



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:25 PM UTC

PDB ID : 7KYP / pdb_00007kyp
Title : PsaBC from Streptococcus pneumoniae in complex with Fab
Authors : Maher, M.J.; Sjöhamn, J.
Deposited on : 2020-12-08
Resolution : 2.90 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:

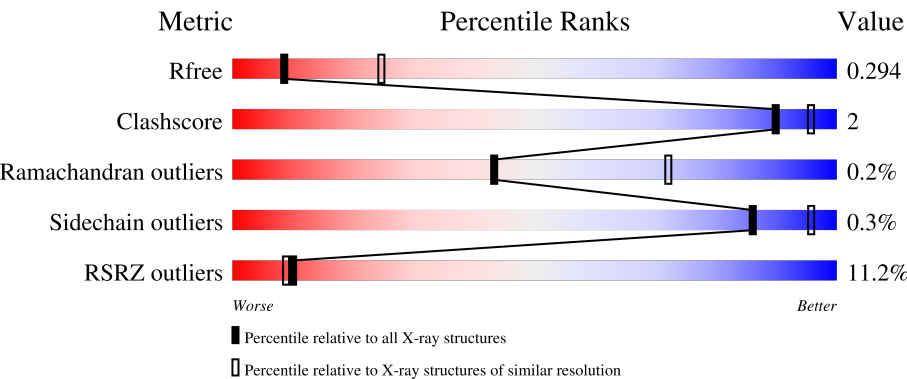
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	240	
1	C	240	
1	E	240	
1	G	240	
1	I	240	

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Mol	Chain	Length	Quality of chain
1	K	240	
1	M	240	
1	O	240	
2	B	290	
2	D	290	
2	F	290	
2	H	290	
2	J	290	
2	L	290	
2	N	290	
2	P	290	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	C	301	-	-	-	X
3	PO4	I	301	-	-	-	X
3	PO4	K	301	-	-	-	X
3	PO4	M	301	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 31187 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 Manganese ABC transporter, ATP-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	0	1	0
			1839	1188	309	337	5			
1	C	232	Total	C	N	O	S	0	1	0
			1793	1155	302	331	5			
1	E	236	Total	C	N	O	S	0	1	0
			1835	1185	308	337	5			
1	G	230	Total	C	N	O	S	0	0	0
			1775	1147	296	327	5			
1	I	234	Total	C	N	O	S	0	0	0
			1831	1182	309	335	5			
1	K	236	Total	C	N	O	S	0	0	0
			1850	1195	312	338	5			
1	M	233	Total	C	N	O	S	0	1	0
			1815	1169	308	333	5			
1	O	234	Total	C	N	O	S	0	1	0
			1810	1170	302	333	5			

- Molecule 2 is a protein called Manganese ABC transporter, permease protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	278	Total	C	N	O	S	0	0	0
			2080	1396	315	357	12			
2	D	272	Total	C	N	O	S	0	0	0
			2027	1362	303	350	12			
2	F	290	Total	C	N	O	S	0	0	0
			2173	1457	329	375	12			
2	H	272	Total	C	N	O	S	0	0	0
			2012	1355	301	345	11			
2	J	274	Total	C	N	O	S	0	0	0
			2015	1352	303	349	11			
2	L	278	Total	C	N	O	S	0	0	0
			2065	1384	313	356	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	277	Total	C	N	O	S	0	0	0
			2060	1381	312	355	12			
2	P	271	Total	C	N	O	S	0	0	0
			2015	1353	302	348	12			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	100	SER	PHE	conflict	UNP A0A0H2ZPI2
B	283	GLY	-	expression tag	UNP A0A0H2ZPI2
B	284	SER	-	expression tag	UNP A0A0H2ZPI2
B	285	GLU	-	expression tag	UNP A0A0H2ZPI2
B	286	ASN	-	expression tag	UNP A0A0H2ZPI2
B	287	LEU	-	expression tag	UNP A0A0H2ZPI2
B	288	TYR	-	expression tag	UNP A0A0H2ZPI2
B	289	PHE	-	expression tag	UNP A0A0H2ZPI2
B	290	GLN	-	expression tag	UNP A0A0H2ZPI2
D	100	SER	PHE	conflict	UNP A0A0H2ZPI2
D	283	GLY	-	expression tag	UNP A0A0H2ZPI2
D	284	SER	-	expression tag	UNP A0A0H2ZPI2
D	285	GLU	-	expression tag	UNP A0A0H2ZPI2
D	286	ASN	-	expression tag	UNP A0A0H2ZPI2
D	287	LEU	-	expression tag	UNP A0A0H2ZPI2
D	288	TYR	-	expression tag	UNP A0A0H2ZPI2
D	289	PHE	-	expression tag	UNP A0A0H2ZPI2
D	290	GLN	-	expression tag	UNP A0A0H2ZPI2
F	100	SER	PHE	conflict	UNP A0A0H2ZPI2
F	283	GLY	-	expression tag	UNP A0A0H2ZPI2
F	284	SER	-	expression tag	UNP A0A0H2ZPI2
F	285	GLU	-	expression tag	UNP A0A0H2ZPI2
F	286	ASN	-	expression tag	UNP A0A0H2ZPI2
F	287	LEU	-	expression tag	UNP A0A0H2ZPI2
F	288	TYR	-	expression tag	UNP A0A0H2ZPI2
F	289	PHE	-	expression tag	UNP A0A0H2ZPI2
F	290	GLN	-	expression tag	UNP A0A0H2ZPI2
H	100	SER	PHE	conflict	UNP A0A0H2ZPI2
H	283	GLY	-	expression tag	UNP A0A0H2ZPI2
H	284	SER	-	expression tag	UNP A0A0H2ZPI2
H	285	GLU	-	expression tag	UNP A0A0H2ZPI2
H	286	ASN	-	expression tag	UNP A0A0H2ZPI2
H	287	LEU	-	expression tag	UNP A0A0H2ZPI2
H	288	TYR	-	expression tag	UNP A0A0H2ZPI2

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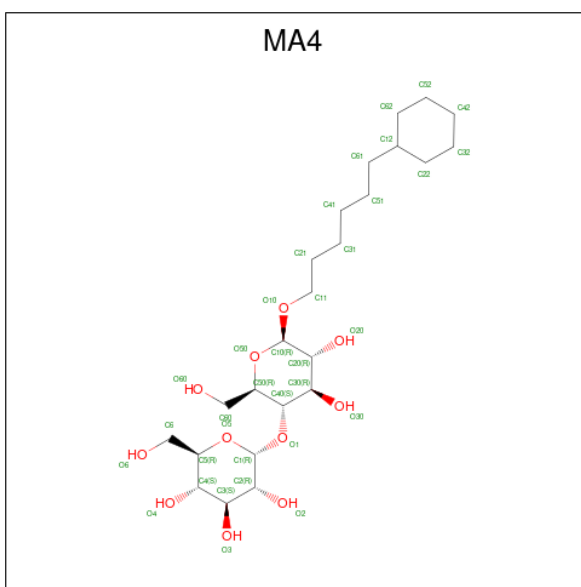
Chain	Residue	Modelled	Actual	Comment	Reference
H	289	PHE	-	expression tag	UNP A0A0H2ZPI2
H	290	GLN	-	expression tag	UNP A0A0H2ZPI2
J	100	SER	PHE	conflict	UNP A0A0H2ZPI2
J	283	GLY	-	expression tag	UNP A0A0H2ZPI2
J	284	SER	-	expression tag	UNP A0A0H2ZPI2
J	285	GLU	-	expression tag	UNP A0A0H2ZPI2
J	286	ASN	-	expression tag	UNP A0A0H2ZPI2
J	287	LEU	-	expression tag	UNP A0A0H2ZPI2
J	288	TYR	-	expression tag	UNP A0A0H2ZPI2
J	289	PHE	-	expression tag	UNP A0A0H2ZPI2
J	290	GLN	-	expression tag	UNP A0A0H2ZPI2
L	100	SER	PHE	conflict	UNP A0A0H2ZPI2
L	283	GLY	-	expression tag	UNP A0A0H2ZPI2
L	284	SER	-	expression tag	UNP A0A0H2ZPI2
L	285	GLU	-	expression tag	UNP A0A0H2ZPI2
L	286	ASN	-	expression tag	UNP A0A0H2ZPI2
L	287	LEU	-	expression tag	UNP A0A0H2ZPI2
L	288	TYR	-	expression tag	UNP A0A0H2ZPI2
L	289	PHE	-	expression tag	UNP A0A0H2ZPI2
L	290	GLN	-	expression tag	UNP A0A0H2ZPI2
N	100	SER	PHE	conflict	UNP A0A0H2ZPI2
N	283	GLY	-	expression tag	UNP A0A0H2ZPI2
N	284	SER	-	expression tag	UNP A0A0H2ZPI2
N	285	GLU	-	expression tag	UNP A0A0H2ZPI2
N	286	ASN	-	expression tag	UNP A0A0H2ZPI2
N	287	LEU	-	expression tag	UNP A0A0H2ZPI2
N	288	TYR	-	expression tag	UNP A0A0H2ZPI2
N	289	PHE	-	expression tag	UNP A0A0H2ZPI2
N	290	GLN	-	expression tag	UNP A0A0H2ZPI2
P	100	SER	PHE	conflict	UNP A0A0H2ZPI2
P	283	GLY	-	expression tag	UNP A0A0H2ZPI2
P	284	SER	-	expression tag	UNP A0A0H2ZPI2
P	285	GLU	-	expression tag	UNP A0A0H2ZPI2
P	286	ASN	-	expression tag	UNP A0A0H2ZPI2
P	287	LEU	-	expression tag	UNP A0A0H2ZPI2
P	288	TYR	-	expression tag	UNP A0A0H2ZPI2
P	289	PHE	-	expression tag	UNP A0A0H2ZPI2
P	290	GLN	-	expression tag	UNP A0A0H2ZPI2

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	G	1	Total	O	P	0	0
			5	4	1		
3	I	1	Total	O	P	0	0
			5	4	1		
3	K	1	Total	O	P	0	0
			5	4	1		
3	M	1	Total	O	P	0	0
			5	4	1		
3	O	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is CYCLOHEXYL-HEXYL-BETA-D-MALTOSIDE (CCD ID: MA4) (formula: $C_{24}H_{44}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total 35	C 24	O 11	0	0
4	F	1	Total 35	C 24	O 11	0	0
4	L	1	Total 35	C 24	O 11	0	0
4	N	1	Total 35	C 24	O 11	0	0

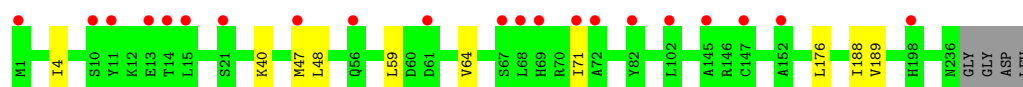
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total O 2 2	0	0
5	C	1	Total O 1 1	0	0
5	D	1	Total O 1 1	0	0
5	E	3	Total O 3 3	0	0
5	H	1	Total O 1 1	0	0
5	O	3	Total O 3 3	0	0
5	P	1	Total O 1 1	0	0

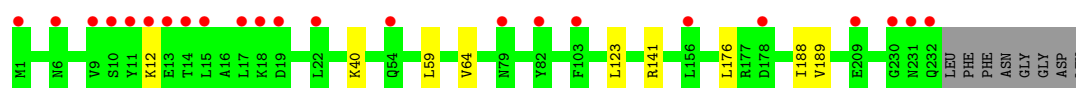
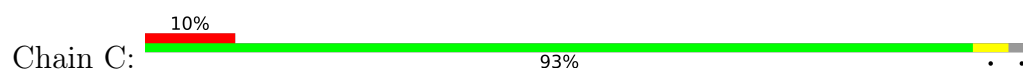
3 Residue-property plots [i](#)

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.

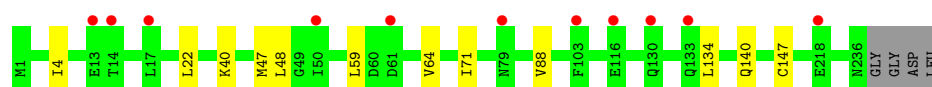
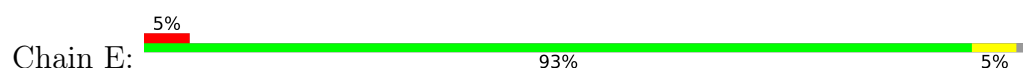
- Molecule 1: Manganese ABC transporter, ATP-binding protein



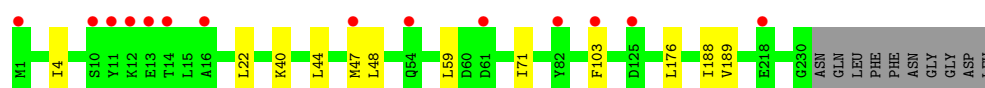
- Molecule 1: Manganese ABC transporter, ATP-binding protein



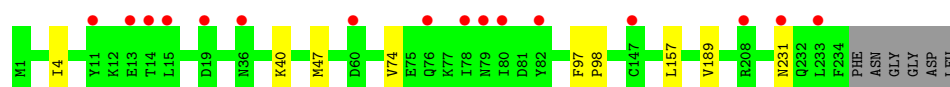
- Molecule 1: Manganese ABC transporter, ATP-binding protein



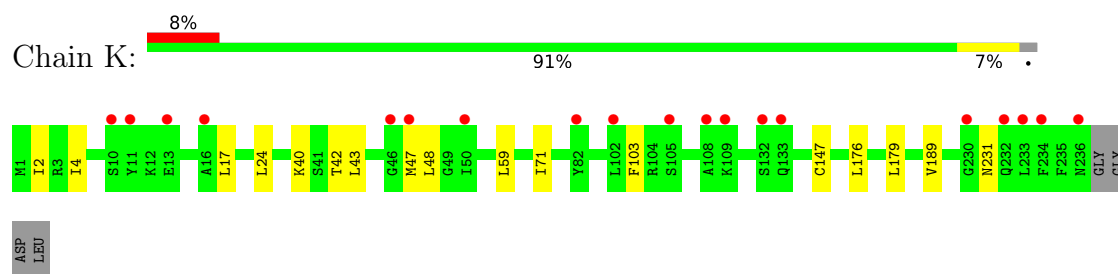
- Molecule 1: Manganese ABC transporter, ATP-binding protein



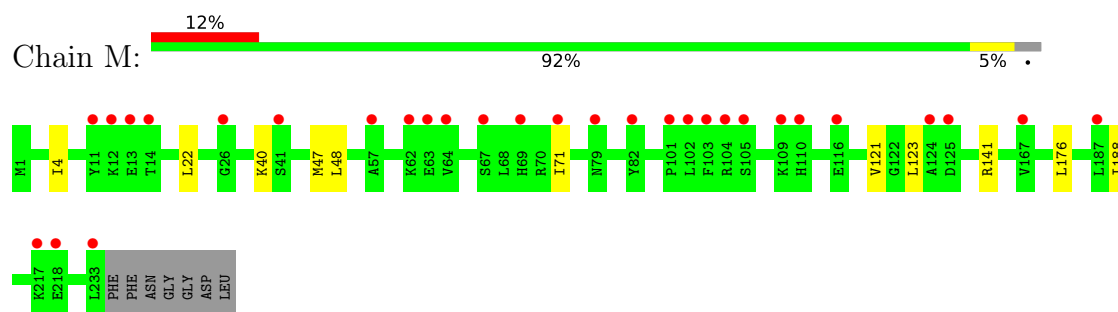
- Molecule 1: Manganese ABC transporter, ATP-binding protein



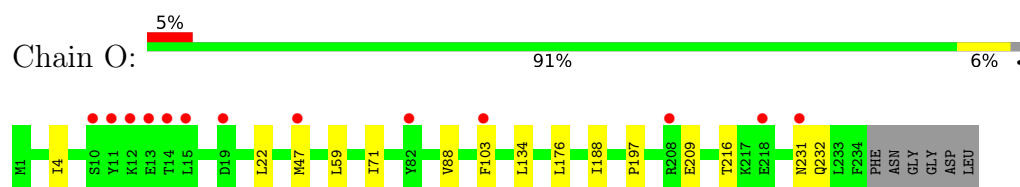
- Molecule 1: Manganese ABC transporter, ATP-binding protein



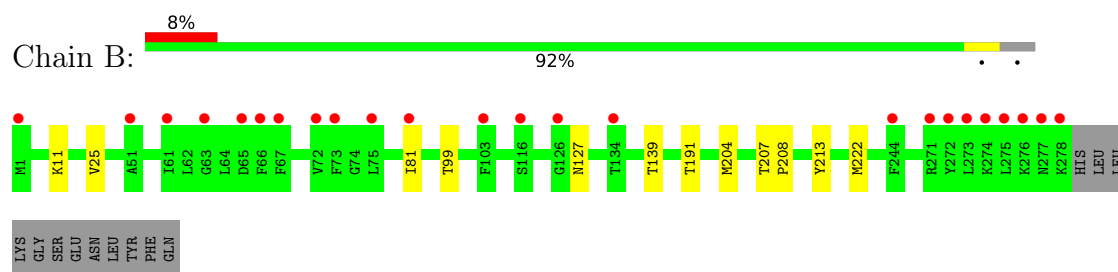
- Molecule 1: Manganese ABC transporter, ATP-binding protein



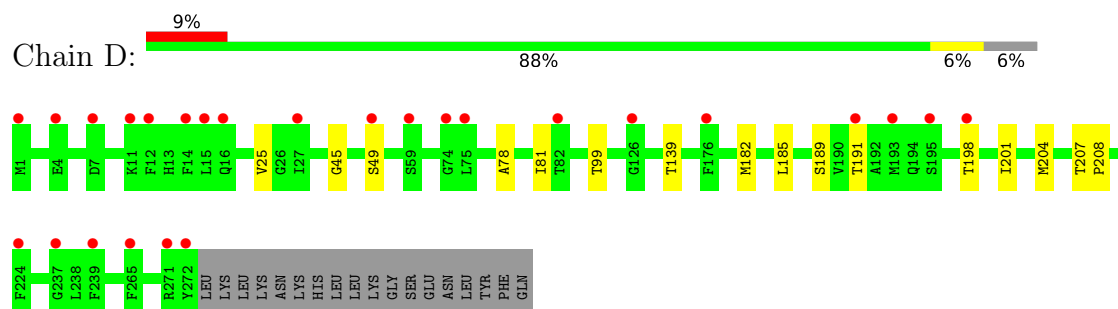
- Molecule 1: Manganese ABC transporter, ATP-binding protein



- Molecule 2: Manganese ABC transporter, permease protein

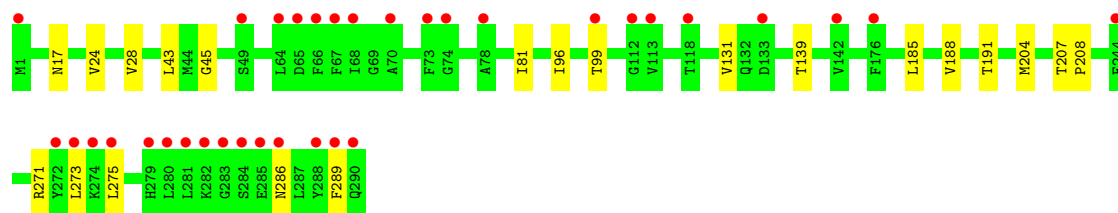


- Molecule 2: Manganese ABC transporter, permease protein

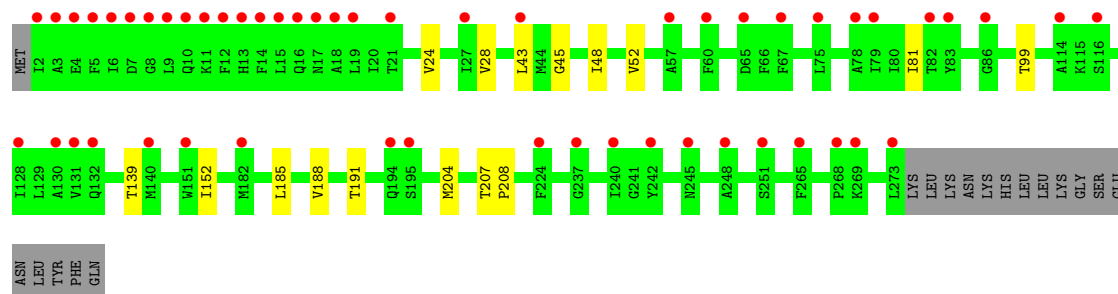
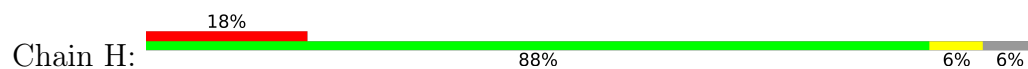


- Molecule 2: Manganese ABC transporter, permease protein

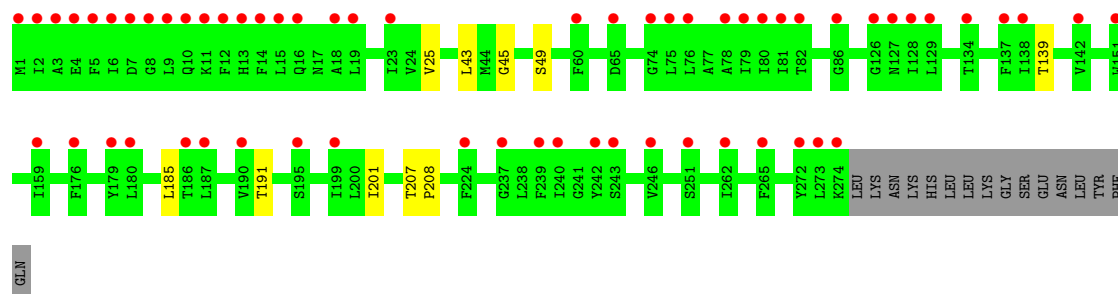
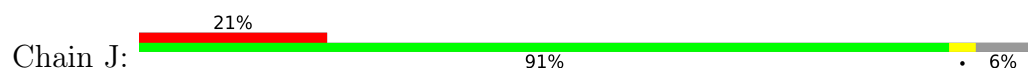




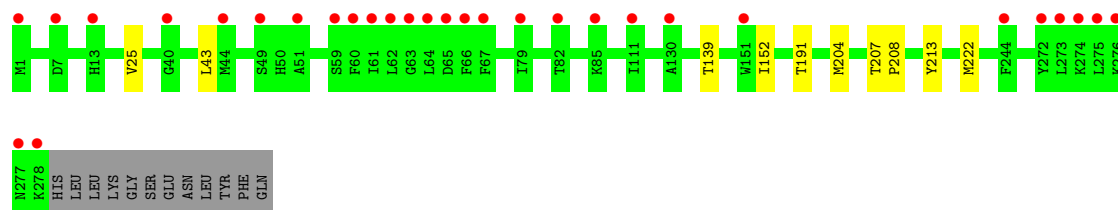
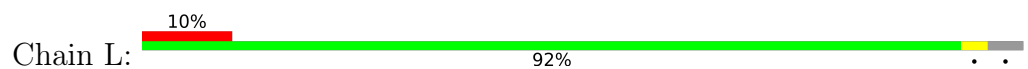
- Molecule 2: Manganese ABC transporter, permease protein



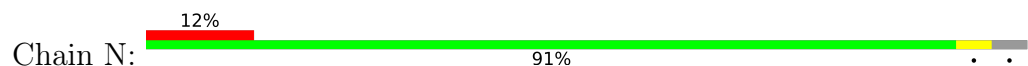
- Molecule 2: Manganese ABC transporter, permease protein

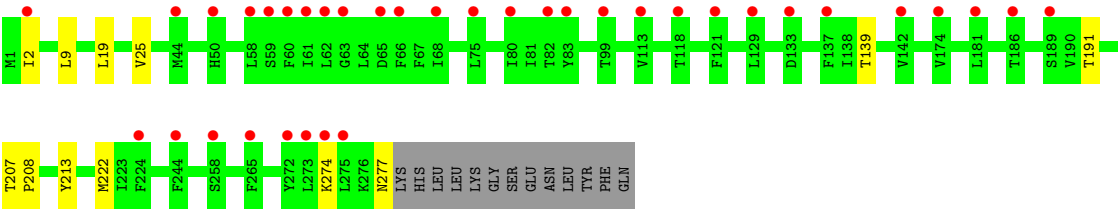


- Molecule 2: Manganese ABC transporter, permease protein

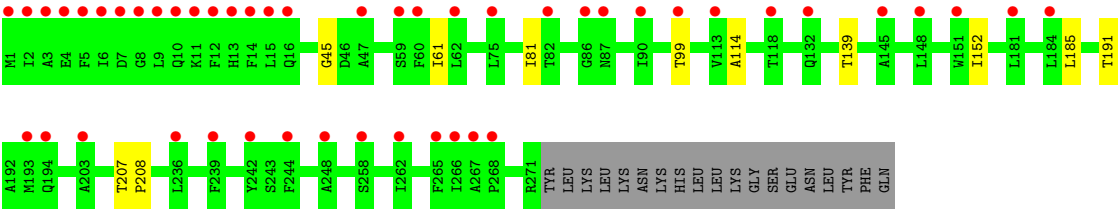
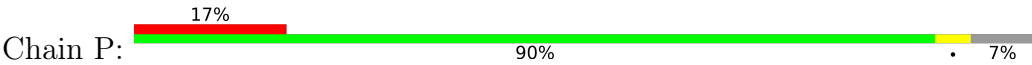


- Molecule 2: Manganese ABC transporter, permease protein





● Molecule 2: Manganese ABC transporter, permease protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	92.55Å 129.56Å 169.62Å 95.35° 92.21° 98.66°	Depositor
Resolution (Å)	49.63 – 2.90 49.63 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.8 (49.63-2.90) 90.9 (49.63-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.260 , 0.281 0.275 , 0.294	Depositor DCC
R_{free} test set	7870 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	55.4	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 13.2	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	0.86	EDS
Total number of atoms	31187	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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: PO4, MA4

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	0.38	0/1876	0.64	0/2539
1	C	0.39	0/1828	0.65	0/2476
1	E	0.39	0/1872	0.65	0/2535
1	G	0.39	0/1808	0.65	0/2450
1	I	0.39	0/1865	0.64	0/2523
1	K	0.38	0/1885	0.63	0/2550
1	M	0.38	0/1850	0.64	0/2503
1	O	0.39	0/1846	0.64	0/2500
2	B	0.39	0/2120	0.75	0/2878
2	D	0.39	0/2067	0.76	0/2809
2	F	0.39	0/2215	0.76	0/3007
2	H	0.39	0/2052	0.75	0/2792
2	J	0.40	0/2054	0.75	0/2796
2	L	0.39	0/2104	0.76	0/2859
2	N	0.39	0/2099	0.76	0/2852
2	P	0.40	0/2054	0.76	0/2791
All	All	0.39	0/31595	0.70	0/42860

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

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	1839	0	1876	6	0
1	C	1793	0	1824	4	0
1	E	1835	0	1865	6	0
1	G	1775	0	1806	7	0
1	I	1831	0	1877	4	0
1	K	1850	0	1892	9	0
1	M	1815	0	1863	5	0
1	O	1810	0	1844	6	0
2	B	2080	0	2206	6	0
2	D	2027	0	2139	12	0
2	F	2173	0	2289	13	0
2	H	2012	0	2115	12	0
2	J	2015	0	2103	6	0
2	L	2065	0	2177	7	0
2	N	2060	0	2175	7	0
2	P	2015	0	2130	6	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0
3	E	5	0	0	0	0
3	G	5	0	0	0	0
3	I	5	0	0	0	0
3	K	5	0	0	1	0
3	M	5	0	0	0	0
3	O	5	0	0	0	0
4	B	35	0	44	0	0
4	F	35	0	44	0	0
4	L	35	0	44	0	0
4	N	35	0	44	0	0
5	A	2	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	3	0	0	0	0
5	H	1	0	0	0	0
5	O	3	0	0	0	0
5	P	1	0	0	0	0
All	All	31187	0	32357	111	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 2.

All (111) 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)
2:D:139:THR:HG23	2:D:191:THR:HG23	1.72	0.69
2:B:25:VAL:HG22	2:B:191:THR:HG22	1.75	0.68
1:E:59:LEU:HD11	1:E:71:ILE:HD11	1.75	0.68
2:D:189:SER:HA	2:D:201:ILE:HD13	1.79	0.64
1:A:4:ILE:HD11	1:A:47:MET:HE2	1.82	0.62
2:J:139:THR:HG23	2:J:191:THR:HG23	1.81	0.61
2:B:139:THR:HG23	2:B:191:THR:HG23	1.80	0.61
1:M:4:ILE:HD11	1:M:47:MET:HE2	1.84	0.60
2:H:139:THR:HG23	2:H:191:THR:HG23	1.85	0.58
2:L:204:MET:HA	2:L:204:MET:HE2	1.85	0.58
2:J:25:VAL:HG22	2:J:191:THR:HG22	1.85	0.58
2:L:139:THR:HG23	2:L:191:THR:HG23	1.86	0.58
2:F:43:LEU:HD13	2:H:43:LEU:HD13	1.86	0.57
1:E:4:ILE:HD11	1:E:47:MET:HE2	1.85	0.57
2:F:28:VAL:HG12	2:F:188:VAL:HG22	1.87	0.56
2:D:45:GLY:HA2	2:D:185:LEU:HD21	1.86	0.56
2:P:139:THR:HG23	2:P:191:THR:HG23	1.88	0.56
2:J:43:LEU:HD13	2:L:43:LEU:HD13	1.88	0.56
2:N:25:VAL:HG22	2:N:191:THR:HG22	1.89	0.55
2:L:25:VAL:HG22	2:L:191:THR:HG22	1.88	0.55
1:K:59:LEU:HD11	1:K:71:ILE:HD11	1.89	0.54
2:F:139:THR:HG23	2:F:191:THR:CG2	2.39	0.53
2:F:24:VAL:HG11	2:F:139:THR:HG22	1.91	0.52
1:M:176:LEU:HD13	1:M:188:ILE:HD11	1.92	0.51
2:J:45:GLY:HA2	2:J:185:LEU:HD21	1.93	0.51
2:F:81:ILE:HG13	2:F:99:THR:HG21	1.92	0.50
2:P:61:ILE:HG22	2:P:114:ALA:HB3	1.93	0.50
1:I:4:ILE:HD11	1:I:47:MET:HE2	1.94	0.49
2:F:28:VAL:CG1	2:F:188:VAL:HG22	2.42	0.49
1:G:4:ILE:HD11	1:G:47:MET:HE2	1.95	0.49
2:D:207:THR:HB	2:D:208:PRO:HD3	1.95	0.49
2:N:139:THR:HG23	2:N:191:THR:HG23	1.93	0.49
2:H:81:ILE:HG13	2:H:99:THR:HG21	1.94	0.49
1:M:48:LEU:HD21	1:M:71:ILE:HG21	1.94	0.48
2:D:78:ALA:HB2	2:D:182:MET:HE1	1.95	0.48
1:O:197:PRO:HA	1:O:216:THR:HG21	1.95	0.48
1:I:40:LYS:HB3	1:I:189:VAL:HG13	1.95	0.48
2:D:25:VAL:HG22	2:D:191:THR:HG22	1.96	0.47
1:O:88:VAL:CG2	1:O:134:LEU:HD11	2.44	0.47
2:F:275:LEU:HD13	2:F:289:PHE:CE2	2.50	0.47
2:H:45:GLY:HA2	2:H:185:LEU:HD21	1.95	0.47
2:N:213:TYR:HB2	2:N:222:MET:HE3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:40:LYS:HB3	1:K:189:VAL:HG13	1.96	0.47
1:A:176:LEU:HD13	1:A:188:ILE:HD11	1.95	0.47
1:K:176:LEU:HD23	1:K:179:LEU:HD12	1.97	0.46
2:B:204:MET:HE2	2:B:204:MET:HA	1.97	0.46
2:D:204:MET:HA	2:D:204:MET:HE2	1.97	0.46
2:H:204:MET:HA	2:H:204:MET:HE2	1.97	0.46
1:K:4:ILE:HD11	1:K:47:MET:HE2	1.97	0.46
1:A:59:LEU:HG	1:A:64:VAL:HG22	1.96	0.46
1:A:59:LEU:HD11	1:A:71:ILE:HD11	1.96	0.46
1:C:176:LEU:HD13	1:C:188:ILE:HD11	1.98	0.46
1:O:4:ILE:HD11	1:O:47:MET:HE2	1.98	0.46
1:K:42:THR:N	3:K:301:PO4:O3	2.46	0.45
2:F:204:MET:HA	2:F:204:MET:HE2	1.98	0.45
1:G:103:PHE:CE1	2:H:152:ILE:HD12	2.51	0.45
2:D:198:THR:HA	2:D:201:ILE:HD11	1.99	0.45
1:G:40:LYS:HB3	1:G:189:VAL:HG13	1.98	0.45
1:A:40:LYS:HB3	1:A:189:VAL:HG13	1.97	0.45
2:F:207:THR:HB	2:F:208:PRO:HD3	1.98	0.45
2:D:49:SER:HB3	2:D:201:ILE:HD12	1.98	0.45
2:H:139:THR:HG23	2:H:191:THR:CG2	2.46	0.45
1:I:74:VAL:HB	1:I:157:LEU:HD23	1.99	0.45
1:E:59:LEU:HG	1:E:64:VAL:HG22	1.98	0.44
2:P:81:ILE:HG13	2:P:99:THR:HG21	2.00	0.44
2:D:81:ILE:CG1	2:D:99:THR:HG21	2.48	0.44
2:D:78:ALA:CB	2:D:182:MET:HE1	2.48	0.44
1:G:44:LEU:HD11	1:G:189:VAL:HG23	1.99	0.44
1:C:40:LYS:HB3	1:C:189:VAL:HG13	2.00	0.43
2:B:213:TYR:HB2	2:B:222:MET:HE3	2.00	0.43
2:N:9:LEU:HD11	2:N:19:LEU:HD22	2.01	0.43
2:F:17:ASN:HB3	2:F:131:VAL:HG23	1.99	0.43
2:H:81:ILE:CG1	2:H:99:THR:HG21	2.49	0.43
1:E:48:LEU:HD21	1:E:71:ILE:HG21	2.00	0.43
2:P:207:THR:HB	2:P:208:PRO:HD3	2.01	0.43
2:H:48:ILE:O	2:H:52:VAL:HG23	2.19	0.43
1:A:48:LEU:HD21	1:A:71:ILE:HG21	2.00	0.42
2:D:81:ILE:HG13	2:D:99:THR:HG21	2.00	0.42
2:H:24:VAL:HG11	2:H:139:THR:HG22	2.00	0.42
2:J:49:SER:HB3	2:J:201:ILE:HD12	2.01	0.42
1:E:88:VAL:HG22	1:E:134:LEU:HD11	2.00	0.42
1:G:176:LEU:HD13	1:G:188:ILE:HD11	2.01	0.42
2:F:45:GLY:HA2	2:F:185:LEU:HD21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:274:LYS:HA	2:N:277:ASN:HD22	1.83	0.42
2:F:43:LEU:HG	2:F:96:ILE:HG21	2.01	0.42
2:H:28:VAL:HG12	2:H:188:VAL:HG22	2.00	0.42
1:M:121:VAL:O	1:M:141:ARG:NE	2.52	0.42
1:K:103:PHE:CE1	2:L:152:ILE:HD12	2.55	0.42
2:H:207:THR:HB	2:H:208:PRO:HD3	2.02	0.42
2:P:45:GLY:HA2	2:P:185:LEU:HD21	2.01	0.42
1:E:88:VAL:CG2	1:E:134:LEU:HD11	2.50	0.42
1:G:59:LEU:CD1	1:G:71:ILE:HD11	2.50	0.42
1:M:123:LEU:HD11	1:M:141:ARG:HB2	2.02	0.42
2:B:207:THR:HB	2:B:208:PRO:HD3	2.02	0.42
2:L:213:TYR:HB2	2:L:222:MET:HE3	2.01	0.42
2:N:2:ILE:HD12	2:N:2:ILE:N	2.35	0.42
1:O:103:PHE:CE1	2:P:152:ILE:HD12	2.55	0.42
1:I:97:PHE:N	1:I:98:PRO:CD	2.83	0.41
1:K:2:ILE:HD12	1:K:24:LEU:HD23	2.01	0.41
1:K:17:LEU:HD21	1:K:43:LEU:HA	2.01	0.41
2:B:81:ILE:HG13	2:B:99:THR:HG21	2.02	0.41
2:L:207:THR:HB	2:L:208:PRO:HD3	2.01	0.41
1:C:59:LEU:HG	1:C:64:VAL:HG22	2.02	0.41
2:J:207:THR:HB	2:J:208:PRO:HD3	2.02	0.41
2:F:271:ARG:C	2:F:273:LEU:H	2.29	0.41
1:C:123:LEU:HD11	1:C:141:ARG:HB2	2.03	0.41
1:K:48:LEU:CD2	1:K:71:ILE:HD13	2.50	0.41
2:N:207:THR:HB	2:N:208:PRO:HD3	2.02	0.41
1:O:176:LEU:HD13	1:O:188:ILE:HD11	2.02	0.41
1:O:59:LEU:HD22	1:O:71:ILE:HD11	2.02	0.40
1:G:48:LEU:HD21	1:G:71:ILE:HG21	2.02	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	235/240 (98%)	218 (93%)	17 (7%)	0	100	100
1	C	231/240 (96%)	218 (94%)	12 (5%)	1 (0%)	30	59
1	E	235/240 (98%)	221 (94%)	13 (6%)	1 (0%)	30	59
1	G	228/240 (95%)	216 (95%)	12 (5%)	0	100	100
1	I	232/240 (97%)	216 (93%)	15 (6%)	1 (0%)	30	59
1	K	234/240 (98%)	221 (94%)	12 (5%)	1 (0%)	30	59
1	M	232/240 (97%)	218 (94%)	13 (6%)	1 (0%)	30	59
1	O	233/240 (97%)	215 (92%)	15 (6%)	3 (1%)	9	32
2	B	276/290 (95%)	266 (96%)	9 (3%)	1 (0%)	30	59
2	D	270/290 (93%)	261 (97%)	9 (3%)	0	100	100
2	F	288/290 (99%)	281 (98%)	7 (2%)	0	100	100
2	H	270/290 (93%)	265 (98%)	5 (2%)	0	100	100
2	J	272/290 (94%)	264 (97%)	8 (3%)	0	100	100
2	L	276/290 (95%)	268 (97%)	8 (3%)	0	100	100
2	N	275/290 (95%)	267 (97%)	8 (3%)	0	100	100
2	P	269/290 (93%)	260 (97%)	9 (3%)	0	100	100
All	All	4056/4240 (96%)	3875 (96%)	172 (4%)	9 (0%)	43	72

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	231	ASN
1	M	40	LYS
1	E	40	LYS
1	K	231	ASN
1	O	232	GLN
1	O	231	ASN
2	B	127	ASN
1	C	12	LYS
1	O	209	GLU

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	199/208 (96%)	199 (100%)	0	100	100
1	C	193/208 (93%)	193 (100%)	0	100	100
1	E	198/208 (95%)	195 (98%)	3 (2%)	57	84
1	G	191/208 (92%)	190 (100%)	1 (0%)	81	93
1	I	199/208 (96%)	199 (100%)	0	100	100
1	K	201/208 (97%)	200 (100%)	1 (0%)	81	93
1	M	197/208 (95%)	196 (100%)	1 (0%)	81	93
1	O	195/208 (94%)	194 (100%)	1 (0%)	81	93
2	B	221/234 (94%)	220 (100%)	1 (0%)	81	93
2	D	215/234 (92%)	215 (100%)	0	100	100
2	F	230/234 (98%)	229 (100%)	1 (0%)	84	94
2	H	211/234 (90%)	211 (100%)	0	100	100
2	J	210/234 (90%)	210 (100%)	0	100	100
2	L	218/234 (93%)	218 (100%)	0	100	100
2	N	218/234 (93%)	218 (100%)	0	100	100
2	P	214/234 (92%)	214 (100%)	0	100	100
All	All	3310/3536 (94%)	3301 (100%)	9 (0%)	86	96

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	11	LYS
1	E	22	LEU
1	E	140	GLN
1	E	147	CYS
2	F	286	ASN
1	G	22	LEU
1	K	147	CYS
1	M	22	LEU
1	O	22	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN

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Mol	Chain	Res	Type
2	B	87	ASN
2	B	122	HIS
2	B	127	ASN
1	C	36	ASN
1	C	56	GLN
1	C	110	HIS
1	C	130	GLN
1	C	140	GLN
1	C	174	ASN
1	E	6	ASN
1	E	36	ASN
1	E	130	GLN
1	E	140	GLN
1	E	198	HIS
2	F	10	GLN
2	F	194	GLN
1	G	36	ASN
1	G	133	GLN
2	J	10	GLN
1	K	140	GLN
1	M	6	ASN
1	M	138	GLN
1	M	202	GLN
2	N	122	HIS
2	N	127	ASN
2	N	277	ASN
1	O	110	HIS
1	O	133	GLN

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

12 ligands are modelled in this entry.

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
3	PO4	C	301	-	4,4,4	0.94	0	6,6,6	0.57	0
3	PO4	K	301	-	4,4,4	1.00	0	6,6,6	0.40	0
3	PO4	A	301	-	4,4,4	0.99	0	6,6,6	0.43	0
3	PO4	G	301	-	4,4,4	0.99	0	6,6,6	0.49	0
3	PO4	E	301	-	4,4,4	0.93	0	6,6,6	0.65	0
3	PO4	O	301	-	4,4,4	1.00	0	6,6,6	0.47	0
3	PO4	I	301	-	4,4,4	0.94	0	6,6,6	0.54	0
4	MA4	N	301	-	37,37,37	0.47	0	50,50,50	0.79	2 (4%)
3	PO4	M	301	-	4,4,4	1.02	0	6,6,6	0.45	0
4	MA4	B	301	-	37,37,37	0.47	0	50,50,50	0.95	2 (4%)
4	MA4	L	301	-	37,37,37	0.44	0	50,50,50	0.63	0
4	MA4	F	301	-	37,37,37	0.47	0	50,50,50	0.67	0

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	MA4	N	301	-	-	3/18/66/66	0/3/3/3
4	MA4	L	301	-	-	8/18/66/66	0/3/3/3
4	MA4	F	301	-	-	6/18/66/66	0/3/3/3
4	MA4	B	301	-	-	3/18/66/66	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	MA4	C20-C30-C40	3.53	117.70	109.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	MA4	C10-C20-C30	3.39	117.13	110.01
4	N	301	MA4	C1-O5-C5	2.42	118.45	113.72
4	N	301	MA4	C3-C4-C5	2.07	113.98	110.23

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	N	301	MA4	C4-C5-C6-O6
4	N	301	MA4	O5-C5-C6-O6
4	F	301	MA4	O5-C1-O1-C40
4	F	301	MA4	C50-C40-O1-C1
4	L	301	MA4	O5-C1-O1-C40
4	F	301	MA4	C30-C40-O1-C1
4	B	301	MA4	C21-C31-C41-C51
4	F	301	MA4	C21-C11-O10-C10
4	F	301	MA4	C11-C21-C31-C41
4	N	301	MA4	C21-C31-C41-C51
4	B	301	MA4	C30-C40-O1-C1
4	L	301	MA4	C30-C40-O1-C1
4	L	301	MA4	C50-C40-O1-C1
4	L	301	MA4	C21-C11-O10-C10
4	B	301	MA4	C50-C40-O1-C1
4	L	301	MA4	C22-C12-C61-C51
4	L	301	MA4	C62-C12-C61-C51
4	L	301	MA4	C41-C51-C61-C12
4	F	301	MA4	C41-C51-C61-C12
4	L	301	MA4	C21-C31-C41-C51

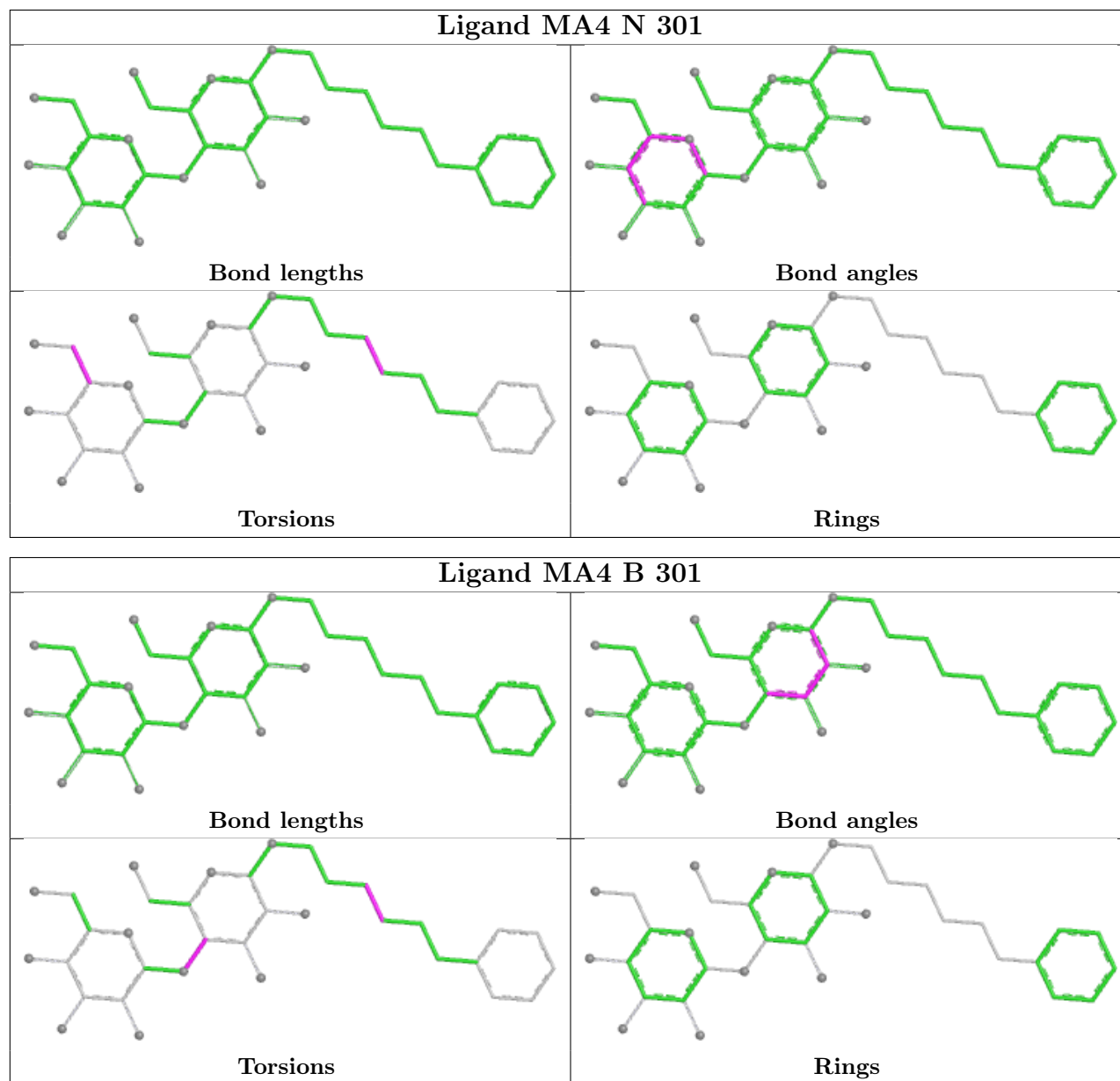
There are no ring outliers.

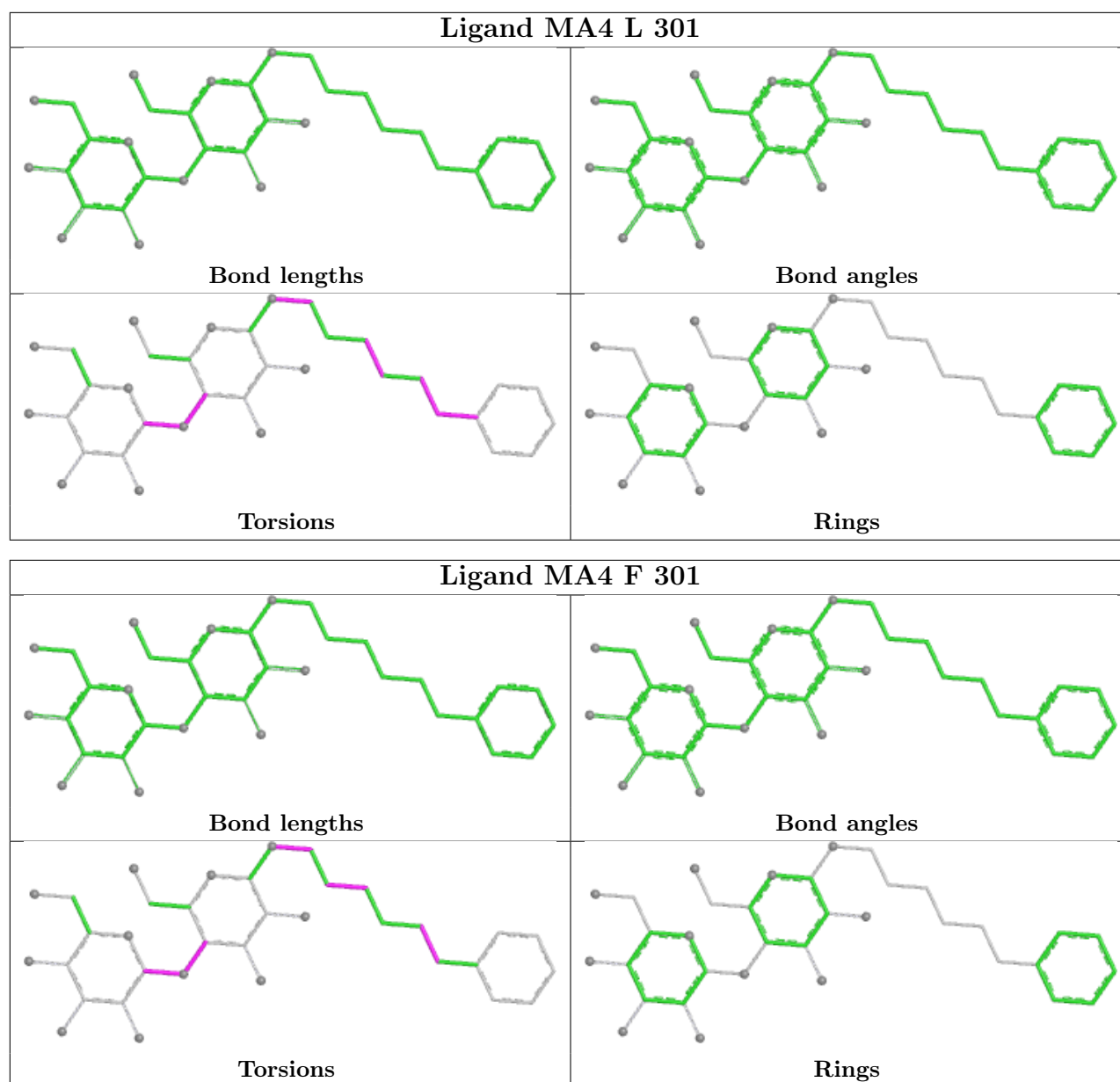
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	K	301	PO4	1	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 [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	236/240 (98%)	0.95	21 (8%)	15 13	34, 40, 53, 55	1 (0%)
1	C	232/240 (96%)	0.93	23 (9%)	13 11	29, 34, 46, 49	1 (0%)
1	E	236/240 (98%)	0.77	11 (4%)	36 28	27, 33, 43, 48	1 (0%)
1	G	230/240 (95%)	0.69	14 (6%)	27 21	24, 33, 41, 45	0
1	I	234/240 (97%)	0.76	16 (6%)	23 18	24, 32, 42, 50	0
1	K	236/240 (98%)	0.84	19 (8%)	18 15	30, 37, 45, 64	0
1	M	233/240 (97%)	1.11	30 (12%)	7 6	31, 42, 51, 57	1 (0%)
1	O	234/240 (97%)	0.79	13 (5%)	30 23	26, 34, 43, 46	1 (0%)
2	B	278/290 (95%)	0.82	24 (8%)	16 13	27, 38, 51, 69	0
2	D	272/290 (93%)	0.93	26 (9%)	13 11	30, 39, 51, 69	0
2	F	290/290 (100%)	1.02	34 (11%)	9 8	29, 41, 61, 79	0
2	H	272/290 (93%)	1.28	53 (19%)	3 2	32, 44, 59, 80	0
2	J	274/290 (94%)	1.48	61 (22%)	2 2	29, 45, 66, 84	0
2	L	278/290 (95%)	0.93	30 (10%)	11 9	31, 40, 62, 77	0
2	N	277/290 (95%)	1.01	36 (12%)	7 6	34, 43, 59, 75	0
2	P	271/290 (93%)	1.34	48 (17%)	4 3	32, 48, 59, 107	0
All	All	4083/4240 (96%)	0.99	459 (11%)	10 9	24, 39, 56, 107	5 (0%)

All (459) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	232	GLN	9.6
2	J	7	ASP	8.8
2	J	14	PHE	8.5
2	J	12	PHE	8.2
2	J	10	GLN	8.1

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Mol	Chain	Res	Type	RSRZ
2	J	11	LYS	8.0
2	P	7	ASP	7.9
2	P	12	PHE	7.8
1	C	13	GLU	7.0
1	I	14	THR	7.0
1	O	14	THR	6.9
2	P	1	MET	6.8
2	N	61	ILE	6.8
2	H	14	PHE	6.3
2	J	9	LEU	6.2
2	H	11	LYS	6.2
1	O	19	ASP	6.0
2	H	10	GLN	5.9
2	P	265	PHE	5.9
2	P	4	GLU	5.9
2	P	14	PHE	5.9
2	J	6	ILE	5.9
2	J	16	GLN	5.8
1	C	14	THR	5.8
1	I	13	GLU	5.7
2	N	59	SER	5.6
2	J	15	LEU	5.5
2	H	2	ILE	5.5
2	J	5	PHE	5.3
2	F	65	ASP	5.3
2	H	12	PHE	5.3
2	H	8	GLY	5.3
2	H	7	ASP	5.1
2	F	284	SER	5.1
2	F	283	GLY	4.9
2	D	7	ASP	4.9
2	J	242	TYR	4.8
2	H	13	HIS	4.8
2	J	13	HIS	4.7
2	H	9	LEU	4.7
2	H	4	GLU	4.7
2	L	61	ILE	4.7
2	L	65	ASP	4.6
2	L	66	PHE	4.6
1	M	109	LYS	4.6
2	J	274	LYS	4.6
2	P	6	ILE	4.6

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Mol	Chain	Res	Type	RSRZ
2	L	51	ALA	4.5
2	F	282	LYS	4.4
2	L	273	LEU	4.4
1	C	103	PHE	4.4
1	C	12	LYS	4.4
1	O	11[A]	TYR	4.4
2	D	14	PHE	4.4
2	B	277	ASN	4.4
2	P	86	GLY	4.4
2	P	10	GLN	4.4
1	M	82	TYR	4.3
2	H	265	PHE	4.3
2	H	237	GLY	4.3
2	N	274	LYS	4.3
1	A	145	ALA	4.3
2	L	244	PHE	4.3
2	N	272	TYR	4.2
2	F	275	LEU	4.2
2	J	127	ASN	4.2
2	N	244	PHE	4.2
2	J	251	SER	4.1
2	B	66	PHE	4.1
1	C	230	GLY	4.1
1	K	46	GLY	4.0
2	J	243	SER	4.0
2	F	67	PHE	4.0
2	B	61	ILE	4.0
2	J	81	ILE	4.0
2	H	75	LEU	4.0
2	B	274	LYS	4.0
2	P	258	SER	4.0
1	I	82	TYR	3.9
2	J	179	TYR	3.9
2	D	11	LYS	3.9
1	O	10	SER	3.9
2	F	281	LEU	3.9
2	J	79	ILE	3.9
2	J	82	THR	3.9
2	N	65	ASP	3.9
2	J	239	PHE	3.9
2	F	274	LYS	3.9
1	C	11[A]	TYR	3.9

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Mol	Chain	Res	Type	RSRZ
2	J	78	ALA	3.9
1	C	10	SER	3.9
1	C	82	TYR	3.8
2	P	267	ALA	3.8
2	L	274	LYS	3.8
2	F	279	HIS	3.8
1	E	79	ASN	3.8
1	I	79	ASN	3.8
2	B	75	LEU	3.7
2	N	62	LEU	3.7
2	H	251	SER	3.7
2	P	11	LYS	3.7
2	B	65	ASP	3.7
2	L	272	TYR	3.7
1	A	13	GLU	3.7
2	J	128	ILE	3.7
1	I	11	TYR	3.7
1	M	11[A]	TYR	3.7
1	M	102	LEU	3.7
2	B	73	PHE	3.7
2	N	60	PHE	3.7
1	C	15	LEU	3.6
1	G	47	MET	3.6
2	F	290	GLN	3.6
1	E	103	PHE	3.6
1	K	11	TYR	3.6
2	N	66	PHE	3.6
1	M	105	SER	3.6
2	P	2	ILE	3.6
1	K	82	TYR	3.6
2	F	273	LEU	3.6
2	L	59	SER	3.5
2	H	5	PHE	3.5
1	M	218	GLU	3.5
2	P	13	HIS	3.5
2	P	5	PHE	3.5
2	F	113	VAL	3.5
1	G	14	THR	3.5
1	G	82	TYR	3.5
2	F	272	TYR	3.5
2	J	3	ALA	3.5
2	J	2	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	231	ASN	3.4
2	J	265	PHE	3.4
2	D	12	PHE	3.4
2	L	1	MET	3.4
1	A	71	ILE	3.4
2	B	273	LEU	3.4
2	F	285	GLU	3.4
2	J	4	GLU	3.4
1	O	13	GLU	3.3
1	O	82	TYR	3.3
1	M	13	GLU	3.3
1	A	102	LEU	3.3
2	L	67	PHE	3.3
1	A	61	ASP	3.3
2	H	240	ILE	3.3
2	P	9	LEU	3.3
2	N	113	VAL	3.3
2	D	126	GLY	3.3
2	L	63	GLY	3.3
1	C	9	VAL	3.2
2	H	16	GLN	3.2
1	I	80	ILE	3.2
2	H	194	GLN	3.2
2	H	86	GLY	3.2
2	B	272	TYR	3.2
2	J	129	LEU	3.2
1	O	103	PHE	3.2
2	H	79	ILE	3.2
1	M	217	LYS	3.2
2	P	82	THR	3.2
1	C	209	GLU	3.2
1	M	110	HIS	3.1
2	B	81	ILE	3.1
2	N	273	LEU	3.1
2	P	203	ALA	3.1
2	J	23	ILE	3.1
2	L	79	ILE	3.1
2	N	58	LEU	3.1
2	B	278	LYS	3.1
2	H	116	SER	3.1
1	A	47	MET	3.1
2	F	286	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
2	J	65	ASP	3.1
1	E	218	GLU	3.1
2	L	278	LYS	3.1
1	A	1	MET	3.0
2	J	1	MET	3.0
1	A	56	GLN	3.0
2	P	16	GLN	3.0
1	O	208	ARG	3.0
2	B	275	LEU	3.0
2	P	15	LEU	3.0
1	O	47	MET	3.0
1	A	10	SER	3.0
1	A	15	LEU	3.0
2	P	75	LEU	3.0
2	B	134	THR	3.0
2	F	99	THR	3.0
2	H	82	THR	3.0
2	B	51	ALA	3.0
2	F	66	PHE	3.0
2	B	103	PHE	3.0
1	G	12	LYS	3.0
1	K	230	GLY	3.0
1	K	109	LYS	2.9
2	F	112	GLY	2.9
2	H	43	LEU	2.9
1	G	61	ASP	2.9
2	P	268	PRO	2.9
1	M	14	THR	2.9
2	H	83	TYR	2.9
2	H	242	TYR	2.9
1	M	12	LYS	2.9
2	H	195	SER	2.9
1	M	103	PHE	2.9
1	A	14	THR	2.9
2	F	68	ILE	2.9
2	F	74	GLY	2.9
2	J	74	GLY	2.9
2	H	78	ALA	2.9
2	L	60	PHE	2.9
2	N	258	SER	2.9
1	K	232	GLN	2.8
1	E	13	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	67	PHE	2.8
2	P	148	LEU	2.8
1	K	233	LEU	2.8
1	M	62	LYS	2.8
1	I	147	CYS	2.8
2	F	289	PHE	2.8
2	J	137	PHE	2.8
2	J	224	PHE	2.8
2	P	242	TYR	2.8
2	L	13	HIS	2.8
2	H	130	ALA	2.7
2	L	64	LEU	2.7
2	B	271	ARG	2.7
1	I	233	LEU	2.7
2	P	3	ALA	2.7
2	L	276	LYS	2.7
1	C	6	ASN	2.7
1	C	178	ASP	2.7
2	J	126	GLY	2.7
2	D	193	MET	2.7
2	D	272	TYR	2.7
2	N	118	THR	2.7
1	M	125	ASP	2.7
1	A	21	SER	2.7
1	M	63	GLU	2.7
2	H	131	VAL	2.7
2	D	15	LEU	2.7
2	L	62	LEU	2.7
2	F	70	ALA	2.7
2	P	60	PHE	2.7
1	C	79	ASN	2.7
1	M	79	ASN	2.7
2	D	74	GLY	2.7
2	N	63	GLY	2.7
1	I	19	ASP	2.7
2	D	265	PHE	2.6
2	H	248	ALA	2.6
1	A	69	HIS	2.6
1	M	104	ARG	2.6
2	N	129	LEU	2.6
2	P	262	ILE	2.6
1	G	10	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	J	18	ALA	2.6
2	J	60	PHE	2.6
2	H	268	PRO	2.6
1	E	14	THR	2.6
2	P	118	THR	2.6
2	H	132	GLN	2.6
1	M	101	PRO	2.6
2	J	237	GLY	2.6
1	O	15	LEU	2.6
2	J	75	LEU	2.6
1	O	12	LYS	2.6
2	H	15	LEU	2.6
2	P	194	GLN	2.6
2	H	224	PHE	2.6
1	M	41	SER	2.5
2	F	280	LEU	2.5
2	J	240	ILE	2.5
2	P	266	ILE	2.5
1	G	13	GLU	2.5
2	F	244	PHE	2.5
2	P	248	ALA	2.5
1	E	61	ASP	2.5
2	B	126	GLY	2.5
2	P	236	LEU	2.5
2	D	27	ILE	2.5
2	D	176	PHE	2.5
2	F	133	ASP	2.5
2	J	195	SER	2.5
1	M	233	LEU	2.5
2	F	1	MET	2.5
1	I	78	ILE	2.5
1	M	167	VAL	2.5
2	L	151	TRP	2.5
2	J	134	THR	2.5
1	A	147	CYS	2.5
2	B	276	LYS	2.5
2	N	75	LEU	2.5
2	B	72	VAL	2.5
1	C	19	ASP	2.4
1	I	15	LEU	2.4
2	L	44	MET	2.4
1	K	105	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	M	67	SER	2.4
2	P	113	VAL	2.4
2	F	176	PHE	2.4
1	A	11[A]	TYR	2.4
1	G	54	GLN	2.4
2	J	272	TYR	2.4
1	I	231	ASN	2.4
2	L	275	LEU	2.4
2	J	151	TRP	2.4
2	F	288	TYR	2.4
2	N	275	LEU	2.4
2	N	2	ILE	2.4
1	M	64	VAL	2.4
1	K	10	SER	2.4
1	M	57	ALA	2.4
2	H	114	ALA	2.4
2	P	151	TRP	2.4
1	G	1	MET	2.4
2	J	19	LEU	2.4
2	J	273	LEU	2.4
2	H	6	ILE	2.4
2	H	245	ASN	2.4
2	H	269	LYS	2.4
1	A	67	SER	2.4
2	F	142	VAL	2.3
2	P	8	GLY	2.3
1	G	125	ASP	2.3
2	H	67	PHE	2.3
2	D	16	GLN	2.3
2	P	132	GLN	2.3
2	B	1	MET	2.3
2	D	49	SER	2.3
1	A	198	HIS	2.3
2	H	273	LEU	2.3
1	E	50	ILE	2.3
2	J	199	ILE	2.3
2	D	239	PHE	2.3
2	F	118	THR	2.3
2	H	27	ILE	2.3
2	P	90	ILE	2.3
2	B	63	GLY	2.3
1	K	16	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	M	124	ALA	2.3
1	O	218	GLU	2.3
2	L	130	ALA	2.3
1	K	133	GLN	2.3
2	P	62	LEU	2.3
1	K	234	PHE	2.3
2	H	60	PHE	2.3
2	P	47	ALA	2.3
1	A	68	LEU	2.3
1	K	102	LEU	2.3
2	L	7	ASP	2.3
2	J	262	ILE	2.3
2	H	21	THR	2.3
2	L	49	SER	2.3
2	N	174	VAL	2.3
2	J	176	PHE	2.2
1	C	1	MET	2.2
2	P	193	MET	2.2
1	K	108	ALA	2.2
1	M	116	GLU	2.2
2	F	78	ALA	2.2
1	C	22	LEU	2.2
1	I	76	GLN	2.2
2	J	80	ILE	2.2
2	D	59	SER	2.2
2	D	195	SER	2.2
2	J	142	VAL	2.2
2	J	246	VAL	2.2
2	J	138	ILE	2.2
2	J	186	THR	2.2
2	J	8	GLY	2.2
2	D	224	PHE	2.2
2	J	180	LEU	2.2
2	P	145	ALA	2.2
1	E	130	GLN	2.2
1	I	208	ARG	2.2
1	M	71	ILE	2.2
2	N	50	HIS	2.2
2	J	190	VAL	2.2
2	D	191	THR	2.2
2	L	82	THR	2.2
2	N	83	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
2	P	59	SER	2.2
1	M	187	LEU	2.2
2	J	187	LEU	2.2
2	N	181	LEU	2.2
1	A	152	ALA	2.2
1	G	218	GLU	2.2
1	K	13	GLU	2.2
2	H	18	ALA	2.2
1	K	50	ILE	2.2
2	N	80	ILE	2.2
1	C	54	GLN	2.2
2	L	277	ASN	2.2
2	N	133	ASP	2.2
2	H	182	MET	2.2
2	N	82	THR	2.2
1	A	82	TYR	2.2
2	F	49	SER	2.2
2	P	244	PHE	2.2
2	H	19	LEU	2.2
2	J	76	LEU	2.2
1	A	72	ALA	2.2
2	D	4	GLU	2.2
2	L	85	LYS	2.1
1	M	26	GLY	2.1
2	L	40	GLY	2.1
2	N	99	THR	2.1
1	C	17	LEU	2.1
2	D	271	ARG	2.1
2	D	1	MET	2.1
2	H	140	MET	2.1
1	I	60	ASP	2.1
2	D	82	THR	2.1
2	J	86	GLY	2.1
1	E	17	LEU	2.1
2	B	244	PHE	2.1
2	H	57	ALA	2.1
2	L	111	ILE	2.1
2	N	68	ILE	2.1
1	C	18	LYS	2.1
1	I	36	ASN	2.1
1	O	231	ASN	2.1
2	N	44	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	103	PHE	2.1
2	N	265	PHE	2.1
1	E	116	GLU	2.1
1	K	132	SER	2.1
2	N	189	SER	2.1
1	E	133	GLN	2.1
1	M	69	HIS	2.1
1	K	236	ASN	2.1
2	P	87	ASN	2.1
2	F	64	LEU	2.1
2	D	237	GLY	2.1
2	F	73	PHE	2.1
2	N	224	PHE	2.1
2	D	198	THR	2.1
1	G	16	ALA	2.1
1	G	11	TYR	2.1
2	D	75	LEU	2.0
2	P	184	LEU	2.0
2	N	186	THR	2.0
2	P	99	THR	2.0
2	H	151	TRP	2.0
2	B	116	SER	2.0
1	C	156	LEU	2.0
2	P	181	LEU	2.0
2	H	17	ASN	2.0
2	N	121	PHE	2.0
2	N	137	PHE	2.0
2	P	239	PHE	2.0
2	H	128	ILE	2.0
2	J	159	ILE	2.0
2	H	3	ALA	2.0
2	H	65	ASP	2.0
1	K	47	MET	2.0
2	N	142	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

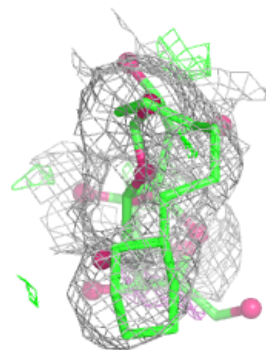
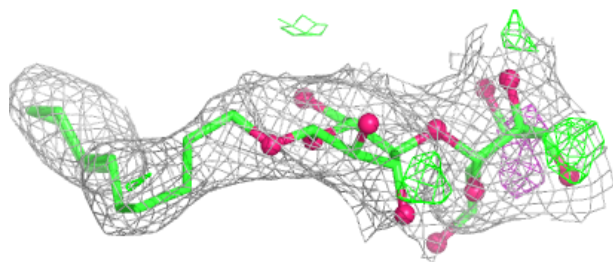
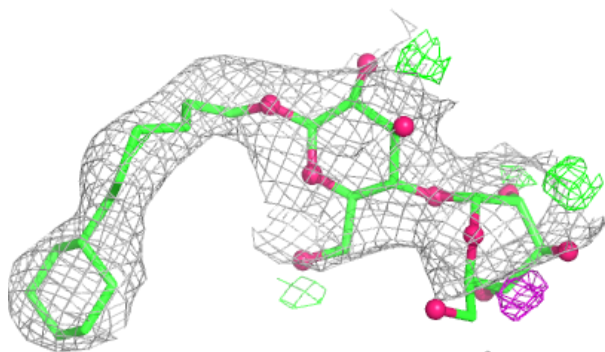
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	PO4	M	301	5/5	0.62	0.42	70,70,70,70	0
3	PO4	K	301	5/5	0.65	0.55	73,74,74,74	0
3	PO4	I	301	5/5	0.65	0.58	75,76,77,77	0
3	PO4	C	301	5/5	0.68	0.46	70,71,71,72	0
3	PO4	O	301	5/5	0.70	0.39	59,60,60,60	0
3	PO4	A	301	5/5	0.72	0.35	66,66,67,67	0
3	PO4	E	301	5/5	0.74	0.37	54,55,55,55	0
4	MA4	B	301	35/35	0.78	0.25	77,83,86,86	0
3	PO4	G	301	5/5	0.79	0.30	60,60,61,61	0
4	MA4	F	301	35/35	0.82	0.21	69,69,72,72	0
4	MA4	L	301	35/35	0.82	0.20	71,73,75,76	0
4	MA4	N	301	35/35	0.84	0.24	88,92,95,95	0

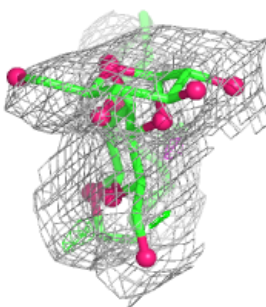
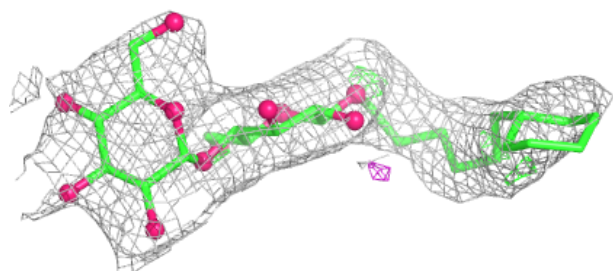
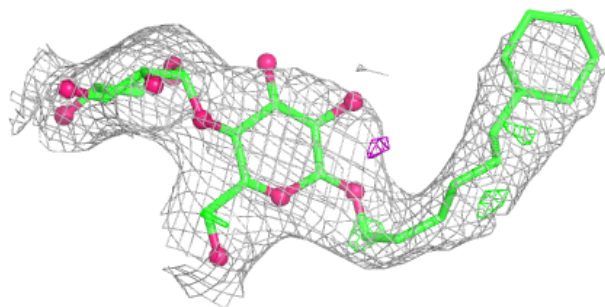
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around MA4 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)

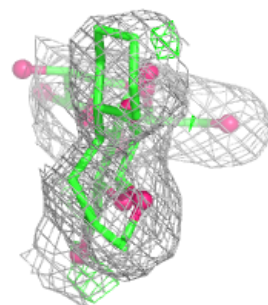
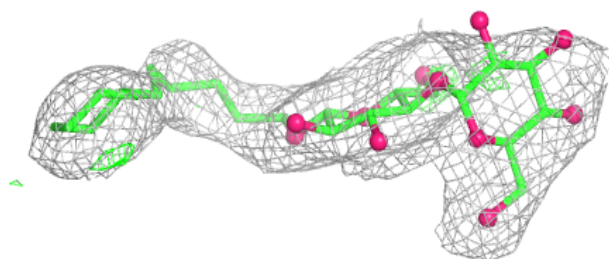
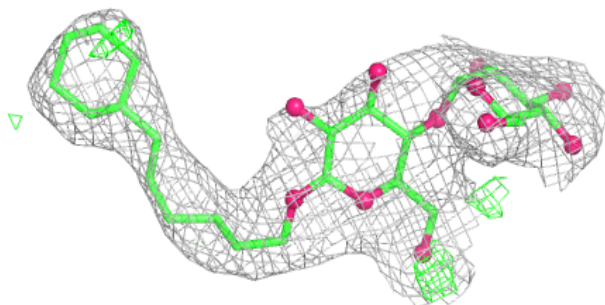
**Electron density around MA4 F 301:**

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

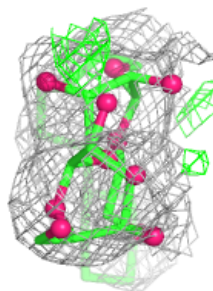
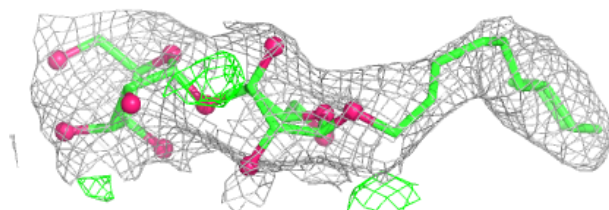
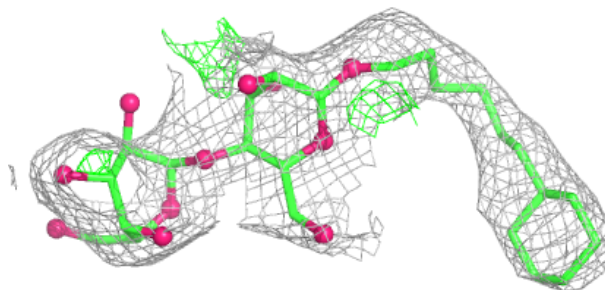


Electron density around MA4 L 301:

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

**Electron density around MA4 N 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 [i](#)

There are no such residues in this entry.