



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:56 PM UTC

PDB ID : 6K1A / pdb_00006k1a
Title : Crystal structure of EXD2 exonuclease domain soaked in Mn and Mg
Authors : Park, J.; Lee, C.
Deposited on : 2019-05-10
Resolution : 2.60 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

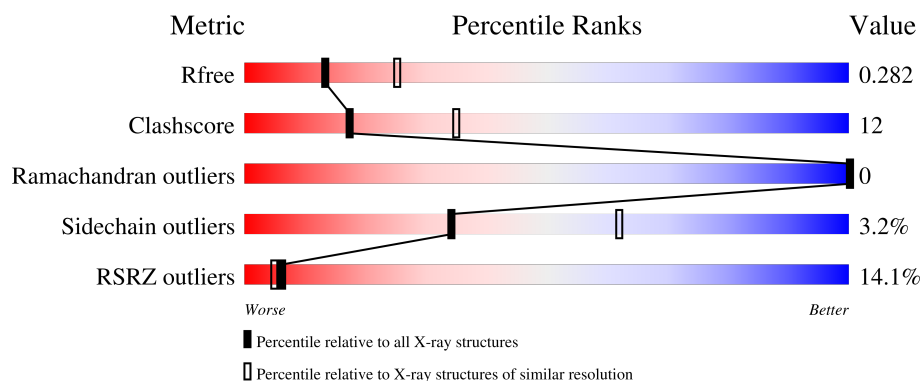
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	220	12% 66% 21% • 11%
1	B	220	13% 69% 19% • 11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3129 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 Exonuclease 3'-5' domain-containing protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1515	968	256	281	10			
1	B	196	Total	C	N	O	S	0	0	0
			1522	972	257	283	10			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

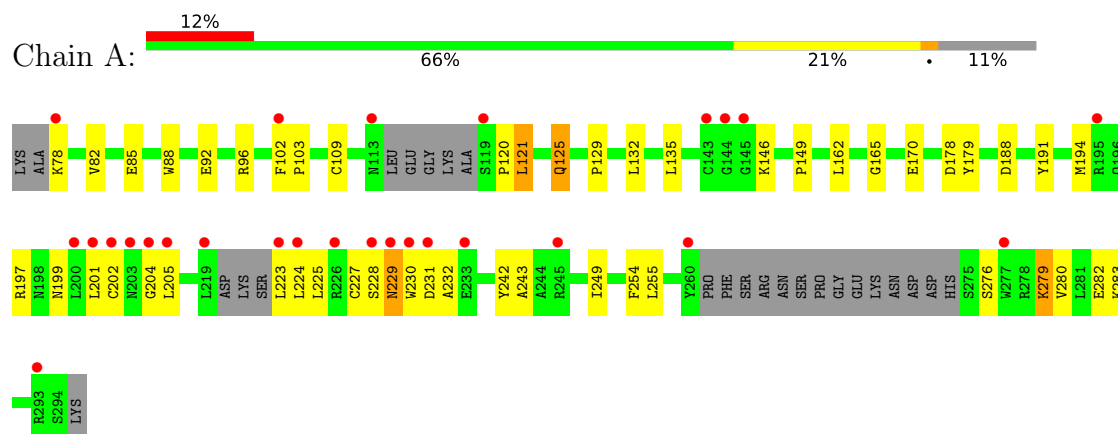
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total	O	0	0
			42	42		
4	B	46	Total	O	0	0
			46	46		

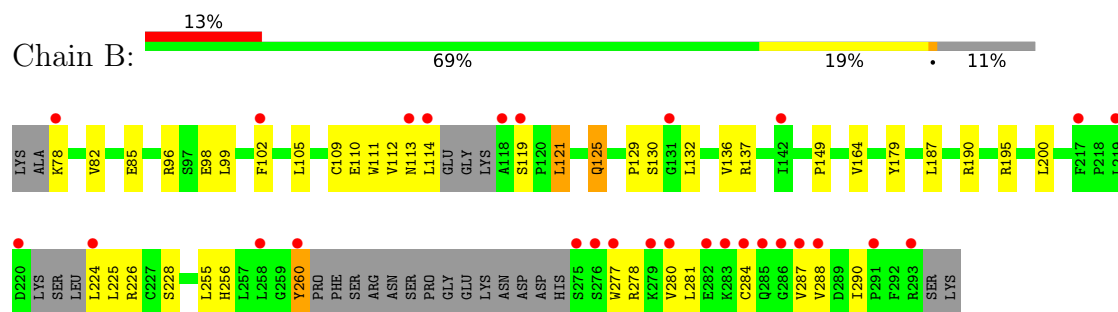
3 Residue-property plots [i](#)

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: Exonuclease 3'-5' domain-containing protein 2



- Molecule 1: Exonuclease 3'-5' domain-containing protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.32Å 73.66Å 131.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.22 – 2.60 39.22 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (39.22-2.60) 98.9 (39.22-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.223 , 0.260 0.264 , 0.282	Depositor DCC
R_{free} test set	714 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3129	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.50% 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: MN, MG

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	0.30	0/1536	0.81	3/2082 (0.1%)
1	B	0.32	0/1543	0.78	1/2092 (0.0%)
All	All	0.31	0/3079	0.80	4/4174 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	229	ASN	N-CA-C	8.81	121.83	109.15
1	A	229	ASN	CB-CA-C	-6.55	100.71	110.67
1	B	278	ARG	N-CA-C	-5.57	105.36	111.82
1	A	204	GLY	N-CA-C	-5.42	103.62	112.83

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	1515	0	1564	53	0
1	B	1522	0	1568	40	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
4	A	42	0	0	1	0
4	B	46	0	0	1	0
All	All	3129	0	3132	77	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 12.

All (77) 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:112:VAL:HG12	1:B:114:LEU:HG	1.56	0.86
1:A:78:LYS:HG3	1:A:132:LEU:HA	1.60	0.83
1:A:78:LYS:HD2	1:A:132:LEU:CA	2.13	0.77
1:B:102:PHE:HB3	1:B:129:PRO:HG2	1.67	0.76
1:A:78:LYS:CG	1:A:132:LEU:HA	2.19	0.72
1:B:288:VAL:O	1:B:290:ILE:HG12	1.91	0.71
1:A:230:TRP:CD1	1:B:110:GLU:HB3	2.26	0.70
1:A:109:CYS:HB3	1:A:121:LEU:HD21	1.74	0.70
1:B:256:HIS:HA	1:B:260:TYR:HE2	1.57	0.69
1:A:78:LYS:CD	1:A:132:LEU:HA	2.22	0.69
1:B:260:TYR:HD2	1:B:260:TYR:N	1.90	0.69
1:B:224:LEU:HG	1:B:224:LEU:O	1.93	0.68
1:B:109:CYS:HB3	1:B:121:LEU:HD21	1.76	0.67
1:B:112:VAL:CG1	1:B:114:LEU:HG	2.24	0.67
1:B:260:TYR:N	1:B:260:TYR:CD2	2.63	0.67
1:A:232:ALA:O	1:B:137:ARG:NH2	2.23	0.66
1:A:194:MET:SD	1:A:197:ARG:NH2	2.70	0.64
1:A:162:LEU:HD11	1:B:284:CYS:SG	2.38	0.64
1:A:255:LEU:HD11	1:B:129:PRO:HB3	1.78	0.64
1:A:85:GLU:HG3	1:A:149:PRO:HG3	1.80	0.63
1:A:201:LEU:HD23	1:A:205:LEU:HD21	1.81	0.62
1:A:78:LYS:HD2	1:A:132:LEU:HB2	1.82	0.61
1:A:125:GLN:HG3	1:A:135:LEU:HD22	1.81	0.60
1:A:78:LYS:HD2	1:A:132:LEU:HA	1.80	0.60
1:A:223:LEU:O	1:A:224:LEU:HB2	2.02	0.60
1:A:191:TYR:CD2	1:B:287:VAL:HG13	2.37	0.59
1:A:229:ASN:OD1	1:A:231:ASP:N	2.25	0.59
1:A:229:ASN:ND2	1:A:232:ALA:HB2	2.18	0.59
1:A:197:ARG:HB2	1:B:200:LEU:HD13	1.86	0.57
1:A:125:GLN:CG	1:A:135:LEU:HD22	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:LYS:HA	1:A:282:GLU:HB3	1.86	0.56
1:A:78:LYS:HD2	1:A:132:LEU:CB	2.35	0.56
1:A:92:GLU:OE2	1:A:96:ARG:HD3	2.06	0.56
1:A:225:LEU:O	1:A:228:SER:OG	2.22	0.53
1:B:164:VAL:HG12	1:B:187:LEU:HB3	1.91	0.53
1:A:162:LEU:CD1	1:B:284:CYS:SG	2.97	0.52
1:B:284:CYS:O	1:B:288:VAL:HG23	2.09	0.52
1:B:256:HIS:HA	1:B:260:TYR:CE2	2.42	0.52
1:A:229:ASN:HD21	1:A:232:ALA:N	2.08	0.52
1:B:121:LEU:O	1:B:179:TYR:OH	2.27	0.51
1:A:78:LYS:HD2	1:A:132:LEU:N	2.25	0.51
1:B:277:TRP:CZ3	1:B:280:VAL:HG11	2.48	0.49
1:A:230:TRP:NE1	1:B:110:GLU:HB3	2.26	0.49
1:A:229:ASN:OD1	1:A:229:ASN:C	2.54	0.48
1:B:85:GLU:HG3	1:B:149:PRO:HG3	1.96	0.48
1:B:78:LYS:HE3	1:B:132:LEU:HD13	1.97	0.47
1:A:102:PHE:HA	1:A:103:PRO:HD3	1.70	0.46
1:A:205:LEU:HB3	1:B:190:ARG:HG2	1.97	0.46
1:A:276:SER:O	1:A:280:VAL:HG23	2.16	0.45
1:A:103:PRO:HD2	1:B:277:TRP:CE2	2.52	0.45
1:B:82:VAL:HG22	1:B:136:VAL:HA	1.99	0.44
1:A:225:LEU:HD12	1:A:242:TYR:HB2	1.98	0.44
1:A:229:ASN:ND2	1:A:232:ALA:N	2.66	0.44
1:A:243:ALA:HB1	1:B:125:GLN:CD	2.43	0.44
1:B:226:ARG:NH2	4:B:408:HOH:O	2.46	0.44
1:A:121:LEU:O	1:A:179:TYR:OH	2.35	0.43
1:B:110:GLU:HG2	1:B:125:GLN:OE1	2.18	0.43
1:A:78:LYS:CD	1:A:132:LEU:HB2	2.49	0.43
1:A:254:PHE:CD1	1:B:164:VAL:HG11	2.54	0.43
1:A:223:LEU:HD12	4:A:434:HOH:O	2.17	0.43
1:A:225:LEU:HD23	1:A:225:LEU:HA	1.83	0.42
1:A:82:VAL:HG11	1:A:88:TRP:HB2	2.01	0.42
1:A:243:ALA:HB1	1:B:125:GLN:NE2	2.34	0.42
1:A:283:LYS:O	1:B:195:ARG:NH1	2.53	0.42
1:A:249:ILE:HD13	1:A:249:ILE:HA	1.87	0.42
1:A:165:GLY:O	1:A:188:ASP:HA	2.19	0.42
1:A:191:TYR:CG	1:B:287:VAL:HG13	2.54	0.42
1:A:199:ASN:HA	1:A:202:CYS:HB3	2.02	0.42
1:B:112:VAL:N	1:B:119:SER:OG	2.53	0.41
1:B:98:GLU:OE1	1:B:130:SER:OG	2.38	0.41
1:B:225:LEU:O	1:B:228:SER:OG	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:LEU:HD21	1:B:105:LEU:HG	2.02	0.41
1:A:129:PRO:HB3	1:B:255:LEU:HD11	2.02	0.41
1:B:111:TRP:HB2	1:B:119:SER:OG	2.21	0.41
1:A:120:PRO:HA	1:A:178:ASP:OD2	2.21	0.40
1:A:78:LYS:CD	1:A:132:LEU:CA	2.84	0.40
1:A:224:LEU:O	1:A:227:CYS:HB2	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	A	187/220 (85%)	186 (100%)	1 (0%)	0	100	100
1	B	188/220 (86%)	187 (100%)	1 (0%)	0	100	100
All	All	375/440 (85%)	373 (100%)	2 (0%)	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	172/193 (89%)	167 (97%)	5 (3%)	37	65
1	B	172/193 (89%)	166 (96%)	6 (4%)	32	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	344/386 (89%)	333 (97%)	11 (3%)	34 62

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	LEU
1	A	125	GLN
1	A	146	LYS
1	A	170	GLU
1	A	279	LYS
1	B	96	ARG
1	B	113	ASN
1	B	121	LEU
1	B	125	GLN
1	B	260	TYR
1	B	281	LEU

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

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 4 ligands modelled in this entry, 4 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	A	195/220 (88%)	0.88	27 (13%) 6 5	28, 48, 95, 154	0
1	B	196/220 (89%)	0.95	28 (14%) 6 5	30, 48, 89, 114	0
All	All	391/440 (88%)	0.92	55 (14%) 6 5	28, 48, 95, 154	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	284	CYS	10.8
1	A	202	CYS	6.4
1	A	223	LEU	6.1
1	B	114	LEU	5.5
1	B	285	GLN	5.4
1	A	260	TYR	5.3
1	B	102	PHE	5.1
1	A	205	LEU	5.0
1	B	282	GLU	4.9
1	B	119	SER	4.7
1	B	286	GLY	4.7
1	B	260	TYR	4.6
1	A	78	LYS	4.3
1	B	279	LYS	4.3
1	B	283	LYS	4.2
1	A	219	LEU	4.2
1	B	219	LEU	4.0
1	B	287	VAL	4.0
1	B	118	ALA	3.8
1	A	228	SER	3.7
1	A	204	GLY	3.6
1	B	113	ASN	3.6
1	B	275	SER	3.4
1	A	102	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	280	VAL	3.4
1	A	195	ARG	3.2
1	A	229	ASN	3.2
1	A	119	SER	3.2
1	A	201	LEU	2.9
1	A	113	ASN	2.8
1	A	224	LEU	2.8
1	A	143	CYS	2.8
1	B	277	TRP	2.7
1	B	142	ILE	2.7
1	A	203	ASN	2.7
1	A	231	ASP	2.6
1	A	144	GLY	2.6
1	B	220	ASP	2.5
1	B	293	ARG	2.4
1	B	291	PRO	2.4
1	B	276	SER	2.3
1	A	200	LEU	2.3
1	A	145	GLY	2.3
1	B	224	LEU	2.3
1	A	277	TRP	2.3
1	A	230	TRP	2.2
1	B	217	PHE	2.2
1	A	245	ARG	2.2
1	A	293	ARG	2.2
1	B	258	LEU	2.1
1	B	78	LYS	2.0
1	B	131	GLY	2.0
1	A	233	GLU	2.0
1	A	226	ARG	2.0
1	B	288	VAL	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

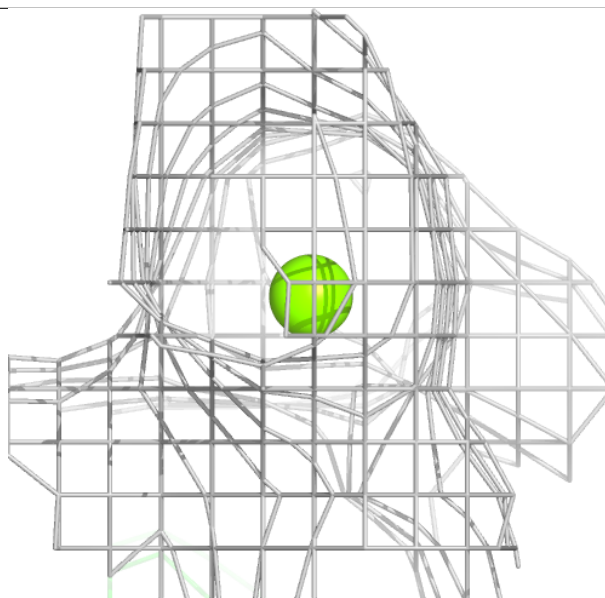
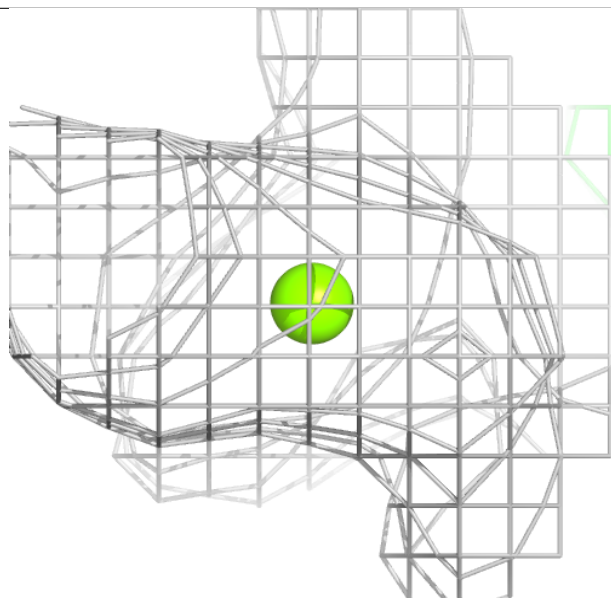
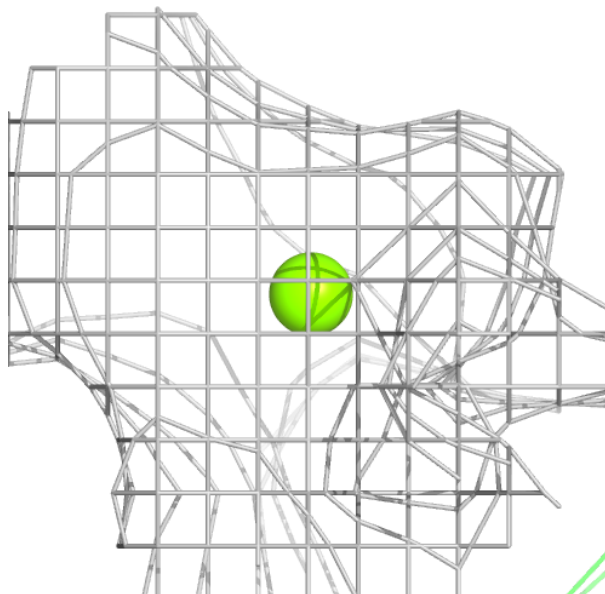
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	B	302	1/1	0.56	0.20	79,79,79,79	0
3	MG	A	302	1/1	0.75	0.17	66,66,66,66	0
2	MN	B	301	1/1	0.91	0.11	52,52,52,52	0
2	MN	A	301	1/1	0.99	0.09	59,59,59,59	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.

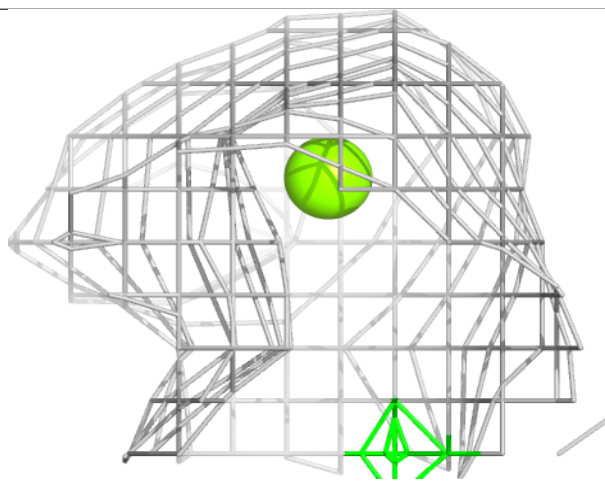
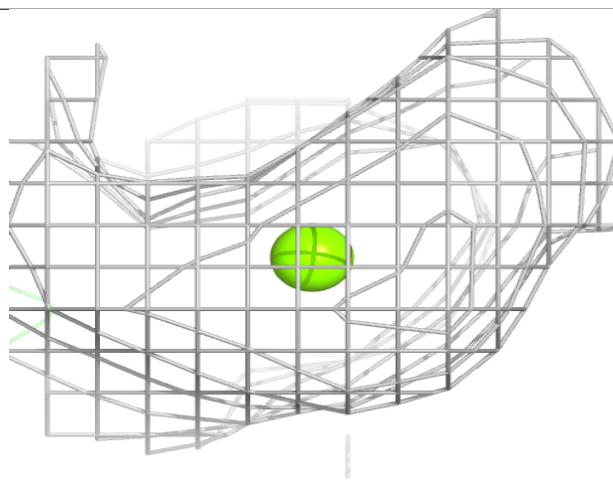
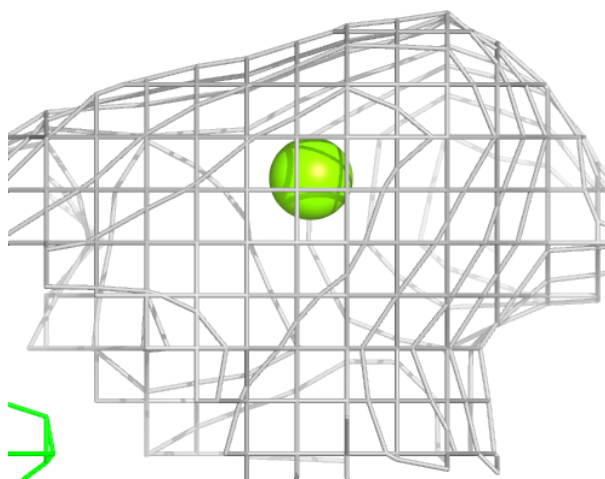
Electron density around MG B 302:

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



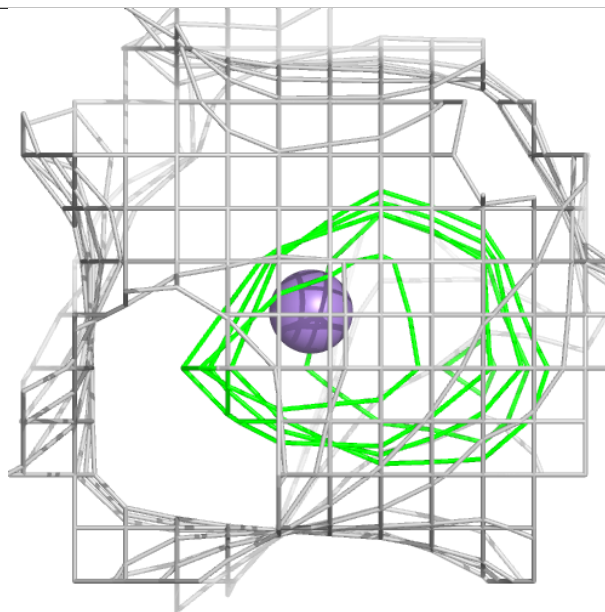
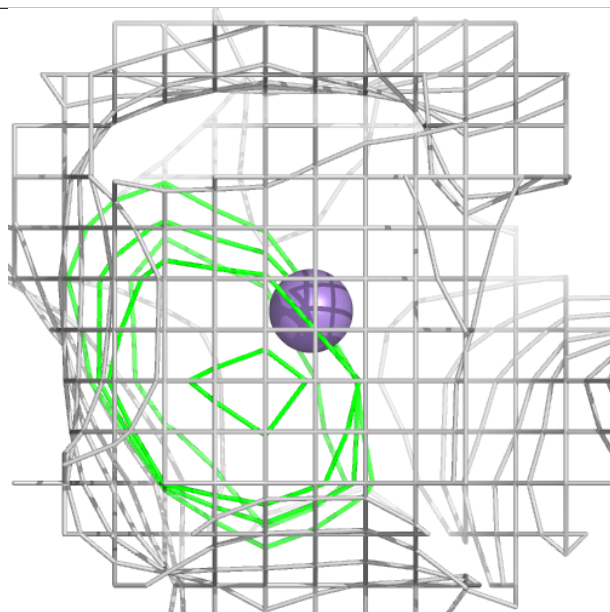
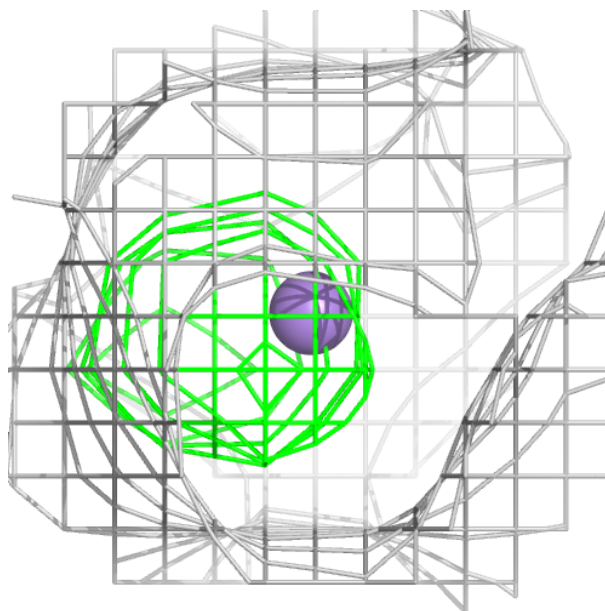
Electron density around MG A 302:

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



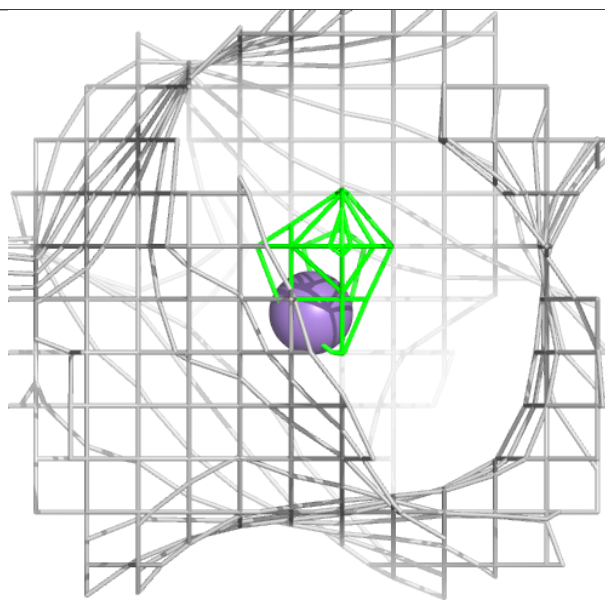
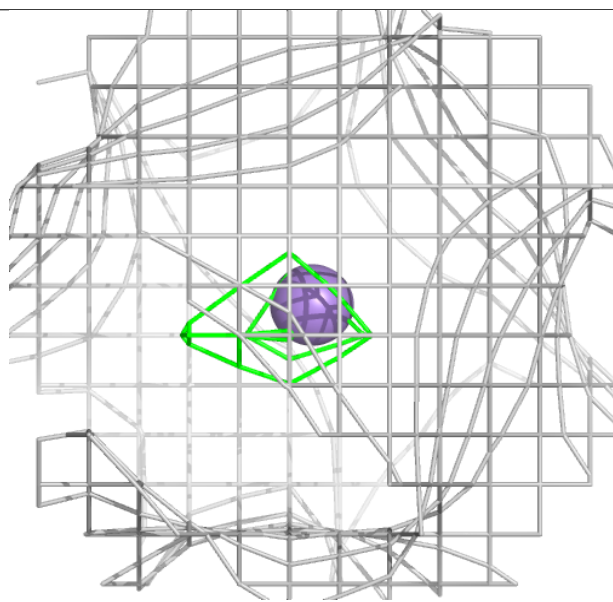
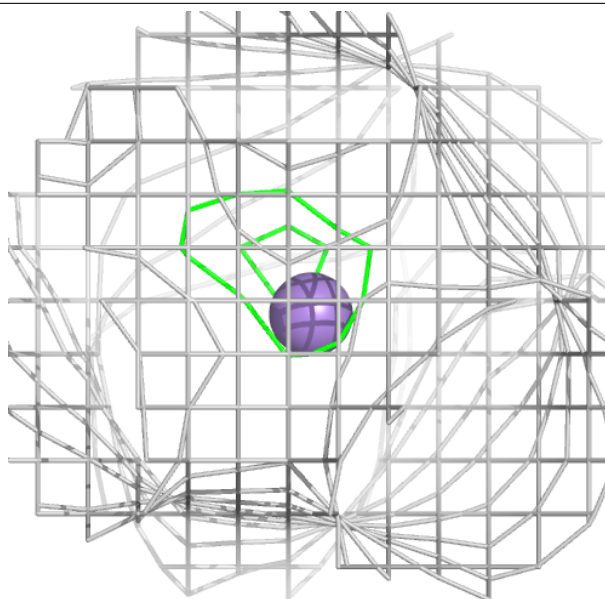
Electron density around MN B 301:

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



Electron density around MN A 301:

$2mF_o-DF_c$ (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.