



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 25, 2024 – 09:37 PM EDT

PDB ID : 7CK1
Title : Crystal structure of arabidopsis CESA3 catalytic domain
Authors : Qiao, Z.; Gao, Y.G.
Deposited on : 2020-07-15
Resolution : 2.35 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
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

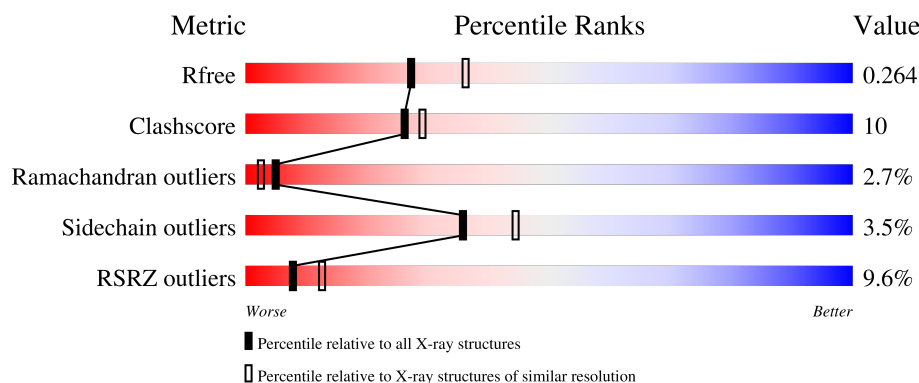
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	410	<div> <div>9%</div> <div>65%</div> <div>19%</div> <div>14%</div> </div>
1	B	410	<div> <div>8%</div> <div>64%</div> <div>19%</div> <div>14%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5573 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 Cellulose synthase A catalytic subunit 3 [UDP-forming], Cellulose synthase A catalytic subunit 3 [UDP-forming].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	0	0
			2788	1781	473	515	19			
1	B	351	Total	C	N	O	S	0	0	0
			2764	1767	469	509	19			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	697	GLY	-	linker	UNP Q941L0
A	698	SER	-	linker	UNP Q941L0
A	699	GLY	-	linker	UNP Q941L0
A	700	SER	-	linker	UNP Q941L0
A	701	GLY	-	linker	UNP Q941L0
B	697	GLY	-	linker	UNP Q941L0
B	698	SER	-	linker	UNP Q941L0
B	699	GLY	-	linker	UNP Q941L0
B	700	SER	-	linker	UNP Q941L0
B	701	GLY	-	linker	UNP Q941L0

- Molecule 2 is MANGANESE (II) ION (three-letter code: 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		

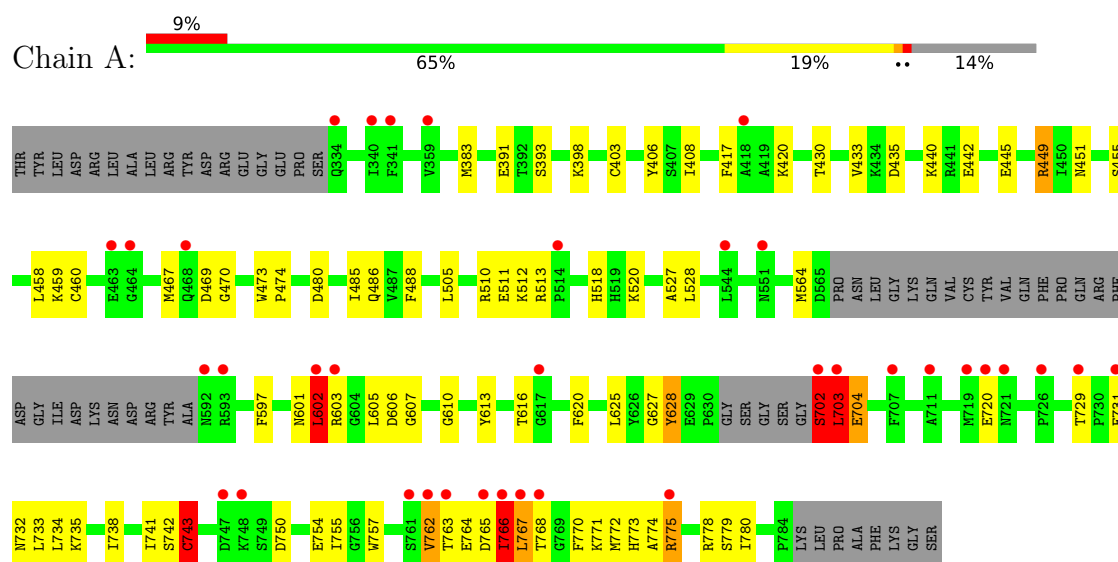
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total 8	O 8	0	0
3	B	11	Total 11	O 11	0	0

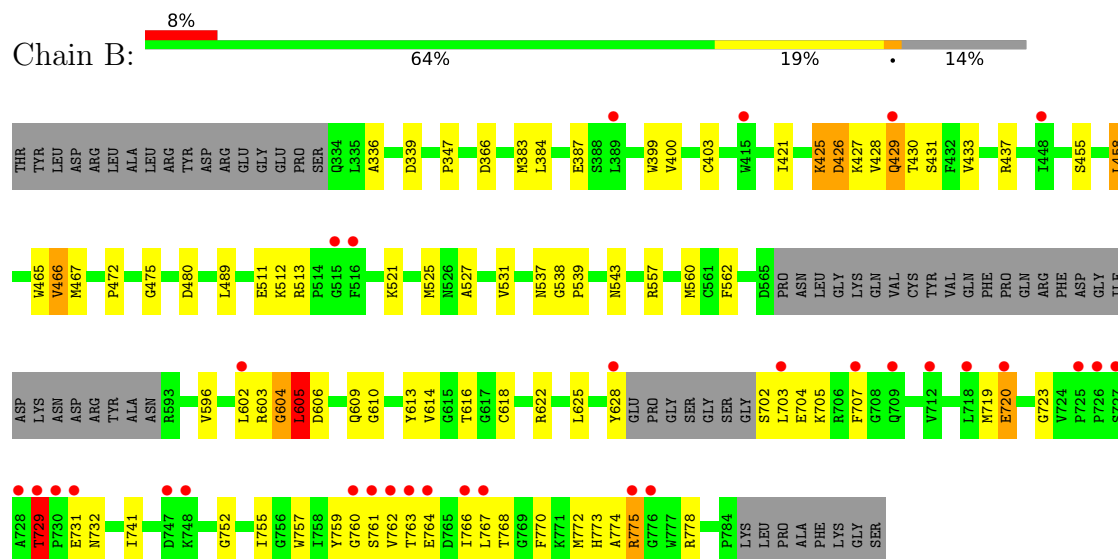
3 Residue-property plots

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: Cellulose synthase A catalytic subunit 3 [UDP-forming], Cellulose synthase A catalytic subunit 3 [UDP-forming]



- Molecule 1: Cellulose synthase A catalytic subunit 3 [UDP-forming], Cellulose synthase A catalytic subunit 3 [UDP-forming]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.62Å 147.03Å 92.51Å 90.00° 102.05° 90.00°	Depositor
Resolution (Å)	48.08 – 2.35 48.08 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.08-2.35) 99.8 (48.08-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.237 , 0.264 0.238 , 0.264	Depositor DCC
R_{free} test set	2961 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5573	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN

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	0.49	1/2856 (0.0%)	0.72	5/3869 (0.1%)
1	B	0.48	0/2831	0.65	2/3834 (0.1%)
All	All	0.48	1/5687 (0.0%)	0.69	7/7703 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	2
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	743	CYS	CB-SG	-5.34	1.73	1.81

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	602	LEU	CA-CB-CG	9.86	137.98	115.30
1	A	767	LEU	CA-CB-CG	8.13	133.99	115.30
1	A	703	LEU	CA-CB-CG	6.51	130.28	115.30
1	B	602	LEU	CB-CG-CD1	6.32	121.74	111.00
1	A	733	LEU	CA-CB-CG	5.86	128.78	115.30
1	B	604	GLY	N-CA-C	5.14	125.95	113.10
1	A	602	LEU	CB-CG-CD2	5.03	119.55	111.00

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	602	LEU	Peptide
1	A	702	SER	Peptide
1	A	742	SER	Peptide
1	A	766	ILE	Peptide
1	B	603	ARG	Peptide
1	B	729	THR	Peptide

5.2 Too-close contacts

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	2788	0	2738	65	0
1	B	2764	0	2719	52	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	8	0	0	2	0
3	B	11	0	0	5	0
All	All	5573	0	5457	115	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 10.

All (115) 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:A:732:ASN:HA	1:A:735:LYS:HG2	1.47	0.94
1:A:766:ILE:HA	1:A:768:THR:HG22	1.54	0.88
1:B:387:GLU:N	3:B:901:HOH:O	2.09	0.85
1:A:764:GLU:HB3	1:A:766:ILE:HD12	1.59	0.83
1:B:336:ALA:H	1:B:560:MET:HE2	1.41	0.83
1:B:429:GLN:O	1:B:431:SER:N	2.17	0.78
1:B:761:SER:O	1:B:763:THR:N	2.19	0.74
1:A:603:ARG:HB3	1:A:606:ASP:OD1	1.89	0.72
1:B:475:GLY:O	3:B:902:HOH:O	2.09	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:527:ALA:HB1	1:B:741:ILE:HG22	1.75	0.68
1:A:430:THR:HG22	1:A:433:VAL:HG13	1.77	0.65
1:A:628:TYR:C	1:A:775:ARG:HD2	2.17	0.65
1:A:627:GLY:O	1:A:628:TYR:HB2	2.00	0.61
1:A:766:ILE:HG22	1:A:767:LEU:N	2.15	0.61
1:B:384:LEU:O	3:B:901:HOH:O	2.17	0.60
1:A:702:SER:O	1:A:704:GLU:N	2.34	0.60
1:A:459:LYS:HD2	1:A:460:CYS:N	2.18	0.59
1:A:602:LEU:HB3	1:A:603:ARG:HG2	1.85	0.59
1:B:625:LEU:HD21	1:B:770:PHE:HB3	1.84	0.59
1:A:628:TYR:O	1:A:775:ARG:NH1	2.35	0.58
1:A:518:HIS:HE1	1:A:520:LYS:HD3	1.67	0.58
1:B:339:ASP:OD2	1:B:622:ARG:NH2	2.37	0.57
1:A:606:ASP:OD1	1:A:606:ASP:N	2.37	0.57
1:A:720:GLU:OE1	1:A:720:GLU:N	2.37	0.57
1:B:537:ASN:HB3	1:B:723:GLY:H	1.69	0.57
1:A:729:THR:HG23	1:A:731:GLU:HG2	1.88	0.56
1:A:750:ASP:HB2	1:A:755:ILE:HD12	1.86	0.56
1:B:336:ALA:H	1:B:560:MET:CE	2.14	0.56
1:B:384:LEU:C	3:B:901:HOH:O	2.44	0.55
1:B:763:THR:HG23	1:B:764:GLU:H	1.71	0.55
1:A:510:ARG:HD2	3:A:902:HOH:O	2.06	0.55
1:A:403:CYS:HA	1:A:408:ILE:HD13	1.89	0.55
1:B:604:GLY:O	1:B:606:ASP:N	2.37	0.54
1:A:398:LYS:HE2	1:A:442:GLU:OE2	2.07	0.54
1:A:625:LEU:HD21	1:A:770:PHE:HB3	1.90	0.54
1:A:602:LEU:HB3	1:A:603:ARG:CG	2.37	0.53
1:A:774:ALA:HA	1:A:778:ARG:CZ	2.38	0.53
1:A:602:LEU:CD2	1:A:603:ARG:H	2.22	0.53
1:B:614:VAL:HB	1:B:772:MET:HE1	1.89	0.52
1:B:760:GLY:O	1:B:767:LEU:HD12	2.10	0.52
1:B:628:TYR:C	1:B:775:ARG:HE	2.13	0.51
1:A:527:ALA:HB1	1:A:741:ILE:HG22	1.93	0.51
1:B:347:PRO:HB2	1:B:384:LEU:HD23	1.91	0.51
1:A:754:GLU:OE1	1:A:771:LYS:NZ	2.41	0.51
1:A:764:GLU:HB3	1:A:766:ILE:CD1	2.36	0.51
1:A:757:TRP:HE1	1:A:767:LEU:C	2.15	0.50
1:B:707:PHE:CE1	1:B:755:ILE:HD13	2.47	0.50
1:A:597:PHE:HB3	1:A:780:ILE:HG12	1.94	0.50
1:A:607:GLY:HA2	1:B:618:CYS:SG	2.52	0.49
1:B:366:ASP:OD1	1:B:557:ARG:NH1	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:562:PHE:CZ	1:B:596:VAL:HG13	2.48	0.49
1:B:426:ASP:O	1:B:428:VAL:N	2.46	0.48
1:B:729:THR:O	1:B:731:GLU:N	2.46	0.48
1:A:772:MET:SD	1:A:780:ILE:HD11	2.52	0.48
1:B:480:ASP:HA	1:B:512:LYS:O	2.14	0.48
1:A:420:LYS:HA	1:A:420:LYS:HD2	1.58	0.47
1:B:521:LYS:O	1:B:525:MET:HG2	2.14	0.47
1:A:485:ILE:HD12	1:A:738:ILE:HG13	1.96	0.47
1:A:613:TYR:HB2	1:B:609:GLN:O	2.14	0.47
1:B:702:SER:O	1:B:704:GLU:N	2.47	0.47
1:B:527:ALA:O	1:B:531:VAL:HG23	2.15	0.47
1:B:610:GLY:O	1:B:774:ALA:HB2	2.15	0.47
1:B:466:VAL:HA	1:B:472:PRO:HA	1.96	0.46
1:B:752:GLY:HA3	1:B:759:TYR:CD2	2.51	0.46
1:A:406:TYR:OH	1:A:435:ASP:OD2	2.23	0.46
1:A:735:LYS:N	1:A:735:LYS:HD3	2.30	0.46
1:A:601:ASN:OD1	1:A:603:ARG:HD2	2.16	0.46
1:A:445:GLU:O	1:A:449:ARG:HG2	2.16	0.46
1:B:383:MET:O	3:B:901:HOH:O	2.20	0.46
1:B:613:TYR:CD1	1:B:766:ILE:HD11	2.51	0.45
1:A:779:SER:C	1:A:780:ILE:HD13	2.36	0.45
1:A:766:ILE:CG2	1:A:768:THR:H	2.30	0.45
1:A:518:HIS:CE1	1:A:520:LYS:HD3	2.50	0.45
1:A:451:ASN:H	1:A:451:ASN:HD22	1.65	0.45
1:A:393:SER:HB3	1:A:505:LEU:HD23	1.98	0.45
1:A:486:GLN:HG2	1:A:488:PHE:CZ	2.51	0.44
1:A:480:ASP:OD1	1:A:513:ARG:NH1	2.48	0.44
1:A:773:HIS:O	1:A:778:ARG:HD3	2.18	0.44
1:B:383:MET:CE	1:B:467:MET:HA	2.48	0.44
1:B:383:MET:HE3	1:B:467:MET:HA	1.99	0.44
1:A:528:LEU:HD23	1:A:528:LEU:HA	1.87	0.44
1:A:610:GLY:O	1:A:774:ALA:HB2	2.18	0.44
1:A:731:GLU:O	1:A:734:LEU:N	2.51	0.44
1:B:752:GLY:HA3	1:B:759:TYR:CE2	2.53	0.44
1:B:702:SER:O	1:B:705:LYS:NZ	2.46	0.43
1:B:543:ASN:OD1	1:B:616:THR:HG22	2.18	0.43
1:A:510:ARG:HG2	3:A:907:HOH:O	2.19	0.43
1:A:391:GLU:CD	1:A:449:ARG:HH21	2.22	0.43
1:A:391:GLU:OE1	1:A:449:ARG:NH2	2.49	0.42
1:A:762:VAL:HG13	1:A:764:GLU:OE1	2.19	0.42
1:B:433:VAL:O	1:B:437:ARG:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:707:PHE:HE1	1:B:755:ILE:HD13	1.82	0.42
1:A:613:TYR:OH	1:A:771:LYS:HD2	2.19	0.42
1:B:704:GLU:HB2	1:B:705:LYS:HZ2	1.84	0.42
1:B:763:THR:HG23	1:B:764:GLU:N	2.34	0.42
1:A:383:MET:CE	1:A:467:MET:HG2	2.50	0.42
1:A:764:GLU:O	1:A:766:ILE:HD12	2.19	0.42
1:A:602:LEU:HD22	1:A:603:ARG:H	1.85	0.42
1:A:473:TRP:HA	1:A:474:PRO:HD3	1.91	0.42
1:A:430:THR:CG2	1:A:433:VAL:HG13	2.45	0.41
1:A:512:LYS:O	1:A:513:ARG:HD3	2.20	0.41
1:A:511:GLU:CG	1:A:743:CYS:HB2	2.50	0.41
1:B:605:LEU:HD13	1:B:605:LEU:HA	1.81	0.41
1:A:417:PHE:CE1	1:A:440:LYS:HB2	2.56	0.41
1:B:383:MET:SD	1:B:465:TRP:HB3	2.61	0.41
1:B:752:GLY:HA2	1:B:757:TRP:O	2.21	0.41
1:B:538:GLY:HA2	1:B:539:PRO:HD3	1.95	0.41
1:B:773:HIS:O	1:B:778:ARG:HD3	2.21	0.41
1:B:399:TRP:CE2	1:B:403:CYS:SG	3.14	0.41
1:B:511:GLU:OE2	1:B:513:ARG:NE	2.50	0.40
1:A:620:PHE:HA	1:A:779:SER:O	2.22	0.40
1:B:425:LYS:O	1:B:426:ASP:O	2.40	0.40
1:A:764:GLU:O	1:A:765:ASP:HB2	2.21	0.40
1:A:511:GLU:HG2	1:A:743:CYS:HB2	2.02	0.40
1:B:455:SER:HA	1:B:458:LEU:HD23	2.04	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	348/410 (85%)	319 (92%)	19 (6%)	10 (3%)	4 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	345/410 (84%)	312 (90%)	24 (7%)	9 (3%)	5	3
All	All	693/820 (84%)	631 (91%)	43 (6%)	19 (3%)	5	2

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	469	ASP
1	A	602	LEU
1	A	628	TYR
1	A	703	LEU
1	A	704	GLU
1	A	743	CYS
1	A	775	ARG
1	B	426	ASP
1	B	430	THR
1	B	703	LEU
1	B	720	GLU
1	B	762	VAL
1	B	427	LYS
1	B	605	LEU
1	A	763	THR
1	A	766	ILE
1	B	429	GLN
1	B	775	ARG
1	A	470	GLY

5.3.2 Protein sidechains ⓘ

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	300/346 (87%)	291 (97%)	9 (3%)	41	50
1	B	297/346 (86%)	285 (96%)	12 (4%)	31	39
All	All	597/692 (86%)	576 (96%)	21 (4%)	36	44

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	449	ARG
1	A	455	SER
1	A	458	LEU
1	A	564	MET
1	A	605	LEU
1	A	616	THR
1	A	702	SER
1	A	703	LEU
1	A	762	VAL
1	B	400	VAL
1	B	421	ILE
1	B	425	LYS
1	B	458	LEU
1	B	466	VAL
1	B	489	LEU
1	B	605	LEU
1	B	719	MET
1	B	720	GLU
1	B	729	THR
1	B	732	ASN
1	B	768	THR

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 monosaccharides 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	A	354/410 (86%)	0.82	36 (10%) 6 11	56, 73, 113, 135	0
1	B	351/410 (85%)	0.73	32 (9%) 9 14	54, 72, 110, 144	0
All	All	705/820 (85%)	0.77	68 (9%) 8 12	54, 72, 111, 144	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	727	SER	10.2
1	B	602	LEU	9.8
1	A	767	LEU	7.0
1	B	762	VAL	6.7
1	A	726	PRO	5.5
1	B	726	PRO	5.3
1	A	762	VAL	4.9
1	B	515	GLY	4.9
1	B	728	ALA	4.9
1	A	766	ILE	4.8
1	B	763	THR	4.6
1	A	765	ASP	4.5
1	A	768	THR	4.4
1	B	730	PRO	4.0
1	B	729	THR	3.9
1	B	761	SER	3.8
1	A	763	THR	3.8
1	A	334	GLN	3.7
1	B	718	LEU	3.6
1	B	748	LYS	3.6
1	A	775	ARG	3.6
1	B	764	GLU	3.5
1	B	429	GLN	3.5
1	B	767	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	628	TYR	3.2
1	A	729	THR	3.2
1	A	514	PRO	3.1
1	A	593	ARG	3.0
1	A	703	LEU	3.0
1	B	725	PRO	2.9
1	A	418	ALA	2.9
1	A	748	LYS	2.8
1	B	731	GLU	2.8
1	B	516	PHE	2.8
1	A	468	GLN	2.8
1	A	761	SER	2.6
1	B	712	VAL	2.6
1	A	592	ASN	2.6
1	B	747	ASP	2.5
1	A	731	GLU	2.5
1	A	544	LEU	2.5
1	A	602	LEU	2.4
1	B	703	LEU	2.4
1	B	389	LEU	2.4
1	A	721	ASN	2.4
1	A	463	GLU	2.4
1	A	603	ARG	2.4
1	A	464	GLY	2.4
1	A	340	ILE	2.3
1	B	766	ILE	2.3
1	B	709	GLN	2.3
1	A	551	ASN	2.3
1	A	711	ALA	2.3
1	B	707	PHE	2.2
1	A	719	MET	2.2
1	A	702	SER	2.2
1	A	720	GLU	2.2
1	B	775	ARG	2.1
1	B	448	ILE	2.1
1	B	415	TRP	2.1
1	A	617	GLY	2.1
1	B	720	GLU	2.1
1	A	747	ASP	2.1
1	B	776	GLY	2.0
1	A	359	VAL	2.0
1	A	341	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	707	PHE	2.0
1	B	760	GLY	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 monosaccharides in this entry.

6.4 Ligands [i](#)

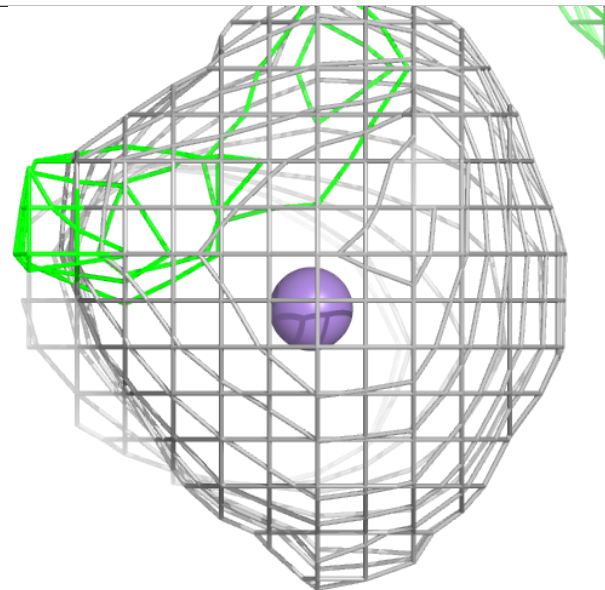
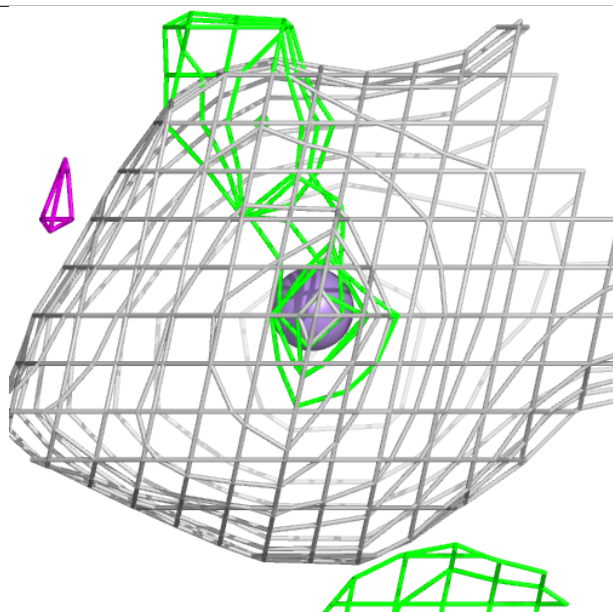
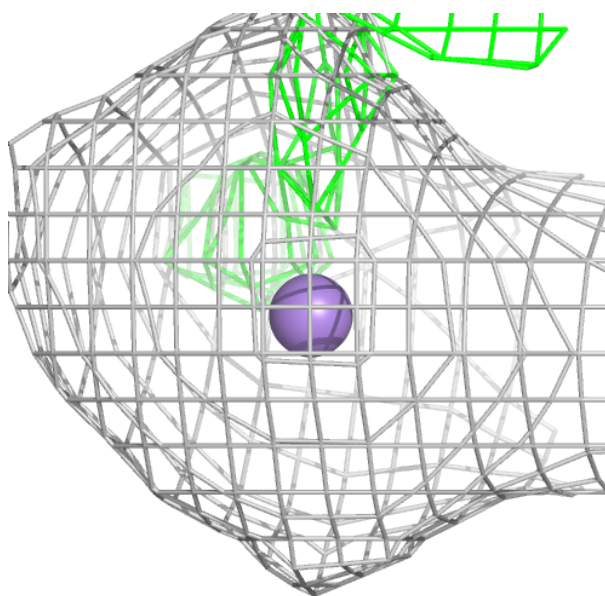
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MN	A	801	1/1	0.97	0.24	85,85,85,85	1
2	MN	B	801	1/1	0.97	0.14	93,93,93,93	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.

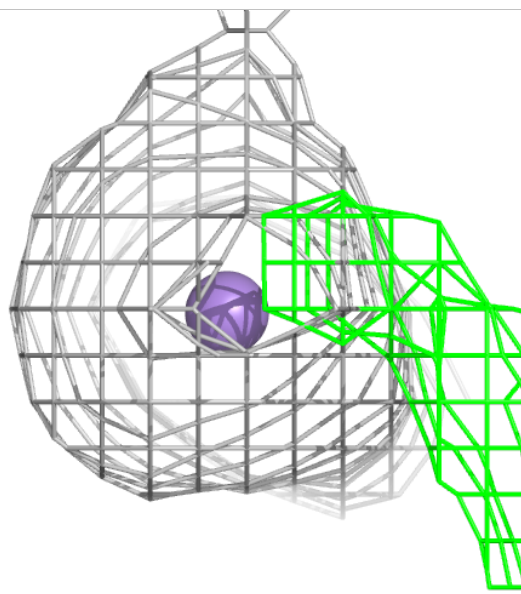
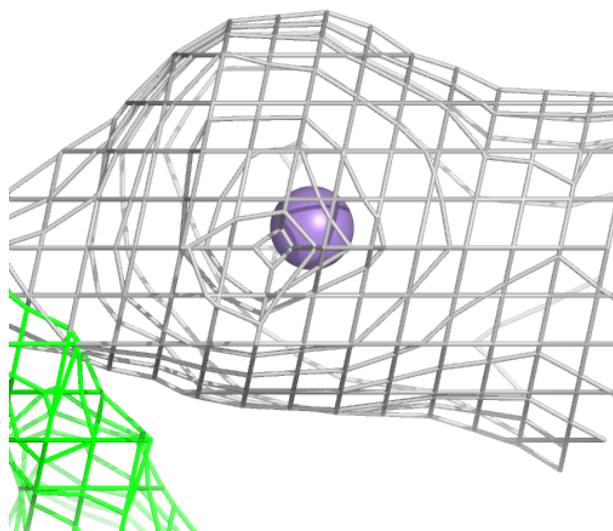
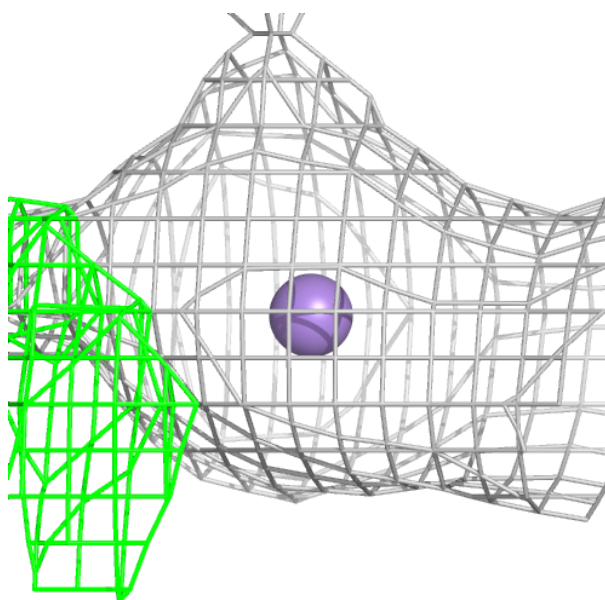
Electron density around MN A 801:

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



Electron density around MN B 801:

$2mF_o-DF_c$ (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.