



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:06 PM UTC

PDB ID : 5DIS / pdb_00005dis
Title : Crystal structure of a CRM1-RanGTP-SPN1 export complex bound to a 113 amino acid FG-repeat containing fragment of Nup214
Authors : Monecke, T.; Port, S.A.; Dickmanns, A.; Kehlenbach, R.H.; Ficner, R.
Deposited on : 2015-09-01
Resolution : 2.85 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

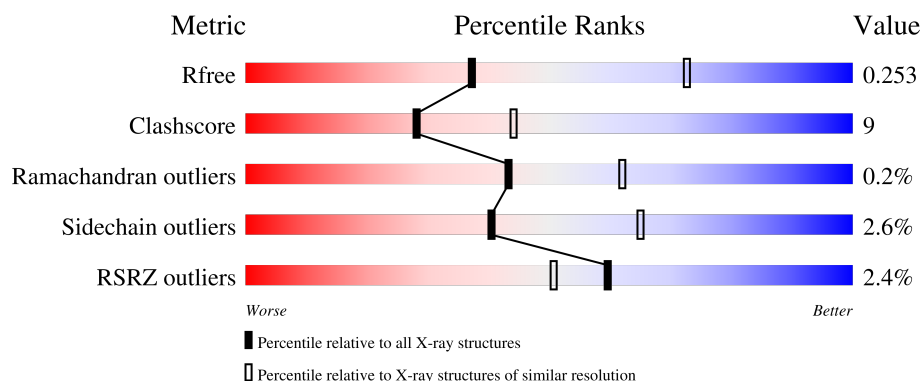
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	1044	77% 22% ..
2	B	172	84% 16% .
3	C	289	3% 67% 23% 9%
4	D	479	6% 62% 17% 18%
5	E	2	50% 50%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14991 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 Exportin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1032	Total	C	N	O	S	0	0	0
			8366	5370	1400	1541	55			

- Molecule 2 is a protein called GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1398	909	245	239	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	69	LEU	GLN	engineered mutation	UNP P62826

- Molecule 3 is a protein called Snurportin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	264	Total	C	N	O	S	0	0	0
			2131	1359	365	393	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	linker	UNP O95149
C	0	SER	-	linker	UNP O95149

- Molecule 4 is a protein called Maltose-binding periplasmic protein,Nuclear pore complex protein Nup214.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	395	Total	C	N	O	S	0	0	0
			3012	1941	489	579	3			

There are 11 discrepancies between the modelled and reference sequences:

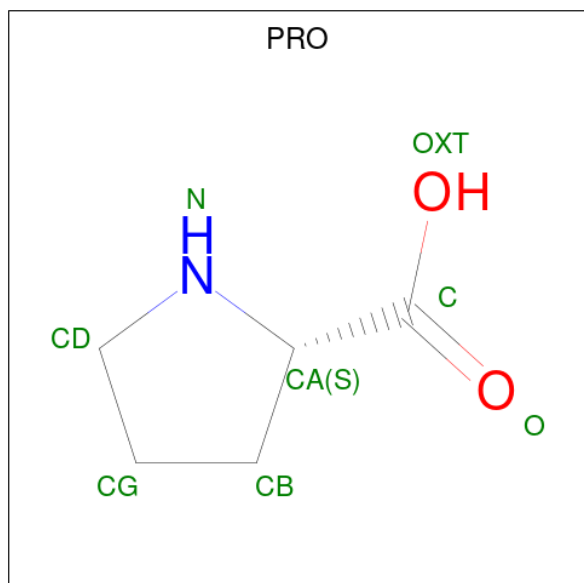
Chain	Residue	Modelled	Actual	Comment	Reference
D	362	ALA	-	linker	UNP P0AEX9
D	363	ALA	-	linker	UNP P0AEX9
D	364	ALA	-	linker	UNP P0AEX9
D	365	GLN	-	linker	UNP P0AEX9
D	366	THR	-	linker	UNP P0AEX9
D	367	ASN	-	linker	UNP P0AEX9
D	368	ALA	-	linker	UNP P0AEX9
D	1912	ALA	-	linker	UNP P0AEX9
D	1913	ALA	-	linker	UNP P0AEX9
D	1914	GLU	-	linker	UNP P0AEX9
D	1915	PHE	-	linker	UNP P0AEX9

- Molecule 5 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



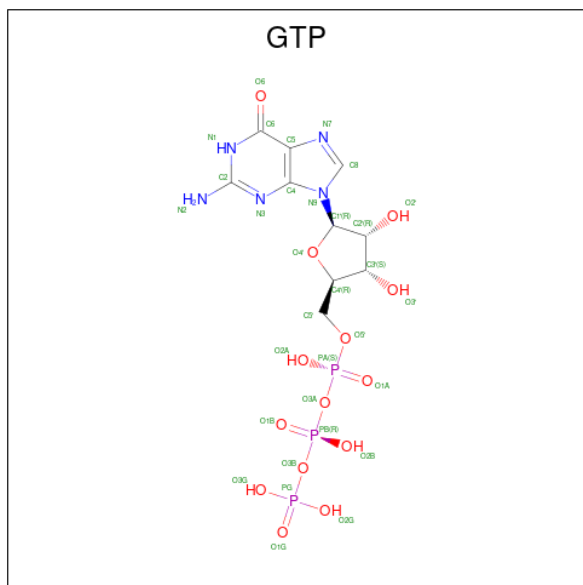
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
5	E	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 6 is PROLINE (CCD ID: PRO) (formula: C₅H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			8	5	1	2		
6	C	1	Total	C	N	O	0	0
			8	5	1	2		

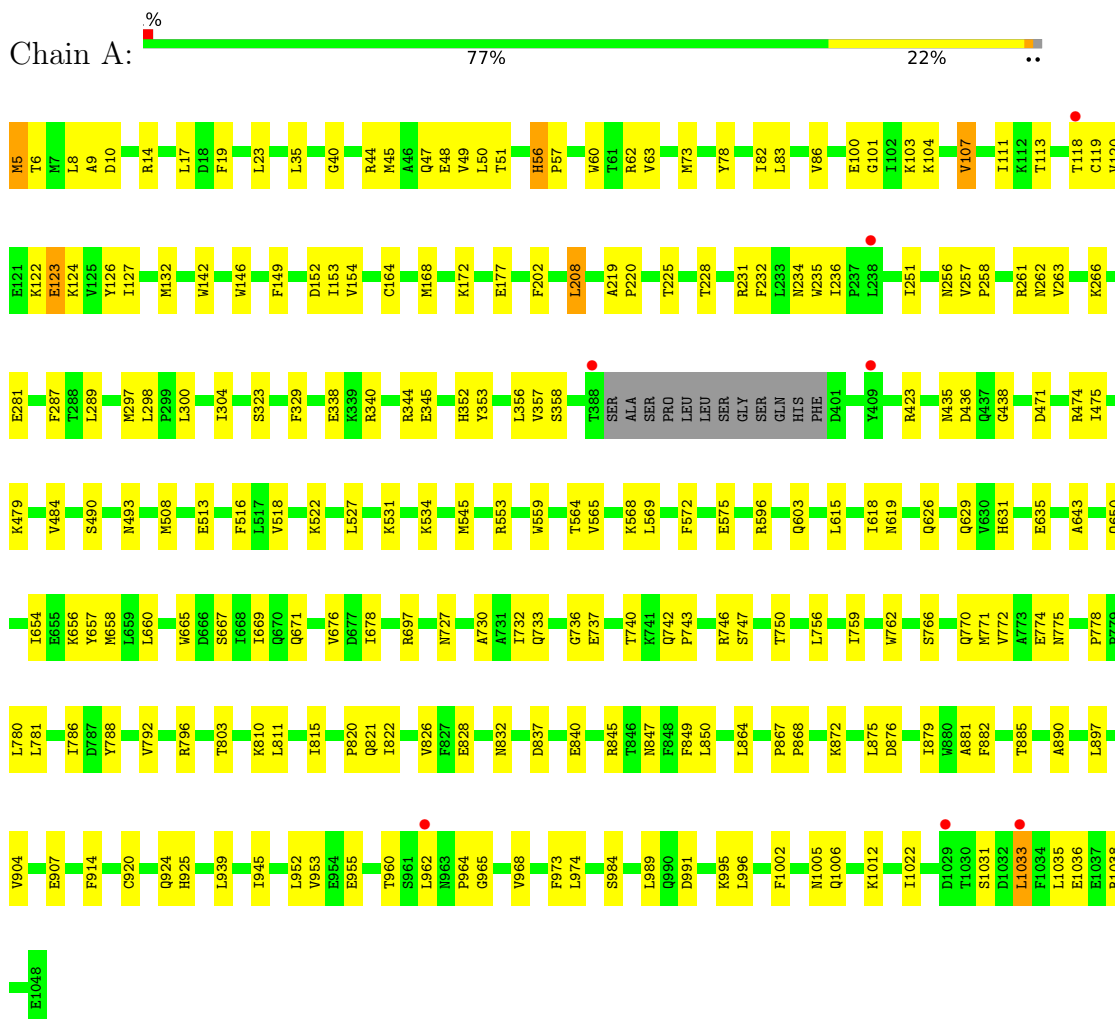
- Molecule 7 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



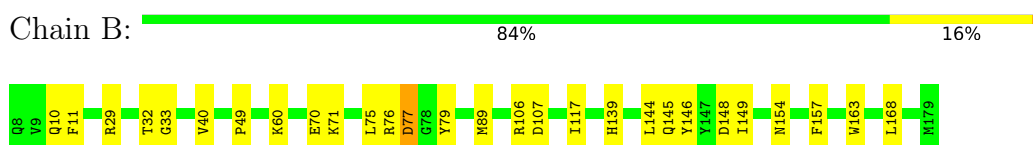
3 Residue-property plots [i](#)

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.

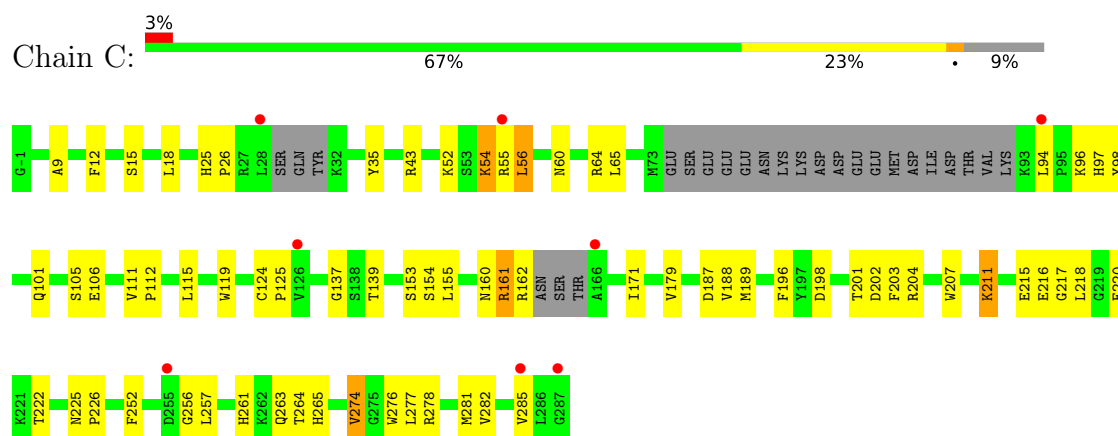
• Molecule 1: Exportin-1



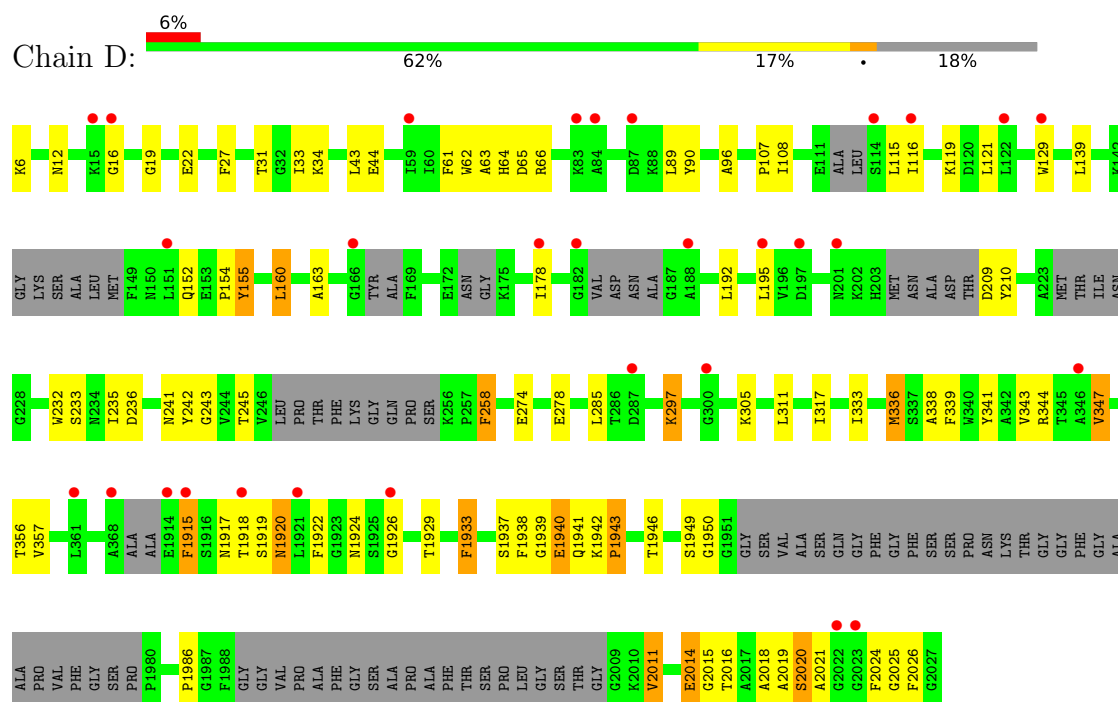
• Molecule 2: GTP-binding nuclear protein Ran



- Molecule 3: Snurportin-1



- Molecule 4: Maltose-binding periplasmic protein, Nuclear pore complex protein Nup214



- Molecule 5: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	112.33Å 248.97Å 210.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.48 – 2.85 48.48 – 2.85	Depositor EDS
% Data completeness (in resolution range)	98.2 (48.48-2.85) 98.2 (48.48-2.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.206 , 0.249 0.213 , 0.253	Depositor DCC
R_{free} test set	3396 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	80.0	Xtriage
Anisotropy	0.811	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 101.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14991	wwPDB-VP
Average B, all atoms (Å ²)	125.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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, GTP, GLC

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	0.41	0/8537	0.82	7/11566 (0.1%)
2	B	0.44	0/1432	0.76	0/1932
3	C	0.37	0/2186	0.84	3/2957 (0.1%)
4	D	0.41	0/3080	0.84	8/4158 (0.2%)
All	All	0.41	0/15235	0.82	18/20613 (0.1%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	336	MET	N-CA-C	8.51	120.63	111.36
3	C	124	CYS	CA-C-N	7.01	126.97	120.03
3	C	124	CYS	C-N-CA	7.01	126.97	120.03
4	D	1940	GLU	N-CA-C	6.14	118.69	108.13
4	D	347	VAL	N-CA-C	5.79	118.33	111.09
1	A	235	TRP	N-CA-C	5.77	122.14	114.12
1	A	778	PRO	N-CA-C	5.67	117.62	110.70
4	D	258	PHE	CA-C-N	-5.49	115.62	122.48
4	D	258	PHE	C-N-CA	-5.49	115.62	122.48
3	C	161	ARG	N-CA-C	-5.37	105.32	111.07
4	D	297	LYS	CA-C-N	5.18	125.19	119.90
4	D	297	LYS	C-N-CA	5.18	125.19	119.90
1	A	618	ILE	N-CA-C	5.17	115.61	110.23
4	D	27	PHE	N-CA-C	-5.16	105.56	111.07
1	A	56	HIS	CA-C-N	5.15	124.87	119.82
1	A	56	HIS	C-N-CA	5.15	124.87	119.82
1	A	742	GLN	CA-C-N	5.01	126.10	119.84
1	A	742	GLN	C-N-CA	5.01	126.10	119.84

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	8366	0	8417	148	0
2	B	1398	0	1425	17	0
3	C	2131	0	2083	46	0
4	D	3012	0	2923	72	0
5	E	23	0	21	1	0
6	A	8	0	7	0	0
6	C	8	0	7	1	0
7	B	32	0	12	0	0
8	B	1	0	0	0	0
9	A	3	0	0	1	0
9	B	9	0	0	0	0
All	All	14991	0	14895	267	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 (267) 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)
4:D:338:ALA:HA	4:D:341:TYR:CE2	2.12	0.84
4:D:2018:ALA:O	4:D:2021:ALA:HB3	1.83	0.78
4:D:155:TYR:HE2	4:D:258:PHE:CE1	2.02	0.77
1:A:251:ILE:HG21	1:A:289:LEU:HB3	1.67	0.77
1:A:471:ASP:OD1	1:A:474:ARG:NH2	2.23	0.72
1:A:19:PHE:HA	1:A:62:ARG:NH2	2.05	0.71
1:A:953:VAL:HG21	1:A:974:LEU:HD11	1.73	0.70
4:D:64:HIS:CE1	4:D:96:ALA:HB1	2.27	0.69
3:C:160:ASN:OD1	3:C:161:ARG:N	2.26	0.69
4:D:2016:THR:O	4:D:2020:SER:OG	2.12	0.68
4:D:339:PHE:O	4:D:343:VAL:HG23	1.94	0.68
1:A:23:LEU:HD22	1:A:62:ARG:HD3	1.76	0.67
4:D:1918:THR:OG1	4:D:1919:SER:N	2.25	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:137:GLY:HA2	3:C:160:ASN:O	1.94	0.67
3:C:65:LEU:HD21	3:C:101:GLN:HE21	1.61	0.66
1:A:358:SER:O	1:A:423:ARG:NH1	2.29	0.65
2:B:70:GLU:O	2:B:76:ARG:NH2	2.30	0.65
1:A:123:GLU:HB2	1:A:126:TYR:HB3	1.79	0.65
4:D:356:THR:HG22	4:D:357:VAL:H	1.61	0.65
1:A:338:GLU:O	1:A:344:ARG:NH2	2.30	0.65
1:A:23:LEU:HB2	1:A:62:ARG:NH1	2.11	0.64
1:A:820:PRO:HB3	4:D:1938:PHE:O	1.97	0.63
4:D:152:GLN:O	4:D:344:ARG:HG3	1.98	0.62
4:D:155:TYR:CE2	4:D:258:PHE:CE1	2.86	0.62
1:A:9:ALA:HA	1:A:49:VAL:HG22	1.80	0.62
1:A:56:HIS:CD2	1:A:57:PRO:HD2	2.34	0.62
2:B:49:PRO:HB3	2:B:60:LYS:HE2	1.81	0.61
4:D:343:VAL:O	4:D:347:VAL:HG23	1.99	0.61
1:A:103:LYS:O	1:A:107:VAL:HG13	2.00	0.61
3:C:54:LYS:HD2	3:C:198:ASP:OD1	2.00	0.61
4:D:258:PHE:O	4:D:258:PHE:CD1	2.54	0.61
1:A:475:ILE:O	1:A:479:LYS:HG2	1.99	0.61
1:A:840:GLU:O	1:A:845:ARG:NH1	2.33	0.60
4:D:2014:GLU:OE1	4:D:2015:GLY:N	2.35	0.60
1:A:5:MET:SD	1:A:6:THR:N	2.72	0.60
1:A:6:THR:HA	1:A:9:ALA:HB3	1.84	0.60
4:D:61:PHE:HE1	4:D:108:ILE:HG13	1.67	0.59
4:D:31:THR:HG22	4:D:33:ILE:HG13	1.84	0.59
1:A:864:LEU:HD11	1:A:907:GLU:HG3	1.85	0.59
1:A:955:GLU:N	1:A:955:GLU:OE1	2.35	0.59
1:A:100:GLU:O	1:A:104:LYS:HG3	2.03	0.59
1:A:733:GLN:HG3	1:A:792:VAL:HG21	1.84	0.59
2:B:10:GLN:HB3	2:B:60:LYS:HB3	1.84	0.58
1:A:939:LEU:HG	1:A:1012:LYS:HG3	1.86	0.58
4:D:154:PRO:HD3	4:D:344:ARG:HB2	1.85	0.58
3:C:155:LEU:HG	3:C:225:ASN:HB2	1.86	0.57
1:A:231:ARG:O	1:A:234:ASN:HB2	2.05	0.57
1:A:756:LEU:HB3	1:A:803:THR:HG21	1.85	0.57
1:A:298:LEU:O	1:A:353:TYR:OH	2.18	0.57
1:A:8:LEU:HD22	1:A:45:MET:SD	2.45	0.57
1:A:828:GLU:O	1:A:832:ASN:ND2	2.38	0.57
4:D:274:GLU:O	4:D:278:GLU:HG2	2.05	0.57
1:A:231:ARG:NH2	9:A:1202:HOH:O	2.37	0.57
3:C:278:ARG:HB2	3:C:281:MET:HG3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:277:LEU:HD21	3:C:285:VAL:HG21	1.88	0.56
1:A:872:LYS:NZ	1:A:876:ASP:OD2	2.38	0.55
1:A:925:HIS:HA	4:D:1986:PRO:HB3	1.86	0.55
4:D:1949:SER:OG	4:D:1950:GLY:N	2.40	0.55
2:B:106:ARG:HG3	2:B:107:ASP:N	2.20	0.55
1:A:73:MET:HE2	1:A:123:GLU:HB3	1.88	0.55
4:D:116:ILE:HD11	4:D:232:TRP:CD1	2.42	0.55
1:A:60:TRP:CG	1:A:86:VAL:HG21	2.43	0.55
4:D:16:GLY:HA2	4:D:297:LYS:HD2	1.89	0.55
1:A:837:ASP:O	1:A:845:ARG:NH2	2.40	0.54
4:D:64:HIS:ND1	4:D:96:ALA:HB1	2.22	0.54
1:A:153:ILE:HG22	4:D:2024:PHE:CG	2.43	0.54
1:A:484:VAL:HA	1:A:527:LEU:HD13	1.90	0.54
3:C:155:LEU:HG	3:C:225:ASN:CB	2.38	0.53
1:A:615:LEU:HB3	1:A:656:LYS:HD2	1.90	0.53
1:A:113:THR:HG23	1:A:122:LYS:HE3	1.91	0.53
1:A:438:GLY:N	1:A:746:ARG:HD3	2.24	0.53
4:D:6:LYS:HD2	4:D:34:LYS:HD2	1.91	0.53
1:A:619:ASN:HB2	1:A:660:LEU:HD21	1.91	0.52
1:A:281:GLU:OE2	1:A:340:ARG:NH2	2.42	0.52
3:C:261:HIS:CD2	3:C:263:GLN:H	2.26	0.52
1:A:152:ASP:HB3	4:D:2025:GLY:H	1.75	0.52
1:A:60:TRP:HB3	1:A:86:VAL:HG21	1.92	0.52
1:A:1033:LEU:HD12	1:A:1033:LEU:H	1.75	0.52
2:B:77:ASP:OD1	2:B:77:ASP:N	2.42	0.52
3:C:160:ASN:CG	3:C:162:ARG:HG2	2.35	0.51
4:D:1939:GLY:C	4:D:1940:GLU:HG3	2.36	0.51
1:A:19:PHE:HA	1:A:62:ARG:HH21	1.71	0.51
4:D:89:LEU:HD13	4:D:107:PRO:HG2	1.91	0.51
4:D:192:LEU:HD23	4:D:357:VAL:HG13	1.91	0.51
3:C:276:TRP:CE3	6:C:301:PRO:HG3	2.46	0.51
4:D:235:ILE:HG22	4:D:235:ILE:O	2.09	0.51
4:D:108:ILE:HD13	4:D:285:LEU:HD21	1.93	0.51
3:C:60:ASN:O	3:C:64:ARG:HG3	2.11	0.50
1:A:40:GLY:O	1:A:44:ARG:HB2	2.11	0.50
3:C:215:GLU:C	3:C:216:GLU:HG2	2.36	0.50
4:D:336:MET:O	4:D:339:PHE:HB3	2.11	0.50
1:A:17:LEU:HD13	1:A:56:HIS:ND1	2.27	0.50
2:B:139:HIS:HB2	2:B:144:LEU:HB2	1.92	0.50
1:A:953:VAL:HG11	1:A:974:LEU:HD12	1.93	0.50
2:B:89:MET:HE2	2:B:149:ILE:HD11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:PHE:HB2	1:A:329:PHE:CZ	2.46	0.50
1:A:882:PHE:HA	1:A:890:ALA:HA	1.93	0.50
4:D:235:ILE:HG21	4:D:242:TYR:HD2	1.75	0.50
4:D:339:PHE:CE2	4:D:343:VAL:HG21	2.47	0.50
4:D:119:LYS:HE3	4:D:243:GLY:HA3	1.93	0.49
4:D:119:LYS:HB2	4:D:241:ASN:HB3	1.94	0.49
1:A:732:ILE:O	1:A:736:GLY:N	2.45	0.49
1:A:811:LEU:HD12	1:A:815:ILE:HD12	1.93	0.49
3:C:257:LEU:HD12	3:C:282:VAL:HG21	1.94	0.49
1:A:879:ILE:HA	1:A:882:PHE:CE2	2.47	0.49
1:A:991:ASP:O	1:A:995:LYS:HG2	2.12	0.49
3:C:202:ASP:OD1	3:C:203:PHE:N	2.46	0.49
1:A:568:LYS:HE2	3:C:12:PHE:HB3	1.95	0.49
1:A:60:TRP:HA	1:A:63:VAL:HG23	1.95	0.48
4:D:338:ALA:HA	4:D:341:TYR:CD2	2.47	0.48
1:A:78:TYR:CE2	1:A:82:ILE:HD11	2.48	0.48
1:A:438:GLY:H	1:A:746:ARG:HD3	1.78	0.48
1:A:676:VAL:HG12	4:D:1922:PHE:HB2	1.95	0.48
3:C:171:ILE:HG22	3:C:189:MET:HE3	1.94	0.48
1:A:202:PHE:CZ	1:A:236:ILE:HG21	2.49	0.48
1:A:172:LYS:HB2	1:A:228:THR:HG23	1.95	0.48
1:A:490:SER:HB3	1:A:493:ASN:OD1	2.14	0.48
4:D:160:LEU:HD23	4:D:163:ALA:HB3	1.96	0.48
2:B:89:MET:CE	2:B:149:ILE:HD11	2.43	0.48
3:C:264:THR:HG22	3:C:265:HIS:O	2.14	0.48
4:D:1915:PHE:CD1	4:D:1915:PHE:N	2.76	0.48
3:C:115:LEU:HD12	3:C:119:TRP:HB2	1.95	0.47
1:A:1002:PHE:O	1:A:1005:ASN:ND2	2.47	0.47
4:D:1920:ASN:N	4:D:1920:ASN:OD1	2.46	0.47
1:A:119:CYS:HA	1:A:122:LYS:HB3	1.95	0.47
1:A:1005:ASN:OD1	1:A:1006:GLN:HG3	2.14	0.47
2:B:32:THR:OG1	2:B:33:GLY:N	2.48	0.47
1:A:657:TYR:CE2	1:A:658:MET:HE3	2.49	0.47
1:A:847:ASN:HA	1:A:850:LEU:HD12	1.95	0.47
3:C:56:LEU:N	3:C:56:LEU:HD12	2.30	0.47
3:C:201:THR:HG23	3:C:264:THR:O	2.14	0.47
4:D:19:GLY:HA2	4:D:22:GLU:CD	2.40	0.47
4:D:1933:PHE:CD1	4:D:1933:PHE:C	2.92	0.47
1:A:132:MET:HE1	1:A:177:GLU:HG3	1.97	0.47
3:C:96:LYS:O	3:C:97:HIS:HB2	2.14	0.47
4:D:311:LEU:HB3	4:D:317:ILE:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:153:SER:O	3:C:226:PRO:HD2	2.15	0.47
4:D:116:ILE:HD11	4:D:232:TRP:HD1	1.80	0.47
1:A:631:HIS:HB3	1:A:697:ARG:HG3	1.97	0.46
1:A:849:PHE:CE1	1:A:881:ALA:HB2	2.50	0.46
2:B:75:LEU:HD12	2:B:79:TYR:CE1	2.51	0.46
1:A:770:GLN:O	1:A:774:GLU:HG2	2.15	0.46
4:D:155:TYR:HE2	4:D:258:PHE:CD1	2.34	0.46
1:A:508:MET:HE1	1:A:516:PHE:CD2	2.51	0.46
1:A:513:GLU:CD	1:A:553:ARG:HH21	2.23	0.46
3:C:105:SER:HB3	3:C:274:VAL:HG22	1.98	0.46
1:A:56:HIS:HD2	1:A:57:PRO:HD2	1.80	0.46
1:A:996:LEU:HD21	1:A:1038:ARG:HG2	1.98	0.46
3:C:207:TRP:CH2	3:C:211:LYS:HD3	2.51	0.46
1:A:8:LEU:HD12	1:A:8:LEU:H	1.81	0.46
1:A:104:LYS:HB3	4:D:2021:ALA:CB	2.46	0.46
1:A:743:PRO:O	1:A:746:ARG:HB3	2.16	0.45
1:A:868:PRO:HG2	4:D:1946:THR:OG1	2.17	0.45
4:D:129:TRP:CZ3	4:D:195:LEU:HD11	2.52	0.45
1:A:60:TRP:CB	1:A:86:VAL:HG21	2.47	0.45
3:C:55:ARG:N	3:C:56:LEU:HD12	2.31	0.45
3:C:125:PRO:HB3	3:C:252:PHE:CG	2.51	0.45
3:C:217:GLY:HA2	3:C:220:GLU:OE1	2.16	0.45
1:A:228:THR:HG22	1:A:232:PHE:CE2	2.51	0.45
2:B:11:PHE:CG	2:B:168:LEU:HD13	2.52	0.45
2:B:117:ILE:HB	2:B:144:LEU:HD22	1.99	0.45
2:B:146:TYR:OH	2:B:148:ASP:OD1	2.24	0.45
1:A:6:THR:O	1:A:10:ASP:HB2	2.16	0.45
1:A:104:LYS:O	4:D:2021:ALA:HB2	2.16	0.45
1:A:1035:LEU:HD12	1:A:1035:LEU:H	1.81	0.45
4:D:2019:ALA:C	4:D:2021:ALA:H	2.25	0.45
1:A:596:ARG:HB2	1:A:643:ALA:HB2	1.99	0.45
1:A:759:ILE:HD13	1:A:780:LEU:HD21	1.99	0.45
1:A:534:LYS:NZ	1:A:575:GLU:OE2	2.50	0.45
1:A:545:MET:HE2	1:A:572:PHE:HD2	1.82	0.45
1:A:727:ASN:HA	4:D:1918:THR:HB	1.99	0.45
3:C:18:LEU:HD23	3:C:35:TYR:HA	1.99	0.45
3:C:125:PRO:HB3	3:C:252:PHE:CD1	2.52	0.45
3:C:139:THR:OG1	3:C:154:SER:HB3	2.17	0.45
4:D:235:ILE:CG2	4:D:242:TYR:HD2	2.30	0.45
4:D:1942:LYS:HA	4:D:1943:PRO:HD3	1.85	0.45
1:A:107:VAL:O	1:A:111:ILE:HG12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:960:THR:HG21	1:A:965:GLY:HA2	1.99	0.44
3:C:196:PHE:O	3:C:204:ARG:HD3	2.17	0.44
3:C:198:ASP:O	3:C:265:HIS:NE2	2.49	0.44
1:A:258:PRO:O	1:A:261:ARG:HG2	2.17	0.44
1:A:788:TYR:CZ	1:A:796:ARG:HD3	2.52	0.44
1:A:1031:SER:HA	1:A:1033:LEU:CD1	2.47	0.44
1:A:564:THR:HG23	3:C:9:ALA:HB2	1.99	0.44
2:B:145:GLN:HE21	2:B:163:TRP:NE1	2.14	0.44
3:C:216:GLU:O	3:C:218:LEU:N	2.49	0.44
4:D:339:PHE:CD2	4:D:339:PHE:C	2.95	0.44
1:A:35:LEU:HD22	1:A:50:LEU:HD12	1.98	0.44
1:A:304:ILE:HG13	1:A:356:LEU:HB3	2.00	0.44
1:A:490:SER:OG	5:E:1:GLC:H61	2.18	0.44
3:C:97:HIS:O	3:C:98:TYR:HB2	2.16	0.44
3:C:112:PRO:HG3	3:C:119:TRP:CE2	2.52	0.44
1:A:63:VAL:HG11	1:A:83:LEU:HD22	2.00	0.44
1:A:897:LEU:HD23	1:A:945:ILE:HD13	1.99	0.44
4:D:62:TRP:CG	4:D:63:ALA:N	2.85	0.44
1:A:635:GLU:OE1	1:A:697:ARG:HD2	2.17	0.44
4:D:65:ASP:OD2	4:D:66:ARG:HG2	2.18	0.44
1:A:146:TRP:CD1	1:A:149:PHE:HB2	2.53	0.43
1:A:565:VAL:O	1:A:569:LEU:HD13	2.18	0.43
1:A:730:ALA:HB3	4:D:1918:THR:HG22	1.99	0.43
1:A:120:VAL:HG13	1:A:124:LYS:HE3	1.98	0.43
1:A:667:SER:O	1:A:671:GLN:HG3	2.18	0.43
1:A:747:SER:O	1:A:750:THR:HB	2.19	0.43
3:C:54:LYS:H	3:C:54:LYS:HG2	1.57	0.43
4:D:90:TYR:CE1	4:D:305:LYS:HG2	2.53	0.43
1:A:172:LYS:HE2	1:A:172:LYS:HB3	1.88	0.43
1:A:766:SER:O	1:A:810:LYS:NZ	2.44	0.43
1:A:518:VAL:O	1:A:522:LYS:HB2	2.18	0.43
1:A:168:MET:HE2	1:A:225:THR:HA	2.00	0.43
1:A:920:CYS:O	1:A:924:GLN:HG2	2.19	0.43
1:A:123:GLU:O	1:A:127:ILE:HG13	2.19	0.43
1:A:154:VAL:HG22	1:A:208:LEU:HD12	2.01	0.43
3:C:256:GLY:O	3:C:257:LEU:HD23	2.19	0.43
4:D:43:LEU:HD12	4:D:44:GLU:N	2.32	0.43
4:D:115:LEU:HD13	4:D:245:THR:HG23	2.00	0.43
1:A:968:VAL:CG1	1:A:973:PHE:HB2	2.49	0.43
2:B:71:LYS:HE3	2:B:71:LYS:HB2	1.84	0.43
1:A:774:GLU:HG3	1:A:775:ASN:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:781:LEU:O	1:A:786:ILE:HG13	2.19	0.42
2:B:145:GLN:HE21	2:B:163:TRP:CD1	2.37	0.42
3:C:43:ARG:NH2	3:C:106:GLU:OE2	2.51	0.42
4:D:336:MET:HE3	4:D:336:MET:HB3	1.90	0.42
1:A:300:LEU:HB3	1:A:352:HIS:CE1	2.55	0.42
1:A:559:TRP:CD2	1:A:603:GLN:HG3	2.55	0.42
4:D:1917:ASN:OD1	4:D:1918:THR:N	2.52	0.42
1:A:19:PHE:CA	1:A:62:ARG:HH21	2.32	0.42
1:A:904:VAL:HG11	1:A:914:PHE:CD2	2.53	0.42
3:C:55:ARG:H	3:C:56:LEU:HD12	1.85	0.42
4:D:1926:GLY:HA2	4:D:1929:THR:HG22	2.00	0.42
1:A:626:GLN:HG2	1:A:629:GLN:HG3	2.02	0.42
1:A:118:THR:O	1:A:122:LYS:HB2	2.20	0.42
1:A:256:ASN:HA	1:A:297:MET:SD	2.60	0.42
1:A:435:ASN:CG	1:A:436:ASP:N	2.78	0.42
3:C:25:HIS:HA	3:C:26:PRO:HD3	1.87	0.42
4:D:155:TYR:CE2	4:D:258:PHE:CD1	3.08	0.42
4:D:233:SER:O	4:D:236:ASP:CG	2.63	0.42
1:A:344:ARG:HA	1:A:344:ARG:HD3	1.75	0.42
1:A:1033:LEU:HD12	1:A:1033:LEU:N	2.33	0.42
3:C:215:GLU:O	3:C:215:GLU:HG3	2.20	0.42
1:A:737:GLU:O	1:A:740:THR:HG22	2.20	0.41
1:A:821:GLN:H	1:A:821:GLN:HG2	1.68	0.41
1:A:104:LYS:HA	4:D:2021:ALA:HB1	2.02	0.41
1:A:531:LYS:HD3	1:A:531:LYS:HA	1.89	0.41
3:C:187:ASP:OD1	3:C:188:VAL:N	2.53	0.41
4:D:1937:SER:O	4:D:1938:PHE:HB2	2.20	0.41
1:A:575:GLU:CD	3:C:15:SER:HG	2.26	0.41
1:A:164:CYS:O	1:A:168:MET:HG2	2.20	0.41
1:A:262:ASN:O	1:A:266:LYS:HG3	2.20	0.41
1:A:650:GLN:O	1:A:654:ILE:HG13	2.20	0.41
3:C:277:LEU:HG	3:C:281:MET:HB2	2.01	0.41
1:A:219:ALA:HB3	1:A:220:PRO:HD3	2.03	0.41
1:A:101:GLY:HA3	4:D:2011:VAL:HG13	2.03	0.41
1:A:107:VAL:HG11	1:A:146:TRP:HH2	1.86	0.41
1:A:142:TRP:HD1	1:A:146:TRP:O	2.04	0.41
1:A:323:SER:HB3	1:A:357:VAL:HG11	2.03	0.41
1:A:665:TRP:O	1:A:669:ILE:HG12	2.21	0.41
1:A:762:TRP:CH2	1:A:772:VAL:HG22	2.55	0.41
1:A:822:ILE:O	1:A:826:VAL:HG22	2.21	0.41
1:A:867:PRO:HA	1:A:868:PRO:HD3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:962:LEU:HB3	1:A:964:PRO:HD2	2.02	0.41
4:D:209:ASP:OD1	4:D:210:TYR:N	2.50	0.41
1:A:104:LYS:CA	4:D:2021:ALA:HB1	2.51	0.41
1:A:953:VAL:HG21	1:A:974:LEU:CD1	2.48	0.41
4:D:121:LEU:HB3	4:D:139:LEU:HD11	2.02	0.41
1:A:44:ARG:O	1:A:48:GLU:HG3	2.21	0.40
2:B:29:ARG:HG2	2:B:157:PHE:CZ	2.57	0.40
1:A:771:MET:HE2	1:A:771:MET:HB3	1.98	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	1028/1044 (98%)	983 (96%)	45 (4%)	0	100	100
2	B	170/172 (99%)	162 (95%)	8 (5%)	0	100	100
3	C	256/289 (89%)	241 (94%)	15 (6%)	0	100	100
4	D	371/479 (78%)	334 (90%)	34 (9%)	3 (1%)	16	31
All	All	1825/1984 (92%)	1720 (94%)	102 (6%)	3 (0%)	43	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	1941	GLN
4	D	1933	PHE
4	D	1943	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	941/951 (99%)	922 (98%)	19 (2%)	48	71
2	B	151/151 (100%)	148 (98%)	3 (2%)	48	71
3	C	237/262 (90%)	228 (96%)	9 (4%)	29	54
4	D	304/360 (84%)	292 (96%)	12 (4%)	28	54
All	All	1633/1724 (95%)	1590 (97%)	43 (3%)	40	65

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	MET
1	A	14	ARG
1	A	47	GLN
1	A	51	THR
1	A	107	VAL
1	A	123	GLU
1	A	208	LEU
1	A	257	VAL
1	A	263	VAL
1	A	345	GLU
1	A	678	ILE
1	A	875	LEU
1	A	885	THR
1	A	952	LEU
1	A	984	SER
1	A	989	LEU
1	A	1022	ILE
1	A	1033	LEU
1	A	1036	GLU
2	B	40	VAL
2	B	77	ASP
2	B	154	ASN
3	C	52	LYS
3	C	54	LYS

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Mol	Chain	Res	Type
3	C	56	LEU
3	C	94	LEU
3	C	111	VAL
3	C	179	VAL
3	C	211	LYS
3	C	222	THR
3	C	274	VAL
4	D	12	ASN
4	D	155	TYR
4	D	160	LEU
4	D	178	ILE
4	D	333	ILE
4	D	1915	PHE
4	D	1920	ASN
4	D	1924	ASN
4	D	2011	VAL
4	D	2014	GLU
4	D	2020	SER
4	D	2026	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	131	ASN
1	A	136	GLN
1	A	145	HIS
1	A	166	ASN
1	A	333	HIS
1	A	809	ASN
1	A	844	HIS
1	A	858	HIS
2	B	10	GLN
2	B	82	GLN
2	B	105	HIS
3	C	100	ASN
3	C	101	GLN
3	C	200	GLN
3	C	225	ASN
3	C	261	HIS
4	D	49	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 ⓘ

2 monosaccharides 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	$\# Z > 2$
5	GLC	E	1	5	12,12,12	0.52	0	17,17,17	1.23	1 (5%)
5	GLC	E	2	5	11,11,12	0.28	0	15,15,17	2.37	3 (20%)

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
5	GLC	E	1	5	-	2/2/22/22	0/1/1/1
5	GLC	E	2	5	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	2	GLC	C1-O5-C5	7.58	122.35	112.19
5	E	1	GLC	O4-C4-C3	3.89	119.54	110.38
5	E	2	GLC	C3-C4-C5	3.11	115.86	110.23
5	E	2	GLC	O5-C5-C4	3.03	118.20	110.83

There are no chirality outliers.

All (3) torsion outliers are listed below:

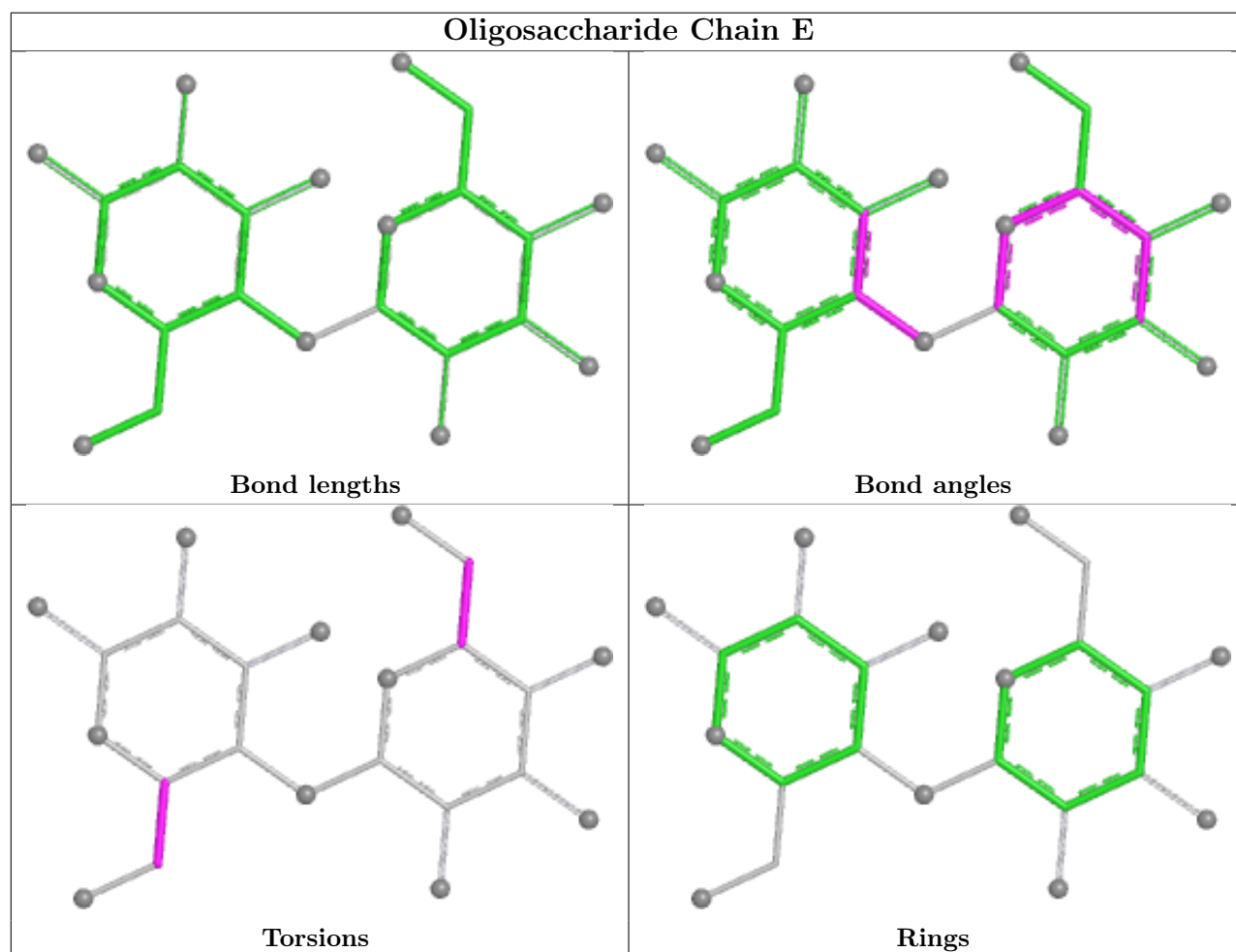
Mol	Chain	Res	Type	Atoms
5	E	1	GLC	O5-C5-C6-O6
5	E	2	GLC	O5-C5-C6-O6
5	E	1	GLC	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	1	GLC	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 oligosaccharide.



5.6 Ligand geometry

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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$
7	GTP	B	201	8	33,34,34	1.46	6 (18%)	50,54,54	1.74	7 (14%)
6	PRO	A	1101	-	8,8,8	0.88	1 (12%)	10,10,10	1.38	2 (20%)
6	PRO	C	301	-	8,8,8	0.83	0	10,10,10	1.05	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	GTP	B	201	8	-	1/22/38/38	0/3/3/3
6	PRO	A	1101	-	-	0/4/11/11	0/1/1/1
6	PRO	C	301	-	-	0/4/11/11	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	201	GTP	PB-O3B	4.20	1.64	1.59
7	B	201	GTP	PB-O3A	2.84	1.62	1.59
7	B	201	GTP	C5-C4	2.69	1.46	1.38
7	B	201	GTP	C6-N1	-2.53	1.34	1.38
7	B	201	GTP	PA-O3A	2.30	1.62	1.59
6	A	1101	PRO	OXT-C	-2.17	1.23	1.30
7	B	201	GTP	C5-N7	-2.06	1.34	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	201	GTP	C5-C4-N3	-6.18	118.56	128.39
7	B	201	GTP	C2-N3-C4	5.41	121.63	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	201	GTP	N9-C4-N3	4.52	135.00	125.95
7	B	201	GTP	C6-C5-N7	3.25	136.20	130.29
7	B	201	GTP	C4-C5-N7	-2.58	106.58	110.67
6	A	1101	PRO	OXT-C-O	-2.56	118.27	124.08
6	A	1101	PRO	OXT-C-CA	2.18	120.88	113.51
7	B	201	GTP	O6-C6-C5	-2.15	120.84	126.53
7	B	201	GTP	C8-N7-C5	2.08	107.96	104.26

There are no chirality outliers.

All (1) torsion outliers are listed below:

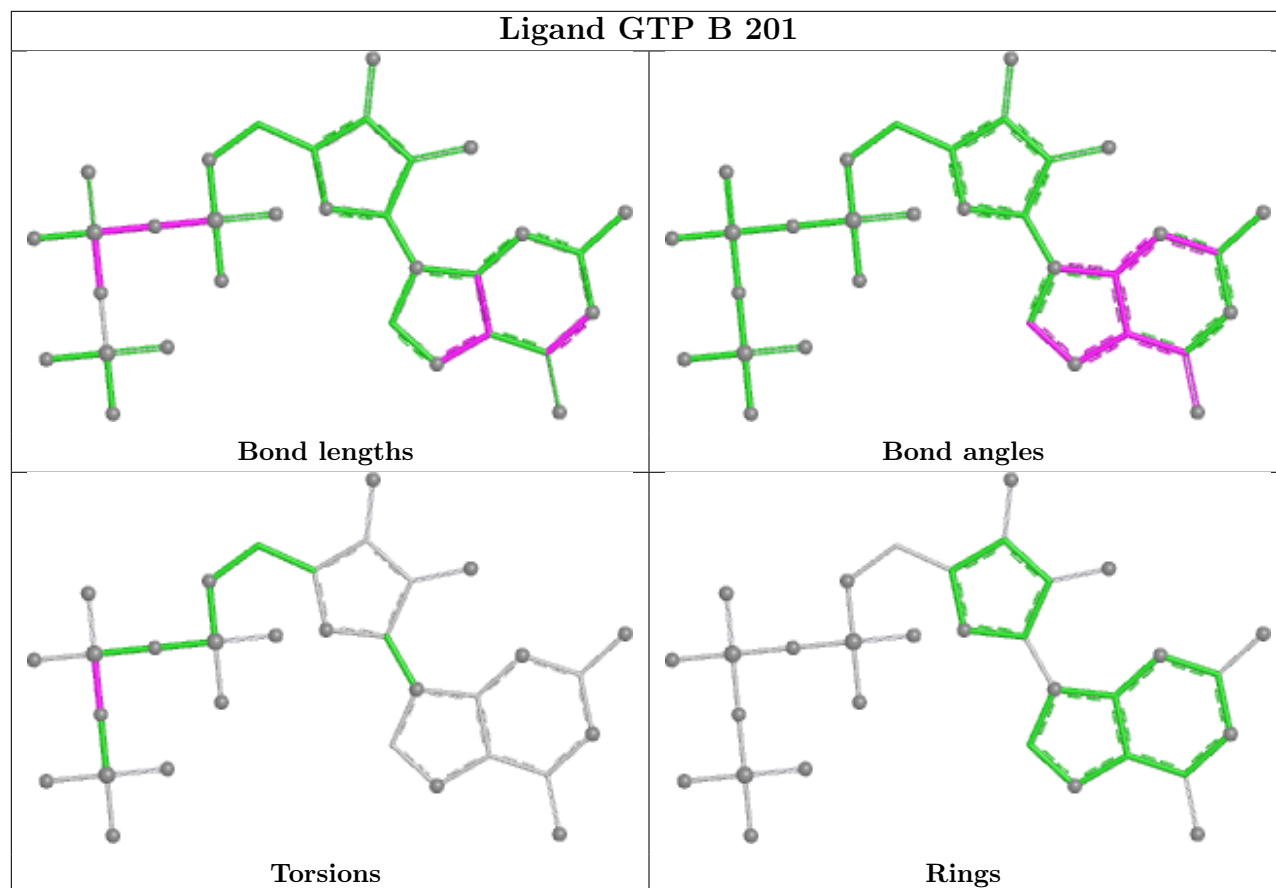
Mol	Chain	Res	Type	Atoms
7	B	201	GTP	PG-O3B-PB-O2B

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	301	PRO	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	1032/1044 (98%)	-0.22	7 (0%) 84 80	59, 92, 145, 259	0
2	B	172/172 (100%)	-0.34	0 100 100	56, 83, 128, 204	0
3	C	264/289 (91%)	0.10	8 (3%) 52 44	72, 125, 190, 230	0
4	D	395/479 (82%)	0.75	30 (7%) 20 14	109, 213, 281, 329	0
All	All	1863/1984 (93%)	0.02	45 (2%) 59 51	56, 105, 246, 329	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	55	ARG	5.1
4	D	16	GLY	4.8
4	D	346	ALA	4.3
4	D	197	ASP	3.7
4	D	182	GLY	3.4
4	D	1915	PHE	3.4
3	C	166	ALA	3.4
4	D	368	ALA	3.2
1	A	962	LEU	3.2
4	D	188	ALA	3.0
4	D	116	ILE	3.0
4	D	59	ILE	2.8
3	C	287	GLY	2.8
4	D	87	ASP	2.8
4	D	114	SER	2.8
4	D	1918	THR	2.8
4	D	1926	GLY	2.7
4	D	1914	GLU	2.7
4	D	195	LEU	2.7
4	D	166	GLY	2.6
4	D	201	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
4	D	2022	GLY	2.6
4	D	361	LEU	2.6
4	D	129	TRP	2.5
4	D	178	ILE	2.5
4	D	151	LEU	2.5
4	D	122	LEU	2.5
3	C	94	LEU	2.4
4	D	2023	GLY	2.4
1	A	238	LEU	2.4
3	C	28	LEU	2.3
4	D	287	ASP	2.3
4	D	83	LYS	2.3
1	A	388	THR	2.3
4	D	1921	LEU	2.2
4	D	300	GLY	2.2
4	D	15	LYS	2.2
4	D	84	ALA	2.2
1	A	1029	ASP	2.2
3	C	126	VAL	2.1
1	A	118	THR	2.1
3	C	255	ASP	2.1
3	C	285	VAL	2.1
1	A	409	TYR	2.1
1	A	1033	LEU	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.

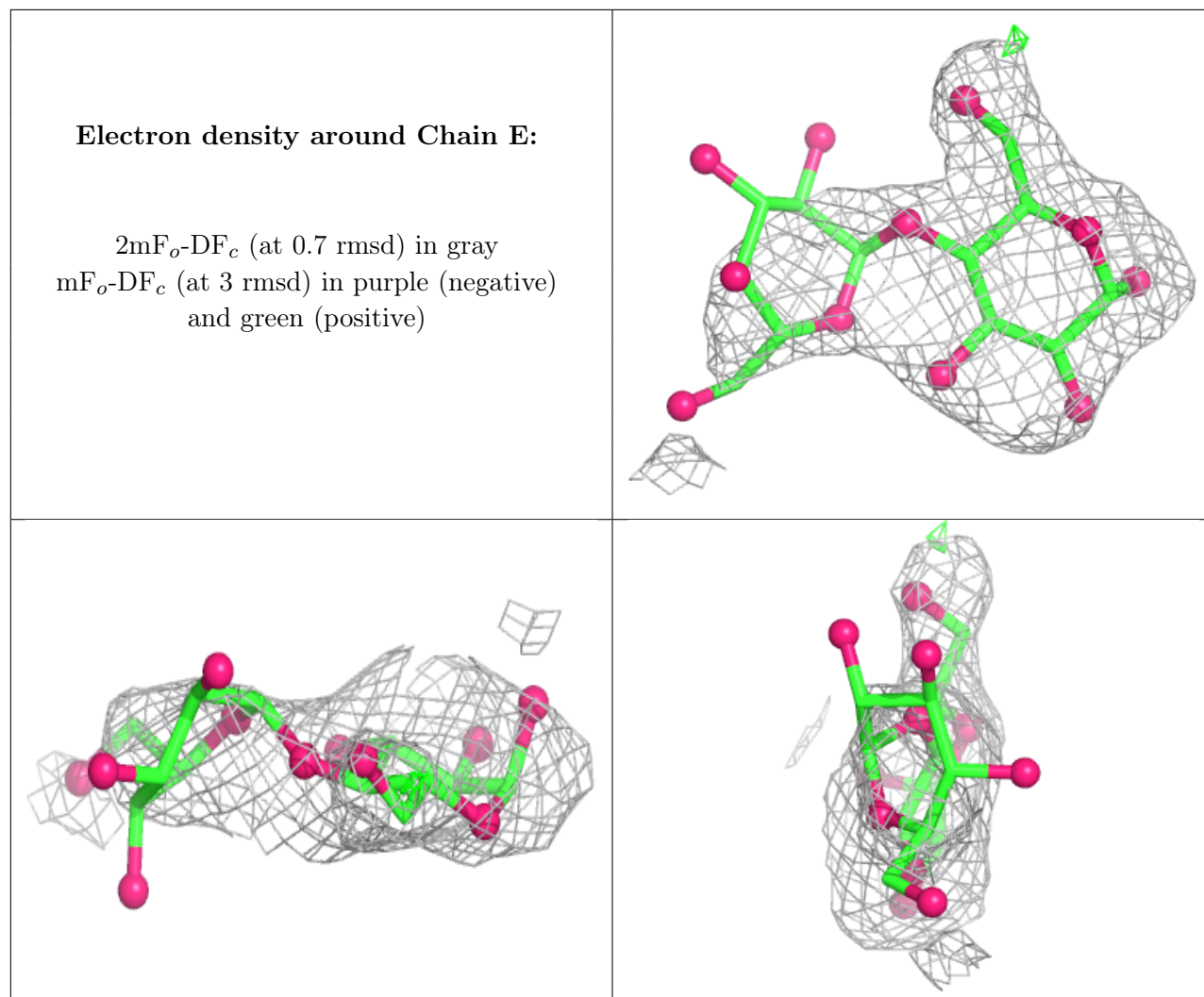
6.3 Carbohydrates [i](#)

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
5	GLC	E	2	11/12	0.28	0.10	224,231,235,236	0
5	GLC	E	1	12/12	0.70	0.11	153,175,193,210	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. Each fit is shown from different orientation to approximate a three-dimensional view.



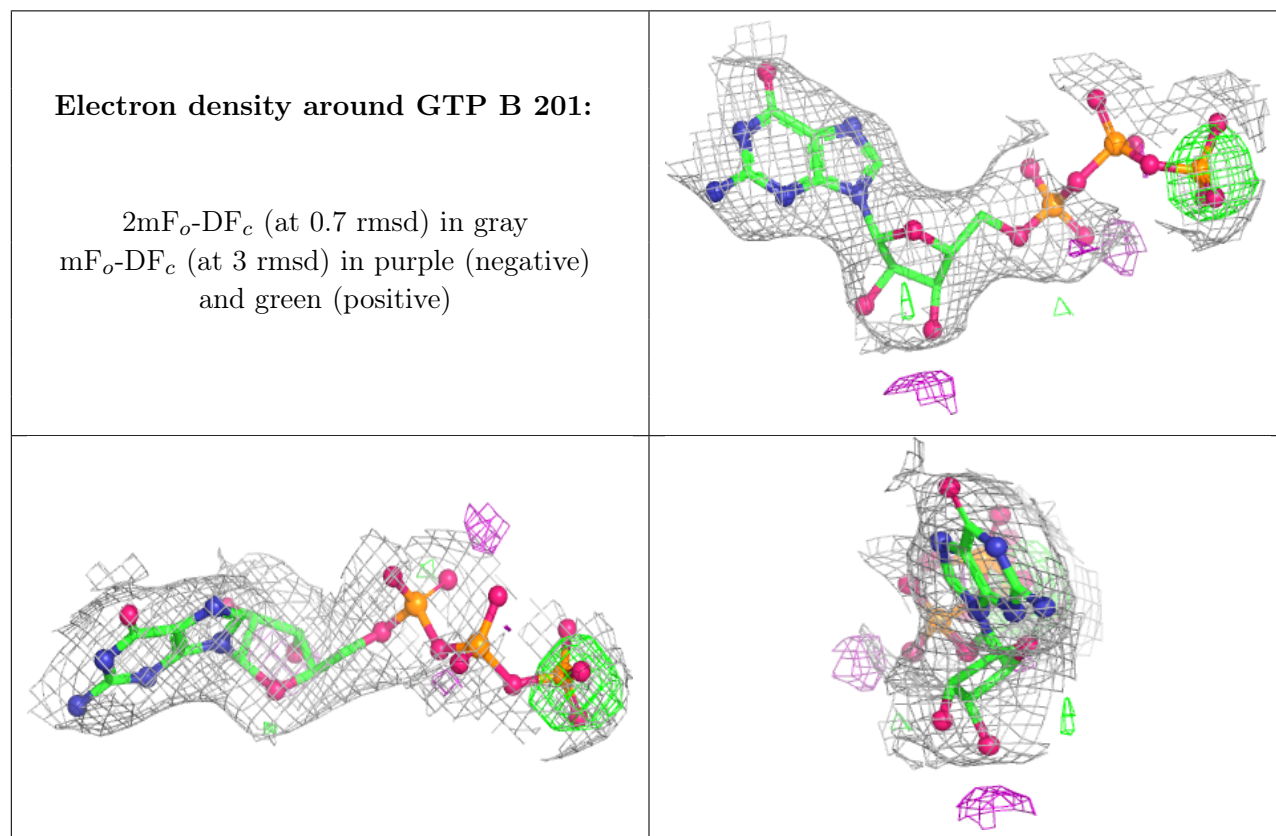
6.4 Ligands [i](#)

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
6	PRO	A	1101	8/8	0.85	0.21	122,126,137,140	0
6	PRO	C	301	8/8	0.93	0.38	166,188,275,329	0
7	GTP	B	201	32/32	0.95	0.07	56,75,106,135	0
8	MG	B	202	1/1	0.98	0.04	94,94,94,94	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.