



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 09:03 PM UTC

PDB ID : 6BL2 / pdb_00006bl2
Title : Novel Modes of Inhibition of Wild-Type IDH1: Direct Covalent Modification of His315 with Cmpd15
Authors : Jakob, C.G.; Qiu, W.
Deposited on : 2017-11-09
Resolution : 1.92 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

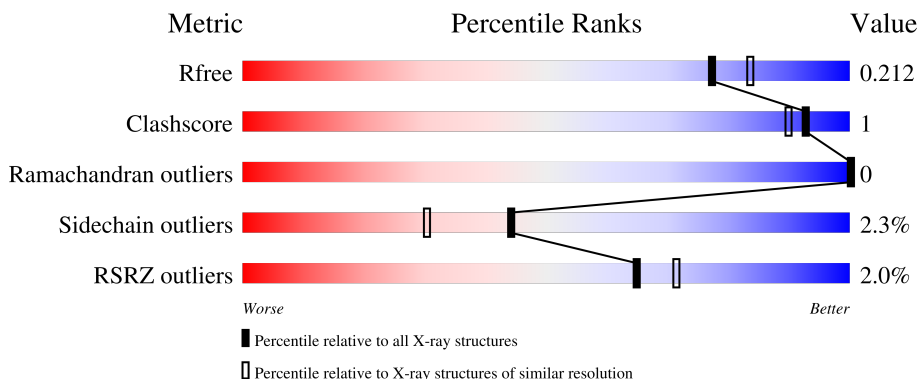
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

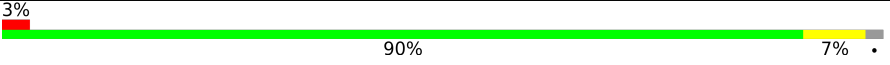
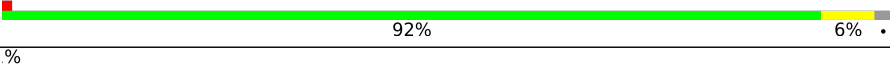
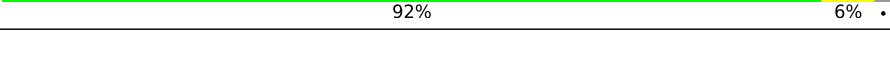
The reported resolution of this entry is 1.92 Å.

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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	425	 3% 90% 7%
1	B	425	 0% 92% 6%
1	C	425	 0% 92% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10885 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 Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3292	2093	559	622	18			
1	B	416	Total	C	N	O	S	0	0	0
			3299	2097	560	624	18			
1	C	416	Total	C	N	O	S	0	0	0
			3292	2093	559	622	18			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	UNP O75874
A	422	HIS	-	expression tag	UNP O75874
A	423	HIS	-	expression tag	UNP O75874
A	424	HIS	-	expression tag	UNP O75874
A	425	HIS	-	expression tag	UNP O75874
B	415	SER	-	expression tag	UNP O75874
B	416	LEU	-	expression tag	UNP O75874
B	417	GLU	-	expression tag	UNP O75874
B	418	HIS	-	expression tag	UNP O75874
B	419	HIS	-	expression tag	UNP O75874
B	420	HIS	-	expression tag	UNP O75874
B	421	HIS	-	expression tag	UNP O75874
B	422	HIS	-	expression tag	UNP O75874
B	423	HIS	-	expression tag	UNP O75874
B	424	HIS	-	expression tag	UNP O75874
B	425	HIS	-	expression tag	UNP O75874
C	415	SER	-	expression tag	UNP O75874

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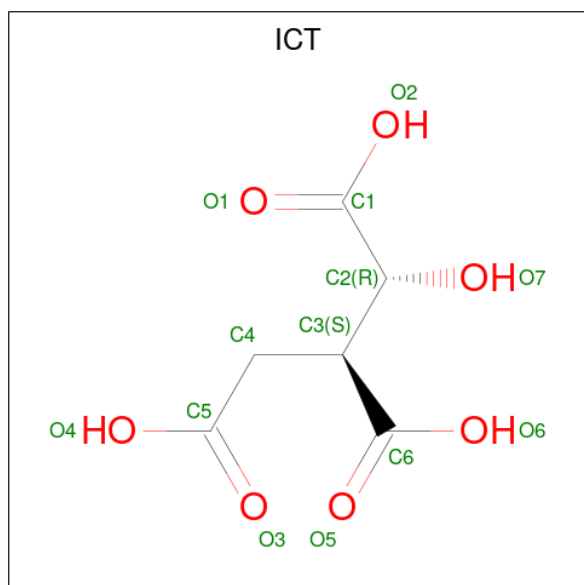
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Chain	Residue	Modelled	Actual	Comment	Reference
C	416	LEU	-	expression tag	UNP O75874
C	417	GLU	-	expression tag	UNP O75874
C	418	HIS	-	expression tag	UNP O75874
C	419	HIS	-	expression tag	UNP O75874
C	420	HIS	-	expression tag	UNP O75874
C	421	HIS	-	expression tag	UNP O75874
C	422	HIS	-	expression tag	UNP O75874
C	423	HIS	-	expression tag	UNP O75874
C	424	HIS	-	expression tag	UNP O75874
C	425	HIS	-	expression tag	UNP O75874

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Ca 2 2	0	0
2	B	1	Total Ca 1 1	0	0

- Molecule 3 is ISOCITRIC ACID (CCD ID: ICT) (formula: C₆H₈O₇).



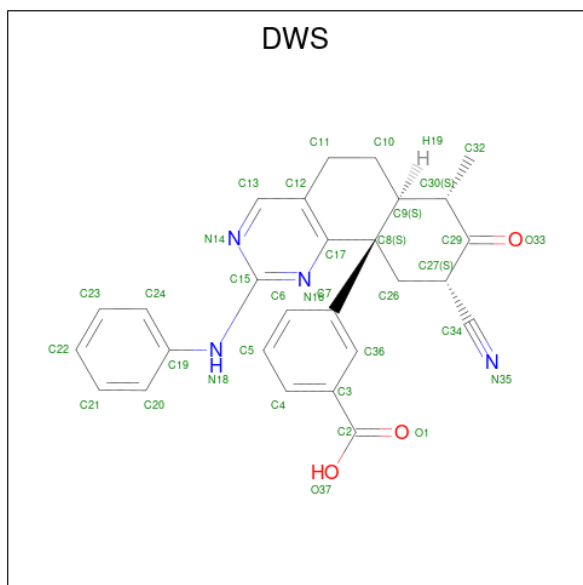
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 13 6 7	0	0
3	A	1	Total C O 13 6 7	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is 3-[(6aS,7S,9S,10aS)-9-cyano-7-methyl-8-oxo-2-(phenylamino)-6,6a,7,8,9,10-hexahydrobenzo[h]quinazolin-10a(5H)-yl]benzoic acid (CCD ID: DWS) (formula: C₂₇H₂₄N₄O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			34	27	4	3		
4	B	1	Total	C	N	O	0	0
			34	27	4	3		
4	C	1	Total	C	N	O	0	0
			34	27	4	3		

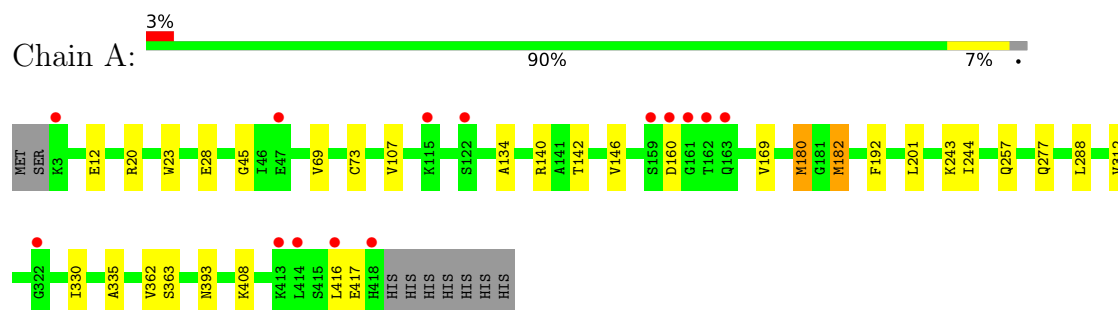
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	230	Total	O	0	0
			230	230		
5	B	331	Total	O	0	0
			331	331		
5	C	297	Total	O	0	0
			297	297		

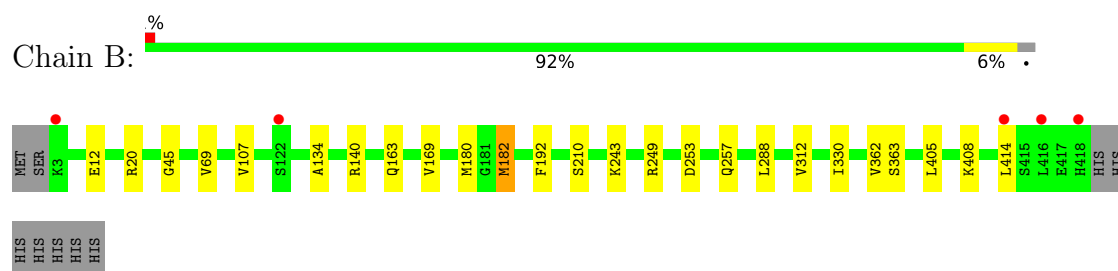
3 Residue-property plots [i](#)

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.

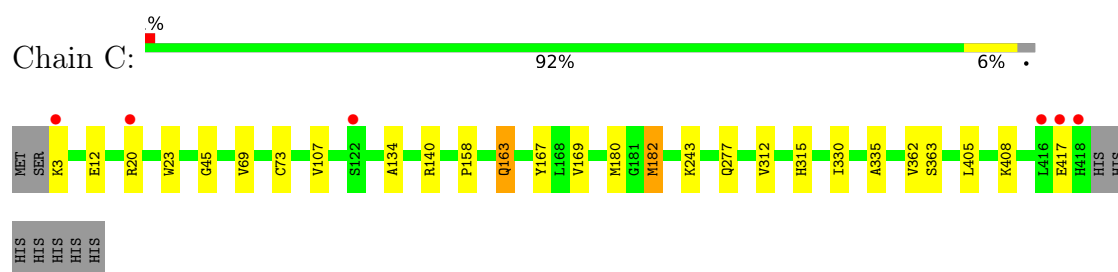
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	95.51Å 274.89Å 116.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.14 – 1.92 28.14 – 1.92	Depositor EDS
% Data completeness (in resolution range)	100.0 (28.14-1.92) 99.9 (28.14-1.92)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 1.92Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.180 , 0.211 0.182 , 0.212	Depositor DCC
R_{free} test set	5851 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.691	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10885	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% of the height of the origin peak. No significant pseudotranslation is detected.*

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: ICT, CA, DWS

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	0.81	2/3361 (0.1%)	1.15	7/4533 (0.2%)
1	B	0.84	1/3368 (0.0%)	1.12	3/4541 (0.1%)
1	C	0.83	1/3361 (0.0%)	1.12	3/4533 (0.1%)
All	All	0.83	4/10090 (0.0%)	1.13	13/13607 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	182	MET	SD-CE	-5.60	1.65	1.79
1	B	182	MET	SD-CE	-5.51	1.65	1.79
1	A	180	MET	SD-CE	-5.49	1.65	1.79
1	C	182	MET	SD-CE	-5.49	1.65	1.79

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	169	VAL	N-CA-C	-7.11	104.84	111.45
1	A	169	VAL	N-CA-C	-6.75	105.17	111.45
1	C	169	VAL	N-CA-C	-6.51	105.39	111.45
1	A	160	ASP	CA-CB-CG	5.83	118.43	112.60
1	C	335	ALA	N-CA-C	-5.75	105.09	111.36
1	A	28	GLU	CB-CG-CD	5.39	121.76	112.60
1	C	417	GLU	CB-CG-CD	5.37	121.73	112.60
1	B	253	ASP	CA-CB-CG	5.30	117.90	112.60
1	A	335	ALA	N-CA-C	-5.21	105.60	111.28
1	A	192	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	146	VAL	N-CA-C	-5.05	102.97	107.56
1	B	192	PHE	CA-CB-CG	5.03	118.83	113.80
1	A	393	ASN	CA-CB-CG	5.03	117.63	112.60

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	3292	0	3255	9	0
1	B	3299	0	3268	9	0
1	C	3292	0	3255	11	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	26	0	8	0	0
3	B	13	0	4	0	0
4	A	34	0	0	0	0
4	B	34	0	0	1	0
4	C	34	0	0	2	0
5	A	230	0	0	0	0
5	B	331	0	0	0	0
5	C	297	0	0	0	0
All	All	10885	0	9790	29	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 1.

All (29) 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:249:ARG:HE	1:B:257:GLN:HE22	1.09	0.99
1:A:180:MET:HE2	1:A:182:MET:HE2	1.67	0.74
1:C:180:MET:HE2	1:C:182:MET:HE2	1.72	0.71
1:B:180:MET:HE2	1:B:182:MET:HE2	1.74	0.69
1:A:142:THR:HG21	1:C:167:TYR:HB3	1.84	0.60
1:A:277:GLN:HE22	1:C:277:GLN:HE22	1.50	0.58
1:B:249:ARG:HE	1:B:257:GLN:NE2	1.93	0.55
1:B:107:VAL:HG23	1:B:134:ALA:HB2	1.94	0.49
1:C:107:VAL:HG23	1:C:134:ALA:HB2	1.96	0.47
4:C:501:DWS:C27	4:C:501:DWS:C6	2.93	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:VAL:HG23	1:B:408:LYS:HD2	1.98	0.46
1:C:20:ARG:HH21	1:C:45:GLY:HA3	1.81	0.45
1:A:20:ARG:HH21	1:A:45:GLY:HA3	1.81	0.45
1:A:107:VAL:HG23	1:A:134:ALA:HB2	1.98	0.45
1:A:362:VAL:HG23	1:A:408:LYS:HD2	1.98	0.45
1:B:20:ARG:HH21	1:B:45:GLY:HA3	1.81	0.45
1:C:362:VAL:HG23	1:C:408:LYS:HD2	2.00	0.44
1:C:330:ILE:HD12	1:C:363:SER:HB3	2.00	0.43
1:A:330:ILE:HD12	1:A:363:SER:HB3	2.01	0.43
4:B:503:DWS:C27	4:B:503:DWS:C6	2.97	0.42
1:A:23:TRP:CD2	1:A:73:CYS:HB2	2.55	0.41
1:C:23:TRP:CD2	1:C:73:CYS:HB2	2.55	0.41
1:B:330:ILE:HD12	1:B:363:SER:HB3	2.02	0.41
1:C:158:PRO:HG2	1:C:163:GLN:HG3	2.03	0.41
1:B:362:VAL:HG21	1:B:405:LEU:HA	2.02	0.41
1:C:315:HIS:CE1	4:C:501:DWS:N16	2.89	0.41
1:B:210:SER:HA	1:B:249:ARG:O	2.21	0.41
1:A:201:LEU:HD23	1:A:244:ILE:HD11	2.04	0.40
1:C:362:VAL:HG21	1:C:405:LEU:HA	2.04	0.40

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	414/425 (97%)	399 (96%)	15 (4%)	0	100	100
1	B	414/425 (97%)	403 (97%)	11 (3%)	0	100	100
1	C	414/425 (97%)	403 (97%)	11 (3%)	0	100	100
All	All	1242/1275 (97%)	1205 (97%)	37 (3%)	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	350/361 (97%)	341 (97%)	9 (3%)	40	24
1	B	352/361 (98%)	344 (98%)	8 (2%)	44	30
1	C	350/361 (97%)	343 (98%)	7 (2%)	48	35
All	All	1052/1083 (97%)	1028 (98%)	24 (2%)	44	30

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLU
1	A	69	VAL
1	A	140	ARG
1	A	243	LYS
1	A	257	GLN
1	A	288	LEU
1	A	312	VAL
1	A	416	LEU
1	A	417	GLU
1	B	12	GLU
1	B	69	VAL
1	B	140	ARG
1	B	163	GLN
1	B	243	LYS
1	B	288	LEU
1	B	312	VAL
1	B	414	LEU
1	C	3	LYS
1	C	12	GLU
1	C	69	VAL
1	C	140	ARG
1	C	163	GLN
1	C	243	LYS
1	C	312	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	68	ASN
1	A	185	GLN
1	A	242	GLN
1	A	277	GLN
1	A	283	GLN
1	A	323	GLN
1	B	14	GLN
1	B	55	GLN
1	B	68	ASN
1	B	163	GLN
1	B	257	GLN
1	B	323	GLN
1	C	14	GLN
1	C	68	ASN
1	C	163	GLN
1	C	228	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](#)

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	$\# Z > 2$
3	ICT	B	502	2	12,12,12	1.15	0	13,16,16	1.57	2 (15%)
3	ICT	A	502	2	12,12,12	1.27	2 (16%)	13,16,16	1.86	5 (38%)
4	DWS	B	503	1	36,38,38	1.73	7 (19%)	44,56,56	1.66	5 (11%)
4	DWS	A	504	1	36,38,38	0.81	2 (5%)	44,56,56	1.81	9 (20%)
3	ICT	A	505	2	12,12,12	1.17	1 (8%)	13,16,16	1.86	5 (38%)
4	DWS	C	501	1	36,38,38	1.90	9 (25%)	44,56,56	1.48	6 (13%)

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	ICT	B	502	2	-	2/16/16/16	-
3	ICT	A	502	2	-	2/16/16/16	-
4	DWS	B	503	1	-	2/14/50/50	0/5/5/5
4	DWS	A	504	1	-	4/14/50/50	0/5/5/5
3	ICT	A	505	2	-	2/16/16/16	-
4	DWS	C	501	1	-	5/14/50/50	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	501	DWS	C26-C8	-5.80	1.50	1.55
4	B	503	DWS	C26-C8	-4.46	1.51	1.55
4	C	501	DWS	C27-C29	-4.31	1.48	1.53
4	B	503	DWS	C27-C29	-4.06	1.48	1.53
4	C	501	DWS	C27-C34	-3.57	1.43	1.47
4	C	501	DWS	C11-C12	-2.75	1.46	1.51
4	B	503	DWS	C27-C34	-2.74	1.44	1.47
4	C	501	DWS	C10-C9	-2.66	1.49	1.53
4	C	501	DWS	O37-C2	-2.52	1.23	1.30
4	B	503	DWS	C8-C7	-2.41	1.49	1.53
4	A	504	DWS	C15-N18	2.34	1.41	1.36
4	B	503	DWS	C11-C12	-2.31	1.47	1.51
4	C	501	DWS	C26-C27	-2.24	1.49	1.54
4	C	501	DWS	C30-C9	-2.22	1.50	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	DWS	C10-C9	-2.20	1.50	1.53
4	A	504	DWS	C27-C29	-2.15	1.50	1.53
3	A	502	ICT	C3-C6	2.13	1.55	1.51
4	C	501	DWS	C30-C29	-2.12	1.48	1.51
3	A	505	ICT	C3-C6	2.10	1.54	1.51
3	A	502	ICT	O5-C6	2.02	1.28	1.22
4	B	503	DWS	C32-C30	-2.02	1.48	1.53

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	504	DWS	C34-C27-C29	6.60	117.05	109.04
4	B	503	DWS	C34-C27-C29	6.05	116.38	109.04
4	C	501	DWS	C34-C27-C29	4.47	114.45	109.04
4	A	504	DWS	O33-C29-C27	-4.41	118.73	123.07
4	A	504	DWS	N14-C15-N16	-4.41	122.16	126.42
3	A	502	ICT	C4-C3-C6	-3.88	105.08	111.25
4	C	501	DWS	N14-C15-N16	-3.87	122.68	126.42
4	B	503	DWS	N14-C15-N16	-3.65	122.89	126.42
3	A	505	ICT	O3-C5-C4	-3.48	112.17	122.84
3	B	502	ICT	O3-C5-C4	-3.24	112.90	122.84
4	C	501	DWS	C32-C30-C29	-3.18	105.08	111.63
3	A	505	ICT	C4-C3-C6	-3.12	106.28	111.25
4	A	504	DWS	C30-C29-C27	2.79	119.15	113.67
4	C	501	DWS	C26-C8-C17	-2.73	107.86	114.77
4	B	503	DWS	C26-C8-C17	-2.58	108.23	114.77
4	A	504	DWS	C10-C9-C30	2.58	118.25	114.15
3	A	502	ICT	C3-C4-C5	-2.55	109.19	113.95
4	B	503	DWS	C7-C8-C17	-2.54	106.54	110.88
4	A	504	DWS	C7-C8-C17	-2.54	106.54	110.88
4	A	504	DWS	C26-C8-C17	-2.52	108.39	114.77
4	B	503	DWS	O33-C29-C27	-2.51	120.60	123.07
4	A	504	DWS	C8-C26-C27	2.49	116.51	111.68
4	C	501	DWS	C23-C24-C19	2.37	122.47	119.73
3	A	502	ICT	O1-C1-C2	-2.32	115.43	121.62
4	A	504	DWS	N18-C15-N14	2.30	122.81	116.29
3	A	502	ICT	O4-C5-O3	2.27	129.16	123.33
3	A	505	ICT	O5-C6-C3	-2.16	117.91	123.01
3	A	505	ICT	O4-C5-O3	2.16	128.88	123.33
3	A	505	ICT	C3-C4-C5	-2.14	109.94	113.95
3	B	502	ICT	O1-C1-C2	-2.11	116.00	121.62
4	C	501	DWS	C32-C30-C9	-2.06	110.75	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	ICT	O5-C6-C3	-2.05	118.17	123.01

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	504	DWS	O1-C2-C3-C4
4	A	504	DWS	O37-C2-C3-C4
3	A	502	ICT	C4-C3-C6-O5
3	A	502	ICT	C4-C3-C6-O6
4	B	503	DWS	C6-C7-C8-C17
4	B	503	DWS	C36-C7-C8-C17
4	C	501	DWS	C36-C7-C8-C17
4	C	501	DWS	O37-C2-C3-C4
4	A	504	DWS	O37-C2-C3-C36
4	A	504	DWS	O1-C2-C3-C36
4	C	501	DWS	O37-C2-C3-C36
3	A	505	ICT	C3-C4-C5-O3
3	B	502	ICT	C3-C4-C5-O4
3	A	505	ICT	C3-C4-C5-O4
3	B	502	ICT	C3-C4-C5-O3
4	C	501	DWS	C6-C7-C8-C17
4	C	501	DWS	O1-C2-C3-C4

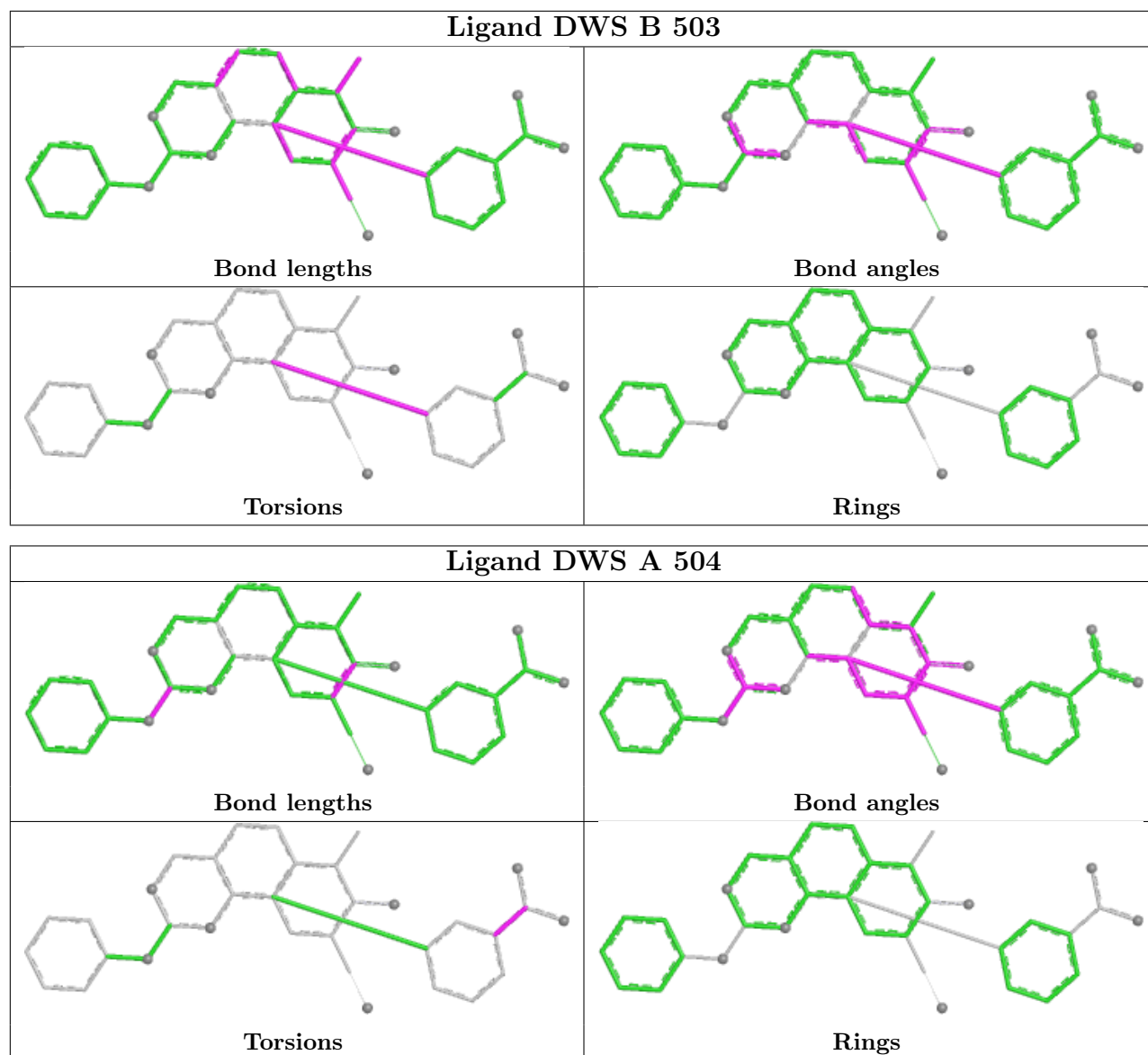
There are no ring outliers.

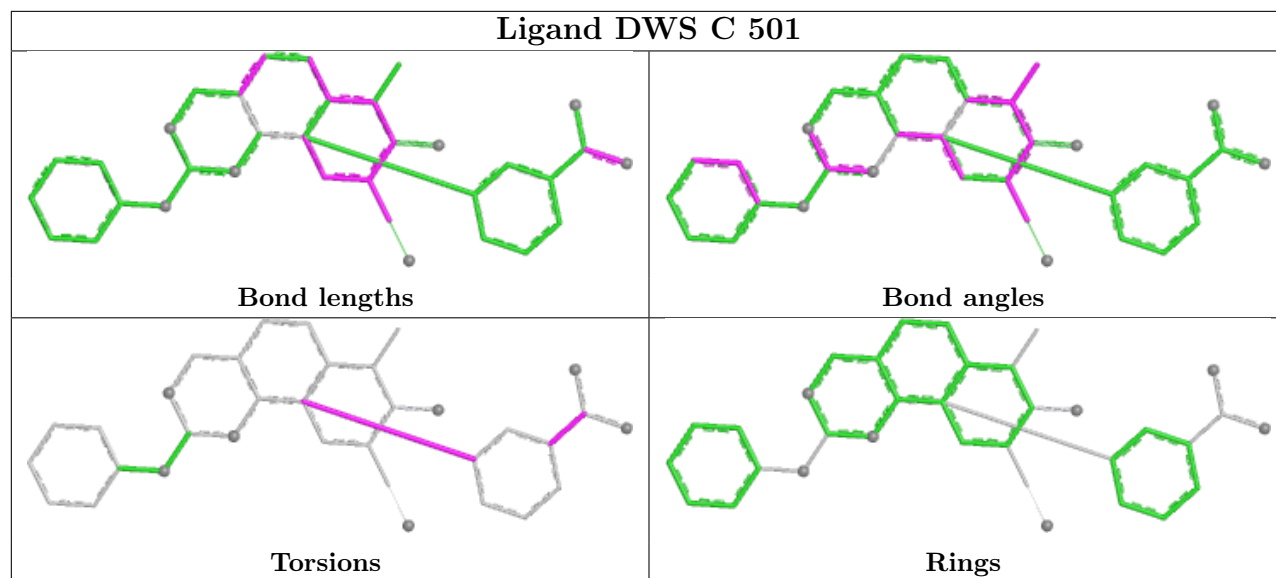
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	DWS	1	0
4	C	501	DWS	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

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	416/425 (97%)	0.25	14 (3%) 48 53	20, 33, 64, 152	0
1	B	416/425 (97%)	-0.01	5 (1%) 76 82	18, 28, 51, 132	0
1	C	416/425 (97%)	0.09	6 (1%) 73 80	21, 31, 52, 112	0
All	All	1248/1275 (97%)	0.11	25 (2%) 65 71	18, 31, 56, 152	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	416	LEU	5.5
1	C	416	LEU	5.4
1	A	161	GLY	4.3
1	C	418	HIS	4.3
1	A	414	LEU	3.9
1	B	418	HIS	3.9
1	A	416	LEU	3.5
1	B	414	LEU	3.2
1	C	417	GLU	3.2
1	A	163	GLN	3.1
1	A	162	THR	2.9
1	A	159	SER	2.9
1	A	413	LYS	2.6
1	A	418	HIS	2.6
1	B	3	LYS	2.6
1	A	47	GLU	2.4
1	C	3	LYS	2.3
1	A	3	LYS	2.2
1	A	115	LYS	2.2
1	B	122	SER	2.1
1	A	322	GLY	2.1
1	A	160	ASP	2.1
1	C	122	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	20	ARG	2.1
1	A	122	SER	2.0

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](#)

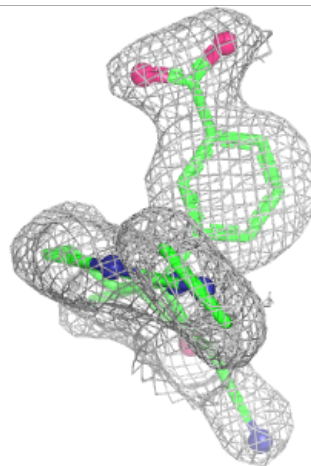
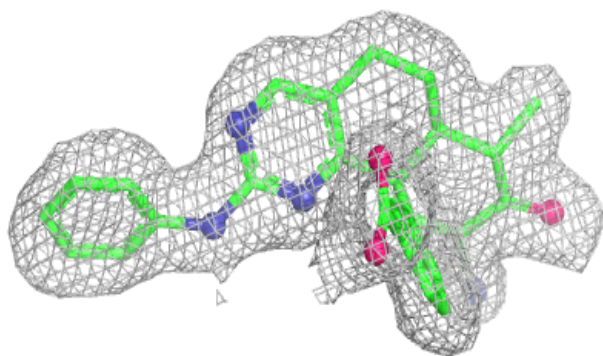
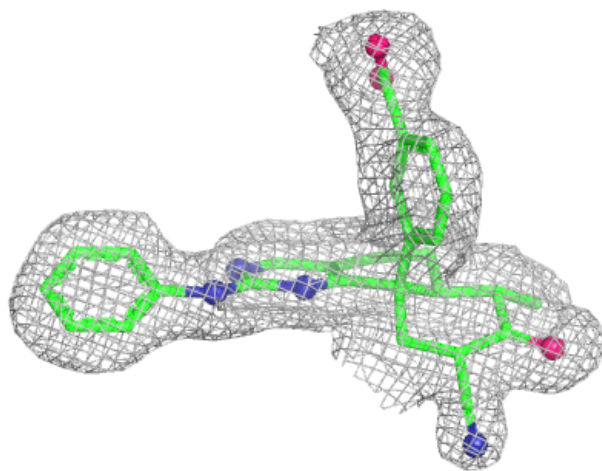
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	DWS	C	501	34/34	0.91	0.08	31,34,43,46	0
4	DWS	A	504	34/34	0.93	0.07	36,39,47,48	0
3	ICT	A	502	13/13	0.94	0.07	27,29,33,35	0
4	DWS	B	503	34/34	0.94	0.07	23,28,36,38	0
3	ICT	A	505	13/13	0.94	0.07	27,29,35,37	0
3	ICT	B	502	13/13	0.96	0.06	21,26,30,30	0
2	CA	A	501	1/1	0.99	0.07	39,39,39,39	0
2	CA	A	503	1/1	0.99	0.07	40,40,40,40	0
2	CA	B	501	1/1	1.00	0.07	35,35,35,35	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.

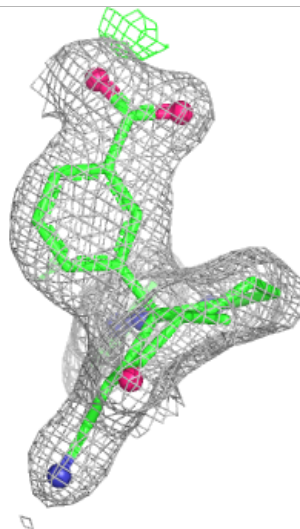
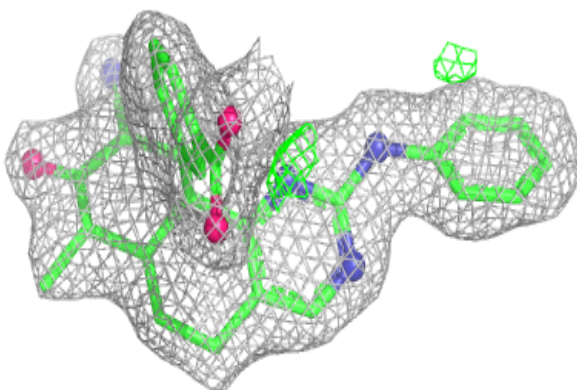
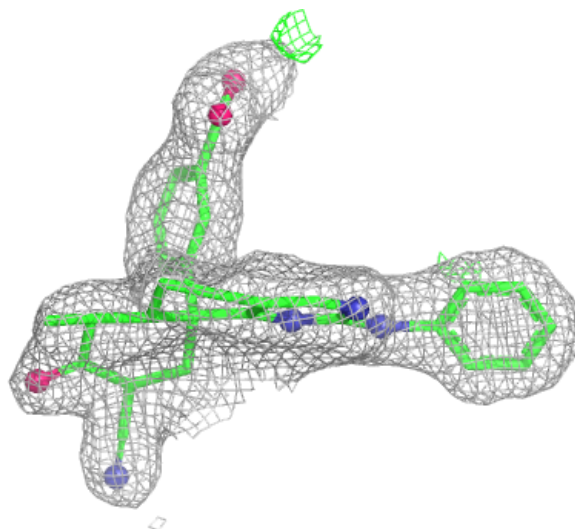
Electron density around DWS C 501:

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



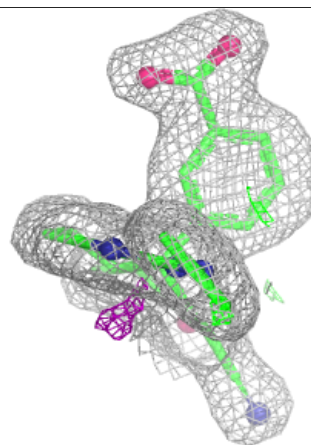
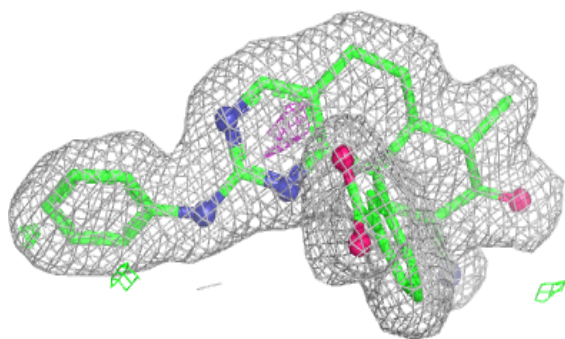
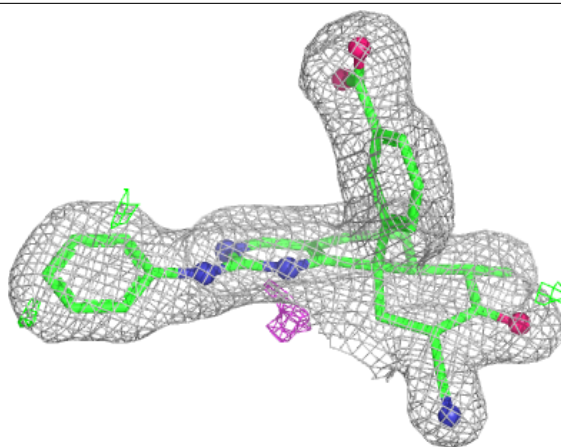
Electron density around DWS A 504:

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



Electron density around DWS B 503:

$2mF_o-DF_c$ (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.