



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 12:00 pm BST

PDB ID : 6BL1
Title : Novel Modes of Inhibition of Wild-Type IDH1: Direct Covalent Modification of His315 with Cmpd13
Authors : Jakob, C.G.; Qiu, W.
Deposited on : 2017-11-09
Resolution : 2.02 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

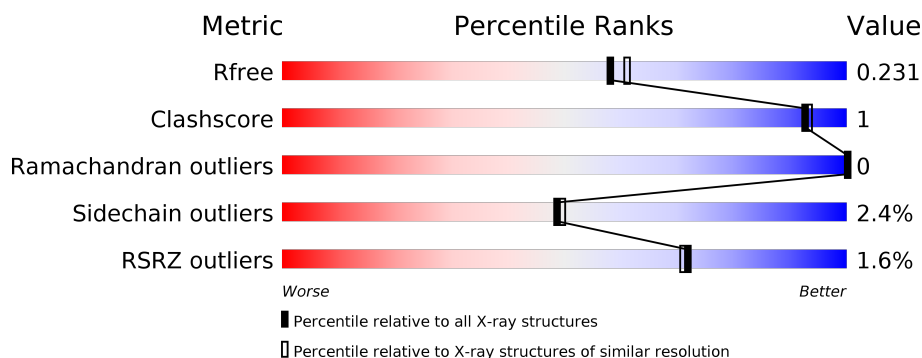
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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}	130704	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 on 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	
1	B	425	
1	C	425	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10744 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	B	416	Total	C	N	O	S	0	2	0
			3303	2101	561	623	18			
1	C	416	Total	C	N	O	S	0	1	0
			3293	2093	559	623	18			
1	A	416	Total	C	N	O	S	0	2	0
			3303	2100	561	624	18			

There are 33 discrepancies between the modelled and reference sequences:

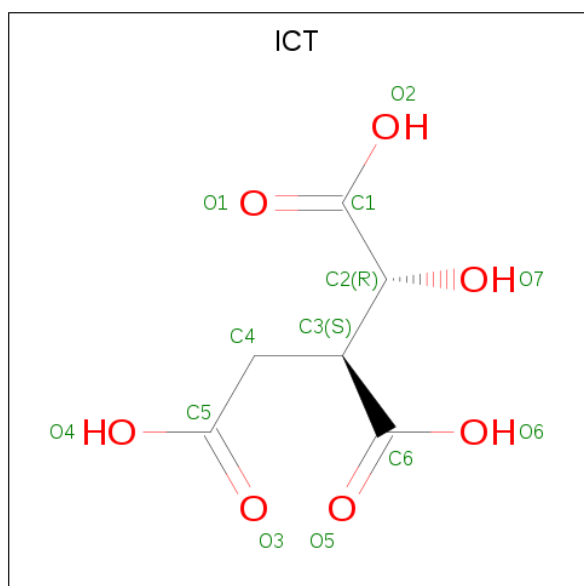
Chain	Residue	Modelled	Actual	Comment	Reference
B	415	SER	-	expression tag	UNP O75874
B	416	LEU	-	expression tag	UNP O75874
B	417	GLU	-	expression tag	UNP O75874
B	418	HIS	-	expression tag	UNP O75874
B	419	HIS	-	expression tag	UNP O75874
B	420	HIS	-	expression tag	UNP O75874
B	421	HIS	-	expression tag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

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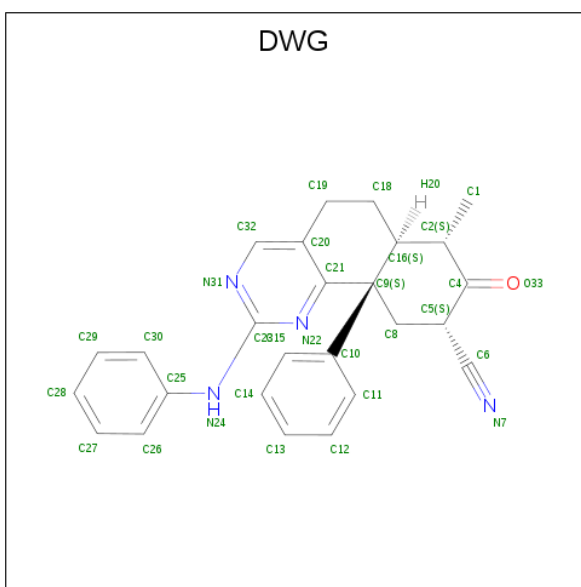
Chain	Residue	Modelled	Actual	Comment	Reference
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	UNP O75874
A	422	HIS	-	expression tag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

- Molecule 2 is ISOCITRIC ACID (three-letter code: ICT) (formula: C₆H₈O₇).



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

- Molecule 3 is (6aS,7S,9S,10aS)-7-methyl-8-oxo-10a-phenyl-2-(phenylamino)-5,6,6a,7,8,9,10,10a-octahydrobenzo[h]quinazoline-9-carbonitrile (three-letter code: DWG) (formula: C₂₆H₂₄N₄O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			31	26	4	1		
3	C	1	Total	C	N	O	0	0
			31	26	4	1		
3	A	1	Total	C	N	O	0	0
			31	26	4	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

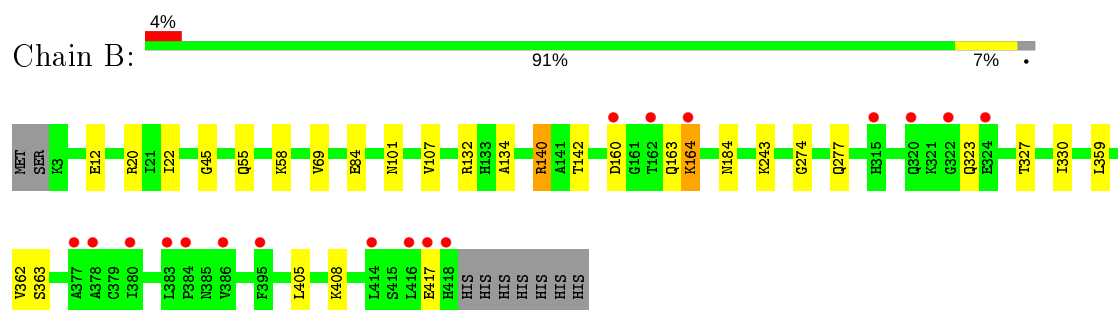
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	193	Total	O	0	0
			193	193		
5	C	313	Total	O	0	0
			313	313		
5	A	204	Total	O	0	0
			204	204		

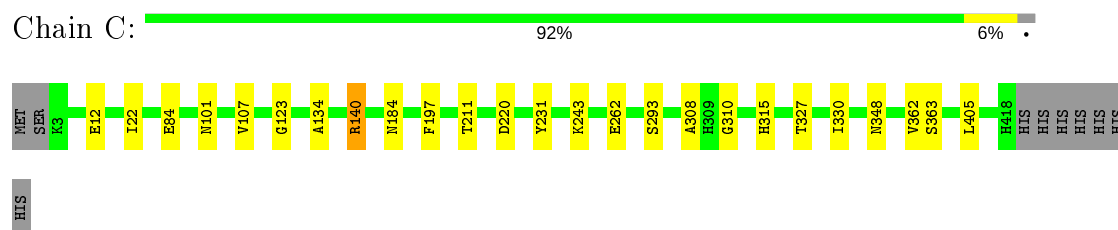
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

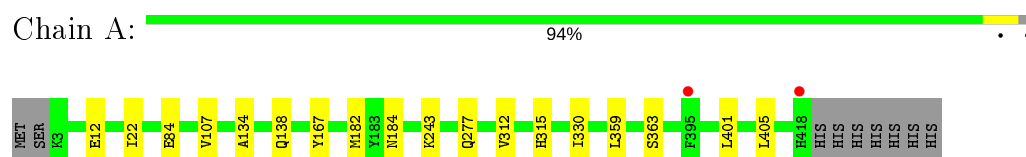
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



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- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	98.61Å 273.81Å 114.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.78 – 2.02 92.78 – 2.02	Depositor EDS
% Data completeness (in resolution range)	99.9 (92.78-2.02) 99.9 (92.78-2.02)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.02Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.183 , 0.223 0.190 , 0.231	Depositor DCC
R_{free} test set	5135 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10744	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, DWG, ICT

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	0/3378	0.65	0/4556
1	B	0.49	0/3378	0.66	0/4555
1	C	0.53	0/3365	0.65	0/4540
All	All	0.50	0/10121	0.66	0/13651

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	3303	0	3269	8	0
1	B	3303	0	3274	12	0
1	C	3293	0	3250	11	0
2	B	26	0	8	0	0
2	C	13	0	4	0	0
3	A	31	0	0	1	0
3	B	31	0	0	0	0
3	C	31	0	0	1	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	204	0	0	0	0
5	B	193	0	0	0	0
5	C	313	0	0	1	0
All	All	10744	0	9805	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:A:315:HIS:NE2	3:A:501:DWG:C8	2.36	0.88
1:C:315:HIS:NE2	3:C:503:DWG:C8	2.46	0.78
1:C:101[B]:ASN:ND2	1:C:140:ARG:HD2	2.03	0.73
1:C:107:VAL:HG23	1:C:134:ALA:HB2	1.78	0.66
1:B:142:THR:HG21	1:A:167:TYR:HB3	1.84	0.60
1:C:293:SER:HB2	1:C:308:ALA:HB2	1.90	0.54
1:B:101[B]:ASN:ND2	1:B:140:ARG:HD2	2.24	0.53
1:A:363:SER:HA	1:A:401:LEU:HD21	1.95	0.48
1:B:277:GLN:HE22	1:A:277:GLN:HE22	1.63	0.47
1:C:362:VAL:HG21	1:C:405:LEU:HA	1.96	0.47
1:B:55:GLN:HA	1:B:58:LYS:HE3	1.97	0.46
1:C:310:GLY:HA3	5:C:761:HOH:O	2.15	0.46
1:C:330:ILE:HD12	1:C:363:SER:HB3	1.97	0.46
1:B:362:VAL:HG21	1:B:405:LEU:HA	1.99	0.45
1:B:362:VAL:HG23	1:B:408:LYS:HD2	1.96	0.45
1:B:132:ARG:HG3	1:B:274:GLY:HA3	1.98	0.45
1:A:138:GLN:HG3	1:A:182:MET:HE1	1.98	0.45
1:C:22:ILE:HD11	1:C:327:THR:HB	1.99	0.45
1:A:107:VAL:HG23	1:A:134:ALA:HB2	1.99	0.45
1:A:330:ILE:HD12	1:A:363:SER:HB3	2.01	0.42
1:B:164[B]:LYS:H	1:B:164[B]:LYS:HG2	1.64	0.42
1:B:330:ILE:HD12	1:B:363:SER:HB3	2.02	0.42
1:C:211:THR:HB	1:C:220:ASP:HB3	2.01	0.41
1:C:197:PHE:CZ	1:C:231:TYR:HB2	2.56	0.41
1:B:359:LEU:HD13	1:B:405:LEU:HD22	2.03	0.41
1:B:20:ARG:HH21	1:B:45:GLY:HA3	1.86	0.41
1:B:107:VAL:HG23	1:B:134:ALA:HB2	2.03	0.41
1:A:359:LEU:HD13	1:A:405:LEU:HD22	2.03	0.41
1:C:123:GLY:O	1:C:262:GLU:HA	2.22	0.40

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	416/425 (98%)	402 (97%)	14 (3%)	0	100	100
1	B	416/425 (98%)	397 (95%)	19 (5%)	0	100	100
1	C	415/425 (98%)	403 (97%)	12 (3%)	0	100	100
All	All	1247/1275 (98%)	1202 (96%)	45 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	352/361 (98%)	346 (98%)	6 (2%)	60	63
1	B	352/361 (98%)	338 (96%)	14 (4%)	31	28
1	C	350/361 (97%)	344 (98%)	6 (2%)	60	63
All	All	1054/1083 (97%)	1028 (98%)	26 (2%)	49	48

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	12	GLU
1	B	22	ILE

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Mol	Chain	Res	Type
1	B	69	VAL
1	B	84	GLU
1	B	140	ARG
1	B	160	ASP
1	B	163	GLN
1	B	164[A]	LYS
1	B	164[B]	LYS
1	B	184	ASN
1	B	243	LYS
1	B	323	GLN
1	B	327	THR
1	B	417	GLU
1	C	12	GLU
1	C	84	GLU
1	C	140	ARG
1	C	184	ASN
1	C	243	LYS
1	C	348	ASN
1	A	12	GLU
1	A	22	ILE
1	A	84	GLU
1	A	184	ASN
1	A	243	LYS
1	A	312	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	277	GLN
1	C	14	GLN
1	C	385	ASN

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 [i](#)

There are no carbohydrates 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	DWG	C	503	-	32,35,35	0.80	2 (6%)	40,51,51	1.76	8 (20%)
2	ICT	C	502	4	2,12,12	0.78	0	5,16,16	3.32	3 (60%)
3	DWG	B	502	1	32,35,35	0.65	1 (3%)	40,51,51	1.62	6 (15%)
2	ICT	B	505	4	2,12,12	0.26	0	5,16,16	2.50	2 (40%)
2	ICT	B	501	4	2,12,12	0.28	0	5,16,16	2.24	3 (60%)
3	DWG	A	501	-	32,35,35	0.72	1 (3%)	40,51,51	1.68	9 (22%)

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	DWG	C	503	-	-	1/10/46/46	0/5/5/5
2	ICT	C	502	4	-	1/6/16/16	-
3	DWG	B	502	1	-	0/10/46/46	0/5/5/5
2	ICT	B	505	4	-	0/6/16/16	-
2	ICT	B	501	4	-	1/6/16/16	-
3	DWG	A	501	-	-	0/10/46/46	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	503	DWG	C23-N24	2.80	1.42	1.36
3	B	502	DWG	C23-N24	2.64	1.41	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	503	DWG	C8-C9	-2.39	1.53	1.55
3	A	501	DWG	C23-N24	2.32	1.41	1.36

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	502	ICT	C4-C3-C6	-5.01	105.91	112.70
2	C	502	ICT	C1-C2-C3	-4.83	106.81	112.25
3	C	503	DWG	C2-C4-C5	4.64	122.20	113.47
3	B	502	DWG	C2-C4-C5	4.37	121.68	113.47
3	A	501	DWG	C9-C8-C5	4.35	120.30	111.68
2	B	505	ICT	C4-C3-C6	-4.16	107.07	112.70
3	C	503	DWG	C9-C8-C5	3.96	119.52	111.68
3	A	501	DWG	C2-C4-C5	3.85	120.71	113.47
3	C	503	DWG	C8-C5-C4	3.65	115.26	108.09
3	B	502	DWG	N31-C23-N22	-3.62	123.12	126.55
3	C	503	DWG	N31-C23-N22	-3.58	123.15	126.55
3	A	501	DWG	C19-C20-C21	3.56	124.52	121.64
3	A	501	DWG	C8-C5-C4	3.54	115.04	108.09
3	C	503	DWG	C10-C9-C16	3.42	118.84	111.75
3	B	502	DWG	C19-C20-C21	3.40	124.39	121.64
2	B	505	ICT	C1-C2-C3	-3.26	108.58	112.25
3	A	501	DWG	C10-C9-C16	3.21	118.41	111.75
2	B	501	ICT	C1-C2-C3	-3.20	108.64	112.25
2	B	501	ICT	C4-C3-C6	-3.10	108.50	112.70
3	C	503	DWG	C19-C20-C21	3.10	124.15	121.64
3	A	501	DWG	N31-C23-N22	-2.81	123.88	126.55
3	B	502	DWG	C8-C5-C4	2.65	113.29	108.09
3	B	502	DWG	C8-C9-C21	-2.52	108.39	114.77
3	B	502	DWG	C1-C2-C16	-2.49	109.99	113.98
3	C	503	DWG	C8-C9-C21	-2.44	108.61	114.77
3	C	503	DWG	C10-C9-C21	-2.38	106.38	110.95
3	A	501	DWG	C8-C9-C10	-2.28	106.54	111.03
2	C	502	ICT	O7-C2-C3	2.26	114.88	108.56
3	A	501	DWG	C19-C20-C32	-2.17	118.38	121.94
3	A	501	DWG	C12-C11-C10	2.12	122.99	120.76
2	B	501	ICT	O7-C2-C3	2.07	114.35	108.56

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	502	ICT	C1-C2-C3-C6
2	B	501	ICT	C1-C2-C3-C6
3	C	503	DWG	C15-C10-C9-C21

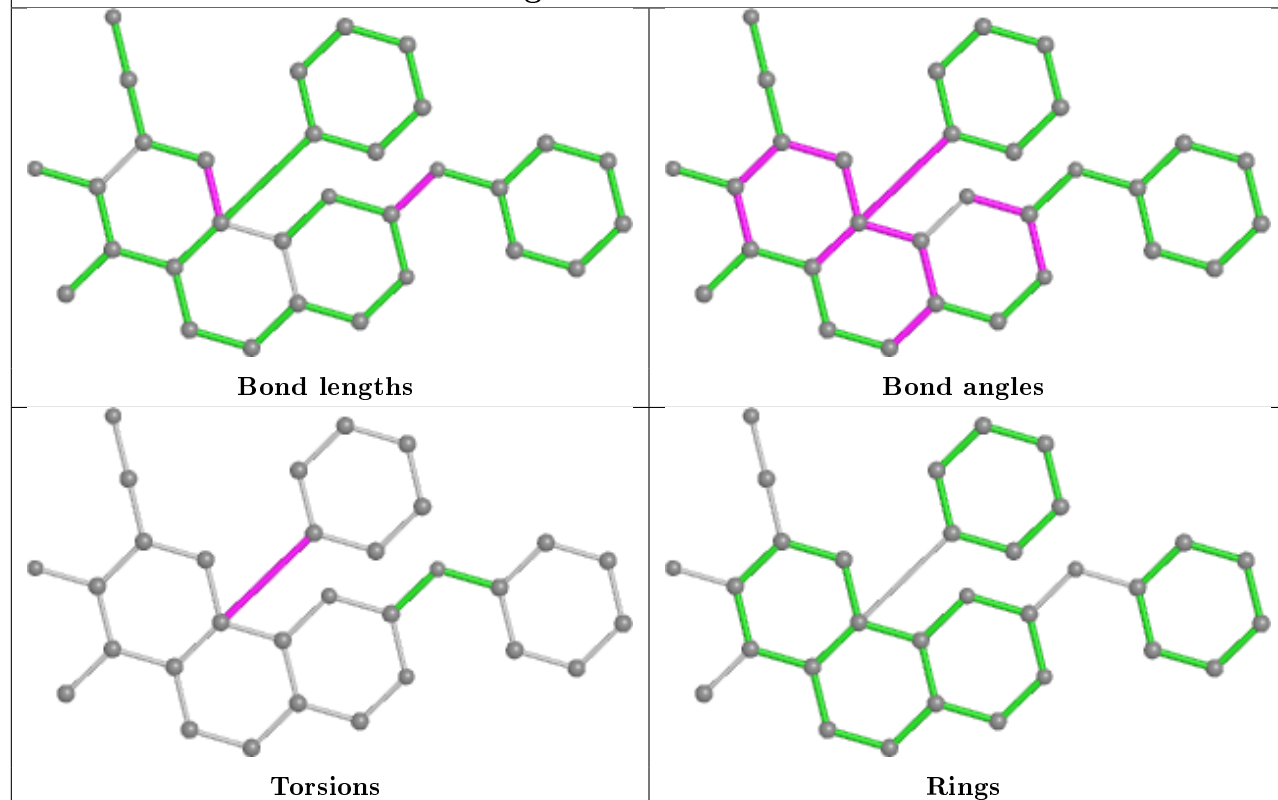
There are no ring outliers.

2 monomers are involved in 2 short contacts:

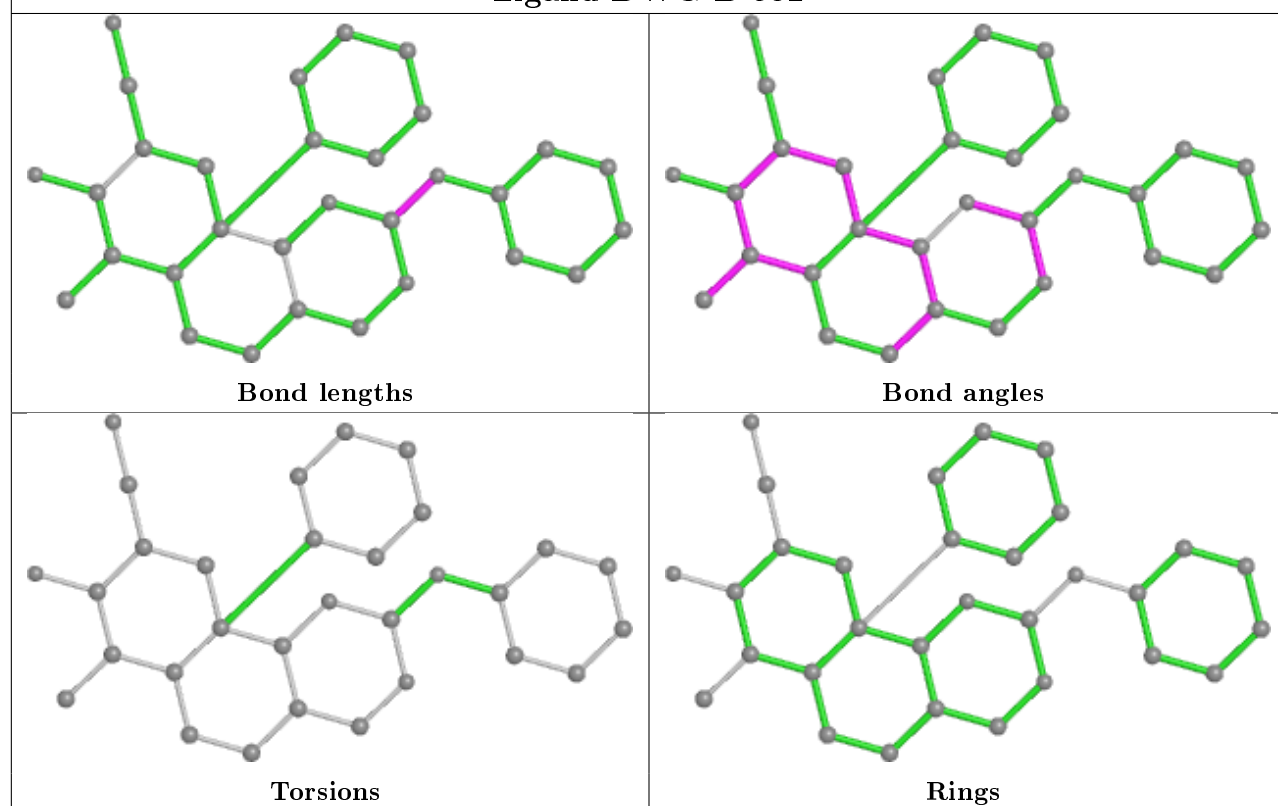
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	DWG	1	0
3	A	501	DWG	1	0

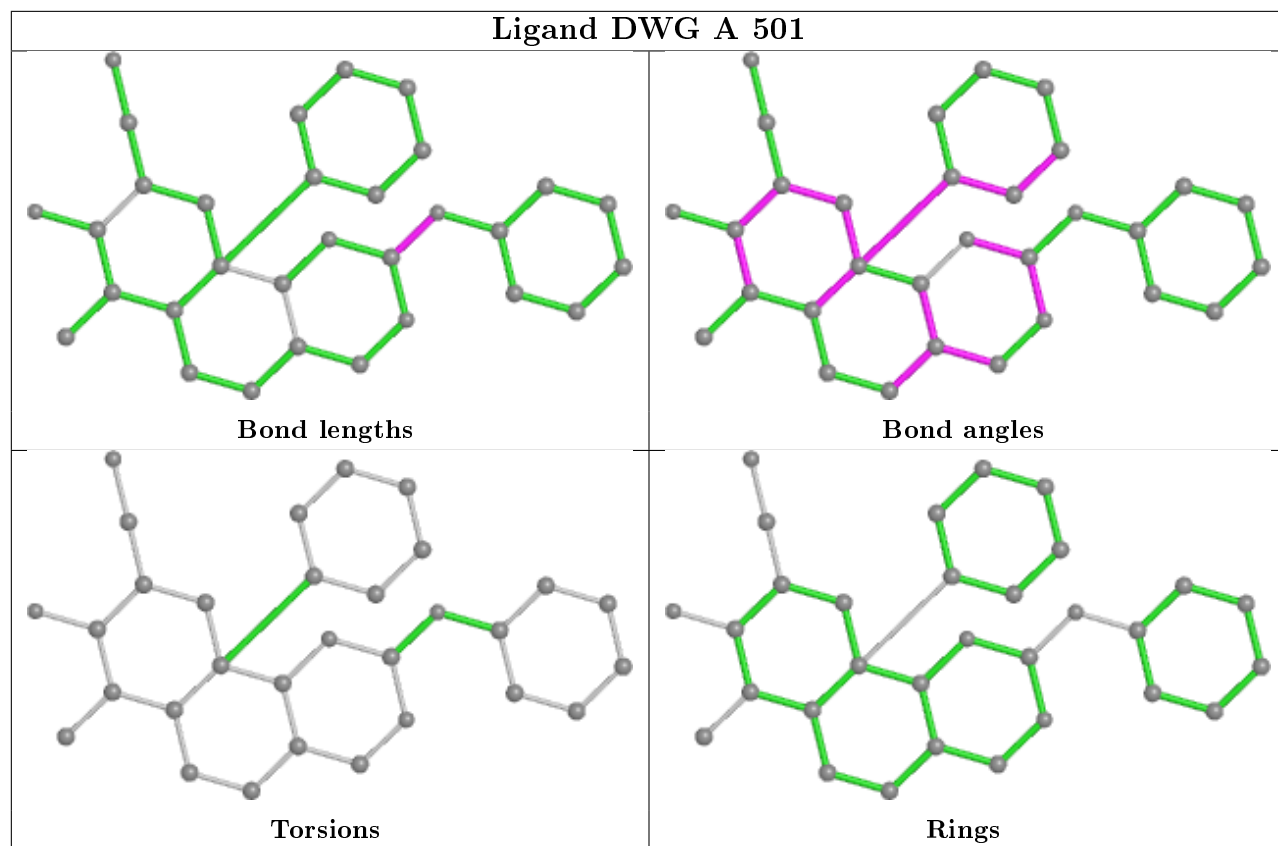
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand DWG C 503



Ligand DWG B 502





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%)	-0.26	2 (0%) 91 91	21, 40, 64, 110	0
1	B	416/425 (97%)	-0.12	18 (4%) 35 34	21, 39, 72, 147	0
1	C	416/425 (97%)	-0.27	0 100 100	19, 29, 53, 91	0
All	All	1248/1275 (97%)	-0.21	20 (1%) 72 71	19, 36, 64, 147	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162	THR	10.0
1	B	418	HIS	5.8
1	B	414	LEU	5.3
1	B	416	LEU	4.3
1	B	384	PRO	3.7
1	A	418	HIS	3.7
1	B	164[A]	LYS	3.1
1	B	383	LEU	3.1
1	B	320	GLN	3.1
1	A	395	PHE	3.0
1	B	160	ASP	2.9
1	B	395	PHE	2.8
1	B	377	ALA	2.7
1	B	386	VAL	2.4
1	B	380	ILE	2.4
1	B	417	GLU	2.4
1	B	324	GLU	2.3
1	B	315	HIS	2.2
1	B	378	ALA	2.2
1	B	322	GLY	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 carbohydrates in this entry.

6.4 Ligands [i](#)

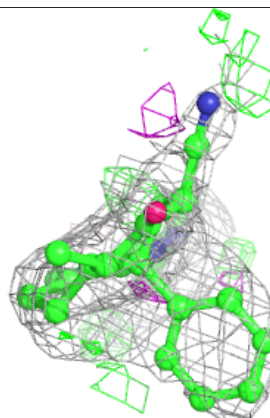
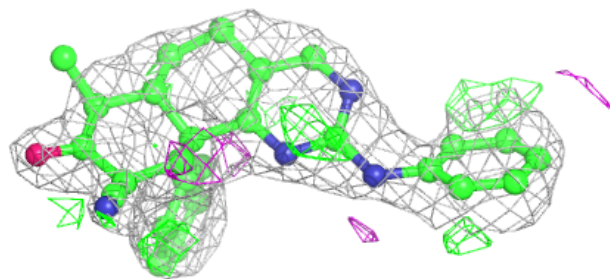
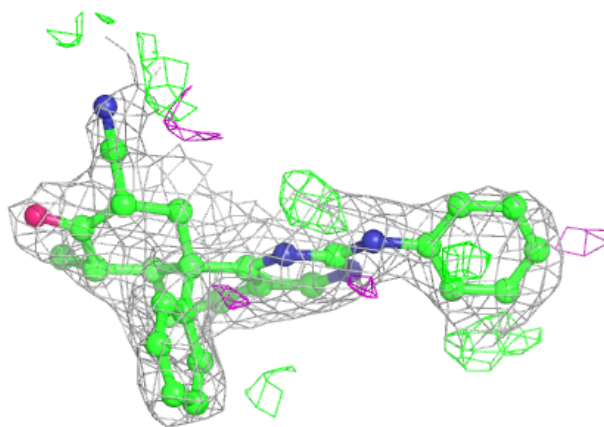
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	DWG	B	502	31/31	0.80	0.24	62,65,68,69	0
3	DWG	C	503	31/31	0.83	0.15	32,45,49,50	0
3	DWG	A	501	31/31	0.87	0.17	46,50,54,56	0
2	ICT	B	505	13/13	0.90	0.17	44,48,55,56	0
2	ICT	C	502	13/13	0.91	0.14	34,40,50,52	0
2	ICT	B	501	13/13	0.92	0.17	41,45,46,47	0
4	CA	C	501	1/1	0.96	0.05	37,37,37,37	0
4	CA	B	504	1/1	0.97	0.04	51,51,51,51	0
4	CA	B	503	1/1	0.99	0.08	45,45,45,45	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.

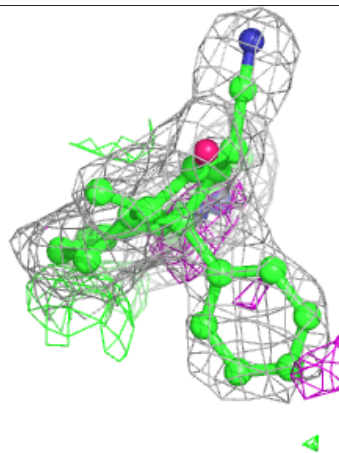
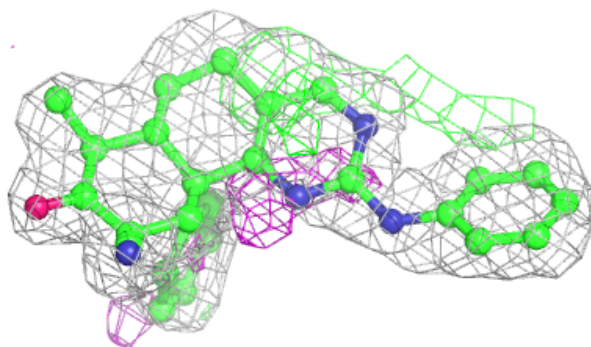
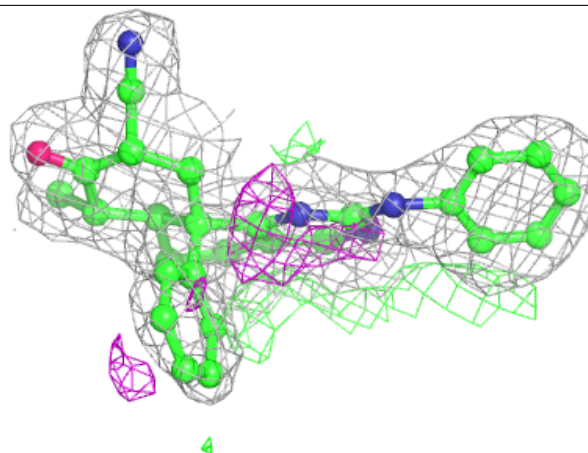
Electron density around DWG B 502:

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



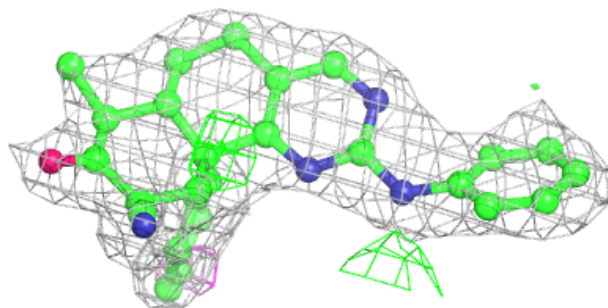
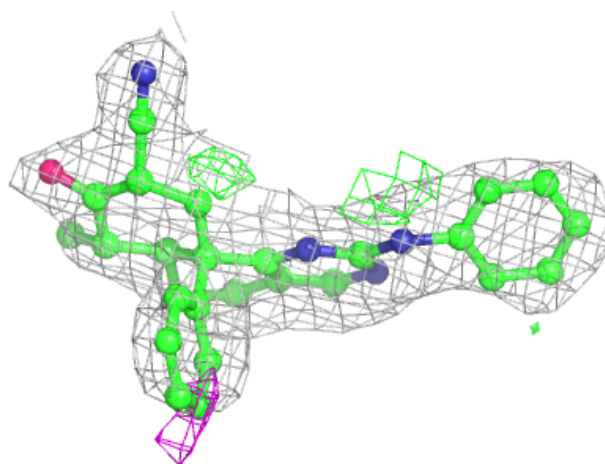
Electron density around DWG C 503:

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



Electron density around DWG A 501:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.