



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:06 PM UTC

PDB ID : 8J56 / pdb_00008j56
Title : Crystal structure of the FlhDC complex from *Cupriavidus necator*
Authors : Cho, S.Y.; Oh, H.B.; Yoon, S.I.
Deposited on : 2023-04-21
Resolution : 3.50 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

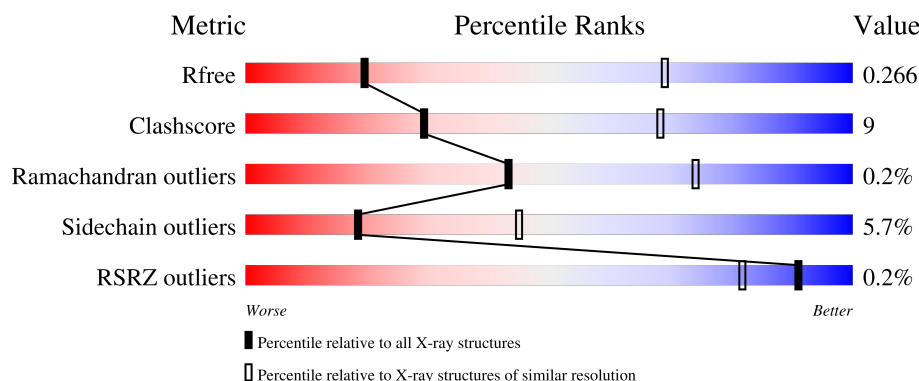
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	111	42% 15% . 41%
1	B	111	71% 17% . 9%
1	D	111	45% 12% . 41%
1	E	111	68% 15% . 15%
2	C	201	63% 14% . 20%

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Mol	Chain	Length	Quality of chain
2	F	201	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment representing 61%, a yellow segment representing 16%, and a grey segment representing 21%. The segments are separated by thin black lines. The percentages are labeled below the bar.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4863 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 Flagellar transcriptional regulator FlhD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	65	Total	C	N	O	S	0	0	0
			503	319	89	92	3			
1	B	101	Total	C	N	O	S	0	0	0
			738	467	128	140	3			
1	D	65	Total	C	N	O	S	0	0	0
			497	314	90	90	3			
1	E	94	Total	C	N	O	S	0	0	0
			679	434	116	125	4			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP A0A7W4VD94
A	-4	SER	-	expression tag	UNP A0A7W4VD94
A	-3	ALA	-	expression tag	UNP A0A7W4VD94
A	-2	LYS	-	expression tag	UNP A0A7W4VD94
A	-1	ASP	-	expression tag	UNP A0A7W4VD94
A	0	PRO	-	expression tag	UNP A0A7W4VD94
A	104	LEU	ILE	conflict	UNP A0A7W4VD94
B	-5	GLY	-	expression tag	UNP A0A7W4VD94
B	-4	SER	-	expression tag	UNP A0A7W4VD94
B	-3	ALA	-	expression tag	UNP A0A7W4VD94
B	-2	LYS	-	expression tag	UNP A0A7W4VD94
B	-1	ASP	-	expression tag	UNP A0A7W4VD94
B	0	PRO	-	expression tag	UNP A0A7W4VD94
B	104	LEU	ILE	conflict	UNP A0A7W4VD94
D	-5	GLY	-	expression tag	UNP A0A7W4VD94
D	-4	SER	-	expression tag	UNP A0A7W4VD94
D	-3	ALA	-	expression tag	UNP A0A7W4VD94
D	-2	LYS	-	expression tag	UNP A0A7W4VD94
D	-1	ASP	-	expression tag	UNP A0A7W4VD94
D	0	PRO	-	expression tag	UNP A0A7W4VD94
D	104	LEU	ILE	conflict	UNP A0A7W4VD94

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-5	GLY	-	expression tag	UNP A0A7W4VD94
E	-4	SER	-	expression tag	UNP A0A7W4VD94
E	-3	ALA	-	expression tag	UNP A0A7W4VD94
E	-2	LYS	-	expression tag	UNP A0A7W4VD94
E	-1	ASP	-	expression tag	UNP A0A7W4VD94
E	0	PRO	-	expression tag	UNP A0A7W4VD94
E	104	LEU	ILE	conflict	UNP A0A7W4VD94

- Molecule 2 is a protein called Flagellar transcriptional regulator FlhC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	160	Total	C	N	O	S	0	0	0
			1230	792	207	223	8			
2	F	158	Total	C	N	O	S	0	0	0
			1214	786	205	215	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	LEU	MET	conflict	UNP A0A7T4G0G8
F	1	LEU	MET	conflict	UNP A0A7T4G0G8

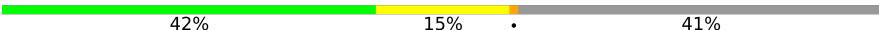
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

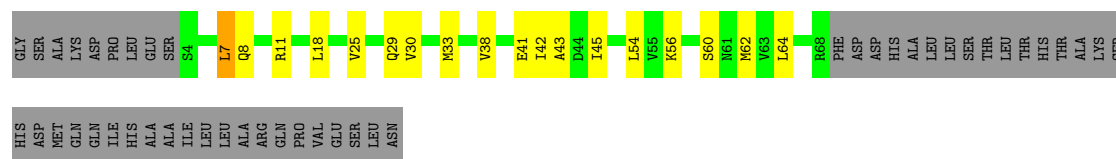
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Flagellar transcriptional regulator FlhD

Chain A: 

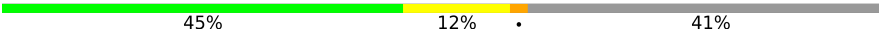


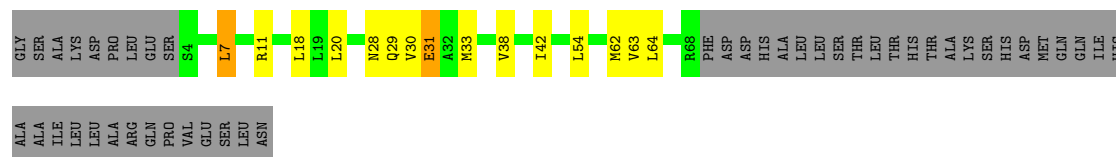
• Molecule 1: Flagellar transcriptional regulator FlhD

Chain B: 



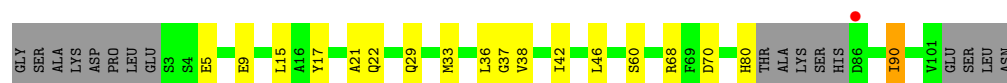
• Molecule 1: Flagellar transcriptional regulator FlhD

Chain D: 



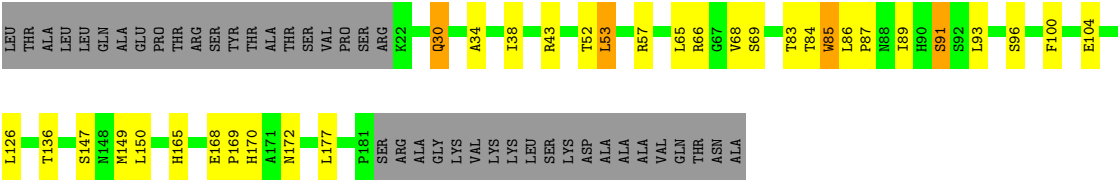
• Molecule 1: Flagellar transcriptional regulator FlhD

Chain E: 

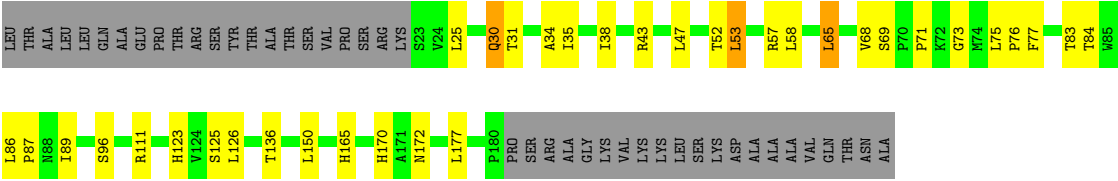


• Molecule 2: Flagellar transcriptional regulator FlhC

Chain C: 



• Molecule 2: Flagellar transcriptional regulator FlhC



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	124.53Å 124.53Å 173.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.91 – 3.50 29.91 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.91-3.50) 99.5 (29.91-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 3.47Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.220 , 0.261 0.221 , 0.266	Depositor DCC
R_{free} test set	949 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	95.4	Xtriage
Anisotropy	0.332	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 85.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.044 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4863	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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:
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	0.17	0/505	0.32	0/680
1	B	0.16	0/745	0.35	0/1015
1	D	0.17	0/499	0.34	0/674
1	E	0.15	0/685	0.33	0/933
2	C	0.17	0/1260	0.35	0/1718
2	F	0.17	0/1244	0.38	0/1696
All	All	0.17	0/4938	0.35	0/6716

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	503	0	525	15	0
1	B	738	0	701	14	0
1	D	497	0	510	13	0
1	E	679	0	650	11	0
2	C	1230	0	1172	23	0
2	F	1214	0	1169	20	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	1	0	0	0	0
All	All	4863	0	4727	82	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 9.

All (82) 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)
2:C:43:ARG:HB2	2:C:136:THR:HG21	1.61	0.82
1:A:29:GLN:HG2	1:A:33:MET:HE2	1.63	0.81
1:D:29:GLN:HG2	1:D:33:MET:HE2	1.67	0.76
2:C:38:ILE:O	2:C:66:ARG:NH1	2.26	0.68
1:B:55:VAL:HG22	1:E:15:LEU:HD23	1.76	0.67
2:F:43:ARG:HB2	2:F:136:THR:HG21	1.79	0.65
1:D:63:VAL:HG11	1:E:17:TYR:HE2	1.62	0.64
2:F:75:LEU:HD13	2:F:77:PHE:HE1	1.62	0.64
2:F:30:GLN:HG2	2:F:52:THR:H	1.63	0.62
1:D:30:VAL:HA	1:D:33:MET:HE3	1.83	0.61
1:B:90:ILE:HG23	2:C:93:LEU:HD22	1.82	0.60
2:C:30:GLN:HG2	2:C:52:THR:H	1.67	0.59
1:A:41:GLU:O	1:A:45:ILE:HG12	2.04	0.56
2:C:85:TRP:O	2:C:89:ILE:HG13	2.06	0.56
1:E:21:ALA:HB1	1:E:46:LEU:HD13	1.86	0.56
2:C:83:THR:HG21	2:C:165:HIS:CE1	2.41	0.55
1:D:18:LEU:HB3	1:D:54:LEU:HD22	1.89	0.55
1:B:30:VAL:HA	1:B:33:MET:HE3	1.88	0.54
1:B:3:SER:HA	1:B:7:LEU:HD23	1.89	0.54
2:F:89:ILE:HD13	2:F:177:LEU:HD21	1.90	0.54
1:A:30:VAL:HA	1:A:33:MET:HE3	1.91	0.53
2:F:84:THR:HB	2:F:87:PRO:HD2	1.91	0.53
1:D:20:LEU:HD13	1:E:9:GLU:HG2	1.91	0.53
2:F:47:LEU:HB2	2:F:58:LEU:HD13	1.90	0.52
2:F:35:ILE:HG23	2:F:65:LEU:HD11	1.92	0.51
2:F:43:ARG:HD3	2:F:73:GLY:O	2.12	0.50
2:F:43:ARG:H	2:F:136:THR:CG2	2.25	0.49
1:A:7:LEU:HG	2:C:126:LEU:HD21	1.93	0.49
2:F:75:LEU:HD13	2:F:77:PHE:CE1	2.43	0.49
2:C:89:ILE:HD13	2:C:177:LEU:HD21	1.94	0.49
2:C:84:THR:HB	2:C:87:PRO:HD2	1.95	0.48
1:B:31:GLU:HB3	1:B:35:ARG:NH2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:VAL:HG22	1:B:42:ILE:HB	1.96	0.48
1:A:18:LEU:HB3	1:A:54:LEU:HD22	1.97	0.47
1:B:98:ARG:NH1	2:C:96:SER:HB3	2.30	0.47
2:C:68:VAL:HG22	2:C:69:SER:H	1.80	0.47
1:E:90:ILE:HD13	2:F:123:HIS:CD2	2.50	0.47
1:E:38:VAL:HG22	1:E:42:ILE:HB	1.97	0.47
2:F:68:VAL:HG22	2:F:69:SER:H	1.80	0.46
1:A:62:MET:HE2	1:A:62:MET:HB3	1.71	0.46
1:E:29:GLN:O	1:E:33:MET:HG3	2.15	0.46
1:D:29:GLN:O	1:D:33:MET:HG3	2.16	0.46
2:F:34:ALA:O	2:F:38:ILE:HG13	2.16	0.46
2:C:93:LEU:O	2:C:96:SER:OG	2.32	0.45
1:D:38:VAL:HG22	1:D:42:ILE:HB	1.99	0.45
1:A:38:VAL:HG22	1:A:42:ILE:HB	1.98	0.45
1:B:91:HIS:CE1	1:B:95:LEU:HD11	2.52	0.45
2:C:43:ARG:H	2:C:136:THR:CG2	2.29	0.45
2:C:100:PHE:CE2	2:C:104:GLU:HB3	2.51	0.45
2:C:30:GLN:HB3	2:C:53:LEU:HD13	1.99	0.44
1:A:64:LEU:HD22	1:B:69:PHE:HE2	1.82	0.44
1:A:29:GLN:O	1:A:33:MET:HG3	2.18	0.44
1:D:64:LEU:HD23	1:D:64:LEU:HA	1.86	0.44
2:F:170:HIS:C	2:F:172:ASN:H	2.25	0.44
1:A:8:GLN:OE1	1:A:11:ARG:NH1	2.50	0.44
1:B:29:GLN:O	1:B:33:MET:HG3	2.18	0.43
2:C:86:LEU:HB3	2:C:87:PRO:HD3	2.00	0.43
1:E:22:GLN:HG3	1:E:46:LEU:O	2.19	0.43
1:A:56:LYS:HE3	1:B:89:GLN:HE21	1.83	0.43
1:D:63:VAL:HG21	1:E:36:LEU:HD23	2.00	0.43
2:C:149:MET:O	2:C:150:LEU:HD12	2.19	0.43
1:A:56:LYS:HE2	1:A:56:LYS:HB3	1.80	0.43
1:D:28:ASN:CG	1:D:31:GLU:HB2	2.44	0.42
2:F:43:ARG:NH2	2:F:71:PRO:HB2	2.34	0.42
1:A:25:VAL:HG11	1:A:43:ALA:O	2.19	0.42
1:D:7:LEU:HG	2:F:126:LEU:HD21	2.02	0.42
2:C:168:GLU:HG3	2:C:169:PRO:HD2	2.01	0.42
2:C:87:PRO:O	2:C:91:SER:OG	2.31	0.42
2:F:111:ARG:HA	2:F:111:ARG:HD3	1.92	0.42
1:A:64:LEU:HD23	1:A:64:LEU:HA	1.81	0.41
2:C:170:HIS:C	2:C:172:ASN:H	2.28	0.41
1:D:62:MET:HE2	1:D:62:MET:HB3	1.77	0.41
1:A:64:LEU:O	1:B:68:ARG:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:147:SER:OG	2:C:149:MET:HE3	2.20	0.41
1:D:42:ILE:HD11	1:E:70:ASP:HA	2.03	0.41
2:F:31:THR:HA	2:F:53:LEU:HD21	2.03	0.41
1:B:39:SER:HB2	1:B:41:GLU:OE1	2.20	0.41
2:C:65:LEU:CD2	2:F:25:LEU:HD21	2.51	0.41
1:B:35:ARG:O	1:B:68:ARG:NH1	2.55	0.40
2:C:34:ALA:O	2:C:38:ILE:HG13	2.21	0.40
1:E:37:GLY:HA3	1:E:68:ARG:NH2	2.37	0.40
2:F:83:THR:HG21	2:F:165:HIS:CE1	2.57	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	63/111 (57%)	61 (97%)	2 (3%)	0	100	100
1	B	97/111 (87%)	88 (91%)	9 (9%)	0	100	100
1	D	63/111 (57%)	61 (97%)	2 (3%)	0	100	100
1	E	90/111 (81%)	87 (97%)	3 (3%)	0	100	100
2	C	158/201 (79%)	149 (94%)	9 (6%)	0	100	100
2	F	156/201 (78%)	147 (94%)	8 (5%)	1 (1%)	21	54
All	All	627/846 (74%)	593 (95%)	33 (5%)	1 (0%)	43	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	76	PRO

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	53/95 (56%)	51 (96%)	2 (4%)	29	55
1	B	71/95 (75%)	65 (92%)	6 (8%)	10	34
1	D	51/95 (54%)	48 (94%)	3 (6%)	18	44
1	E	64/95 (67%)	60 (94%)	4 (6%)	16	42
2	C	127/171 (74%)	122 (96%)	5 (4%)	28	54
2	F	125/171 (73%)	117 (94%)	8 (6%)	16	42
All	All	491/722 (68%)	463 (94%)	28 (6%)	18	45

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	60	SER
1	B	7	LEU
1	B	35	ARG
1	B	42	ILE
1	B	51	SER
1	B	63	VAL
1	B	80	HIS
2	C	30	GLN
2	C	53	LEU
2	C	57	ARG
2	C	85	TRP
2	C	91	SER
1	D	7	LEU
1	D	11	ARG
1	D	31	GLU
1	E	5	GLU
1	E	60	SER
1	E	80	HIS
1	E	90	ILE
2	F	30	GLN
2	F	53	LEU

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Mol	Chain	Res	Type
2	F	57	ARG
2	F	65	LEU
2	F	86	LEU
2	F	96	SER
2	F	125	SER
2	F	150	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	89	GLN
2	F	148	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å²)	Q<0.9
1	A	65/111 (58%)	-0.50	0	100100	51, 74, 101, 129	0
1	B	101/111 (90%)	-0.01	0	100100	63, 99, 142, 189	0
1	D	65/111 (58%)	-0.41	0	100100	59, 88, 131, 151	0
1	E	94/111 (84%)	-0.09	1 (1%)	7854	57, 100, 157, 194	0
2	C	160/201 (79%)	-0.23	0	100100	60, 92, 134, 154	0
2	F	158/201 (78%)	-0.21	0	100100	64, 93, 136, 151	0
All	All	643/846 (76%)	-0.21	1 (0%)	9182	51, 92, 138, 194	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	86	ASP	2.3

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](#)

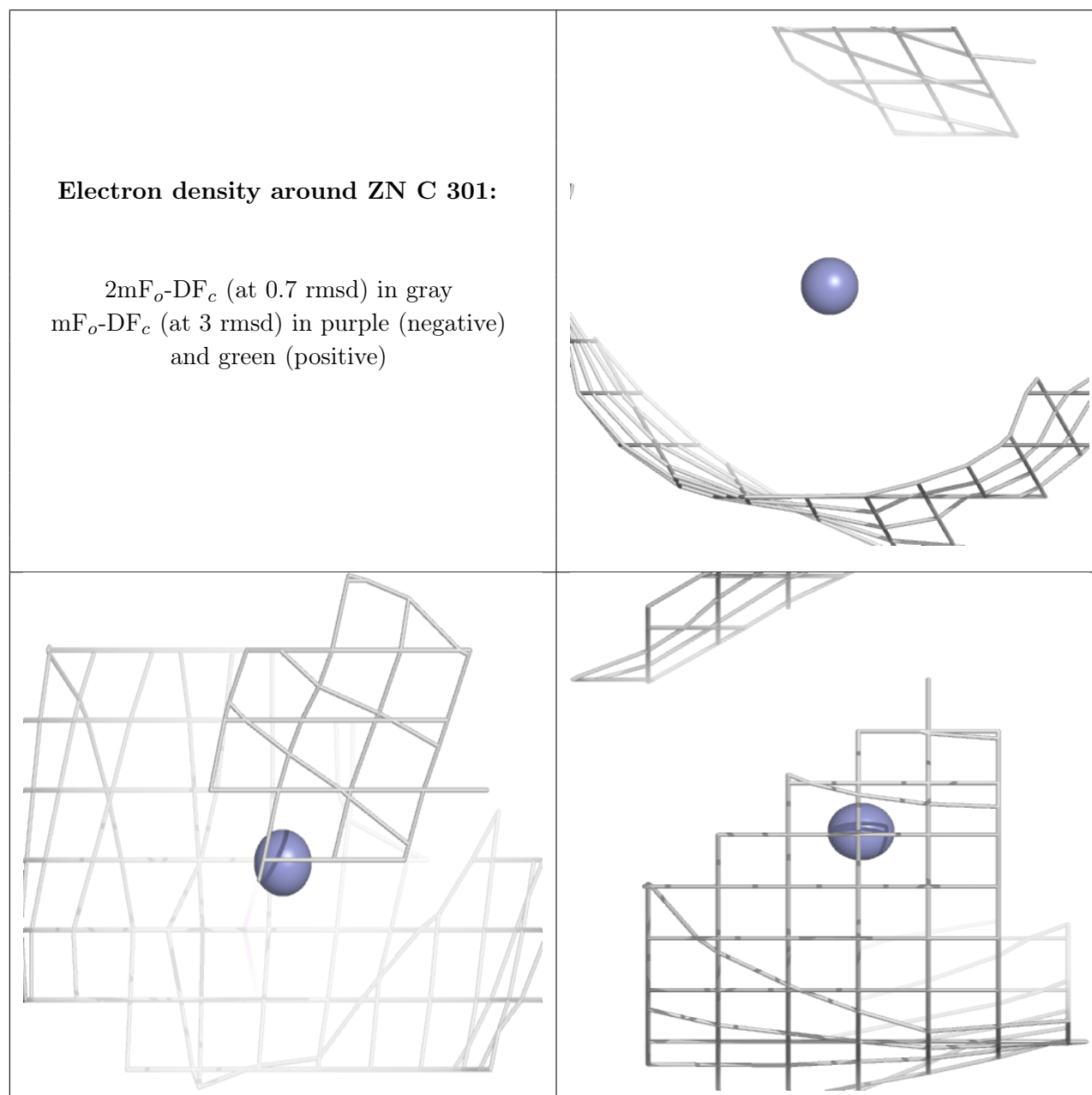
There are no oligosaccharides in this entry.

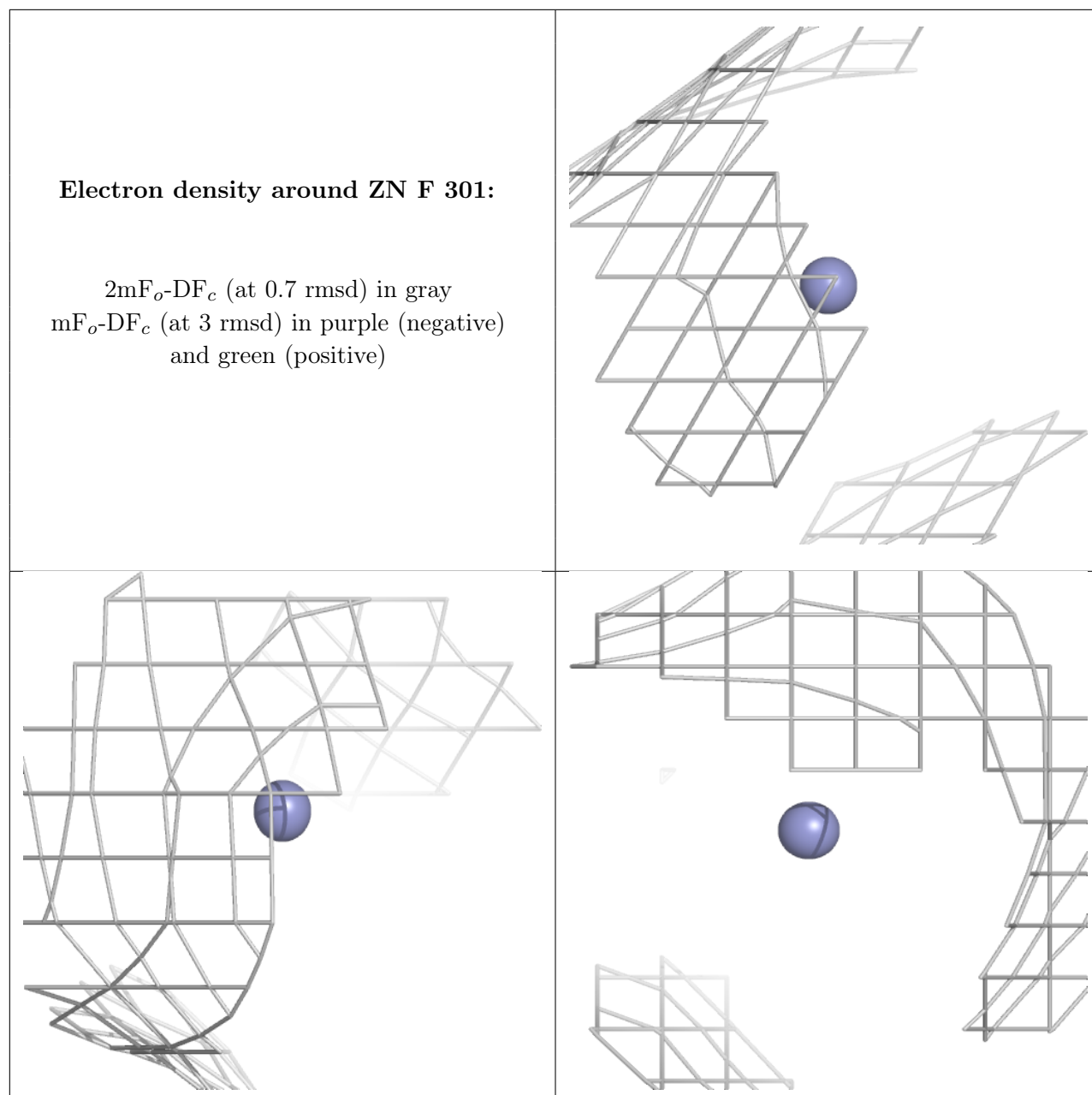
6.4 Ligands [i](#)

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
3	ZN	C	301	1/1	1.00	0.02	88,88,88,88	0
3	ZN	F	301	1/1	1.00	0.01	90,90,90,90	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.





6.5 Other polymers ⓘ

There are no such residues in this entry.