



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:16 AM UTC

PDB ID : 7Q39 / pdb_00007q39
Title : Ribonucleotide Reductase R2_genomic protein from Aquifex aeolicus
Authors : Scaletti, E.; Rehling, D.; Stenmark, P.
Deposited on : 2021-10-27
Resolution : 2.10 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

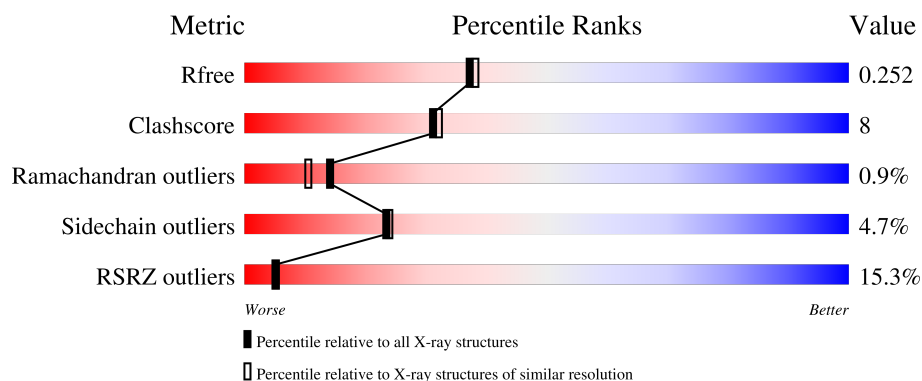
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	AAA	350	14% 73% 16% • 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2758 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 Ribonucleoside-diphosphate reductase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	320	Total	C	N	O	S	0	3	0
			2687	1746	430	504	7			

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	2	Total	Fe	0	0
			2	2		

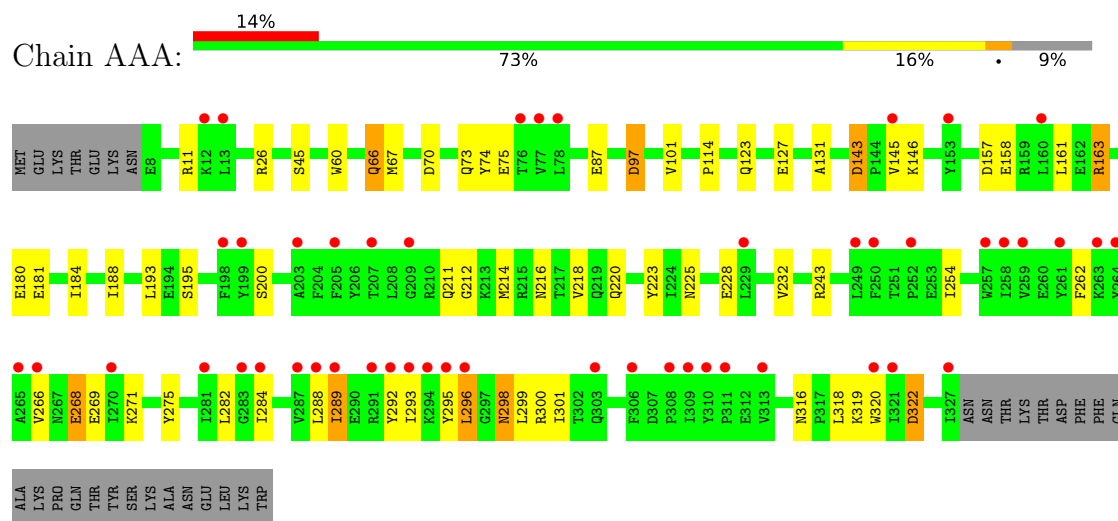
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	69	Total	O	0	0
			69	69		

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: Ribonucleoside-diphosphate reductase subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	69.25Å 69.25Å 177.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.25 – 2.10 47.25 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.25-2.10) 100.0 (47.25-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.238 , 0.297 (Not available) , 0.252	Depositor DCC
R_{free} test set	1284 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2758	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: FE

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	AAA	0.98	0/2753	1.48	3/3736 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	97	ASP	CB-CA-C	5.24	120.31	110.63
1	AAA	296	LEU	CA-C-N	5.11	125.77	119.99
1	AAA	296	LEU	C-N-CA	5.11	125.77	119.99

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	AAA	2687	0	2557	43	0
2	AAA	2	0	0	0	0
3	AAA	69	0	0	7	0
All	All	2758	0	2557	43	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 8.

All (43) 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:AAA:200:SER:HB2	1:AAA:269:GLU:OE2	1.89	0.72
1:AAA:200:SER:CB	1:AAA:269:GLU:OE2	2.37	0.72
1:AAA:228[B]:GLU:OE2	1:AAA:228[B]:GLU:HA	1.90	0.71
1:AAA:163:ARG:NH2	1:AAA:269:GLU:OE1	2.26	0.67
1:AAA:97:ASP:HB3	3:AAA:542:HOH:O	1.95	0.67
1:AAA:266:VAL:HG13	1:AAA:293:ILE:HG22	1.77	0.66
1:AAA:143:ASP:HB3	3:AAA:562:HOH:O	1.96	0.65
1:AAA:284:ILE:N	3:AAA:503:HOH:O	2.30	0.64
1:AAA:211:GLN:NE2	1:AAA:282:LEU:O	2.32	0.62
1:AAA:268:GLU:OE2	1:AAA:268:GLU:O	2.18	0.61
1:AAA:214:MET:O	1:AAA:218:VAL:HG23	2.02	0.59
1:AAA:195:SER:HB3	1:AAA:300:ARG:HG3	1.83	0.59
1:AAA:200:SER:OG	1:AAA:269:GLU:OE2	2.22	0.57
1:AAA:66:GLN:O	3:AAA:501:HOH:O	2.17	0.57
1:AAA:228[B]:GLU:O	1:AAA:232:VAL:HG23	2.03	0.57
1:AAA:193:LEU:C	1:AAA:193:LEU:HD23	2.30	0.57
1:AAA:131:ALA:HB2	3:AAA:542:HOH:O	2.05	0.56
1:AAA:322:ASP:OD1	1:AAA:322:ASP:N	2.33	0.54
1:AAA:298:ASN:O	1:AAA:301:ILE:HG22	2.08	0.53
1:AAA:75:GLU:HA	3:AAA:527:HOH:O	2.09	0.52
1:AAA:181:GLU:HG3	1:AAA:254:ILE:HD11	1.91	0.52
1:AAA:181:GLU:CG	1:AAA:254:ILE:HD11	2.41	0.50
1:AAA:180:GLU:O	1:AAA:184:ILE:HG12	2.11	0.50
1:AAA:296:LEU:O	1:AAA:299:LEU:HB3	2.11	0.50
1:AAA:67:MET:HE3	1:AAA:220:GLN:HE21	1.77	0.49
1:AAA:292:TYR:CD1	1:AAA:318:LEU:HD12	2.49	0.46
1:AAA:318:LEU:O	1:AAA:320:TRP:N	2.48	0.46
1:AAA:157:ASP:OD2	1:AAA:275:TYR:OH	2.29	0.46
1:AAA:74:TYR:HB2	1:AAA:214:MET:HE1	1.97	0.46
1:AAA:295:TYR:CE2	1:AAA:316:ASN:HB2	2.51	0.45
1:AAA:101[B]:VAL:HG22	1:AAA:127:GLU:HB2	1.99	0.45
1:AAA:161:LEU:HD12	1:AAA:161:LEU:O	2.17	0.45
1:AAA:87:GLU:OE1	1:AAA:146:LYS:NZ	2.49	0.44
1:AAA:288:LEU:O	1:AAA:289:ILE:C	2.60	0.44
1:AAA:268:GLU:HG3	3:AAA:523:HOH:O	2.17	0.44
1:AAA:60:TRP:CE2	1:AAA:223:TYR:HB3	2.53	0.43
1:AAA:26:ARG:HD3	1:AAA:114:PRO:HD2	2.01	0.43
1:AAA:73:GLN:NE2	1:AAA:212:GLY:O	2.45	0.43
1:AAA:228[A]:GLU:O	1:AAA:232:VAL:HG23	2.20	0.42
1:AAA:70:ASP:CG	1:AAA:216:ASN:HB2	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:184:ILE:O	1:AAA:188:ILE:HG12	2.20	0.41
1:AAA:318:LEU:C	1:AAA:320:TRP:H	2.29	0.41
1:AAA:123:GLN:O	1:AAA:127:GLU:HG2	2.20	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	AAA	321/350 (92%)	284 (88%)	34 (11%)	3 (1%)	14 10

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	289	ILE
1	AAA	319	LYS
1	AAA	298	ASN

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	AAA	282/323 (87%)	269 (95%)	13 (5%)	24 25

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	11	ARG
1	AAA	45	SER
1	AAA	66	GLN
1	AAA	143	ASP
1	AAA	145	VAL
1	AAA	158	GLU
1	AAA	163	ARG
1	AAA	225	ASN
1	AAA	243	ARG
1	AAA	262	PHE
1	AAA	268	GLU
1	AAA	271	LYS
1	AAA	322	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 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 ⓘ

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	AAA	320/350 (91%)	1.00	49 (15%) 5 5	18, 65, 96, 113	3 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	270	ILE	3.5
1	AAA	160	LEU	3.4
1	AAA	292	TYR	3.2
1	AAA	281	ILE	3.1
1	AAA	327	ILE	3.1
1	AAA	229	LEU	3.0
1	AAA	284	ILE	3.0
1	AAA	313	VAL	3.0
1	AAA	289	ILE	2.9
1	AAA	265	ALA	2.9
1	AAA	145	VAL	2.8
1	AAA	266	VAL	2.8
1	AAA	309	ILE	2.8
1	AAA	308	PRO	2.8
1	AAA	293	ILE	2.8
1	AAA	288	LEU	2.8
1	AAA	199	TYR	2.7
1	AAA	259	VAL	2.7
1	AAA	321	ILE	2.7
1	AAA	198	PHE	2.6
1	AAA	311	PRO	2.5
1	AAA	320	TRP	2.5
1	AAA	258	ILE	2.4
1	AAA	263	LYS	2.4
1	AAA	291	ARG	2.4
1	AAA	283	GLY	2.4
1	AAA	261	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	AAA	295	TYR	2.3
1	AAA	203	ALA	2.3
1	AAA	12	LYS	2.3
1	AAA	287	VAL	2.3
1	AAA	249	LEU	2.3
1	AAA	264	TYR	2.3
1	AAA	296	LEU	2.3
1	AAA	13	LEU	2.3
1	AAA	209	GLY	2.2
1	AAA	310	TYR	2.2
1	AAA	303	GLN	2.2
1	AAA	306	PHE	2.2
1	AAA	77	VAL	2.2
1	AAA	76	THR	2.2
1	AAA	252	PRO	2.1
1	AAA	257	TRP	2.1
1	AAA	205	PHE	2.1
1	AAA	250	PHE	2.1
1	AAA	207	THR	2.1
1	AAA	153	TYR	2.1
1	AAA	78	LEU	2.0
1	AAA	294	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(Å ²)	Q<0.9
2	FE	AAA	401	1/1	0.95	0.13	96,96,96,96	0

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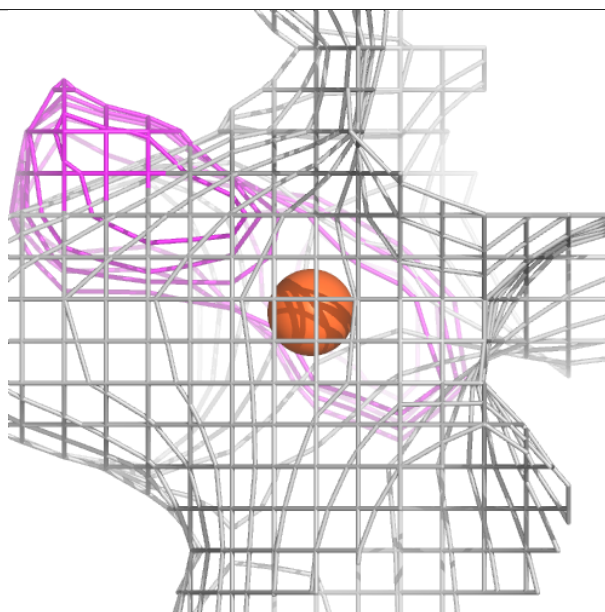
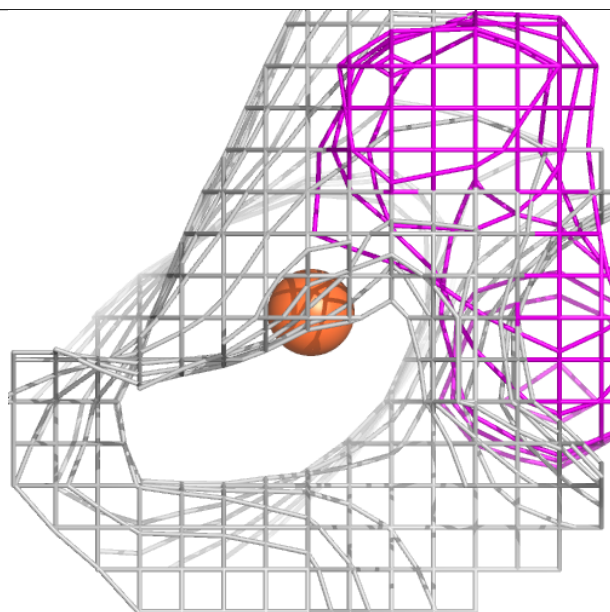
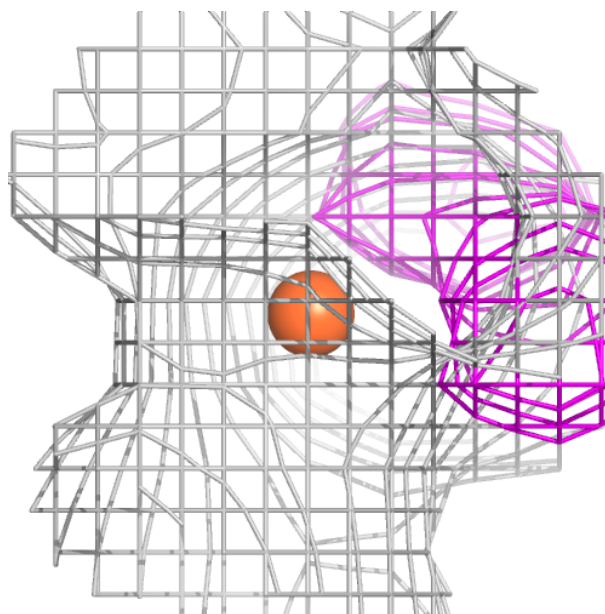
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FE	AAA	402	1/1	0.99	0.03	54,54,54,54	0

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.

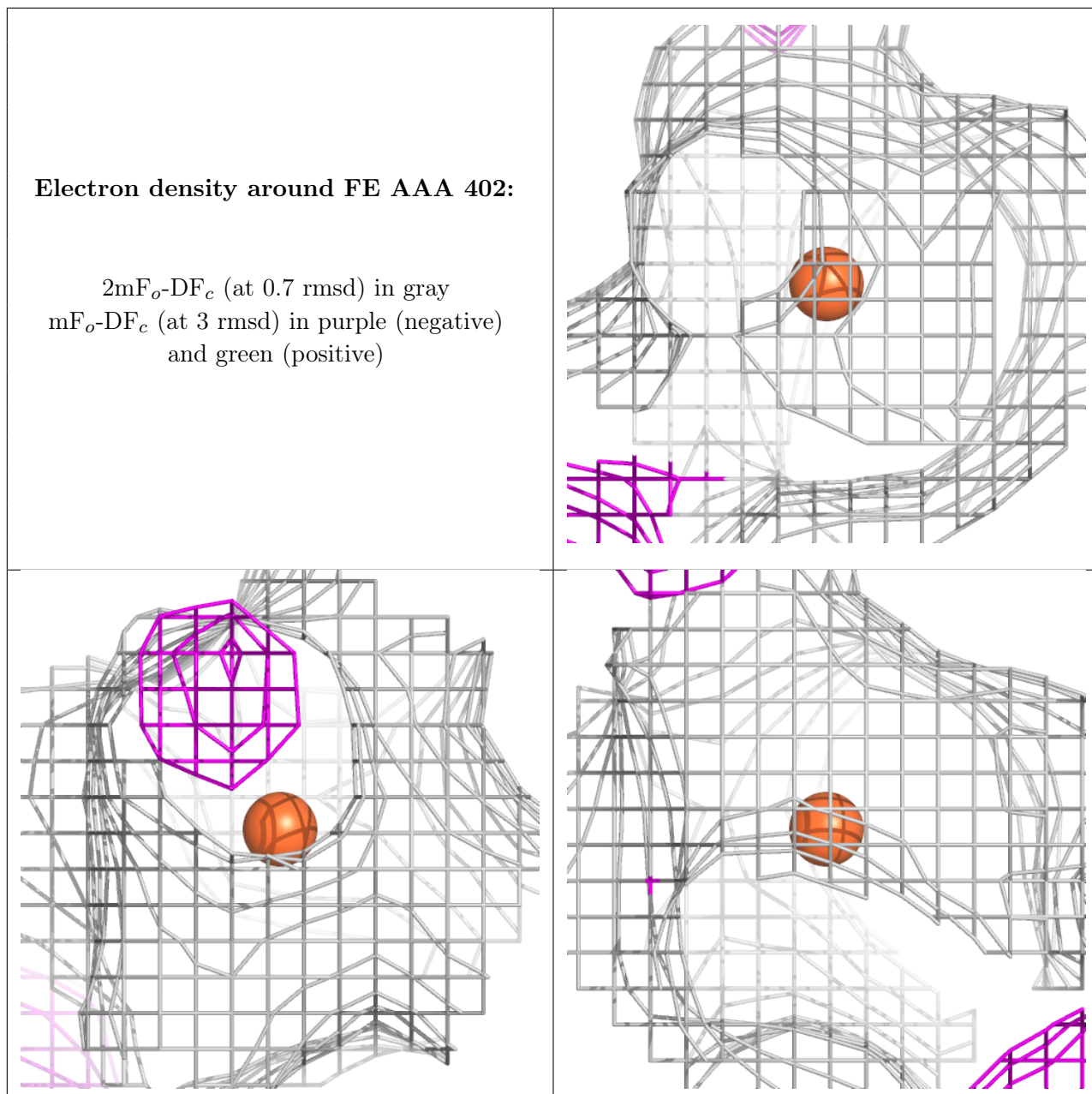
Electron density around FE AAA 401:

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 FE AAA 402:

$2mF_o - DF_c$ (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.