



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:04 PM UTC

PDB ID : 4LWZ / pdb_00004lwz
Title : Crystal structure of Myo5b globular tail domain in complex with inactive Rab11a
Authors : Pylypenko, O.; Attanda, W.; Gauquelin, C.; Houdusse, A.
Deposited on : 2013-07-29
Resolution : 2.55 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

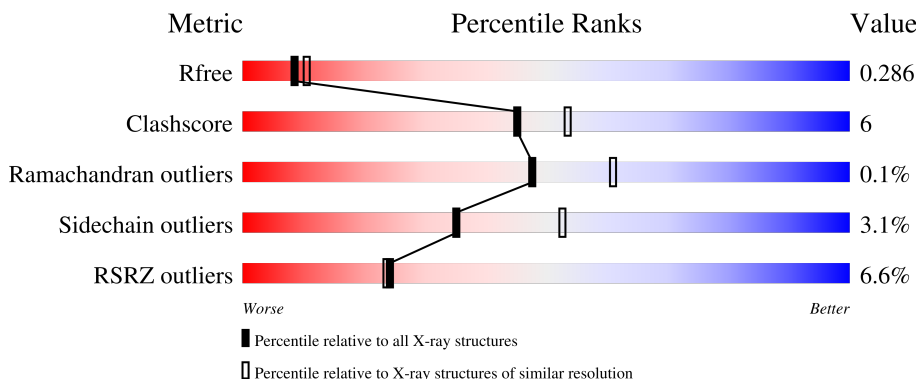
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	177	% 79% 11% • 8%
1	C	177	11% 71% 15% • 12%
2	B	427	% 78% 7% • 14%
2	D	427	10% 68% 12% 19%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8137 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 Ras-related protein Rab-11A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	0
			1265	803	215	246	1			
1	C	155	Total	C	N	O	S	0	0	0
			1174	739	206	228	1			

- Molecule 2 is a protein called Unconventional myosin-Vb.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	367	Total	C	N	O	S	0	1	0
			2903	1847	492	539	25			
2	D	346	Total	C	N	O	S	0	0	0
			2612	1651	447	493	21			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1422	MET	-	expression tag	UNP Q9ULV0
B	1423	ARG	-	expression tag	UNP Q9ULV0
B	1424	SER	-	expression tag	UNP Q9ULV0
B	1425	GLU	-	expression tag	UNP Q9ULV0
B	1426	THR	-	expression tag	UNP Q9ULV0
B	1427	MET	-	expression tag	UNP Q9ULV0
B	1428	SER	-	expression tag	UNP Q9ULV0
B	1429	TYR	-	expression tag	UNP Q9ULV0
B	1430	TYR	-	expression tag	UNP Q9ULV0
B	1431	HIS	-	expression tag	UNP Q9ULV0
B	1432	HIS	-	expression tag	UNP Q9ULV0
B	1433	HIS	-	expression tag	UNP Q9ULV0
B	1434	HIS	-	expression tag	UNP Q9ULV0
B	1435	HIS	-	expression tag	UNP Q9ULV0
B	1436	HIS	-	expression tag	UNP Q9ULV0
B	1437	ASP	-	expression tag	UNP Q9ULV0

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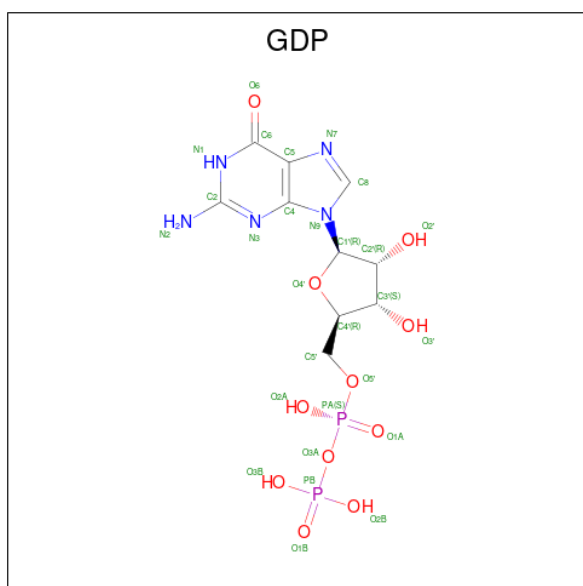
Chain	Residue	Modelled	Actual	Comment	Reference
B	1438	TYR	-	expression tag	UNP Q9ULV0
B	1439	ASP	-	expression tag	UNP Q9ULV0
B	1440	ILE	-	expression tag	UNP Q9ULV0
B	1441	PRO	-	expression tag	UNP Q9ULV0
B	1442	THR	-	expression tag	UNP Q9ULV0
B	1443	THR	-	expression tag	UNP Q9ULV0
B	1444	GLU	-	expression tag	UNP Q9ULV0
B	1445	ASN	-	expression tag	UNP Q9ULV0
B	1446	LEU	-	expression tag	UNP Q9ULV0
B	1447	TYR	-	expression tag	UNP Q9ULV0
B	1448	PHE	-	expression tag	UNP Q9ULV0
B	1449	GLN	-	expression tag	UNP Q9ULV0
B	1450	GLY	-	expression tag	UNP Q9ULV0
B	1451	ALA	-	expression tag	UNP Q9ULV0
B	1452	MET	-	expression tag	UNP Q9ULV0
B	1453	GLY	-	expression tag	UNP Q9ULV0
B	1454	SER	-	expression tag	UNP Q9ULV0
B	1455	MET	-	expression tag	UNP Q9ULV0
D	1422	MET	-	expression tag	UNP Q9ULV0
D	1423	ARG	-	expression tag	UNP Q9ULV0
D	1424	SER	-	expression tag	UNP Q9ULV0
D	1425	GLU	-	expression tag	UNP Q9ULV0
D	1426	THR	-	expression tag	UNP Q9ULV0
D	1427	MET	-	expression tag	UNP Q9ULV0
D	1428	SER	-	expression tag	UNP Q9ULV0
D	1429	TYR	-	expression tag	UNP Q9ULV0
D	1430	TYR	-	expression tag	UNP Q9ULV0
D	1431	HIS	-	expression tag	UNP Q9ULV0
D	1432	HIS	-	expression tag	UNP Q9ULV0
D	1433	HIS	-	expression tag	UNP Q9ULV0
D	1434	HIS	-	expression tag	UNP Q9ULV0
D	1435	HIS	-	expression tag	UNP Q9ULV0
D	1436	HIS	-	expression tag	UNP Q9ULV0
D	1437	ASP	-	expression tag	UNP Q9ULV0
D	1438	TYR	-	expression tag	UNP Q9ULV0
D	1439	ASP	-	expression tag	UNP Q9ULV0
D	1440	ILE	-	expression tag	UNP Q9ULV0
D	1441	PRO	-	expression tag	UNP Q9ULV0
D	1442	THR	-	expression tag	UNP Q9ULV0
D	1443	THR	-	expression tag	UNP Q9ULV0
D	1444	GLU	-	expression tag	UNP Q9ULV0
D	1445	ASN	-	expression tag	UNP Q9ULV0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1446	LEU	-	expression tag	UNP Q9ULV0
D	1447	TYR	-	expression tag	UNP Q9ULV0
D	1448	PHE	-	expression tag	UNP Q9ULV0
D	1449	GLN	-	expression tag	UNP Q9ULV0
D	1450	GLY	-	expression tag	UNP Q9ULV0
D	1451	ALA	-	expression tag	UNP Q9ULV0
D	1452	MET	-	expression tag	UNP Q9ULV0
D	1453	GLY	-	expression tag	UNP Q9ULV0
D	1454	SER	-	expression tag	UNP Q9ULV0
D	1455	MET	-	expression tag	UNP Q9ULV0

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



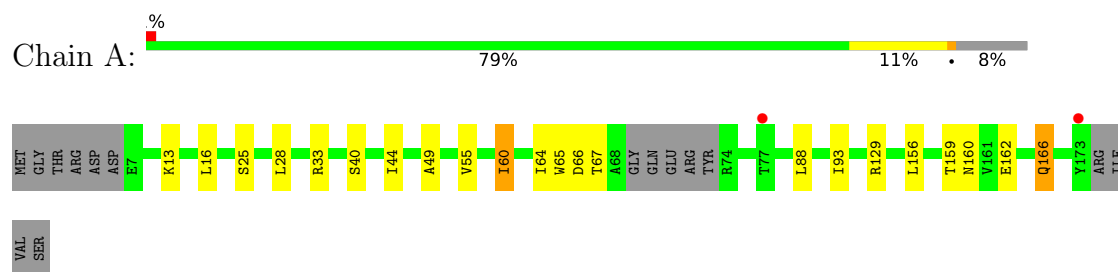
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total 33	O 33	0	0
5	B	49	Total 49	O 49	0	0
5	C	17	Total 17	O 17	0	0
5	D	26	Total 26	O 26	0	0

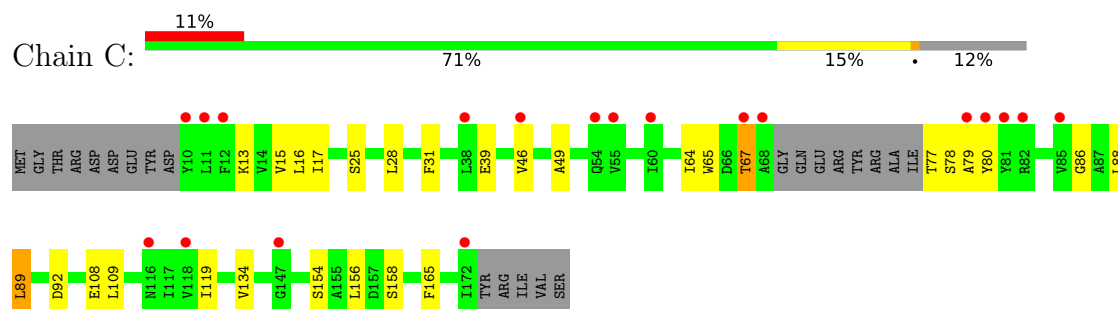
3 Residue-property plots [i](#)

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.

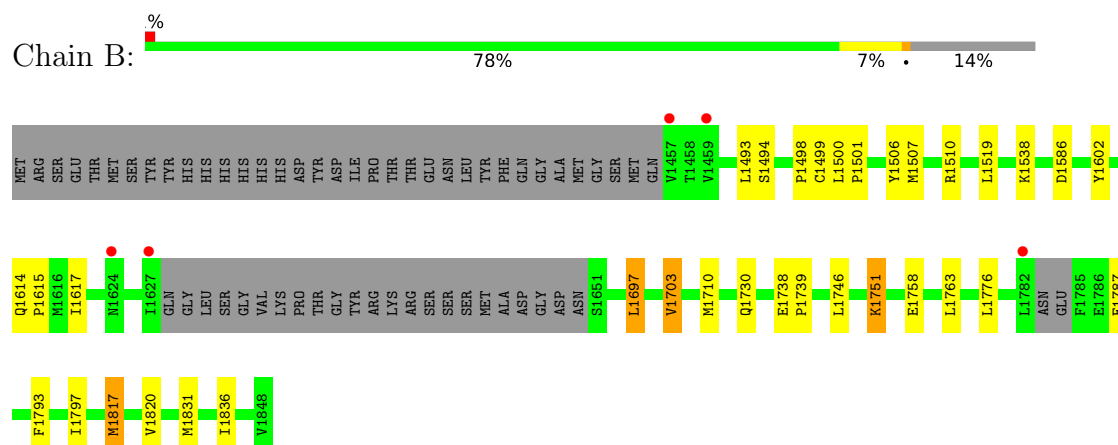
- Molecule 1: Ras-related protein Rab-11A



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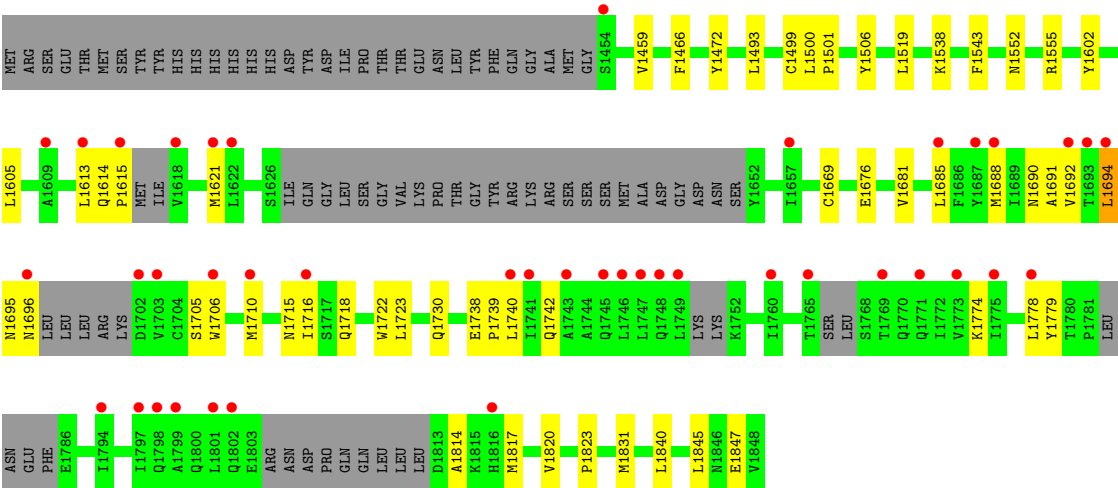


- Molecule 2: Unconventional myosin-Vb



- Molecule 2: Unconventional myosin-Vb





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.68Å 143.14Å 162.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.97 – 2.55 24.97 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.7 (24.97-2.55) 99.5 (24.97-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.231 , 0.295 0.226 , 0.286	Depositor DCC
R_{free} test set	2606 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	66.8	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 63.9	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	0.95	EDS
Total number of atoms	8137	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: MG, GDP

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	0.53	0/1285	0.79	1/1742 (0.1%)
1	C	0.54	0/1191	0.79	1/1614 (0.1%)
2	B	0.50	0/2955	0.74	1/4000 (0.0%)
2	D	0.51	0/2653	0.77	0/3598
All	All	0.51	0/8084	0.77	3/10954 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	67	THR	N-CA-C	5.51	117.47	108.76
2	B	1617	ILE	N-CA-C	5.39	115.83	110.23
1	C	134	VAL	CB-CA-C	-5.19	106.15	111.08

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	1265	0	1236	16	0
1	C	1174	0	1123	24	0
2	B	2903	0	2882	22	0
2	D	2612	0	2435	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	28	0	12	1	0
3	C	28	0	12	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	33	0	0	0	0
5	B	49	0	0	0	0
5	C	17	0	0	0	0
5	D	26	0	0	0	0
All	All	8137	0	7700	92	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 6.

All (92) 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:1602:TYR:CZ	2:D:1820:VAL:HG11	2.19	0.78
2:B:1602:TYR:CZ	2:B:1820:VAL:HG11	2.26	0.71
2:B:1507:MET:HG3	2:B:1831:MET:CE	2.25	0.67
1:C:13:LYS:HE3	1:C:65:TRP:CE2	2.32	0.64
2:D:1602:TYR:CE2	2:D:1820:VAL:HG11	2.33	0.63
1:A:55:VAL:HB	1:A:60:ILE:CD1	2.30	0.61
2:B:1507:MET:CG	2:B:1831:MET:CE	2.79	0.60
1:C:13:LYS:HE3	1:C:65:TRP:CD2	2.37	0.59
2:D:1506:TYR:CD1	2:D:1831:MET:HE2	2.38	0.58
1:C:156:LEU:C	1:C:156:LEU:HD23	2.29	0.57
2:D:1705:SER:HA	2:D:1779:TYR:CE1	2.39	0.56
1:C:109:LEU:HD23	1:C:119:ILE:HD13	1.88	0.56
2:D:1706:TRP:O	2:D:1710:MET:HG2	2.07	0.55
2:D:1614:GLN:HB3	2:D:1615:PRO:HD3	1.89	0.55
1:C:77:THR:C	1:C:79:ALA:N	2.63	0.55
2:D:1614:GLN:HB3	2:D:1615:PRO:CD	2.38	0.54
1:C:16:LEU:HD23	1:C:88:LEU:HB2	1.90	0.54
1:C:13:LYS:HG3	1:C:65:TRP:CD1	2.43	0.54
1:C:89:LEU:HD12	1:C:89:LEU:C	2.34	0.53
2:D:1688:MET:C	2:D:1690:ASN:N	2.65	0.53
2:B:1738:GLU:HA	2:B:1738:GLU:OE2	2.08	0.53
1:C:77:THR:O	1:C:78:SER:C	2.51	0.52
2:D:1555:ARG:O	2:D:1555:ARG:HD3	2.10	0.51
1:C:15:VAL:HB	1:C:80:TYR:CE2	2.46	0.51
2:D:1738:GLU:O	2:D:1742:GLN:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1493:LEU:HD21	2:B:1501:PRO:HD3	1.94	0.50
2:B:1614:GLN:N	2:B:1615:PRO:HD2	2.27	0.50
2:D:1543:PHE:CE1	2:D:1605:LEU:HD11	2.47	0.50
1:A:28:LEU:HD11	1:A:49:ALA:HB3	1.94	0.49
1:C:13:LYS:HG3	1:C:65:TRP:NE1	2.27	0.49
2:D:1506:TYR:HD1	2:D:1831:MET:CE	2.25	0.49
2:D:1774:LYS:O	2:D:1778:LEU:HG	2.11	0.49
2:B:1697:LEU:HD12	2:B:1703:VAL:HG13	1.94	0.49
1:C:92:ASP:OD1	1:C:92:ASP:C	2.56	0.49
1:C:109:LEU:CD2	1:C:119:ILE:HD13	2.44	0.48
2:D:1506:TYR:CD1	2:D:1506:TYR:C	2.91	0.48
2:B:1500:LEU:HB3	2:B:1501:PRO:HD3	1.95	0.48
1:A:159:THR:O	1:A:160:ASN:HB2	2.14	0.47
2:D:1506:TYR:CD1	2:D:1831:MET:CE	2.97	0.47
2:D:1681:VAL:O	2:D:1685:LEU:HG	2.14	0.47
1:A:28:LEU:HD13	1:A:64:ILE:HG21	1.96	0.47
1:A:156:LEU:C	1:A:156:LEU:HD23	2.40	0.47
2:D:1738:GLU:N	2:D:1739:PRO:CD	2.78	0.47
2:D:1692:VAL:O	2:D:1696:ASN:ND2	2.49	0.46
1:A:65:TRP:HZ3	2:B:1710:MET:HE1	1.81	0.46
1:A:40:SER:HB3	3:A:201:GDP:H3'	1.98	0.46
1:A:55:VAL:HB	1:A:60:ILE:HD13	1.97	0.46
2:B:1500:LEU:N	2:B:1501:PRO:CD	2.79	0.46
2:D:1715:ASN:O	2:D:1718:GLN:HB2	2.16	0.46
1:C:77:THR:O	1:C:80:TYR:N	2.49	0.45
2:B:1498:PRO:O	2:B:1499:CYS:HB2	2.17	0.45
2:D:1716:ILE:HD11	2:D:1740:LEU:HG	1.98	0.45
1:C:31:PHE:CE1	1:C:165:PHE:HB2	2.52	0.45
1:C:46:VAL:HG22	1:C:67:THR:CG2	2.47	0.43
2:D:1500:LEU:N	2:D:1501:PRO:CD	2.81	0.43
2:B:1602:TYR:CE2	2:B:1820:VAL:HG11	2.52	0.43
2:B:1817:MET:HE2	2:B:1817:MET:HB3	1.90	0.43
2:D:1613:LEU:O	2:D:1614:GLN:C	2.61	0.43
1:A:13:LYS:HE3	1:A:65:TRP:CE2	2.53	0.43
1:C:15:VAL:HG22	1:C:86:GLY:O	2.19	0.43
2:D:1694:LEU:O	2:D:1695:ASN:C	2.62	0.43
1:A:25:SER:OG	1:A:66:ASP:OD2	2.33	0.43
1:C:28:LEU:HD11	1:C:49:ALA:HB3	2.01	0.43
2:D:1621:MET:O	2:D:1715:ASN:ND2	2.51	0.43
1:C:31:PHE:CE1	1:C:165:PHE:CB	3.01	0.43
1:C:46:VAL:HG22	1:C:67:THR:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:1738:GLU:HB3	2:D:1739:PRO:HD3	2.00	0.43
1:A:16:LEU:HD23	1:A:88:LEU:HB2	2.01	0.42
1:A:162:GLU:O	1:A:166:GLN:HB2	2.20	0.42
2:D:1472:TYR:HB3	2:D:1845:LEU:CD2	2.49	0.42
2:B:1493:LEU:O	2:B:1494:SER:C	2.62	0.42
2:B:1507:MET:CG	2:B:1831:MET:HE3	2.49	0.42
1:A:55:VAL:HB	1:A:60:ILE:HD12	2.02	0.42
2:D:1499:CYS:HA	2:D:1552:ASN:OD1	2.20	0.42
1:A:129:ARG:HD2	2:D:1669:CYS:HB3	2.00	0.42
2:B:1586:ASP:OD1	2:B:1586:ASP:C	2.63	0.42
2:D:1466:PHE:CD1	2:D:1847:GLU:HB3	2.54	0.42
2:B:1738:GLU:HB3	2:B:1739:PRO:HD3	2.02	0.42
2:B:1506:TYR:CE2	2:B:1510:ARG:HG3	2.55	0.41
2:B:1751:LYS:HD2	2:B:1776:LEU:CD2	2.51	0.41
1:C:154:SER:O	1:C:158:SER:N	2.50	0.41
1:C:108:GLU:O	1:C:108:GLU:HG2	2.19	0.41
1:C:28:LEU:HD13	1:C:64:ILE:CG2	2.51	0.41
2:D:1722:TRP:CZ3	2:D:1723:LEU:HD23	2.56	0.41
2:D:1676:GLU:HB2	2:D:1823:PRO:HG3	2.03	0.41
1:A:13:LYS:HE3	1:A:65:TRP:CD2	2.55	0.41
2:B:1793:PHE:CE2	2:B:1797:ILE:HD11	2.56	0.41
1:A:44:ILE:HD13	2:B:1763:LEU:HD13	2.02	0.41
1:C:13:LYS:CE	1:C:65:TRP:CE2	3.01	0.41
2:B:1507:MET:HG3	2:B:1831:MET:HE2	2.00	0.40
2:D:1691:ALA:CB	2:D:1814:ALA:HB2	2.51	0.40
2:D:1459:VAL:O	2:D:1459:VAL:HG13	2.22	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	158/177 (89%)	153 (97%)	5 (3%)	0	100	100
1	C	151/177 (85%)	141 (93%)	10 (7%)	0	100	100
2	B	362/427 (85%)	345 (95%)	17 (5%)	0	100	100
2	D	330/427 (77%)	312 (94%)	17 (5%)	1 (0%)	36	45
All	All	1001/1208 (83%)	951 (95%)	49 (5%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	1694	LEU

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	132/150 (88%)	128 (97%)	4 (3%)	36	53
1	C	118/150 (79%)	113 (96%)	5 (4%)	26	40
2	B	321/391 (82%)	310 (97%)	11 (3%)	32	48
2	D	266/391 (68%)	260 (98%)	6 (2%)	44	62
All	All	837/1082 (77%)	811 (97%)	26 (3%)	35	53

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	A	60	ILE
1	A	93	ILE
1	A	166	GLN
2	B	1519	LEU
2	B	1538	LYS
2	B	1697	LEU
2	B	1703	VAL
2	B	1730	GLN
2	B	1746	LEU

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Mol	Chain	Res	Type
2	B	1751	LYS
2	B	1758	GLU
2	B	1787	GLU
2	B	1817	MET
2	B	1836	ILE
1	C	17	ILE
1	C	25	SER
1	C	39	GLU
1	C	67	THR
1	C	89	LEU
2	D	1493	LEU
2	D	1519	LEU
2	D	1538	LYS
2	D	1730	GLN
2	D	1817	MET
2	D	1840	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	1511	HIS
2	B	1516	ASN
2	B	1539	HIS
2	B	1604	GLN
2	B	1624	ASN
2	B	1727	ASN
2	B	1730	GLN
2	D	1467	GLN
2	D	1511	HIS
2	D	1539	HIS
2	D	1558	HIS
2	D	1584	ASN
2	D	1624	ASN
2	D	1730	GLN
2	D	1816	HIS

5.3.3 RNA ⓘ

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	GDP	C	201	4	29,30,30	1.13	3 (10%)	45,47,47	1.74	8 (17%)
3	GDP	A	201	4	29,30,30	1.17	4 (13%)	45,47,47	1.84	8 (17%)

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	GDP	C	201	4	-	0/16/32/32	0/3/3/3
3	GDP	A	201	4	-	1/16/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	201	GDP	C5-C4	2.76	1.46	1.38
3	C	201	GDP	C6-N1	-2.60	1.34	1.38
3	A	201	GDP	C5-C4	2.49	1.45	1.38
3	C	201	GDP	C4-N9	-2.38	1.32	1.38
3	A	201	GDP	C6-N1	-2.37	1.34	1.38
3	A	201	GDP	C4-N9	-2.33	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	201	GDP	C5-N7	-2.02	1.35	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201	GDP	C5-C4-N3	-5.34	119.89	128.39
3	C	201	GDP	C5-C4-N3	-5.28	119.99	128.39
3	A	201	GDP	C2-N3-C4	5.17	121.20	112.30
3	C	201	GDP	C2-N3-C4	4.88	120.71	112.30
3	C	201	GDP	N9-C4-N3	4.02	133.99	125.95
3	A	201	GDP	C6-C5-N7	3.96	137.50	130.29
3	C	201	GDP	C6-C5-N7	3.56	136.77	130.29
3	A	201	GDP	N9-C4-N3	3.32	132.59	125.95
3	A	201	GDP	C4-C5-N7	-2.98	105.95	110.67
3	A	201	GDP	C2'-C3'-C4'	2.40	107.24	102.61
3	A	201	GDP	O2A-PA-O3A	2.34	113.59	107.27
3	C	201	GDP	C4-C5-N7	-2.15	107.26	110.67
3	C	201	GDP	C2'-C1'-N9	-2.15	107.27	113.25
3	C	201	GDP	O2A-PA-O3A	2.09	112.92	107.27
3	C	201	GDP	O6-C6-C5	-2.07	121.07	126.53
3	A	201	GDP	C8-N7-C5	2.04	107.90	104.26

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	201	GDP	PA-O3A-PB-O2B

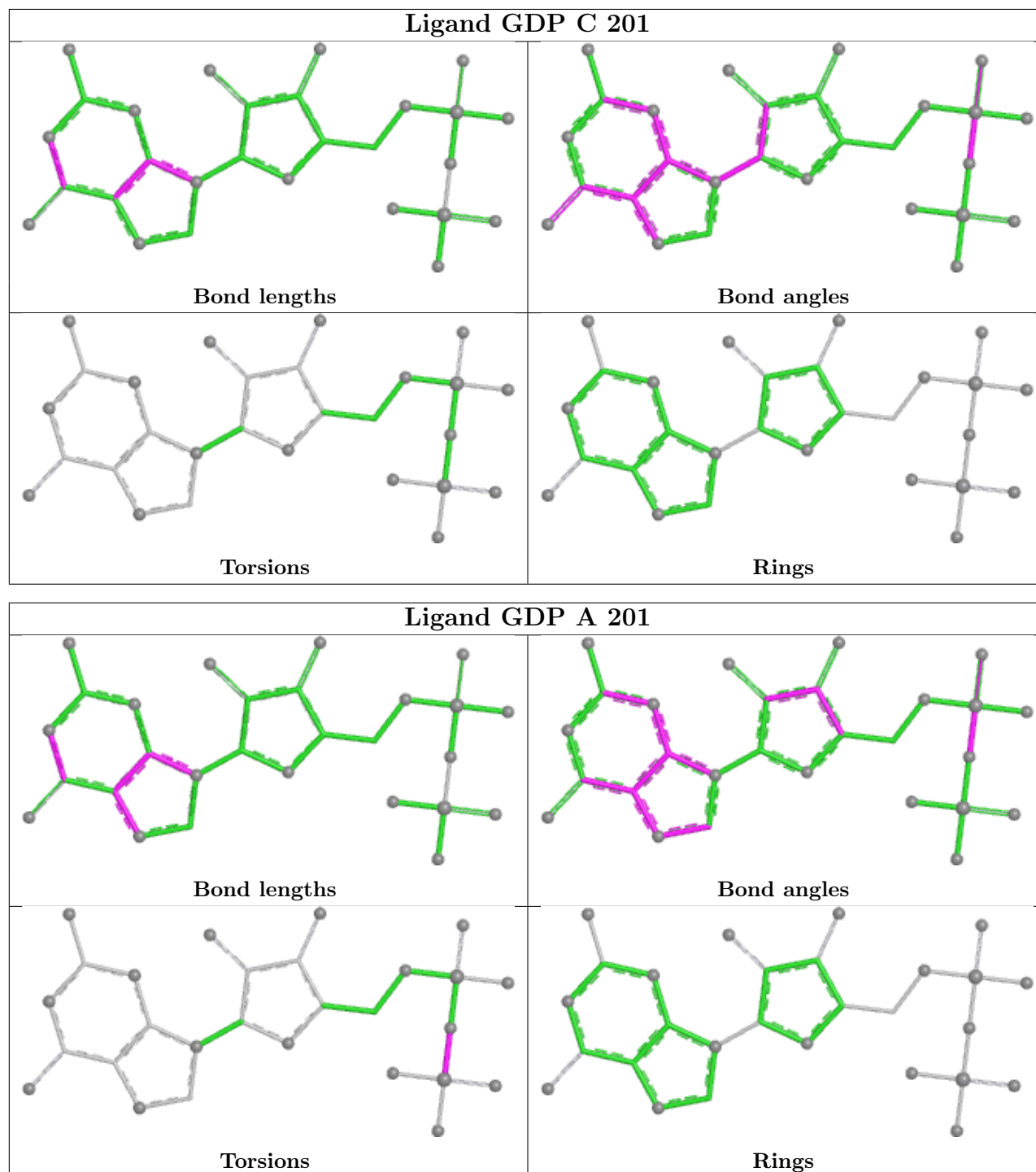
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	201	GDP	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	162/177 (91%)	-0.07	2 (1%) 76 77	52, 76, 128, 162	0
1	C	155/177 (87%)	0.75	19 (12%) 8 7	60, 97, 141, 181	0
2	B	367/427 (85%)	-0.10	5 (1%) 73 74	42, 81, 140, 188	1 (0%)
2	D	346/427 (81%)	0.62	42 (12%) 8 7	59, 98, 146, 178	0
All	All	1030/1208 (85%)	0.28	68 (6%) 24 24	42, 90, 142, 188	1 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	1710	MET	5.4
2	D	1741	ILE	5.2
1	C	79	ALA	5.0
2	D	1747	LEU	4.2
2	D	1797	ILE	4.1
2	D	1771	GLN	3.8
2	B	1782	LEU	3.6
2	D	1778	LEU	3.4
2	D	1748	GLN	3.4
2	D	1622	LEU	3.4
2	D	1696	ASN	3.2
1	C	55	VAL	3.2
1	C	60	ILE	3.1
2	D	1794	ILE	3.1
2	D	1743	ALA	3.1
2	D	1749	LEU	3.1
1	C	118	VAL	3.0
2	D	1613	LEU	3.0
1	C	68	ALA	3.0
2	D	1702	ASP	2.9
2	D	1454	SER	2.9

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Mol	Chain	Res	Type	RSRZ
2	D	1760	ILE	2.8
2	D	1746	LEU	2.8
2	D	1801	LEU	2.8
2	B	1624	ASN	2.8
1	C	11	LEU	2.8
2	D	1694	LEU	2.8
2	D	1706	TRP	2.7
1	C	46	VAL	2.7
2	D	1618	VAL	2.7
2	D	1769	THR	2.7
1	C	172	ILE	2.7
2	B	1459	VAL	2.6
2	D	1687	TYR	2.6
2	D	1609	ALA	2.6
2	D	1816	HIS	2.6
2	D	1799	ALA	2.6
1	C	82	ARG	2.5
1	A	173	TYR	2.4
2	D	1745	GLN	2.4
2	B	1457	VAL	2.4
2	D	1775	ILE	2.4
1	C	12	PHE	2.4
2	D	1765	THR	2.3
2	D	1740	LEU	2.3
1	C	85	VAL	2.3
2	D	1621	MET	2.3
2	D	1685	LEU	2.2
1	C	10	TYR	2.2
1	C	147	GLY	2.2
2	B	1627	ILE	2.2
1	C	54	GLN	2.2
1	A	77	THR	2.1
1	C	67	THR	2.1
1	C	81	TYR	2.1
2	D	1692	VAL	2.1
1	C	38	LEU	2.1
2	D	1657	ILE	2.1
2	D	1716	ILE	2.1
1	C	80	TYR	2.1
2	D	1773	VAL	2.0
1	C	116	ASN	2.0
2	D	1615	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
2	D	1798	GLN	2.0
2	D	1802	GLN	2.0
2	D	1703	VAL	2.0
2	D	1693	THR	2.0
2	D	1688	MET	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](#)

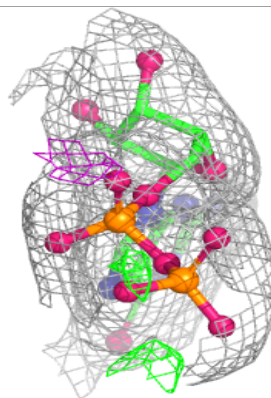
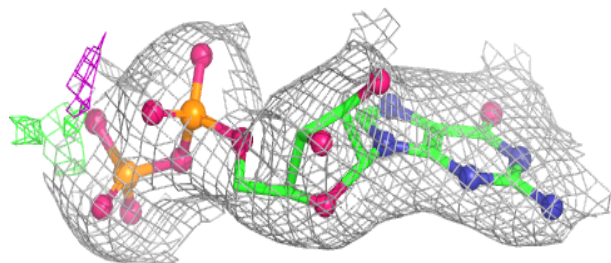
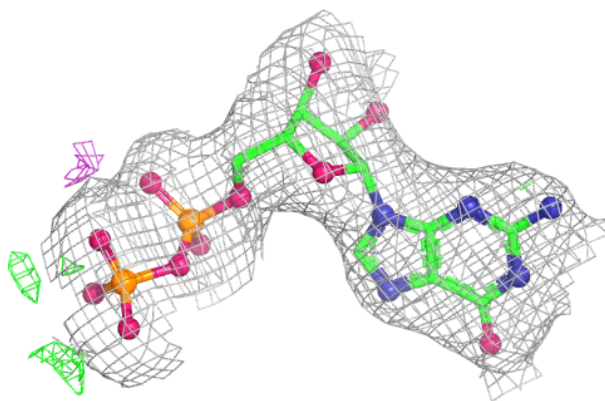
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	C	202	1/1	0.81	0.11	90,90,90,90	0
3	GDP	C	201	28/28	0.97	0.06	56,70,84,96	0
3	GDP	A	201	28/28	0.98	0.05	47,61,72,83	0
4	MG	A	202	1/1	0.99	0.08	74,74,74,74	0

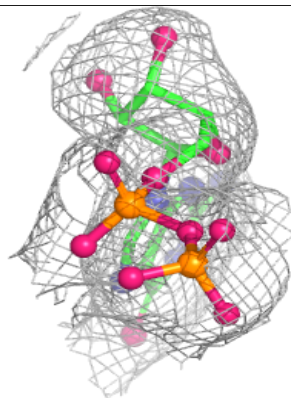
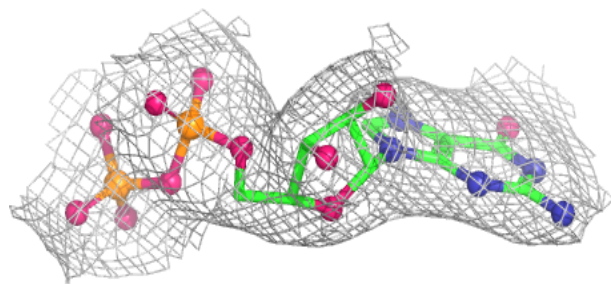
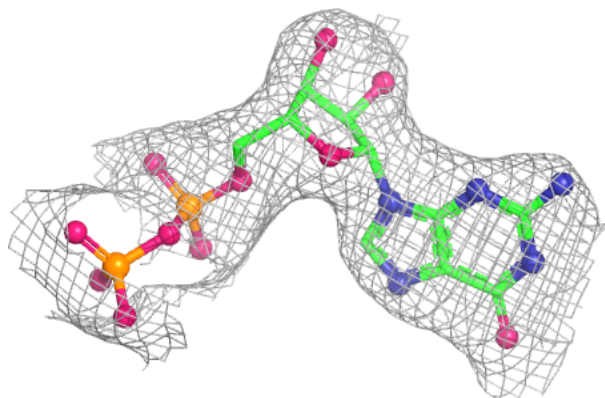
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GDP C 201:

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

**Electron density around GDP A 201:**

$2mF_o-DF_c$ (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.