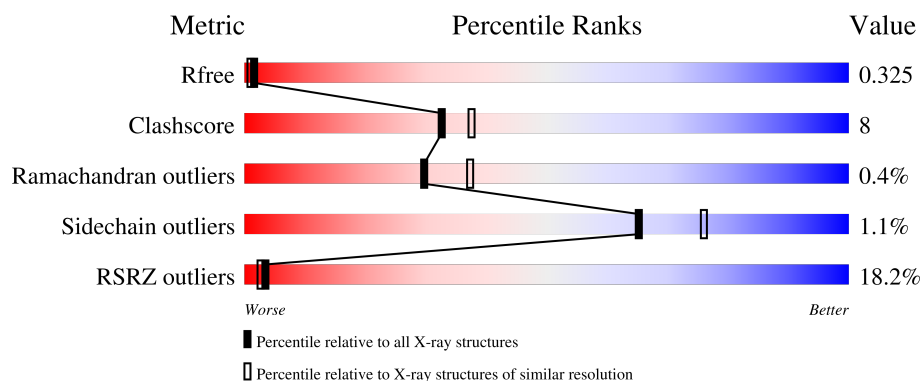
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	235	7% 83%10%6%
1	B	235	5% 82%13%5%
1	C	235	14% 77%18%. .
1	D	235	17% 83%12%. .

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Mol	Chain	Length	Quality of chain
1	E	235	
1	F	235	
1	G	235	
1	H	235	
1	I	235	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 14268 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 4'-phosphopantetheinyl transferase PptT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1686	1077	299	303	7			
1	B	223	Total	C	N	O	S	0	0	0
			1708	1093	302	306	7			
1	C	225	Total	C	N	O	S	0	0	0
			1712	1093	301	311	7			
1	D	225	Total	C	N	O	S	0	0	0
			1659	1061	296	297	5			
1	E	222	Total	C	N	O	S	0	0	0
			1672	1072	295	298	7			
1	F	97	Total	C	N	O	S	0	0	0
			701	455	121	123	2			
1	G	197	Total	C	N	O	S	0	0	0
			1508	968	267	268	5			
1	H	216	Total	C	N	O	S	0	0	0
			1617	1039	279	293	6			
1	I	217	Total	C	N	O	S	0	0	0
			1642	1049	284	302	7			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	ALA	-	expression tag	UNP O33336
A	229	GLY	-	expression tag	UNP O33336
A	230	GLU	-	expression tag	UNP O33336
A	231	ASN	-	expression tag	UNP O33336
A	232	LEU	-	expression tag	UNP O33336
A	233	TYR	-	expression tag	UNP O33336
A	234	PHE	-	expression tag	UNP O33336
A	235	GLN	-	expression tag	UNP O33336
B	228	ALA	-	expression tag	UNP O33336
B	229	GLY	-	expression tag	UNP O33336
B	230	GLU	-	expression tag	UNP O33336

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Chain	Residue	Modelled	Actual	Comment	Reference
B	231	ASN	-	expression tag	UNP O33336
B	232	LEU	-	expression tag	UNP O33336
B	233	TYR	-	expression tag	UNP O33336
B	234	PHE	-	expression tag	UNP O33336
B	235	GLN	-	expression tag	UNP O33336
C	228	ALA	-	expression tag	UNP O33336
C	229	GLY	-	expression tag	UNP O33336
C	230	GLU	-	expression tag	UNP O33336
C	231	ASN	-	expression tag	UNP O33336
C	232	LEU	-	expression tag	UNP O33336
C	233	TYR	-	expression tag	UNP O33336
C	234	PHE	-	expression tag	UNP O33336
C	235	GLN	-	expression tag	UNP O33336
D	228	ALA	-	expression tag	UNP O33336
D	229	GLY	-	expression tag	UNP O33336
D	230	GLU	-	expression tag	UNP O33336
D	231	ASN	-	expression tag	UNP O33336
D	232	LEU	-	expression tag	UNP O33336
D	233	TYR	-	expression tag	UNP O33336
D	234	PHE	-	expression tag	UNP O33336
D	235	GLN	-	expression tag	UNP O33336
E	228	ALA	-	expression tag	UNP O33336
E	229	GLY	-	expression tag	UNP O33336
E	230	GLU	-	expression tag	UNP O33336
E	231	ASN	-	expression tag	UNP O33336
E	232	LEU	-	expression tag	UNP O33336
E	233	TYR	-	expression tag	UNP O33336
E	234	PHE	-	expression tag	UNP O33336
E	235	GLN	-	expression tag	UNP O33336
F	228	ALA	-	expression tag	UNP O33336
F	229	GLY	-	expression tag	UNP O33336
F	230	GLU	-	expression tag	UNP O33336
F	231	ASN	-	expression tag	UNP O33336
F	232	LEU	-	expression tag	UNP O33336
F	233	TYR	-	expression tag	UNP O33336
F	234	PHE	-	expression tag	UNP O33336
F	235	GLN	-	expression tag	UNP O33336
G	228	ALA	-	expression tag	UNP O33336
G	229	GLY	-	expression tag	UNP O33336
G	230	GLU	-	expression tag	UNP O33336
G	231	ASN	-	expression tag	UNP O33336
G	232	LEU	-	expression tag	UNP O33336

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Chain	Residue	Modelled	Actual	Comment	Reference
G	233	TYR	-	expression tag	UNP O33336
G	234	PHE	-	expression tag	UNP O33336
G	235	GLN	-	expression tag	UNP O33336
H	228	ALA	-	expression tag	UNP O33336
H	229	GLY	-	expression tag	UNP O33336
H	230	GLU	-	expression tag	UNP O33336
H	231	ASN	-	expression tag	UNP O33336
H	232	LEU	-	expression tag	UNP O33336
H	233	TYR	-	expression tag	UNP O33336
H	234	PHE	-	expression tag	UNP O33336
H	235	GLN	-	expression tag	UNP O33336
I	228	ALA	-	expression tag	UNP O33336
I	229	GLY	-	expression tag	UNP O33336
I	230	GLU	-	expression tag	UNP O33336
I	231	ASN	-	expression tag	UNP O33336
I	232	LEU	-	expression tag	UNP O33336
I	233	TYR	-	expression tag	UNP O33336
I	234	PHE	-	expression tag	UNP O33336
I	235	GLN	-	expression tag	UNP O33336

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- Chemical structure of D16, a complex molecule featuring a thiazole ring, a pyrimidine ring, and a carboxylic acid group. The structure is labeled with various atoms and bonds, including a chiral center marked with a wedge bond.

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	C	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	D	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	E	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	F	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	G	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	H	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
2	I	1	Total	C	N	O	S	0	0
			32	21	4	6	1		

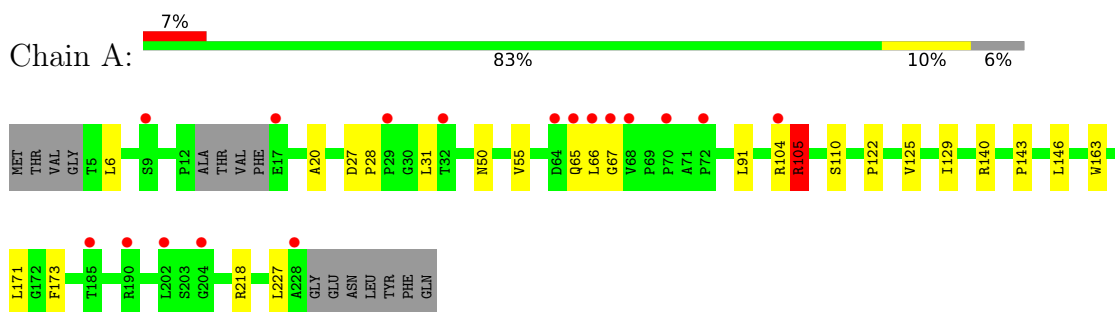
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	20	Total	O	0	0
			20	20		
3	C	6	Total	O	0	0
			6	6		
3	D	7	Total	O	0	0
			7	7		
3	E	8	Total	O	0	0
			8	8		
3	F	3	Total	O	0	0
			3	3		
3	G	9	Total	O	0	0
			9	9		
3	H	10	Total	O	0	0
			10	10		
3	I	1	Total	O	0	0
			1	1		

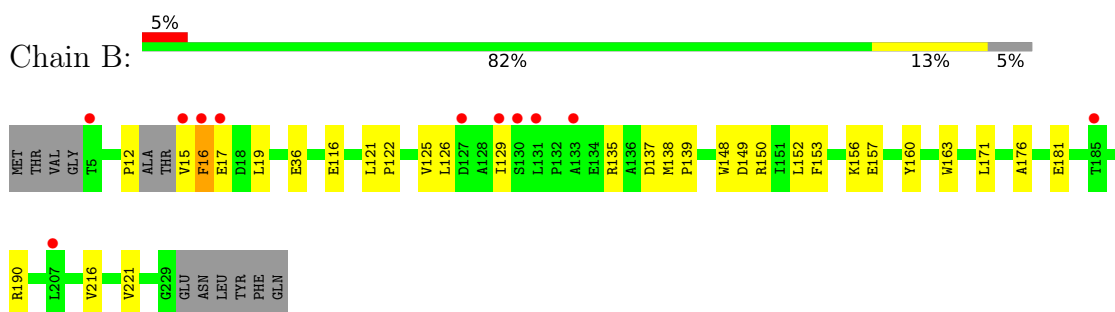
3 Residue-property plots [i](#)

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.

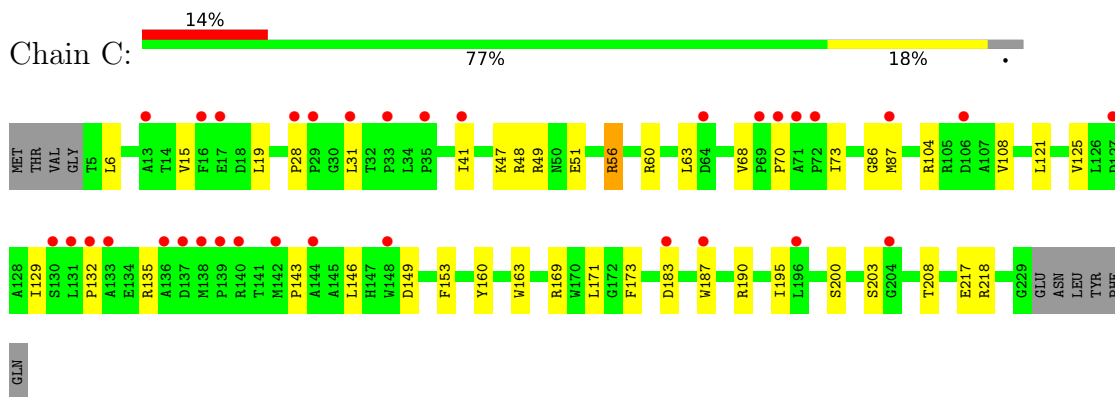
- Molecule 1: 4'-phosphopantetheinyl transferase PptT



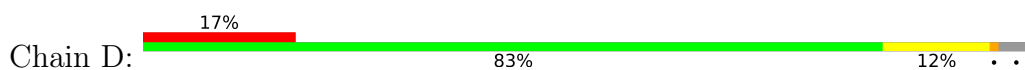
- Molecule 1: 4'-phosphopantetheinyl transferase PptT

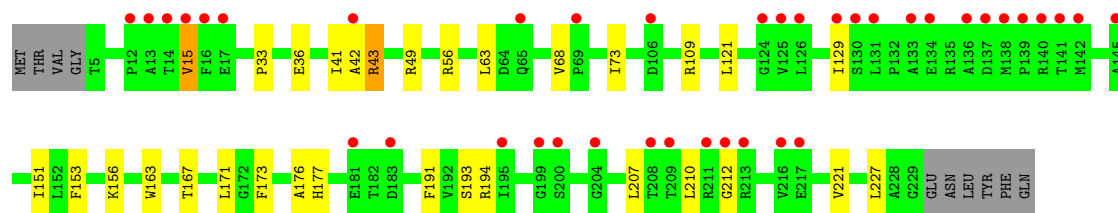


- Molecule 1: 4'-phosphopantetheinyl transferase PptT

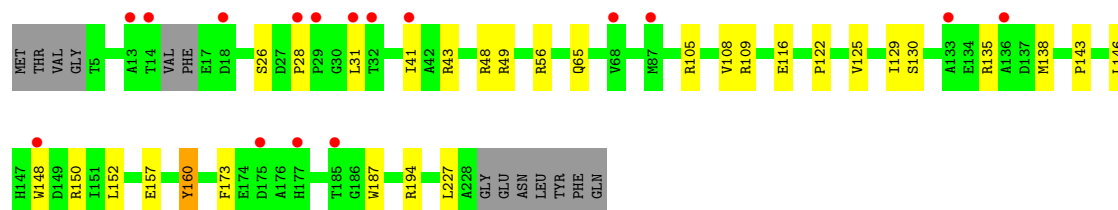
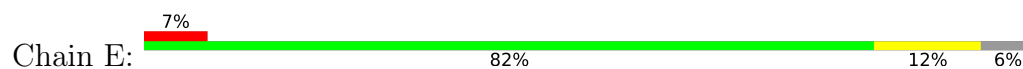


- Molecule 1: 4'-phosphopantetheinyl transferase PptT

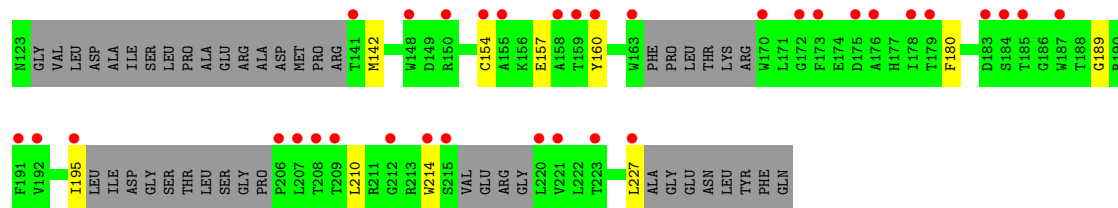
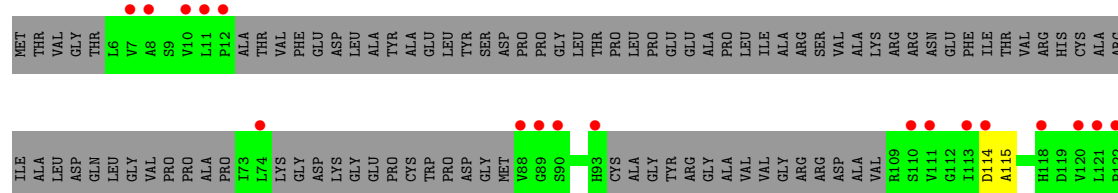
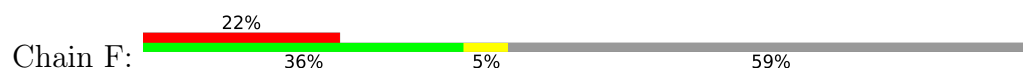




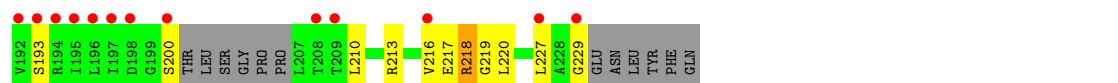
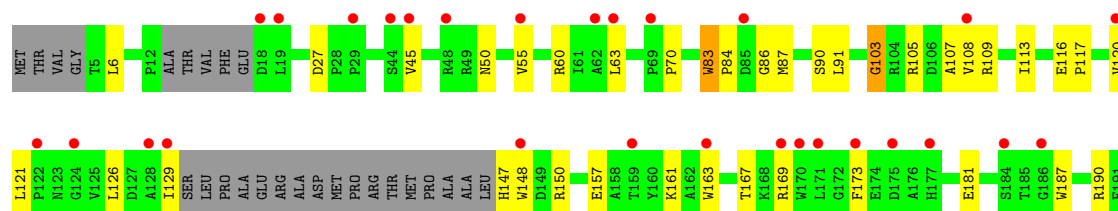
• Molecule 1: 4'-phosphopantetheinyl transferase PptT



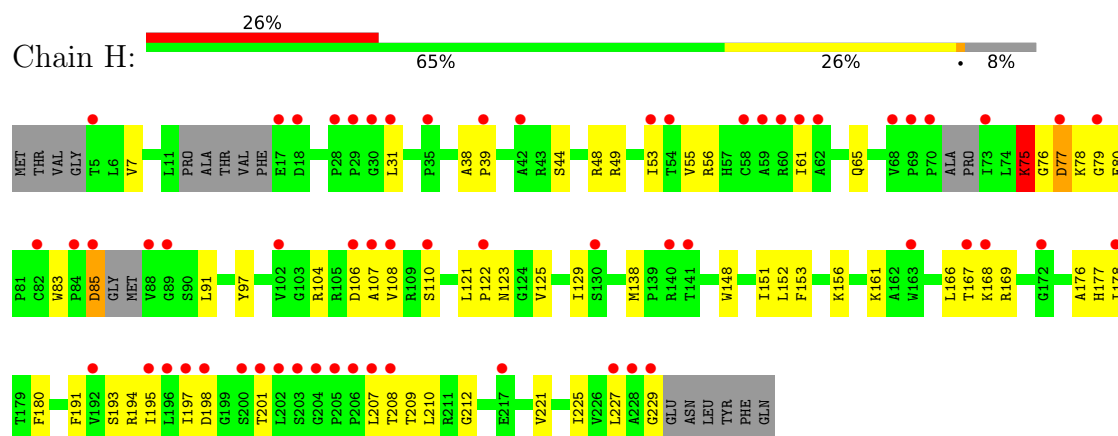
• Molecule 1: 4'-phosphopantetheinyl transferase PptT



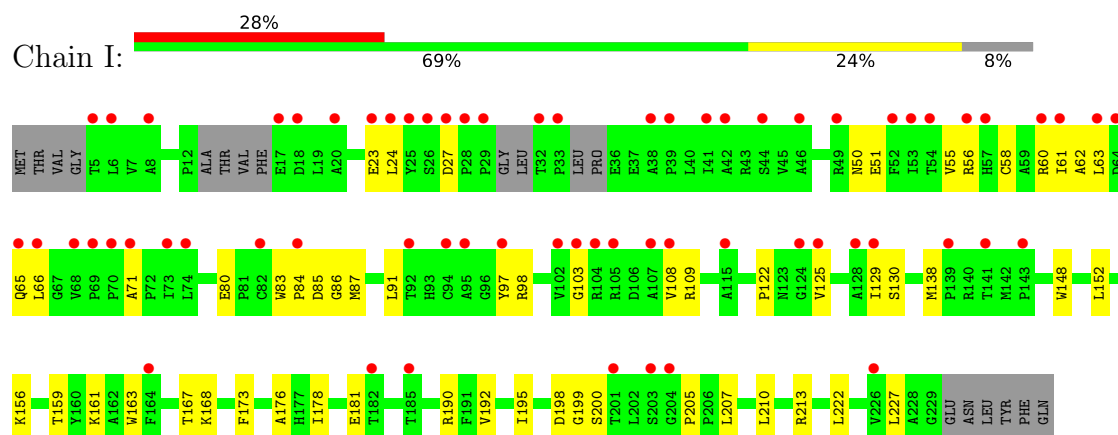
• Molecule 1: 4'-phosphopantetheinyl transferase PptT



• Molecule 1: 4'-phosphopantetheinyl transferase PptT



• Molecule 1: 4'-phosphopantetheinyl transferase PptT



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	141.83Å 141.83Å 209.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.08 – 2.45 21.08 – 2.45	Depositor EDS
% Data completeness (in resolution range)	91.0 (21.08-2.45) 90.9 (21.08-2.45)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.44Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.281 , 0.326 0.282 , 0.325	Depositor DCC
R_{free} test set	3619 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	14268	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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 38.92 % 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 3.4281e-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: D16

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	0.25	0/1730	0.50	1/2363 (0.0%)
1	B	0.34	1/1753 (0.1%)	0.57	4/2394 (0.2%)
1	C	0.28	0/1757	0.53	0/2403
1	D	0.39	0/1702	0.60	0/2330
1	E	0.43	0/1716	0.57	1/2348 (0.0%)
1	F	0.22	0/712	0.57	0/965
1	G	0.41	0/1545	0.70	4/2106 (0.2%)
1	H	0.43	1/1657 (0.1%)	0.65	3/2264 (0.1%)
1	I	0.39	0/1683	0.64	2/2299 (0.1%)
All	All	0.36	2/14255 (0.0%)	0.60	15/19472 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	126	LEU	CA-C	-5.56	1.45	1.52
1	H	123	ASN	CA-C	-5.47	1.45	1.52

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	105	ARG	N-CA-C	7.20	119.13	111.28
1	G	103	GLY	O-C-N	-6.85	117.68	123.23
1	E	130	SER	N-CA-C	6.56	120.21	109.72
1	H	77	ASP	N-CA-C	6.00	120.36	109.32
1	G	83	TRP	O-C-N	5.99	126.44	121.35

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	1686	0	1695	17	0
1	B	1708	0	1716	21	0
1	C	1712	0	1715	29	0
1	D	1659	0	1628	22	0
1	E	1672	0	1673	18	0
1	F	701	0	653	7	0
1	G	1508	0	1508	36	0
1	H	1617	0	1586	45	0
1	I	1642	0	1622	33	0
2	A	32	0	20	0	0
2	B	32	0	20	0	0
2	C	32	0	20	2	0
2	D	32	0	20	0	0
2	E	32	0	20	2	0
2	F	32	0	20	0	0
2	G	32	0	20	1	0
2	H	32	0	20	2	0
2	I	32	0	20	0	0
3	A	11	0	0	0	0
3	B	20	0	0	0	0
3	C	6	0	0	0	0
3	D	7	0	0	0	0
3	E	8	0	0	0	0
3	F	3	0	0	0	0
3	G	9	0	0	0	0
3	H	10	0	0	0	0
3	I	1	0	0	0	0
All	All	14268	0	13976	222	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

The worst 5 of 222 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:G:169:ARG:HH22	1:G:200:SER:HB2	1.27	0.98
1:G:169:ARG:NH2	1:G:200:SER:HB2	1.95	0.81
1:I:66:LEU:HD22	1:I:87:MET:HE1	1.64	0.76
1:A:122:PRO:HG2	1:A:125:VAL:HG21	1.66	0.76
1:G:181:GLU:HG3	1:G:190:ARG:HE	1.50	0.75

There are no symmetry-related clashes.

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	216/235 (92%)	211 (98%)	5 (2%)	0	100	100
1	B	219/235 (93%)	214 (98%)	5 (2%)	0	100	100
1	C	223/235 (95%)	215 (96%)	8 (4%)	0	100	100
1	D	223/235 (95%)	213 (96%)	8 (4%)	2 (1%)	14	17
1	E	218/235 (93%)	217 (100%)	1 (0%)	0	100	100
1	F	81/235 (34%)	81 (100%)	0	0	100	100
1	G	189/235 (80%)	172 (91%)	14 (7%)	3 (2%)	7	7
1	H	208/235 (88%)	195 (94%)	11 (5%)	2 (1%)	12	14
1	I	209/235 (89%)	194 (93%)	14 (7%)	1 (0%)	24	32
All	All	1786/2115 (84%)	1712 (96%)	66 (4%)	8 (0%)	30	37

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	42	ALA
1	G	218	ARG
1	G	219	GLY
1	G	45	VAL
1	H	83	TRP

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	177/190 (93%)	175 (99%)	2 (1%)	65	76
1	B	179/190 (94%)	178 (99%)	1 (1%)	78	86
1	C	179/190 (94%)	176 (98%)	3 (2%)	53	67
1	D	163/190 (86%)	161 (99%)	2 (1%)	63	74
1	E	172/190 (90%)	169 (98%)	3 (2%)	53	67
1	F	65/190 (34%)	65 (100%)	0	100	100
1	G	156/190 (82%)	156 (100%)	0	100	100
1	H	163/190 (86%)	160 (98%)	3 (2%)	51	65
1	I	171/190 (90%)	169 (99%)	2 (1%)	63	74
All	All	1425/1710 (83%)	1409 (99%)	16 (1%)	65	76

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	56	ARG
1	H	85	ASP
1	E	65	GLN
1	H	78	LYS
1	D	43	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	123	ASN
1	B	123	ASN
1	E	65	GLN
1	I	177	HIS

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 ⓘ

9 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	D16	I	301	-	33,34,34	0.73	0	46,48,48	0.89	2 (4%)
2	D16	F	301	-	33,34,34	0.73	0	46,48,48	0.95	2 (4%)
2	D16	A	301	-	33,34,34	0.72	0	46,48,48	0.88	1 (2%)
2	D16	B	301	-	33,34,34	0.74	0	46,48,48	0.91	1 (2%)
2	D16	D	301	-	33,34,34	0.74	0	46,48,48	0.89	1 (2%)
2	D16	C	301	-	33,34,34	0.75	0	46,48,48	0.88	2 (4%)
2	D16	H	301	-	33,34,34	0.75	0	46,48,48	0.90	1 (2%)
2	D16	G	301	-	33,34,34	0.75	0	46,48,48	0.95	1 (2%)
2	D16	E	301	-	33,34,34	0.74	0	46,48,48	0.90	1 (2%)

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	D16	I	301	-	-	4/25/25/25	0/3/3/3
2	D16	F	301	-	-	4/25/25/25	0/3/3/3
2	D16	A	301	-	-	4/25/25/25	0/3/3/3
2	D16	B	301	-	-	4/25/25/25	0/3/3/3
2	D16	D	301	-	-	3/25/25/25	0/3/3/3
2	D16	C	301	-	-	4/25/25/25	0/3/3/3
2	D16	H	301	-	-	3/25/25/25	0/3/3/3
2	D16	G	301	-	-	7/25/25/25	0/3/3/3
2	D16	E	301	-	-	3/25/25/25	0/3/3/3

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	D16	C8A-C4A-C4	-3.06	116.81	120.11
2	I	301	D16	C8A-C4A-C4	-3.05	116.81	120.11
2	F	301	D16	C8A-C4A-C4	-3.05	116.82	120.11
2	H	301	D16	C8A-C4A-C4	-3.02	116.85	120.11
2	A	301	D16	C8A-C4A-C4	-3.02	116.85	120.11

There are no chirality outliers.

5 of 36 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	301	D16	CB-CA-CT-O2
2	G	301	D16	CB-CA-CT-O1
2	H	301	D16	CT-CA-CB-CG
2	C	301	D16	CB-CA-CT-O2
2	B	301	D16	CB-CA-CT-O1

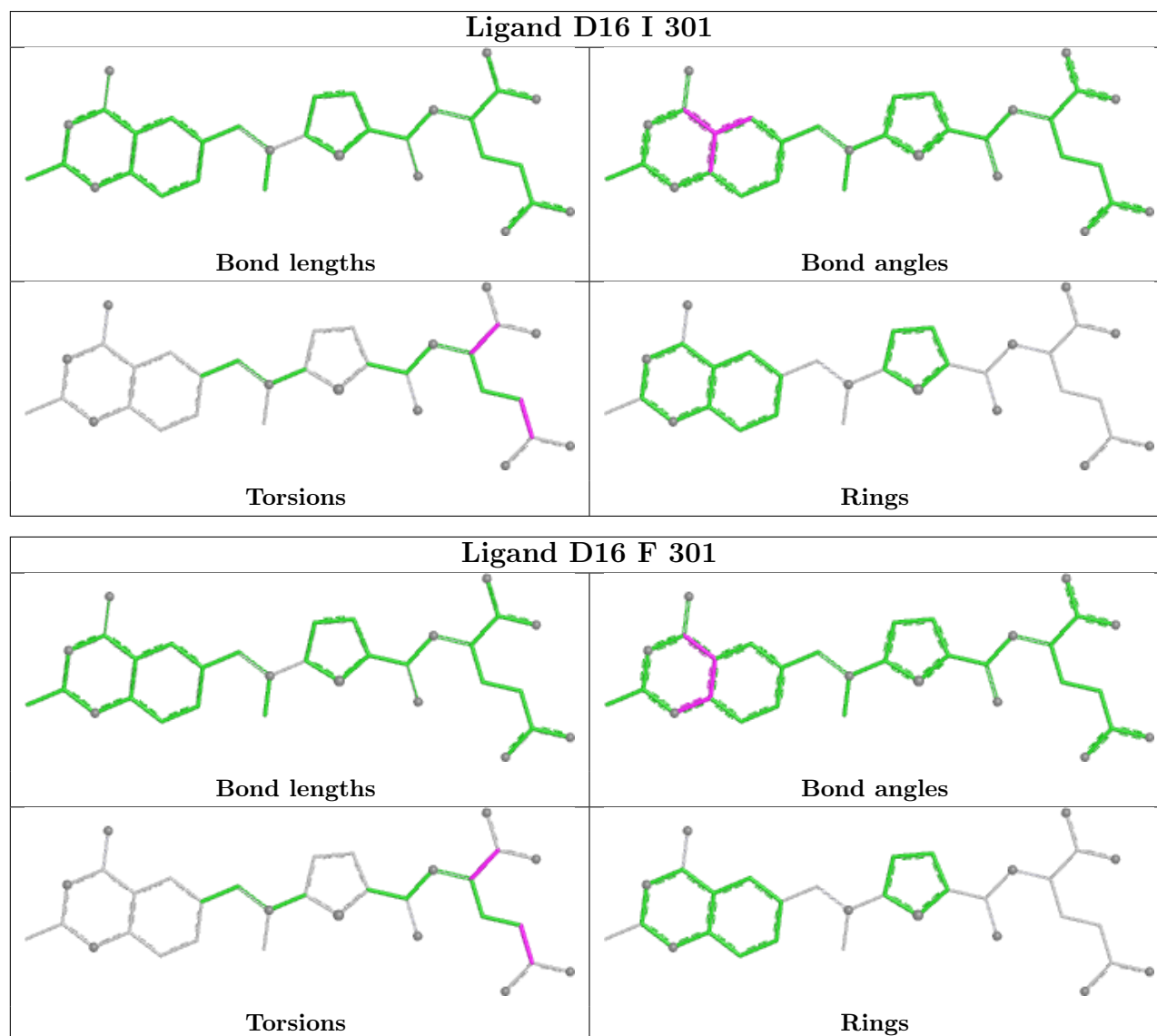
There are no ring outliers.

4 monomers are involved in 7 short contacts:

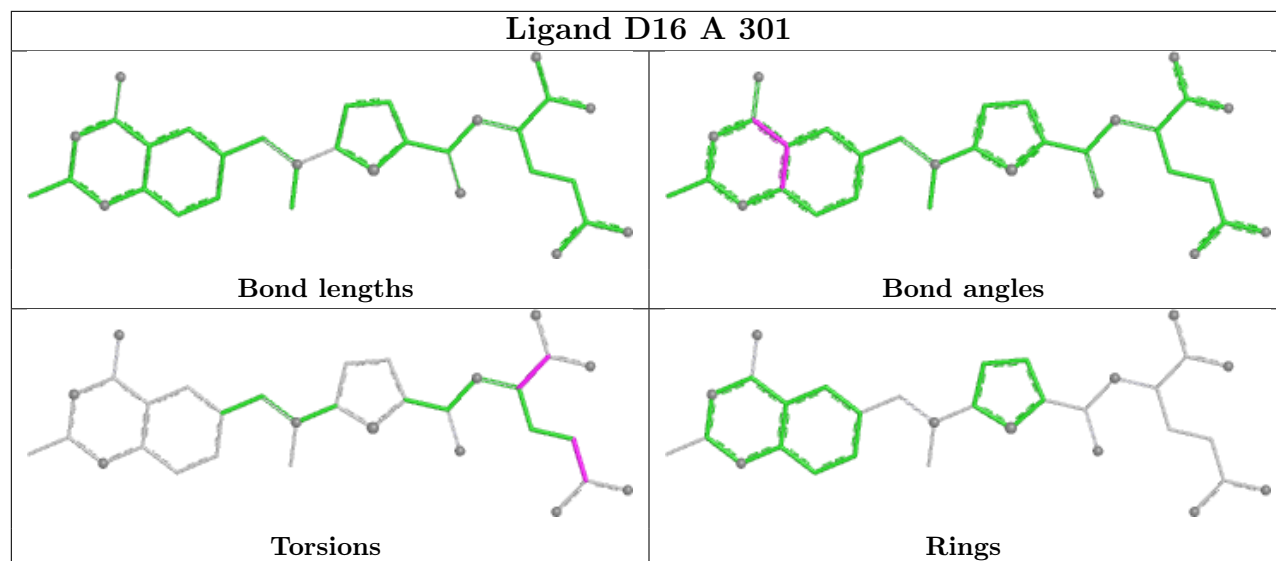
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	D16	2	0
2	H	301	D16	2	0
2	G	301	D16	1	0
2	E	301	D16	2	0

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

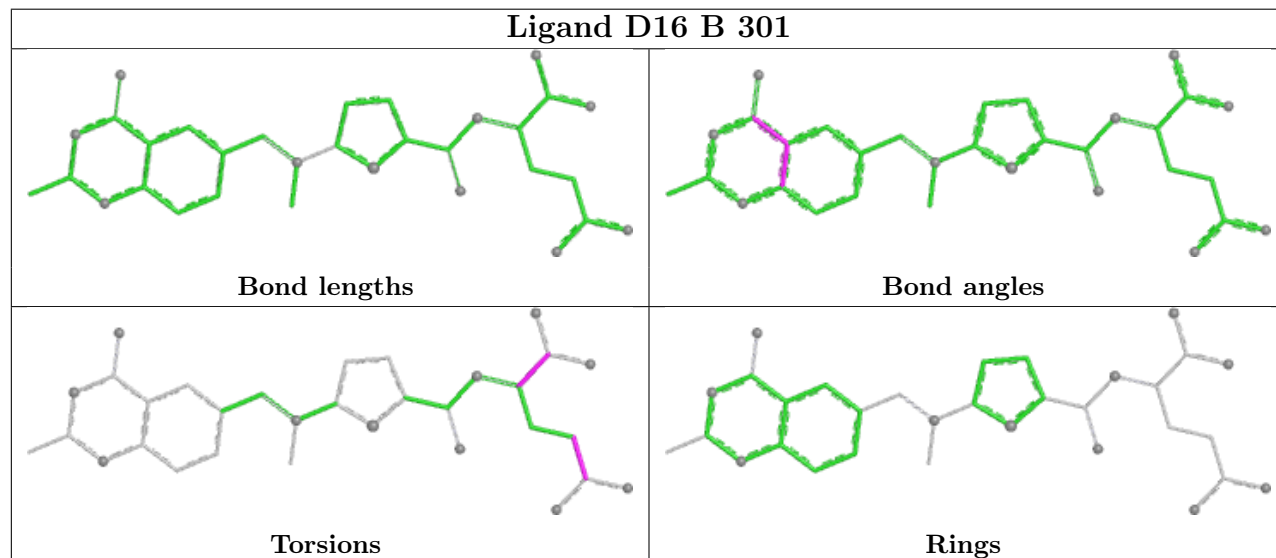
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.



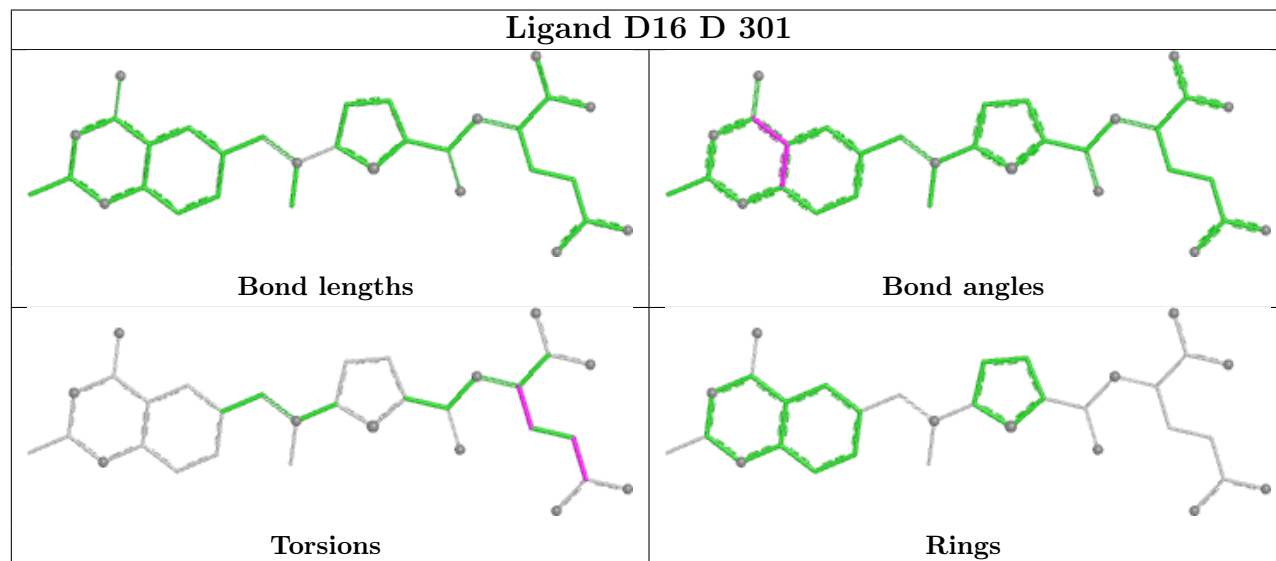
Ligand D16 A 301



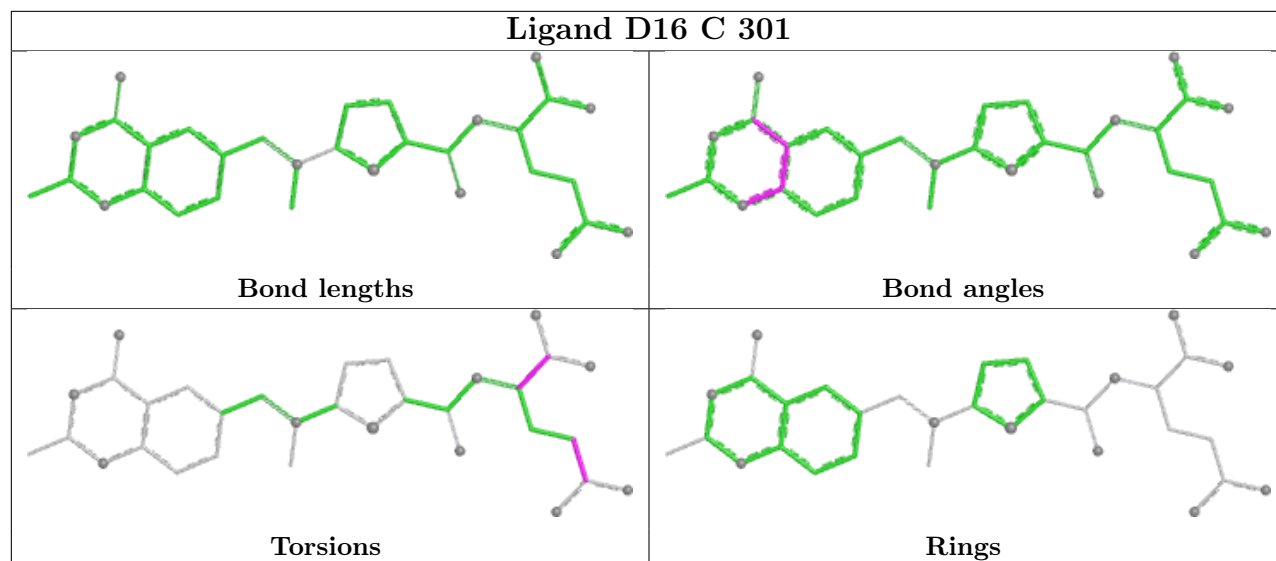
Ligand D16 B 301



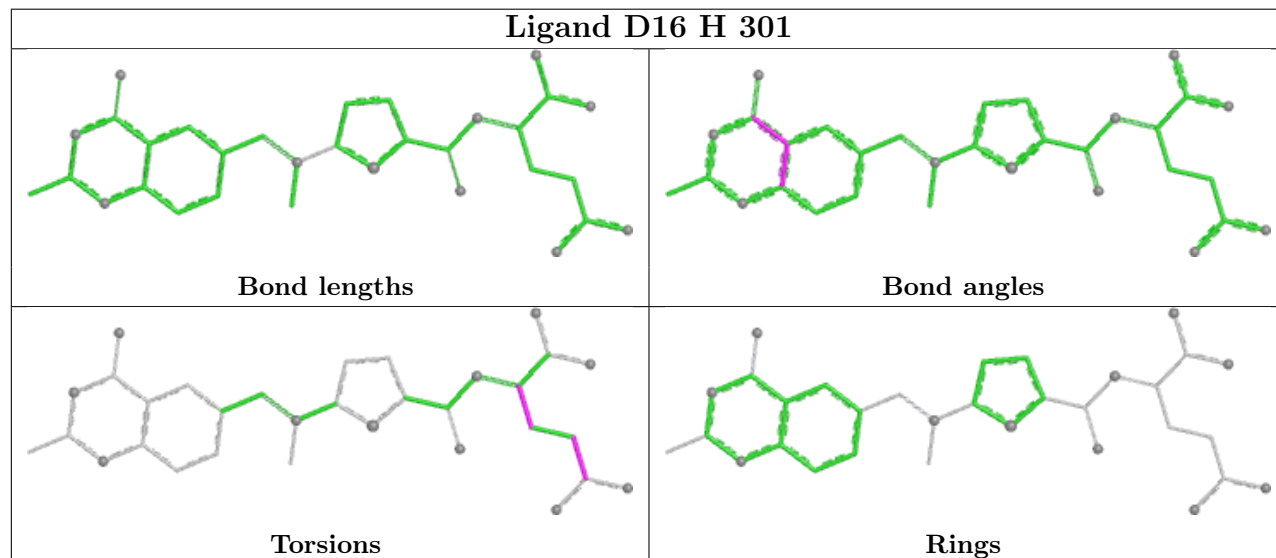
Ligand D16 D 301



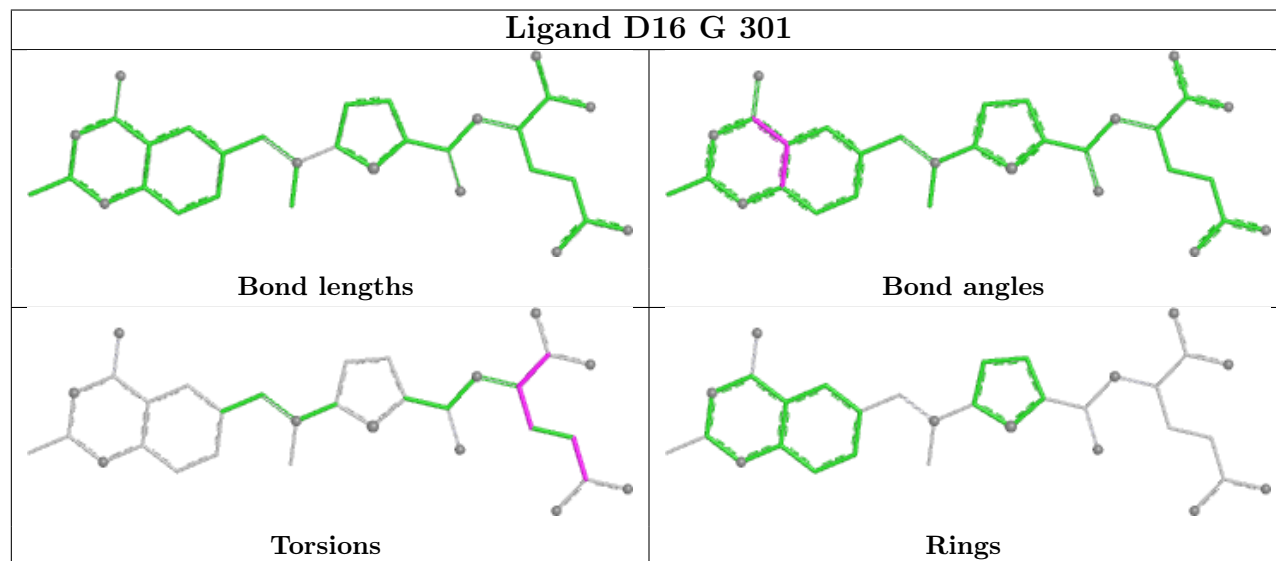
Ligand D16 C 301

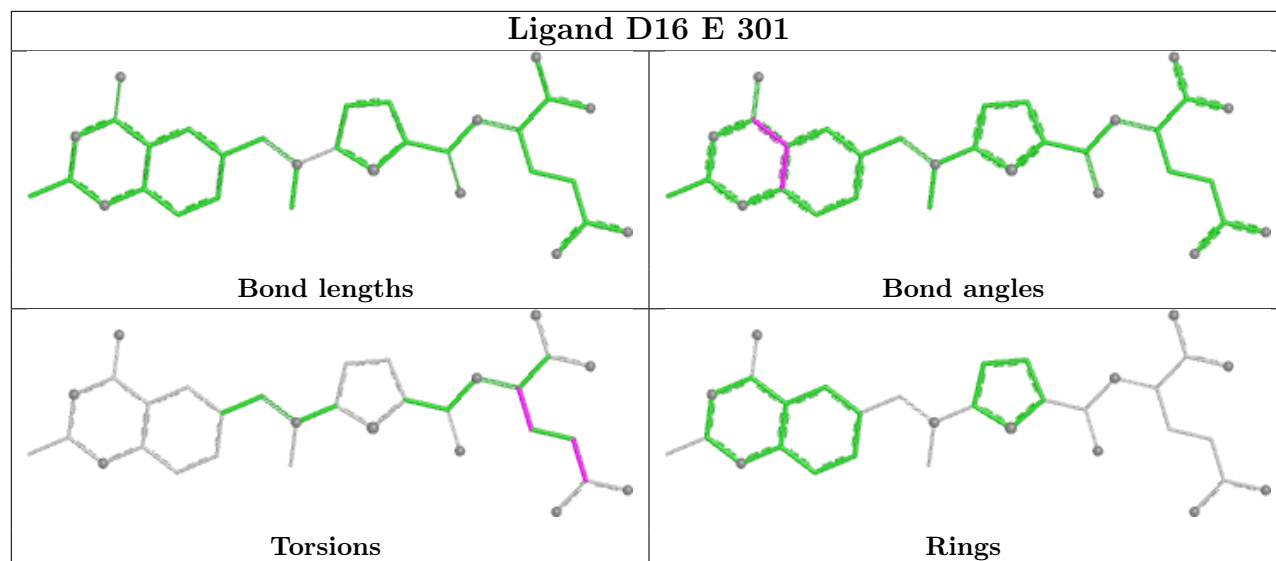


Ligand D16 H 301



Ligand D16 G 301





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	220/235 (93%)	0.59	17 (7%)	19 17	25, 40, 55, 71	0
1	B	223/235 (94%)	0.51	11 (4%)	35 33	21, 37, 54, 73	0
1	C	225/235 (95%)	0.92	33 (14%)	6 4	21, 43, 67, 88	0
1	D	225/235 (95%)	1.17	39 (17%)	4 3	25, 46, 75, 83	0
1	E	222/235 (94%)	0.66	16 (7%)	21 19	24, 42, 60, 69	0
1	F	97/235 (41%)	2.38	52 (53%)	0 0	42, 60, 71, 78	0
1	G	197/235 (83%)	1.32	41 (20%)	2 2	35, 53, 69, 80	0
1	H	216/235 (91%)	1.43	60 (27%)	1 1	34, 51, 74, 82	0
1	I	217/235 (92%)	1.51	66 (30%)	1 1	33, 51, 74, 82	0
All	All	1842/2115 (87%)	1.08	335 (18%)	3 3	21, 46, 71, 88	0

The worst 5 of 335 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	60	ARG	6.8
1	F	178	ILE	6.6
1	F	185	THR	5.6
1	I	53	ILE	5.6
1	F	175	ASP	5.6

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](#)

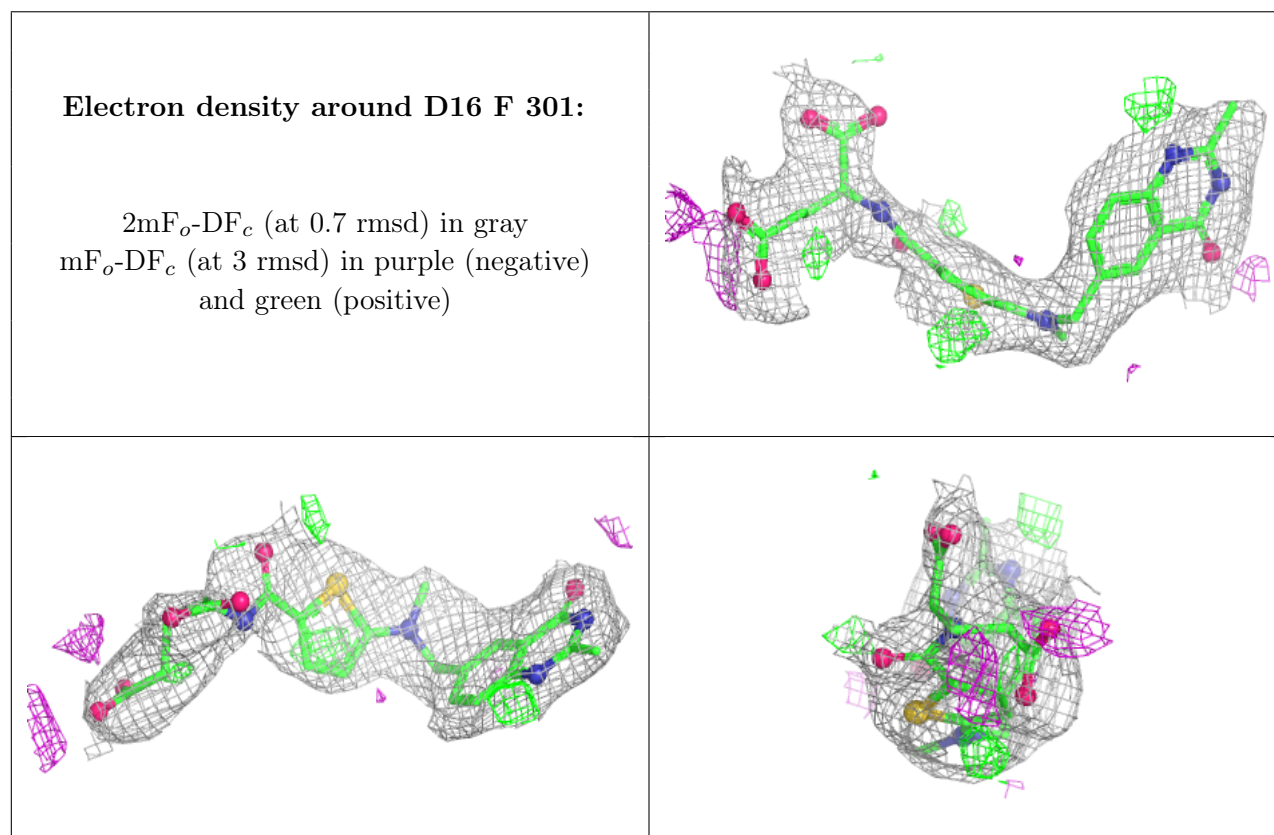
There are no oligosaccharides in this entry.

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.

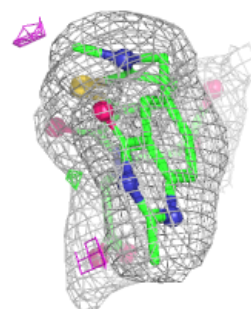
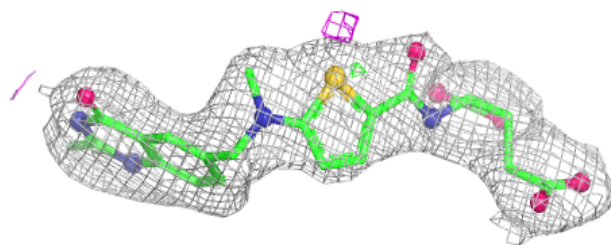
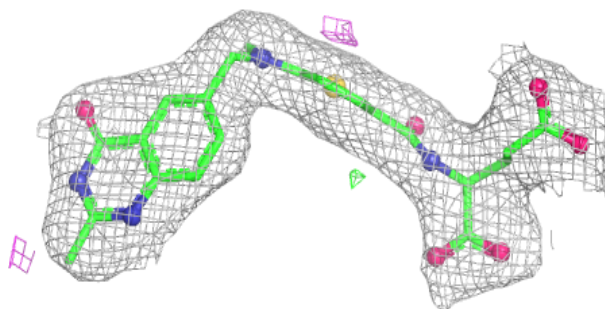
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	D16	F	301	32/32	0.78	0.16	36,48,58,61	0
2	D16	G	301	32/32	0.91	0.10	33,38,46,47	0
2	D16	H	301	32/32	0.91	0.12	32,41,48,51	0
2	D16	I	301	32/32	0.91	0.11	30,43,54,58	0
2	D16	D	301	32/32	0.92	0.09	24,30,36,41	0
2	D16	E	301	32/32	0.92	0.10	25,30,39,40	0
2	D16	C	301	32/32	0.93	0.09	22,27,32,38	0
2	D16	A	301	32/32	0.94	0.08	19,28,33,38	0
2	D16	B	301	32/32	0.95	0.08	17,26,36,40	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.

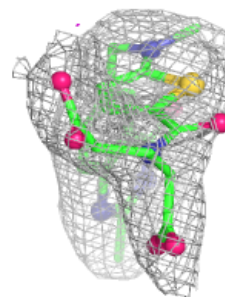
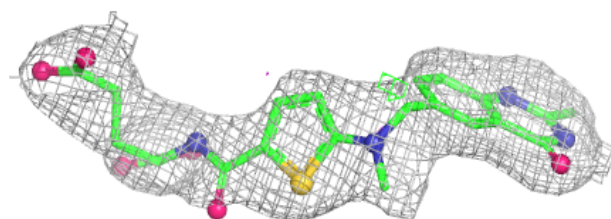
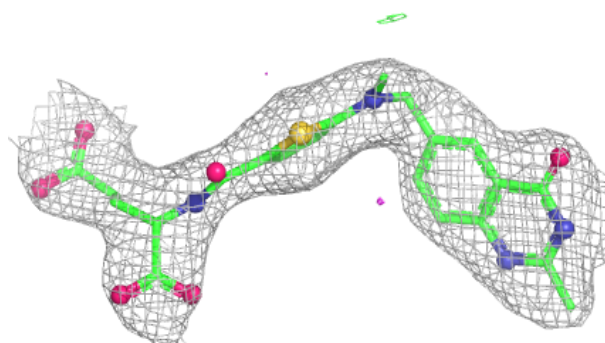


Electron density around D16 G 301:

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

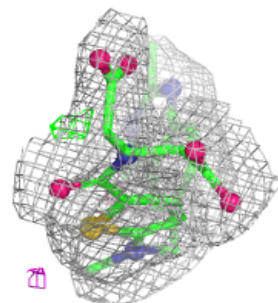
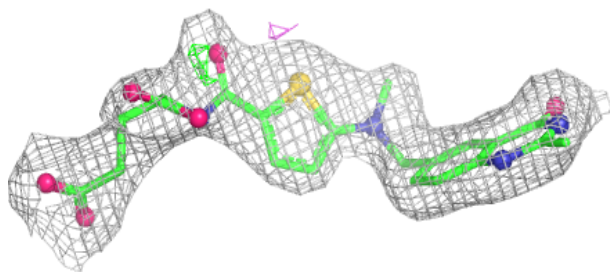
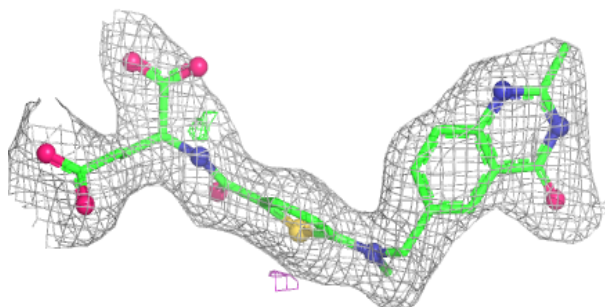
**Electron density around D16 H 301:**

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

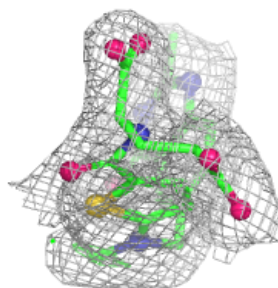
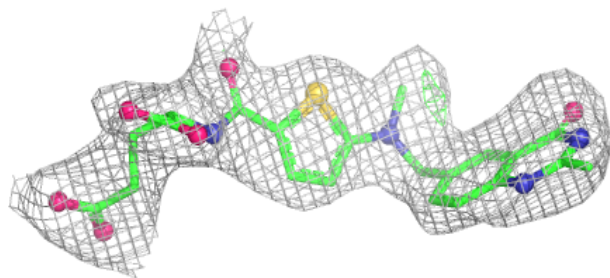
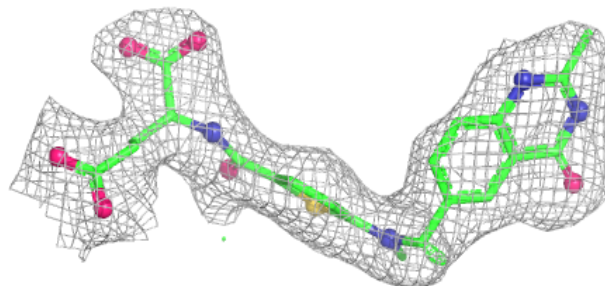


Electron density around D16 I 301:

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

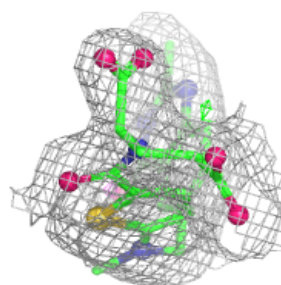
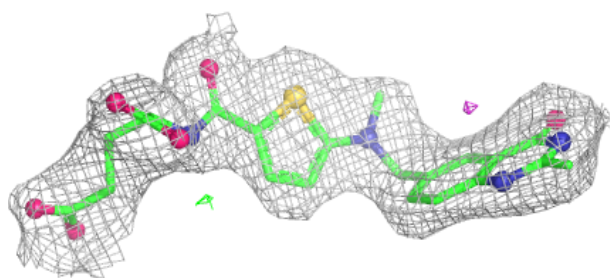
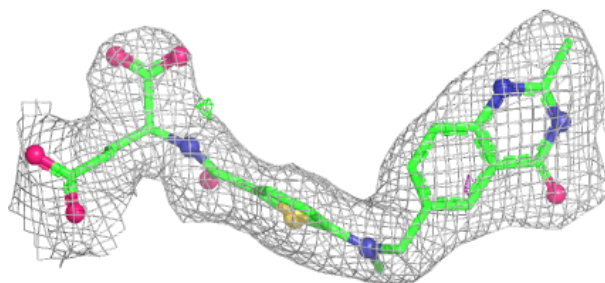
**Electron density around D16 D 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
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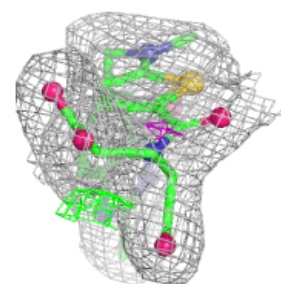
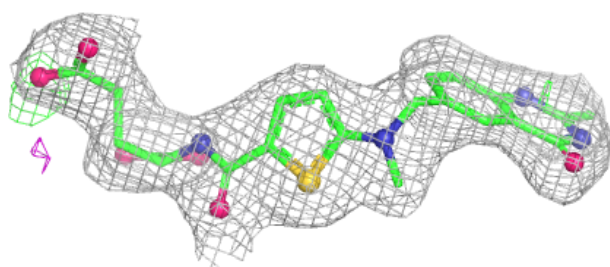
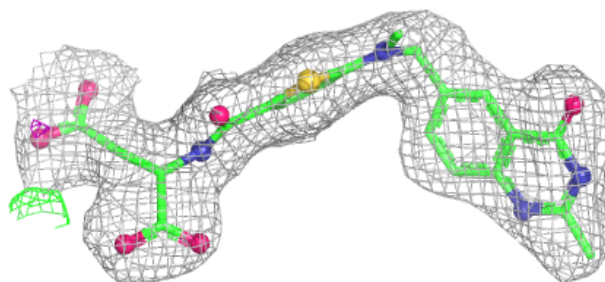


Electron density around D16 E 301:

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

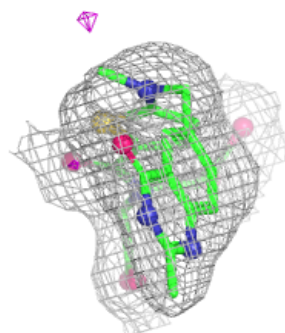
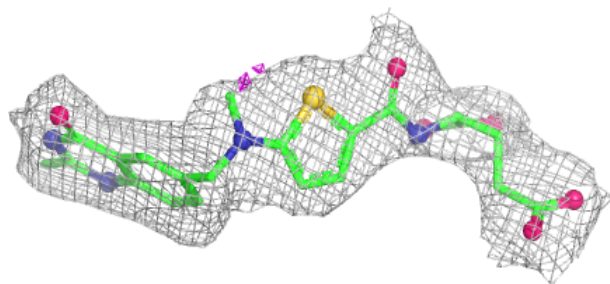
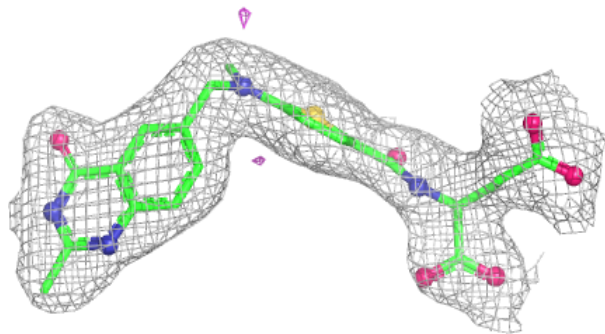
**Electron density around D16 C 301:**

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

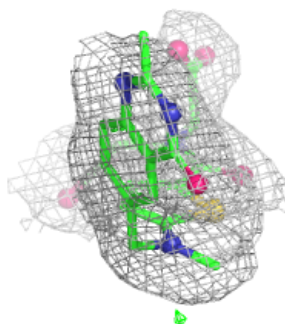
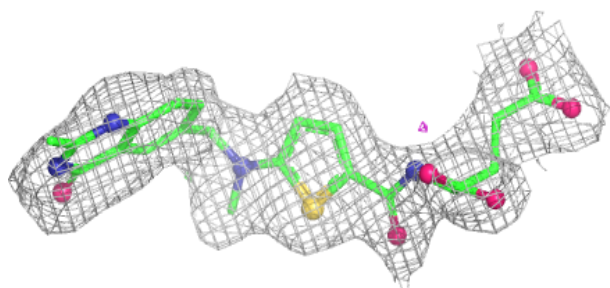
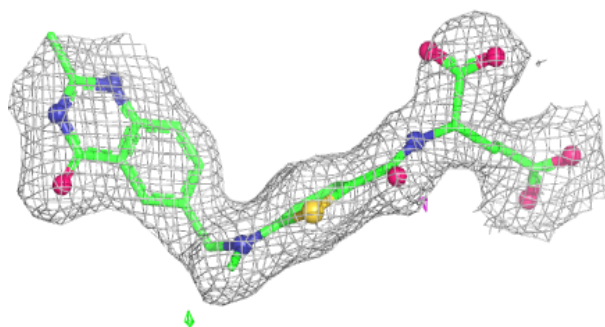


Electron density around D16 A 301:

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

**Electron density around D16 B 301:**

$2mF_o-DF_c$ (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.