



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:19 AM UTC

PDB ID : 9LS7 / pdb_00009ls7
Title : The crystal structure of human RyR3 Repeat12 domain complex with RI-2 and AMP-PCP
Authors : Hadiatullah, H.; Lin, L.; Yuchi, Z.
Deposited on : 2025-02-04
Resolution : 2.49 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

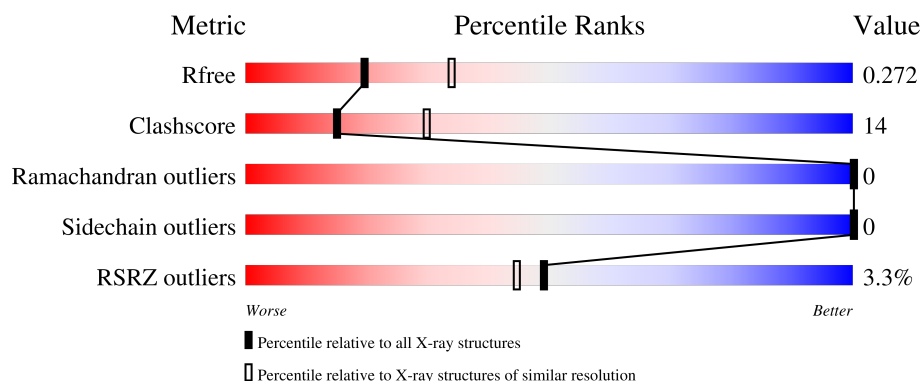
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	218	2% 71% 23% 6%
1	B	218	0% 67% 27% 6%
1	C	218	2% 68% 25% 6%
1	D	218	7% 68% 19% 13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6340 atoms, of which 66 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.

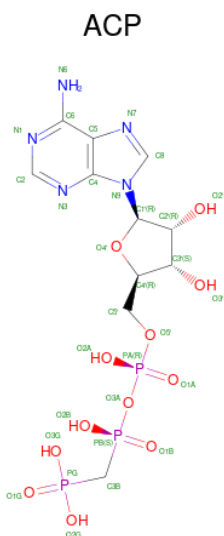
- Molecule 1 is a protein called Ryanodine receptor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1598	1016	280	296	6			
1	B	205	Total	C	N	O	S	0	0	0
			1582	1016	277	283	6			
1	C	206	Total	C	N	O	S	0	0	0
			1543	992	266	279	6			
1	D	190	Total	C	N	O	S	0	0	0
			1378	886	242	245	5			

There are 4 discrepancies between the modelled and reference sequences:

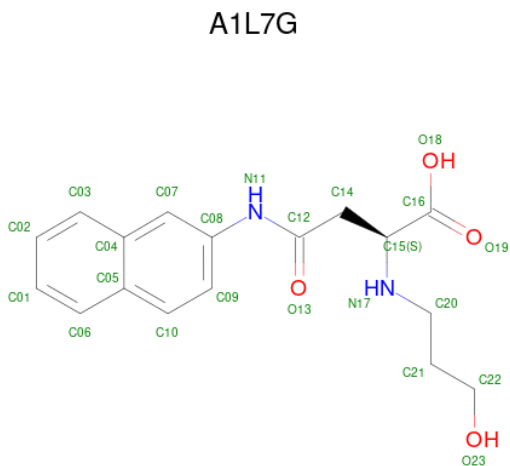
Chain	Residue	Modelled	Actual	Comment	Reference
A	846	ASN	GLN	conflict	UNP Q15413
B	846	ASN	GLN	conflict	UNP Q15413
C	846	ASN	GLN	conflict	UNP Q15413
D	846	ASN	GLN	conflict	UNP Q15413

- Molecule 2 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (CCD ID: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 45	C 11	H 14	N 5	O 12	P 3	0	0
2	B	1	Total 45	C 11	H 14	N 5	O 12	P 3	0	0

- Molecule 3 is (2 {S})-4-(naphthalen-2-ylamino)-4-oxidanylidene-2-(3-oxidanylpropylamino)butanoic acid (CCD ID: A1L7G) (formula: $C_{17}H_{20}N_2O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			42	17	19	2	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	0	0
			42	17	19	2	4		

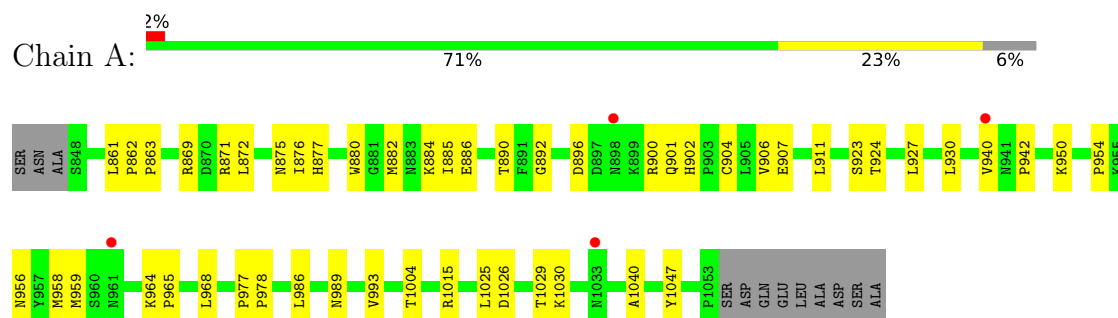
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		
4	B	15	Total	O	0	0
			15	15		
4	C	14	Total	O	0	0
			14	14		
4	D	9	Total	O	0	0
			9	9		

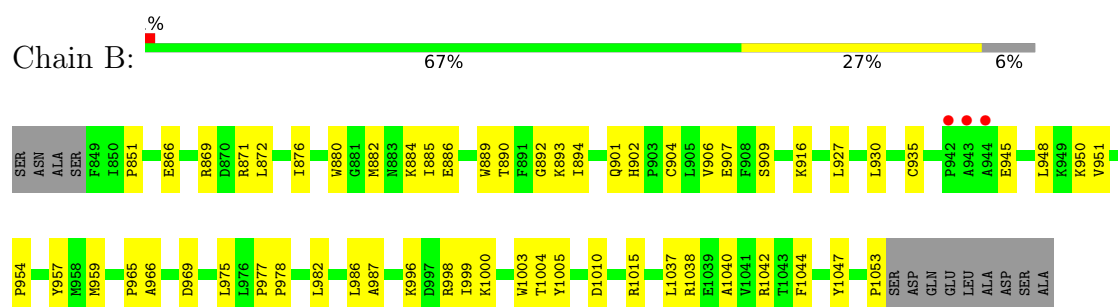
3 Residue-property plots [i](#)

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.

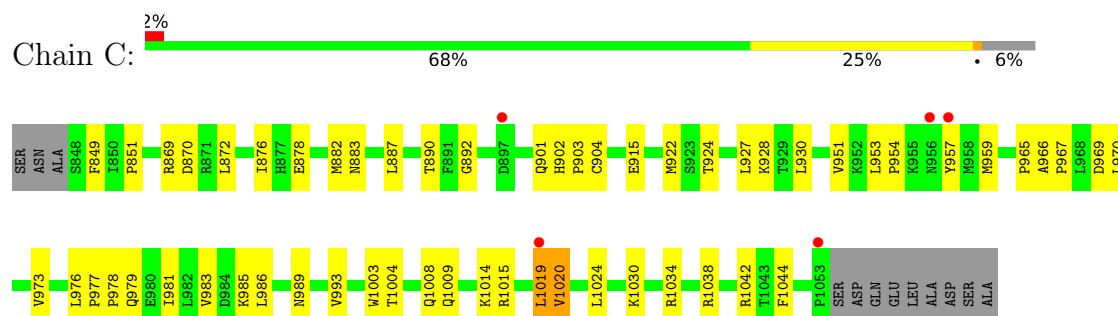
• Molecule 1: Ryanodine receptor 3



• Molecule 1: Ryanodine receptor 3

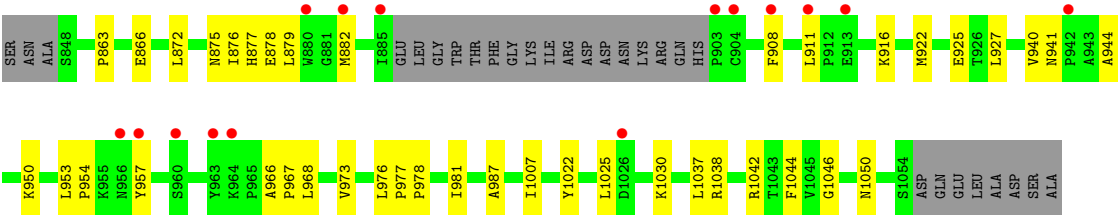


• Molecule 1: Ryanodine receptor 3



• Molecule 1: Ryanodine receptor 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.36Å 61.62Å 277.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	16.96 – 2.49 16.96 – 2.49	Depositor EDS
% Data completeness (in resolution range)	92.1 (16.96-2.49) 78.3 (16.96-2.49)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.49Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, R_{free}	0.234 , 0.271 0.235 , 0.272	Depositor DCC
R_{free} test set	1485 reflections (3.97%)	wwPDB-VP
Wilson B-factor (Å ²)	49.5	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.033 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6340	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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: ACP, A1L7G

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	0.07	0/1637	0.20	0/2235
1	B	0.08	0/1621	0.21	0/2212
1	C	0.07	0/1582	0.32	2/2166 (0.1%)
1	D	0.07	0/1410	0.25	0/1932
All	All	0.07	0/6250	0.25	2/8545 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1019	LEU	N-CA-C	-8.52	96.89	109.15
1	C	1020	VAL	N-CA-CB	6.76	120.67	111.21

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	1598	0	1529	32	0
1	B	1582	0	1539	44	0
1	C	1543	0	1451	52	0
1	D	1378	0	1276	37	0
2	A	31	14	14	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	31	14	14	0	0
3	A	23	19	0	2	0
3	B	23	19	0	1	0
4	A	27	0	0	0	0
4	B	15	0	0	0	0
4	C	14	0	0	0	0
4	D	9	0	0	0	0
All	All	6274	66	5823	165	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 14.

All (165) 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:C:882:MET:HE2	1:C:953:LEU:HD11	1.46	0.95
1:C:1019:LEU:O	1:C:1020:VAL:HG13	1.74	0.88
1:C:977:PRO:HG2	1:C:978:PRO:HD3	1.54	0.88
1:C:930:LEU:HD21	1:C:986:LEU:HD11	1.56	0.86
1:B:987:ALA:HA	1:B:1037:LEU:HD12	1.57	0.85
1:C:882:MET:HE3	1:C:967:PRO:HG2	1.61	0.83
1:C:1019:LEU:O	1:C:1020:VAL:CG1	2.28	0.81
1:C:922:MET:HA	1:C:922:MET:HE2	1.65	0.78
1:A:892:GLY:HA3	1:A:901:GLN:HB3	1.66	0.78
1:C:959:MET:HA	1:C:959:MET:HE2	1.66	0.78
1:C:849:PHE:CZ	1:C:851:PRO:HG3	2.22	0.75
1:A:861:LEU:HD22	1:A:924:THR:HG23	1.71	0.73
1:B:892:GLY:HA3	1:B:901:GLN:HB3	1.71	0.73
1:D:954:PRO:HD2	1:D:957:TYR:CE1	2.27	0.69
1:C:927:LEU:HD21	1:C:1044:PHE:CE1	2.27	0.69
1:C:1003:TRP:HA	1:C:1014:LYS:HB3	1.77	0.67
1:C:890:THR:O	1:C:901:GLN:HA	1.94	0.67
1:A:930:LEU:HD21	1:A:986:LEU:HD11	1.76	0.66
1:D:882:MET:SD	1:D:953:LEU:HD11	2.35	0.66
1:C:977:PRO:CG	1:C:978:PRO:HD3	2.23	0.66
1:A:885:ILE:HD12	1:A:965:PRO:HG3	1.77	0.66
1:C:927:LEU:HD21	1:C:1044:PHE:HE1	1.60	0.65
1:A:959:MET:HA	1:A:959:MET:HE2	1.79	0.64
1:C:976:LEU:HB3	1:C:977:PRO:HD2	1.79	0.64
1:D:882:MET:HG2	1:D:953:LEU:HD11	1.80	0.63
1:C:970:LEU:HB3	1:C:1042:ARG:HG2	1.81	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:882:MET:CG	1:D:953:LEU:HD11	2.29	0.63
1:B:987:ALA:HA	1:B:1037:LEU:CD1	2.29	0.62
1:D:927:LEU:HD21	1:D:1044:PHE:CE1	2.34	0.62
1:C:989:ASN:O	1:C:993:VAL:HG23	1.99	0.62
1:B:890:THR:O	1:B:901:GLN:HA	2.00	0.61
1:A:940:VAL:O	1:A:942:PRO:HD3	2.02	0.60
1:D:1007:ILE:HD12	1:D:1007:ILE:H	1.67	0.59
1:B:884:LYS:O	1:B:889:TRP:HB2	2.02	0.58
1:A:923:SER:O	1:A:927:LEU:HD13	2.04	0.58
1:B:959:MET:HA	1:B:959:MET:HE2	1.86	0.57
1:C:930:LEU:CD2	1:C:986:LEU:HD11	2.32	0.57
1:C:882:MET:CE	1:C:953:LEU:HD11	2.29	0.57
1:C:882:MET:CE	1:C:967:PRO:HG2	2.33	0.57
1:B:872:LEU:HD11	1:B:1040:ALA:HB1	1.87	0.57
1:B:975:LEU:HD11	1:B:1042:ARG:CG	2.35	0.56
1:A:977:PRO:N	1:A:978:PRO:HD2	2.20	0.56
1:A:880:TRP:CZ2	1:A:884:LYS:HD2	2.40	0.56
1:C:1008:GLN:HG2	1:C:1009:GLN:H	1.68	0.56
1:D:863:PRO:HA	1:D:866:GLU:OE2	2.05	0.56
1:A:871:ARG:HD3	1:A:1047:TYR:OH	2.05	0.56
1:C:922:MET:HE2	1:C:922:MET:CA	2.36	0.55
1:C:970:LEU:HD13	1:C:1042:ARG:HG2	1.87	0.55
1:D:950:LYS:HD3	1:D:966:ALA:O	2.05	0.55
1:B:902:HIS:CE1	1:B:904:CYS:HB2	2.42	0.55
1:B:975:LEU:HD11	1:B:1042:ARG:HG2	1.87	0.55
1:D:977:PRO:N	1:D:978:PRO:HD2	2.21	0.55
1:B:951:VAL:HG12	1:B:969:ASP:HB2	1.89	0.55
1:B:977:PRO:N	1:B:978:PRO:HD2	2.22	0.55
1:A:906:VAL:HG23	1:A:907:GLU:O	2.08	0.54
1:C:1034:ARG:HG2	1:C:1038:ARG:HH21	1.73	0.54
1:C:1019:LEU:C	1:C:1020:VAL:HG13	2.32	0.53
1:B:930:LEU:HD21	1:B:986:LEU:HD11	1.90	0.53
1:A:954:PRO:HB2	1:A:956:ASN:OD1	2.10	0.51
1:A:877:HIS:CE1	1:A:911:LEU:HD11	2.46	0.51
1:C:882:MET:HE2	1:C:953:LEU:HD21	1.93	0.51
1:D:987:ALA:HA	1:D:1037:LEU:CD1	2.41	0.51
1:A:890:THR:O	1:A:901:GLN:HA	2.11	0.51
1:A:989:ASN:O	1:A:993:VAL:HG23	2.10	0.51
1:D:1022:TYR:O	1:D:1030:LYS:HE2	2.11	0.51
1:D:944:ALA:HB1	1:D:1046:GLY:O	2.11	0.51
1:C:878:GLU:OE1	1:C:965:PRO:HA	2.12	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1102:A1L7G:N11	3:A:1102:A1L7G:O19	2.44	0.50
1:B:950:LYS:HD3	1:B:966:ALA:O	2.10	0.50
1:B:893:LYS:C	1:B:894:ILE:HD12	2.36	0.50
1:B:935:CYS:SG	1:B:1053:PRO:HA	2.52	0.50
1:D:872:LEU:O	1:D:876:ILE:HG13	2.12	0.50
1:C:979:GLN:O	1:C:983:VAL:HG23	2.11	0.49
1:D:908:PHE:O	1:D:916:LYS:HE2	2.12	0.49
1:B:882:MET:O	1:B:886:GLU:HG3	2.13	0.49
1:B:906:VAL:HG23	1:B:907:GLU:O	2.12	0.49
1:C:882:MET:HE2	1:C:953:LEU:CD1	2.30	0.49
1:B:998:ARG:O	1:B:1003:TRP:HB2	2.13	0.48
1:B:885:ILE:HD12	1:B:965:PRO:HG2	1.95	0.48
1:C:959:MET:HE2	1:C:959:MET:CA	2.42	0.48
1:D:1007:ILE:HD12	1:D:1007:ILE:N	2.29	0.48
1:B:876:ILE:HG13	3:B:1102:A1L7G:C22	2.43	0.48
1:D:922:MET:HA	1:D:922:MET:HE2	1.94	0.48
1:C:872:LEU:O	1:C:876:ILE:HG12	2.13	0.48
1:B:945:GLU:O	1:B:948:LEU:HB2	2.13	0.48
1:C:951:VAL:HG22	1:C:969:ASP:HB2	1.95	0.48
1:B:954:PRO:HD2	1:B:957:TYR:CE1	2.49	0.47
1:A:872:LEU:HD11	1:A:1040:ALA:HB1	1.95	0.47
1:C:869:ARG:HB2	1:C:927:LEU:HD12	1.96	0.47
1:C:902:HIS:ND1	1:C:904:CYS:HB2	2.29	0.47
1:A:958:MET:HB2	1:A:964:LYS:HE3	1.97	0.47
1:A:1025:LEU:HB3	1:A:1030:LYS:HD2	1.97	0.47
1:A:1004:THR:O	1:A:1015:ARG:HA	2.15	0.47
1:B:871:ARG:HD3	1:B:1047:TYR:OH	2.15	0.47
1:C:902:HIS:CE1	1:C:904:CYS:HB2	2.50	0.47
1:B:1038:ARG:HD3	1:B:1042:ARG:NH2	2.30	0.46
1:D:940:VAL:HG21	1:D:1050:ASN:HB2	1.98	0.46
1:A:875:ASN:OD1	1:A:968:LEU:HG	2.16	0.46
1:B:872:LEU:HD11	1:B:1040:ALA:CB	2.45	0.46
1:C:973:VAL:CG1	1:C:1042:ARG:HG3	2.46	0.46
1:A:890:THR:HA	1:A:959:MET:SD	2.56	0.46
1:B:975:LEU:HD13	1:B:1042:ARG:HE	1.80	0.46
1:B:951:VAL:CG1	1:B:969:ASP:HB2	2.46	0.45
1:B:996:LYS:HG2	1:B:1000:LYS:NZ	2.30	0.45
1:B:866:GLU:HG3	1:B:869:ARG:HH21	1.81	0.45
1:B:996:LYS:HG2	1:B:1000:LYS:HZ3	1.81	0.45
1:B:1004:THR:HG22	1:B:1005:TYR:N	2.32	0.45
1:C:1019:LEU:O	1:C:1020:VAL:HG12	2.14	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:849:PHE:CE1	1:C:851:PRO:HG3	2.52	0.45
1:D:875:ASN:ND2	1:D:968:LEU:HD12	2.32	0.45
1:C:869:ARG:HB2	1:C:927:LEU:CD1	2.47	0.45
1:C:981:ILE:CG2	1:C:985:LYS:HE3	2.47	0.45
1:B:1010:ASP:H	1:B:1015:ARG:HB2	1.81	0.45
1:C:924:THR:CG2	1:C:928:LYS:HE3	2.47	0.45
1:B:996:LYS:HE2	1:B:1000:LYS:HZ3	1.82	0.44
1:A:872:LEU:HD11	1:A:1040:ALA:CB	2.47	0.44
1:C:849:PHE:CE2	1:C:851:PRO:HG3	2.52	0.44
1:D:976:LEU:HB3	1:D:977:PRO:HD2	1.99	0.44
1:D:1025:LEU:O	1:D:1030:LYS:HE3	2.18	0.44
1:C:1024:LEU:N	1:C:1024:LEU:HD12	2.33	0.44
1:D:922:MET:HE2	1:D:922:MET:CA	2.48	0.44
1:B:927:LEU:HD21	1:B:1044:PHE:CE1	2.53	0.44
1:B:851:PRO:HG3	1:B:999:ILE:HD12	1.99	0.44
1:B:986:LEU:C	1:B:1037:LEU:HD13	2.43	0.44
1:C:954:PRO:HD2	1:C:957:TYR:CE1	2.53	0.43
1:A:902:HIS:CE1	1:A:904:CYS:HB2	2.54	0.43
1:D:940:VAL:HG23	1:D:941:ASN:N	2.34	0.43
1:B:975:LEU:HD22	1:B:1038:ARG:HG3	2.00	0.43
1:C:883:ASN:O	1:C:887:LEU:HG	2.19	0.43
1:D:973:VAL:HB	1:D:1042:ARG:HD3	2.01	0.43
1:B:909:SER:O	1:B:916:LYS:NZ	2.42	0.43
1:B:982:LEU:O	1:B:986:LEU:HG	2.19	0.43
1:B:977:PRO:CD	1:B:978:PRO:HD2	2.49	0.42
1:C:927:LEU:HG	1:C:1044:PHE:HZ	1.84	0.42
1:D:879:LEU:CD2	1:D:967:PRO:HB3	2.49	0.42
1:D:954:PRO:HD2	1:D:957:TYR:HE1	1.77	0.42
1:A:1026:ASP:OD1	1:A:1029:THR:HB	2.19	0.42
1:D:922:MET:HE2	1:D:922:MET:N	2.35	0.42
1:D:977:PRO:CD	1:D:978:PRO:HD2	2.49	0.42
1:A:886:GLU:HB3	1:D:981:ILE:HG13	2.00	0.42
1:C:869:ARG:HD2	1:C:870:ASP:OD1	2.19	0.42
1:D:877:HIS:CE1	1:D:911:LEU:HD21	2.54	0.42
1:A:950:LYS:HG2	1:A:968:LEU:HD23	2.01	0.42
1:A:876:ILE:HG21	3:A:1102:A1L7G:C21	2.49	0.42
1:B:894:ILE:HD12	1:B:894:ILE:N	2.34	0.42
1:C:915:GLU:OE1	1:C:915:GLU:N	2.51	0.42
1:D:882:MET:HG2	1:D:953:LEU:CD1	2.48	0.42
1:A:1026:ASP:OD1	1:A:1026:ASP:N	2.53	0.42
1:C:1004:THR:O	1:C:1015:ARG:HA	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:925:GLU:HA	1:D:925:GLU:OE1	2.21	0.41
1:C:1030:LYS:HB3	1:C:1034:ARG:HH12	1.85	0.41
1:D:1038:ARG:HG2	1:D:1042:ARG:NH1	2.36	0.41
1:A:862:PRO:HA	1:A:863:PRO:HD3	1.97	0.41
1:C:966:ALA:N	1:C:967:PRO:HD3	2.35	0.41
1:A:869:ARG:HD2	1:A:869:ARG:C	2.46	0.41
1:A:882:MET:O	1:A:886:GLU:HG3	2.20	0.41
1:A:896:ASP:O	1:A:900:ARG:N	2.54	0.41
1:B:977:PRO:HD2	1:B:978:PRO:HD2	2.03	0.41
1:D:976:LEU:HB3	1:D:977:PRO:CD	2.51	0.40
1:B:880:TRP:CZ2	1:B:884:LYS:HG3	2.57	0.40
1:D:878:GLU:O	1:D:967:PRO:HG3	2.21	0.40
1:D:879:LEU:HD23	1:D:967:PRO:HB3	2.04	0.40
1:C:892:GLY:O	1:C:903:PRO:HA	2.22	0.40
1:D:879:LEU:HD23	1:D:967:PRO:CB	2.52	0.40
1:D:977:PRO:HD2	1:D:978:PRO:HD2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	204/218 (94%)	197 (97%)	7 (3%)	0	100	100
1	B	203/218 (93%)	199 (98%)	4 (2%)	0	100	100
1	C	204/218 (94%)	195 (96%)	9 (4%)	0	100	100
1	D	186/218 (85%)	182 (98%)	4 (2%)	0	100	100
All	All	797/872 (91%)	773 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	166/195 (85%)	166 (100%)	0	100	100
1	B	162/195 (83%)	162 (100%)	0	100	100
1	C	151/195 (77%)	151 (100%)	0	100	100
1	D	128/195 (66%)	128 (100%)	0	100	100
All	All	607/780 (78%)	607 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	864	HIS
1	B	1009	GLN
1	C	1009	GLN

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](#)

4 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	A1L7G	A	1102	-	24,24,24	1.37	2 (8%)	30,31,31	1.28	3 (10%)
3	A1L7G	B	1102	-	24,24,24	1.38	3 (12%)	30,31,31	1.25	4 (13%)
2	ACP	B	1101	-	31,33,33	0.73	1 (3%)	47,52,52	1.08	2 (4%)
2	ACP	A	1101	-	31,33,33	0.64	1 (3%)	47,52,52	1.12	2 (4%)

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	A1L7G	A	1102	-	-	12/17/17/17	0/2/2/2
3	A1L7G	B	1102	-	-	10/17/17/17	0/2/2/2
2	ACP	B	1101	-	-	0/19/38/38	0/3/3/3
2	ACP	A	1101	-	-	6/19/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1102	A1L7G	C12-N11	4.68	1.45	1.35
3	A	1102	A1L7G	C12-N11	4.66	1.45	1.35
2	B	1101	ACP	PB-O3A	3.40	1.62	1.58
2	A	1101	ACP	PB-O3A	3.20	1.62	1.58
3	B	1102	A1L7G	C08-N11	2.16	1.46	1.41
3	A	1102	A1L7G	O13-C12	-2.09	1.19	1.23
3	B	1102	A1L7G	O13-C12	-2.01	1.19	1.23

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	ACP	PB-O3A-PA	-6.10	112.46	132.37
2	B	1101	ACP	PB-O3A-PA	-5.97	112.90	132.37
3	A	1102	A1L7G	C14-C12-N11	3.38	118.80	114.62

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1102	A1L7G	C14-C12-N11	3.32	118.73	114.62
3	A	1102	A1L7G	C08-N11-C12	-3.12	122.00	127.52
3	B	1102	A1L7G	C14-C15-N17	-2.98	105.46	110.69
3	A	1102	A1L7G	C14-C15-C16	2.32	115.42	110.79
3	B	1102	A1L7G	C08-N11-C12	-2.23	123.58	127.52
3	B	1102	A1L7G	O18-C16-C15	2.07	120.53	113.51
2	B	1101	ACP	O3'-C3'-C2'	-2.05	105.25	111.82
2	A	1101	ACP	O3'-C3'-C2'	-2.04	105.26	111.82

There are no chirality outliers.

All (28) torsion outliers are listed below:

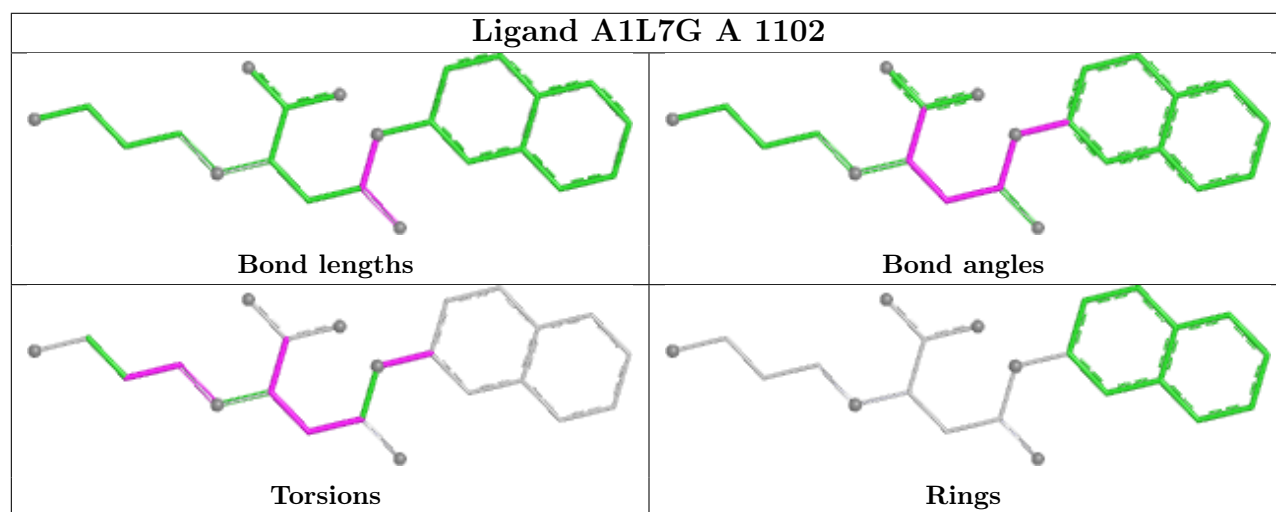
Mol	Chain	Res	Type	Atoms
2	A	1101	ACP	C5'-O5'-PA-O2A
2	A	1101	ACP	C5'-O5'-PA-O3A
2	A	1101	ACP	O4'-C4'-C5'-O5'
3	A	1102	A1L7G	C12-C14-C15-C16
3	A	1102	A1L7G	C12-C14-C15-N17
3	A	1102	A1L7G	N17-C20-C21-C22
3	B	1102	A1L7G	C14-C15-N17-C20
3	B	1102	A1L7G	C16-C15-N17-C20
3	A	1102	A1L7G	C07-C08-N11-C12
3	A	1102	A1L7G	C09-C08-N11-C12
2	A	1101	ACP	C3'-C4'-C5'-O5'
3	A	1102	A1L7G	C21-C20-N17-C15
3	B	1102	A1L7G	C21-C20-N17-C15
3	A	1102	A1L7G	N17-C15-C16-O19
3	B	1102	A1L7G	N17-C15-C16-O19
3	B	1102	A1L7G	C14-C15-C16-O19
3	A	1102	A1L7G	N17-C15-C16-O18
3	B	1102	A1L7G	N17-C15-C16-O18
3	B	1102	A1L7G	C14-C15-C16-O18
3	B	1102	A1L7G	N11-C12-C14-C15
2	A	1101	ACP	C5'-O5'-PA-O1A
3	A	1102	A1L7G	C14-C15-C16-O18
3	A	1102	A1L7G	C14-C15-C16-O19
3	A	1102	A1L7G	O13-C12-C14-C15
3	B	1102	A1L7G	O13-C12-C14-C15
3	A	1102	A1L7G	N11-C12-C14-C15
2	A	1101	ACP	PG-C3B-PB-O1B
3	B	1102	A1L7G	C09-C08-N11-C12

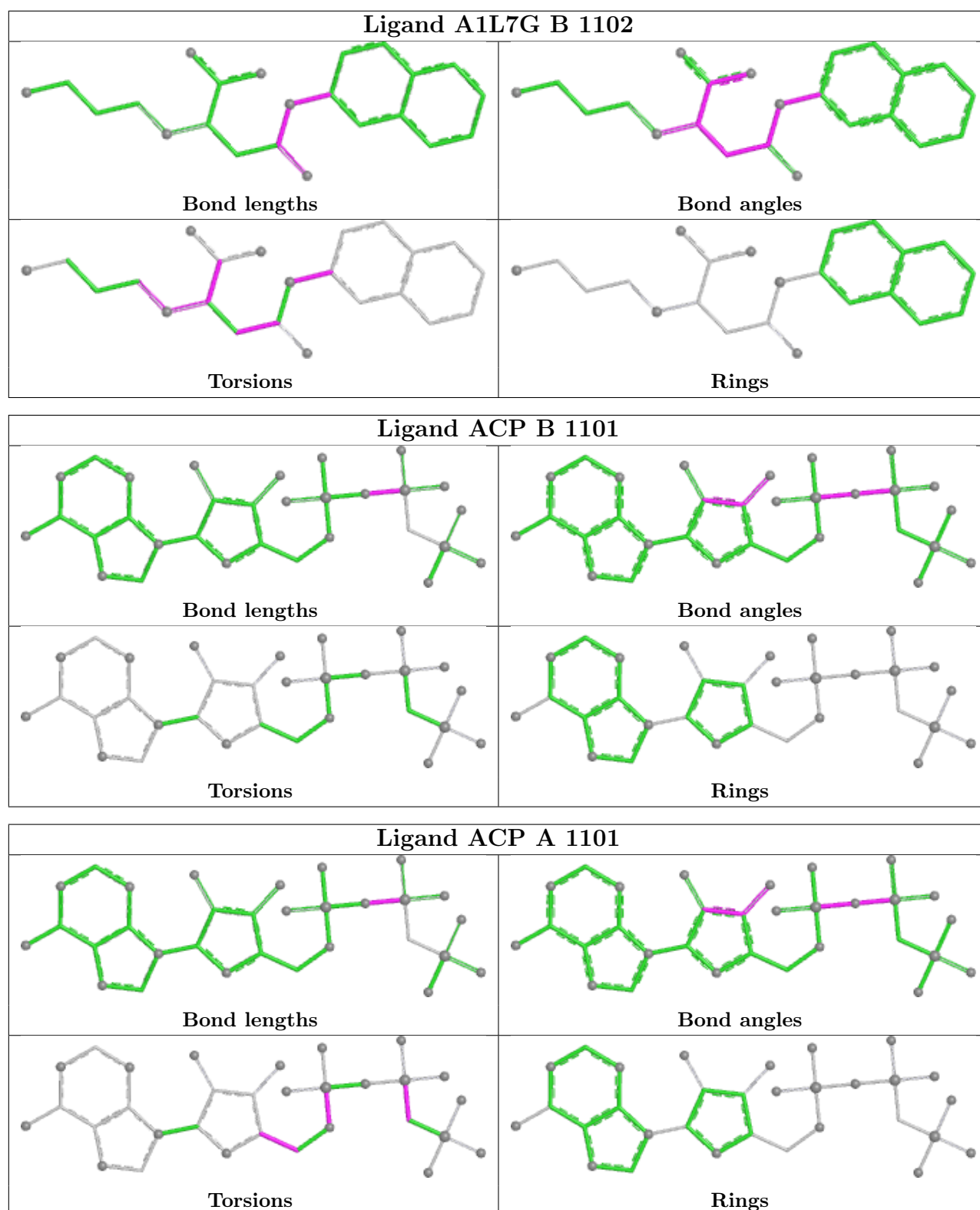
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1102	A1L7G	2	0
3	B	1102	A1L7G	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 ⓘ

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	206/218 (94%)	0.12	4 (1%) 66 62	25, 47, 68, 80	0
1	B	205/218 (94%)	0.23	3 (1%) 72 68	28, 50, 70, 87	0
1	C	206/218 (94%)	0.34	5 (2%) 59 55	33, 58, 74, 88	0
1	D	190/218 (87%)	0.44	15 (7%) 18 16	28, 57, 80, 89	0
All	All	807/872 (92%)	0.28	27 (3%) 49 45	25, 53, 75, 89	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	943	ALA	3.6
1	D	903	PRO	3.4
1	C	956	ASN	3.2
1	D	957	TYR	3.1
1	A	898	ASN	3.0
1	D	913	GLU	3.0
1	D	964	LYS	2.9
1	C	957	TYR	2.9
1	C	1019	LEU	2.8
1	D	911	LEU	2.6
1	D	904	CYS	2.5
1	D	882	MET	2.5
1	D	885	ILE	2.4
1	D	1026	ASP	2.4
1	D	908	PHE	2.4
1	C	897	ASP	2.4
1	A	961	ASN	2.4
1	D	960	SER	2.3
1	B	942	PRO	2.2
1	C	1053	PRO	2.2
1	B	944	ALA	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	956	ASN	2.1
1	D	963	TYR	2.1
1	D	880	TRP	2.1
1	A	1033	ASN	2.1
1	A	940	VAL	2.0
1	D	942	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

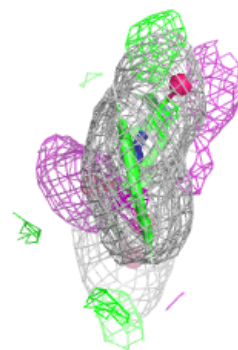
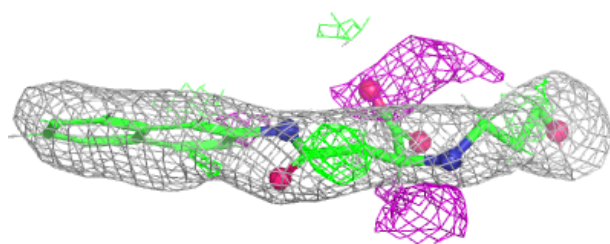
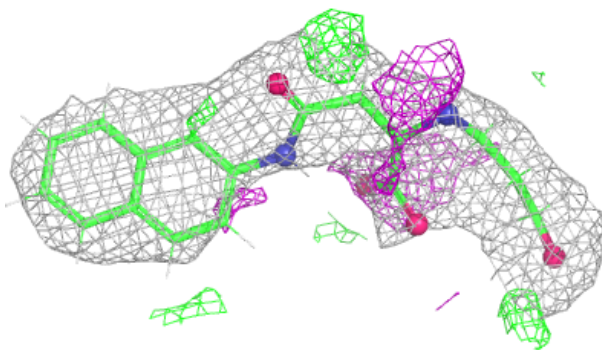
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	A1L7G	B	1102	23/23	0.83	0.14	34,50,61,76	0
3	A1L7G	A	1102	23/23	0.84	0.17	21,49,62,70	0
2	ACP	B	1101	31/31	0.91	0.08	40,61,77,91	0
2	ACP	A	1101	31/31	0.94	0.07	33,50,72,79	0

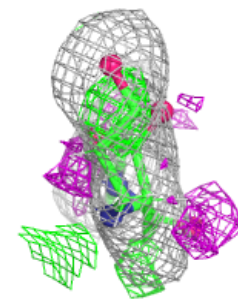
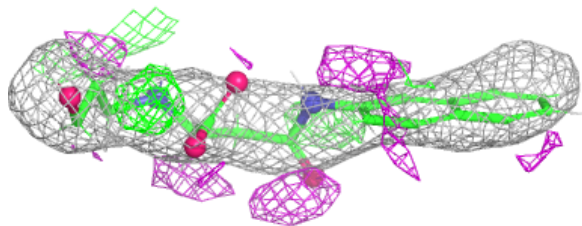
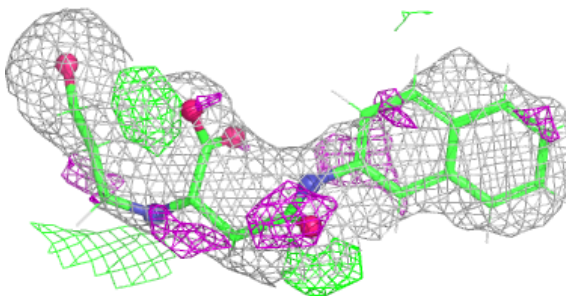
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A1L7G B 1102:

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

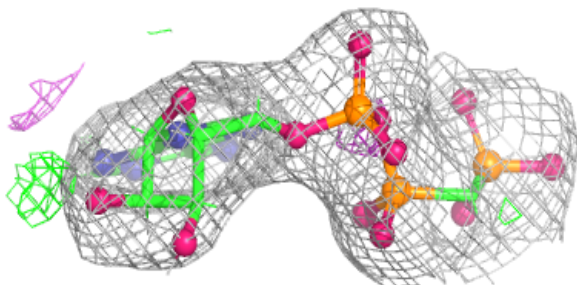
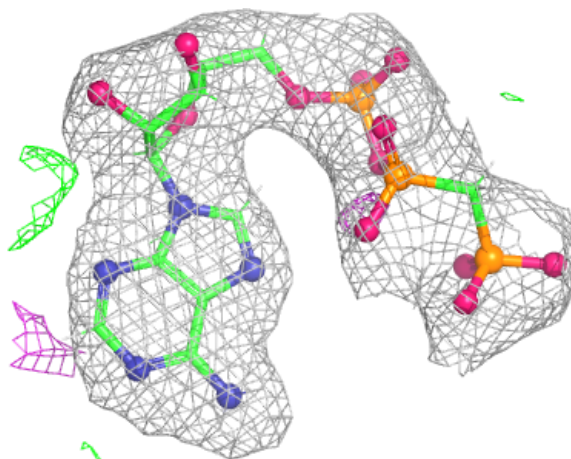
**Electron density around A1L7G A 1102:**

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



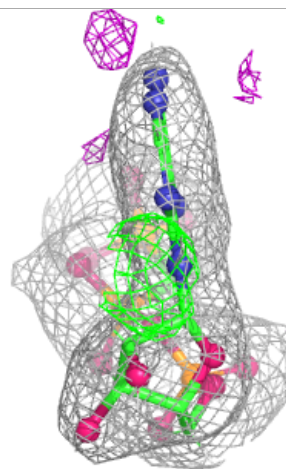
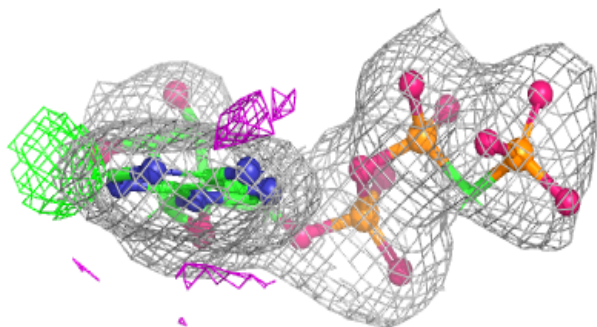
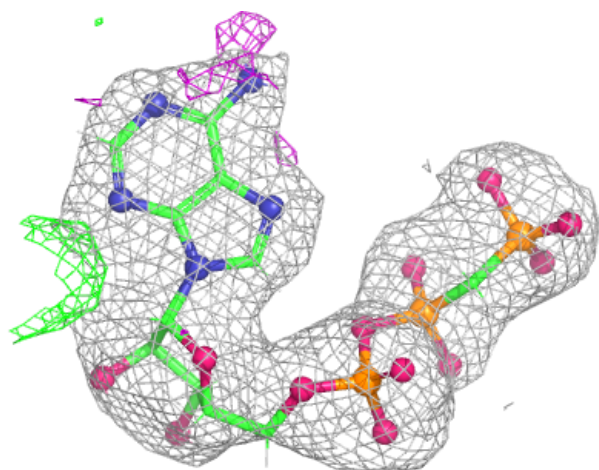
Electron density around ACP B 1101:

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



Electron density around ACP A 1101:

$2mF_o-DF_c$ (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.