



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2024 – 04:50 pm BST

PDB ID : 8ROL
Title : Human cohesin SMC3-HD(EQ)/RAD21-N complex - ATP-Mg-bound conformation - Form 2
Authors : Vitoria Gomes, M.; Romier, C.
Deposited on : 2024-01-11
Resolution : 3.11 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.002 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.38.2

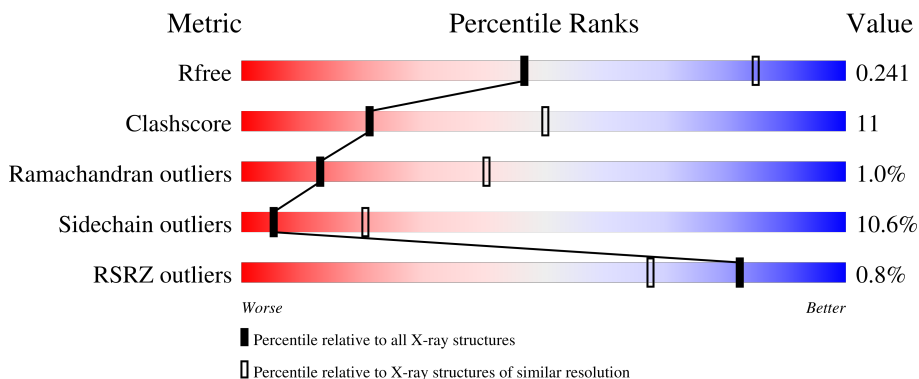
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.11 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	1668 (3.14-3.10)
Clashscore	180529	1788 (3.14-3.10)
Ramachandran outliers	177936	1696 (3.14-3.10)
Sidechain outliers	177891	1696 (3.14-3.10)
RSRZ outliers	164620	1668 (3.14-3.10)

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	A	462	 3% 58% 28% 11%
2	B	110	 3% 49% 20% 29%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3987 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 Structural maintenance of chromosomes protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	411	3333	2119	576	624	14	0	0	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	967	GLY	-	linker	UNP Q9UQE7
A	968	SER	-	linker	UNP Q9UQE7
A	969	GLY	-	linker	UNP Q9UQE7
A	970	SER	-	linker	UNP Q9UQE7
A	971	LEU	-	linker	UNP Q9UQE7
A	972	VAL	-	linker	UNP Q9UQE7
A	973	PRO	-	linker	UNP Q9UQE7
A	974	ARG	-	linker	UNP Q9UQE7
A	975	GLY	-	linker	UNP Q9UQE7
A	976	SER	-	linker	UNP Q9UQE7
A	977	GLY	-	linker	UNP Q9UQE7
A	978	SER	-	linker	UNP Q9UQE7
A	1144	GLN	GLU	engineered mutation	UNP Q9UQE7

- Molecule 2 is a protein called Double-strand-break repair protein rad21 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	78	622	408	109	101	4	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	103	GLY	-	expression tag	UNP O60216
B	104	SER	-	expression tag	UNP O60216
B	105	LEU	-	expression tag	UNP O60216

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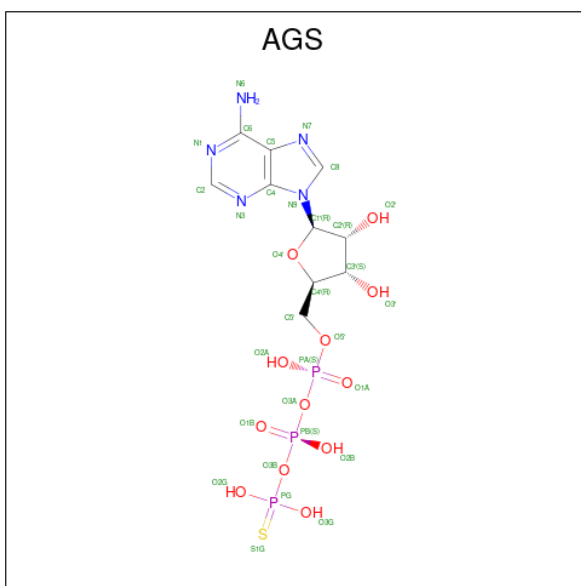
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Chain	Residue	Modelled	Actual	Comment	Reference
B	106	GLU	-	expression tag	UNP O60216
B	107	VAL	-	expression tag	UNP O60216
B	108	LEU	-	expression tag	UNP O60216
B	109	PHE	-	expression tag	UNP O60216
B	110	GLN	-	expression tag	UNP O60216

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

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

- Molecule 4 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (three-letter code: AGS) (formula: C₁₀H₁₆N₅O₁₂P₃S) (labeled as "Ligand of Interest" by depositor).

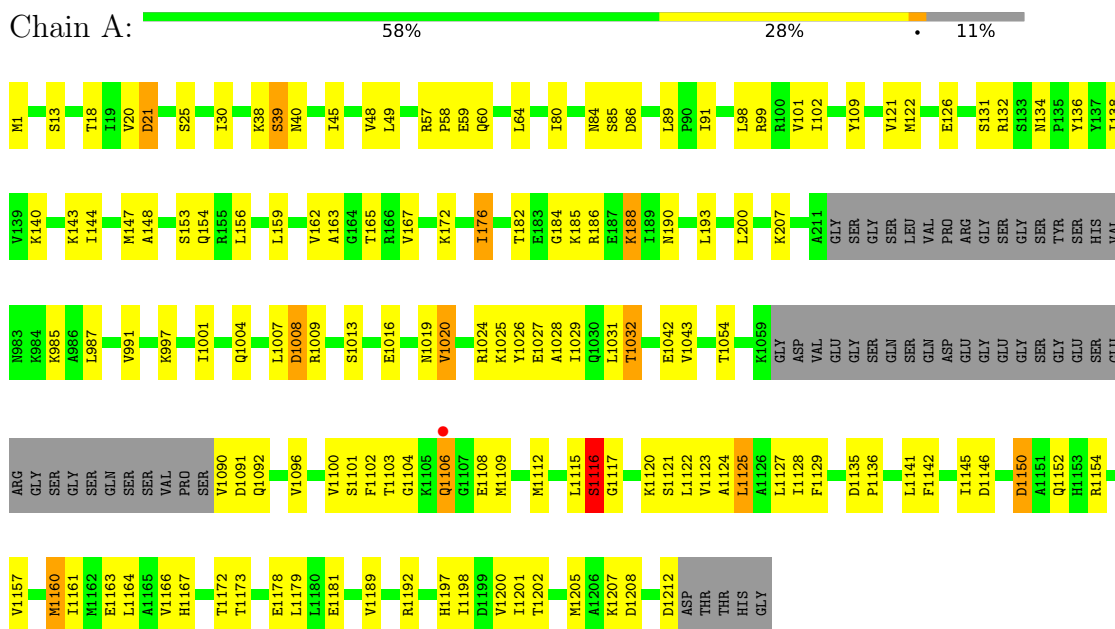


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
4	A	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		

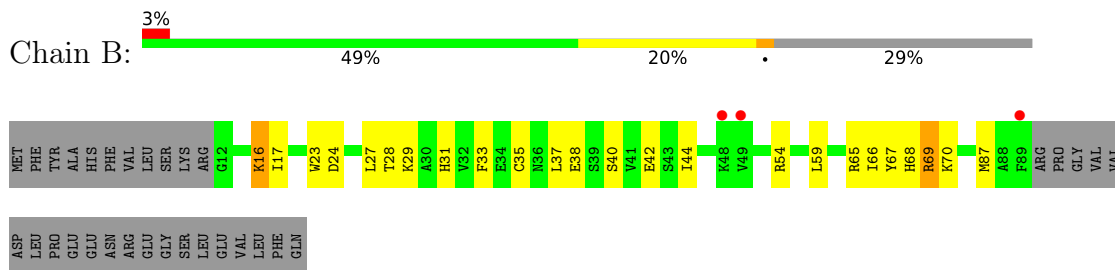
3 Residue-property plots [i](#)

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: Structural maintenance of chromosomes protein 3



- Molecule 2: Double-strand-break repair protein rad21 homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	201.07Å 201.07Å 92.38Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.64 – 3.11 44.64 – 3.11	Depositor EDS
% Data completeness (in resolution range)	99.6 (44.64-3.11) 99.7 (44.64-3.11)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 3.12Å)	Xtrriage
Refinement program	PHENIX 1.17_3644	Depositor
R, R_{free}	0.187 , 0.238 0.192 , 0.241	Depositor DCC
R_{free} test set	1011 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	104.9	Xtrriage
Anisotropy	0.365	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 82.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3987	wwPDB-VP
Average B, all atoms (Å ²)	112.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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: MG, AGS

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	0.42	0/3384	0.57	0/4532
2	B	0.41	0/636	0.51	0/855
All	All	0.42	0/4020	0.56	0/5387

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	3333	0	3388	75	1
2	B	622	0	661	14	0
3	A	1	0	0	0	0
4	A	31	0	11	2	0
All	All	3987	0	4060	87	1

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 11.

All (87) 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)
4:A:2001:AGS:O4'	4:A:2001:AGS:C1'	1.64	1.17
1:A:148:ALA:HB1	1:A:1120:LYS:HB3	1.65	0.77
1:A:1154:ARG:NH1	1:A:1178:GLU:OE1	2.19	0.76
1:A:186:ARG:HH21	1:A:1008:ASP:HA	1.55	0.70
2:B:38:GLU:O	2:B:42:GLU:HG2	1.94	0.67
2:B:16:LYS:NZ	2:B:35:CYS:SG	2.61	0.66
1:A:1142:PHE:HB2	1:A:1172:THR:HG22	1.80	0.64
1:A:39:SER:OG	4:A:2001:AGS:O2B	2.17	0.63
1:A:163:ALA:HA	1:A:1032:THR:HG21	1.79	0.63
2:B:27:LEU:O	2:B:69:ARG:NH2	2.28	0.62
1:A:172:LYS:O	1:A:176:ILE:HG23	2.00	0.61
1:A:1029:ILE:HD12	1:A:1029:ILE:H	1.65	0.61
1:A:38:LYS:HB3	1:A:1173:THR:HG21	1.84	0.60
1:A:1028:ALA:O	1:A:1032:THR:HG23	2.01	0.60
1:A:48:VAL:HG12	1:A:49:LEU:HD23	1.84	0.58
1:A:1116:SER:OG	1:A:1117:GLY:N	2.38	0.57
1:A:207:LYS:HG3	1:A:987:LEU:HD22	1.86	0.56
1:A:186:ARG:NH2	1:A:1008:ASP:OD1	2.39	0.56
2:B:37:LEU:HD11	2:B:70:LYS:HD2	1.87	0.56
1:A:1027:GLU:O	1:A:1031:LEU:HB2	2.05	0.55
1:A:186:ARG:NH2	1:A:1008:ASP:HA	2.22	0.54
1:A:185:LYS:HE3	2:B:68:HIS:CE1	2.42	0.54
2:B:17:ILE:HD11	2:B:40:SER:HB3	1.89	0.54
1:A:156:LEU:HB2	1:A:1096:VAL:HG23	1.90	0.54
1:A:1091:ASP:OD1	1:A:1092:GLN:N	2.35	0.53
1:A:109:TYR:CD2	1:A:121:VAL:HG21	2.45	0.52
1:A:1166:VAL:HG23	1:A:1167:HIS:CE1	2.45	0.52
1:A:182:THR:HG21	2:B:67:TYR:CE2	2.45	0.51
1:A:122:MET:HB3	1:A:132:ARG:NH1	2.26	0.51
1:A:167:VAL:HG21	2:B:54:ARG:HG3	1.92	0.51
1:A:30:ILE:HB	1:A:1173:THR:HG22	1.92	0.50
1:A:49:LEU:HD21	1:A:80:ILE:HD12	1.93	0.50
1:A:98:LEU:HD11	1:A:109:TYR:HD2	1.76	0.50
1:A:1025:LYS:HE2	1:A:1026:TYR:HE1	1.75	0.49
1:A:1007:LEU:O	1:A:1009:ARG:N	2.45	0.49
2:B:28:THR:HG22	2:B:31:HIS:ND1	2.28	0.48
2:B:29:LYS:HZ1	2:B:33:PHE:HE2	1.63	0.47
1:A:1166:VAL:HG23	1:A:1167:HIS:ND1	2.29	0.47
2:B:37:LEU:HD21	2:B:66:ILE:HG22	1.96	0.47
1:A:131:SER:HB2	1:A:134:ASN:HB2	1.97	0.47
1:A:144:ILE:HD13	1:A:1125:LEU:CD1	2.44	0.47
1:A:136:TYR:HE1	1:A:143:LYS:HD3	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1160:MET:O	1:A:1164:LEU:HG	2.14	0.46
1:A:1007:LEU:C	1:A:1009:ARG:H	2.20	0.46
1:A:1125:LEU:HD21	1:A:1145:ILE:HD12	1.98	0.45
1:A:1150:ASP:OD1	1:A:1152:GLN:N	2.49	0.45
1:A:193:LEU:HD13	1:A:1001:ILE:HD13	1.99	0.45
1:A:985:LYS:HA	1:A:985:LYS:HD2	1.77	0.45
1:A:13:SER:H	1:A:40:ASN:ND2	2.15	0.45
1:A:1202:THR:OG1	1:A:1205:MET:HG3	2.16	0.44
1:A:45:ILE:HG21	1:A:1141:LEU:HD11	1.98	0.44
1:A:1124:ALA:O	1:A:1128:ILE:HG13	2.17	0.44
1:A:997:LYS:HB3	1:A:997:LYS:HE2	1.77	0.44
1:A:1142:PHE:O	1:A:1172:THR:HA	2.17	0.44
1:A:147:MET:CE	1:A:1127:LEU:HD23	2.48	0.44
1:A:1101:SER:HB2	1:A:1109:MET:SD	2.57	0.44
1:A:1106:GLN:C	1:A:1108:GLU:H	2.21	0.44
1:A:1207:LYS:HE2	1:A:1207:LYS:HB3	1.89	0.44
1:A:1201:ILE:HB	1:A:1205:MET:HB2	2.00	0.44
1:A:60:GLN:O	1:A:64:LEU:HG	2.17	0.43
1:A:89:LEU:O	1:A:91:ILE:N	2.51	0.43
1:A:144:ILE:HD11	1:A:1145:ILE:HG22	2.00	0.43
1:A:1100:VAL:HB	1:A:1115:LEU:HD11	2.00	0.43
1:A:163:ALA:CA	1:A:1032:THR:HG21	2.45	0.43
1:A:20:VAL:HG12	1:A:21:ASP:O	2.19	0.43
1:A:1135:ASP:N	1:A:1136:PRO:HD3	2.33	0.43
1:A:1166:VAL:HG23	1:A:1167:HIS:CG	2.52	0.43
1:A:57:ARG:HB3	1:A:58:PRO:HD2	2.01	0.43
1:A:1212:ASP:N	1:A:1212:ASP:OD1	2.52	0.43
2:B:44:ILE:HD13	2:B:59:LEU:HB3	2.00	0.42
1:A:1043:VAL:HG11	1:A:1129:PHE:HB2	2.01	0.42
2:B:59:LEU:HD23	2:B:59:LEU:HA	1.89	0.42
1:A:84:ASN:C	1:A:86:ASP:H	2.23	0.42
1:A:1146:ASP:HB2	1:A:1154:ARG:HG2	2.02	0.41
1:A:1157:VAL:O	1:A:1161:ILE:HG13	2.21	0.41
1:A:1146:ASP:OD2	1:A:1179:LEU:HD13	2.19	0.41
1:A:1163:GLU:O	1:A:1166:VAL:HG22	2.20	0.41
1:A:1122:LEU:HD12	1:A:1122:LEU:HA	1.88	0.41
1:A:1192:ARG:HB2	1:A:1197:HIS:CD2	2.56	0.41
1:A:30:ILE:O	1:A:1173:THR:HA	2.21	0.41
1:A:1100:VAL:HG21	1:A:1123:VAL:HG21	2.02	0.41
2:B:27:LEU:HD23	2:B:27:LEU:HA	1.85	0.41
1:A:1102:PHE:O	1:A:1104:GLY:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:GLY:O	1:A:188:LYS:HD3	2.21	0.40
1:A:1025:LYS:HE2	1:A:1026:TYR:CE1	2.54	0.40
1:A:1020:VAL:O	1:A:1024:ARG:HG3	2.21	0.40
1:A:138:ILE:HA	1:A:1141:LEU:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1016:GLU:OE2	1:A:1024:ARG:NH1[7_554]	2.11	0.09

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	405/462 (88%)	370 (91%)	30 (7%)	5 (1%)	11	35
2	B	76/110 (69%)	74 (97%)	2 (3%)	0	100	100
All	All	481/572 (84%)	444 (92%)	32 (7%)	5 (1%)	13	40

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1150	ASP
1	A	1008	ASP
1	A	1103	THR
1	A	85	SER
1	A	1116	SER

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	366/405 (90%)	326 (89%)	40 (11%)	5	20
2	B	66/95 (70%)	60 (91%)	6 (9%)	7	27
All	All	432/500 (86%)	386 (89%)	46 (11%)	5	20

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	18	THR
1	A	21	ASP
1	A	25	SER
1	A	39	SER
1	A	59	GLU
1	A	99	ARG
1	A	101	VAL
1	A	102	ILE
1	A	126	GLU
1	A	140	LYS
1	A	153	SER
1	A	154	GLN
1	A	159	LEU
1	A	162	VAL
1	A	165	THR
1	A	176	ILE
1	A	188	LYS
1	A	190	ASN
1	A	200	LEU
1	A	991	VAL
1	A	1004	GLN
1	A	1013	SER
1	A	1019	ASN
1	A	1020	VAL
1	A	1032	THR
1	A	1042	GLU

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Mol	Chain	Res	Type
1	A	1054	THR
1	A	1090	VAL
1	A	1106	GLN
1	A	1112	MET
1	A	1116	SER
1	A	1121	SER
1	A	1125	LEU
1	A	1160	MET
1	A	1181	GLU
1	A	1189	VAL
1	A	1198	ILE
1	A	1200	VAL
1	A	1208	ASP
2	B	16	LYS
2	B	23	TRP
2	B	24	ASP
2	B	65	ARG
2	B	69	ARG
2	B	87	MET

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, 1 is monoatomic - leaving 1 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	AGS	A	2001	3	26,33,33	4.63	9 (34%)	26,52,52	1.58	4 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	AGS	A	2001	3	-	4/17/38/38	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2001	AGS	O4'-C1'	17.04	1.64	1.41
4	A	2001	AGS	C2'-C1'	-12.49	1.34	1.53
4	A	2001	AGS	O4'-C4'	-6.17	1.31	1.45
4	A	2001	AGS	PG-O2G	3.79	1.67	1.54
4	A	2001	AGS	O3'-C3'	-3.28	1.35	1.43
4	A	2001	AGS	C3'-C4'	3.27	1.61	1.53
4	A	2001	AGS	C6-N6	3.13	1.45	1.34
4	A	2001	AGS	PG-O3G	2.54	1.63	1.54
4	A	2001	AGS	O2'-C2'	2.20	1.48	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2001	AGS	N3-C2-N1	-5.10	120.71	128.68
4	A	2001	AGS	PA-O3A-PB	-3.51	120.78	132.83
4	A	2001	AGS	C4-C5-N7	-3.00	106.27	109.40
4	A	2001	AGS	O4'-C1'-C2'	-2.29	103.58	106.93

There are no chirality outliers.

All (4) torsion outliers are listed below:

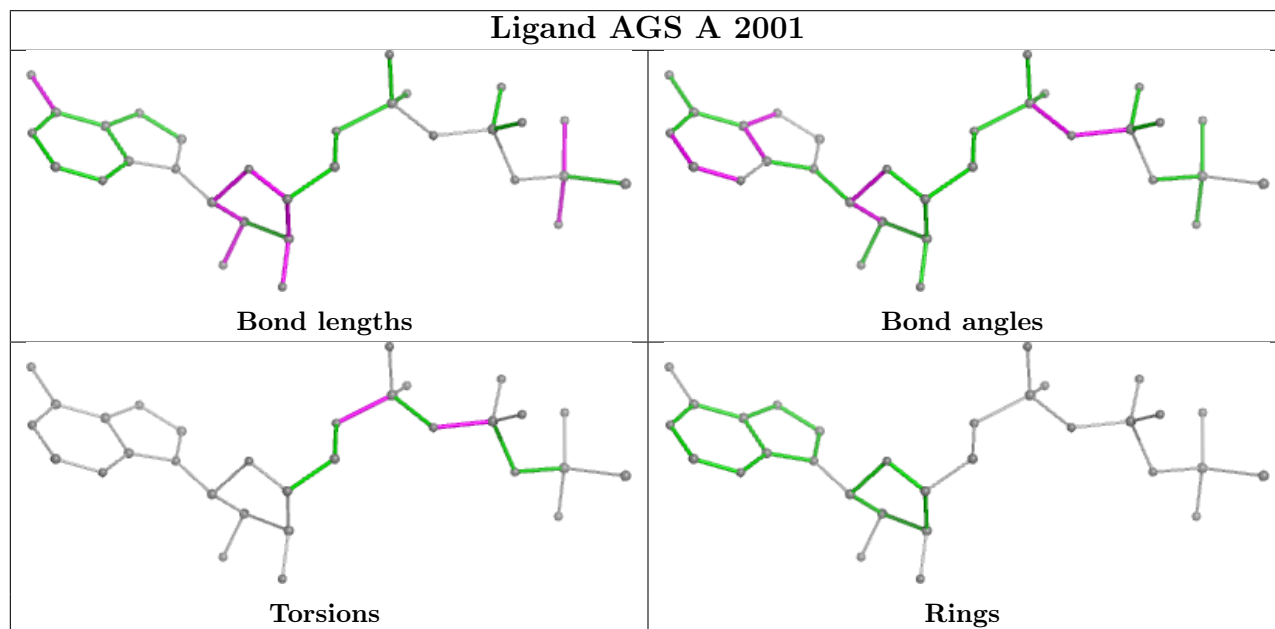
Mol	Chain	Res	Type	Atoms
4	A	2001	AGS	C5'-O5'-PA-O3A
4	A	2001	AGS	C5'-O5'-PA-O1A
4	A	2001	AGS	C5'-O5'-PA-O2A
4	A	2001	AGS	PA-O3A-PB-O2B

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2001	AGS	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	411/462 (88%)	-0.40	1 (0%) 92 84	74, 106, 159, 193	0
2	B	78/110 (70%)	-0.20	3 (3%) 44 28	78, 116, 162, 195	0
All	All	489/572 (85%)	-0.37	4 (0%) 82 68	74, 107, 160, 195	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	48	LYS	4.3
2	B	49	VAL	2.8
2	B	89	PHE	2.6
1	A	1106	GLN	2.1

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.

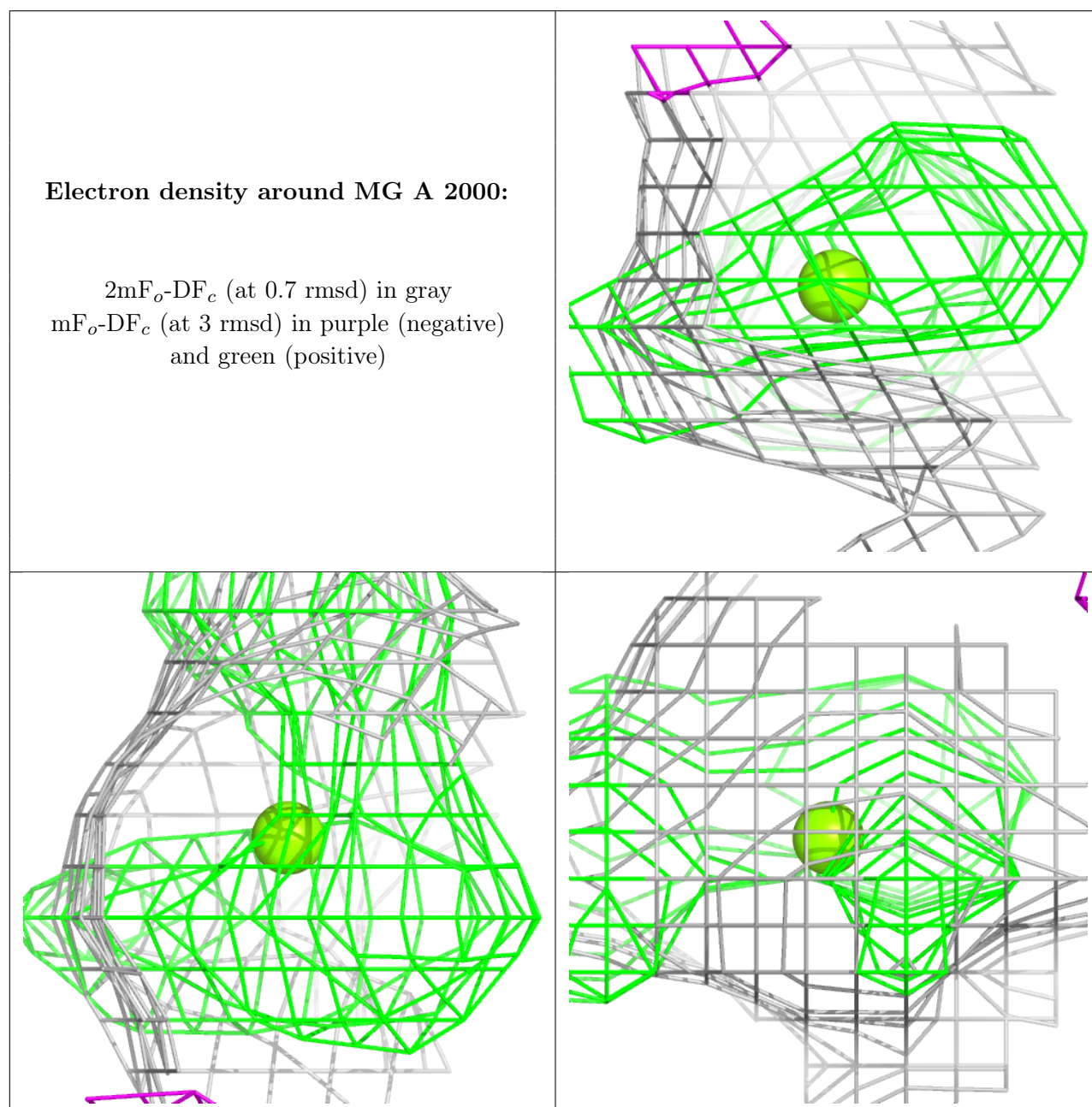
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	A	2000	1/1	0.92	0.17	79,79,79,79	0

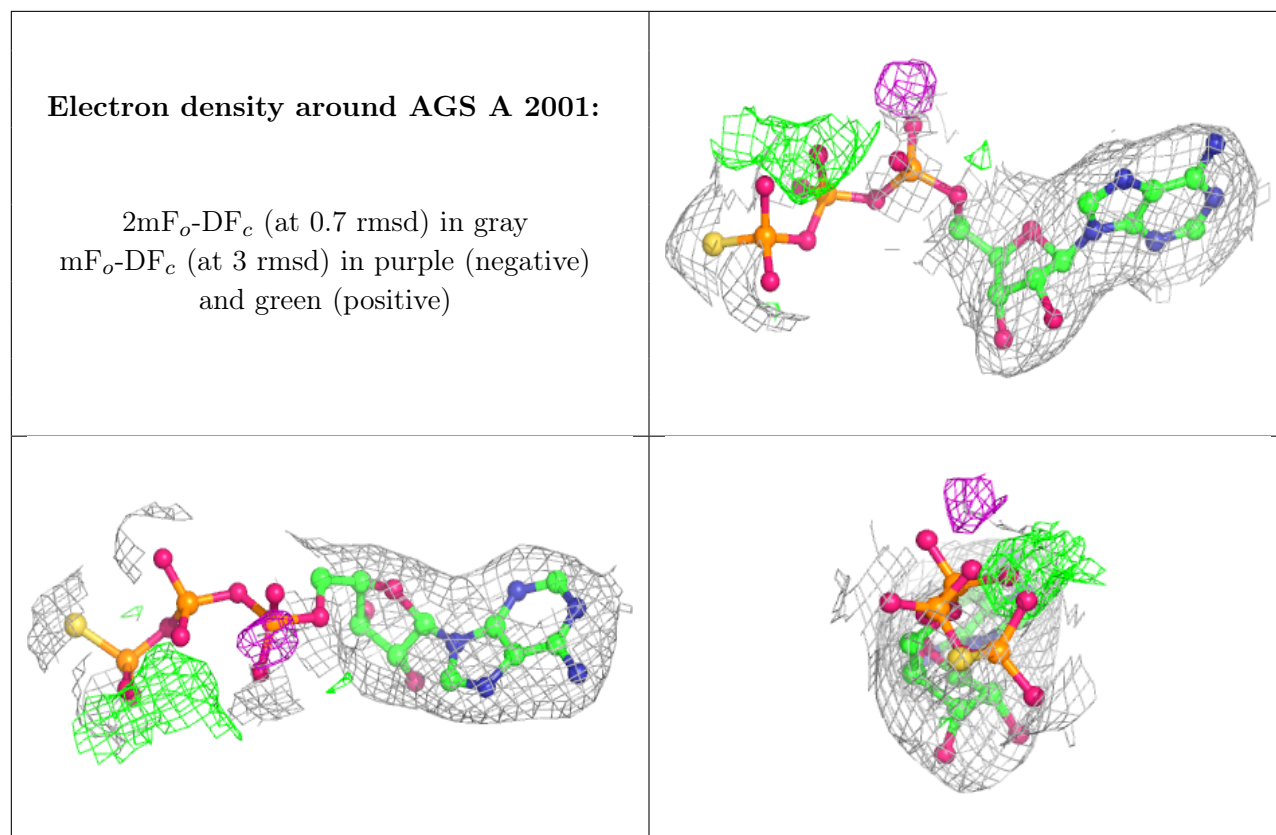
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	AGS	A	2001	31/31	0.97	0.06	68,95,106,115	0

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 [i](#)

There are no such residues in this entry.