



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2026 – 12:16 PM JST

PDB ID : 9V5Y / pdb_00009v5y
Title : PfDXR - Mn2+ - SHOK314 ternary complex
Authors : Takada, S.; Sakamoto, Y.; Tanaka, N.
Deposited on : 2025-05-26
Resolution : 1.57 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 : 1.8.5 (274361), CSD as541be (2020)
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

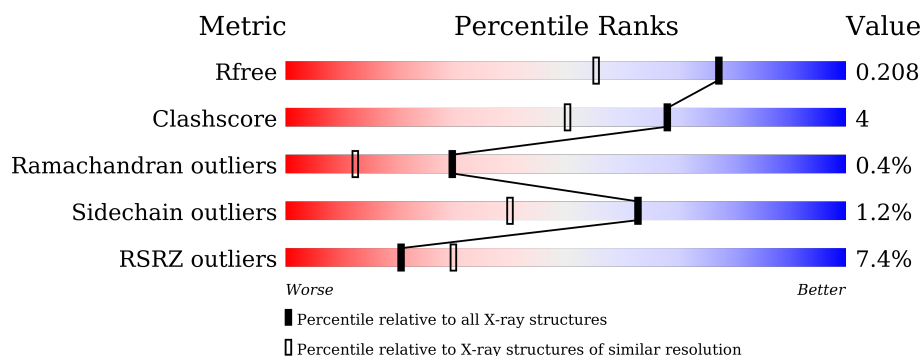
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	488	<div> <div>9%</div> <div> <div></div> <div>77%</div> <div>7%</div> <div>•</div> <div>16%</div> </div> </div>
1	B	488	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>7%</div> <div>•</div> <div>16%</div> </div> </div>

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.

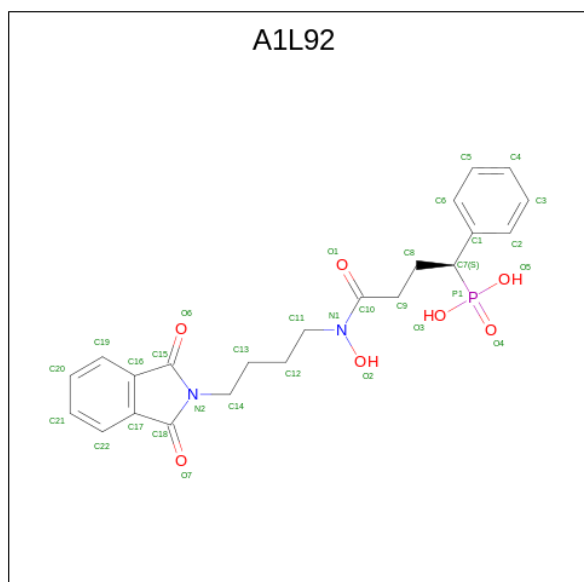
- Molecule 1 is a protein called 1-deoxy-D-xylulose 5-phosphate reductoisomerase, apicoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total 3285	C 2109	N 539	O 617	S 20	0	0	0
1	B	411	Total 3285	C 2109	N 539	O 617	S 20	0	0	0

- 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 1 1	0	0
2	B	1	Total Mn 1 1	0	0

- Molecule 3 is [(1 {S})-4-[4-[1,3-bis(oxidanylidene)isoindol-2-yl]butyl-oxidanyl-amino]-4-oxidanylidene-1-phenyl-butyl]phosphonic acid (CCD ID: A1L92) (formula: C₂₂H₂₅N₂O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	22	2	7	1		
3	B	1	Total	C	N	O	P	0	0
			32	22	2	7	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	3	Total	Ca	0	0
			3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	265	Total	O	0	0
			265	265		
6	B	296	Total	O	0	0
			296	296		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.76Å 77.55Å 111.49Å 90.00° 92.85° 90.00°	Depositor
Resolution (Å)	63.64 – 1.57 63.64 – 1.57	Depositor EDS
% Data completeness (in resolution range)	99.8 (63.64-1.57) 99.9 (63.64-1.57)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.167 , 0.195 0.181 , 0.208	Depositor DCC
R_{free} test set	6216 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.020 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7212	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.51% 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: GOL, MN, CA, A1L92

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	0.70	0/3348	1.15	5/4521 (0.1%)
1	B	0.74	0/3348	1.16	8/4521 (0.2%)
All	All	0.72	0/6696	1.15	13/9042 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	360	MET	CG-SD-CE	8.70	120.04	100.90
1	A	360	MET	CG-SD-CE	6.82	115.89	100.90
1	B	305	ASP	CA-CB-CG	6.45	119.05	112.60
1	A	234	HIS	CA-CB-CG	-6.40	107.40	113.80
1	B	245	LYS	CG-CD-CE	-5.35	99.00	111.30
1	B	175	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	279	MET	CG-SD-CE	-5.33	89.18	100.90
1	B	152	LYS	CB-CA-C	5.16	118.25	109.53
1	A	355	MET	CG-SD-CE	-5.15	89.57	100.90
1	A	286	THR	CA-CB-OG1	-5.14	101.88	109.60
1	A	305	ASP	CA-CB-CG	5.05	117.66	112.60
1	B	213	PHE	CA-CB-CG	-5.05	108.75	113.80
1	B	266	PHE	CA-CB-CG	5.04	118.84	113.80

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	3285	0	3334	21	0
1	B	3285	0	3334	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	32	0	0	0	0
3	B	32	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
5	B	3	0	0	0	0
6	A	265	0	0	5	0
6	B	296	0	0	13	0
All	All	7212	0	6684	48	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 4.

All (48) 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:360:MET:HG3	6:B:869:HOH:O	1.68	0.94
1:B:372:ASP:HB3	6:B:675:HOH:O	1.71	0.88
1:B:328:ASN:HB3	6:B:838:HOH:O	1.84	0.75
1:B:204:ASN:HB3	6:B:614:HOH:O	1.87	0.74
1:B:202:LEU:HD11	6:B:614:HOH:O	1.88	0.72
1:A:122:TYR:OH	1:A:126:ARG:NH1	2.22	0.72
1:B:204:ASN:ND2	1:B:207:SER:OG	2.27	0.67
1:B:153:ASP:C	6:B:601:HOH:O	2.37	0.66
1:B:372:ASP:CB	6:B:675:HOH:O	2.37	0.65
1:A:292:LYS:O	1:A:294:PRO:HD3	1.99	0.61
6:A:733:HOH:O	1:B:393:LYS:HE3	2.03	0.59
1:B:279:MET:HE1	1:B:438:PHE:CZ	2.38	0.58
1:A:360:MET:HG3	6:A:836:HOH:O	2.04	0.57
1:A:340:ILE:HG12	1:A:355:MET:HG2	1.85	0.57
1:A:328:ASN:HB3	6:A:823:HOH:O	2.05	0.56
1:A:138:LYS:HA	1:A:160:CYS:SG	2.47	0.55
1:B:279:MET:CA	1:B:279:MET:HE3	2.37	0.55
1:B:393:LYS:HG2	6:B:710:HOH:O	2.08	0.54
1:B:292:LYS:NZ	6:B:603:HOH:O	2.37	0.54
1:B:403:LYS:HE3	1:B:407:GLN:OE1	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:LYS:HG2	6:A:656:HOH:O	2.10	0.51
1:A:141:TYR:CE2	1:A:145:LYS:HD2	2.46	0.51
1:B:84:GLY:HA3	1:B:181:ILE:HG23	1.93	0.51
1:A:481:ILE:O	1:A:484:LYS:HG3	2.10	0.51
1:A:403:LYS:HE3	1:A:407:GLN:NE2	2.25	0.50
1:A:230:VAL:O	1:A:360:MET:HE1	2.11	0.49
1:B:122:TYR:CZ	1:B:126:ARG:HD2	2.48	0.49
1:B:227:ILE:HD12	1:B:252:LEU:HD13	1.96	0.48
1:B:328:ASN:CB	6:B:838:HOH:O	2.53	0.48
1:B:279:MET:HE3	1:B:279:MET:HA	1.95	0.48
1:A:87:GLY:O	1:A:91:THR:HG23	2.14	0.48
1:A:280:ASP:OD1	1:A:283:LYS:NZ	2.47	0.48
1:B:254:ASP:O	1:B:255:ASN:HB2	2.14	0.47
1:B:360:MET:CG	6:B:869:HOH:O	2.44	0.46
1:A:290:ALA:C	1:A:292:LYS:H	2.25	0.45
1:B:424:GLU:HG3	6:B:882:HOH:O	2.17	0.45
1:B:292:LYS:HA	1:B:292:LYS:HD3	1.80	0.44
1:B:76:LYS:HB3	1:B:77:PRO:HD3	2.00	0.44
1:A:131:GLU:OE1	1:A:131:GLU:N	2.50	0.44
1:B:113:TYR:C	1:B:113:TYR:CD1	2.96	0.43
1:A:484:LYS:HD2	1:A:485:HIS:N	2.34	0.42
1:A:293:HIS:CD2	1:A:293:HIS:O	2.72	0.42
1:A:141:TYR:CD1	1:A:160:CYS:HB2	2.54	0.42
1:A:372:ASP:OD2	6:A:601:HOH:O	2.22	0.42
1:B:449:GLU:CD	6:B:602:HOH:O	2.63	0.42
1:A:104:GLU:HG3	1:A:106:VAL:CG1	2.50	0.41
1:A:481:ILE:O	1:A:484:LYS:HE3	2.20	0.41
1:B:181:ILE:HD12	1:B:185:GLN:HB3	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	409/488 (84%)	398 (97%)	9 (2%)	2 (0%)	24	9
1	B	409/488 (84%)	400 (98%)	8 (2%)	1 (0%)	43	26
All	All	818/976 (84%)	798 (98%)	17 (2%)	3 (0%)	30	12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	342	SER
1	A	294	PRO
1	B	342	SER

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	378/449 (84%)	373 (99%)	5 (1%)	61	37
1	B	378/449 (84%)	374 (99%)	4 (1%)	65	43
All	All	756/898 (84%)	747 (99%)	9 (1%)	63	40

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

Mol	Chain	Res	Type
1	A	117	SER
1	A	293	HIS
1	A	296	TRP
1	A	298	MET
1	A	349	LYS
1	B	163	GLU
1	B	292	LYS
1	B	387	SER
1	B	460	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	ASN
1	A	220	ASN
1	A	224	ASN
1	A	260	ASN
1	A	261	ASN
1	A	293	HIS
1	A	377	ASN
1	A	407	GLN
1	B	92	ASN
1	B	105	ASN
1	B	119	ASN
1	B	204	ASN
1	B	255	ASN
1	B	293	HIS
1	B	433	ASN
1	B	452	ASN
1	B	486	ASN

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 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 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
3	A1L92	A	502	2	33,34,34	1.97	8 (24%)	39,48,48	2.95	16 (41%)
3	A1L92	B	502	2	33,34,34	2.33	8 (24%)	39,48,48	2.34	13 (33%)
4	GOL	B	506	-	5,5,5	0.18	0	5,5,5	0.61	0
4	GOL	A	503	-	5,5,5	0.18	0	5,5,5	0.99	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
3	A1L92	A	502	2	-	4/26/42/42	0/3/3/3
3	A1L92	B	502	2	-	3/26/42/42	0/3/3/3
4	GOL	B	506	-	-	0/4/4/4	-
4	GOL	A	503	-	-	0/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	A1L92	O7-C18	6.90	1.36	1.22
3	B	502	A1L92	O6-C15	6.45	1.35	1.22
3	B	502	A1L92	O7-C18	5.57	1.33	1.22
3	B	502	A1L92	C18-N2	-5.19	1.33	1.39
3	B	502	A1L92	C11-N1	4.31	1.54	1.46
3	A	502	A1L92	O6-C15	3.72	1.29	1.22
3	B	502	A1L92	C15-N2	-3.63	1.35	1.39
3	A	502	A1L92	C15-N2	-3.52	1.35	1.39
3	A	502	A1L92	P1-C7	3.25	1.86	1.82
3	B	502	A1L92	P1-O4	3.08	1.54	1.49
3	B	502	A1L92	C3-C2	3.01	1.45	1.38
3	A	502	A1L92	O2-N1	2.93	1.42	1.40
3	B	502	A1L92	C10-N1	-2.53	1.31	1.34
3	A	502	A1L92	P1-O4	2.48	1.53	1.49
3	A	502	A1L92	C9-C10	2.26	1.56	1.51
3	A	502	A1L92	C6-C1	2.04	1.42	1.39

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	A1L92	C16-C15-N2	9.56	112.67	105.88
3	A	502	A1L92	C18-N2-C15	-6.65	106.61	112.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	A1L92	C17-C18-N2	6.30	110.36	105.88
3	B	502	A1L92	C16-C15-N2	6.19	110.28	105.88
3	A	502	A1L92	C14-N2-C18	5.37	129.28	123.88
3	A	502	A1L92	C17-C18-N2	5.30	109.64	105.88
3	B	502	A1L92	C18-N2-C15	-4.10	108.69	112.03
3	A	502	A1L92	C12-C11-N1	4.03	119.42	111.06
3	A	502	A1L92	C17-C16-C15	-3.92	104.86	108.26
3	B	502	A1L92	C17-C16-C15	-3.91	104.87	108.26
3	A	502	A1L92	O7-C18-N2	3.87	128.58	124.81
3	A	502	A1L92	C8-C7-C1	-3.81	107.15	113.42
3	A	502	A1L92	O7-C18-C17	-3.55	121.77	128.68
3	B	502	A1L92	C5-C6-C1	-3.42	116.42	120.65
3	A	502	A1L92	O6-C15-C16	-3.34	122.19	128.68
3	A	502	A1L92	C22-C17-C18	3.18	134.82	129.63
3	B	502	A1L92	O1-C10-C9	-3.15	112.34	121.31
3	B	502	A1L92	C22-C17-C18	2.94	134.42	129.63
3	A	502	A1L92	O2-N1-C11	2.93	120.91	113.59
3	B	502	A1L92	C19-C16-C15	2.87	134.31	129.63
3	B	502	A1L92	C16-C17-C18	-2.84	105.80	108.26
3	B	502	A1L92	C4-C3-C2	-2.71	116.06	120.19
3	A	502	A1L92	C16-C17-C18	-2.37	106.20	108.26
3	A	502	A1L92	O4-P1-C7	-2.35	106.78	113.35
3	A	502	A1L92	C19-C16-C15	2.25	133.30	129.63
3	A	502	A1L92	C6-C1-C7	2.20	124.37	120.82
3	B	502	A1L92	C2-C1-C6	2.13	120.95	118.29
3	B	502	A1L92	O3-P1-O4	-2.03	108.34	113.45
3	B	502	A1L92	O7-C18-N2	-2.01	122.86	124.81

There are no chirality outliers.

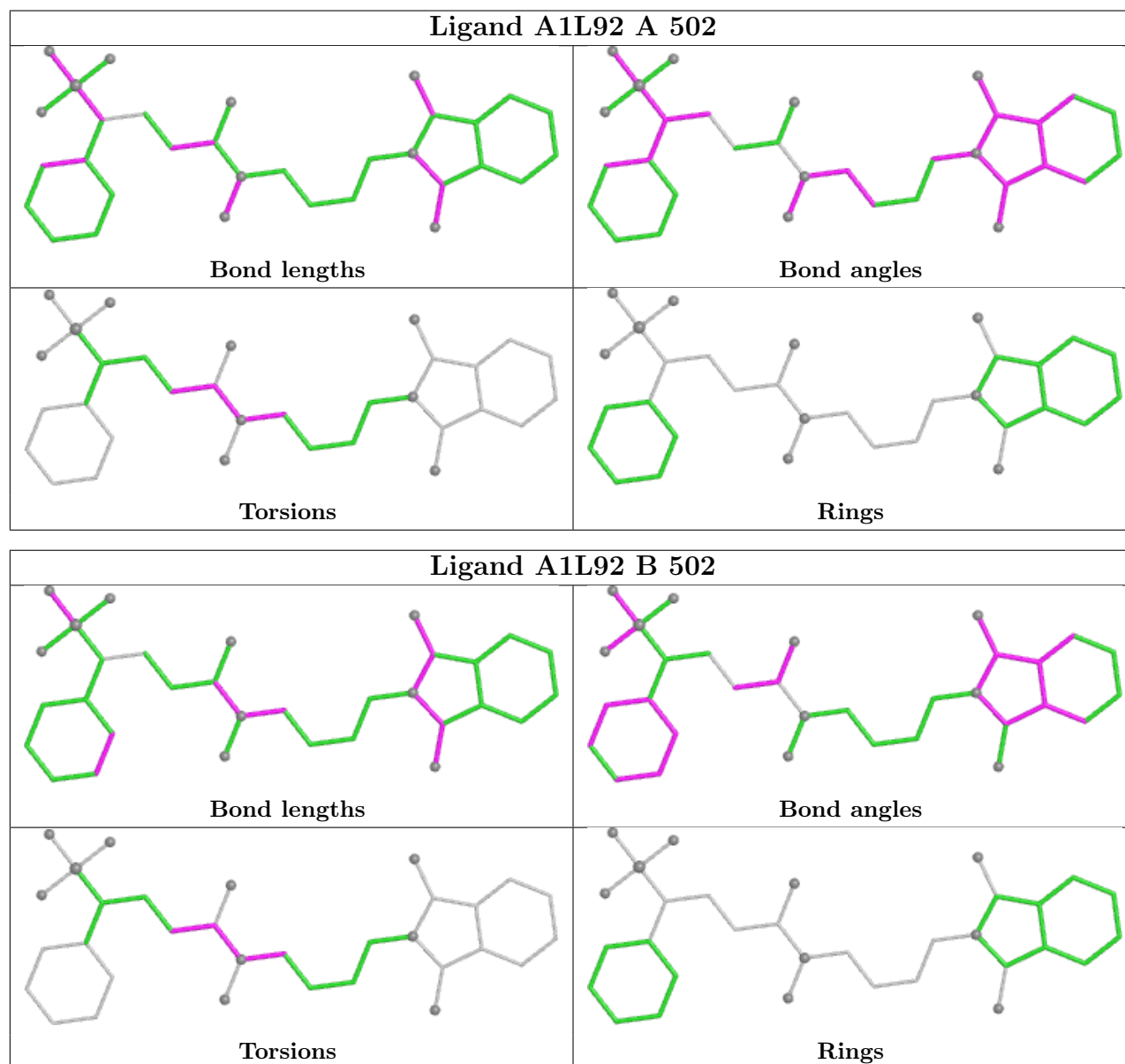
All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	A1L92	C12-C11-N1-O2
3	B	502	A1L92	N1-C10-C9-C8
3	A	502	A1L92	C12-C11-N1-C10
3	B	502	A1L92	C12-C11-N1-C10
3	A	502	A1L92	O1-C10-N1-O2
3	A	502	A1L92	N1-C10-C9-C8
3	B	502	A1L92	O1-C10-N1-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	411/488 (84%)	0.46	42 (10%) 12 19	11, 22, 58, 100	0
1	B	411/488 (84%)	0.16	19 (4%) 37 49	11, 20, 41, 76	0
All	All	822/976 (84%)	0.31	61 (7%) 20 30	11, 20, 52, 100	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	294	PRO	8.0
1	A	296	TRP	6.3
1	B	296	TRP	5.7
1	A	292	LYS	5.1
1	A	293	HIS	4.5
1	A	151	ILE	4.5
1	A	141	TYR	4.2
1	A	140	VAL	3.8
1	A	147	LEU	3.7
1	B	294	PRO	3.7
1	A	76	LYS	3.6
1	B	76	LYS	3.5
1	A	184	PHE	3.5
1	A	295	LYS	3.5
1	A	297	LYS	3.5
1	B	485	HIS	3.4
1	A	122	TYR	3.2
1	A	77	PRO	3.2
1	B	295	LYS	3.2
1	A	148	VAL	3.2
1	A	154	TYR	3.1
1	B	297	LYS	2.9
1	A	485	HIS	2.9
1	A	144	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	213	PHE	2.8
1	A	157	ILE	2.8
1	A	158	ILE	2.8
1	A	160	CYS	2.7
1	A	133	LEU	2.6
1	B	486	ASN	2.6
1	A	119	ASN	2.6
1	B	153	ASP	2.5
1	A	328	ASN	2.5
1	A	106	VAL	2.4
1	A	118	VAL	2.4
1	B	141	TYR	2.4
1	B	293	HIS	2.4
1	B	460	SER	2.4
1	A	135	ILE	2.3
1	A	172	ASN	2.3
1	A	117	SER	2.3
1	A	138	LYS	2.2
1	B	484	LYS	2.2
1	A	155	LYS	2.2
1	B	292	LYS	2.2
1	A	185	GLN	2.2
1	B	182	ASP	2.2
1	A	291	LEU	2.1
1	A	192	TYR	2.1
1	A	173	SER	2.1
1	B	377	ASN	2.1
1	A	113	TYR	2.1
1	A	149	LYS	2.1
1	A	152	LYS	2.1
1	A	290	ALA	2.1
1	B	184	PHE	2.0
1	B	148	VAL	2.0
1	A	129	LEU	2.0
1	A	115	ASN	2.0
1	A	162	ASP	2.0
1	B	105	ASN	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

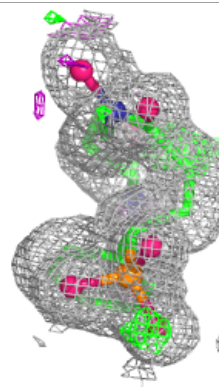
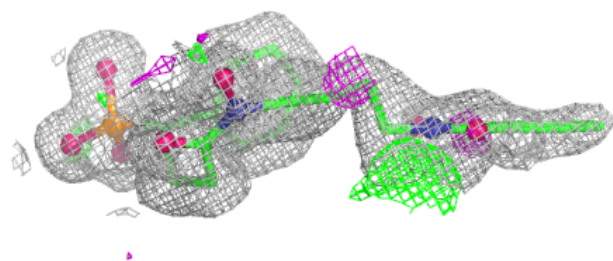
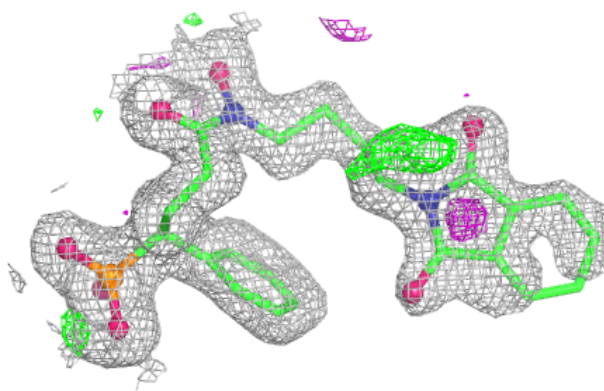
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	A	503	6/6	0.94	0.08	20,22,26,27	0
4	GOL	B	506	6/6	0.95	0.07	20,23,26,27	0
3	A1L92	A	502	32/32	0.97	0.10	15,21,55,59	0
5	CA	B	504	1/1	0.97	0.07	23,23,23,23	0
5	CA	B	505	1/1	0.97	0.08	24,24,24,24	0
3	A1L92	B	502	32/32	0.98	0.09	12,19,48,53	0
5	CA	B	503	1/1	0.99	0.06	16,16,16,16	0
2	MN	B	501	1/1	1.00	0.02	12,12,12,12	0
2	MN	A	501	1/1	1.00	0.02	13,13,13,13	0

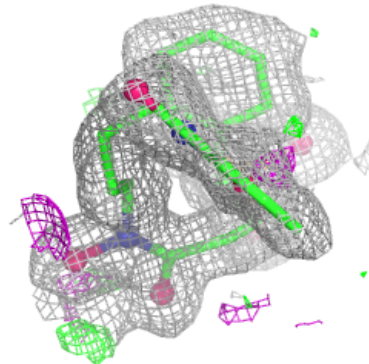
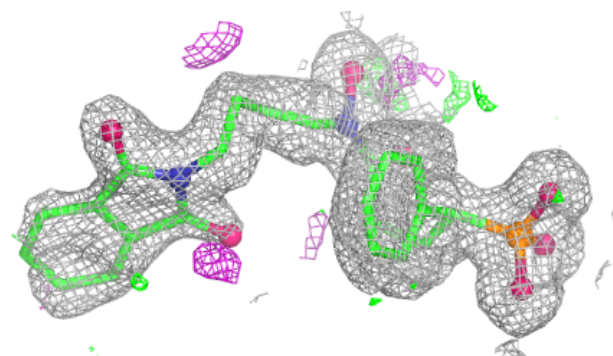
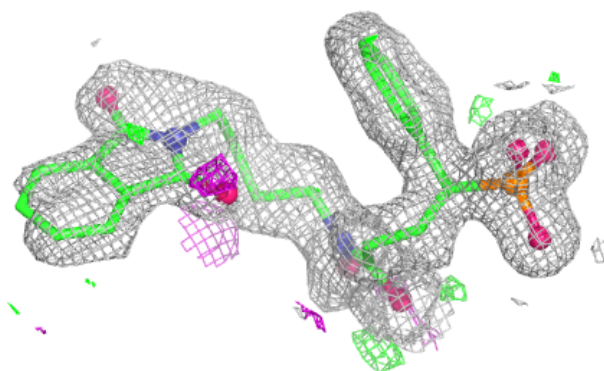
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A1L92 A 502:

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

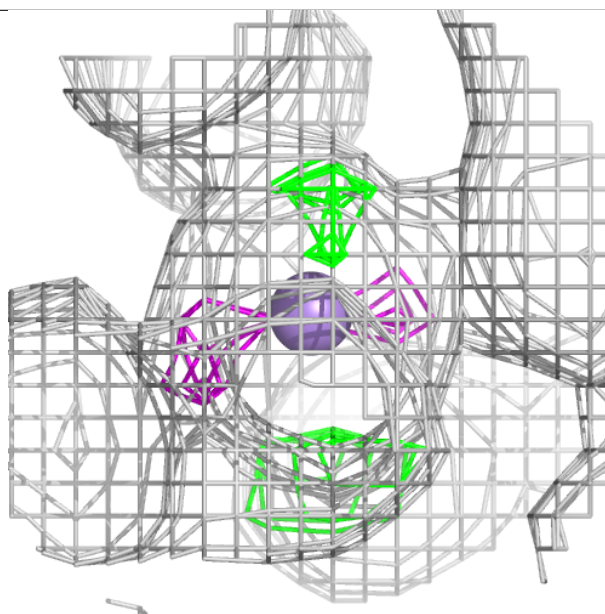
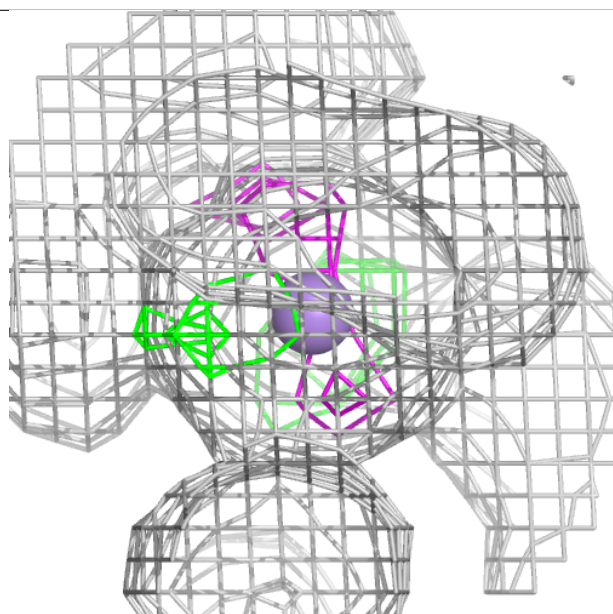
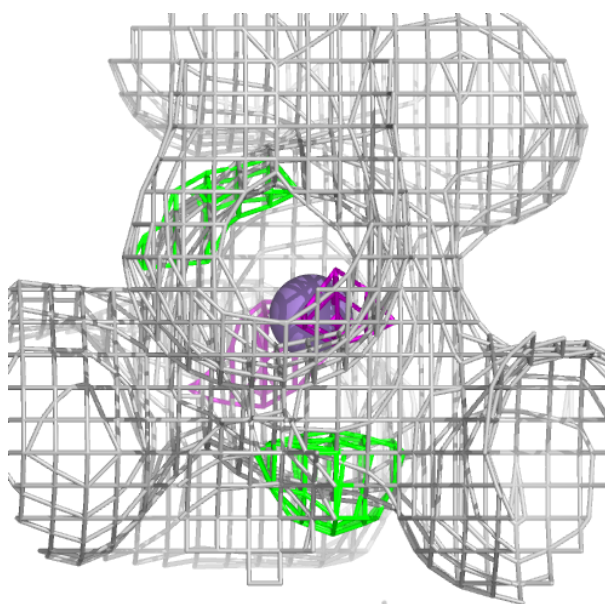
**Electron density around A1L92 B 502:**

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



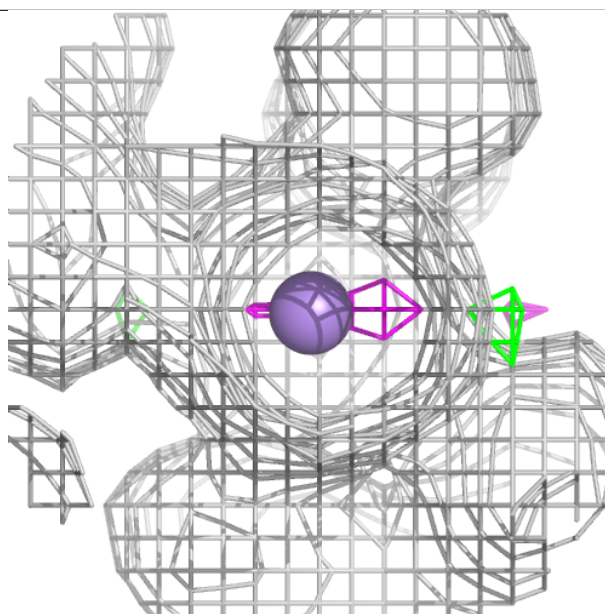
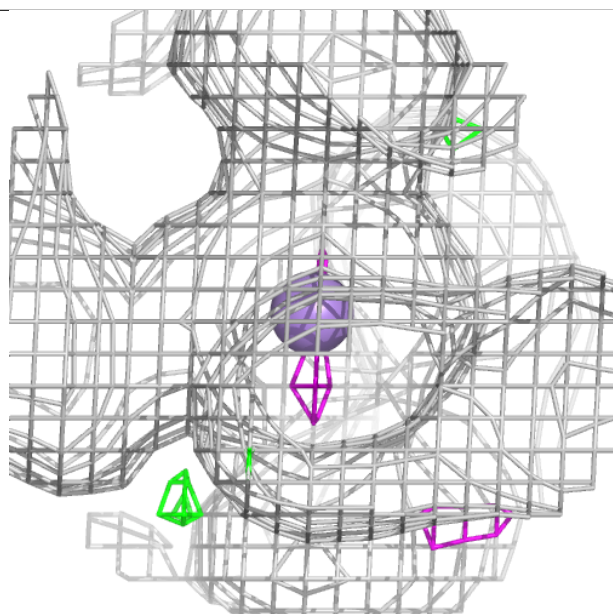
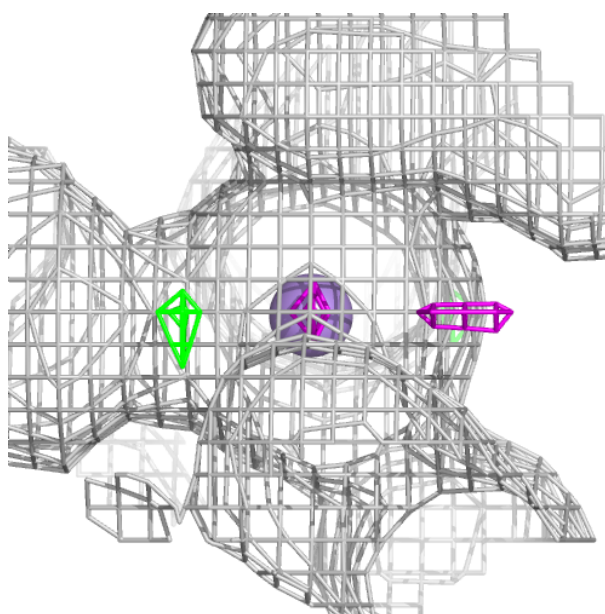
Electron density around MN B 501:

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



Electron density around MN A 501:

$2mF_o-DF_c$ (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.