



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 12:11 PM UTC

PDB ID : 9HNE / pdb\_00009hne  
Title : Cereblon in complex with DDB1, GSPT1 and Compound-1  
Authors : Klejnot, M.; Walczak, M.J.  
Deposited on : 2024-12-10  
Resolution : 3.90 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](mailto: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>.

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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

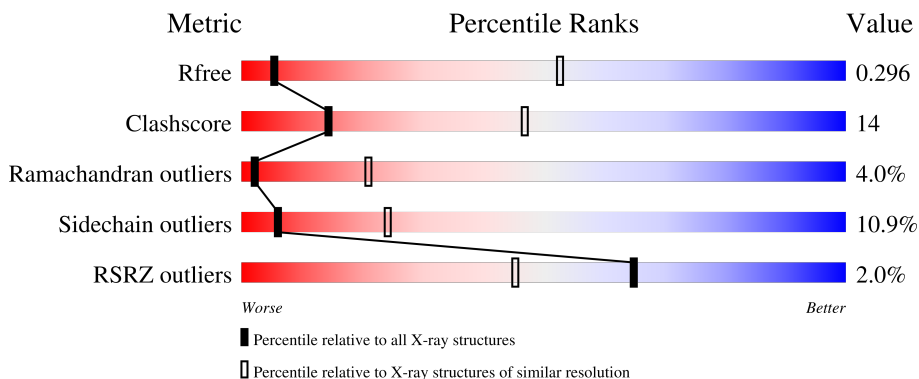
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


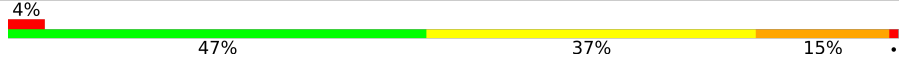



The reported resolution of this entry is 3.90 Å.

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	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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  $\geq 3$ , 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	197	
1	D	197	
2	B	1140	
2	E	1140	
3	C	404	

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Mol	Chain	Length	Quality of chain
3	F	404	<div><div></div><div>2%</div><div>57%</div><div>28%</div><div>6%</div><div>9%</div></div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 26021 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 Eukaryotic peptide chain release factor GTP-binding subunit ERF3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1513	960	261	280	12			
1	D	196	Total	C	N	O	S	0	0	0
			1520	965	262	281	12			

- Molecule 2 is a protein called DNA damage-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1081	Total	C	N	O	S	0	0	0
			8491	5397	1431	1617	46			
2	E	1082	Total	C	N	O	S	0	0	0
			8501	5402	1432	1621	46			

- Molecule 3 is a protein called Protein cereblon.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	371	Total	C	N	O	S	0	0	0
			2996	1910	511	552	23			
3	F	367	Total	C	N	O	S	0	0	0
			2954	1892	502	536	24			

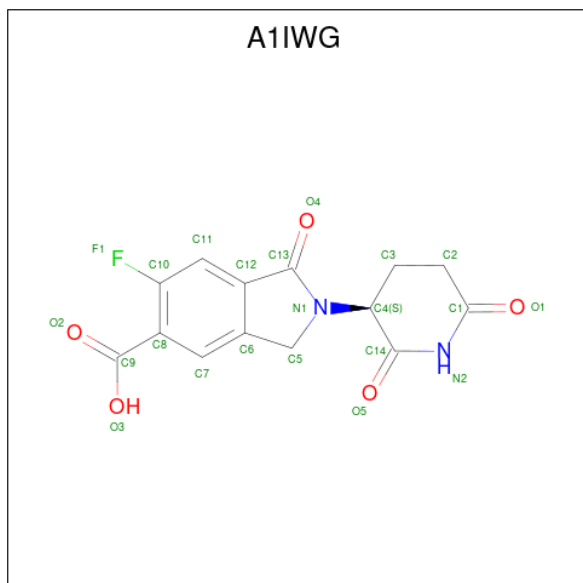
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	39	SER	-	expression tag	UNP Q96SW2
F	39	SER	-	expression tag	UNP Q96SW2

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Zn	0	0
			1	1		
4	F	1	Total	Zn	0	0
			1	1		

- Molecule 5 is 2-[(3 {S})-2,6-bis(oxidanylidene)piperidin-3-yl]-6-fluoranyl-1-oxidanylidene-3 {H}-isoindole-5-carboxylic acid (CCD ID: A1IWG) (formula: C<sub>14</sub>H<sub>11</sub>FN<sub>2</sub>O<sub>5</sub>) (labeled as "Ligand of Interest" by depositor).

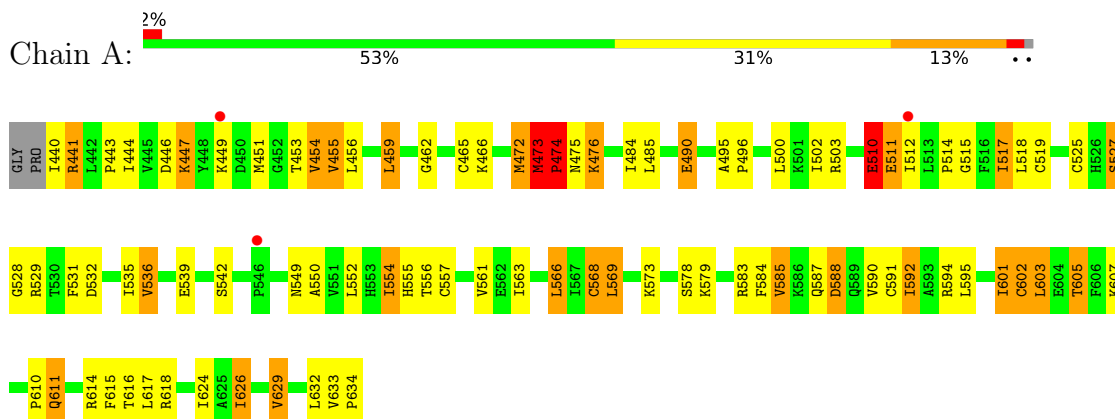


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total	C	F	N	O	0	0
			22	14	1	2	5		
5	F	1	Total	C	F	N	O	0	0
			22	14	1	2	5		

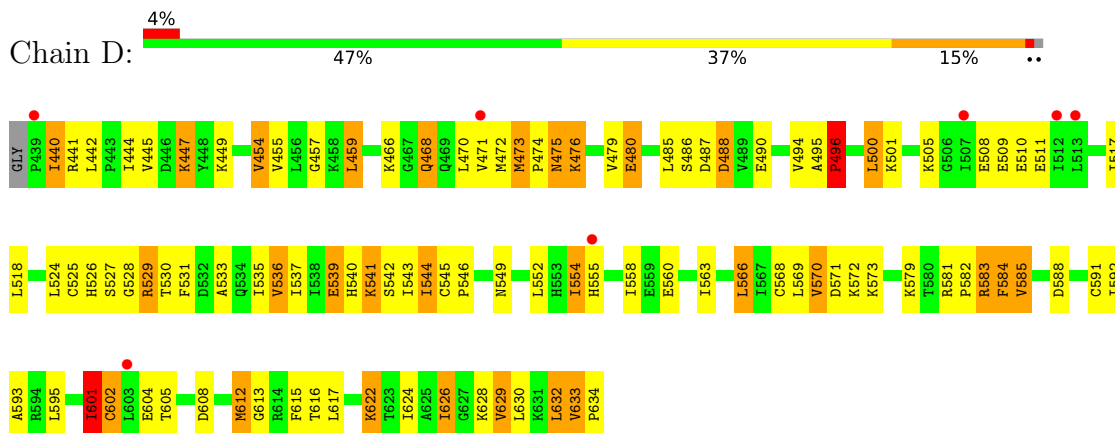
### 3 Residue-property plots [i](#)

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.

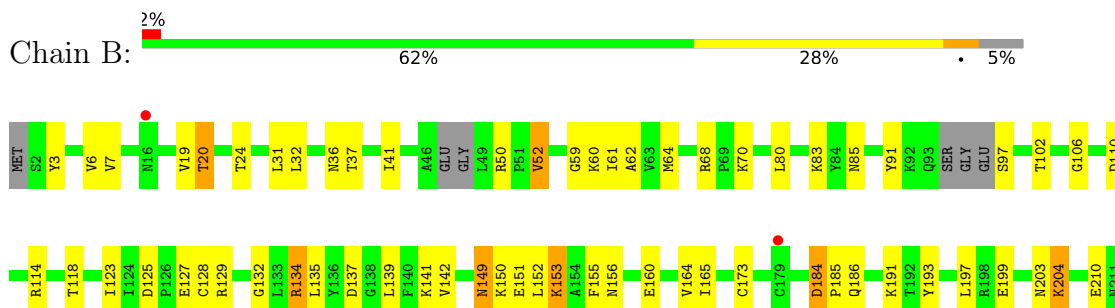
- Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit ERF3A

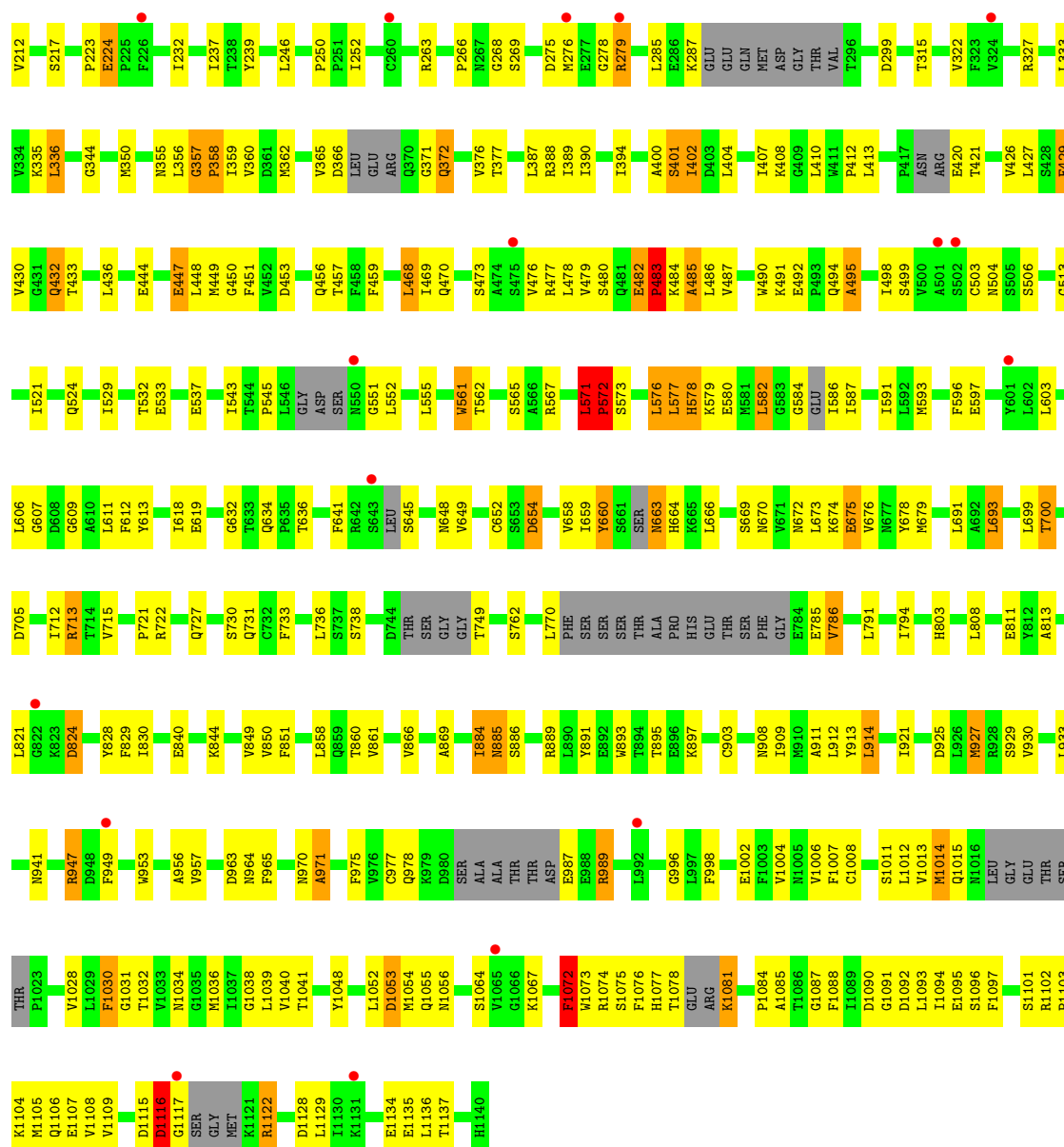


- Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit ERF3A

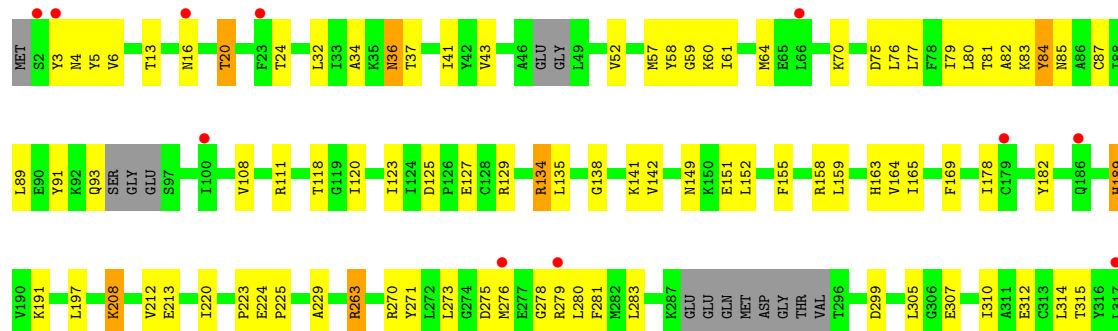


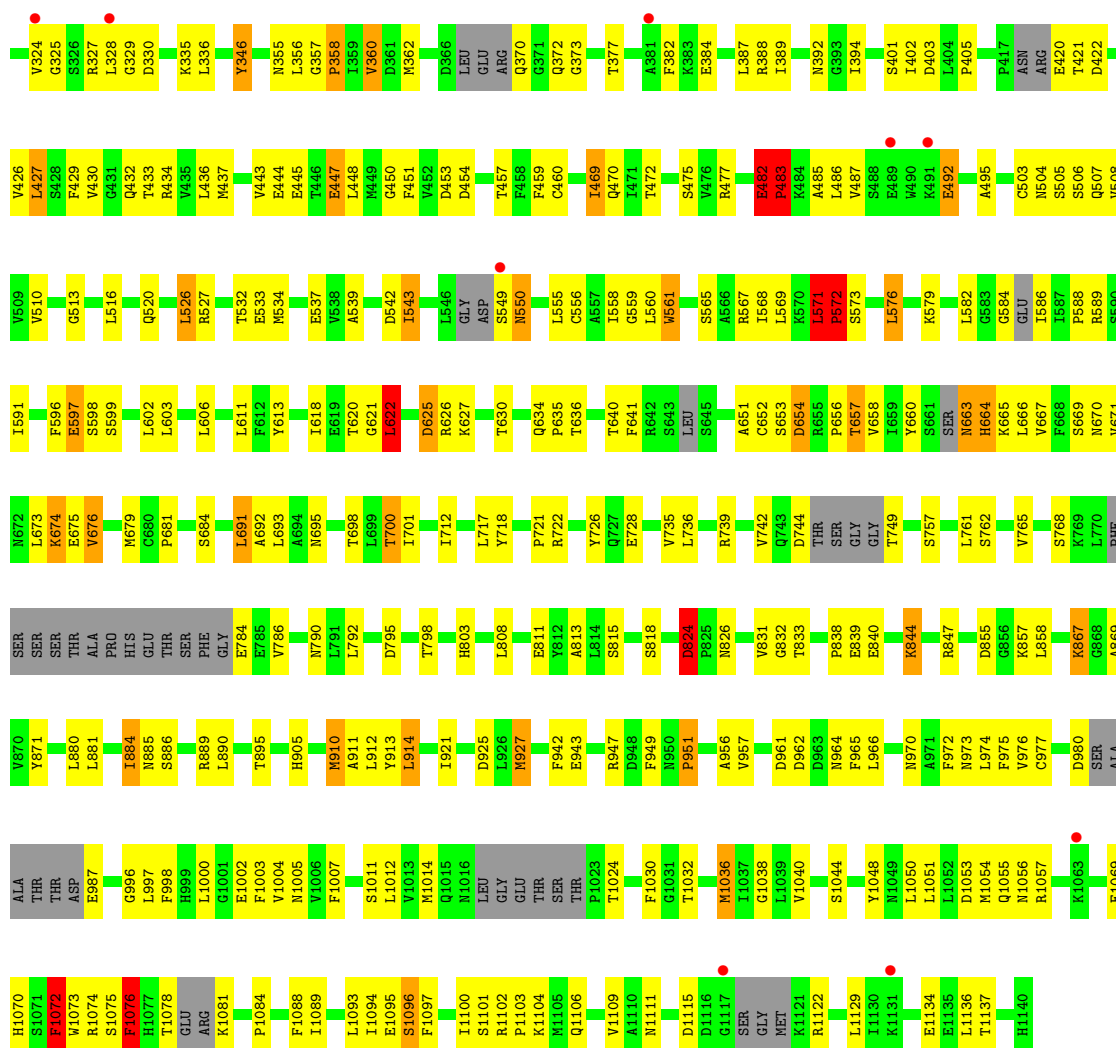
- Molecule 2: DNA damage-binding protein 1



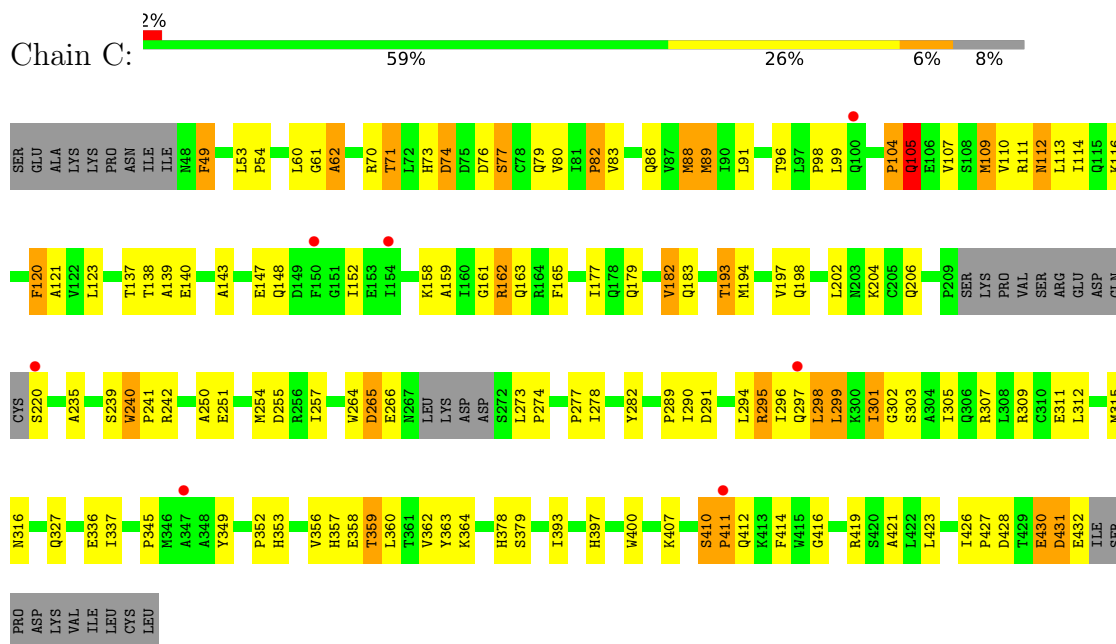


• Molecule 2: DNA damage-binding protein 1

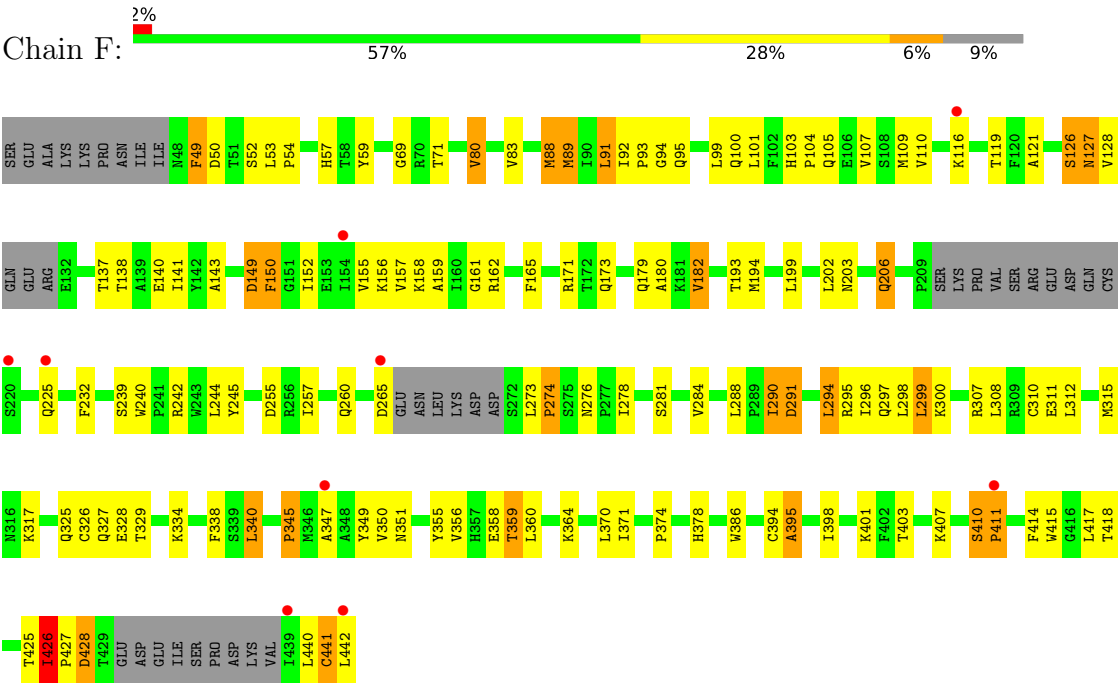




• Molecule 3: Protein cereblon



● Molecule 3: Protein cereblon



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	158.34Å 112.38Å 177.94Å 90.00° 95.25° 90.00°	Depositor
Resolution (Å)	49.49 – 3.90 49.49 – 3.90	Depositor EDS
% Data completeness (in resolution range)	96.5 (49.49-3.90) 96.5 (49.49-3.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.28 (at 3.88Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, $R_{free}$	0.259 , 0.336 (Not available) , 0.296	Depositor DCC
$R_{free}$ test set	2100 reflections (3.81%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	107.4	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 183.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	26021	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	135.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: A1IWG, ZN

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	0.49	0/1534	1.11	1/2068 (0.0%)
1	D	0.49	0/1542	1.11	2/2079 (0.1%)
2	B	0.49	0/8637	1.01	17/11683 (0.1%)
2	E	0.49	0/8647	1.00	14/11696 (0.1%)
3	C	0.48	0/3067	1.06	10/4160 (0.2%)
3	F	0.48	0/3023	1.04	7/4099 (0.2%)
All	All	0.49	0/26450	1.03	51/35785 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6
1	D	0	2
2	B	0	4
2	E	0	3
3	C	0	3
All	All	0	18

There are no bond length outliers.

The worst 5 of 51 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	572	PRO	N-CA-CB	-8.46	94.36	103.25
2	B	572	PRO	N-CA-CB	-7.99	94.86	103.25
2	B	360	VAL	N-CA-CB	-7.69	105.33	111.64
2	B	1072	PHE	CA-CB-CG	7.39	121.19	113.80
2	B	1076	PHE	N-CA-CB	-7.30	99.30	110.04

There are no chirality outliers.

5 of 18 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	472	MET	Peptide
1	A	473	MET	Peptide
1	A	510	GLU	Peptide
1	A	588	ASP	Peptide
1	A	601	ILE	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1513	0	1584	64	0
1	D	1520	0	1592	81	0
2	B	8491	0	8487	202	0
2	E	8501	0	8496	232	0
3	C	2996	0	2971	83	0
3	F	2954	0	2952	89	0
4	C	1	0	0	0	0
4	F	1	0	0	0	0
5	C	22	0	0	2	0
5	F	22	0	0	1	0
All	All	26021	0	26082	734	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 14.

The worst 5 of 734 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)
3:C:74:ASP:O	3:C:77:SER:OG	1.89	0.90
2:B:451:PHE:CE2	2:B:470:GLN:HB2	2.13	0.81
3:F:257:ILE:HG23	3:F:315:MET:HE1	1.62	0.81
2:B:433:THR:OG1	2:B:457:THR:OG1	1.99	0.80
2:E:1078:THR:HG1	2:E:1081:LYS:N	1.81	0.79

There are no symmetry-related clashes.

## 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	193/197 (98%)	152 (79%)	30 (16%)	11 (6%)	1	16
1	D	194/197 (98%)	149 (77%)	36 (19%)	9 (5%)	2	19
2	B	1049/1140 (92%)	884 (84%)	126 (12%)	39 (4%)	2	22
2	E	1050/1140 (92%)	880 (84%)	135 (13%)	35 (3%)	3	24
3	C	365/404 (90%)	287 (79%)	61 (17%)	17 (5%)	2	19
3	F	357/404 (88%)	287 (80%)	53 (15%)	17 (5%)	2	18
All	All	3208/3482 (92%)	2639 (82%)	441 (14%)	128 (4%)	2	21

5 of 128 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	473	MET
1	A	474	PRO
1	A	511	GLU
1	A	555	HIS
1	A	584	PHE

### 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	173/174 (99%)	140 (81%)	33 (19%)	1	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	174/174 (100%)	137 (79%)	37 (21%)	1	6
2	B	951/999 (95%)	865 (91%)	86 (9%)	9	31
2	E	953/999 (95%)	857 (90%)	96 (10%)	7	27
3	C	335/367 (91%)	298 (89%)	37 (11%)	6	24
3	F	331/367 (90%)	301 (91%)	30 (9%)	9	31
All	All	2917/3080 (95%)	2598 (89%)	319 (11%)	6	24

5 of 319 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	533	GLU
2	E	1106	GLN
2	E	582	LEU
2	E	790	ASN
3	F	182	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 59 such sidechains are listed below:

Mol	Chain	Res	Type
2	E	262	ASN
3	F	206	GLN
2	E	578	HIS
3	F	198	GLN
3	F	103	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

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z  > 2$
5	A1IWG	F	502	-	24,24,24	0.73	0	36,36,36	0.96	3 (8%)
5	A1IWG	C	502	-	24,24,24	0.72	0	36,36,36	0.87	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
5	A1IWG	F	502	-	-	4/8/33/33	0/3/3/3
5	A1IWG	C	502	-	-	0/8/33/33	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
5	F	502	A1IWG	C11-C10-C8	-2.36	120.87	123.48
5	F	502	A1IWG	C14-C4-N1	-2.32	108.18	110.32
5	F	502	A1IWG	C7-C8-C10	2.23	119.59	116.83
5	C	502	A1IWG	C7-C8-C10	2.00	119.31	116.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

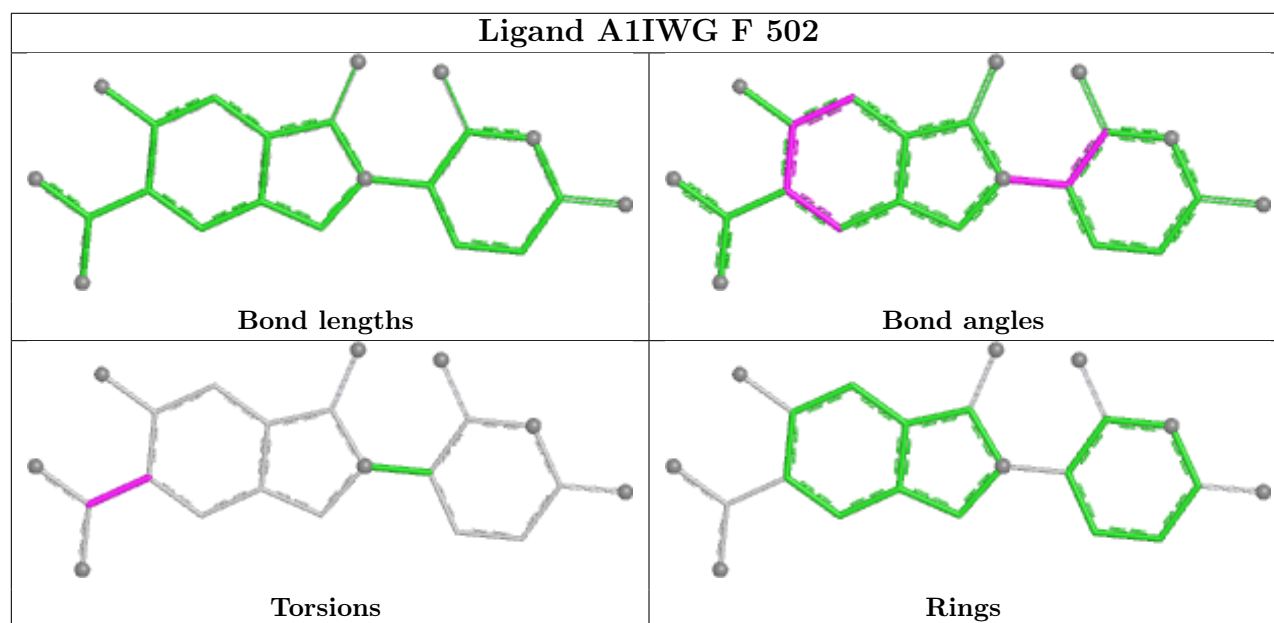
Mol	Chain	Res	Type	Atoms
5	F	502	A1IWG	C10-C8-C9-O2
5	F	502	A1IWG	C10-C8-C9-O3
5	F	502	A1IWG	C7-C8-C9-O2
5	F	502	A1IWG	C7-C8-C9-O3

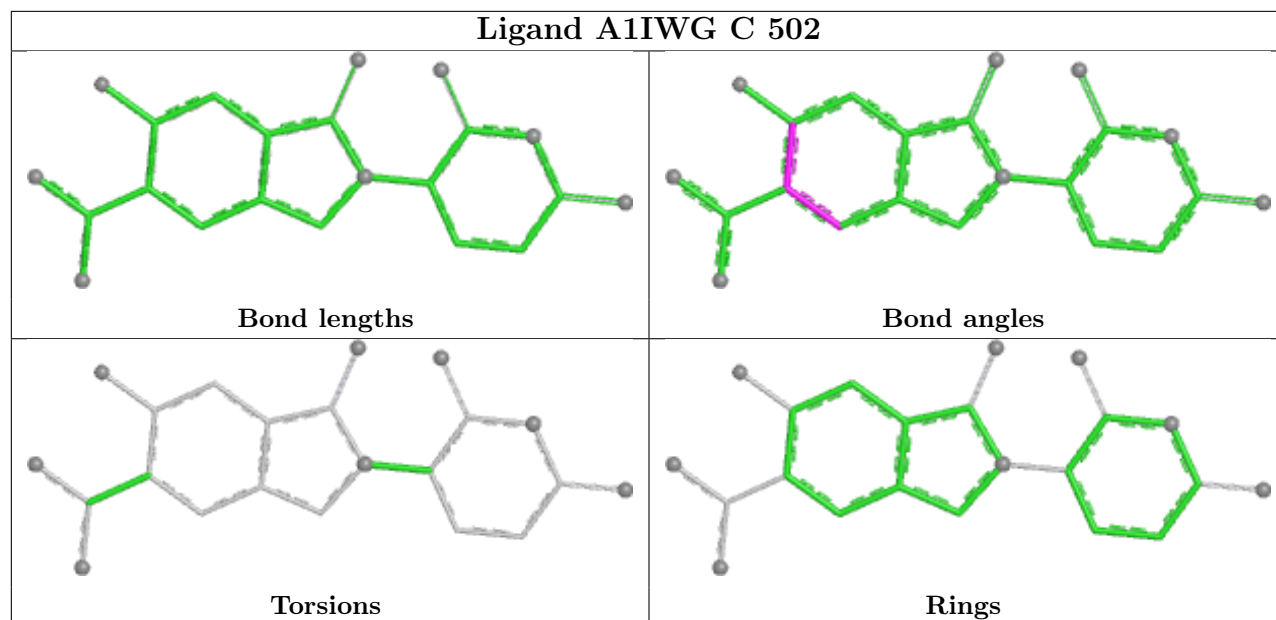
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	502	A1IWG	1	0
5	C	502	A1IWG	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.





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	195/197 (98%)	0.19	3 (1%) 72 50	69, 145, 271, 323	0
1	D	196/197 (99%)	0.25	7 (3%) 46 33	60, 143, 239, 326	0
2	B	1081/1140 (94%)	0.01	19 (1%) 67 47	31, 125, 203, 284	5 (0%)
2	E	1082/1140 (94%)	0.08	20 (1%) 67 47	31, 130, 221, 308	5 (0%)
3	C	371/404 (91%)	0.05	7 (1%) 66 46	63, 122, 211, 291	0
3	F	367/404 (90%)	0.09	9 (2%) 58 41	59, 122, 198, 262	0
All	All	3292/3482 (94%)	0.07	65 (1%) 65 45	31, 128, 219, 326	10 (0%)

The worst 5 of 65 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	226	PHE	4.4
2	B	260	CYS	4.0
2	E	179	CYS	4.0
1	D	555	HIS	3.9
2	E	276	MET	3.9

### 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](#)

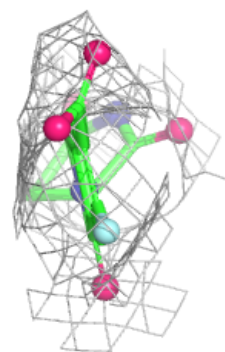
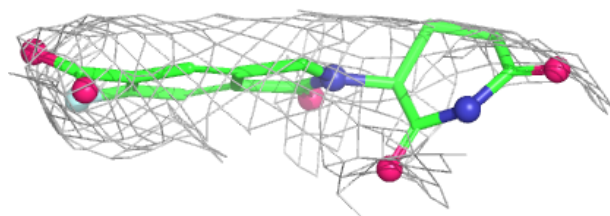
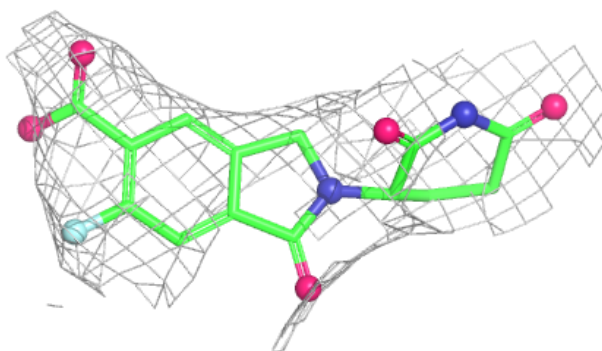
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	A1IWG	F	502	22/22	0.95	0.08	72,88,105,113	0
5	A1IWG	C	502	22/22	0.96	0.08	69,88,107,115	0
4	ZN	C	501	1/1	1.00	0.03	113,113,113,113	0
4	ZN	F	501	1/1	1.00	0.02	100,100,100,100	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.

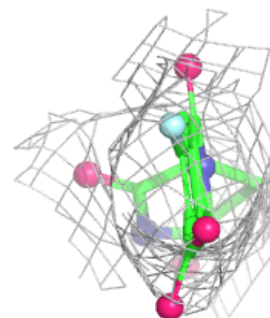
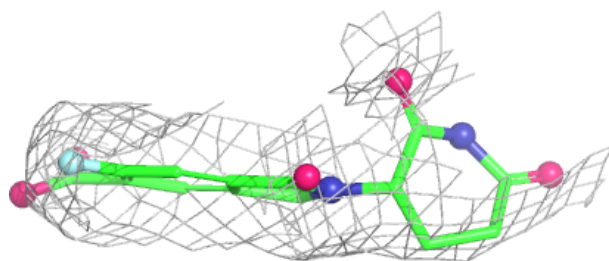
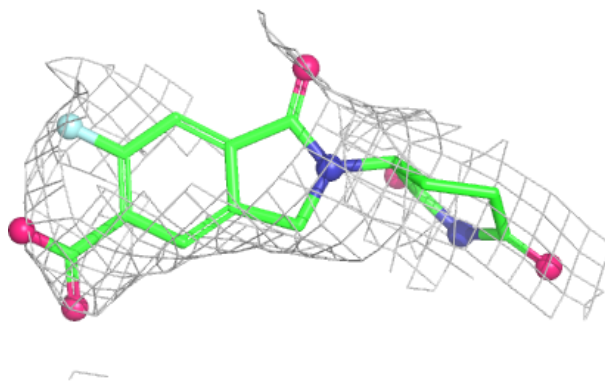
### Electron density around A1IWG F 502:

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around A1IWG C 502:**

$2mF_o - DF_c$  (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.