



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:51 PM UTC

PDB ID : 4M8N / pdb_00004m8n
Title : Crystal Structure of PlexinC1/Rap1B Complex
Authors : Pascoe, H.G.; Wang, Y.; Brautigam, C.A.; He, H.; Zhang, X.
Deposited on : 2013-08-13
Resolution : 3.29 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

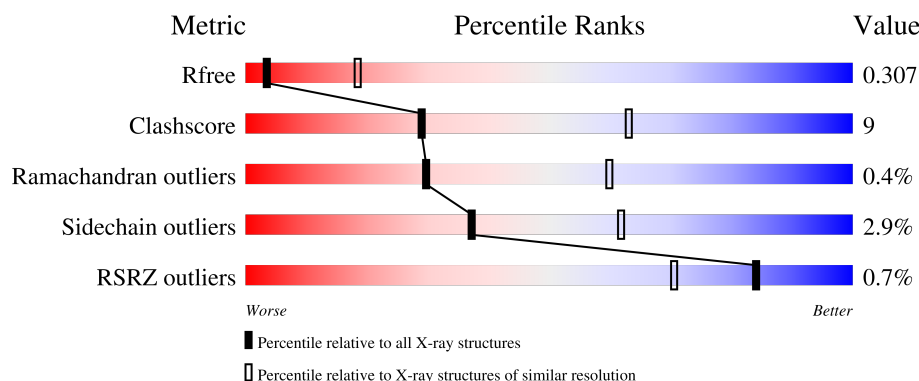
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	599	% 77% 18% • •
1	B	599	71% 23% • 5%
1	C	599	73% 19% • 7%
1	D	599	% 72% 20% • 7%
2	E	199	% 70% 14% • 16%

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Mol	Chain	Length	Quality of chain
2	F	199	63%19%17%
2	G	199	% 66%18%16%
2	H	199	% 69%14%16%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22228 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 PlexinC1 Intracellular Region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	574	Total	C	N	O	S	0	0	0
			4356	2818	717	801	20			
1	B	571	Total	C	N	O	S	0	0	0
			4262	2756	702	784	20			
1	C	560	Total	C	N	O	S	0	0	0
			4218	2719	699	780	20			
1	D	560	Total	C	N	O	S	0	0	0
			4230	2738	695	777	20			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	549	GLY	-	expression tag	UNP Q5RGW1
A	550	GLY	-	expression tag	UNP Q5RGW1
A	551	SER	-	expression tag	UNP Q5RGW1
B	549	GLY	-	expression tag	UNP Q5RGW1
B	550	GLY	-	expression tag	UNP Q5RGW1
B	551	SER	-	expression tag	UNP Q5RGW1
C	549	GLY	-	expression tag	UNP Q5RGW1
C	550	GLY	-	expression tag	UNP Q5RGW1
C	551	SER	-	expression tag	UNP Q5RGW1
D	549	GLY	-	expression tag	UNP Q5RGW1
D	550	GLY	-	expression tag	UNP Q5RGW1
D	551	SER	-	expression tag	UNP Q5RGW1

- Molecule 2 is a protein called Ras-related protein Rap-1b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	168	Total	C	N	O	S	0	0	0
			1277	805	215	250	7			
2	F	166	Total	C	N	O	S	0	0	0
			1262	797	213	245	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	168	Total	C	N	O	S	0	0	0
			1258	793	211	247	7			
2	H	167	Total	C	N	O	S	0	0	0
			1223	779	206	232	6			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	expression tag	UNP P61224
E	-1	PRO	-	expression tag	UNP P61224
E	0	HIS	-	expression tag	UNP P61224
E	167	SER	-	expression tag	UNP P61224
E	168	GLY	-	expression tag	UNP P61224
E	169	GLY	-	expression tag	UNP P61224
E	170	SER	-	expression tag	UNP P61224
E	171	GLY	-	expression tag	UNP P61224
E	172	SER	-	expression tag	UNP P61224
E	173	GLY	-	expression tag	UNP P61224
E	174	SER	-	expression tag	UNP P61224
E	175	SER	-	expression tag	UNP P61224
E	176	GLY	-	expression tag	UNP P61224
E	177	GLY	-	expression tag	UNP P61224
E	178	SER	-	expression tag	UNP P61224
E	179	GLY	-	expression tag	UNP P61224
E	180	SER	-	expression tag	UNP P61224
E	181	GLY	-	expression tag	UNP P61224
E	182	GLY	-	expression tag	UNP P61224
E	183	GLY	-	expression tag	UNP P61224
E	184	SER	-	expression tag	UNP P61224
E	185	GLY	-	expression tag	UNP P61224
E	186	SER	-	expression tag	UNP P61224
E	187	GLY	-	expression tag	UNP P61224
E	188	SER	-	expression tag	UNP P61224
E	189	SER	-	expression tag	UNP P61224
E	190	GLY	-	expression tag	UNP P61224
E	191	LEU	-	expression tag	UNP P61224
E	192	PRO	-	expression tag	UNP P61224
E	193	GLU	-	expression tag	UNP P61224
E	194	THR	-	expression tag	UNP P61224
E	195	GLY	-	expression tag	UNP P61224
E	196	GLY	-	expression tag	UNP P61224
F	-2	GLY	-	expression tag	UNP P61224

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	PRO	-	expression tag	UNP P61224
F	0	HIS	-	expression tag	UNP P61224
F	167	SER	-	expression tag	UNP P61224
F	168	GLY	-	expression tag	UNP P61224
F	169	GLY	-	expression tag	UNP P61224
F	170	SER	-	expression tag	UNP P61224
F	171	GLY	-	expression tag	UNP P61224
F	172	SER	-	expression tag	UNP P61224
F	173	GLY	-	expression tag	UNP P61224
F	174	SER	-	expression tag	UNP P61224
F	175	SER	-	expression tag	UNP P61224
F	176	GLY	-	expression tag	UNP P61224
F	177	GLY	-	expression tag	UNP P61224
F	178	SER	-	expression tag	UNP P61224
F	179	GLY	-	expression tag	UNP P61224
F	180	SER	-	expression tag	UNP P61224
F	181	GLY	-	expression tag	UNP P61224
F	182	GLY	-	expression tag	UNP P61224
F	183	GLY	-	expression tag	UNP P61224
F	184	SER	-	expression tag	UNP P61224
F	185	GLY	-	expression tag	UNP P61224
F	186	SER	-	expression tag	UNP P61224
F	187	GLY	-	expression tag	UNP P61224
F	188	SER	-	expression tag	UNP P61224
F	189	SER	-	expression tag	UNP P61224
F	190	GLY	-	expression tag	UNP P61224
F	191	LEU	-	expression tag	UNP P61224
F	192	PRO	-	expression tag	UNP P61224
F	193	GLU	-	expression tag	UNP P61224
F	194	THR	-	expression tag	UNP P61224
F	195	GLY	-	expression tag	UNP P61224
F	196	GLY	-	expression tag	UNP P61224
G	-2	GLY	-	expression tag	UNP P61224
G	-1	PRO	-	expression tag	UNP P61224
G	0	HIS	-	expression tag	UNP P61224
G	167	SER	-	expression tag	UNP P61224
G	168	GLY	-	expression tag	UNP P61224
G	169	GLY	-	expression tag	UNP P61224
G	170	SER	-	expression tag	UNP P61224
G	171	GLY	-	expression tag	UNP P61224
G	172	SER	-	expression tag	UNP P61224
G	173	GLY	-	expression tag	UNP P61224

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Chain	Residue	Modelled	Actual	Comment	Reference
G	174	SER	-	expression tag	UNP P61224
G	175	SER	-	expression tag	UNP P61224
G	176	GLY	-	expression tag	UNP P61224
G	177	GLY	-	expression tag	UNP P61224
G	178	SER	-	expression tag	UNP P61224
G	179	GLY	-	expression tag	UNP P61224
G	180	SER	-	expression tag	UNP P61224
G	181	GLY	-	expression tag	UNP P61224
G	182	GLY	-	expression tag	UNP P61224
G	183	GLY	-	expression tag	UNP P61224
G	184	SER	-	expression tag	UNP P61224
G	185	GLY	-	expression tag	UNP P61224
G	186	SER	-	expression tag	UNP P61224
G	187	GLY	-	expression tag	UNP P61224
G	188	SER	-	expression tag	UNP P61224
G	189	SER	-	expression tag	UNP P61224
G	190	GLY	-	expression tag	UNP P61224
G	191	LEU	-	expression tag	UNP P61224
G	192	PRO	-	expression tag	UNP P61224
G	193	GLU	-	expression tag	UNP P61224
G	194	THR	-	expression tag	UNP P61224
G	195	GLY	-	expression tag	UNP P61224
G	196	GLY	-	expression tag	UNP P61224
H	-2	GLY	-	expression tag	UNP P61224
H	-1	PRO	-	expression tag	UNP P61224
H	0	HIS	-	expression tag	UNP P61224
H	167	SER	-	expression tag	UNP P61224
H	168	GLY	-	expression tag	UNP P61224
H	169	GLY	-	expression tag	UNP P61224
H	170	SER	-	expression tag	UNP P61224
H	171	GLY	-	expression tag	UNP P61224
H	172	SER	-	expression tag	UNP P61224
H	173	GLY	-	expression tag	UNP P61224
H	174	SER	-	expression tag	UNP P61224
H	175	SER	-	expression tag	UNP P61224
H	176	GLY	-	expression tag	UNP P61224
H	177	GLY	-	expression tag	UNP P61224
H	178	SER	-	expression tag	UNP P61224
H	179	GLY	-	expression tag	UNP P61224
H	180	SER	-	expression tag	UNP P61224
H	181	GLY	-	expression tag	UNP P61224
H	182	GLY	-	expression tag	UNP P61224

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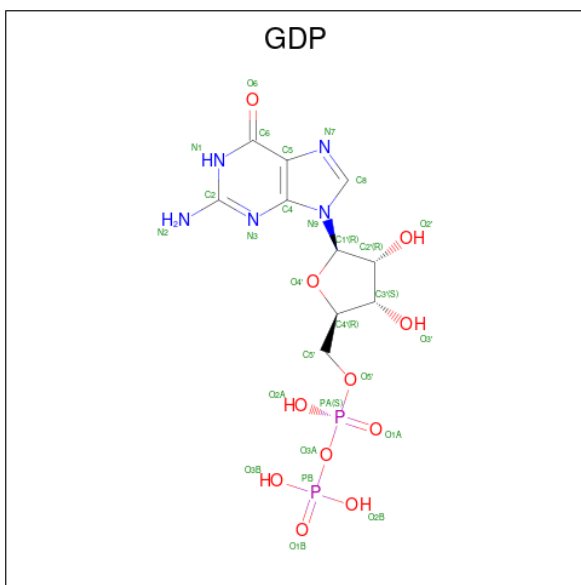
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Chain	Residue	Modelled	Actual	Comment	Reference
H	183	GLY	-	expression tag	UNP P61224
H	184	SER	-	expression tag	UNP P61224
H	185	GLY	-	expression tag	UNP P61224
H	186	SER	-	expression tag	UNP P61224
H	187	GLY	-	expression tag	UNP P61224
H	188	SER	-	expression tag	UNP P61224
H	189	SER	-	expression tag	UNP P61224
H	190	GLY	-	expression tag	UNP P61224
H	191	LEU	-	expression tag	UNP P61224
H	192	PRO	-	expression tag	UNP P61224
H	193	GLU	-	expression tag	UNP P61224
H	194	THR	-	expression tag	UNP P61224
H	195	GLY	-	expression tag	UNP P61224
H	196	GLY	-	expression tag	UNP P61224

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

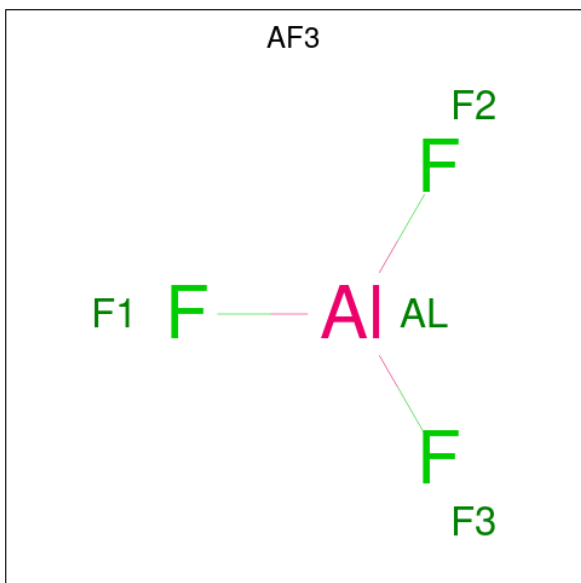
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total Mg 1 1	0	0
3	F	1	Total Mg 1 1	0	0
3	G	1	Total Mg 1 1	0	0
3	H	1	Total Mg 1 1	0	0

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	E	1	Total 28	C 10	N 5	O 11	P 2	0	0
4	F	1	Total 28	C 10	N 5	O 11	P 2	0	0
4	G	1	Total 28	C 10	N 5	O 11	P 2	0	0
4	H	1	Total 28	C 10	N 5	O 11	P 2	0	0

- Molecule 5 is ALUMINUM FLUORIDE (CCD ID: AF3) (formula: AlF_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	Al	F	0	0
			4	1	3		
5	F	1	Total	Al	F	0	0
			4	1	3		
5	G	1	Total	Al	F	0	0
			4	1	3		
5	H	1	Total	Al	F	0	0
			4	1	3		

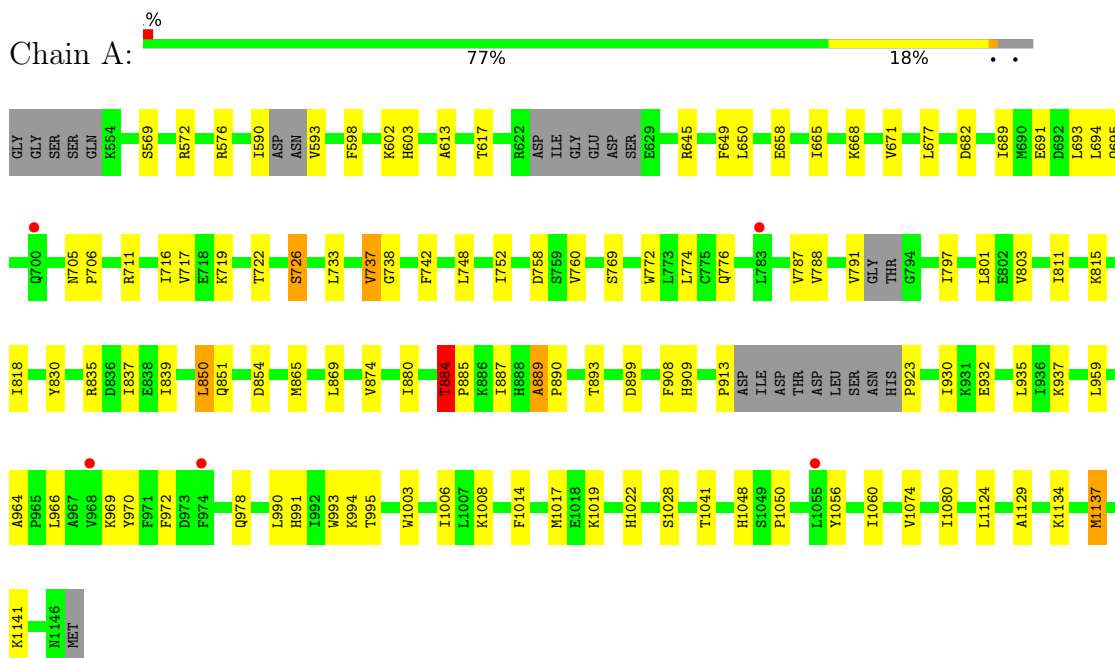
- Molecule 6 is water.

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

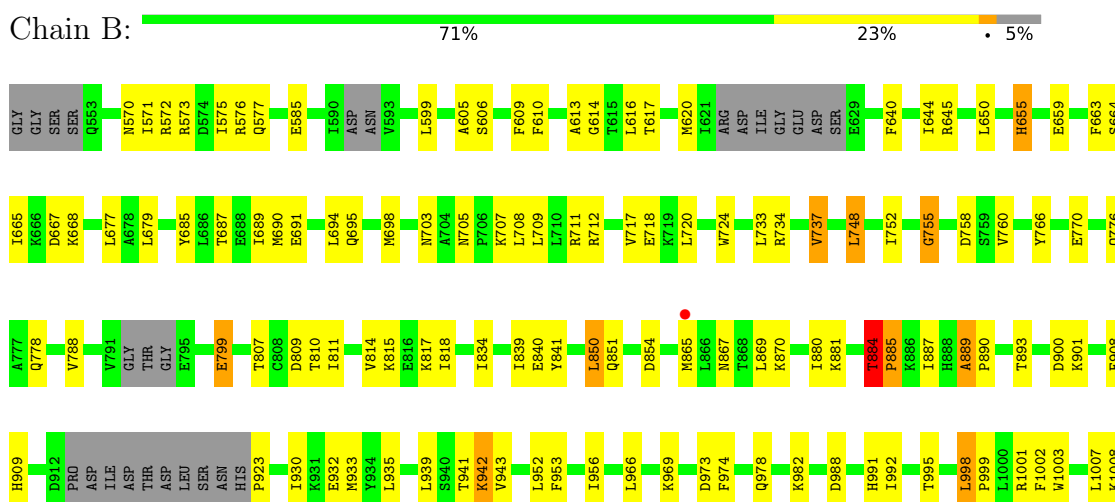
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PlexinC1 Intracellular Region



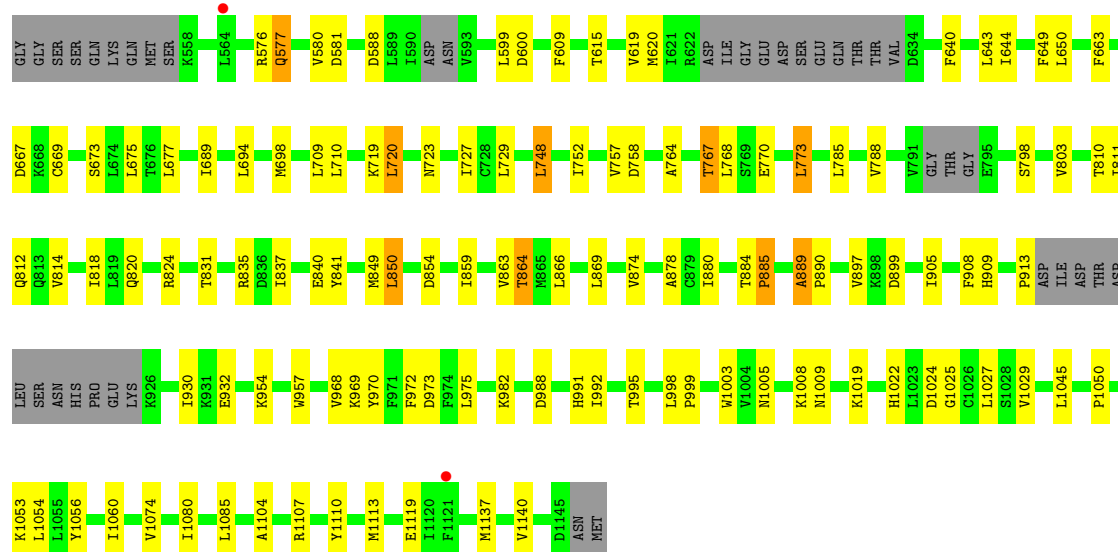
• Molecule 1: PlexinC1 Intracellular Region





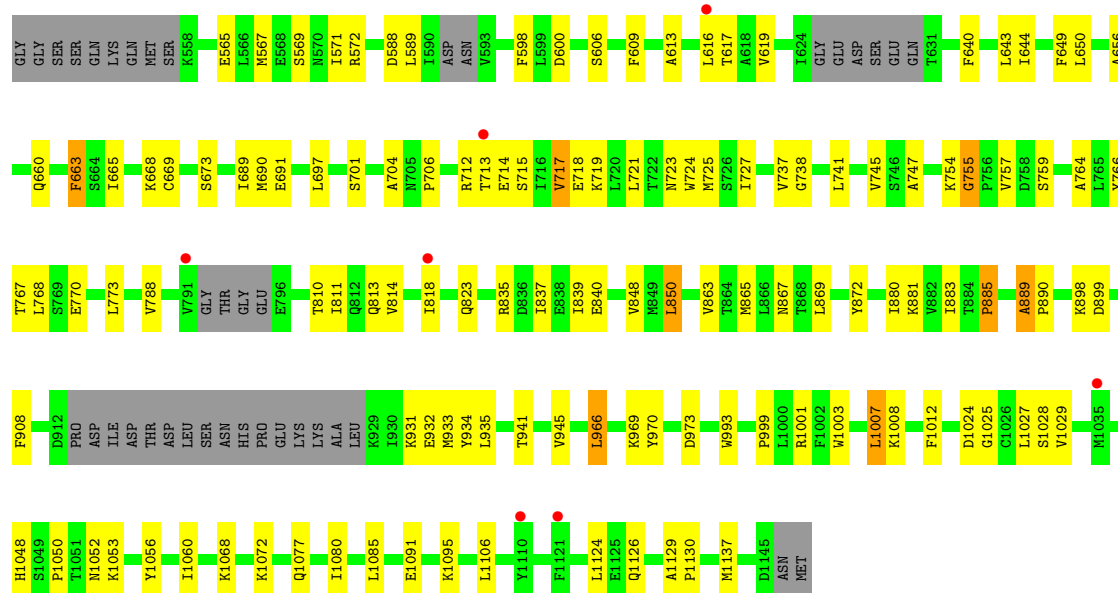
• Molecule 1: PlexinC1 Intracellular Region

Chain C: 73% 19% 7%



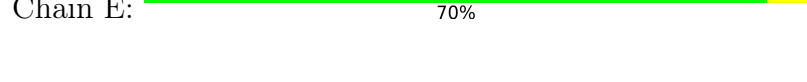
• Molecule 1: PlexinC1 Intracellular Region

Chain D: 72% 20% 7%



• Molecule 2: Ras-related protein Rap-1b

Chain E: 70% 14% 16%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	76.28Å 84.73Å 138.75Å 91.09° 95.15° 90.32°	Depositor
Resolution (Å)	41.05 – 3.29 41.05 – 3.29	Depositor EDS
% Data completeness (in resolution range)	90.2 (41.05-3.29) 90.2 (41.05-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 3.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.243 , 0.299 (Not available) , 0.307	Depositor DCC
R_{free} test set	1804 reflections (3.45%)	wwPDB-VP
Wilson B-factor (Å ²)	111.8	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 100.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.033 for -h,k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	22228	wwPDB-VP
Average B, all atoms (Å ²)	143.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.70% 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: AF3, 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	0.34	0/4437	0.77	6/6023 (0.1%)
1	B	0.34	0/4341	0.79	9/5905 (0.2%)
1	C	0.33	0/4298	0.75	3/5845 (0.1%)
1	D	0.31	0/4312	0.73	5/5861 (0.1%)
2	E	0.30	0/1293	0.75	1/1750 (0.1%)
2	F	0.29	0/1278	0.73	1/1730 (0.1%)
2	G	0.28	0/1274	0.71	1/1729 (0.1%)
2	H	0.29	0/1239	0.73	1/1682 (0.1%)
All	All	0.32	0/22472	0.76	27/30525 (0.1%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	755	GLY	CA-C-N	7.84	127.89	119.90
1	B	755	GLY	C-N-CA	7.84	127.89	119.90
1	C	890	PRO	N-CA-CB	7.38	110.59	103.51
1	C	885	PRO	N-CA-CB	7.33	110.27	103.46
1	A	890	PRO	N-CA-CB	7.19	110.41	103.51
1	D	890	PRO	N-CA-CB	7.08	110.68	103.25
1	D	885	PRO	N-CA-CB	7.00	110.19	103.19
1	C	913	PRO	N-CA-CB	6.80	110.48	103.00
1	A	885	PRO	N-CA-CB	6.79	110.38	103.25
1	B	923	PRO	N-CA-CB	6.79	110.47	103.00
1	B	885	PRO	N-CA-CB	6.76	110.35	103.25
1	A	913	PRO	N-CA-CB	6.75	110.43	103.00
1	A	923	PRO	N-CA-CB	6.74	110.42	103.00
1	B	890	PRO	N-CA-CB	6.34	110.55	103.44
2	H	63	GLN	N-CA-C	-6.13	105.96	113.50
1	D	755	GLY	CA-C-N	6.05	126.08	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	755	GLY	C-N-CA	6.05	126.08	119.90
1	B	998	LEU	CA-C-N	-6.01	112.33	119.84
1	B	998	LEU	C-N-CA	-6.01	112.33	119.84
1	A	705	ASN	CA-C-N	5.99	127.33	119.84
1	A	705	ASN	C-N-CA	5.99	127.33	119.84
1	B	705	ASN	CA-C-N	5.58	126.82	119.84
1	B	705	ASN	C-N-CA	5.58	126.82	119.84
2	G	63	GLN	N-CA-C	-5.55	106.64	113.41
2	E	66	ALA	N-CA-C	5.39	117.60	109.41
2	F	63	GLN	N-CA-C	-5.10	106.56	112.89
1	D	1012	PHE	N-CA-C	-5.04	106.91	113.16

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	4356	0	4143	62	0
1	B	4262	0	3980	98	0
1	C	4218	0	3952	71	0
1	D	4230	0	3971	80	0
2	E	1277	0	1229	16	0
2	F	1262	0	1212	29	0
2	G	1258	0	1188	23	0
2	H	1223	0	1153	22	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	E	28	0	12	0	0
4	F	28	0	12	3	0
4	G	28	0	12	2	0
4	H	28	0	12	5	0
5	E	4	0	0	1	0
5	F	4	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	4	0	0	0	0
5	H	4	0	0	0	0
6	A	4	0	0	0	0
6	B	1	0	0	0	0
6	C	3	0	0	0	0
6	D	1	0	0	0	0
6	H	1	0	0	0	0
All	All	22228	0	20876	375	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 9.

All (375) 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:577:GLN:NE2	1:C:581:ASP:OD2	2.21	0.74
1:C:811:ILE:HD12	1:C:850:LEU:HD13	1.71	0.73
1:A:835:ARG:NH1	1:A:899:ASP:OD2	2.22	0.72
1:B:811:ILE:HD12	1:B:850:LEU:HD13	1.70	0.72
2:H:60:GLY:O	2:H:68:ARG:NH1	2.22	0.71
2:G:45:GLU:HA	2:G:50:GLN:HA	1.73	0.71
1:C:859:ILE:HG12	1:C:864:THR:HG23	1.72	0.71
1:A:691:GLU:OE2	1:A:970:TYR:OH	2.06	0.70
1:A:932:GLU:HA	1:A:935:LEU:HG	1.74	0.69
1:D:835:ARG:NH2	1:D:899:ASP:OD2	2.24	0.69
1:B:1110:TYR:HA	1:B:1113:MET:HB3	1.76	0.68
1:B:695:GLN:HG3	1:B:978:GLN:HE22	1.59	0.68
1:D:1091:GLU:HG2	1:D:1095:LYS:HD3	1.76	0.67
1:B:1095:LYS:HE3	1:D:1130:PRO:HG3	1.78	0.66
1:D:1008:LYS:NZ	1:D:1024:ASP:OD1	2.28	0.66
1:A:760:VAL:HG21	1:A:908:PHE:HA	1.75	0.66
1:B:576:ARG:NH2	1:C:770:GLU:OE2	2.28	0.66
1:A:811:ILE:HD12	1:A:850:LEU:HD13	1.78	0.66
1:B:841:TYR:HB2	1:B:850:LEU:HD23	1.78	0.65
2:H:82:TYR:HB3	2:H:93:LEU:HD11	1.79	0.65
1:D:932:GLU:HA	1:D:935:LEU:HD13	1.78	0.65
2:F:82:TYR:HB3	2:F:93:LEU:HD11	1.78	0.65
1:D:1050:PRO:HG2	1:D:1053:LYS:HB2	1.77	0.64
1:D:706:PRO:O	1:D:993:TRP:NE1	2.28	0.64
1:D:718:GLU:OE1	1:D:1001:ARG:NH1	2.31	0.64
1:A:803:VAL:HG11	1:A:818:ILE:HG22	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:885:PRO:HA	1:C:889:ALA:HB3	1.80	0.64
1:B:809:ASP:OD2	1:B:817:LYS:NZ	2.22	0.64
1:D:650:LEU:HD23	1:D:689:ILE:HG12	1.79	0.64
1:B:570:ASN:OD1	1:B:573:ARG:NH2	2.30	0.63
1:D:669:CYS:HA	1:D:719:LYS:HG2	1.80	0.63
1:A:811:ILE:HG22	1:A:815:LYS:HE3	1.81	0.63
1:B:885:PRO:HA	1:B:889:ALA:HB3	1.79	0.63
2:E:25:GLN:HB2	2:E:27:ILE:HG22	1.81	0.63
2:H:40:TYR:HB2	2:H:55:ILE:HB	1.79	0.63
1:C:640:PHE:HD1	1:C:1140:VAL:HG11	1.64	0.62
1:B:999:PRO:HA	1:B:1003:TRP:HB2	1.81	0.62
2:F:9:LEU:HB3	2:F:96:LEU:HD23	1.81	0.62
2:F:77:GLY:HA2	2:F:110:PRO:HB2	1.82	0.61
1:C:767:THR:OG1	1:C:768:LEU:N	2.33	0.61
1:A:966:LEU:HD12	1:A:1080:ILE:HD13	1.81	0.61
1:A:969:LYS:NZ	1:A:1074:VAL:O	2.24	0.61
1:B:650:LEU:HD23	1:B:689:ILE:HG12	1.81	0.61
1:D:840:GLU:HB2	1:D:881:LYS:HB2	1.81	0.61
1:A:1056:TYR:O	1:A:1060:ILE:HG12	2.01	0.60
1:C:999:PRO:HA	1:C:1003:TRP:HB2	1.83	0.60
2:G:141:CYS:SG	2:G:142:ALA:N	2.75	0.60
1:B:606:SER:HB2	1:B:617:THR:HB	1.83	0.60
1:D:606:SER:HB2	1:D:617:THR:HB	1.83	0.60
1:D:1001:ARG:NH2	2:H:62:GLU:O	2.33	0.60
1:B:690:MET:HE1	1:B:724:TRP:CD2	2.36	0.60
1:B:1054:LEU:HD21	1:C:580:VAL:HG22	1.84	0.60
1:D:668:LYS:HZ3	1:D:714:GLU:C	2.10	0.60
1:B:640:PHE:HD2	1:B:1106:LEU:HD22	1.67	0.60
2:E:82:TYR:HB3	2:E:93:LEU:HD11	1.83	0.59
2:G:93:LEU:HD13	2:G:113:LEU:HD11	1.85	0.59
2:G:85:THR:HG23	2:G:123:GLU:HB3	1.84	0.59
1:D:770:GLU:HG2	1:D:931:LYS:HE2	1.84	0.58
1:D:613:ALA:HB1	1:D:616:LEU:HB2	1.85	0.58
1:D:588:ASP:OD1	1:D:754:LYS:NZ	2.30	0.58
1:B:810:THR:HG23	1:B:909:HIS:CE1	2.39	0.58
1:D:598:PHE:HE1	1:D:738:GLY:HA3	1.69	0.58
1:C:863:VAL:HG12	1:C:908:PHE:HD1	1.69	0.58
1:B:748:LEU:O	1:B:752:ILE:HG13	2.03	0.58
1:D:811:ILE:HD12	1:D:850:LEU:HD13	1.85	0.58
1:B:734:ARG:HH21	1:D:1126:GLN:HA	1.69	0.58
1:C:600:ASP:OD1	1:C:600:ASP:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:807:THR:HB	1:B:870:LYS:HG3	1.85	0.57
2:F:97:ARG:HH11	2:F:111:MET:HE3	1.69	0.57
2:F:61:THR:O	2:F:68:ARG:NH2	2.37	0.57
1:A:590:ILE:HG23	1:A:593:VAL:HG23	1.87	0.57
1:D:759:SER:O	1:D:898:LYS:NZ	2.37	0.57
1:A:649:PHE:HE1	1:A:1137:MET:HE3	1.70	0.57
1:B:576:ARG:HB2	1:C:932:GLU:HG2	1.86	0.57
1:B:1036:ASP:OD2	1:B:1053:LYS:NZ	2.37	0.57
2:G:84:ILE:HD12	2:G:124:ARG:HE	1.68	0.56
1:D:1056:TYR:O	1:D:1060:ILE:HG12	2.05	0.56
1:A:576:ARG:HB2	1:D:932:GLU:HG2	1.86	0.56
1:C:988:ASP:O	1:C:992:ILE:HG12	2.06	0.56
1:D:565:GLU:O	1:D:569:SER:OG	2.22	0.56
1:B:707:LYS:HA	1:B:992:ILE:HG21	1.87	0.56
1:A:990:LEU:HD22	1:A:994:LYS:HE3	1.87	0.56
1:C:1008:LYS:NZ	1:C:1024:ASP:OD1	2.39	0.56
2:F:84:ILE:HD11	2:F:118:CYS:HA	1.88	0.56
1:B:644:ILE:O	1:B:685:TYR:OH	2.22	0.55
1:B:703:ASN:HD22	1:B:712:ARG:NH1	2.04	0.55
2:E:106:THR:OG1	2:E:107:ASP:N	2.27	0.55
1:A:737:VAL:HG21	1:A:1003:TRP:HH2	1.70	0.55
1:A:1129:ALA:HB3	1:A:1134:LYS:HE3	1.88	0.55
1:C:1019:LYS:HE3	1:C:1024:ASP:HB2	1.88	0.55
1:D:966:LEU:HD23	1:D:1080:ILE:HD13	1.89	0.55
1:A:851:GLN:HE21	1:A:854:ASP:HB3	1.71	0.55
2:E:45:GLU:HA	2:E:50:GLN:HA	1.89	0.55
2:G:53:LEU:HD11	2:G:162:VAL:HG21	1.89	0.54
1:C:972:PHE:HD2	1:C:1074:VAL:HG11	1.72	0.54
2:H:117:LYS:HG2	4:H:201:GDP:C5	2.43	0.54
2:E:41:ARG:NH1	2:E:54:GLU:OE2	2.39	0.54
1:C:748:LEU:O	1:C:752:ILE:HG13	2.08	0.54
1:D:867:ASN:ND2	1:D:872:TYR:OH	2.40	0.54
2:E:11:SER:O	2:E:16:LYS:NZ	2.41	0.54
2:H:53:LEU:HD23	2:H:55:ILE:HD11	1.89	0.54
1:D:713:THR:HG21	2:H:62:GLU:HB3	1.90	0.53
2:F:35:THR:HB	5:F:202:AF3:F2	1.97	0.53
2:F:99:GLN:HG3	2:F:102:ARG:HH21	1.73	0.53
1:B:605:ALA:HB1	1:B:609:PHE:CE2	2.43	0.53
2:E:19:LEU:HD23	2:E:148:ALA:HB2	1.90	0.53
2:F:118:CYS:HB3	2:F:145:GLU:HG2	1.91	0.53
1:D:885:PRO:HA	1:D:889:ALA:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1009:ASN:HD21	2:F:65:THR:HB	1.74	0.53
1:D:609:PHE:HB3	1:D:673:SER:HB3	1.90	0.52
1:D:767:THR:OG1	1:D:768:LEU:N	2.42	0.52
1:B:900:ASP:OD1	1:B:901:LYS:N	2.43	0.52
1:B:942:LYS:HG3	1:B:1055:LEU:HD22	1.91	0.52
2:H:119:ASP:OD1	2:H:119:ASP:N	2.43	0.52
1:C:991:HIS:O	1:C:995:THR:HG23	2.09	0.52
1:D:643:LEU:HD11	1:D:1137:MET:HA	1.92	0.52
2:H:119:ASP:OD2	4:H:201:GDP:N2	2.40	0.52
1:A:645:ARG:NE	1:A:682:ASP:OD2	2.41	0.52
1:B:942:LYS:HE3	1:B:1056:TYR:CE2	2.45	0.51
1:A:1050:PRO:HD3	2:E:40:TYR:CE2	2.45	0.51
1:B:1007:LEU:HD23	1:B:1031:ALA:HA	1.92	0.51
1:A:671:VAL:HG11	1:A:716:ILE:HD13	1.93	0.51
1:C:620:MET:HB3	1:C:677:LEU:HD21	1.92	0.51
1:A:706:PRO:O	1:A:993:TRP:NE1	2.42	0.51
1:A:1124:LEU:HG	1:A:1129:ALA:HB2	1.92	0.51
1:C:729:LEU:HD21	1:C:968:VAL:HG23	1.91	0.51
1:C:810:THR:O	1:C:814:VAL:HG23	2.10	0.51
1:A:1008:LYS:O	1:A:1019:LYS:NZ	2.39	0.51
1:B:969:LYS:HE3	1:B:973:ASP:OD2	2.10	0.51
2:F:5:LYS:HB2	2:F:76:GLN:HG3	1.92	0.51
1:B:655:HIS:O	1:B:659:GLU:HG2	2.11	0.51
1:D:589:LEU:HD13	1:D:747:ALA:HB1	1.93	0.51
1:B:737:VAL:HG11	1:B:1003:TRP:CZ3	2.46	0.51
1:B:620:MET:HB3	1:B:677:LEU:HD21	1.91	0.50
1:D:837:ILE:HA	1:D:883:ILE:O	2.11	0.50
1:A:752:ILE:HD13	1:A:1022:HIS:CE1	2.46	0.50
1:B:884:THR:H	1:B:887:ILE:HD12	1.76	0.50
1:B:889:ALA:O	1:B:893:THR:OG1	2.28	0.50
1:B:1035:MET:HE3	2:G:34:PRO:HB2	1.92	0.50
1:B:953:PHE:CZ	1:B:1034:PHE:HB2	2.47	0.50
1:D:721:LEU:HD11	1:D:725:MET:HE2	1.93	0.50
1:A:748:LEU:O	1:A:752:ILE:HG13	2.11	0.50
1:B:770:GLU:OE2	1:C:576:ARG:NH2	2.44	0.50
1:B:760:VAL:HG21	1:B:908:PHE:HA	1.94	0.50
2:G:69:ASP:OD1	2:G:69:ASP:N	2.42	0.50
1:A:884:THR:H	1:A:887:ILE:HD12	1.77	0.50
1:D:941:THR:O	1:D:945:VAL:HG22	2.12	0.50
1:D:1048:HIS:HB3	2:H:24:VAL:HG12	1.94	0.50
1:A:1048:HIS:HB3	2:E:25:GLN:HG2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:737:VAL:HG21	1:B:1003:TRP:HH2	1.76	0.49
1:D:715:SER:HB2	1:D:717:VAL:HG23	1.93	0.49
2:G:19:LEU:HD23	2:G:148:ALA:HB2	1.93	0.49
1:A:791:VAL:H	1:A:797:ILE:HB	1.76	0.49
2:F:118:CYS:SG	2:F:147:SER:HB2	2.52	0.49
1:C:835:ARG:NH1	1:C:899:ASP:OD2	2.33	0.49
1:C:841:TYR:HB2	1:C:850:LEU:HD23	1.95	0.49
1:C:954:LYS:HG2	1:C:957:TRP:HE1	1.77	0.49
1:C:1025:GLY:O	1:C:1029:VAL:HG23	2.13	0.49
1:B:752:ILE:HD13	1:B:1022:HIS:CE1	2.47	0.48
1:B:1056:TYR:O	1:B:1060:ILE:HG12	2.13	0.48
1:C:969:LYS:NZ	1:C:1074:VAL:O	2.27	0.48
2:F:100:ILE:HG21	2:F:111:MET:HE1	1.94	0.48
1:B:865:MET:HB2	1:B:908:PHE:CZ	2.48	0.48
1:D:1068:LYS:O	1:D:1072:LYS:HG2	2.13	0.48
1:A:602:LYS:HG3	1:A:603:HIS:CD2	2.49	0.48
2:E:124:ARG:NH1	2:E:145:GLU:OE2	2.41	0.48
1:B:854:ASP:OD1	1:B:867:ASN:ND2	2.36	0.48
1:C:650:LEU:HD23	1:C:689:ILE:HG12	1.96	0.48
1:C:969:LYS:HE3	1:C:973:ASP:OD2	2.13	0.48
2:G:15:GLY:HA3	2:G:116:ASN:HD22	1.77	0.48
2:G:9:LEU:HD21	2:G:71:TYR:CE1	2.49	0.48
1:B:815:LYS:HB3	1:B:834:ILE:HD11	1.96	0.48
1:D:704:ALA:HA	1:D:712:ARG:HH22	1.79	0.48
1:A:711:ARG:NE	5:E:202:AF3:F3	2.37	0.48
1:A:737:VAL:HG11	1:A:1003:TRP:CZ3	2.49	0.48
1:B:952:LEU:HD21	1:B:1007:LEU:HD11	1.95	0.48
1:B:969:LYS:NZ	1:B:1074:VAL:O	2.34	0.48
1:D:600:ASP:N	1:D:600:ASP:OD1	2.46	0.48
2:H:16:LYS:NZ	4:H:201:GDP:O1B	2.41	0.48
1:C:1110:TYR:HA	1:C:1113:MET:HB3	1.95	0.48
1:B:930:ILE:HB	1:B:933:MET:HE3	1.95	0.47
1:B:758:ASP:OD1	1:B:909:HIS:HB3	2.13	0.47
1:A:722:THR:O	1:A:726:SER:OG	2.32	0.47
1:C:752:ILE:HD13	1:C:1022:HIS:CE1	2.49	0.47
1:B:687:THR:O	1:B:691:GLU:HG3	2.13	0.47
1:C:1104:ALA:HA	1:C:1107:ARG:HD2	1.97	0.47
2:F:8:VAL:O	2:F:58:THR:OG1	2.22	0.47
1:C:669:CYS:SG	2:F:62:GLU:HG2	2.55	0.47
1:D:814:VAL:HG21	1:D:869:LEU:HD21	1.95	0.47
1:B:839:ILE:HG23	1:B:880:ILE:HG23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:119:ASP:OD1	2:F:119:ASP:N	2.47	0.47
1:D:660:GLN:HB2	1:D:663:PHE:HB2	1.97	0.47
1:D:1025:GLY:O	1:D:1028:SER:OG	2.25	0.47
1:B:613:ALA:O	1:B:617:THR:HG23	2.15	0.47
1:C:609:PHE:HB3	1:C:673:SER:HB3	1.96	0.47
1:A:733:LEU:HD21	1:A:1006:ILE:HG12	1.97	0.46
1:A:865:MET:HE2	1:A:865:MET:HB3	1.88	0.46
1:D:773:LEU:HD21	1:D:934:TYR:HB2	1.96	0.46
1:B:733:LEU:HD12	1:B:737:VAL:HB	1.96	0.46
1:B:1050:PRO:HD3	2:G:40:TYR:CE2	2.51	0.46
1:D:640:PHE:O	1:D:644:ILE:HG12	2.16	0.46
1:B:998:LEU:O	1:B:1002:PHE:HB3	2.16	0.46
2:H:79:ALA:HB2	2:H:161:LEU:HD11	1.96	0.46
1:D:613:ALA:O	1:D:617:THR:HG23	2.15	0.46
2:E:84:ILE:HD12	2:E:124:ARG:HG3	1.96	0.46
1:A:658:GLU:OE2	1:A:716:ILE:N	2.34	0.46
1:A:889:ALA:O	1:A:893:THR:OG1	2.34	0.46
1:D:644:ILE:HG23	1:D:650:LEU:HD13	1.98	0.46
1:D:999:PRO:HA	1:D:1003:TRP:HB2	1.98	0.46
2:H:117:LYS:HE2	4:H:201:GDP:C8	2.51	0.46
1:A:665:ILE:HA	1:A:668:LYS:HD2	1.98	0.46
1:D:970:TYR:HB2	1:D:1080:ILE:HD12	1.98	0.46
1:C:1119:GLU:OE1	1:C:1119:GLU:N	2.49	0.45
1:D:1124:LEU:HG	1:D:1129:ALA:HB2	1.97	0.45
1:B:644:ILE:HG21	1:B:679:LEU:HD21	1.98	0.45
1:B:663:PHE:CE1	1:B:667:ASP:HB3	2.52	0.45
2:H:135:ALA:HB2	2:H:143:PHE:HB2	1.97	0.45
1:D:788:VAL:N	1:D:880:ILE:O	2.40	0.45
1:A:650:LEU:HD23	1:A:689:ILE:HG12	1.99	0.45
1:D:724:TRP:HA	1:D:727:ILE:HD12	1.99	0.45
2:F:136:ARG:HD3	2:F:140:ASN:HA	1.98	0.45
1:A:830:TYR:HD2	1:A:837:ILE:HD11	1.82	0.45
2:F:119:ASP:OD2	4:F:201:GDP:N2	2.42	0.45
1:A:758:ASP:OD1	1:A:909:HIS:HB3	2.17	0.45
2:F:161:LEU:O	2:F:165:ILE:HG13	2.17	0.45
1:B:664:SER:O	1:B:668:LYS:HG3	2.16	0.44
1:C:785:LEU:HD13	1:C:880:ILE:HD12	1.99	0.44
1:C:840:GLU:HA	1:C:849:MET:HA	1.99	0.44
1:C:1056:TYR:O	1:C:1060:ILE:HG12	2.17	0.44
1:D:690:MET:HE1	1:D:724:TRP:CD2	2.52	0.44
1:D:713:THR:HA	1:D:718:GLU:OE1	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:972:PHE:HD2	1:A:1074:VAL:HG11	1.82	0.44
1:B:694:LEU:HD22	1:B:974:PHE:HE2	1.83	0.44
2:E:96:LEU:O	2:E:100:ILE:HG12	2.17	0.44
2:F:116:ASN:OD1	2:F:117:LYS:N	2.48	0.44
1:A:613:ALA:O	1:A:617:THR:HG23	2.18	0.44
1:A:839:ILE:HG23	1:A:880:ILE:HG23	2.00	0.44
1:C:788:VAL:HG13	1:C:798:SER:HA	1.99	0.44
2:G:96:LEU:O	2:G:100:ILE:HG12	2.17	0.44
1:A:930:ILE:HG23	1:D:572:ARG:HG2	1.99	0.44
1:A:937:LYS:HA	1:A:937:LYS:HD3	1.66	0.44
1:B:708:LEU:HD23	1:B:712:ARG:HH21	1.82	0.44
1:B:869:LEU:HA	1:B:869:LEU:HD23	1.85	0.44
1:C:669:CYS:HA	1:C:719:LYS:HG2	2.00	0.44
1:C:675:LEU:HD23	1:C:720:LEU:HD11	1.99	0.44
1:B:571:ILE:O	1:B:575:ILE:HG13	2.18	0.44
1:C:663:PHE:CE1	1:C:667:ASP:HB3	2.53	0.44
2:E:104:LYS:HD3	2:E:104:LYS:HA	1.82	0.44
1:A:835:ARG:HH12	1:A:899:ASP:CG	2.25	0.44
1:B:991:HIS:O	1:B:995:THR:HG23	2.17	0.44
1:D:1050:PRO:HB2	1:D:1052:ASN:OD1	2.16	0.44
1:D:865:MET:HB2	1:D:908:PHE:CZ	2.53	0.43
1:B:814:VAL:O	1:B:818:ILE:HG13	2.18	0.43
1:C:1050:PRO:HD3	2:F:40:TYR:CE1	2.53	0.43
2:F:139:ASN:O	2:F:139:ASN:ND2	2.51	0.43
1:D:723:ASN:O	1:D:727:ILE:HG13	2.17	0.43
1:B:952:LEU:O	1:B:956:ILE:HG13	2.19	0.43
1:B:1008:LYS:O	1:B:1019:LYS:NZ	2.30	0.43
1:C:820:GLN:O	1:C:824:ARG:HB2	2.19	0.43
1:D:567:MET:O	1:D:571:ILE:HG13	2.18	0.43
2:F:141:CYS:SG	2:F:142:ALA:N	2.91	0.43
1:A:991:HIS:O	1:A:995:THR:HG23	2.18	0.43
1:C:698:MET:HE2	1:C:709:LEU:HD21	2.00	0.43
2:G:35:THR:HG23	2:G:38:ASP:OD2	2.18	0.43
2:H:96:LEU:O	2:H:100:ILE:HG12	2.19	0.43
1:B:988:ASP:O	1:B:992:ILE:HG12	2.18	0.43
1:C:803:VAL:HG11	1:C:818:ILE:HG12	2.00	0.43
1:C:970:TYR:CE1	1:C:1085:LEU:HD11	2.54	0.43
1:C:1045:LEU:HD13	1:C:1054:LEU:HD12	2.01	0.43
2:F:117:LYS:HG2	4:F:201:GDP:C6	2.53	0.43
1:B:1025:GLY:O	1:B:1029:VAL:HG23	2.19	0.43
1:C:1027:LEU:HD23	1:C:1027:LEU:HA	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:6:LEU:HD22	2:G:161:LEU:HD13	2.01	0.43
1:B:734:ARG:NH2	1:D:1126:GLN:HA	2.33	0.43
1:A:774:LEU:C	1:A:776:GLN:H	2.27	0.42
1:B:1009:ASN:HD21	2:G:65:THR:HB	1.84	0.42
1:D:691:GLU:OE2	1:D:970:TYR:OH	2.28	0.42
2:H:40:TYR:N	2:H:55:ILE:O	2.48	0.42
1:A:869:LEU:HD22	1:A:874:VAL:HG21	2.01	0.42
1:C:841:TYR:HE1	1:C:878:ALA:HB1	1.84	0.42
1:D:656:ALA:O	1:D:660:GLN:NE2	2.30	0.42
1:D:697:LEU:O	1:D:701:SER:HB3	2.19	0.42
1:D:757:VAL:HG22	1:D:764:ALA:HB2	2.00	0.42
2:F:9:LEU:HD21	2:F:71:TYR:HE1	1.83	0.42
1:B:1008:LYS:HE3	1:B:1008:LYS:HB3	1.91	0.42
1:B:1050:PRO:HB2	1:B:1052:ASN:OD1	2.20	0.42
1:C:694:LEU:HD11	1:C:710:LEU:HD21	2.01	0.42
2:F:9:LEU:HD21	2:F:71:TYR:CE1	2.54	0.42
2:F:138:TRP:HB2	2:F:141:CYS:HB2	2.01	0.42
2:H:4:TYR:CE2	2:H:162:VAL:HG13	2.55	0.42
1:A:598:PHE:HZ	1:A:738:GLY:HA3	1.84	0.42
1:B:703:ASN:O	1:B:712:ARG:NH1	2.53	0.42
1:C:640:PHE:O	1:C:644:ILE:HG12	2.20	0.42
2:E:80:LEU:HB3	2:E:93:LEU:HD22	2.00	0.42
2:F:117:LYS:HG2	4:F:201:GDP:C5	2.54	0.42
2:G:117:LYS:HE2	4:G:201:GDP:C4	2.54	0.42
1:A:959:LEU:HD22	1:A:964:ALA:HA	2.02	0.42
1:B:613:ALA:HB1	1:B:616:LEU:HB2	2.02	0.42
1:B:711:ARG:HH21	2:G:34:PRO:HA	1.85	0.42
1:C:599:LEU:HD23	1:C:599:LEU:HA	1.91	0.42
1:C:643:LEU:HD11	1:C:1137:MET:HA	2.01	0.42
1:C:758:ASP:OD1	1:C:909:HIS:HB3	2.20	0.42
1:D:588:ASP:OD1	1:D:588:ASP:N	2.53	0.42
1:A:695:GLN:HG3	1:A:978:GLN:HE22	1.84	0.42
1:A:787:VAL:HB	1:A:801:LEU:HB3	2.01	0.42
1:B:1102:GLU:N	1:B:1102:GLU:OE1	2.52	0.42
1:A:1141:LYS:HB3	1:A:1141:LYS:HE2	1.79	0.42
1:B:932:GLU:HA	1:B:935:LEU:HG	2.01	0.42
1:D:766:TYR:HB3	1:D:941:THR:OG1	2.20	0.42
2:G:161:LEU:O	2:G:165:ILE:HG13	2.20	0.42
1:B:640:PHE:CE2	1:B:644:ILE:HD11	2.55	0.42
1:C:854:ASP:OD1	1:C:854:ASP:N	2.52	0.42
1:C:869:LEU:HD22	1:C:874:VAL:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:84:ILE:HD11	2:H:118:CYS:HA	2.02	0.42
1:A:966:LEU:HD11	1:A:1080:ILE:HG21	2.01	0.42
1:B:572:ARG:HG2	1:C:930:ILE:HG23	2.00	0.42
1:D:741:LEU:O	1:D:745:VAL:HG23	2.19	0.42
1:A:769:SER:HB3	1:A:772:TRP:HB2	2.02	0.42
1:B:694:LEU:O	1:B:698:MET:HG2	2.20	0.41
1:D:640:PHE:HD2	1:D:1106:LEU:HD22	1.85	0.41
1:D:755:GLY:HA3	1:D:766:TYR:CD1	2.54	0.41
2:H:90:PHE:HE2	2:H:130:GLN:HE21	1.66	0.41
1:A:966:LEU:CD1	1:A:1080:ILE:HD13	2.48	0.41
1:C:850:LEU:HD12	1:C:850:LEU:O	2.20	0.41
1:D:810:THR:O	1:D:814:VAL:HG23	2.20	0.41
1:D:933:MET:HE3	1:D:933:MET:HB2	1.88	0.41
1:B:788:VAL:HA	1:B:799:GLU:O	2.21	0.41
1:C:972:PHE:HD1	1:C:975:LEU:HD12	1.86	0.41
1:A:719:LYS:HA	1:A:719:LYS:HD2	1.84	0.41
1:B:614:GLY:O	1:B:617:THR:OG1	2.34	0.41
1:D:1007:LEU:O	1:D:1027:LEU:HD13	2.21	0.41
1:A:1017:MET:HE2	1:A:1017:MET:HB2	1.85	0.41
1:B:776:GLN:HG2	1:B:778:GLN:HE21	1.86	0.41
1:C:757:VAL:HG22	1:C:764:ALA:HB2	2.02	0.41
1:C:1053:LYS:NZ	2:F:38:ASP:OD1	2.53	0.41
2:G:117:LYS:HG2	4:G:201:GDP:C6	2.56	0.41
2:H:35:THR:HG21	2:H:57:ASP:OD2	2.21	0.41
1:B:645:ARG:HA	1:B:685:TYR:CZ	2.55	0.41
1:C:588:ASP:OD1	1:C:588:ASP:N	2.53	0.41
1:D:1024:ASP:OD2	2:H:67:MET:HB3	2.20	0.41
2:H:13:GLY:HA2	4:H:201:GDP:O3A	2.21	0.41
1:A:742:PHE:HB2	1:A:1014:PHE:CZ	2.55	0.41
1:A:569:SER:HA	1:A:572:ARG:HD2	2.03	0.41
1:B:665:ILE:HD12	2:G:61:THR:HG21	2.03	0.41
1:B:766:TYR:HB3	1:B:941:THR:OG1	2.21	0.41
1:C:970:TYR:HB2	1:C:1080:ILE:HG13	2.03	0.41
1:D:814:VAL:O	1:D:818:ILE:HG13	2.21	0.41
1:A:737:VAL:HG21	1:A:1003:TRP:CH2	2.53	0.41
1:B:1021:PRO:O	1:B:1024:ASP:HB3	2.21	0.41
1:B:1035:MET:HE2	2:G:36:ILE:HG13	2.02	0.41
1:C:810:THR:HG21	1:C:866:LEU:HD12	2.02	0.41
1:D:969:LYS:HD3	1:D:1077:GLN:HB3	2.03	0.41
1:D:1008:LYS:HE3	1:D:1008:LYS:HB3	1.89	0.41
1:B:610:PHE:HB2	1:B:617:THR:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:982:LYS:HE3	1:B:982:LYS:HB3	1.85	0.41
1:C:723:ASN:O	1:C:727:ILE:HG13	2.21	0.41
2:E:53:LEU:HD11	2:E:162:VAL:HG21	2.03	0.41
1:B:599:LEU:HD23	1:B:599:LEU:HA	1.93	0.40
1:B:1003:TRP:O	1:B:1007:LEU:HB2	2.21	0.40
1:B:1055:LEU:HD23	1:B:1055:LEU:HA	1.86	0.40
1:C:677:LEU:HD12	1:C:677:LEU:HA	1.95	0.40
1:D:737:VAL:HG11	1:D:1003:TRP:CZ3	2.55	0.40
2:E:66:ALA:HB3	2:E:68:ARG:CZ	2.51	0.40
1:B:840:GLU:HB2	1:B:881:LYS:HB3	2.04	0.40
1:D:1027:LEU:HD23	1:D:1027:LEU:HA	1.89	0.40
2:G:119:ASP:OD1	2:G:119:ASP:N	2.52	0.40
1:B:616:LEU:HD23	1:B:616:LEU:HA	1.96	0.40
1:C:643:LEU:HB3	1:C:649:PHE:CE2	2.57	0.40
1:C:982:LYS:HE2	1:C:982:LYS:HB3	1.81	0.40
1:D:839:ILE:HG23	1:D:880:ILE:HG23	2.03	0.40
1:B:718:GLU:HB3	1:B:1001:ARG:CZ	2.51	0.40
1:B:755:GLY:HA3	1:B:766:TYR:CE1	2.56	0.40
1:B:939:LEU:O	1:B:943:VAL:HG23	2.22	0.40
1:C:767:THR:HG22	1:C:773:LEU:HD13	2.04	0.40
1:D:643:LEU:HD13	1:D:649:PHE:CZ	2.56	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	552/599 (92%)	510 (92%)	40 (7%)	2 (0%)	30	60
1	B	561/599 (94%)	527 (94%)	31 (6%)	3 (0%)	24	55
1	C	550/599 (92%)	514 (94%)	34 (6%)	2 (0%)	30	60
1	D	550/599 (92%)	505 (92%)	43 (8%)	2 (0%)	30	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	166/199 (83%)	152 (92%)	13 (8%)	1 (1%)	21	52
2	F	164/199 (82%)	153 (93%)	10 (6%)	1 (1%)	21	52
2	G	166/199 (83%)	157 (95%)	9 (5%)	0	100	100
2	H	165/199 (83%)	151 (92%)	14 (8%)	0	100	100
All	All	2874/3192 (90%)	2669 (93%)	194 (7%)	11 (0%)	30	60

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	889	ALA
1	D	889	ALA
1	A	889	ALA
1	A	884	THR
1	B	585	GLU
1	D	663	PHE
1	C	889	ALA
2	E	30	GLU
1	B	884	THR
1	C	998	LEU
2	F	110	PRO

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	439/552 (80%)	427 (97%)	12 (3%)	39	63
1	B	418/552 (76%)	404 (97%)	14 (3%)	33	59
1	C	420/552 (76%)	404 (96%)	16 (4%)	29	57
1	D	421/552 (76%)	408 (97%)	13 (3%)	35	60
2	E	133/163 (82%)	130 (98%)	3 (2%)	44	66
2	F	131/163 (80%)	129 (98%)	2 (2%)	57	72
2	G	128/163 (78%)	126 (98%)	2 (2%)	55	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	118/163 (72%)	117 (99%)	1 (1%)	73	79
All	All	2208/2860 (77%)	2145 (97%)	63 (3%)	37	62

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	677	LEU
1	A	693	LEU
1	A	694	LEU
1	A	717	VAL
1	A	726	SER
1	A	737	VAL
1	A	788	VAL
1	A	850	LEU
1	A	884	THR
1	A	1028	SER
1	A	1041	THR
1	A	1137	MET
1	B	577	GLN
1	B	655	HIS
1	B	709	LEU
1	B	717	VAL
1	B	720	LEU
1	B	737	VAL
1	B	748	LEU
1	B	799	GLU
1	B	850	LEU
1	B	851	GLN
1	B	884	THR
1	B	942	LYS
1	B	966	LEU
1	B	1054	LEU
1	C	577	GLN
1	C	615	THR
1	C	619	VAL
1	C	720	LEU
1	C	748	LEU
1	C	767	THR
1	C	773	LEU
1	C	812	GLN
1	C	831	THR
1	C	837	ILE

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Mol	Chain	Res	Type
1	C	850	LEU
1	C	864	THR
1	C	884	THR
1	C	897	VAL
1	C	905	ILE
1	C	1005	ASN
1	D	619	VAL
1	D	665	ILE
1	D	717	VAL
1	D	813	GLN
1	D	823	GLN
1	D	848	VAL
1	D	850	LEU
1	D	863	VAL
1	D	966	LEU
1	D	973	ASP
1	D	1007	LEU
1	D	1029	VAL
1	D	1085	LEU
2	E	67	MET
2	E	97	ARG
2	E	106	THR
2	F	67	MET
2	F	136	ARG
2	G	52	MET
2	G	67	MET
2	H	53	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	563	GLN
1	A	705	ASN
1	A	813	GLN
1	A	823	GLN
1	A	946	HIS
1	A	1009	ASN
1	A	1083	GLN
1	A	1136	ASN
1	B	703	ASN
1	B	978	GLN
1	B	1009	ASN

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Mol	Chain	Res	Type
1	B	1138	HIS
1	C	577	GLN
1	C	695	GLN
1	C	867	ASN
1	C	978	GLN
1	C	996	ASN
1	C	1101	ASN
1	D	705	ASN
1	D	867	ASN
1	D	946	HIS
1	D	978	GLN
1	D	1048	HIS
2	E	22	GLN
2	F	22	GLN
2	F	76	GLN
2	F	91	ASN
2	F	94	GLN
2	F	139	ASN
2	G	49	GLN
2	G	91	ASN
2	G	116	ASN
2	H	22	GLN
2	H	153	ASN
2	H	164	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 ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
5	AF3	G	202	-	0,3,3	-	-	-		
5	AF3	E	202	-	0,3,3	-	-	-		
4	GDP	H	201	3	29,30,30	1.20	4 (13%)	45,47,47	1.81	7 (15%)
5	AF3	H	202	-	0,3,3	-	-	-		
4	GDP	E	201	3	29,30,30	1.17	3 (10%)	45,47,47	1.78	7 (15%)
5	AF3	F	202	-	0,3,3	-	-	-		
4	GDP	F	201	3	29,30,30	1.13	2 (6%)	45,47,47	1.84	7 (15%)
4	GDP	G	201	3	29,30,30	1.14	2 (6%)	45,47,47	1.75	7 (15%)

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	GDP	E	201	3	-	3/16/32/32	0/3/3/3
4	GDP	H	201	3	-	0/16/32/32	0/3/3/3
4	GDP	F	201	3	-	5/16/32/32	0/3/3/3
4	GDP	G	201	3	-	8/16/32/32	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	201	GDP	C5-C4	3.13	1.47	1.38
4	G	201	GDP	C5-C4	3.12	1.47	1.38
4	H	201	GDP	C5-C4	3.11	1.47	1.38
4	E	201	GDP	C5-C4	3.10	1.47	1.38
4	E	201	GDP	C6-N1	-2.45	1.34	1.38
4	G	201	GDP	C6-N1	-2.42	1.34	1.38
4	F	201	GDP	C6-N1	-2.39	1.34	1.38
4	H	201	GDP	C5-N7	-2.27	1.34	1.39
4	H	201	GDP	C6-N1	-2.26	1.34	1.38
4	E	201	GDP	C5-N7	-2.12	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	201	GDP	PA-O3A	2.11	1.61	1.59

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	201	GDP	C5-C4-N3	-6.25	118.44	128.39
4	F	201	GDP	C5-C4-N3	-6.19	118.54	128.39
4	E	201	GDP	C5-C4-N3	-6.09	118.70	128.39
4	G	201	GDP	C5-C4-N3	-5.95	118.93	128.39
4	F	201	GDP	C2-N3-C4	5.12	121.12	112.30
4	H	201	GDP	C2-N3-C4	5.09	121.07	112.30
4	E	201	GDP	C2-N3-C4	4.87	120.69	112.30
4	G	201	GDP	C2-N3-C4	4.87	120.69	112.30
4	E	201	GDP	N9-C4-N3	4.77	135.48	125.95
4	H	201	GDP	N9-C4-N3	4.66	135.26	125.95
4	F	201	GDP	N9-C4-N3	4.65	135.24	125.95
4	G	201	GDP	N9-C4-N3	4.54	135.03	125.95
4	F	201	GDP	C6-C5-N7	3.22	136.14	130.29
4	G	201	GDP	C6-C5-N7	3.18	136.09	130.29
4	E	201	GDP	C6-C5-N7	2.99	135.73	130.29
4	H	201	GDP	C6-C5-N7	2.98	135.71	130.29
4	F	201	GDP	C4-C5-N7	-2.52	106.67	110.67
4	H	201	GDP	C4-C5-N7	-2.47	106.75	110.67
4	G	201	GDP	C4-C5-N7	-2.47	106.75	110.67
4	H	201	GDP	C3'-C2'-C1'	2.41	106.02	101.46
4	E	201	GDP	C4-C5-N7	-2.37	106.92	110.67
4	H	201	GDP	O6-C6-C5	-2.32	120.42	126.53
4	E	201	GDP	O6-C6-C5	-2.30	120.45	126.53
4	F	201	GDP	O6-C6-C5	-2.19	120.75	126.53
4	E	201	GDP	C3'-C2'-C1'	2.17	105.56	101.46
4	G	201	GDP	O6-C6-C5	-2.13	120.91	126.53
4	F	201	GDP	C3'-C2'-C1'	2.09	105.42	101.46
4	G	201	GDP	C3'-C2'-C1'	2.05	105.34	101.46

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	201	GDP	PA-O3A-PB-O2B
4	F	201	GDP	PA-O3A-PB-O2B
4	G	201	GDP	PA-O3A-PB-O2B
4	G	201	GDP	C5'-O5'-PA-O3A

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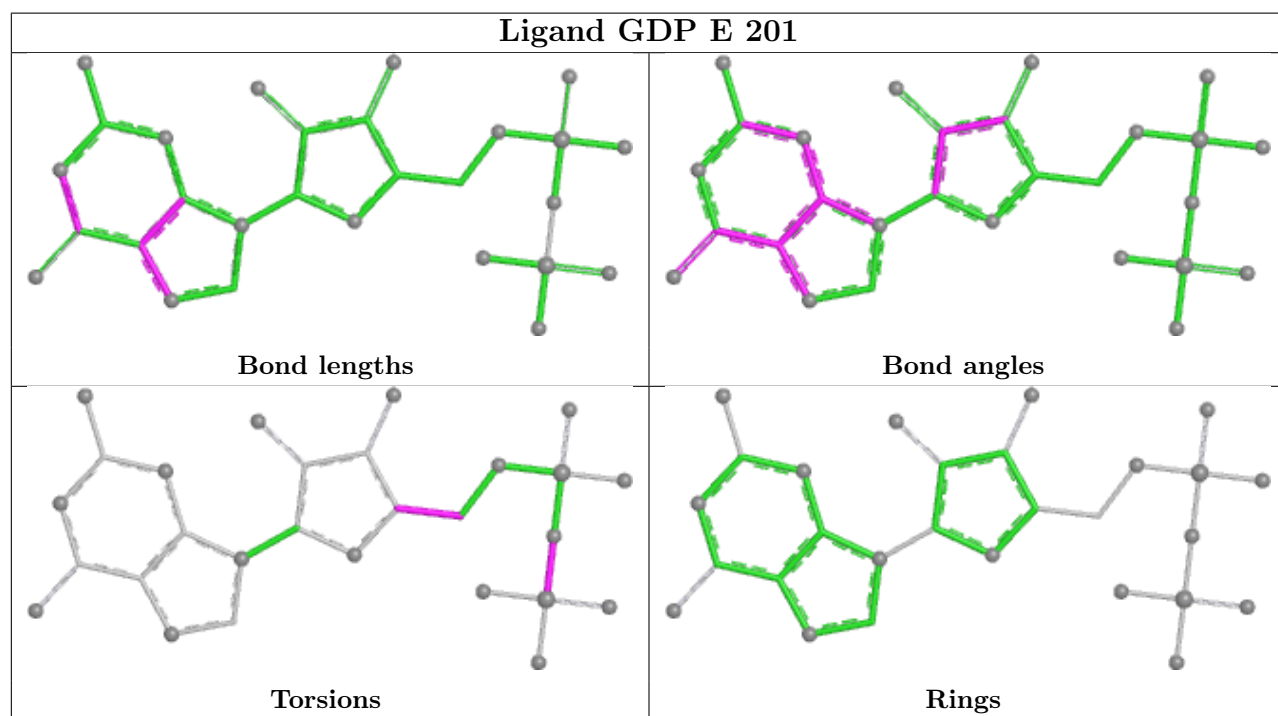
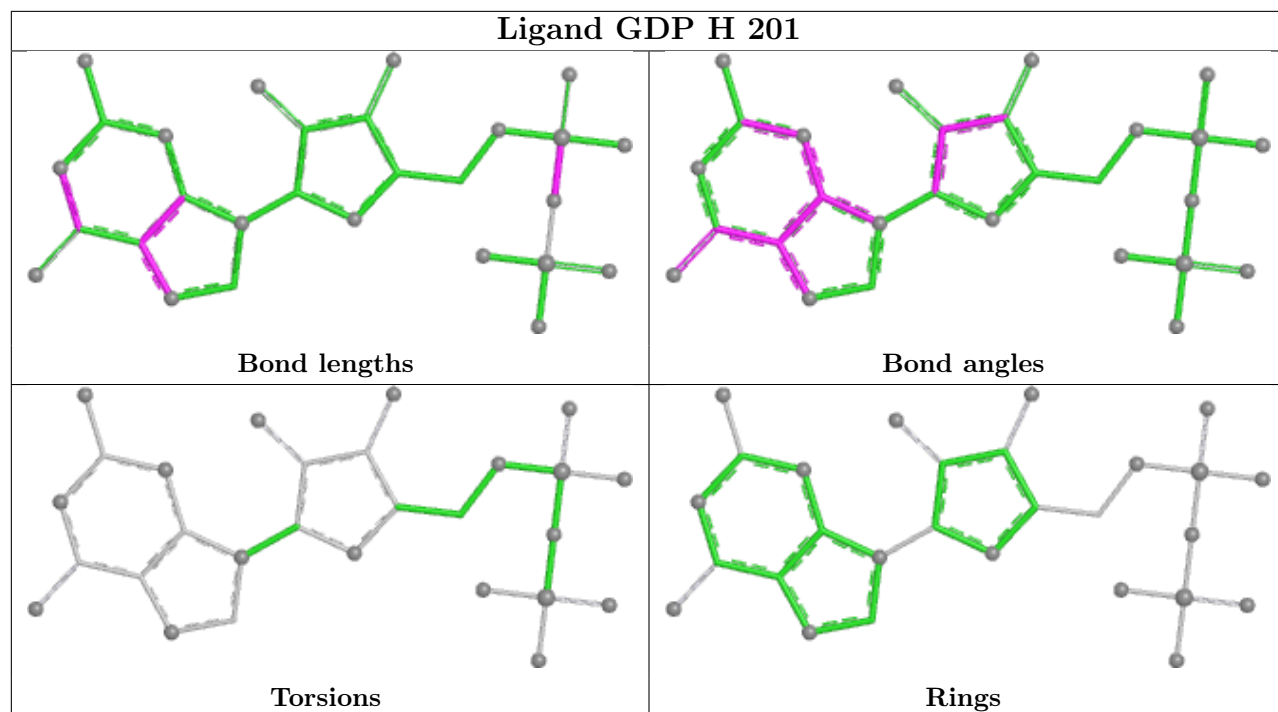
Mol	Chain	Res	Type	Atoms
4	G	201	GDP	C5'-O5'-PA-O1A
4	G	201	GDP	C5'-O5'-PA-O2A
4	G	201	GDP	O4'-C4'-C5'-O5'
4	G	201	GDP	C3'-C4'-C5'-O5'
4	F	201	GDP	C5'-O5'-PA-O3A
4	F	201	GDP	C5'-O5'-PA-O1A
4	E	201	GDP	PA-O3A-PB-O1B
4	F	201	GDP	PA-O3A-PB-O1B
4	F	201	GDP	PA-O3A-PB-O3B
4	G	201	GDP	PA-O3A-PB-O3B
4	G	201	GDP	PA-O3A-PB-O1B
4	E	201	GDP	O4'-C4'-C5'-O5'

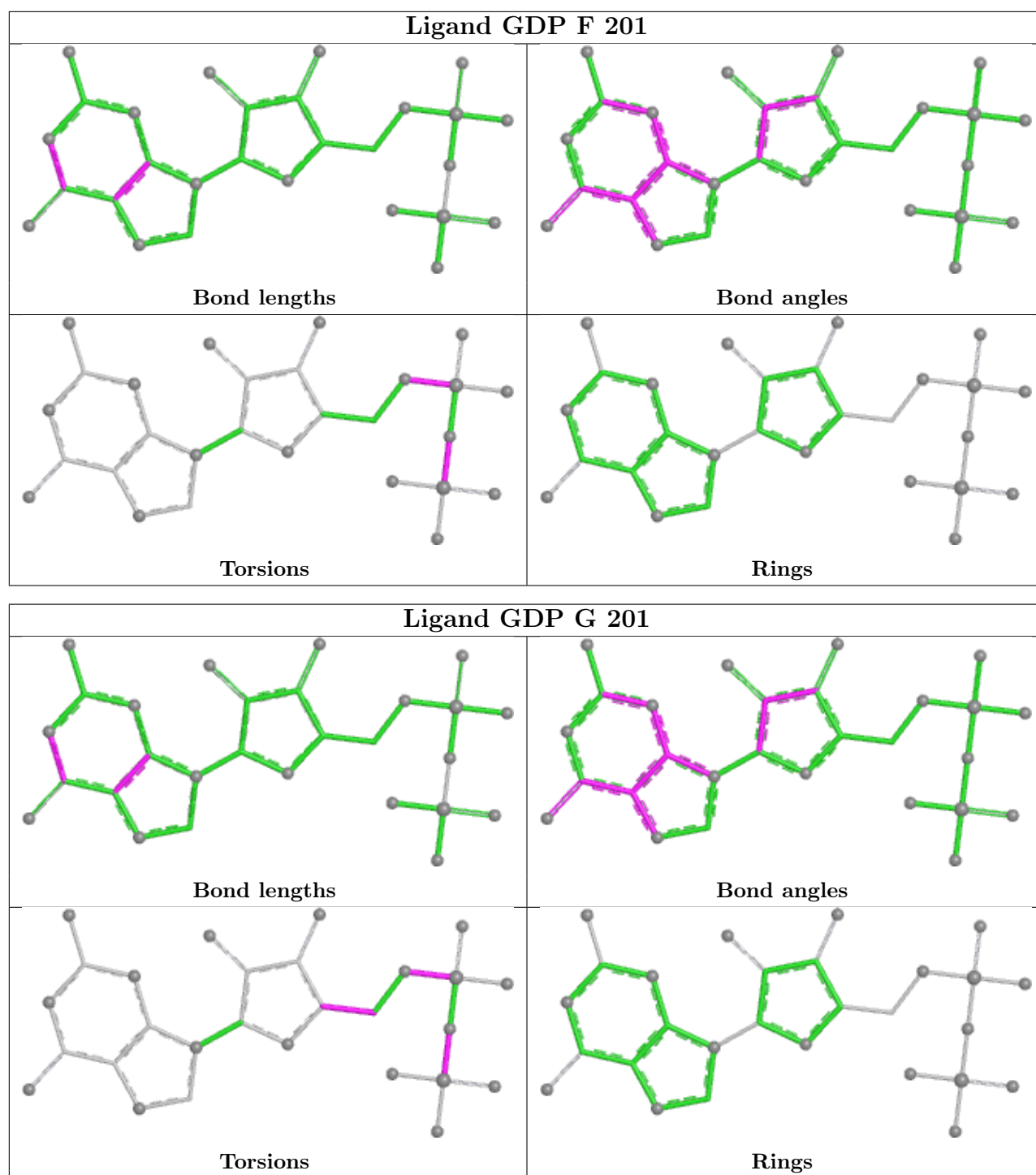
There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	202	AF3	1	0
4	H	201	GDP	5	0
5	F	202	AF3	1	0
4	F	201	GDP	3	0
4	G	201	GDP	2	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	574/599 (95%)	-0.21	5 (0%) 81 65	66, 132, 194, 219	0
1	B	571/599 (95%)	-0.23	1 (0%) 91 86	62, 135, 220, 277	0
1	C	560/599 (93%)	-0.29	2 (0%) 88 79	70, 134, 189, 214	0
1	D	560/599 (93%)	-0.19	7 (1%) 75 56	83, 153, 223, 253	0
2	E	168/199 (84%)	-0.29	1 (0%) 85 73	91, 139, 179, 208	0
2	F	166/199 (83%)	-0.22	0 100 100	94, 147, 186, 205	0
2	G	168/199 (84%)	-0.13	2 (1%) 76 58	95, 166, 214, 234	0
2	H	167/199 (83%)	-0.07	2 (1%) 76 58	110, 168, 222, 243	0
All	All	2934/3192 (91%)	-0.22	20 (0%) 84 70	62, 143, 207, 277	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1035	MET	3.8
1	A	700	GLN	3.4
2	H	67	MET	3.3
2	E	6	LEU	3.0
1	C	1121	PHE	2.9
1	A	1055	LEU	2.6
1	D	616	LEU	2.6
1	D	1110	TYR	2.6
2	H	11	SER	2.6
1	D	791	VAL	2.4
1	B	865	MET	2.3
1	C	564	LEU	2.2
1	D	1121	PHE	2.2
1	D	713	THR	2.2
1	A	968	VAL	2.1
1	A	783	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	G	95	ASP	2.1
2	G	37	GLU	2.1
1	A	974	PHE	2.0
1	D	818	ILE	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](#)

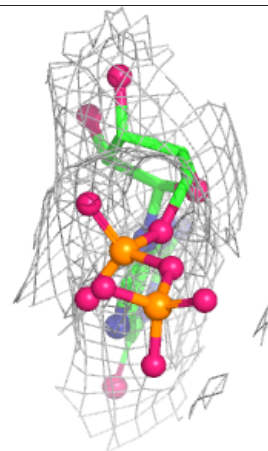
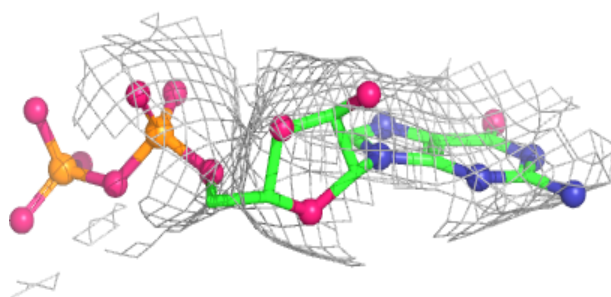
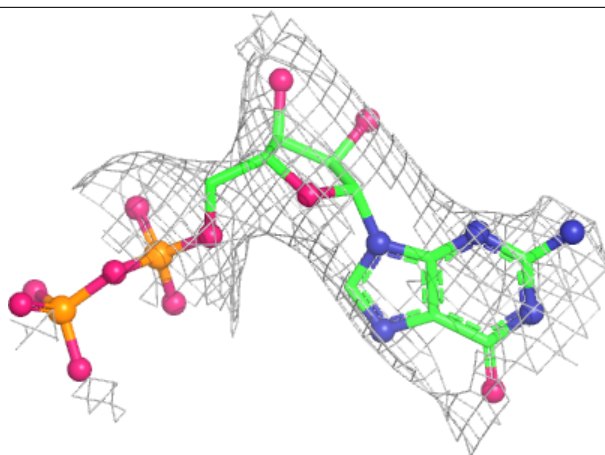
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	AF3	E	202	4/4	0.92	0.12	89,128,128,129	0
5	AF3	H	202	4/4	0.93	0.10	115,121,159,159	0
5	AF3	G	202	4/4	0.94	0.10	87,91,97,125	0
4	GDP	H	201	28/28	0.94	0.08	123,139,162,166	0
4	GDP	E	201	28/28	0.95	0.06	91,127,141,183	0
4	GDP	F	201	28/28	0.96	0.06	97,120,134,138	0
4	GDP	G	201	28/28	0.96	0.05	107,139,151,162	0
3	MG	E	200	1/1	0.96	0.05	88,88,88,88	0
3	MG	G	200	1/1	0.97	0.04	108,108,108,108	0
5	AF3	F	202	4/4	0.97	0.09	91,103,128,169	0
3	MG	F	200	1/1	0.98	0.06	88,88,88,88	0
3	MG	H	200	1/1	0.98	0.07	113,113,113,113	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.

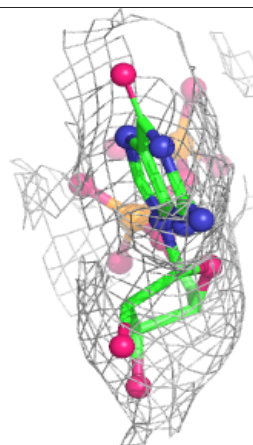
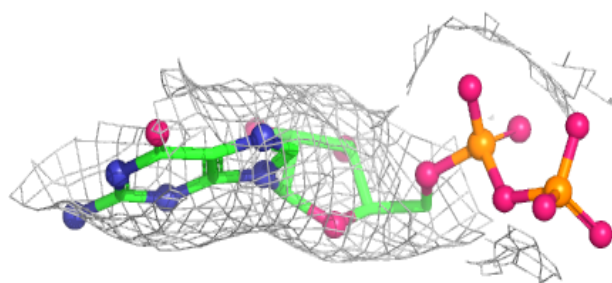
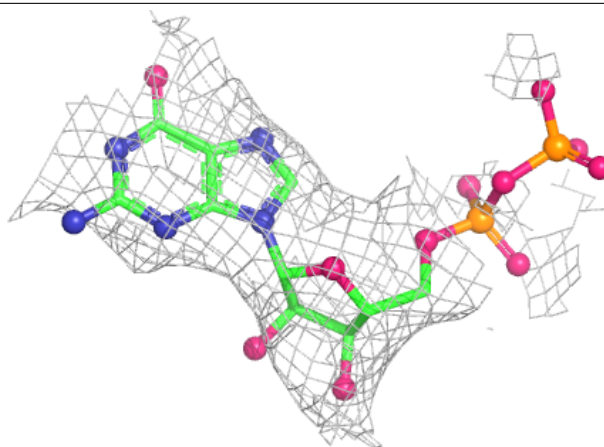
Electron density around GDP H 201:

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

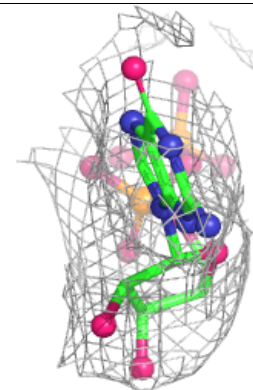
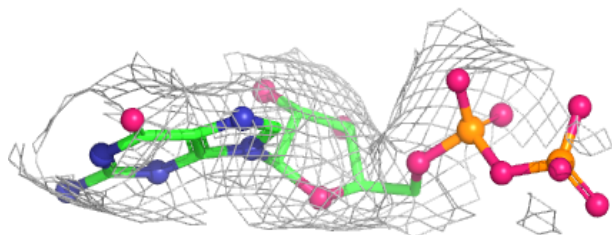
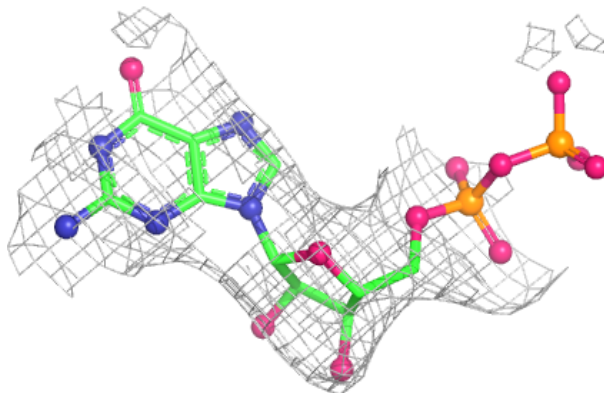


Electron density around GDP E 201:

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

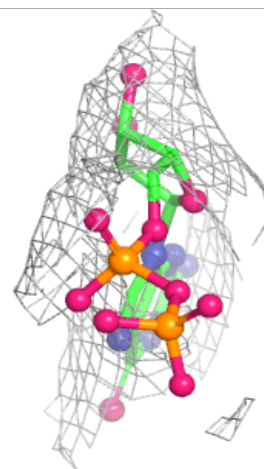
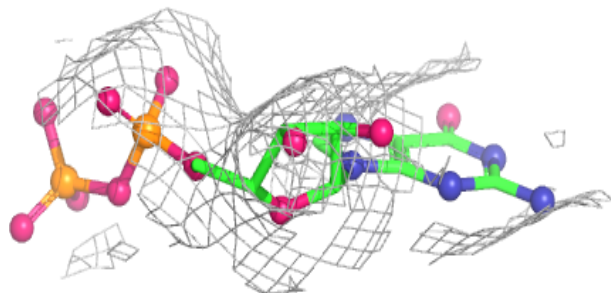
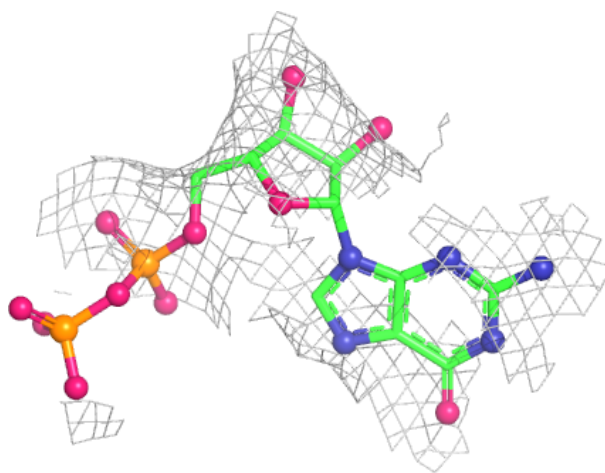
**Electron density around GDP F 201:**

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



Electron density around GDP G 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 [i](#)

There are no such residues in this entry.