



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 11:43 AM BST

PDB ID : 2NZY
Title : Crystal Structure of alpha1,3-Fucosyltransferase with GDP-fucose
Authors : Sun, H.Y.; Ko, T.P.
Deposited on : 2006-11-27
Resolution : 2.05 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

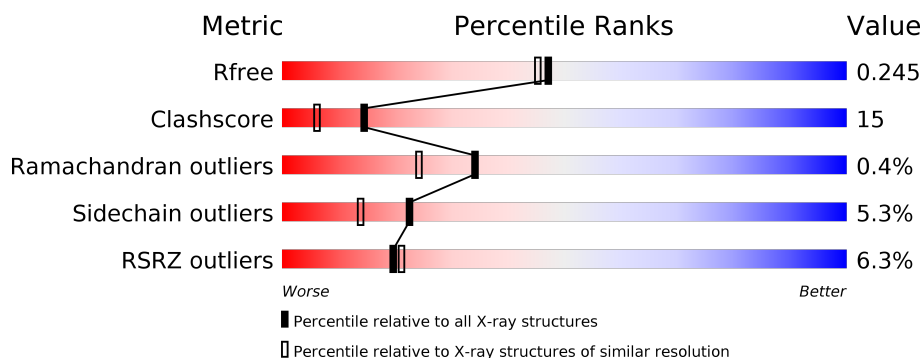
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 on 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	371	<div> <div>7%</div> <div> <div></div> <div>65%</div> <div>23%</div> <div>• • 8%</div> </div> </div>
1	B	371	<div> <div>6%</div> <div> <div></div> <div>64%</div> <div>23%</div> <div>• 9%</div> </div> </div>
1	C	371	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>20%</div> <div>• 5%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FUC	A	4001	X	-	-	-
2	FUC	B	4002	X	-	-	-
2	FUC	C	4003	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9905 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 Alpha1,3-Fucosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	0	0
			2825	1835	457	526	7			
1	B	338	Total	C	N	O	S	0	0	0
			2776	1797	452	520	7			
1	C	351	Total	C	N	O	S	0	0	0
			2879	1867	469	536	7			

There are 24 discrepancies between the modelled and reference sequences:

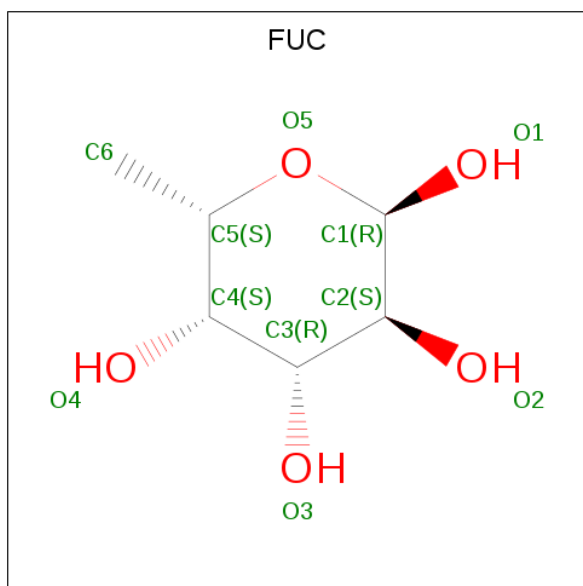
Chain	Residue	Modelled	Actual	Comment	Reference
A	364	LEU	-	expression tag	UNP O30511
A	365	GLU	-	expression tag	UNP O30511
A	366	HIS	-	expression tag	UNP O30511
A	367	HIS	-	expression tag	UNP O30511
A	368	HIS	-	expression tag	UNP O30511
A	369	HIS	-	expression tag	UNP O30511
A	370	HIS	-	expression tag	UNP O30511
A	371	HIS	-	expression tag	UNP O30511
B	364	LEU	-	expression tag	UNP O30511
B	365	GLU	-	expression tag	UNP O30511
B	366	HIS	-	expression tag	UNP O30511
B	367	HIS	-	expression tag	UNP O30511
B	368	HIS	-	expression tag	UNP O30511
B	369	HIS	-	expression tag	UNP O30511
B	370	HIS	-	expression tag	UNP O30511
B	371	HIS	-	expression tag	UNP O30511
C	364	LEU	-	expression tag	UNP O30511
C	365	GLU	-	expression tag	UNP O30511
C	366	HIS	-	expression tag	UNP O30511
C	367	HIS	-	expression tag	UNP O30511
C	368	HIS	-	expression tag	UNP O30511
C	369	HIS	-	expression tag	UNP O30511
C	370	HIS	-	expression tag	UNP O30511

Continued on next page...

Continued from previous page...

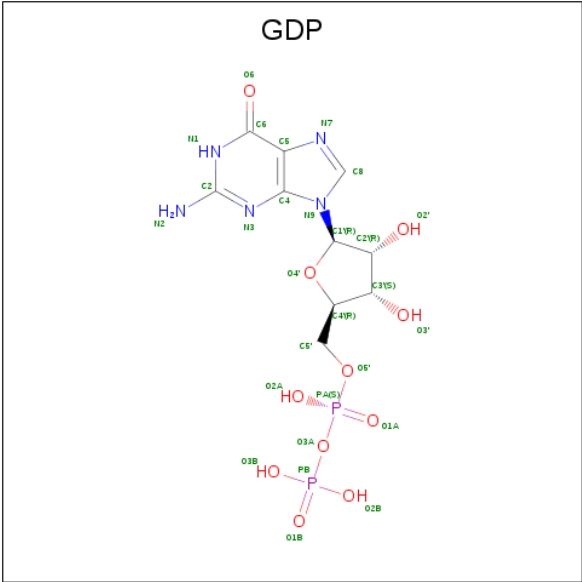
Chain	Residue	Modelled	Actual	Comment	Reference
C	371	HIS	-	expression tag	UNP O30511

- Molecule 2 is alpha-L-fucopyranose (three-letter code: FUC) (formula: $C_6H_{12}O_5$).



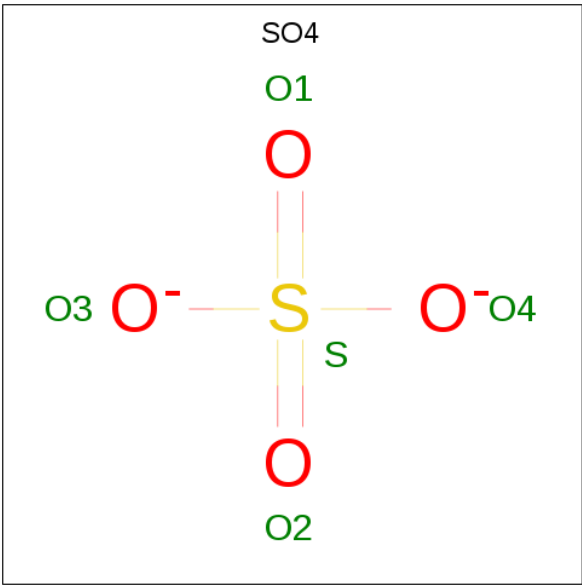
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	C	1	Total	C	O	0	0
			10	6	4		

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
3	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		

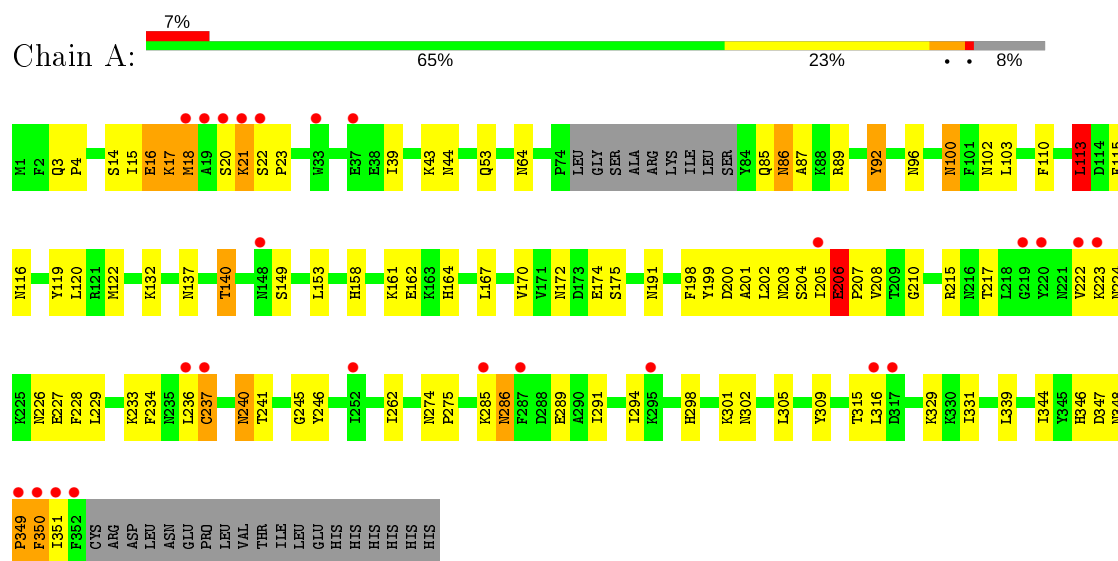
- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	432	Total	O		0	0
			432	432			
5	B	396	Total	O		0	0
			396	396			
5	C	473	Total	O		0	0
			473	473			

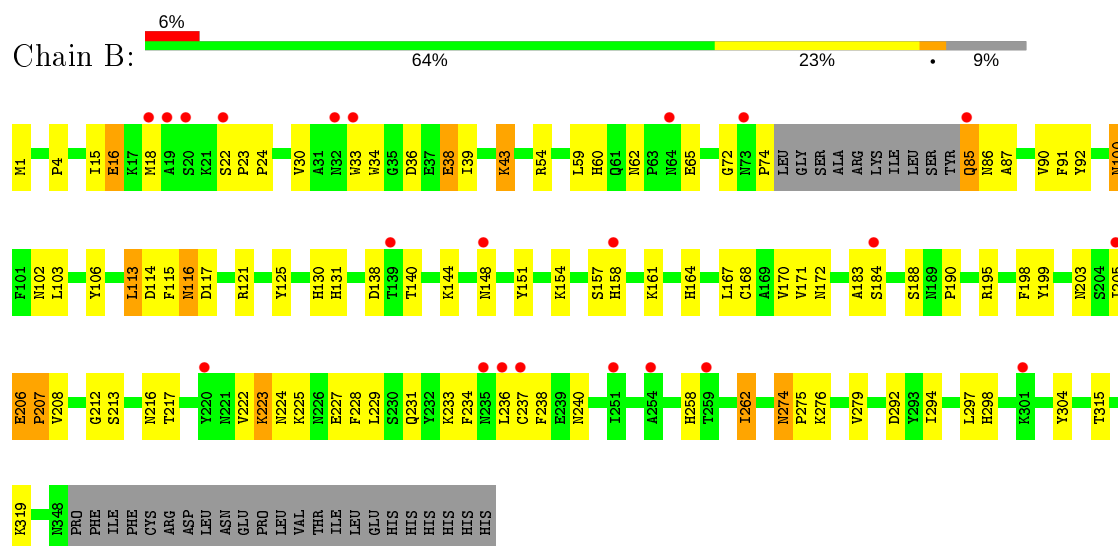
3 Residue-property plots

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: Alpha1,3-Fucosyltransferase

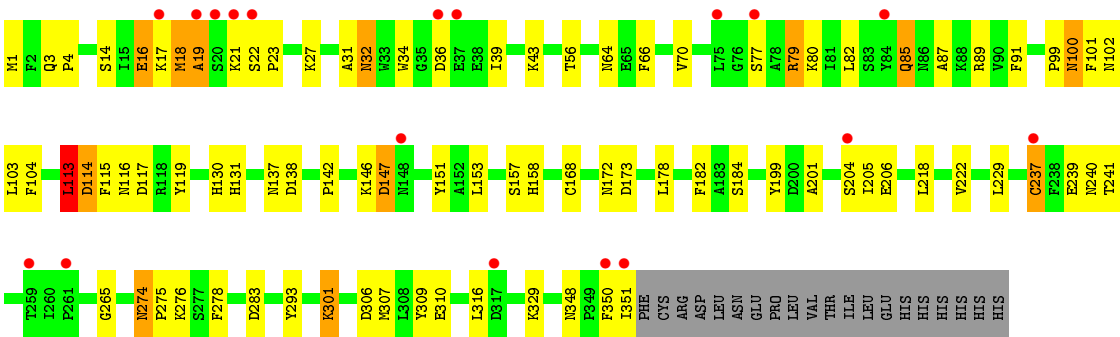


• Molecule 1: Alpha1,3-Fucosyltransferase



• Molecule 1: Alpha1,3-Fucosyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	103.96 Å 135.74 Å 96.19 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.05 29.89 – 2.05	Depositor EDS
% Data completeness (in resolution range)	92.0 (30.00-2.05) 92.2 (29.89-2.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 2.04 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.190 , 0.246 0.188 , 0.245	Depositor DCC
R_{free} test set	3989 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.6	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	0.96	EDS
Total number of atoms	9905	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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: GDP, FUC, SO4

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	0.90	1/2912 (0.0%)	0.87	6/3953 (0.2%)
1	B	0.98	3/2859 (0.1%)	0.89	3/3880 (0.1%)
1	C	1.02	3/2966 (0.1%)	0.92	7/4026 (0.2%)
All	All	0.97	7/8737 (0.1%)	0.90	16/11859 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	237	CYS	CB-SG	-12.10	1.61	1.82
1	B	237	CYS	CB-SG	-8.31	1.68	1.82
1	A	237	CYS	CB-SG	-7.05	1.70	1.82
1	B	319	LYS	CD-CE	5.99	1.66	1.51
1	B	92	TYR	CD1-CE1	5.91	1.48	1.39
1	C	278	PHE	CE2-CZ	5.50	1.47	1.37
1	C	278	PHE	CE1-CZ	5.01	1.46	1.37

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121	ARG	NE-CZ-NH2	-7.44	116.58	120.30
1	A	350	PHE	N-CA-C	-6.69	92.94	111.00
1	A	113	LEU	CA-CB-CG	6.65	130.59	115.30
1	A	113	LEU	CB-CG-CD1	-6.12	100.59	111.00
1	B	262	ILE	N-CA-C	-6.09	94.56	111.00
1	A	339	LEU	CA-CB-CG	-6.02	101.45	115.30
1	C	237	CYS	N-CA-CB	-5.93	99.93	110.60
1	A	229	LEU	CA-CB-CG	5.79	128.61	115.30
1	A	137	ASN	N-CA-C	-5.46	96.27	111.00
1	C	113	LEU	CA-CB-CG	5.45	127.83	115.30
1	C	137	ASN	N-CA-C	-5.45	96.30	111.00
1	C	114	ASP	CB-CG-OD1	5.43	123.19	118.30
1	B	319	LYS	CD-CE-NZ	5.10	123.43	111.70
1	C	113	LEU	CB-CG-CD1	-5.04	102.43	111.00
1	C	350	PHE	N-CA-C	-5.03	97.41	111.00
1	C	79	ARG	NE-CZ-NH1	-5.03	117.79	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	92	TYR	Sidechain
1	B	125	TYR	Sidechain
1	C	293	TYR	Sidechain

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	2825	0	2698	93	0
1	B	2776	0	2653	95	0
1	C	2879	0	2767	76	0
2	A	10	0	10	0	0
2	B	10	0	10	0	0
2	C	10	0	10	1	0
3	A	28	0	12	0	0
3	B	28	0	12	0	0
3	C	28	0	12	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	A	432	0	0	15	0
5	B	396	0	0	18	0
5	C	473	0	0	10	0
All	All	9905	0	8184	258	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 15.

All (258) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:HIS:HD2	1:B:62:ASN:H	1.10	0.98
1:A:286:ASN:HB2	5:A:4201:HOH:O	1.64	0.97
1:C:146:LYS:HG2	5:C:4263:HOH:O	1.70	0.92
1:A:200:ASP:OD1	1:A:215:ARG:NH1	2.05	0.90
1:A:85:GLN:HG2	5:A:4165:HOH:O	1.73	0.88
1:C:130:HIS:HD2	1:C:151:TYR:OH	1.59	0.85
1:B:130:HIS:HD2	1:B:151:TYR:OH	1.60	0.84
1:B:223:LYS:N	1:B:223:LYS:HD3	1.94	0.82
1:B:100:ASN:HD22	1:B:100:ASN:C	1.82	0.82
1:B:224:ASN:HD22	1:B:227:GLU:CB	1.93	0.81
1:B:223:LYS:HD3	1:B:223:LYS:H	1.43	0.81
1:A:100:ASN:C	1:A:100:ASN:HD22	1.84	0.81
1:B:60:HIS:CD2	1:B:62:ASN:H	1.98	0.81
1:A:116:ASN:HB3	5:A:4295:HOH:O	1.81	0.80
1:A:286:ASN:HD21	1:A:289:GLU:HG3	1.45	0.80
1:A:348:ASN:HD22	1:C:64:ASN:ND2	1.81	0.78
1:B:16:GLU:OE1	1:B:16:GLU:N	2.17	0.78
1:C:153:LEU:HD11	5:C:4146:HOH:O	1.85	0.77
1:A:22:SER:HA	5:A:4014:HOH:O	1.85	0.74
1:B:62:ASN:OD1	1:B:65:GLU:HG3	1.87	0.74
1:A:222:VAL:HG21	1:A:228:PHE:HB2	1.68	0.74
1:A:348:ASN:HD22	1:C:64:ASN:HD22	1.36	0.73
1:A:17:LYS:NZ	1:A:17:LYS:HB3	2.04	0.72
1:B:158:HIS:HB2	5:B:4125:HOH:O	1.90	0.72
1:B:33:TRP:CZ2	5:B:4298:HOH:O	2.42	0.72
1:B:222:VAL:HG11	1:B:228:PHE:HB2	1.70	0.72
1:B:224:ASN:HD22	1:B:227:GLU:HB2	1.55	0.70
1:B:100:ASN:ND2	1:B:102:ASN:H	1.89	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:PHE:CD2	1:B:236:LEU:HD11	2.27	0.69
1:A:100:ASN:ND2	1:A:102:ASN:H	1.91	0.69
1:C:201:ALA:O	1:C:205:ILE:HD13	1.93	0.68
1:A:164:HIS:CD2	1:A:224:ASN:HD21	2.12	0.68
1:B:223:LYS:H	1:B:223:LYS:CD	2.06	0.68
1:B:315:THR:HG23	5:B:4183:HOH:O	1.93	0.68
1:A:17:LYS:HB3	1:A:17:LYS:HZ3	1.59	0.67
1:A:198:PHE:CD2	1:A:236:LEU:HD21	2.29	0.67
1:A:14:SER:HA	5:A:4006:HOH:O	1.93	0.67
1:C:85:GLN:HA	1:C:85:GLN:HE21	1.59	0.67
1:A:204:SER:C	1:A:205:ILE:HD13	2.15	0.67
1:B:117:ASP:HB2	5:B:4210:HOH:O	1.95	0.67
1:C:1:MET:HE2	1:C:309:TYR:HD1	1.60	0.67
1:C:283:ASP:OD2	5:C:4126:HOH:O	2.13	0.66
1:B:100:ASN:HD21	1:B:102:ASN:HB2	1.60	0.66
1:B:198:PHE:HD2	1:B:236:LEU:HD11	1.59	0.66
1:A:274:ASN:HD22	1:A:275:PRO:HD2	1.60	0.65
1:B:38:GLU:HG2	5:B:4302:HOH:O	1.96	0.65
1:C:16:GLU:H	1:C:16:GLU:CD	2.00	0.64
1:B:203:ASN:HD22	1:B:208:VAL:H	1.43	0.64
1:B:203:ASN:ND2	1:B:208:VAL:H	1.94	0.64
1:A:316:LEU:HD22	5:A:4119:HOH:O	1.97	0.64
1:C:113:LEU:HD22	1:C:119:TYR:CD2	2.32	0.64
1:A:200:ASP:CG	1:A:215:ARG:HH12	2.01	0.64
1:C:206:GLU:OE2	1:C:206:GLU:HA	1.98	0.64
1:B:224:ASN:HD22	1:B:227:GLU:HB3	1.62	0.63
1:B:22:SER:OG	5:B:4288:HOH:O	2.16	0.63
1:A:301:LYS:HG3	1:A:302:ASN:H	1.64	0.62
1:A:286:ASN:ND2	1:A:289:GLU:HG3	2.15	0.62
1:C:100:ASN:HD22	1:C:102:ASN:H	1.47	0.62
1:C:147:ASP:N	1:C:147:ASP:OD1	2.32	0.62
1:A:301:LYS:HG3	1:A:302:ASN:N	2.14	0.62
1:A:100:ASN:HD22	1:A:102:ASN:H	1.47	0.62
1:A:172:ASN:O	1:A:174:GLU:HG3	2.00	0.62
1:B:140:THR:HB	1:B:161:LYS:HB2	1.82	0.62
1:A:116:ASN:ND2	5:A:4178:HOH:O	2.25	0.62
1:A:285:LYS:HE2	1:A:289:GLU:OE1	2.00	0.61
1:A:198:PHE:HD2	1:A:236:LEU:HD21	1.65	0.61
1:B:39:ILE:HG22	1:B:43:LYS:NZ	2.16	0.61
1:C:113:LEU:C	1:C:113:LEU:HD23	2.21	0.61
1:A:286:ASN:N	1:A:286:ASN:HD22	1.97	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:ASN:HD22	1:C:100:ASN:C	2.03	0.60
1:C:158:HIS:HB2	5:C:4149:HOH:O	2.00	0.60
1:B:22:SER:O	1:B:24:PRO:HD3	2.01	0.60
1:C:205:ILE:HD12	1:C:205:ILE:N	2.16	0.60
1:B:100:ASN:HD22	1:B:102:ASN:H	1.48	0.60
1:C:1:MET:CE	1:C:309:TYR:HD1	2.14	0.60
1:C:100:ASN:ND2	1:C:102:ASN:H	2.00	0.59
1:B:33:TRP:CH2	5:B:4298:HOH:O	2.56	0.59
1:B:224:ASN:ND2	1:B:227:GLU:H	2.01	0.58
1:A:222:VAL:HG21	1:A:228:PHE:CB	2.33	0.58
1:A:348:ASN:ND2	1:C:64:ASN:HD22	2.01	0.58
1:A:226:ASN:HB2	5:A:4308:HOH:O	2.02	0.57
1:C:182:PHE:CD2	1:C:206:GLU:HG3	2.40	0.57
1:C:172:ASN:O	1:C:173:ASP:HB2	2.04	0.57
1:A:291:ILE:HA	1:A:294:ILE:HD12	1.87	0.57
1:B:258:HIS:HA	5:B:4175:HOH:O	2.04	0.57
1:A:202:LEU:HD23	1:A:208:VAL:HG21	1.85	0.56
1:B:262:ILE:HD13	1:B:294:ILE:HD13	1.88	0.56
1:C:130:HIS:CD2	1:C:151:TYR:OH	2.50	0.56
1:C:131:HIS:CD2	5:C:4190:HOH:O	2.58	0.56
1:C:204:SER:C	1:C:205:ILE:HD12	2.26	0.56
1:B:144:LYS:NZ	5:B:4275:HOH:O	2.38	0.56
1:B:224:ASN:ND2	1:B:227:GLU:HB2	2.20	0.56
1:B:22:SER:OG	1:B:23:PRO:HD2	2.06	0.56
1:B:15:ILE:HB	1:B:16:GLU:OE1	2.06	0.56
1:C:70:VAL:HG23	1:C:87:ALA:HB3	1.88	0.56
1:A:15:ILE:HG13	1:A:16:GLU:O	2.06	0.56
1:C:1:MET:CE	1:C:309:TYR:CD1	2.89	0.55
1:C:85:GLN:CA	1:C:85:GLN:HE21	2.20	0.55
1:B:231:GLN:NE2	5:B:4321:HOH:O	2.39	0.55
1:C:1:MET:HE2	1:C:309:TYR:CD1	2.41	0.55
1:C:39:ILE:O	1:C:43:LYS:HG3	2.06	0.55
1:B:274:ASN:HD22	1:B:275:PRO:HD2	1.71	0.55
1:A:100:ASN:C	1:A:100:ASN:ND2	2.57	0.55
1:A:87:ALA:HA	5:C:4008:HOH:O	2.07	0.55
1:B:184:SER:OG	1:B:229:LEU:HD22	2.07	0.55
1:A:17:LYS:NZ	1:A:17:LYS:CB	2.69	0.54
1:B:205:ILE:O	1:B:205:ILE:HG22	2.08	0.54
1:B:224:ASN:ND2	1:B:227:GLU:CB	2.69	0.54
1:C:22:SER:HA	5:C:4272:HOH:O	2.08	0.54
1:A:191:ASN:HB2	5:A:4413:HOH:O	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:TYR:OH	1:A:122:MET:CE	2.56	0.53
1:B:203:ASN:ND2	1:B:207:PRO:HA	2.23	0.53
1:B:183:ALA:HA	1:B:234:PHE:O	2.08	0.53
1:B:164:HIS:CD2	1:B:224:ASN:HD21	2.27	0.53
1:C:82:LEU:HD11	1:C:89:ARG:HD3	1.91	0.53
1:B:100:ASN:ND2	1:B:100:ASN:C	2.55	0.53
1:B:18:MET:HG2	1:B:54:ARG:NH1	2.23	0.52
1:C:82:LEU:O	1:C:85:GLN:HG2	2.10	0.52
1:B:16:GLU:CD	1:B:16:GLU:H	1.94	0.52
1:C:274:ASN:ND2	1:C:276:LYS:H	2.07	0.52
1:C:178:LEU:HD11	1:C:301:LYS:HB2	1.93	0.51
1:C:205:ILE:CD1	1:C:205:ILE:N	2.73	0.51
1:A:100:ASN:HD21	1:A:102:ASN:HB2	1.76	0.51
1:C:19:ALA:N	5:C:4458:HOH:O	2.43	0.51
1:C:276:LYS:O	1:C:307:MET:HG3	2.11	0.51
1:A:140:THR:HG21	1:A:158:HIS:O	2.12	0.50
1:A:164:HIS:HA	1:A:227:GLU:OE1	2.11	0.50
1:B:60:HIS:HD2	1:B:62:ASN:N	1.94	0.50
1:A:206:GLU:OE1	1:A:207:PRO:HD2	2.11	0.50
1:C:23:PRO:HG2	1:C:56:THR:OG1	2.12	0.50
1:A:149:SER:O	1:A:153:LEU:HG	2.11	0.50
1:A:286:ASN:HD22	1:A:286:ASN:H	1.59	0.50
1:A:64:ASN:HD22	1:C:348:ASN:HD22	1.60	0.50
1:A:86:ASN:ND2	1:A:346:HIS:NE2	2.58	0.49
1:B:274:ASN:ND2	1:B:276:LYS:H	2.10	0.49
1:B:233:LYS:HD3	1:B:298:HIS:CD2	2.47	0.49
1:A:89:ARG:HD3	1:A:103:LEU:O	2.13	0.49
1:B:233:LYS:HB3	1:B:304:TYR:CE2	2.48	0.49
1:B:203:ASN:HD22	1:B:208:VAL:N	2.10	0.49
1:C:91:PHE:CD2	1:C:99:PRO:HB3	2.48	0.49
1:B:130:HIS:CD2	1:B:151:TYR:OH	2.52	0.48
1:B:199:TYR:HE2	1:B:208:VAL:HG12	1.78	0.48
1:B:240:ASN:HD22	1:B:240:ASN:H	1.60	0.48
1:C:274:ASN:HD22	1:C:275:PRO:HD2	1.77	0.48
1:C:113:LEU:CD2	1:C:113:LEU:C	2.81	0.48
1:B:167:LEU:O	1:B:170:VAL:HG12	2.14	0.48
1:A:110:PHE:HB2	1:A:245:GLY:O	2.14	0.48
1:A:21:LYS:O	1:A:21:LYS:HG2	2.14	0.48
1:B:138:ASP:OD2	1:B:140:THR:OG1	2.32	0.48
1:B:131:HIS:CD2	5:B:4145:HOH:O	2.67	0.48
1:C:138:ASP:HA	1:C:157:SER:O	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:ASN:OD1	1:B:217:THR:N	2.47	0.47
1:C:31:ALA:HB3	1:C:34:TRP:CD1	2.49	0.47
1:A:286:ASN:HD21	1:A:289:GLU:CG	2.22	0.47
1:C:113:LEU:HD23	1:C:114:ASP:N	2.29	0.47
1:A:233:LYS:HD3	1:A:298:HIS:CE1	2.50	0.47
1:A:18:MET:HG2	1:A:53:GLN:O	2.15	0.47
1:A:113:LEU:HD22	1:A:119:TYR:CG	2.50	0.47
1:C:32:ASN:ND2	5:C:4123:HOH:O	2.47	0.47
1:B:72:GLY:HA3	5:B:4166:HOH:O	2.14	0.47
1:C:79:ARG:O	1:C:82:LEU:HB3	2.15	0.46
1:B:90:VAL:HA	1:B:106:TYR:O	2.15	0.46
1:B:100:ASN:ND2	1:B:102:ASN:N	2.60	0.46
1:A:347:ASP:OD2	5:A:4079:HOH:O	2.20	0.46
1:B:148:ASN:HB2	5:B:4375:HOH:O	2.16	0.46
1:B:1:MET:O	1:B:4:PRO:HD2	2.16	0.46
1:A:100:ASN:ND2	1:A:102:ASN:N	2.61	0.46
1:A:234:PHE:HB3	1:A:262:ILE:CD1	2.46	0.46
1:B:43:LYS:HE3	1:B:59:LEU:HD23	1.98	0.46
1:A:15:ILE:HA	1:A:329:LYS:HZ2	1.80	0.46
1:A:64:ASN:ND2	1:C:348:ASN:HD22	2.14	0.46
1:A:286:ASN:N	1:A:286:ASN:ND2	2.64	0.45
1:A:164:HIS:HB3	1:A:167:LEU:HB3	1.98	0.45
1:A:349:PRO:HB3	5:A:4165:HOH:O	2.16	0.45
1:C:218:LEU:CD2	1:C:222:VAL:HG13	2.46	0.45
1:B:223:LYS:N	1:B:223:LYS:CD	2.68	0.45
1:A:113:LEU:HD22	1:A:119:TYR:CD2	2.51	0.45
1:A:22:SER:N	1:A:23:PRO:HD3	2.31	0.45
1:C:1:MET:HE1	1:C:309:TYR:CD1	2.52	0.45
1:C:14:SER:O	1:C:329:LYS:HE3	2.16	0.45
1:A:3:GLN:HB3	1:A:4:PRO:HD3	1.98	0.45
1:C:99:PRO:HG2	1:C:101:PHE:CE2	2.52	0.45
1:B:23:PRO:HA	1:B:24:PRO:HD2	1.78	0.45
1:B:36:ASP:O	5:B:4206:HOH:O	2.21	0.45
1:C:205:ILE:O	1:C:205:ILE:HG22	2.16	0.45
1:C:39:ILE:CG2	1:C:43:LYS:HE3	2.47	0.45
1:C:205:ILE:O	1:C:205:ILE:CG2	2.65	0.45
5:A:4276:HOH:O	1:C:351:ILE:HD12	2.17	0.45
1:B:212:GLY:O	1:B:216:ASN:HB2	2.17	0.44
1:A:132:LYS:NZ	5:A:4272:HOH:O	2.50	0.44
1:B:205:ILE:O	1:B:206:GLU:HB2	2.18	0.44
1:B:33:TRP:HZ2	5:B:4298:HOH:O	1.92	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ASN:HD22	1:B:102:ASN:N	2.14	0.44
1:A:223:LYS:HG2	5:A:4060:HOH:O	2.17	0.44
1:A:86:ASN:HA	1:A:346:HIS:NE2	2.33	0.44
1:B:86:ASN:CG	1:B:86:ASN:O	2.55	0.44
1:A:3:GLN:HB3	1:A:4:PRO:CD	2.47	0.43
1:C:240:ASN:OD1	1:C:241:THR:HG23	2.18	0.43
1:A:199:TYR:CE2	1:A:208:VAL:HB	2.53	0.43
1:B:199:TYR:CE2	1:B:208:VAL:CG1	3.01	0.43
1:A:3:GLN:N	1:A:4:PRO:HD2	2.33	0.43
1:C:239:GLU:CG	1:C:265:GLY:HA3	2.49	0.43
1:B:113:LEU:HD12	1:B:113:LEU:C	2.38	0.43
1:B:199:TYR:CE2	1:B:208:VAL:HG12	2.54	0.43
1:A:203:ASN:OD1	1:A:207:PRO:HA	2.18	0.43
1:A:205:ILE:HD13	1:A:205:ILE:N	2.33	0.43
1:A:20:SER:C	1:A:22:SER:H	2.22	0.43
1:A:92:TYR:OH	1:A:122:MET:HE1	2.19	0.42
1:B:170:VAL:HG13	1:B:171:VAL:N	2.34	0.42
1:C:100:ASN:HD21	1:C:102:ASN:HB2	1.83	0.42
1:A:301:LYS:HE2	1:A:302:ASN:ND2	2.34	0.42
1:A:305:LEU:HD22	1:A:309:TYR:CE2	2.54	0.42
1:A:201:ALA:O	1:A:205:ILE:HG12	2.19	0.42
1:B:188:SER:HA	1:B:213:SER:HB2	2.01	0.42
1:B:262:ILE:HA	1:B:279:VAL:HB	2.02	0.42
1:A:161:LYS:HD2	5:A:4108:HOH:O	2.19	0.42
1:C:306:ASP:O	1:C:310:GLU:HG2	2.20	0.42
1:A:210:GLY:H	1:A:217:THR:HB	1.85	0.42
1:B:30:VAL:HG13	1:B:34:TRP:CE3	2.55	0.42
1:C:205:ILE:O	1:C:206:GLU:HB2	2.20	0.42
1:B:297:LEU:O	1:B:304:TYR:HB2	2.20	0.42
1:A:39:ILE:O	1:A:43:LYS:HG3	2.19	0.41
1:B:116:ASN:HB2	1:B:117:ASP:H	1.54	0.41
1:B:154:LYS:HE3	5:B:4192:HOH:O	2.20	0.41
1:B:274:ASN:HD22	1:B:275:PRO:CD	2.33	0.41
1:C:77:SER:HA	1:C:80:LYS:HD2	2.02	0.41
1:A:344:ILE:HG12	1:C:66:PHE:CZ	2.55	0.41
1:B:161:LYS:HA	1:B:168:CYS:SG	2.59	0.41
1:B:74:PRO:HD2	1:B:91:PHE:HE1	1.85	0.41
1:B:222:VAL:HG11	1:B:228:PHE:CB	2.44	0.41
1:B:100:ASN:ND2	1:B:103:LEU:H	2.17	0.41
1:C:1:MET:O	1:C:4:PRO:HD2	2.20	0.41
1:A:240:ASN:OD1	1:A:241:THR:HG23	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:ASN:ND2	1:A:289:GLU:H	2.18	0.41
1:C:142:PRO:HD3	1:C:168:CYS:SG	2.60	0.41
1:B:195:ARG:HG3	1:B:238:PHE:HB2	2.02	0.41
1:B:85:GLN:NE2	5:B:4096:HOH:O	2.51	0.41
1:A:140:THR:HG23	1:A:158:HIS:HA	2.03	0.41
1:A:204:SER:O	1:A:205:ILE:HD13	2.20	0.41
1:C:199:TYR:CD1	1:C:199:TYR:C	2.94	0.41
1:A:351:ILE:HA	1:A:351:ILE:HD13	1.88	0.41
1:C:240:ASN:ND2	2:C:4003:FUC:O2	2.53	0.41
1:C:85:GLN:CA	1:C:85:GLN:NE2	2.84	0.41
1:A:86:ASN:HD22	1:A:86:ASN:HA	1.69	0.41
1:A:96:ASN:HB2	1:A:246:TYR:CZ	2.55	0.40
1:C:17:LYS:HG2	1:C:18:MET:N	2.36	0.40
1:C:91:PHE:HB2	1:C:104:PHE:CG	2.56	0.40
1:B:225:LYS:HE2	1:B:229:LEU:HD11	2.03	0.40
1:A:120:LEU:HD21	1:A:331:ILE:HG12	2.01	0.40
1:A:349:PRO:O	1:A:350:PHE:CD2	2.74	0.40
1:B:113:LEU:HD12	1:B:114:ASP:N	2.37	0.40
1:C:274:ASN:HD22	1:C:275:PRO:CD	2.34	0.40
1:A:39:ILE:CG2	1:A:43:LYS:HE3	2.51	0.40
1:C:117:ASP:HB3	5:C:4178:HOH:O	2.21	0.40
1:B:168:CYS:O	1:B:172:ASN:HB2	2.22	0.40
1:B:87:ALA:HB2	5:B:4291:HOH:O	2.21	0.40
1:C:184:SER:OG	1:C:229:LEU:HD22	2.22	0.40
1:C:3:GLN:N	1:C:4:PRO:CD	2.84	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	339/371 (91%)	318 (94%)	19 (6%)	2 (1%)	25 15

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	334/371 (90%)	323 (97%)	10 (3%)	1 (0%)	41	31
1	C	349/371 (94%)	337 (97%)	11 (3%)	1 (0%)	41	31
All	All	1022/1113 (92%)	978 (96%)	40 (4%)	4 (0%)	34	24

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	19	ALA
1	A	21	LYS
1	A	206	GLU
1	B	206	GLU

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	308/334 (92%)	290 (94%)	18 (6%)	20	11
1	B	303/334 (91%)	289 (95%)	14 (5%)	27	19
1	C	314/334 (94%)	297 (95%)	17 (5%)	22	13
All	All	925/1002 (92%)	876 (95%)	49 (5%)	22	14

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	17	LYS
1	A	18	MET
1	A	44	ASN
1	A	86	ASN
1	A	100	ASN
1	A	113	LEU
1	A	115	PHE
1	A	140	THR
1	A	162	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	170	VAL
1	A	175	SER
1	A	206	GLU
1	A	237	CYS
1	A	240	ASN
1	A	286	ASN
1	A	315	THR
1	A	349	PRO
1	B	16	GLU
1	B	38	GLU
1	B	43	LYS
1	B	85	GLN
1	B	100	ASN
1	B	113	LEU
1	B	115	PHE
1	B	116	ASN
1	B	157	SER
1	B	190	PRO
1	B	207	PRO
1	B	223	LYS
1	B	274	ASN
1	B	292	ASP
1	C	16	GLU
1	C	18	MET
1	C	21	LYS
1	C	27	LYS
1	C	32	ASN
1	C	36	ASP
1	C	85	GLN
1	C	100	ASN
1	C	103	LEU
1	C	113	LEU
1	C	115	PHE
1	C	116	ASN
1	C	147	ASP
1	C	237	CYS
1	C	274	ASN
1	C	301	LYS
1	C	316	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	64	ASN
1	A	86	ASN
1	A	100	ASN
1	A	131	HIS
1	A	172	ASN
1	A	196	ASN
1	A	274	ASN
1	A	286	ASN
1	A	300	HIS
1	A	302	ASN
1	B	60	HIS
1	B	61	GLN
1	B	73	ASN
1	B	100	ASN
1	B	130	HIS
1	B	131	HIS
1	B	172	ASN
1	B	196	ASN
1	B	203	ASN
1	B	224	ASN
1	B	231	GLN
1	B	240	ASN
1	B	274	ASN
1	B	298	HIS
1	B	300	HIS
1	C	60	HIS
1	C	64	ASN
1	C	85	GLN
1	C	86	ASN
1	C	100	ASN
1	C	130	HIS
1	C	166	ASN
1	C	172	ASN
1	C	240	ASN
1	C	274	ASN
1	C	300	HIS
1	C	302	ASN
1	C	314	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GDP	A	3001	2	24,30,30	1.60	6 (25%)	31,47,47	2.22	9 (29%)
2	FUC	C	4003	3	10,10,11	1.34	2 (20%)	14,14,16	1.28	2 (14%)
4	SO4	C	2002	-	4,4,4	0.33	0	6,6,6	0.24	0
3	GDP	B	3002	2	24,30,30	1.48	3 (12%)	31,47,47	2.16	7 (22%)
2	FUC	B	4002	3	10,10,11	0.99	0	14,14,16	1.30	2 (14%)
4	SO4	B	2001	-	4,4,4	0.38	0	6,6,6	0.12	0
3	GDP	C	3003	2	24,30,30	1.47	3 (12%)	31,47,47	2.08	8 (25%)
2	FUC	A	4001	3	10,10,11	1.03	1 (10%)	14,14,16	0.96	1 (7%)

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	A	3001	2	-	1/12/32/32	0/3/3/3
2	FUC	C	4003	3	1/1/4/5	-	0/1/1/1
3	GDP	B	3002	2	-	0/12/32/32	0/3/3/3
2	FUC	B	4002	3	1/1/4/5	-	0/1/1/1
3	GDP	C	3003	2	-	1/12/32/32	0/3/3/3
2	FUC	A	4001	3	1/1/4/5	-	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3002	GDP	O4'-C1'	4.12	1.46	1.41
3	C	3003	GDP	C6-N1	3.91	1.39	1.33
3	A	3001	GDP	C6-N1	3.83	1.39	1.33
3	C	3003	GDP	O4'-C1'	3.73	1.46	1.41
3	C	3003	GDP	C2-N1	3.21	1.41	1.35
3	A	3001	GDP	C2-N1	2.94	1.40	1.35
3	A	3001	GDP	C8-N7	-2.91	1.29	1.34
2	A	4001	FUC	C1-C2	2.76	1.58	1.52
2	C	4003	FUC	C1-C2	2.62	1.58	1.52
3	A	3001	GDP	C2-N2	2.61	1.39	1.33
3	A	3001	GDP	O4'-C1'	2.57	1.44	1.41
3	B	3002	GDP	C6-N1	2.48	1.37	1.33
3	A	3001	GDP	C6-C5	2.21	1.45	1.41
3	B	3002	GDP	C6-C5	2.19	1.45	1.41
2	C	4003	FUC	C4-C3	2.17	1.57	1.52

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	3001	GDP	C2-N3-C4	5.75	121.92	115.36
3	B	3002	GDP	C2-N3-C4	5.69	121.85	115.36
3	C	3003	GDP	C2-N3-C4	5.44	121.57	115.36
3	A	3001	GDP	PA-O3A-PB	-5.30	114.63	132.83
3	C	3003	GDP	N3-C2-N1	-5.09	120.43	127.22
3	A	3001	GDP	N3-C2-N1	-4.99	120.57	127.22
3	B	3002	GDP	PA-O3A-PB	-4.90	116.02	132.83
3	C	3003	GDP	PA-O3A-PB	-4.87	116.11	132.83
3	B	3002	GDP	N3-C2-N1	-4.76	120.87	127.22
3	A	3001	GDP	C5-C6-N1	-3.74	118.32	123.43
3	B	3002	GDP	C5-C6-N1	-3.62	118.48	123.43
3	C	3003	GDP	C5-C6-N1	-3.42	118.75	123.43
3	B	3002	GDP	C6-C5-C4	-3.42	117.53	120.80
3	B	3002	GDP	O3B-PB-O3A	3.36	115.91	104.64
3	B	3002	GDP	C6-N1-C2	3.31	121.19	115.93
3	A	3001	GDP	O3B-PB-O3A	3.22	115.42	104.64
3	A	3001	GDP	C6-N1-C2	2.98	120.66	115.93
3	A	3001	GDP	C6-C5-C4	-2.93	118.00	120.80
3	C	3003	GDP	C6-N1-C2	2.86	120.47	115.93
3	C	3003	GDP	O3B-PB-O3A	2.77	113.91	104.64
2	C	4003	FUC	C2-C3-C4	-2.62	106.35	110.89
2	A	4001	FUC	C2-C3-C4	-2.52	106.53	110.89
2	B	4002	FUC	C1-C2-C3	2.46	112.69	109.67

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	3003	GDP	C6-C5-C4	-2.46	118.45	120.80
2	B	4002	FUC	O2-C2-C3	-2.42	105.29	110.14
2	C	4003	FUC	O2-C2-C3	-2.16	105.80	110.14
3	A	3001	GDP	O2B-PB-O3A	2.08	111.61	104.64
3	A	3001	GDP	PA-O5'-C5'	-2.05	109.67	121.68
3	C	3003	GDP	O2B-PB-O3A	2.03	111.45	104.64

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	4003	FUC	C1
2	B	4002	FUC	C1
2	A	4001	FUC	C1

All (2) torsion outliers are listed below:

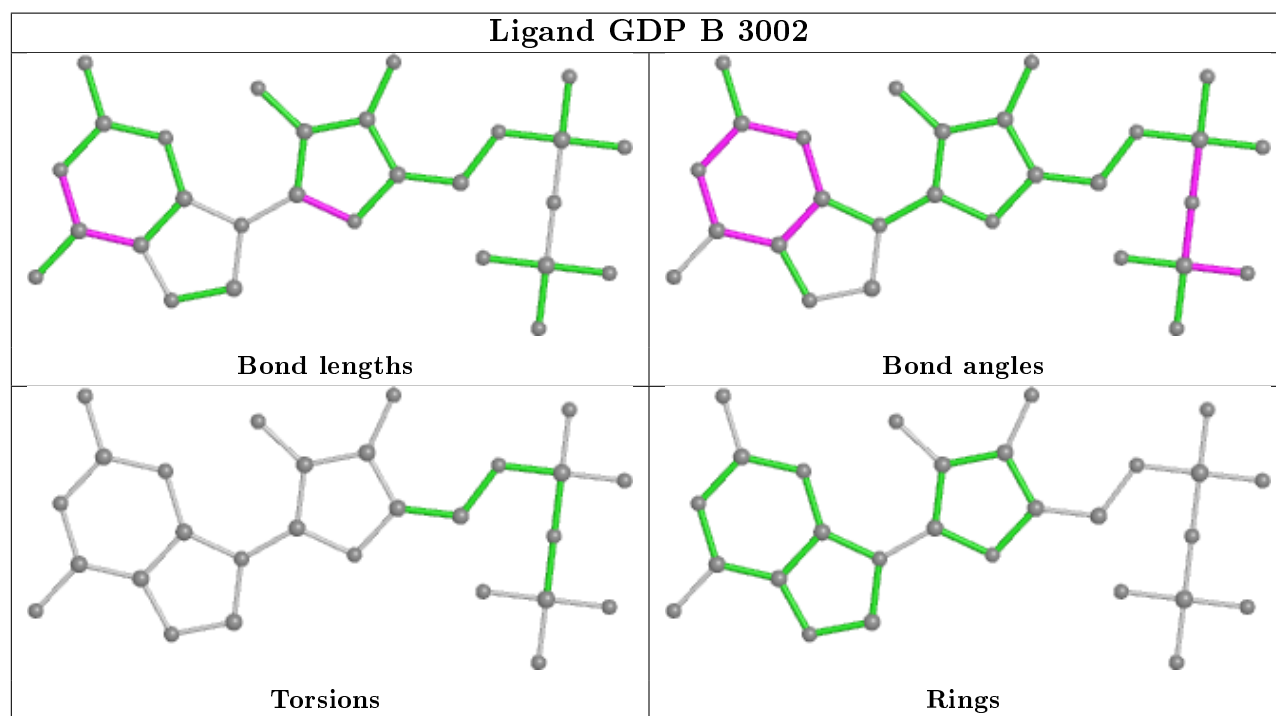
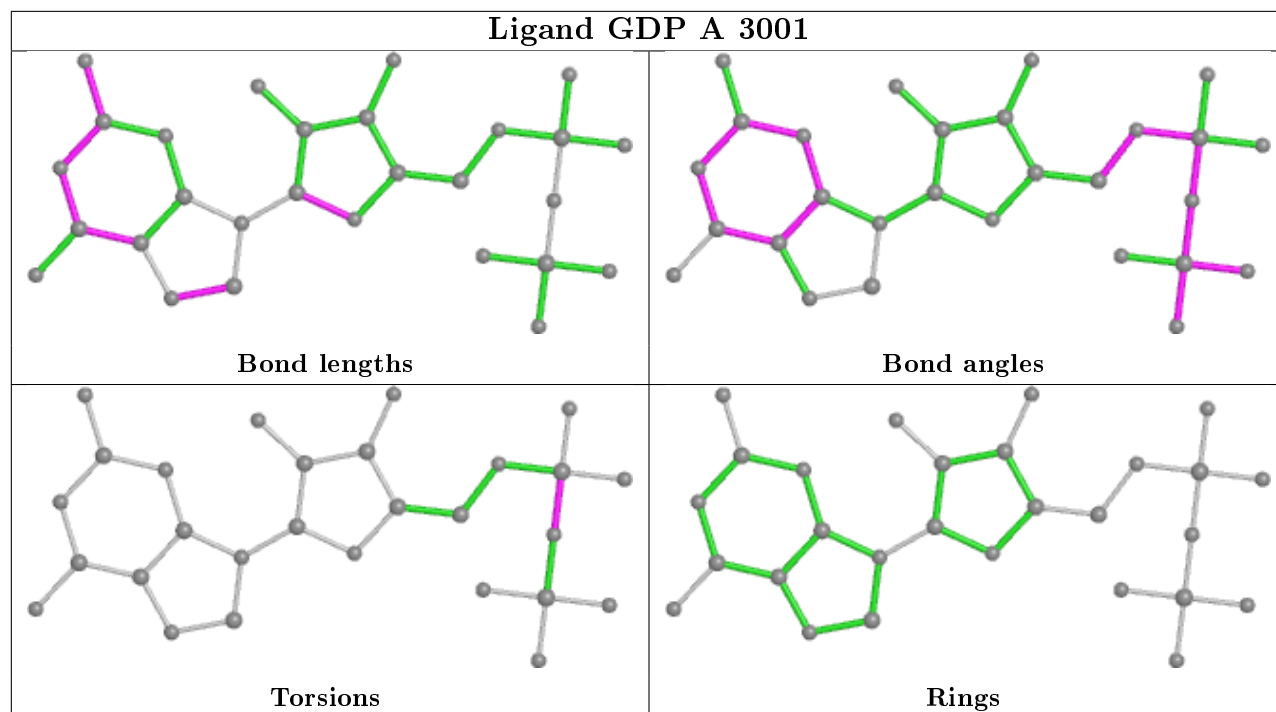
Mol	Chain	Res	Type	Atoms
3	A	3001	GDP	PB-O3A-PA-O2A
3	C	3003	GDP	PB-O3A-PA-O2A

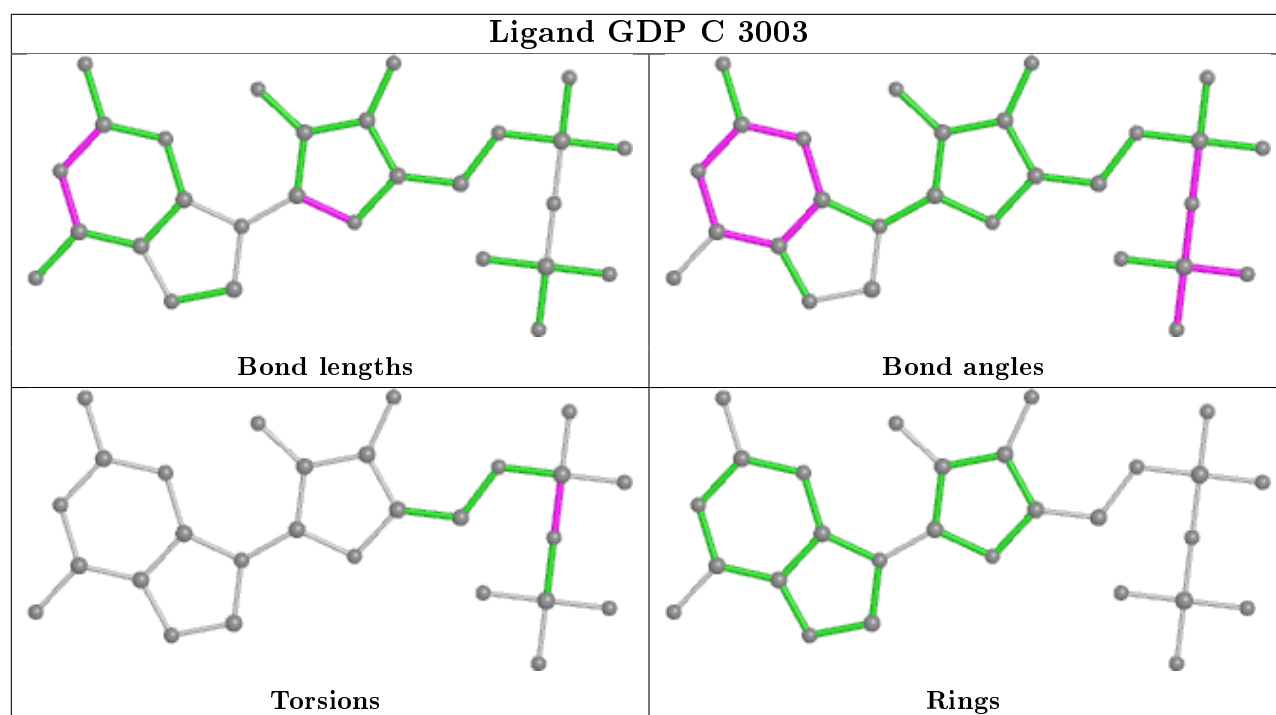
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4003	FUC	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	343/371 (92%)	0.24	25 (7%)	15 16	23, 44, 68, 87	0
1	B	338/371 (91%)	0.16	22 (6%)	18 20	21, 38, 66, 76	0
1	C	351/371 (94%)	-0.01	18 (5%)	28 30	22, 31, 57, 83	0
All	All	1032/1113 (92%)	0.13	65 (6%)	20 21	21, 37, 66, 87	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	19	ALA	6.5
1	A	19	ALA	6.2
1	A	20	SER	5.9
1	B	20	SER	5.1
1	C	20	SER	4.6
1	C	19	ALA	4.4
1	C	17	LYS	4.2
1	A	351	ILE	4.1
1	A	18	MET	4.0
1	A	22	SER	3.9
1	A	352	PHE	3.8
1	C	77	SER	3.7
1	A	33	TRP	3.5
1	C	75	LEU	3.5
1	A	21	LYS	3.3
1	B	18	MET	3.2
1	A	287	PHE	3.2
1	A	223	LYS	3.2
1	A	316	LEU	3.1
1	B	73	ASN	3.1
1	B	158	HIS	3.1
1	B	301	LYS	3.0
1	C	21	LYS	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	259	THR	3.0
1	C	148	ASN	2.9
1	C	350	PHE	2.8
1	A	350	PHE	2.7
1	C	351	ILE	2.7
1	A	222	VAL	2.6
1	A	236	LEU	2.6
1	A	285	LYS	2.5
1	C	204	SER	2.5
1	A	252	ILE	2.5
1	A	205	ILE	2.5
1	B	22	SER	2.4
1	C	22	SER	2.4
1	B	64	ASN	2.4
1	A	148	ASN	2.3
1	A	37	GLU	2.3
1	A	237	CYS	2.3
1	B	251	ILE	2.3
1	A	220	TYR	2.3
1	B	32	ASN	2.3
1	B	205	ILE	2.3
1	B	33	TRP	2.3
1	B	237	CYS	2.3
1	C	37	GLU	2.3
1	A	317	ASP	2.3
1	B	236	LEU	2.3
1	C	84	TYR	2.2
1	A	295	LYS	2.2
1	B	235	ASN	2.2
1	A	349	PRO	2.2
1	C	36	ASP	2.2
1	B	254	ALA	2.2
1	C	261	PRO	2.2
1	C	317	ASP	2.1
1	B	184	SER	2.1
1	C	237	CYS	2.1
1	A	219	GLY	2.1
1	B	148	ASN	2.1
1	B	85	GLN	2.0
1	B	220	TYR	2.0
1	B	139	THR	2.0
1	C	259	THR	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 monosaccharides in this entry.

6.4 Ligands [i](#)

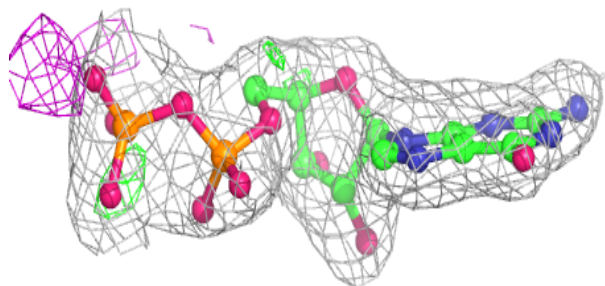
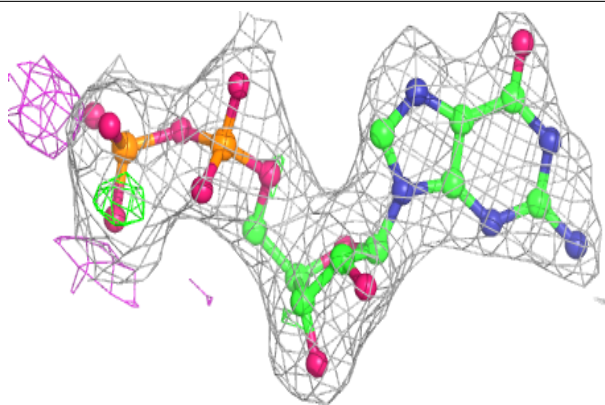
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FUC	A	4001	10/11	0.85	0.37	61,61,62,64	0
3	GDP	A	3001	28/28	0.92	0.14	42,49,57,58	0
2	FUC	C	4003	10/11	0.93	0.18	33,36,38,39	0
4	SO4	B	2001	5/5	0.94	0.37	88,88,88,89	0
2	FUC	B	4002	10/11	0.94	0.15	34,36,38,40	0
3	GDP	B	3002	28/28	0.96	0.10	35,44,49,51	0
4	SO4	C	2002	5/5	0.97	0.17	67,67,68,69	0
3	GDP	C	3003	28/28	0.98	0.07	24,27,36,38	0

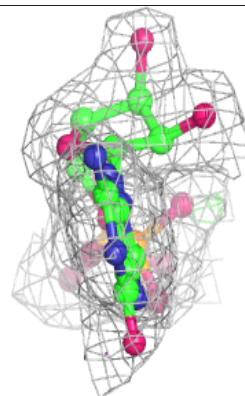
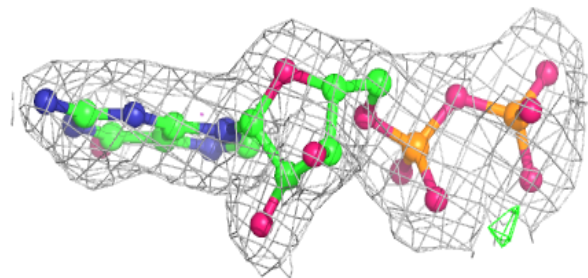
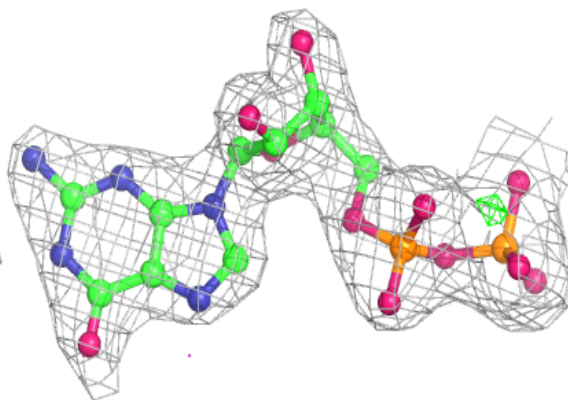
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

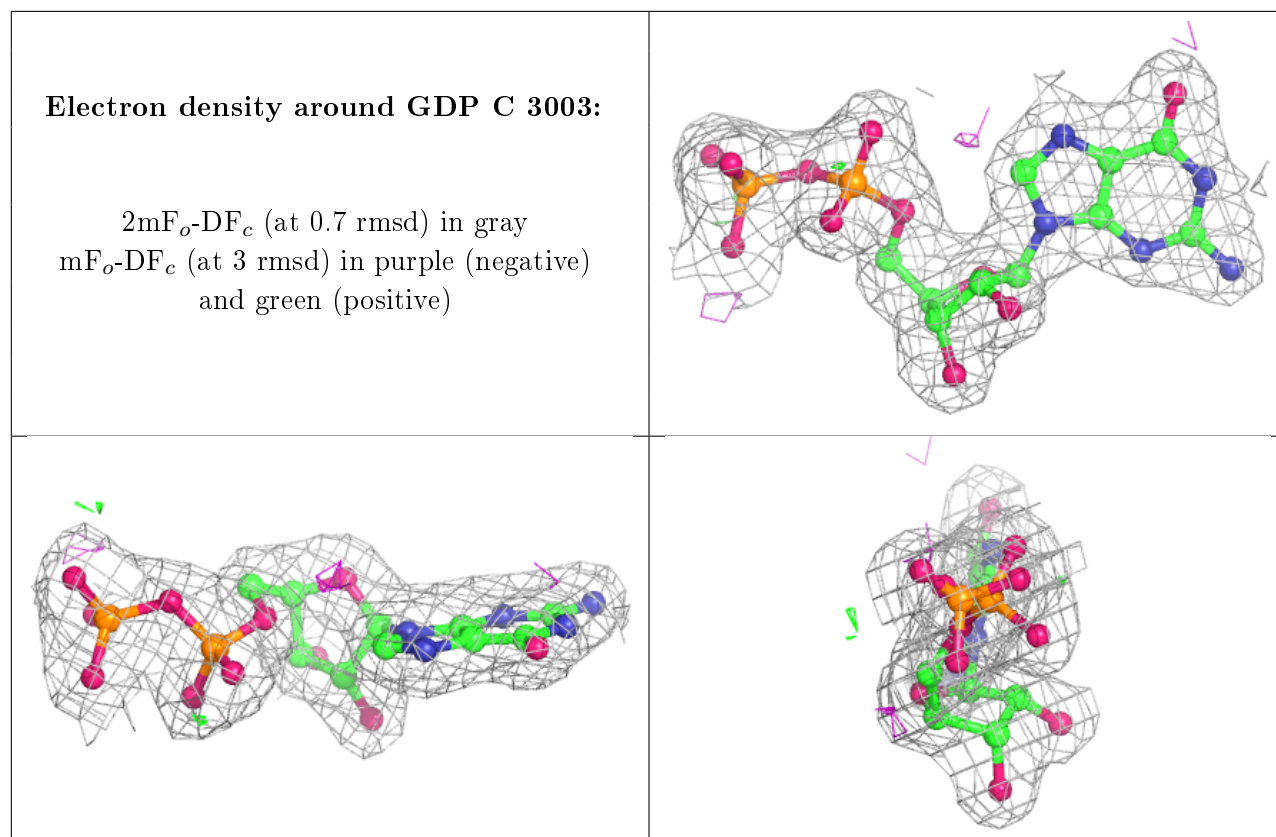
Electron density around GDP A 3001:

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

**Electron density around GDP B 3002:**

$2mF_o-DF_c$ (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.