



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 10:02 AM UTC

PDB ID : 1RL4 / pdb_00001rl4
Title : Plasmodium falciparum peptide deformylase complex with inhibitor
Authors : Robien, M.A.; Nguyen, K.T.; Kumar, A.; Hirsh, I.; Turley, S.; Pei, D.; Hol, W.G.J.
Deposited on : 2003-11-24
Resolution : 2.18 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

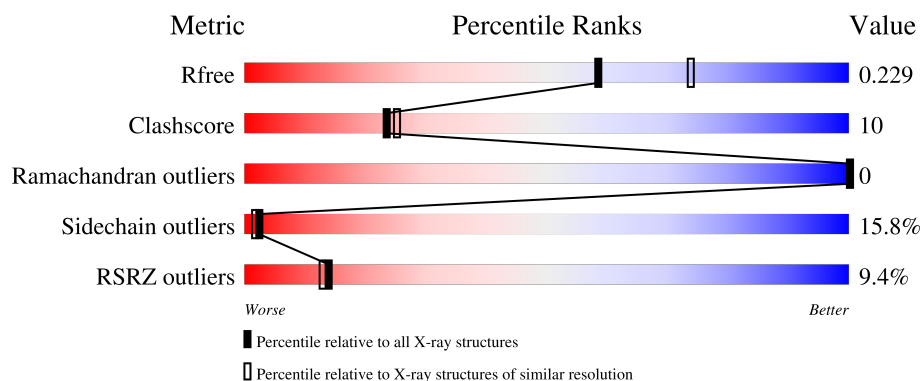
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	188	6% 59% 19% 5% 18%
1	B	188	9% 61% 18% • 17%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2822 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 formylmethionine deformylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1269	819	225	221	4			
1	B	156	Total	C	N	O	S	0	0	0
			1286	831	227	224	4			

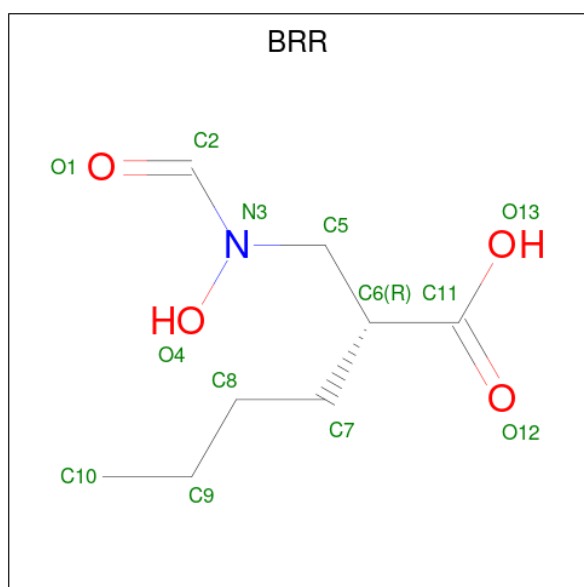
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	57	MET	-	cloning artifact	UNP Q8I372
A	58	SER	-	cloning artifact	UNP Q8I372
A	237	LEU	-	cloning artifact	UNP Q8I372
A	238	GLU	-	cloning artifact	UNP Q8I372
A	239	HIS	-	cloning artifact	UNP Q8I372
A	240	HIS	-	cloning artifact	UNP Q8I372
A	241	HIS	-	cloning artifact	UNP Q8I372
A	242	HIS	-	cloning artifact	UNP Q8I372
A	243	HIS	-	cloning artifact	UNP Q8I372
A	244	HIS	-	cloning artifact	UNP Q8I372
B	57	MET	-	cloning artifact	UNP Q8I372
B	58	SER	-	cloning artifact	UNP Q8I372
B	237	LEU	-	cloning artifact	UNP Q8I372
B	238	GLU	-	cloning artifact	UNP Q8I372
B	239	HIS	-	cloning artifact	UNP Q8I372
B	240	HIS	-	cloning artifact	UNP Q8I372
B	241	HIS	-	cloning artifact	UNP Q8I372
B	242	HIS	-	cloning artifact	UNP Q8I372
B	243	HIS	-	cloning artifact	UNP Q8I372
B	244	HIS	-	cloning artifact	UNP Q8I372

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

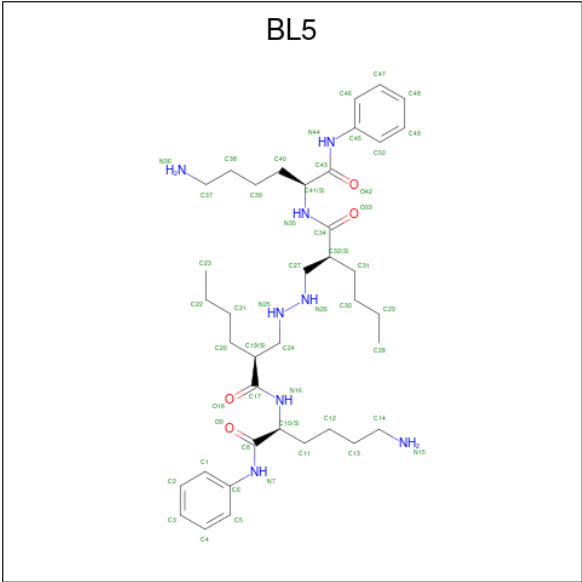
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Co 1 1	0	0
2	B	1	Total Co 1 1	0	0

- Molecule 3 is (2R)-2-{[FORMYL(HYDROXY)AMINO]METHYL}HEXANOIC ACID (CCD ID: BRR) (formula: C₈H₁₅NO₄).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 13 8 1 4	0	0
3	B	1	Total C N O 13 8 1 4	0	0

- Molecule 4 is 2-{N'-[2-(5-AMINO-1-PHENYLCARBAMOYL-PENTYLCARBAMOYL)-HEXYL]-HYDRAZINOMETHYL}-HEXANOIC ACID(5-AMINO-1-PHENYLCARBAMOYL-PENTYL)-AMIDE (CCD ID: BL5) (formula: C₃₈H₆₂N₈O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			50	38	8	4		
4	B	1	Total	C	N	O	0	0
			50	38	8	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	72	Total	O	0	0
			72	72		
5	B	67	Total	O	0	0
			67	67		

4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	102.32Å 102.32Å 118.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.18 20.00 – 2.18	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.18) 99.8 (20.00-2.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.197 , 0.212 0.208 , 0.229	Depositor DCC
R_{free} test set	3566 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.045 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2822	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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: BRR, CO, BL5

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	0.61	0/1292	0.84	0/1733
1	B	0.61	0/1310	0.87	1/1758 (0.1%)
All	All	0.61	0/2602	0.85	1/3491 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	168	ARG	NE-CZ-NH2	-5.77	114.01	119.20

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	1269	0	1334	31	0
1	B	1286	0	1345	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	13	0	13	4	0
3	B	13	0	13	1	0
4	A	50	0	62	2	0
4	B	50	0	62	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	72	0	0	2	1
5	B	67	0	0	1	1
All	All	2822	0	2829	56	1

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 10.

All (56) 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:166:VAL:O	1:A:194:ARG:NH2	2.02	0.92
1:B:166:VAL:O	1:B:194:ARG:NH2	2.12	0.81
1:A:162:ILE:HG23	1:A:220:VAL:HG21	1.71	0.71
1:A:135:ARG:HG2	1:A:135:ARG:HH11	1.56	0.70
1:A:162:ILE:HD11	1:A:227:LEU:HD22	1.72	0.70
1:B:168:ARG:NH2	1:B:201:ASP:OD2	2.29	0.65
1:A:162:ILE:HD13	1:A:224:LEU:HD23	1.78	0.65
1:A:162:ILE:CD1	1:A:227:LEU:HD22	2.30	0.62
1:B:69:LYS:HE2	1:B:231:TYR:OH	2.00	0.60
1:A:162:ILE:HD13	1:A:224:LEU:CD2	2.32	0.59
1:A:199:GLU:OE2	3:A:401:BRR:H2	2.04	0.58
1:A:195:ILE:O	1:A:199:GLU:HG3	2.03	0.58
4:B:1501:BL5:O9	4:B:1501:BL5:H1	2.04	0.58
1:A:219:LYS:O	1:A:219:LYS:HD2	2.05	0.56
1:A:122:TRP:CD1	1:A:122:TRP:C	2.84	0.56
1:A:135:ARG:HG2	1:A:135:ARG:NH1	2.14	0.55
1:B:69:LYS:HG3	1:B:231:TYR:CE2	2.42	0.55
1:A:162:ILE:HD12	1:A:227:LEU:CD2	2.37	0.54
1:A:162:ILE:CD1	1:A:227:LEU:CD2	2.85	0.54
1:B:122:TRP:C	1:B:122:TRP:CD1	2.85	0.53
1:B:218:LYS:HG3	1:B:219:LYS:N	2.24	0.53
1:A:162:ILE:CD1	1:A:224:LEU:HD23	2.38	0.53
1:A:122:TRP:HZ3	1:A:196:PHE:HB2	1.74	0.52
3:A:401:BRR:H103	4:A:501:BL5:H4	1.91	0.51
1:B:107:GLY:HA3	1:B:199:GLU:OE2	2.11	0.50
1:A:157:LEU:HD22	3:A:401:BRR:H52	1.93	0.50
1:B:192:HIS:CD2	5:B:1741:HOH:O	2.66	0.48
1:B:89:LEU:HD13	1:B:179:ILE:HG13	1.96	0.47
1:A:122:TRP:C	1:A:122:TRP:HD1	2.24	0.46
1:A:221:ARG:HB3	1:A:222:PRO:HD3	1.98	0.46
1:B:151:LYS:HA	1:B:166:VAL:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:LYS:HA	1:A:66:LYS:HD2	1.49	0.46
1:B:220:VAL:HG13	1:B:224:LEU:HD22	1.98	0.46
1:A:162:ILE:HG23	1:A:220:VAL:CG2	2.41	0.45
1:A:226:GLU:O	1:A:230:ASP:OD2	2.34	0.45
1:A:211:LYS:HE2	5:A:809:HOH:O	2.17	0.45
1:B:101:TYR:HE1	1:B:123:ASN:HD22	1.64	0.44
1:A:72:ASP:OD1	1:A:73:PRO:HD2	2.17	0.44
1:B:122:TRP:C	1:B:122:TRP:HD1	2.26	0.44
1:A:120:ILE:HD11	1:A:200:PHE:HA	1.99	0.44
1:A:135:ARG:HH11	1:A:135:ARG:CG	2.29	0.43
1:A:189:LYS:HB3	1:A:189:LYS:HE3	1.43	0.43
3:A:401:BRR:C2	3:A:401:BRR:H71	2.49	0.42
4:A:501:BL5:H141	5:A:818:HOH:O	2.19	0.42
3:B:1401:BRR:C2	3:B:1401:BRR:H71	2.48	0.42
4:B:1501:BL5:O42	4:B:1501:BL5:H50	2.19	0.42
1:A:140:PRO:HA	1:A:175:SER:O	2.20	0.42
1:B:189:LYS:HE3	1:B:189:LYS:HB3	1.55	0.42
1:B:100:MET:CE	1:B:123:ASN:HB2	2.50	0.41
1:B:72:ASP:HA	1:B:73:PRO:HD3	1.83	0.41
1:A:110:ALA:N	1:A:111:PRO:CD	2.84	0.41
1:A:142:ILE:HD12	1:A:200:PHE:CD2	2.56	0.41
1:B:110:ALA:HB3	1:B:111:PRO:HD3	2.02	0.41
1:B:110:ALA:N	1:B:111:PRO:CD	2.83	0.40
1:A:157:LEU:HD12	1:A:157:LEU:HA	1.93	0.40
4:B:1501:BL5:O9	4:B:1501:BL5:C1	2.69	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:751:HOH:O	5:B:820:HOH:O[6_554]	1.79	0.41

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	150/188 (80%)	148 (99%)	2 (1%)	0	100	100
1	B	152/188 (81%)	151 (99%)	1 (1%)	0	100	100
All	All	302/376 (80%)	299 (99%)	3 (1%)	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	145/177 (82%)	121 (83%)	24 (17%)	2	1
1	B	146/177 (82%)	124 (85%)	22 (15%)	3	2
All	All	291/354 (82%)	245 (84%)	46 (16%)	2	2

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LYS
1	A	67	ILE
1	A	68	VAL
1	A	75	LEU
1	A	88	ASN
1	A	89	LEU
1	A	90	LYS
1	A	91	ARG
1	A	104	LYS
1	A	114	ASN
1	A	135	ARG
1	A	150	LEU
1	A	157	LEU
1	A	179	ILE
1	A	185	LEU

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Mol	Chain	Res	Type
1	A	186	LYS
1	A	189	LYS
1	A	199	GLU
1	A	207	LEU
1	A	217	LYS
1	A	219	LYS
1	A	220	VAL
1	A	227	LEU
1	A	230	ASP
1	B	66	LYS
1	B	67	ILE
1	B	68	VAL
1	B	75	LEU
1	B	88	ASN
1	B	89	LEU
1	B	90	LYS
1	B	94	ARG
1	B	95	LYS
1	B	106	ILE
1	B	114	ASN
1	B	118	ARG
1	B	135	ARG
1	B	150	LEU
1	B	179	ILE
1	B	185	LEU
1	B	186	LYS
1	B	189	LYS
1	B	207	LEU
1	B	218	LYS
1	B	219	LYS
1	B	224	LEU

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

Mol	Chain	Res	Type
1	B	145	GLN
1	B	180	ASN
1	B	192	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	BRR	B	1401	2	11,12,12	0.98	1 (9%)	11,14,14	2.29	5 (45%)
3	BRR	A	401	2	11,12,12	0.96	1 (9%)	11,14,14	2.29	3 (27%)
4	BL5	A	501	-	49,51,51	1.03	4 (8%)	56,62,62	0.93	2 (3%)
4	BL5	B	1501	-	49,51,51	1.02	4 (8%)	56,62,62	1.01	3 (5%)

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	BRR	B	1401	2	-	1/10/14/14	-
3	BRR	A	401	2	-	1/10/14/14	-
4	BL5	A	501	-	-	9/56/57/57	0/2/2/2
4	BL5	B	1501	-	-	4/56/57/57	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1501	BL5	C45-N44	-3.82	1.33	1.41
4	A	501	BL5	C19-C17	3.40	1.57	1.51
4	B	1501	BL5	C6-N7	-3.27	1.35	1.41
4	A	501	BL5	C45-N44	-3.23	1.35	1.41
4	A	501	BL5	C6-N7	-3.20	1.35	1.41
4	B	1501	BL5	C19-C17	2.76	1.56	1.51
3	A	401	BRR	O13-C11	-2.36	1.23	1.30
4	A	501	BL5	C32-C34	2.28	1.55	1.51
3	B	1401	BRR	O13-C11	-2.19	1.23	1.30
4	B	1501	BL5	C32-C34	2.16	1.55	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	BRR	O4-N3-C5	5.43	126.39	113.81
3	B	1401	BRR	O4-N3-C5	5.43	126.38	113.81
4	B	1501	BL5	C24-N25-N26	-4.56	106.65	111.32
4	A	501	BL5	C24-N25-N26	-3.76	107.47	111.32
3	A	401	BRR	C6-C5-N3	-3.52	107.77	112.77
4	A	501	BL5	C27-N26-N25	-3.20	108.04	111.32
3	B	1401	BRR	C6-C5-N3	-3.03	108.46	112.77
3	B	1401	BRR	O13-C11-C6	2.58	120.84	114.16
4	B	1501	BL5	C20-C19-C17	-2.48	105.60	109.63
4	B	1501	BL5	C27-N26-N25	2.36	113.73	111.32
3	B	1401	BRR	O13-C11-O12	-2.34	118.77	124.08
3	A	401	BRR	O13-C11-C6	2.30	120.13	114.16
3	B	1401	BRR	O1-C2-N3	-2.22	119.81	125.49

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	BL5	C20-C19-C24-N25
4	A	501	BL5	C19-C24-N25-N26
4	A	501	BL5	C30-C31-C32-C34
4	B	1501	BL5	C20-C19-C24-N25
4	B	1501	BL5	C19-C24-N25-N26
4	B	1501	BL5	C29-C30-C31-C32
4	A	501	BL5	C29-C30-C31-C32
3	A	401	BRR	C6-C7-C8-C9
4	A	501	BL5	C28-C29-C30-C31

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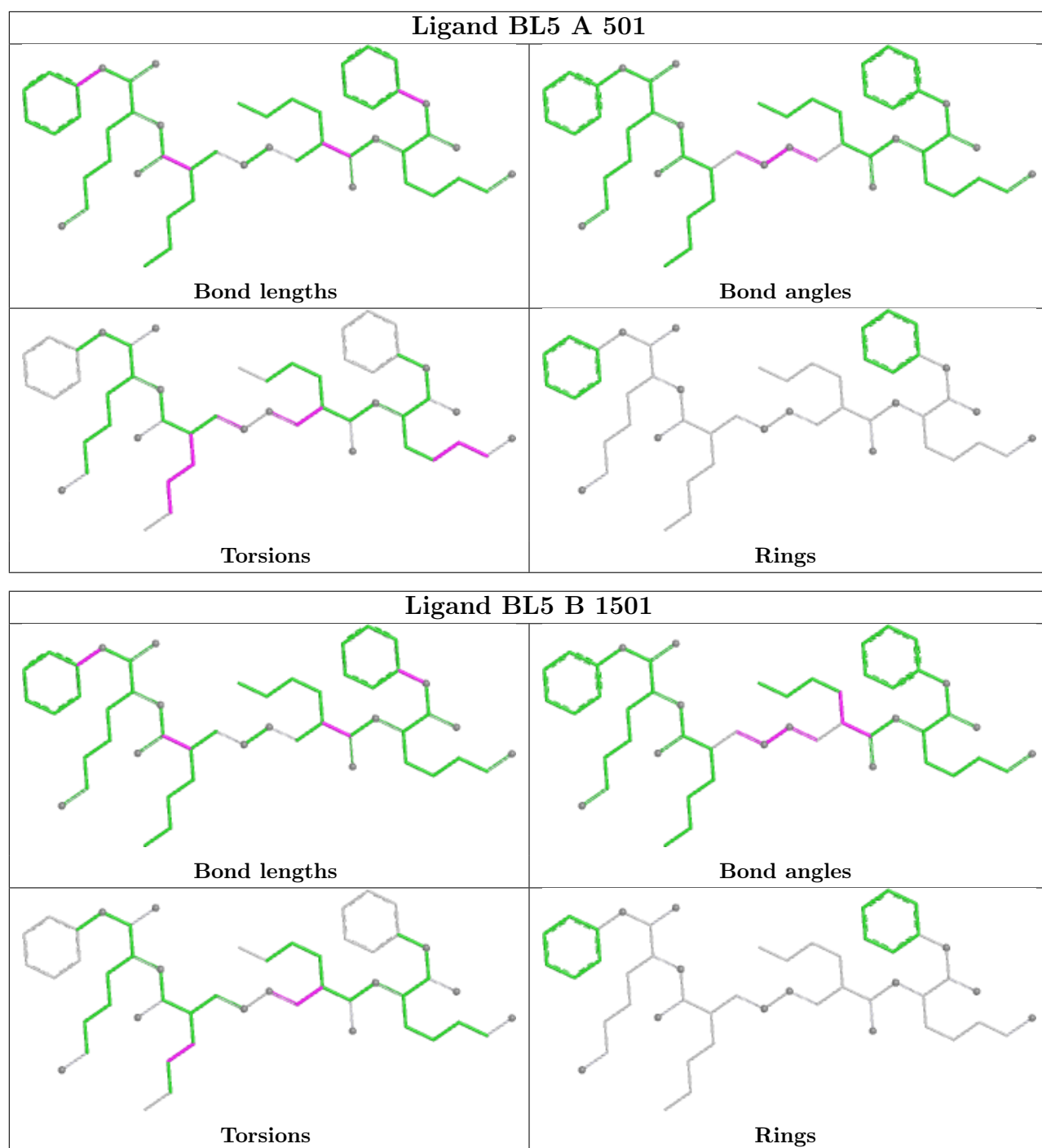
Mol	Chain	Res	Type	Atoms
4	A	501	BL5	C11-C12-C13-C14
4	B	1501	BL5	C17-C19-C24-N25
4	A	501	BL5	C12-C13-C14-N15
4	A	501	BL5	C32-C27-N26-N25
3	B	1401	BRR	C7-C8-C9-C10
4	A	501	BL5	C30-C31-C32-C27

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1401	BRR	1	0
3	A	401	BRR	4	0
4	A	501	BL5	2	0
4	B	1501	BL5	3	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	154/188 (81%)	0.23	12 (7%) 19 18	21, 34, 69, 82	0
1	B	156/188 (82%)	0.43	17 (10%) 10 9	20, 36, 72, 82	0
All	All	310/376 (82%)	0.33	29 (9%) 14 13	20, 35, 71, 82	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	232	LYS	4.9
1	A	66	LYS	4.4
1	B	66	LYS	4.4
1	B	231	TYR	4.2
1	A	230	ASP	3.7
1	B	83	THR	3.3
1	B	67	ILE	3.3
1	A	87	ASP	3.3
1	A	101	TYR	3.1
1	B	101	TYR	3.1
1	B	91	ARG	3.0
1	A	83	THR	2.9
1	B	88	ASN	2.7
1	A	88	ASN	2.7
1	A	67	ILE	2.7
1	B	123	ASN	2.6
1	A	214	GLN	2.6
1	B	86	ASP	2.5
1	B	122	TRP	2.4
1	A	89	LEU	2.3
1	A	86	ASP	2.3
1	B	214	GLN	2.3
1	B	87	ASP	2.3
1	A	106	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	215	VAL	2.2
1	B	106	ILE	2.2
1	A	123	ASN	2.2
1	B	135	ARG	2.1
1	B	104	LYS	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](#)

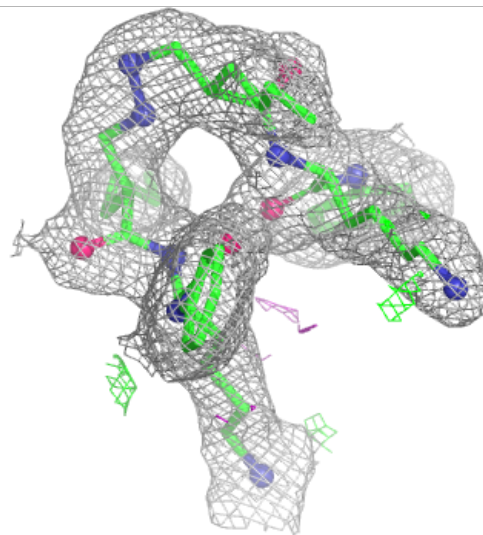
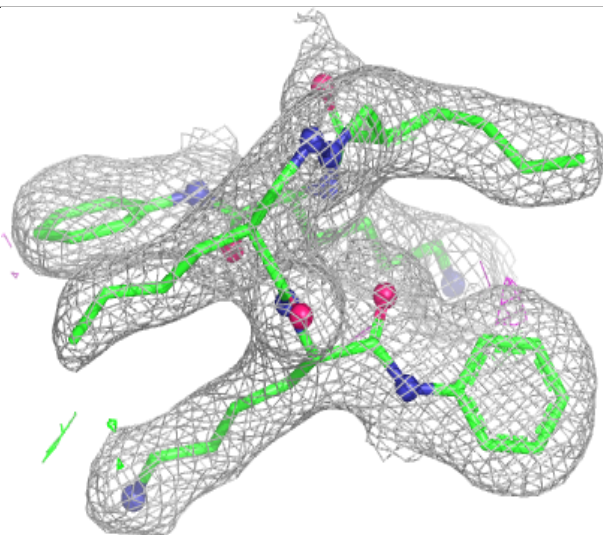
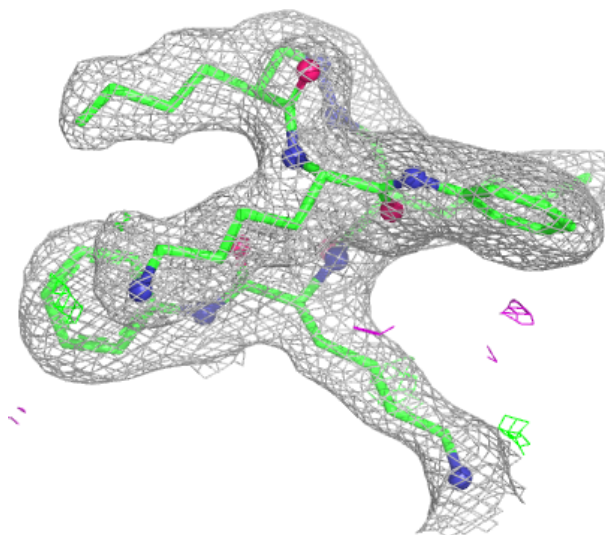
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	BL5	B	1501	50/50	0.95	0.09	24,35,56,68	0
3	BRR	B	1401	13/13	0.96	0.08	29,39,49,50	0
4	BL5	A	501	50/50	0.96	0.07	23,32,56,68	0
3	BRR	A	401	13/13	0.96	0.09	26,41,47,48	0
2	CO	B	1301	1/1	0.99	0.02	30,30,30,30	0
2	CO	A	301	1/1	1.00	0.01	27,27,27,27	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.

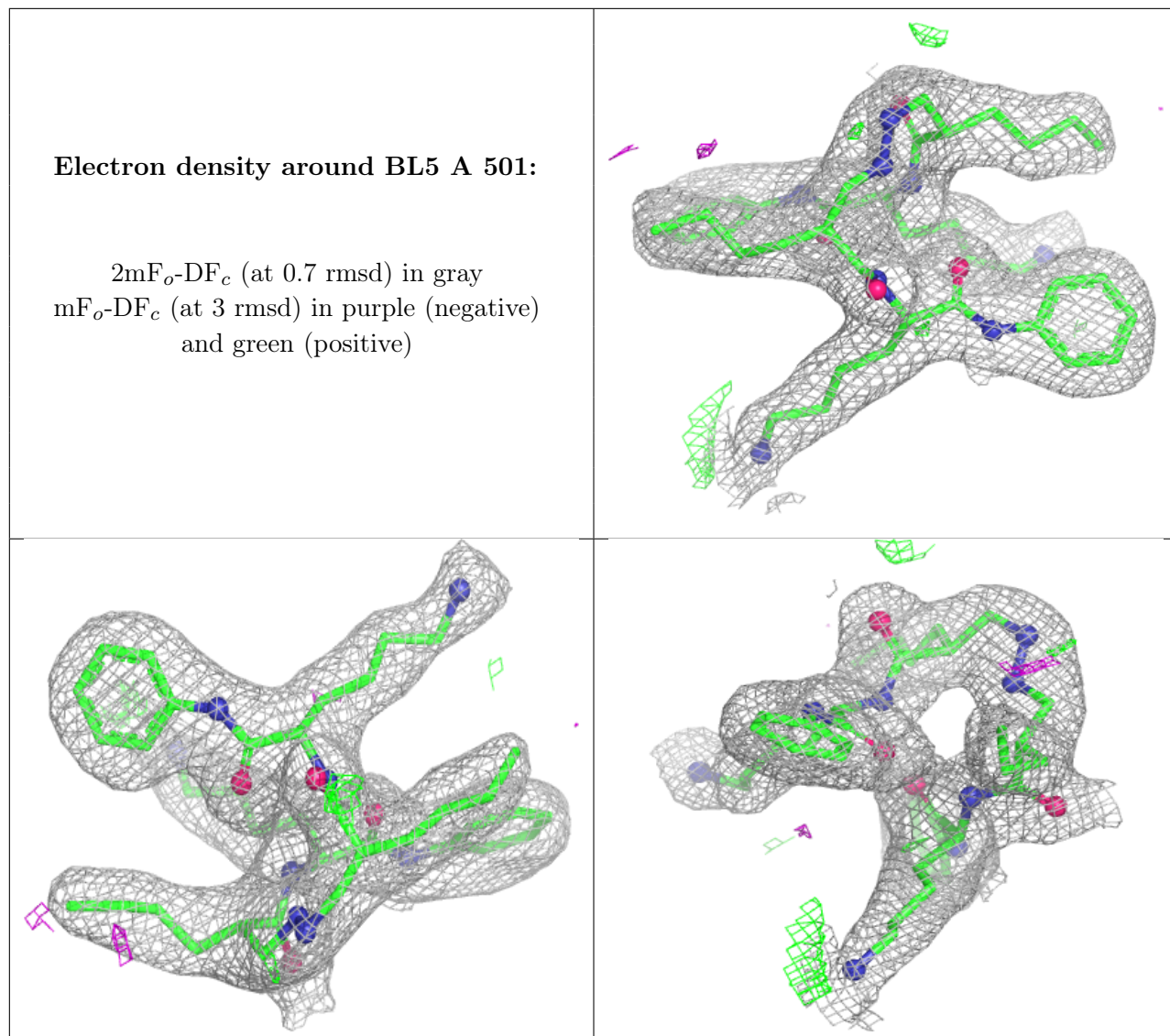
Electron density around BL5 B 1501:

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



Electron density around BL5 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.