



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:41 AM UTC

PDB ID : 7ULT / pdb_00007ult
Title : Crystal Structure of SARS-CoV-2 Nsp16/10 Heterodimer Apo-Form.
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Deposited on : 2022-04-05
Resolution : 1.90 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
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

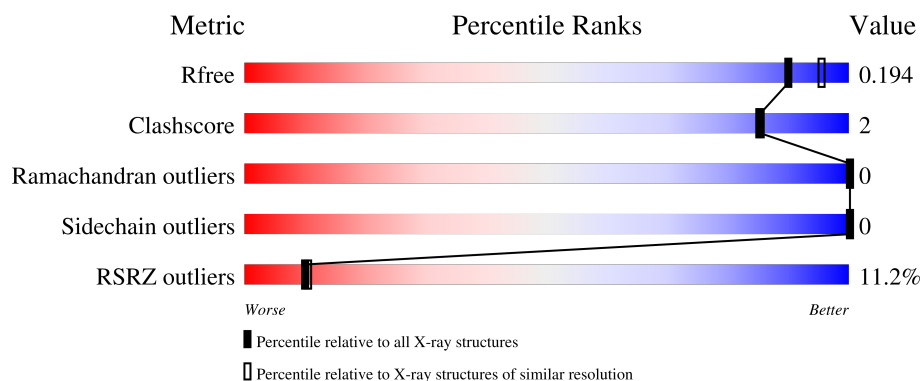
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	300	12% 90% 7% .
1	C	300	9% 96% . ..
2	B	141	14% 79% 7% 13%
2	D	141	6% 78% . 18%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7433 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 2'-O-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	14	0
			2412	1543	403	448	18			
1	C	298	Total	C	N	O	S	0	9	0
			2403	1528	403	454	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6797	SER	-	expression tag	UNP P0DTD1
A	6798	ASN	-	expression tag	UNP P0DTD1
C	6797	SER	-	expression tag	UNP P0DTD1
C	6798	ASN	-	expression tag	UNP P0DTD1

- Molecule 2 is a protein called Non-structural protein 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	122	Total	C	N	O	S	0	1	0
			916	569	155	176	16			
2	D	115	Total	C	N	O	S	0	1	0
			855	531	144	165	15			

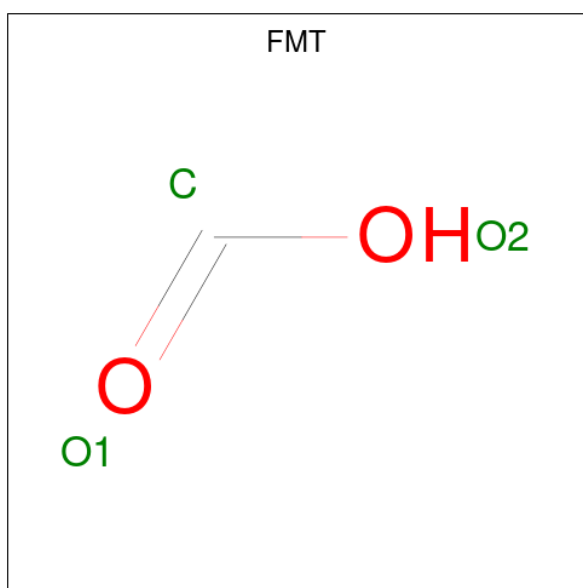
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	4252	SER	-	expression tag	UNP P0DTD1
B	4253	ASN	-	expression tag	UNP P0DTD1
D	4252	SER	-	expression tag	UNP P0DTD1
D	4253	ASN	-	expression tag	UNP P0DTD1

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Na 2 2	0	0
3	B	2	Total Na 2 2	0	0
3	C	3	Total Na 3 3	0	0
3	D	1	Total Na 1 1	0	0

- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	C	1	Total C O 3 1 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Zn	0	0
			2	2		
5	D	2	Total	Zn	0	0
			2	2		

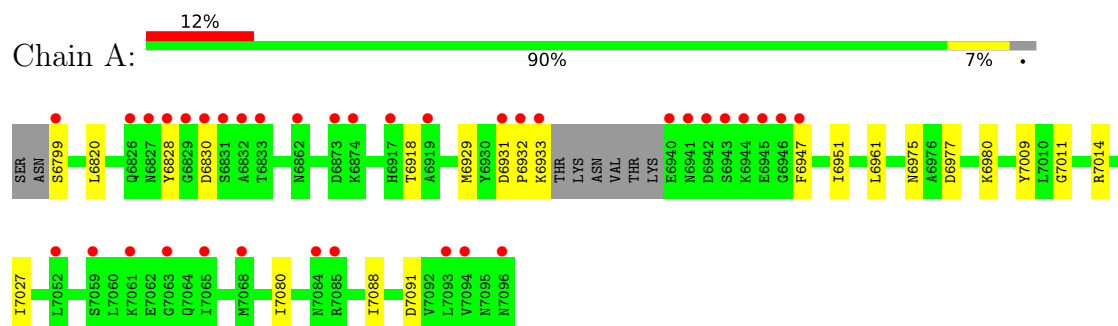
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	263	Total	O	0	9
			266	266		
6	B	112	Total	O	0	2
			113	113		
6	C	287	Total	O	0	13
			296	296		
6	D	106	Total	O	0	3
			109	109		

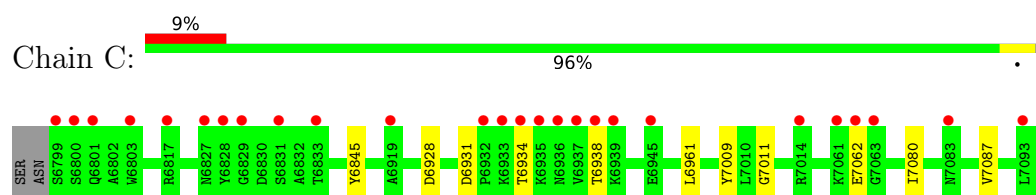
3 Residue-property plots [i](#)

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.

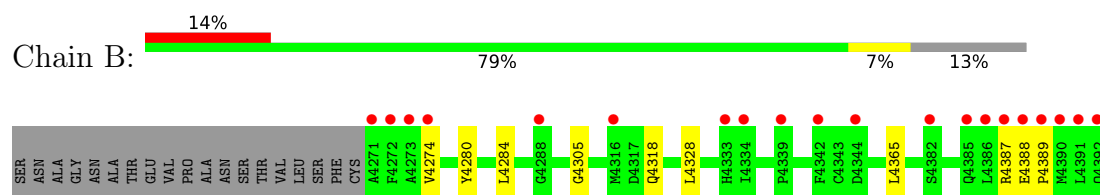
• Molecule 1: 2'-O-methyltransferase



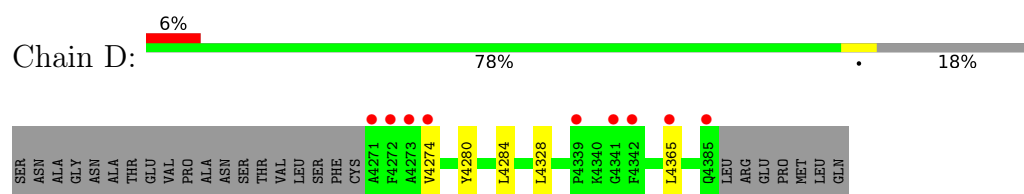
• Molecule 1: 2'-O-methyltransferase



• Molecule 2: Non-structural protein 10



• Molecule 2: Non-structural protein 10



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	166.68Å 166.68Å 98.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.90 – 1.90 29.90 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.90-1.90) 99.7 (29.90-1.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.164 , 0.188 0.171 , 0.194	Depositor DCC
R_{free} test set	6205 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7433	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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: FMT, ZN, NA

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	0.96	0/2465	1.13	0/3345
1	C	0.96	0/2454	1.15	0/3327
2	B	1.02	0/936	1.20	0/1268
2	D	1.01	0/874	1.19	0/1186
All	All	0.98	0/6729	1.16	0/9126

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	2412	0	2381	15	0
1	C	2403	0	2373	7	0
2	B	916	0	876	6	0
2	D	855	0	811	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	3	0	0	0	0
3	D	1	0	0	0	0
4	A	21	0	7	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	27	0	9	0	0
4	D	3	0	1	0	0
5	B	2	0	0	0	0
5	D	2	0	0	0	0
6	A	266	0	0	2	0
6	B	113	0	0	0	0
6	C	296	0	0	0	0
6	D	109	0	0	0	0
All	All	7433	0	6458	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 2.

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)
2:B:4388:GLU:HB3	2:B:4389:PRO:HD3	1.75	0.66
2:B:4328:LEU:HD22	2:B:4365:LEU:HD11	1.78	0.65
2:D:4328:LEU:HD22	2:D:4365:LEU:HD11	1.83	0.60
1:A:6918:THR:HG22	1:A:7088:ILE:HG22	1.86	0.58
1:A:6929[A]:MET:HE3	1:A:6947:PHE:CD1	2.40	0.56
2:B:4387:ARG:HG2	2:B:4387:ARG:O	2.07	0.55
1:A:6980:LYS:NZ	6:A:7206[B]:HOH:O	2.44	0.51
1:C:6931:ASP:O	1:C:6934:THR:HG22	2.10	0.51
1:A:6931:ASP:HB3	1:A:6947:PHE:CE2	2.47	0.49
2:D:4280:TYR:CE2	2:D:4284:LEU:HD11	2.48	0.49
1:C:7062:GLU:OE1	1:C:7062:GLU:N	2.44	0.48
2:B:4274:VAL:CG1	2:D:4274:VAL:HG11	2.46	0.45
1:C:6961:LEU:HB2	1:C:7080:ILE:HB	1.97	0.45
1:C:7009:TYR:CZ	1:C:7011:GLY:HA2	2.52	0.45
1:A:6961:LEU:HB2	1:A:7080:ILE:HB	1.99	0.45
1:C:6845:TYR:OH	1:C:6928:ASP:HB2	2.18	0.44
2:B:4280:TYR:CE2	2:B:4284:LEU:HD11	2.53	0.44
1:A:6975:ASN:OD1	1:A:6977:ASP:HB2	2.18	0.43
1:A:7091:ASP:HB3	1:C:7087:VAL:HG13	2.00	0.43
1:A:6951[A]:ILE:HD12	1:A:6951[A]:ILE:HA	1.89	0.43
1:C:6938:THR:HG22	1:C:6938:THR:O	2.18	0.42
2:B:4305:GLY:HA2	2:B:4318:GLN:OE1	2.20	0.42
1:A:6820:LEU:HD11	1:A:7027[B]:ILE:HD12	2.02	0.41
1:A:6799:SER:N	6:A:7219:HOH:O	2.53	0.41
1:A:6828:TYR:C	1:A:6830:ASP:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6932:PRO:O	1:A:6933:LYS:HB2	2.19	0.41
1:A:6947:PHE:CD1	1:A:6947:PHE:C	2.99	0.41
1:A:7014:ARG:NH1	4:A:7104:FMT:O2	2.45	0.40
1:A:7009:TYR:CZ	1:A:7011:GLY:HA2	2.56	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	302/300 (101%)	293 (97%)	9 (3%)	0	100	100
1	C	305/300 (102%)	297 (97%)	8 (3%)	0	100	100
2	B	121/141 (86%)	116 (96%)	5 (4%)	0	100	100
2	D	114/141 (81%)	112 (98%)	2 (2%)	0	100	100
All	All	842/882 (96%)	818 (97%)	24 (3%)	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	266/260 (102%)	266 (100%)	0	100	100
1	C	266/260 (102%)	266 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	101/115 (88%)	101 (100%)	0	100	100
2	D	94/115 (82%)	94 (100%)	0	100	100
All	All	727/750 (97%)	727 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	4293	ASN
2	B	4367	ASN
2	D	4293	ASN
2	D	4367	ASN

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 29 ligands modelled in this entry, 12 are monoatomic - leaving 17 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
4	FMT	C	7110	-	2,2,2	0.21	0	1,1,1	0.17	0
4	FMT	A	7105	-	2,2,2	0.32	0	1,1,1	0.14	0
4	FMT	D	4404	-	2,2,2	0.24	0	1,1,1	0.16	0
4	FMT	A	7106	-	2,2,2	0.22	0	1,1,1	0.16	0
4	FMT	A	7104	-	2,2,2	0.25	0	1,1,1	0.16	0
4	FMT	A	7109	-	2,2,2	0.23	0	1,1,1	0.15	0
4	FMT	A	7108	-	2,2,2	0.25	0	1,1,1	0.12	0
4	FMT	C	7108	-	2,2,2	0.25	0	1,1,1	0.15	0
4	FMT	C	7112	-	2,2,2	0.20	0	1,1,1	0.15	0
4	FMT	A	7103	-	2,2,2	0.22	0	1,1,1	0.19	0
4	FMT	C	7105	-	2,2,2	0.24	0	1,1,1	0.19	0
4	FMT	C	7109	-	2,2,2	0.26	0	1,1,1	0.16	0
4	FMT	A	7107	-	2,2,2	0.25	0	1,1,1	0.17	0
4	FMT	C	7104	-	2,2,2	0.24	0	1,1,1	0.18	0
4	FMT	C	7107	-	2,2,2	0.22	0	1,1,1	0.17	0
4	FMT	C	7106	-	2,2,2	0.23	0	1,1,1	0.17	0
4	FMT	C	7111	-	2,2,2	0.27	0	1,1,1	0.14	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	7104	FMT	1	0

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	292/300 (97%)	0.34	36 (12%) 8 9	11, 30, 63, 102	14 (4%)
1	C	298/300 (99%)	0.18	28 (9%) 14 14	12, 28, 69, 120	9 (3%)
2	B	122/141 (86%)	0.69	20 (16%) 4 4	18, 42, 76, 120	1 (0%)
2	D	115/141 (81%)	0.37	9 (7%) 19 20	21, 38, 64, 104	1 (0%)
All	All	827/882 (93%)	0.34	93 (11%) 10 10	11, 31, 69, 120	25 (3%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	6937	VAL	9.7
1	A	6944	LYS	8.8
1	C	6938	THR	8.4
2	B	4386	LEU	6.1
1	C	6803	TRP	5.9
2	B	4271	ALA	5.7
2	D	4271	ALA	5.7
1	A	6933	LYS	5.3
2	B	4391	LEU	5.0
2	B	4390	MET	5.0
2	B	4387	ARG	4.9
1	A	6799	SER	4.9
1	A	6942	ASP	4.8
2	B	4389	PRO	4.8
1	A	6940	GLU	4.7
1	C	6934	THR	4.6
1	C	6799	SER	4.6
2	B	4392	GLN	4.6
2	B	4388	GLU	4.5
1	C	6939	LYS	4.5
1	C	6935	LYS	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	7096	ASN	4.1
1	C	6932	PRO	4.0
1	A	6828	TYR	3.9
2	B	4333	HIS	3.9
1	A	6941	ASN	3.7
1	C	6828	TYR	3.7
1	C	7062	GLU	3.7
1	C	6933	LYS	3.7
1	A	6832	ALA	3.6
1	A	6931	ASP	3.6
1	C	6936	ASN	3.5
1	C	6800	SER	3.5
1	A	7085	ARG	3.4
1	A	6947	PHE	3.4
2	D	4273	ALA	3.4
1	A	6945	GLU	3.4
1	A	6943	SER	3.4
1	A	6833	THR	3.4
1	A	7068	MET	3.4
1	A	6827	ASN	3.4
1	C	6945	GLU	3.3
1	A	7063	GLY	3.3
1	A	7052	LEU	3.3
1	A	6829	GLY	3.2
1	C	6829	GLY	3.2
2	B	4274	VAL	3.1
1	C	6831[A]	SER	2.9
1	C	7095	ASN	2.9
1	A	7065	ILE	2.9
2	B	4334	ILE	2.9
1	C	6827	ASN	2.9
2	B	4273	ALA	2.9
1	C	6833	THR	2.8
2	D	4385	GLN	2.8
1	A	6932	PRO	2.8
1	A	6830	ASP	2.8
1	A	6919	ALA	2.8
1	A	7059	SER	2.8
1	C	7061	LYS	2.8
2	D	4342	PHE	2.7
2	B	4339	PRO	2.7
2	B	4272	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	4342	PHE	2.7
1	C	6817	ARG	2.6
1	C	7063	GLY	2.6
1	A	7094	VAL	2.4
1	A	6946	GLY	2.4
1	C	7083	ASN	2.4
2	B	4344	ASP	2.3
2	D	4339	PRO	2.3
1	A	6873	ASP	2.3
1	A	6826	GLN	2.3
1	A	6862	ASN	2.3
1	A	6874	LYS	2.3
2	D	4274	VAL	2.3
1	C	6919	ALA	2.2
2	D	4341	GLY	2.2
1	A	6831	SER	2.2
2	B	4288	GLY	2.2
1	A	7061	LYS	2.2
1	A	7096	ASN	2.2
1	A	6917	HIS	2.1
1	A	7084	ASN	2.1
1	C	6801	GLN	2.1
2	B	4385	GLN	2.1
2	D	4272	PHE	2.1
1	C	7014	ARG	2.1
2	B	4316	MET	2.1
2	D	4365	LEU	2.1
1	C	7093	LEU	2.1
2	B	4382	SER	2.0
1	A	7093	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

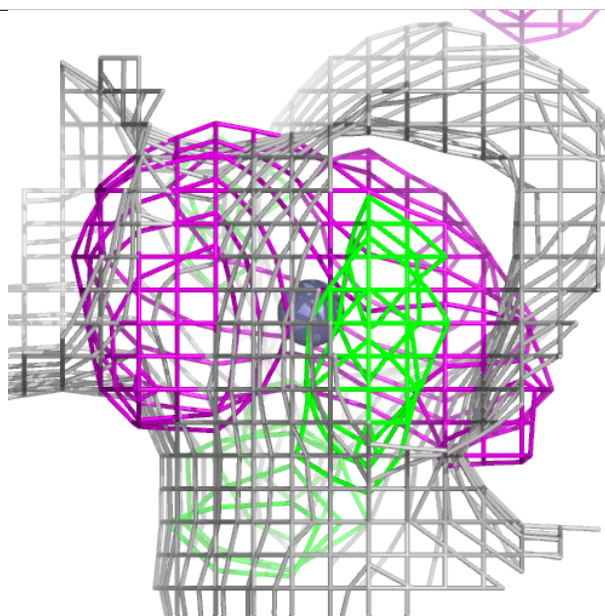
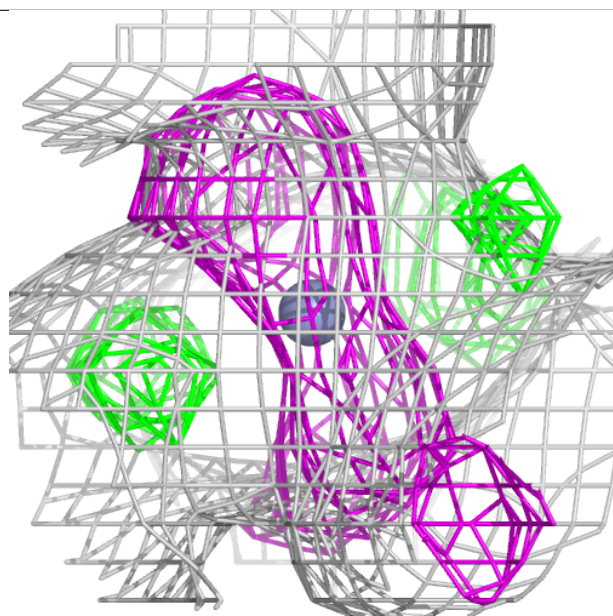
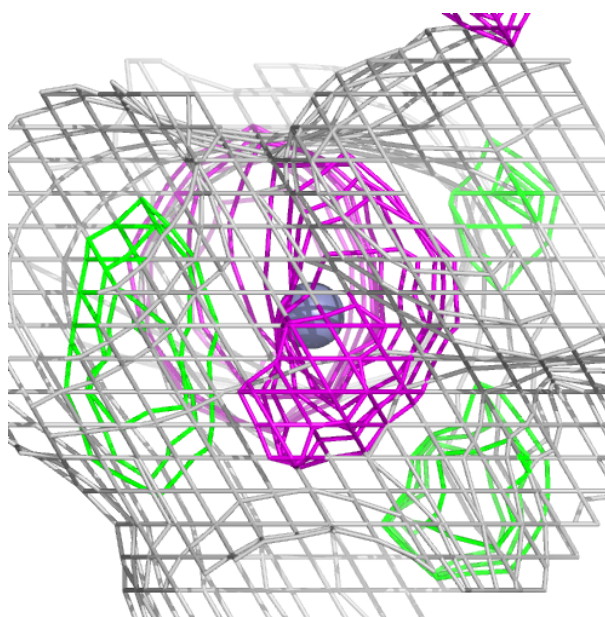
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	FMT	C	7112	3/3	0.78	0.23	57,57,63,71	0
4	FMT	A	7107	3/3	0.82	0.24	67,67,69,70	0
4	FMT	C	7111	3/3	0.87	0.18	46,46,53,57	0
4	FMT	A	7108	3/3	0.89	0.35	60,60,60,61	0
3	NA	B	4403	1/1	0.90	0.12	57,57,57,57	0
4	FMT	A	7105	3/3	0.90	0.19	54,54,59,62	0
4	FMT	C	7107	3/3	0.91	0.14	52,52,53,54	0
4	FMT	C	7109	3/3	0.91	0.23	49,49,53,59	0
3	NA	B	4404	1/1	0.91	0.11	54,54,54,54	0
4	FMT	A	7109	3/3	0.91	0.15	45,45,50,56	0
3	NA	D	4403	1/1	0.92	0.09	65,65,65,65	0
4	FMT	C	7106	3/3	0.92	0.14	45,45,52,53	0
4	FMT	A	7106	3/3	0.93	0.22	53,53,55,59	0
4	FMT	C	7110	3/3	0.93	0.21	38,38,43,52	0
4	FMT	D	4404	3/3	0.93	0.12	53,53,59,62	0
4	FMT	C	7108	3/3	0.94	0.10	53,53,53,56	0
4	FMT	C	7105	3/3	0.95	0.11	31,31,31,34	0
3	NA	C	7101	1/1	0.95	0.20	39,39,39,39	0
3	NA	C	7103	1/1	0.95	0.08	56,56,56,56	0
3	NA	A	7102	1/1	0.96	0.14	55,55,55,55	0
5	ZN	B	4402	1/1	0.96	0.08	45,45,45,45	0
5	ZN	D	4402	1/1	0.96	0.07	46,46,46,46	0
4	FMT	A	7103	3/3	0.97	0.10	29,29,30,31	0
4	FMT	A	7104	3/3	0.97	0.08	40,40,47,48	0
4	FMT	C	7104	3/3	0.98	0.05	35,35,36,37	0
5	ZN	B	4401	1/1	0.98	0.04	34,34,34,34	0
3	NA	C	7102	1/1	0.98	0.12	29,29,29,29	0
5	ZN	D	4401	1/1	0.98	0.04	30,30,30,30	0
3	NA	A	7101	1/1	0.98	0.10	37,37,37,37	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.

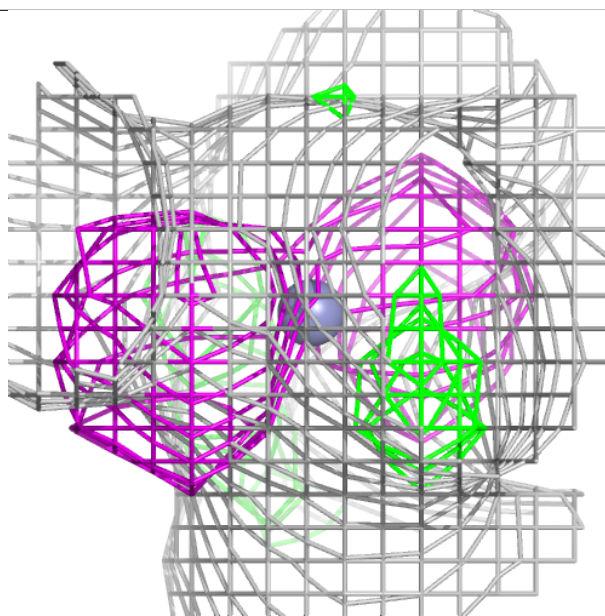
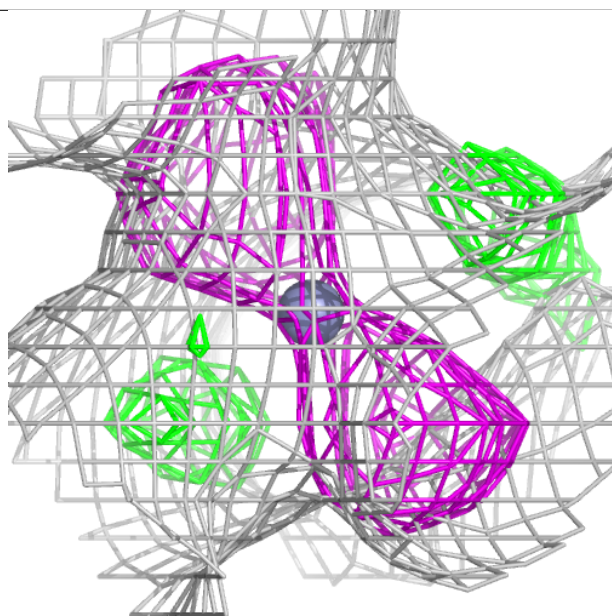
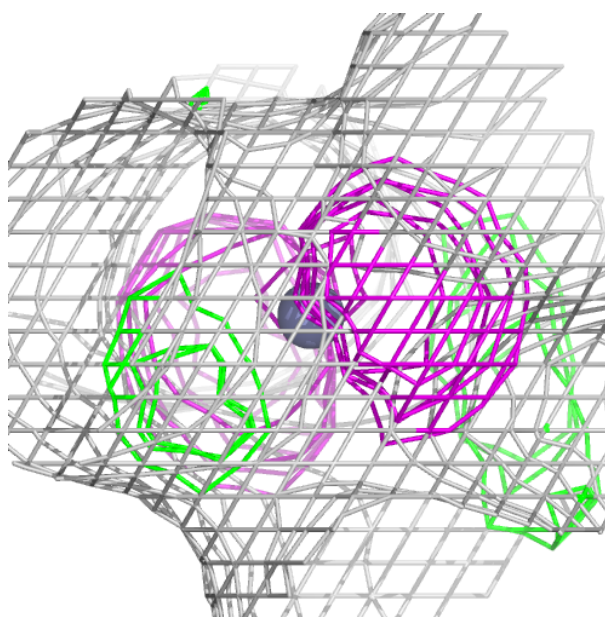
Electron density around ZN B 4402:

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



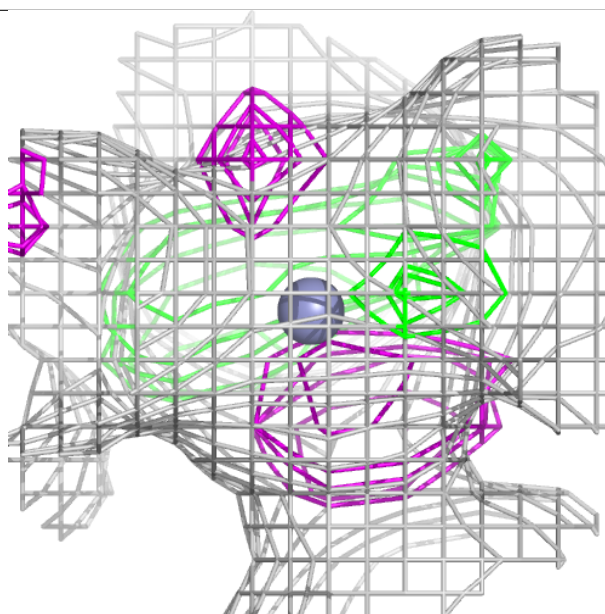
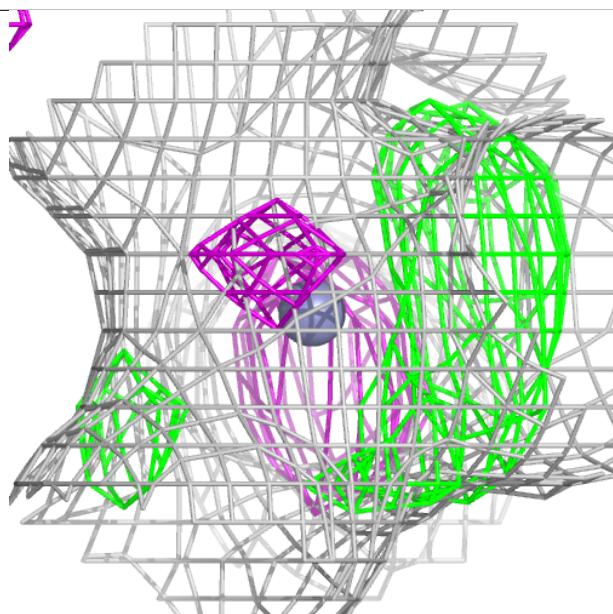
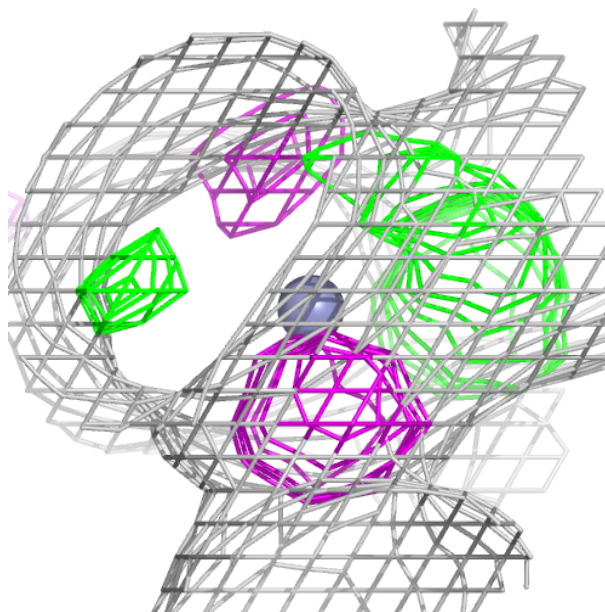
Electron density around ZN D 4402:

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



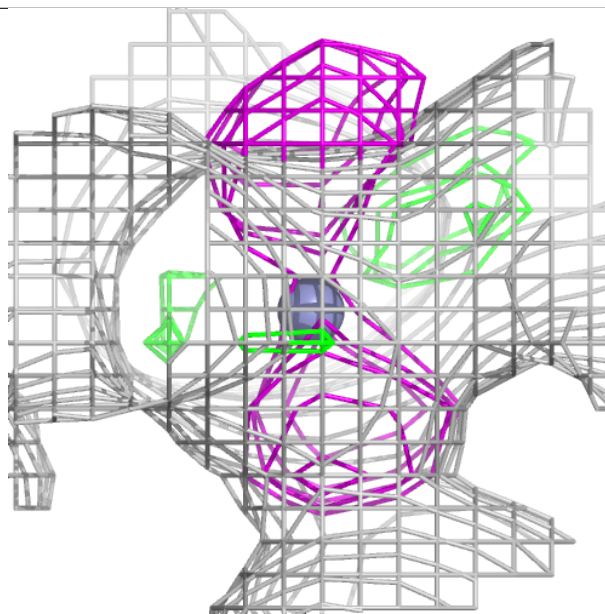
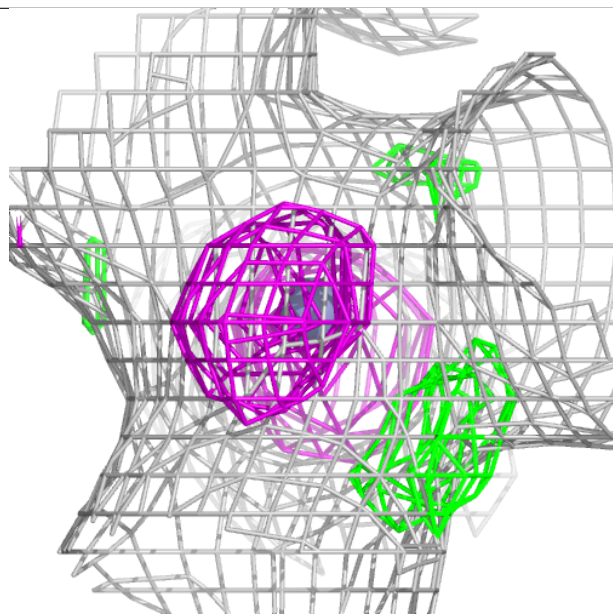
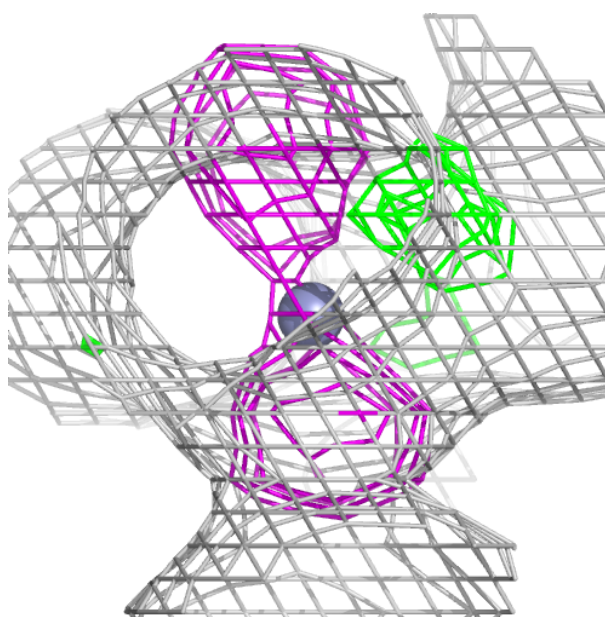
Electron density around ZN B 4401:

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



Electron density around ZN D 4401:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.