



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 31, 2024 – 12:30 pm GMT

PDB ID : 9GL0
Title : Crystal Structure of Acetylpolyamine aminohydrolase (ApaH) from Legionella pneumophila
Authors : Graf, L.G.; Schulze, S.; Palm, G.J.; Lammers, M.
Deposited on : 2024-08-26
Resolution : 2.70 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

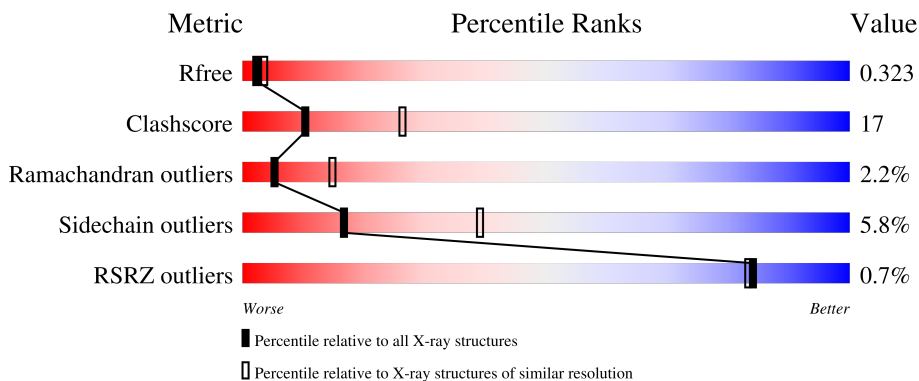
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

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	437	 57% 31% • 10%
1	B	437	 60% 28% • 10%
1	C	437	 54% 32% • 10%
1	D	437	 49% 37% •• 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12712 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 Acetylpolyamine aminohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	393	3182	2037	538	590	17	0	1	0
1	B	392	3175	2032	537	589	17	0	1	0
1	C	392	3164	2027	535	585	17	0	0	0
1	D	391	3167	2029	535	585	18	0	1	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP A0A2S6EWW0
A	-9	ALA	-	expression tag	UNP A0A2S6EWW0
A	-8	HIS	-	expression tag	UNP A0A2S6EWW0
A	-7	HIS	-	expression tag	UNP A0A2S6EWW0
A	-6	HIS	-	expression tag	UNP A0A2S6EWW0
A	-5	HIS	-	expression tag	UNP A0A2S6EWW0
A	-4	HIS	-	expression tag	UNP A0A2S6EWW0
A	-3	HIS	-	expression tag	UNP A0A2S6EWW0
A	-2	VAL	-	expression tag	UNP A0A2S6EWW0
A	-1	GLY	-	expression tag	UNP A0A2S6EWW0
A	0	THR	-	expression tag	UNP A0A2S6EWW0
B	-10	MET	-	initiating methionine	UNP A0A2S6EWW0
B	-9	ALA	-	expression tag	UNP A0A2S6EWW0
B	-8	HIS	-	expression tag	UNP A0A2S6EWW0
B	-7	HIS	-	expression tag	UNP A0A2S6EWW0
B	-6	HIS	-	expression tag	UNP A0A2S6EWW0
B	-5	HIS	-	expression tag	UNP A0A2S6EWW0
B	-4	HIS	-	expression tag	UNP A0A2S6EWW0
B	-3	HIS	-	expression tag	UNP A0A2S6EWW0
B	-2	VAL	-	expression tag	UNP A0A2S6EWW0
B	-1	GLY	-	expression tag	UNP A0A2S6EWW0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	THR	-	expression tag	UNP A0A2S6EWW0
C	-10	MET	-	initiating methionine	UNP A0A2S6EWW0
C	-9	ALA	-	expression tag	UNP A0A2S6EWW0
C	-8	HIS	-	expression tag	UNP A0A2S6EWW0
C	-7	HIS	-	expression tag	UNP A0A2S6EWW0
C	-6	HIS	-	expression tag	UNP A0A2S6EWW0
C	-5	HIS	-	expression tag	UNP A0A2S6EWW0
C	-4	HIS	-	expression tag	UNP A0A2S6EWW0
C	-3	HIS	-	expression tag	UNP A0A2S6EWW0
C	-2	VAL	-	expression tag	UNP A0A2S6EWW0
C	-1	GLY	-	expression tag	UNP A0A2S6EWW0
C	0	THR	-	expression tag	UNP A0A2S6EWW0
D	-10	MET	-	initiating methionine	UNP A0A2S6EWW0
D	-9	ALA	-	expression tag	UNP A0A2S6EWW0
D	-8	HIS	-	expression tag	UNP A0A2S6EWW0
D	-7	HIS	-	expression tag	UNP A0A2S6EWW0
D	-6	HIS	-	expression tag	UNP A0A2S6EWW0
D	-5	HIS	-	expression tag	UNP A0A2S6EWW0
D	-4	HIS	-	expression tag	UNP A0A2S6EWW0
D	-3	HIS	-	expression tag	UNP A0A2S6EWW0
D	-2	VAL	-	expression tag	UNP A0A2S6EWW0
D	-1	GLY	-	expression tag	UNP A0A2S6EWW0
D	0	THR	-	expression tag	UNP A0A2S6EWW0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total K 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total K 1 1	0	0

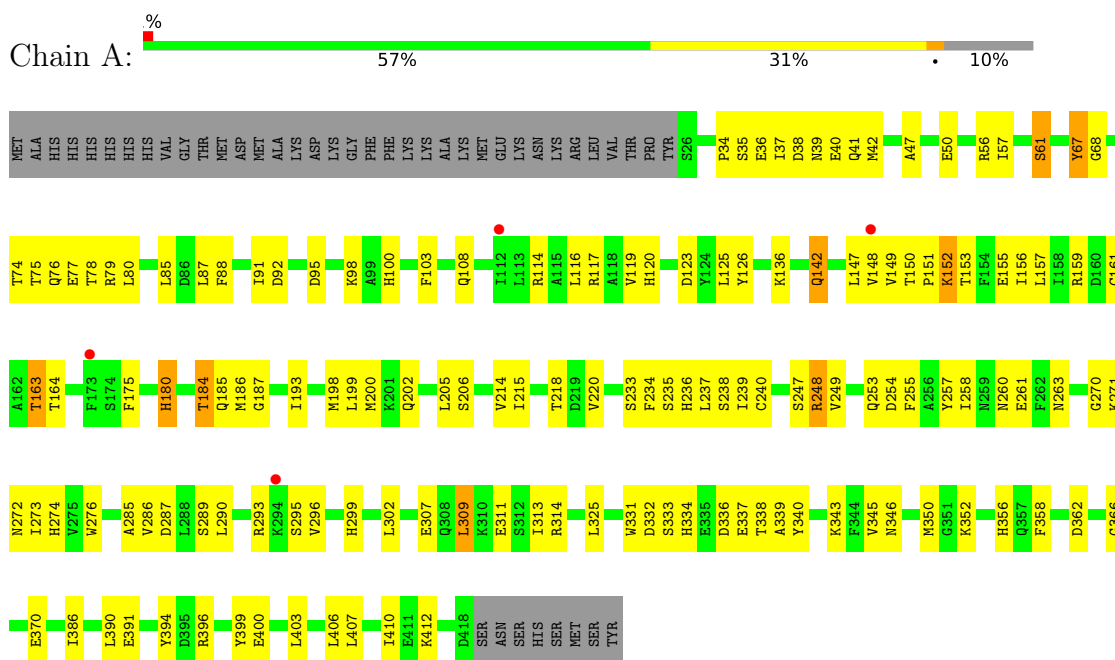
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	B	4	Total O 4 4	0	0
4	C	4	Total O 4 4	0	0
4	D	6	Total O 6 6	0	0

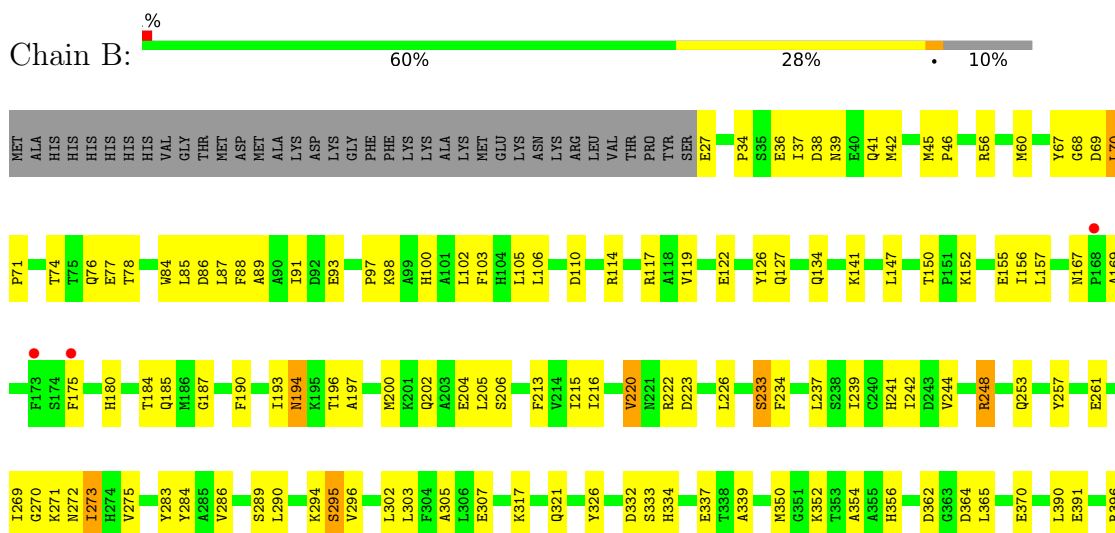
3 Residue-property plots [i](#)

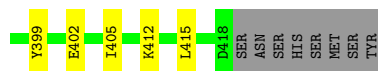
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Acetylpolyamine aminohydrolase



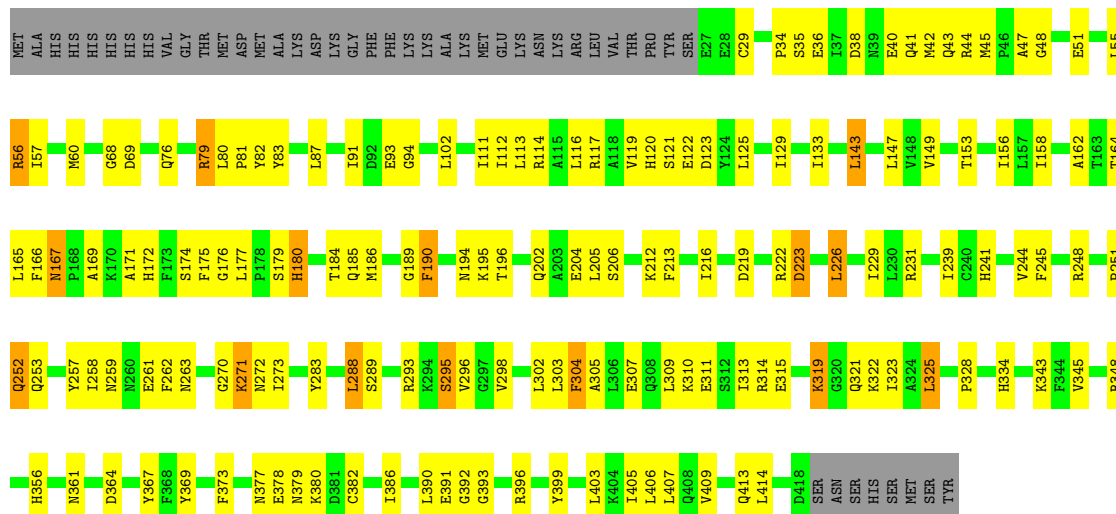
- Molecule 1: Acetylpolyamine aminohydrolase





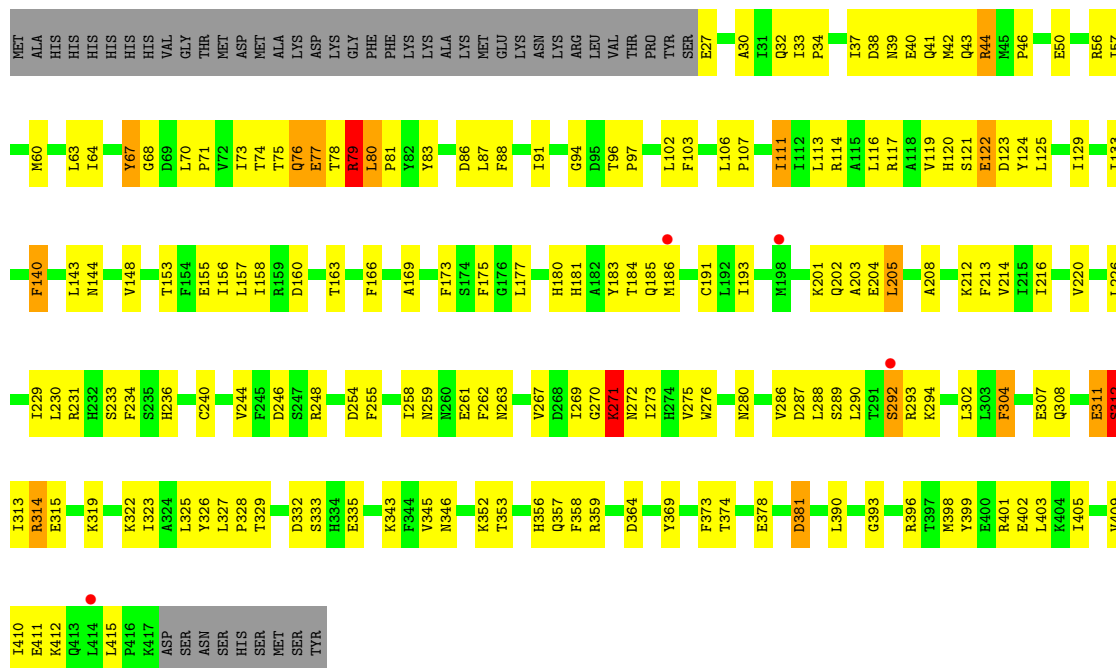
- Molecule 1: Acetylpolyamine aminohydrolase

Chain C: 54% 32% 10%



- Molecule 1: Acetylpolyamine aminohydrolase

Chain D: 49% 37% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	109.12Å 109.12Å 400.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.96 – 2.70 45.96 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.6 (45.96-2.70) 94.5 (45.96-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.87 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.240 , 0.316 0.251 , 0.323	Depositor DCC
R_{free} test set	3452 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	92.2	Xtrriage
Anisotropy	0.512	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 79.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12712	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.67 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2778e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, K

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	0.45	0/3261	0.65	1/4413 (0.0%)
1	B	0.47	1/3254 (0.0%)	0.65	1/4406 (0.0%)
1	C	0.47	0/3243	0.67	1/4390 (0.0%)
1	D	0.45	0/3246	0.63	0/4393
All	All	0.46	1/13004 (0.0%)	0.65	3/17602 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	69	ASP	CA-CB	6.89	1.69	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	220	VAL	CG1-CB-CG2	5.77	120.14	110.90
1	C	325	LEU	CA-CB-CG	5.69	128.39	115.30
1	A	309	LEU	CA-CB-CG	5.68	128.37	115.30

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	3182	0	3107	104	0
1	B	3175	0	3096	78	0
1	C	3164	0	3088	114	0
1	D	3167	0	3094	138	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	6	0	0	0	0
All	All	12712	0	12385	422	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 17.

The worst 5 of 422 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:MET:HE1	1:B:390:LEU:HD11	1.28	1.12
1:A:41:GLN:OE1	1:A:152:LYS:HE3	1.57	1.02
1:C:213:PHE:HB2	1:C:239:ILE:HG12	1.53	0.90
1:C:114:ARG:HH11	1:C:205:LEU:HD23	1.34	0.90
1:A:41:GLN:HB2	1:A:156:ILE:HD11	1.54	0.89

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	392/437 (90%)	341 (87%)	43 (11%)	8 (2%)	6	16
1	B	391/437 (90%)	347 (89%)	38 (10%)	6 (2%)	8	22
1	C	390/437 (89%)	341 (87%)	42 (11%)	7 (2%)	7	18
1	D	390/437 (89%)	346 (89%)	30 (8%)	14 (4%)	3	6
All	All	1563/1748 (89%)	1375 (88%)	153 (10%)	35 (2%)	5	15

5 of 35 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	68	GLY
1	B	295	SER
1	C	252	GLN
1	C	271	LYS
1	D	122	GLU

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	343/381 (90%)	324 (94%)	19 (6%)	18	41
1	B	342/381 (90%)	323 (94%)	19 (6%)	17	41
1	C	340/381 (89%)	318 (94%)	22 (6%)	14	34
1	D	341/381 (90%)	321 (94%)	20 (6%)	16	38
All	All	1366/1524 (90%)	1286 (94%)	80 (6%)	17	38

5 of 80 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	319	LYS
1	D	271	LYS
1	C	348	ARG
1	D	111	ILE
1	D	312	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	167	ASN
1	C	180	HIS
1	D	43	GLN
1	C	379	ASN
1	C	413	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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

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(Å ²)	Q<0.9
1	A	393/437 (89%)	-0.35	4 (1%) 79 79	57, 108, 131, 164	1 (0%)
1	B	392/437 (89%)	-0.37	3 (0%) 82 82	61, 111, 134, 159	1 (0%)
1	C	392/437 (89%)	-0.35	0 100 100	90, 113, 138, 156	0
1	D	391/437 (89%)	-0.36	4 (1%) 79 79	59, 115, 143, 158	1 (0%)
All	All	1568/1748 (89%)	-0.36	11 (0%) 84 83	57, 112, 138, 164	3 (0%)

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	186[A]	MET	5.8
1	A	294[A]	LYS	4.0
1	B	173	PHE	3.1
1	D	414	LEU	3.0
1	A	112	ILE	2.5

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 monosaccharides in this entry.

6.4 Ligands [i](#)

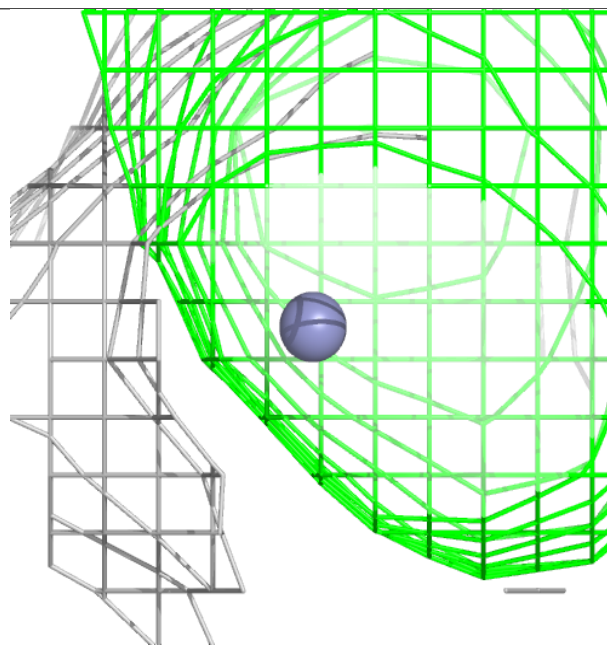
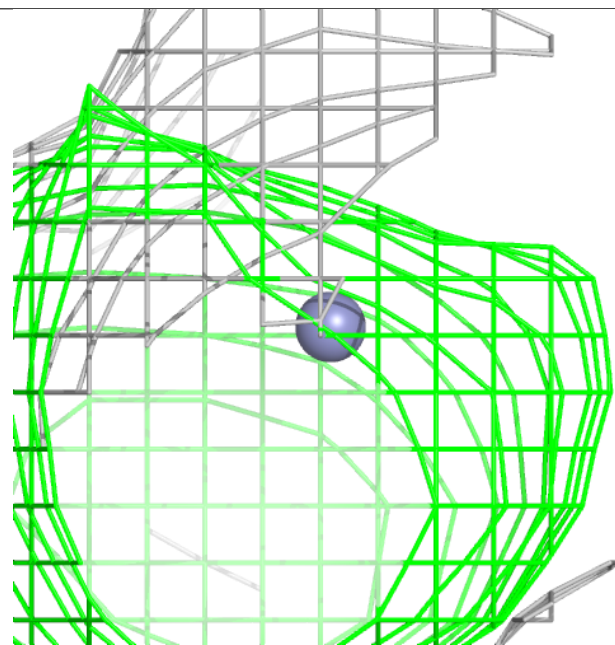
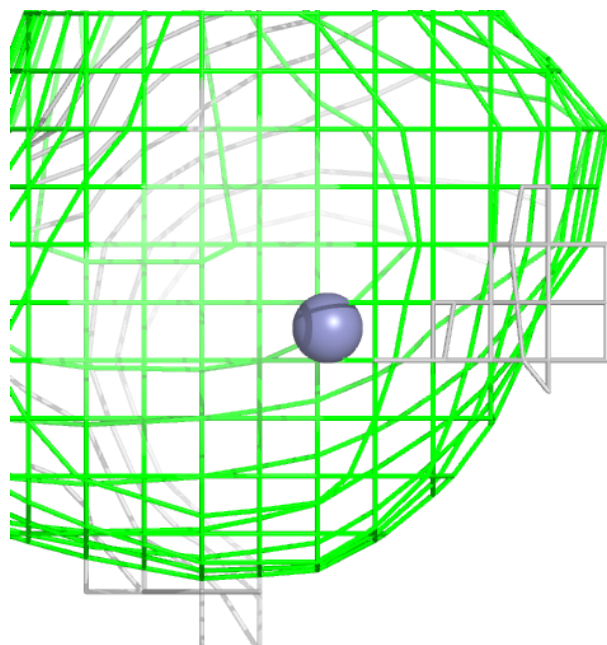
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	B	501	1/1	0.64	0.20	101,101,101,101	1
2	ZN	A	501	1/1	0.66	0.18	103,103,103,103	1
2	ZN	D	501	1/1	0.76	0.20	111,111,111,111	1
3	K	C	502	1/1	0.90	0.06	113,113,113,113	0
3	K	A	502	1/1	0.96	0.04	100,100,100,100	0
2	ZN	C	501	1/1	0.99	0.09	113,113,113,113	1

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.

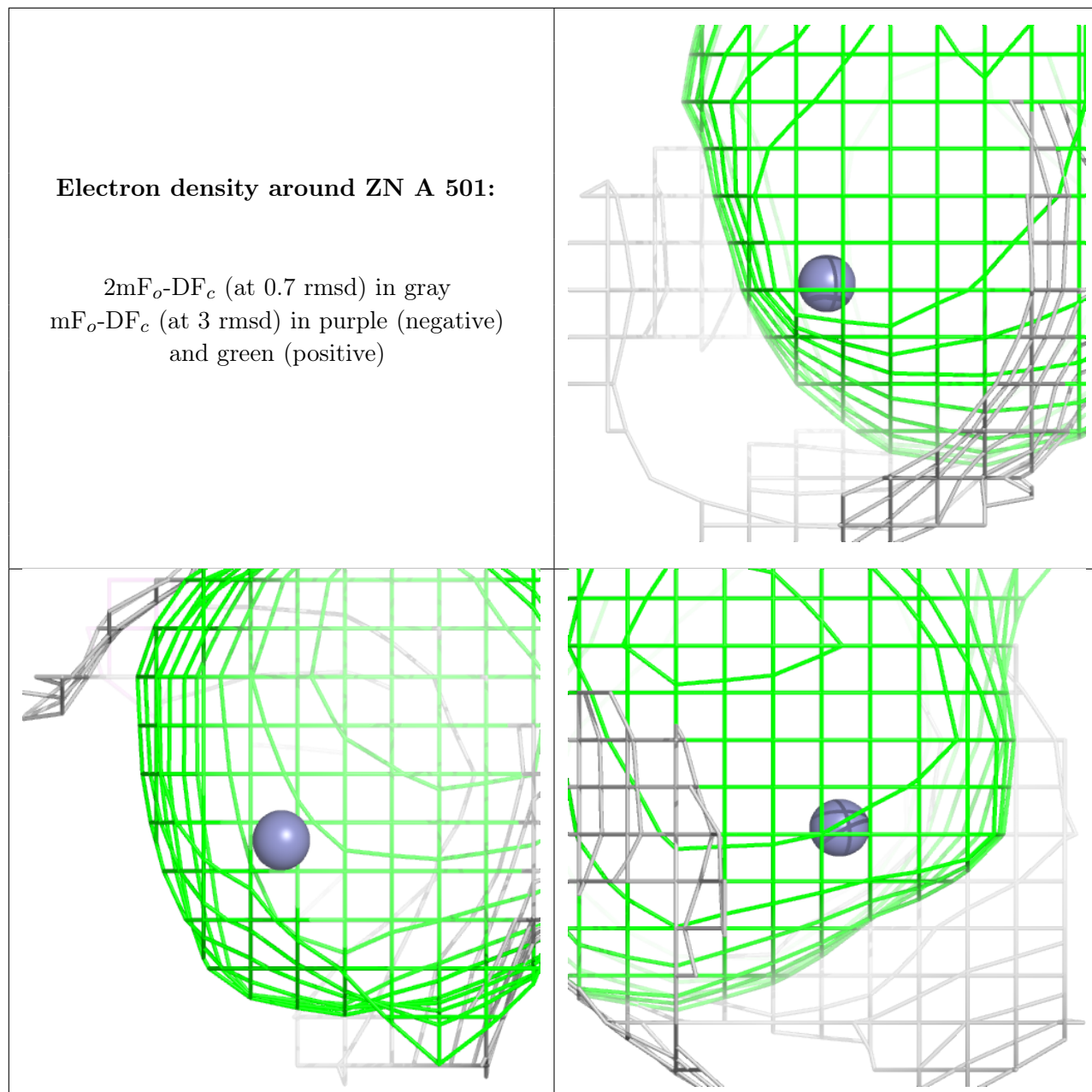
Electron density around ZN B 501:

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



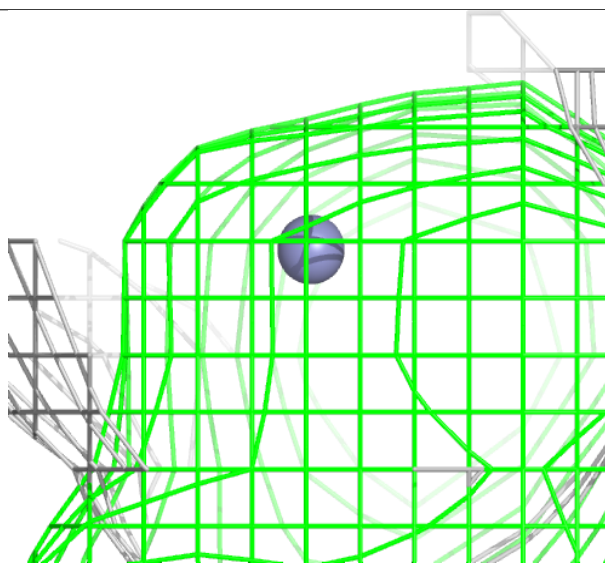
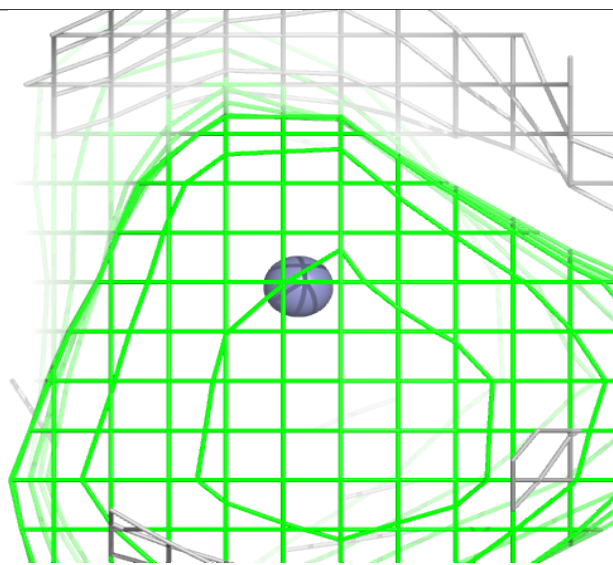
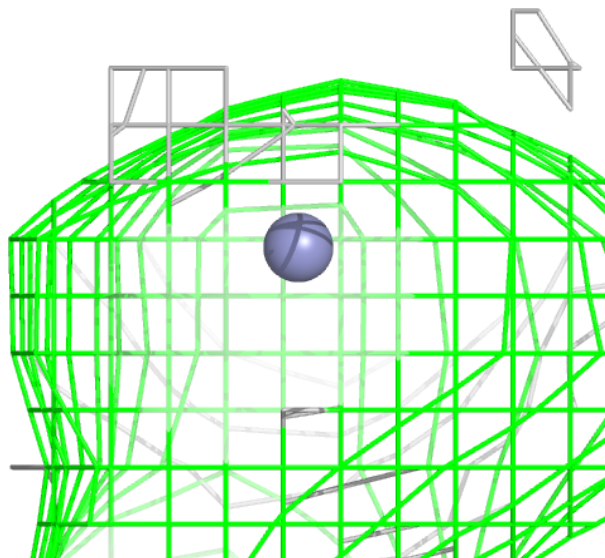
Electron density around ZN A 501:

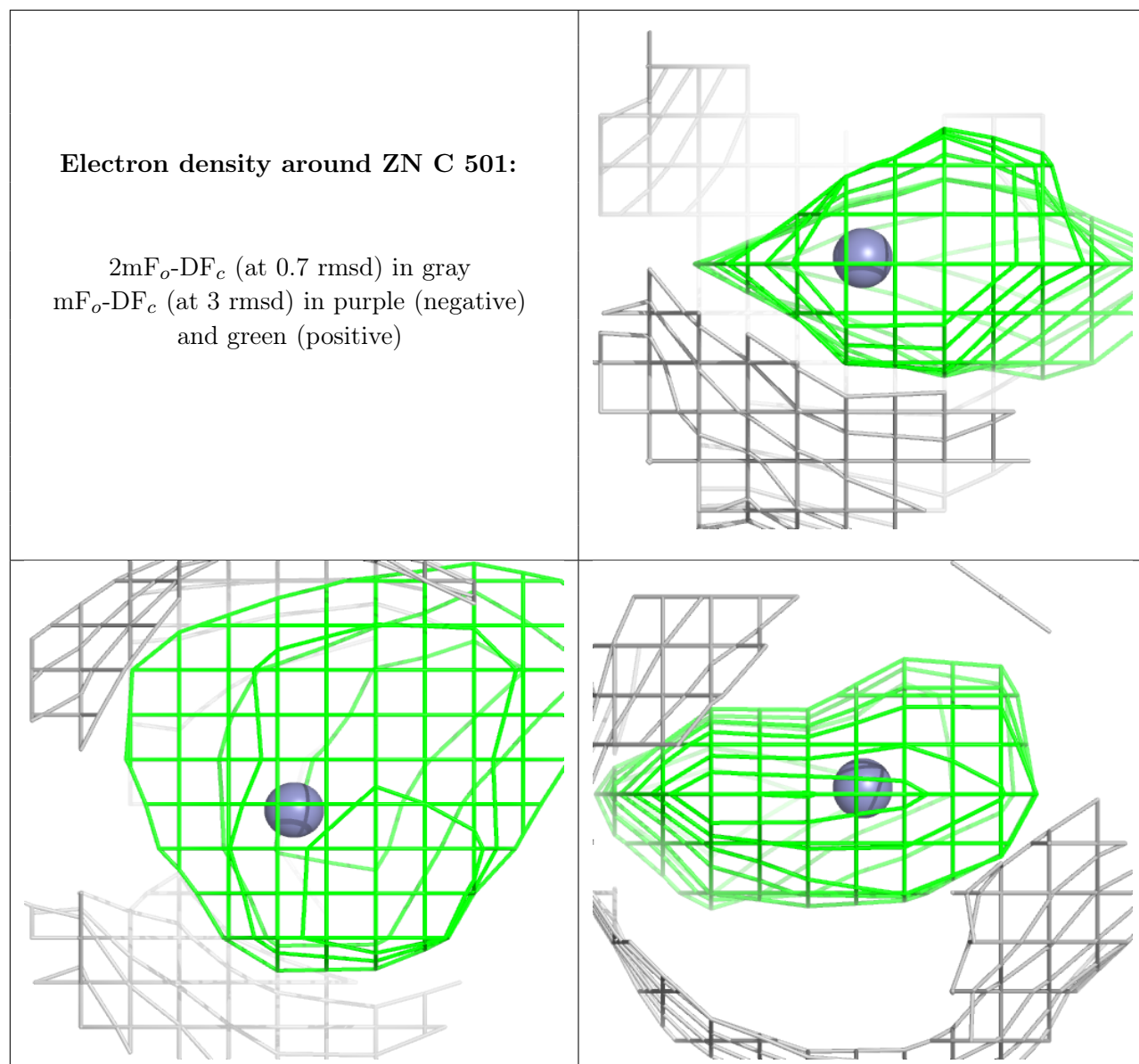
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN D 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.