



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:20 AM UTC

PDB ID : 3U6K / pdb_00003u6k
Title : Ef-tu (escherichia coli) in complex with nvp-ldk733
Authors : Palestrant, D.J.
Deposited on : 2011-10-12
Resolution : 2.45 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:

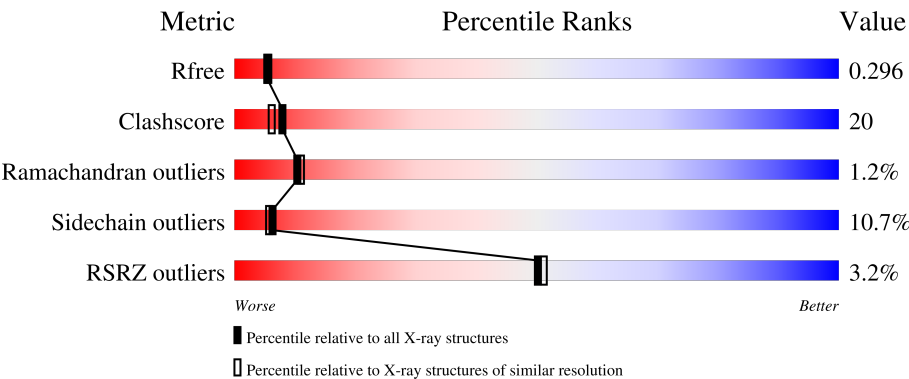
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	394	3% 62% 31% • • •
1	B	394	4% 55% 37% 5% •
2	C	12	17% 75% 8%
2	D	12	17% 75% 8%

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	MH6	C	11	-	X	-	-
2	BB9	C	9	-	X	-	-
2	MH6	D	11	-	X	-	-
2	BB9	D	9	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6311 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 Elongation factor Tu 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	386	Total	C	N	O	S	0	0	0
			2969	1877	511	568	13			
1	B	386	Total	C	N	O	S	0	0	0
			2969	1877	511	568	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP P0CE47
A	1	ALA	-	expression tag	UNP P0CE47
B	0	MET	-	expression tag	UNP P0CE47
B	1	ALA	-	expression tag	UNP P0CE47

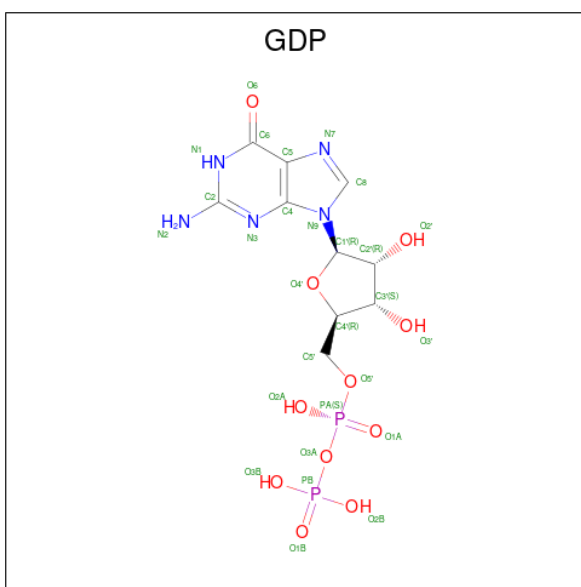
- Molecule 2 is a protein called Thiocillin GE2270 analogue NVP-LDK733.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	S	0	0	0
			85	55	13	11	6			
2	D	12	Total	C	N	O	S	0	0	0
			85	55	13	11	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	12	9BB	CYS	SEE REMARK 999	UNP Q7M0J8
D	12	9BB	CYS	SEE REMARK 999	UNP Q7M0J8

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		

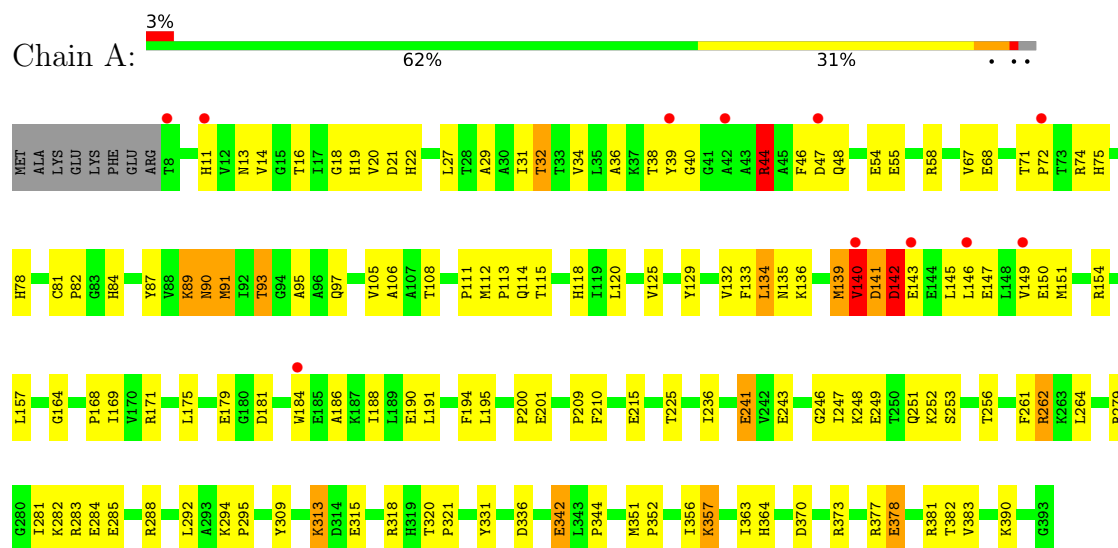
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	69	Total	O	0	0
			69	69		
5	B	72	Total	O	0	0
			72	72		
5	C	1	Total	O	0	0
			1	1		
5	D	3	Total	O	0	0
			3	3		

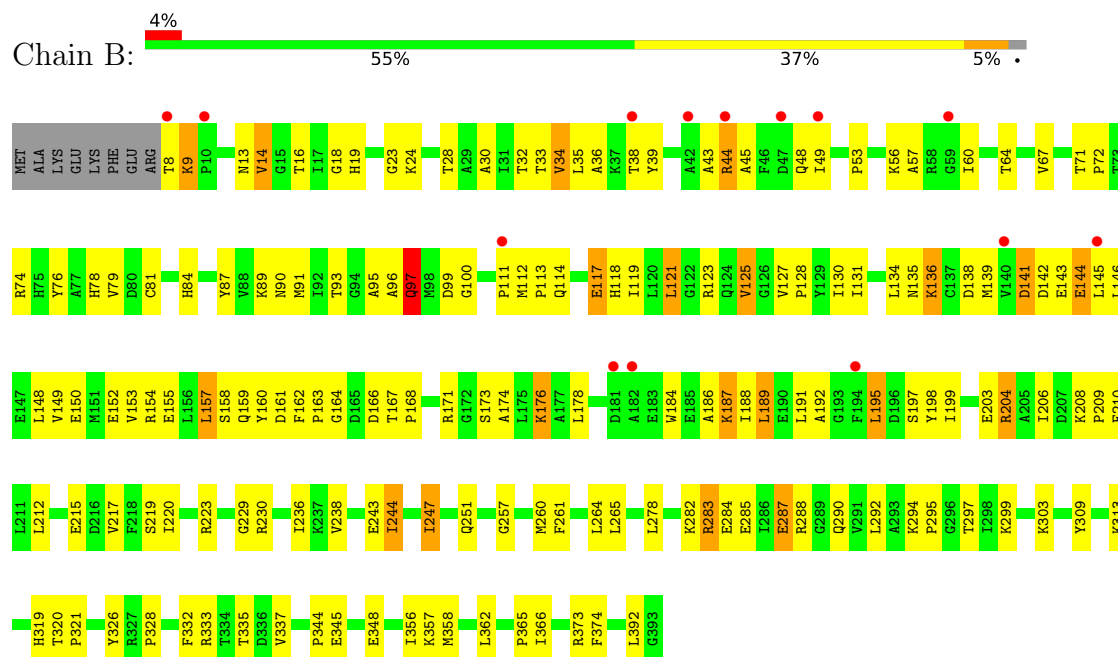
3 Residue-property plots [i](#)

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.

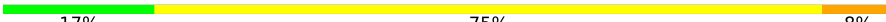
• Molecule 1: Elongation factor Tu 1

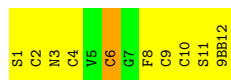


• Molecule 1: Elongation factor Tu 1

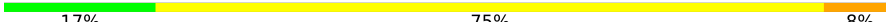


- Molecule 2: Thiocillin GE2270 analogue NVP-LDK733

Chain C:  17% 75% 8%



- Molecule 2: Thiocillin GE2270 analogue NVP-LDK733

Chain D:  17% 75% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.37Å 82.22Å 129.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.04 – 2.45 43.04 – 2.45	Depositor EDS
% Data completeness (in resolution range)	(Not available) (43.04-2.45) 94.9 (43.04-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 2.45Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.224 , 0.298 0.222 , 0.296	Depositor DCC
R_{free} test set	1590 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtriage
Anisotropy	0.228	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6311	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.10% 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: BB8, BB6, MH6, MG, BB7, 9BB, MEN, BB9, 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	0.64	0/3024	0.94	6/4096 (0.1%)
1	B	0.81	6/3024 (0.2%)	0.89	2/4096 (0.0%)
2	C	1.87	1/10 (10.0%)	2.08	0/9
2	D	2.02	1/10 (10.0%)	2.06	0/9
All	All	0.73	8/6068 (0.1%)	0.92	8/8210 (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	2
2	C	1	0
2	D	1	0
All	All	2	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	230	ARG	C-O	-21.92	0.97	1.24
1	B	230	ARG	CA-C	-14.98	1.34	1.52
1	B	230	ARG	N-CA	-6.69	1.37	1.46
1	B	230	ARG	CZ-NH1	-6.45	1.23	1.32
1	B	229	GLY	C-N	-5.58	1.25	1.33
2	C	1	SER	CA-CB	-5.21	1.43	1.53
1	B	230	ARG	CZ-NH2	-5.16	1.26	1.33
2	D	1	SER	CA-CB	-5.03	1.43	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ARG	N-CA-C	9.60	123.67	109.07
1	A	140	VAL	N-CA-C	9.04	120.82	108.17
1	A	141	ASP	N-CA-C	-6.44	96.75	108.02
1	A	139	MET	CA-C-N	-5.64	116.20	123.19
1	A	139	MET	C-N-CA	-5.64	116.20	123.19
1	B	230	ARG	CA-C-O	-5.58	114.23	120.54
1	B	335	THR	N-CA-C	5.51	117.35	109.14
1	A	139	MET	N-CA-C	-5.39	106.55	113.23

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	SER	CA
2	D	1	SER	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	140	VAL	Peptide
1	A	142	ASP	Peptide

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	2969	0	2981	112	0
1	B	2969	0	2982	128	0
2	C	85	0	54	1	0
2	D	85	0	54	1	0
3	A	28	0	12	1	0
3	B	28	0	12	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	69	0	0	3	0
5	B	72	0	0	5	0
5	C	1	0	0	0	0
5	D	3	0	0	0	0
All	All	6311	0	6095	241	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 20.

All (241) 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:A:140:VAL:HG11	1:A:146:LEU:HD21	1.44	0.97
1:B:18:GLY:H	1:B:118:HIS:HD2	1.22	0.87
1:B:282:LYS:HD2	1:B:283:ARG:H	1.40	0.87
1:A:93:THR:HG22	1:A:95:ALA:H	1.44	0.82
1:B:131:ILE:HD11	1:B:198:TYR:CD2	2.13	0.82
1:A:18:GLY:H	1:A:118:HIS:HD2	1.28	0.79
1:B:30:ALA:HA	1:B:178:LEU:HD23	1.67	0.77
1:B:294:LYS:O	1:B:297:THR:HG22	1.87	0.75
1:B:178:LEU:HD13	1:B:178:LEU:O	1.86	0.75
1:B:348:GLU:CD	1:B:348:GLU:H	1.96	0.74
1:A:44:ARG:HD3	1:A:67:VAL:HG22	1.70	0.74
1:A:18:GLY:H	1:A:118:HIS:CD2	2.06	0.73
1:B:71:THR:HB	1:B:72:PRO:HD2	1.72	0.71
1:B:189:LEU:HD22	1:B:189:LEU:H	1.55	0.71
1:A:140:VAL:CG1	1:A:146:LEU:HD21	2.19	0.70
1:B:34:VAL:HG21	1:B:188:ILE:HG21	1.73	0.70
1:A:318:ARG:HH12	1:A:378:GLU:CD	2.00	0.70
1:B:91:MET:O	1:B:125:VAL:HG21	1.93	0.69
1:B:134:LEU:HD11	1:B:149:VAL:HG11	1.75	0.68
1:A:19:HIS:CD2	1:A:20:VAL:H	2.14	0.66
1:A:150:GLU:O	1:A:154:ARG:HG3	1.96	0.66
1:B:155:GLU:HG2	5:B:408:HOH:O	1.95	0.66
1:A:22:HIS:CD2	1:A:106:ALA:H	2.14	0.66
1:A:34:VAL:HG21	1:A:188:ILE:HG21	1.78	0.66
1:B:16:THR:HG23	1:B:78:HIS:CE1	2.30	0.66
1:B:18:GLY:H	1:B:118:HIS:CD2	2.10	0.65
1:A:32:THR:HG23	1:A:44:ARG:H	1.60	0.65
1:A:32:THR:CG2	1:A:44:ARG:H	2.09	0.65
1:A:91:MET:O	1:A:125:VAL:HG21	1.96	0.65
1:A:282:LYS:O	1:A:285:GLU:HG2	1.96	0.64
1:B:189:LEU:HD22	1:B:189:LEU:N	2.12	0.64
1:B:158:SER:HA	1:B:162:PHE:O	1.98	0.64
1:A:19:HIS:CD2	1:A:20:VAL:HG12	2.32	0.63
1:A:89:LYS:O	1:A:93:THR:HB	1.99	0.63
1:B:81:CYS:HB2	1:B:87:TYR:CE2	2.34	0.63
1:B:89:LYS:O	1:B:93:THR:HG23	1.99	0.62
1:A:54:GLU:CD	1:A:54:GLU:H	2.07	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:THR:HB	1:A:72:PRO:CD	2.29	0.62
1:A:112:MET:HE2	1:A:112:MET:HA	1.81	0.62
1:A:16:THR:HG23	1:A:78:HIS:CE1	2.34	0.62
1:B:30:ALA:CA	1:B:178:LEU:HD23	2.29	0.62
1:A:93:THR:HG22	1:A:95:ALA:N	2.14	0.61
1:B:91:MET:HE3	1:B:121:LEU:HB3	1.81	0.61
1:B:44:ARG:NH2	1:B:67:VAL:HA	2.15	0.61
1:A:14:VAL:O	1:A:78:HIS:HA	2.01	0.61
1:B:57:ALA:HB3	1:B:60:ILE:HB	1.81	0.60
1:A:150:GLU:OE2	1:A:171:ARG:HD2	2.01	0.60
1:B:264:LEU:HD21	5:B:399:HOH:O	2.01	0.60
1:A:168:PRO:HB3	1:A:194:PHE:CD1	2.35	0.60
1:B:99:ASP:O	1:B:128:PRO:HG2	2.02	0.60
1:A:84:HIS:CD2	1:A:118:HIS:HE1	2.18	0.60
1:A:146:LEU:O	1:A:150:GLU:HG3	2.02	0.60
1:A:32:THR:HG21	1:A:44:ARG:HG2	1.84	0.59
1:A:44:ARG:HB2	1:A:44:ARG:CZ	2.33	0.59
1:B:96:ALA:O	1:B:97:GLN:HB2	2.01	0.59
1:B:192:ALA:O	1:B:195:LEU:HB2	2.03	0.59
1:B:148:LEU:O	1:B:152:GLU:HG3	2.04	0.58
1:A:225:THR:HG23	1:A:281:ILE:O	2.04	0.58
1:B:173:SER:OG	1:B:176:LYS:HB2	2.04	0.58
1:A:90:ASN:HD22	1:A:95:ALA:HB3	1.69	0.57
1:A:283:ARG:HD3	1:A:283:ARG:C	2.29	0.57
1:B:32:THR:O	1:B:36:ALA:HB2	2.05	0.57
1:A:19:HIS:CD2	1:A:20:VAL:N	2.72	0.57
1:A:34:VAL:CG2	1:A:188:ILE:HG21	2.35	0.57
1:B:64:THR:HG23	1:B:79:VAL:HG23	1.85	0.57
1:A:36:ALA:O	1:A:40:GLY:HA2	2.05	0.57
1:B:8:THR:O	1:B:8:THR:HG23	2.04	0.57
1:B:243:GLU:HG3	1:B:295:PRO:HB3	1.85	0.57
1:B:167:THR:HG22	1:B:168:PRO:O	2.04	0.57
1:A:89:LYS:HD3	1:A:309:TYR:CE2	2.39	0.56
1:A:294:LYS:HG2	1:A:295:PRO:HD2	1.88	0.56
1:B:53:PRO:HD2	1:B:64:THR:O	2.06	0.56
1:B:244:ILE:HD11	1:B:278:LEU:HD13	1.86	0.56
1:B:189:LEU:HA	1:B:192:ALA:HB3	1.88	0.56
1:B:174:ALA:O	1:B:178:LEU:HB2	2.06	0.56
1:A:132:VAL:HB	1:A:169:ILE:HG12	1.89	0.55
1:B:38:THR:HG21	1:B:189:LEU:HD12	1.89	0.55
1:B:89:LYS:HE2	1:B:309:TYR:CZ	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:TYR:O	1:B:161:ASP:HB2	2.06	0.55
1:A:84:HIS:HD2	1:A:118:HIS:HE1	1.54	0.54
1:A:141:ASP:O	1:A:142:ASP:HB2	2.07	0.54
1:A:29:ALA:O	1:A:32:THR:HG22	2.07	0.54
1:A:19:HIS:HA	1:A:114:GLN:HB2	1.90	0.54
1:B:204:ARG:HB2	1:B:206:ILE:HG22	1.90	0.53
1:A:133:PHE:CD1	1:A:191:LEU:HD22	2.44	0.53
1:A:71:THR:HB	1:A:72:PRO:HD3	1.90	0.53
1:A:111:PRO:O	1:A:112:MET:HE2	2.09	0.53
1:A:320:THR:HB	1:A:321:PRO:HD2	1.90	0.53
1:B:35:LEU:HD22	1:B:39:TYR:CE2	2.44	0.52
1:A:225:THR:HG21	1:A:283:ARG:N	2.25	0.52
1:B:260:MET:O	1:B:261:PHE:C	2.53	0.52
1:A:154:ARG:HD3	1:A:164:GLY:O	2.09	0.52
1:A:209:PRO:HB3	1:A:294:LYS:HE2	1.91	0.52
1:B:45:ALA:O	1:B:49:ILE:HG12	2.09	0.52
1:B:189:LEU:HA	1:B:192:ALA:CB	2.40	0.52
1:B:251:GLN:OE1	1:B:285:GLU:HB3	2.10	0.52
1:B:319:HIS:HE1	5:B:449:HOH:O	1.91	0.52
1:B:119:ILE:HD13	1:B:157:LEU:HD13	1.92	0.51
1:A:89:LYS:HD3	1:A:309:TYR:CZ	2.46	0.51
1:B:163:PRO:HB2	1:B:166:ASP:HB2	1.91	0.51
1:B:303:LYS:HG2	1:B:392:LEU:HB2	1.92	0.51
1:A:210:PHE:CE1	1:A:236:ILE:HB	2.46	0.51
1:A:68:GLU:CD	1:A:75:HIS:HE2	2.20	0.51
1:B:35:LEU:HD22	1:B:39:TYR:HE2	1.76	0.51
1:A:84:HIS:HD2	1:A:118:HIS:CE1	2.29	0.50
1:A:168:PRO:HB3	1:A:194:PHE:CE1	2.47	0.50
1:B:34:VAL:HG21	1:B:188:ILE:HG13	1.92	0.50
1:B:138:ASP:O	1:B:139:MET:HE2	2.10	0.50
1:B:320:THR:HB	1:B:321:PRO:HD2	1.93	0.50
1:A:370:ASP:OD1	1:A:390:LYS:HA	2.12	0.50
1:B:145:LEU:O	1:B:148:LEU:HB2	2.12	0.50
1:B:178:LEU:HD13	1:B:178:LEU:C	2.37	0.50
1:B:356:ILE:HD11	1:B:358:MET:HE3	1.94	0.50
1:B:14:VAL:HG12	1:B:100:GLY:O	2.12	0.50
1:B:264:LEU:N	1:B:264:LEU:HD22	2.27	0.50
1:A:19:HIS:CE1	1:A:113:PRO:HD2	2.47	0.49
1:B:28:THR:HG23	1:B:78:HIS:CD2	2.47	0.49
1:B:74:ARG:NH2	1:B:199:ILE:O	2.44	0.49
1:B:294:LYS:O	1:B:297:THR:CG2	2.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:TRP:C	1:A:186:ALA:N	2.70	0.49
1:A:44:ARG:CZ	1:A:44:ARG:CB	2.85	0.49
1:B:84:HIS:CD2	1:B:118:HIS:HE1	2.31	0.49
1:B:149:VAL:HG12	1:B:153:VAL:HG23	1.95	0.49
1:A:344:PRO:HD2	1:A:356:ILE:HD11	1.94	0.49
1:B:257:GLY:HA3	2:D:12:9BB:O	2.13	0.49
1:B:282:LYS:CD	1:B:283:ARG:H	2.21	0.49
1:A:21:ASP:OD1	3:A:500:GDP:H5''	2.13	0.49
1:A:246:GLY:C	1:A:248:LYS:H	2.21	0.49
1:A:209:PRO:HB3	1:A:294:LYS:CE	2.43	0.48
1:A:215:GLU:OE1	1:A:288:ARG:NH1	2.46	0.48
1:B:141:ASP:N	1:B:141:ASP:OD1	2.44	0.48
1:A:377:ARG:NH1	1:A:382:THR:OG1	2.47	0.48
1:B:130:ILE:O	1:B:131:ILE:HD13	2.14	0.47
1:B:13:ASN:ND2	5:B:405:HOH:O	2.41	0.47
1:B:76:TYR:CZ	1:B:195:LEU:HB3	2.49	0.47
1:A:11:HIS:HE1	1:A:13:ASN:OD1	1.98	0.47
1:B:112:MET:HB3	1:B:113:PRO:CD	2.43	0.47
1:B:139:MET:HE2	1:B:139:MET:HA	1.97	0.47
1:A:175:LEU:O	1:A:179:GLU:HG3	2.14	0.47
5:A:402:HOH:O	2:C:6:BB7:H31	2.15	0.47
1:B:44:ARG:HE	1:B:67:VAL:HG12	1.80	0.47
1:B:348:GLU:OE1	1:B:348:GLU:N	2.45	0.47
1:A:34:VAL:HG21	1:A:188:ILE:HG13	1.97	0.47
1:A:135:ASN:CG	1:A:136:LYS:N	2.72	0.47
1:B:19:HIS:ND1	1:B:112:MET:HB3	2.29	0.47
1:B:287:GLU:O	1:B:290:GLN:HG3	2.15	0.47
1:A:251:GLN:HG3	5:A:455:HOH:O	2.14	0.46
1:B:30:ALA:HA	1:B:178:LEU:CD2	2.41	0.46
1:A:184:TRP:C	1:A:186:ALA:H	2.22	0.46
1:A:331:TYR:HD2	1:A:336:ASP:OD1	1.98	0.46
1:B:9:LYS:HD2	1:B:74:ARG:N	2.30	0.46
1:A:243:GLU:HG3	1:A:295:PRO:HA	1.97	0.46
1:B:144:GLU:O	1:B:148:LEU:HD23	2.15	0.46
1:B:210:PHE:CE1	1:B:236:ILE:HB	2.50	0.46
1:B:171:ARG:O	1:B:187:LYS:HG3	2.16	0.46
1:B:23:GLY:O	1:B:24:LYS:C	2.59	0.45
1:A:135:ASN:CG	1:A:136:LYS:H	2.24	0.45
1:B:114:GLN:HA	1:B:117:GLU:HB2	1.98	0.45
1:B:138:ASP:OD2	1:B:139:MET:HE3	2.16	0.45
1:B:176:LYS:HB3	1:B:184:TRP:CD1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:LYS:HB3	1:B:209:PRO:HD2	1.99	0.45
1:B:212:LEU:C	1:B:212:LEU:HD23	2.41	0.45
1:A:342:GLU:O	1:A:342:GLU:HG2	2.17	0.45
1:A:27:LEU:O	1:A:31:ILE:HG13	2.16	0.45
1:A:90:ASN:ND2	1:A:95:ALA:HB3	2.31	0.45
1:A:34:VAL:CB	1:A:188:ILE:HG21	2.47	0.45
1:A:46:PHE:O	1:A:47:ASP:C	2.60	0.45
1:A:318:ARG:NH1	1:A:378:GLU:CD	2.72	0.45
1:A:120:LEU:O	1:A:120:LEU:HD12	2.16	0.45
1:A:145:LEU:N	1:A:145:LEU:HD12	2.32	0.45
1:A:125:VAL:HG22	1:A:373:ARG:NH2	2.32	0.45
1:B:282:LYS:O	1:B:283:ARG:C	2.59	0.45
1:B:19:HIS:CE1	1:B:112:MET:HB3	2.52	0.45
1:B:135:ASN:CG	1:B:136:LYS:H	2.25	0.45
1:B:184:TRP:C	1:B:186:ALA:N	2.73	0.45
1:B:74:ARG:NH1	1:B:199:ILE:O	2.49	0.44
1:B:112:MET:HE3	1:B:113:PRO:HD3	1.99	0.44
1:B:189:LEU:N	1:B:189:LEU:CD2	2.79	0.44
1:B:344:PRO:HG3	1:B:357:LYS:O	2.16	0.44
1:A:105:VAL:O	1:A:134:LEU:HD13	2.18	0.44
1:A:129:TYR:CE2	1:A:200:PRO:HD2	2.53	0.44
1:B:19:HIS:HA	1:B:114:GLN:HB2	1.98	0.44
1:B:187:LYS:HA	1:B:187:LYS:HD2	1.74	0.44
1:A:112:MET:HB3	1:A:113:PRO:CD	2.48	0.44
1:A:241:GLU:HG3	1:A:253:SER:O	2.17	0.43
1:B:19:HIS:ND1	1:B:112:MET:CB	2.81	0.43
1:B:247:ILE:N	1:B:365:PRO:O	2.41	0.43
1:B:319:HIS:CE1	5:B:449:HOH:O	2.67	0.43
1:B:143:GLU:O	1:B:146:LEU:HB2	2.19	0.43
1:A:81:CYS:HB2	1:A:87:TYR:CE2	2.52	0.43
1:B:76:TYR:OH	1:B:195:LEU:HB3	2.17	0.43
1:B:96:ALA:O	1:B:97:GLN:CB	2.65	0.43
1:A:147:GLU:HG2	1:A:151:MET:HE3	2.00	0.43
1:A:256:THR:CG2	1:A:279:ARG:HB2	2.49	0.43
1:B:9:LYS:HE2	1:B:71:THR:O	2.19	0.43
1:B:188:ILE:O	1:B:191:LEU:HB3	2.18	0.43
1:A:74:ARG:NH1	1:A:195:LEU:O	2.52	0.43
1:A:149:VAL:HG13	1:A:150:GLU:N	2.33	0.43
1:B:348:GLU:CD	1:B:348:GLU:N	2.70	0.43
1:B:244:ILE:HG22	1:B:290:GLN:NE2	2.34	0.42
1:B:33:THR:HA	1:B:43:ALA:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:LEU:C	1:A:149:VAL:HG12	2.44	0.42
1:B:34:VAL:CG2	1:B:188:ILE:HG21	2.45	0.42
1:A:201:GLU:O	1:A:201:GLU:HG3	2.20	0.42
1:B:91:MET:O	1:B:125:VAL:CG2	2.64	0.42
1:B:337:VAL:HG21	1:B:366:ILE:HD11	2.02	0.42
1:A:34:VAL:HB	1:A:188:ILE:HG21	2.02	0.42
1:B:16:THR:HG23	1:B:78:HIS:HE1	1.79	0.42
1:A:134:LEU:HD13	1:A:134:LEU:HA	1.71	0.41
1:B:168:PRO:HD3	1:B:198:TYR:CE1	2.55	0.41
1:A:90:ASN:HD22	1:A:90:ASN:HA	1.75	0.41
1:A:38:THR:HG22	1:A:39:TYR:N	2.35	0.41
1:A:84:HIS:CD2	1:A:118:HIS:CE1	3.03	0.41
1:A:282:LYS:HD2	1:A:282:LYS:N	2.36	0.41
1:A:71:THR:CB	1:A:72:PRO:CD	2.98	0.41
1:A:261:PHE:O	1:A:262:ARG:HB2	2.21	0.41
1:B:34:VAL:HG21	1:B:188:ILE:CG2	2.48	0.41
1:B:87:TYR:O	1:B:91:MET:HG2	2.20	0.41
1:A:357:LYS:HZ2	1:A:357:LYS:HG3	1.75	0.41
1:B:238:VAL:HG22	1:B:257:GLY:HA2	2.02	0.41
1:B:326:TYR:CE2	1:B:328:PRO:HG3	2.54	0.41
1:A:32:THR:HG21	1:A:44:ARG:CG	2.49	0.41
1:A:378:GLU:HG3	1:A:383:VAL:HG11	2.03	0.41
1:B:123:ARG:HH21	1:B:160:TYR:HD1	1.67	0.41
1:A:112:MET:O	1:A:115:THR:N	2.51	0.41
1:A:363:ILE:HD11	5:A:462:HOH:O	2.21	0.41
1:B:131:ILE:HD11	1:B:198:TYR:CG	2.53	0.41
1:B:164:GLY:HA2	1:B:167:THR:OG1	2.20	0.41
1:A:46:PHE:CD1	1:A:46:PHE:C	2.99	0.40
1:A:112:MET:HB3	1:A:113:PRO:HD2	2.03	0.40
1:A:351:MET:HE3	1:A:352:PRO:HD2	2.03	0.40
1:A:313:LYS:C	1:A:313:LYS:HD2	2.46	0.40
1:A:140:VAL:HG11	1:A:146:LEU:CD2	2.32	0.40
1:B:90:ASN:OD1	1:B:95:ALA:HB3	2.21	0.40
1:B:125:VAL:CG1	1:B:127:VAL:HG23	2.52	0.40
1:B:153:VAL:O	1:B:157:LEU:HD22	2.21	0.40
1:B:215:GLU:O	1:B:288:ARG:HD2	2.21	0.40
1:A:141:ASP:O	1:A:142:ASP:CB	2.66	0.40
1:B:150:GLU:O	1:B:154:ARG:HG3	2.22	0.40
1:B:282:LYS:HD2	1:B:283:ARG:N	2.22	0.40
1:B:332:PHE:CE1	1:B:374:PHE:HB3	2.57	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	384/394 (98%)	357 (93%)	24 (6%)	3 (1%)	16	20
1	B	384/394 (98%)	354 (92%)	24 (6%)	6 (2%)	7	7
2	C	2/12 (17%)	1 (50%)	1 (50%)	0	100	100
2	D	2/12 (17%)	1 (50%)	1 (50%)	0	100	100
All	All	772/812 (95%)	713 (92%)	50 (6%)	9 (1%)	10	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	97	GLN
1	A	247	ILE
1	B	34	VAL
1	B	142	ASP
1	A	108	THR
1	B	333	ARG
1	B	247	ILE
1	B	111	PRO
1	A	82	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	318/325 (98%)	286 (90%)	32 (10%)	7	7
1	B	318/325 (98%)	282 (89%)	36 (11%)	5	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	1/2 (50%)	1 (100%)	0	100	100
2	D	1/2 (50%)	1 (100%)	0	100	100
All	All	638/654 (98%)	570 (89%)	68 (11%)	6	6

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	THR
1	A	44	ARG
1	A	48	GLN
1	A	55	GLU
1	A	58	ARG
1	A	89	LYS
1	A	90	ASN
1	A	91	MET
1	A	93	THR
1	A	97	GLN
1	A	134	LEU
1	A	139	MET
1	A	140	VAL
1	A	142	ASP
1	A	143	GLU
1	A	157	LEU
1	A	181	ASP
1	A	190	GLU
1	A	241	GLU
1	A	249	GLU
1	A	252	LYS
1	A	262	ARG
1	A	264	LEU
1	A	284	GLU
1	A	292	LEU
1	A	313	LYS
1	A	315	GLU
1	A	342	GLU
1	A	357	LYS
1	A	364	HIS
1	A	378	GLU
1	A	381	ARG
1	B	9	LYS
1	B	14	VAL

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Mol	Chain	Res	Type
1	B	44	ARG
1	B	48	GLN
1	B	56	LYS
1	B	97	GLN
1	B	117	GLU
1	B	121	LEU
1	B	125	VAL
1	B	136	LYS
1	B	141	ASP
1	B	144	GLU
1	B	157	LEU
1	B	159	GLN
1	B	176	LYS
1	B	187	LYS
1	B	189	LEU
1	B	195	LEU
1	B	197	SER
1	B	203	GLU
1	B	204	ARG
1	B	217	VAL
1	B	219	SER
1	B	220	ILE
1	B	223	ARG
1	B	244	ILE
1	B	265	LEU
1	B	283	ARG
1	B	284	GLU
1	B	287	GLU
1	B	292	LEU
1	B	299	LYS
1	B	313	LYS
1	B	345	GLU
1	B	362	LEU
1	B	373	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	22	HIS
1	A	51	ASN
1	A	84	HIS

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Mol	Chain	Res	Type
1	A	90	ASN
1	A	114	GLN
1	A	118	HIS
1	A	124	GLN
1	A	159	GLN
1	A	301	HIS
1	A	329	GLN
1	B	11	HIS
1	B	63	ASN
1	B	84	HIS
1	B	118	HIS
1	B	159	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

18 non-standard protein/DNA/RNA residues 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
2	BB6	C	4	2	4,6,7	2.83	2 (50%)	3,7,9	1.20	0
2	MH6	C	11	2	3,3,6	3.30	2 (66%)	1,3,7	2.53	1 (100%)
2	BB9	C	9	2	2,4,6	3.59	1 (50%)	3,4,7	2.65	3 (100%)
2	MEN	D	3	2	7,7,9	1.10	1 (14%)	8,8,11	1.52	1 (12%)
2	9BB	C	12	2	15,17,17	2.29	3 (20%)	18,22,22	1.91	2 (11%)
2	BB8	C	8	2	11,11,13	2.02	5 (45%)	12,14,17	1.46	1 (8%)
2	MEN	C	3	2	7,7,9	0.93	0	8,8,11	1.41	1 (12%)
2	MH6	D	11	2	3,3,6	4.31	3 (100%)	1,3,7	2.29	1 (100%)
2	BB9	C	2	2	2,5,6	2.75	1 (50%)	1,5,7	2.85	1 (100%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BB9	D	10	2	2,4,6	3.69	1 (50%)	3,4,7	5.40	2 (66%)
2	BB9	D	9	2	2,4,6	3.30	1 (50%)	3,4,7	2.76	3 (100%)
2	BB7	C	6	2	6,8,9	2.18	2 (33%)	4,9,11	1.60	1 (25%)
2	BB9	D	2	2	2,5,6	2.82	1 (50%)	1,5,7	2.97	1 (100%)
2	BB8	D	8	2	11,11,13	1.95	6 (54%)	12,14,17	1.35	1 (8%)
2	9BB	D	12	2	15,17,17	2.38	4 (26%)	18,22,22	1.74	4 (22%)
2	BB9	C	10	2	2,4,6	4.06	1 (50%)	3,4,7	4.57	2 (66%)
2	BB7	D	6	2	6,8,9	2.13	2 (33%)	4,9,11	1.84	1 (25%)
2	BB6	D	4	2	4,6,7	2.84	2 (50%)	3,7,9	0.82	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
2	BB6	C	4	2	-	0/0/6/8	-
2	9BB	C	12	2	-	0/10/24/24	0/1/1/1
2	BB9	C	9	2	-	0/0/2/6	-
2	MEN	D	3	2	-	2/6/6/10	-
2	BB8	C	8	2	-	5/8/8/12	0/1/1/1
2	MEN	C	3	2	-	2/6/6/10	-
2	BB9	C	2	2	-	0/0/4/6	-
2	BB9	D	10	2	-	0/0/2/6	-
2	BB9	D	9	2	-	0/0/2/6	-
2	BB7	C	6	2	-	1/1/9/11	-
2	BB9	D	2	2	-	0/0/4/6	-
2	BB8	D	8	2	-	3/8/8/12	0/1/1/1
2	9BB	D	12	2	-	2/10/24/24	0/1/1/1
2	BB9	C	10	2	-	0/0/2/6	-
2	BB7	D	6	2	-	1/1/9/11	-
2	BB6	D	4	2	-	0/0/6/8	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	12	9BB	CA-N	5.88	1.47	1.34
2	D	11	MH6	CB-CA	-5.80	1.41	1.49
2	C	10	BB9	CA-N	5.52	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	12	9BB	CA-N	5.15	1.45	1.34
2	D	4	BB6	CA-N	5.08	1.48	1.36
2	D	10	BB9	CA-N	4.94	1.45	1.33
2	C	11	MH6	CB-CA	-4.76	1.42	1.49
2	C	9	BB9	CA-N	4.74	1.44	1.33
2	D	12	9BB	CA-N13	4.63	1.49	1.40
2	C	12	9BB	CA-N13	4.49	1.49	1.40
2	C	4	BB6	CA-N	4.49	1.47	1.36
2	D	9	BB9	CA-N	4.45	1.44	1.33
2	D	6	BB7	CA-N	4.40	1.47	1.36
2	C	6	BB7	CA-N	4.29	1.46	1.36
2	C	2	BB9	C-CA	3.69	1.51	1.45
2	D	2	BB9	C-CA	3.67	1.51	1.45
2	D	12	9BB	C52-C55	3.62	1.57	1.51
2	D	11	MH6	C-CA	-3.46	1.44	1.49
2	C	4	BB6	C-CA	3.34	1.50	1.45
2	C	8	BB8	CD1-CG	3.30	1.44	1.39
2	C	8	BB8	CE2-CD2	3.18	1.44	1.38
2	D	11	MH6	CA-N	3.17	1.35	1.27
2	D	8	BB8	CD1-CG	3.04	1.43	1.39
2	D	8	BB8	CE2-CD2	2.99	1.44	1.38
2	C	6	BB7	C-CA	2.75	1.49	1.45
2	C	11	MH6	CA-N	2.73	1.33	1.27
2	D	6	BB7	C-CA	2.73	1.49	1.45
2	D	12	9BB	OXT-C49	-2.64	1.40	1.46
2	D	4	BB6	C-CA	2.51	1.49	1.45
2	C	8	BB8	CE1-CD1	2.50	1.43	1.38
2	D	8	BB8	CE2-CZ	2.49	1.43	1.38
2	D	8	BB8	CD2-CG	2.49	1.43	1.39
2	C	8	BB8	CE2-CZ	2.40	1.43	1.38
2	C	12	9BB	OXT-C49	-2.33	1.41	1.46
2	D	8	BB8	CZ-CE1	2.23	1.43	1.38
2	D	8	BB8	CE1-CD1	2.20	1.42	1.38
2	D	3	MEN	CB-CA	2.11	1.57	1.53
2	C	8	BB8	CD2-CG	2.11	1.42	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	10	BB9	C-CA-N	8.72	125.96	116.44
2	C	10	BB9	C-CA-N	7.14	124.25	116.44
2	C	12	9BB	C49-OXT-C	6.66	126.03	116.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	12	9BB	OXT-C-N13	4.85	115.34	107.97
2	C	8	BB8	C-CA-CB	-4.20	105.75	112.04
2	D	8	BB8	C-CA-CB	-3.85	106.27	112.04
2	D	10	BB9	CB-CA-N	-3.37	112.16	122.53
2	C	10	BB9	CB-CA-N	-3.27	112.46	122.53
2	D	6	BB7	O-C-CA	-3.12	119.68	125.53
2	D	2	BB9	O-C-CA	-2.97	121.66	125.39
2	D	3	MEN	CB-CG-ND2	-2.92	111.72	115.53
2	C	3	MEN	CB-CG-ND2	-2.86	111.80	115.53
2	C	2	BB9	O-C-CA	-2.85	121.82	125.39
2	D	9	BB9	C-CA-CB	2.81	126.62	121.45
2	D	12	9BB	C49-OXT-C	2.78	120.49	116.52
2	D	9	BB9	CB-CA-N	-2.78	113.97	122.53
2	D	9	BB9	C-CA-N	2.70	119.40	116.44
2	C	9	BB9	C-CA-N	2.69	119.39	116.44
2	C	9	BB9	CB-CA-N	-2.65	114.36	122.53
2	C	6	BB7	O-C-CA	-2.65	120.56	125.53
2	C	9	BB9	C-CA-CB	2.61	126.25	121.45
2	C	11	MH6	C-CA-CB	2.53	122.85	117.62
2	D	12	9BB	OXT-C-O	-2.53	120.83	124.55
2	D	12	9BB	C53-C54-C49	-2.34	107.20	110.96
2	C	12	9BB	C53-C54-C49	2.32	114.69	110.96
2	D	11	MH6	C-CA-CB	2.29	122.37	117.62

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	6	BB7	CB-CB1-OB2-CB3
2	D	12	9BB	O-C-OXT-C49
2	D	12	9BB	N13-C-OXT-C49
2	C	8	BB8	C-CA-CB-CG
2	D	8	BB8	C-CA-CB-CG
2	C	8	BB8	OB-CB-CG-CD1
2	C	8	BB8	OB-CB-CG-CD2
2	C	6	BB7	CB-CB1-OB2-CB3
2	C	8	BB8	CA-CB-CG-CD1
2	C	8	BB8	CA-CB-CG-CD2
2	C	3	MEN	CA-CB-CG-OD1
2	D	3	MEN	CA-CB-CG-OD1
2	D	8	BB8	OB-CB-CG-CD2
2	D	8	BB8	OB-CB-CG-CD1

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Mol	Chain	Res	Type	Atoms
2	C	3	MEN	CA-CB-CG-ND2
2	D	3	MEN	CA-CB-CG-ND2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	6	BB7	1	0
2	D	12	9BB	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GDP	A	500	4	29,30,30	1.89	4 (13%)	45,47,47	1.99	11 (24%)
3	GDP	B	500	4	29,30,30	1.87	5 (17%)	45,47,47	2.00	11 (24%)

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	500	4	-	6/16/32/32	0/3/3/3
3	GDP	B	500	4	-	3/16/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	GDP	C4-N3	6.49	1.49	1.34
3	B	500	GDP	C4-N3	6.27	1.48	1.34
3	A	500	GDP	C5-N7	4.55	1.48	1.39
3	B	500	GDP	C5-N7	4.12	1.47	1.39
3	B	500	GDP	PA-O3A	-3.59	1.55	1.59
3	A	500	GDP	C4-N9	3.46	1.46	1.38
3	B	500	GDP	C4-N9	2.87	1.45	1.38
3	A	500	GDP	C8-N9	2.37	1.42	1.37
3	B	500	GDP	C8-N9	2.26	1.42	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	GDP	C1'-N9-C4	-5.80	109.37	126.49
3	A	500	GDP	C1'-N9-C4	-5.67	109.75	126.49
3	A	500	GDP	O4'-C1'-N9	5.00	119.68	108.36
3	B	500	GDP	O4'-C1'-N9	4.67	118.93	108.36
3	A	500	GDP	C5-C4-N3	-4.26	121.61	128.39
3	B	500	GDP	C5-C4-N3	-4.20	121.71	128.39
3	A	500	GDP	C2-N3-C4	3.83	118.90	112.30
3	B	500	GDP	C2'-C1'-N9	-3.71	102.93	113.25
3	B	500	GDP	C2-N3-C4	3.66	118.61	112.30
3	A	500	GDP	C2'-C1'-N9	-3.64	103.11	113.25
3	A	500	GDP	C5-C4-N9	3.31	111.57	105.66
3	B	500	GDP	C5-C4-N9	3.23	111.44	105.66
3	A	500	GDP	C8-N9-C4	-3.08	100.25	106.03
3	B	500	GDP	C8-N9-C4	-2.99	100.42	106.03
3	B	500	GDP	O6-C6-C5	-2.72	119.35	126.53
3	A	500	GDP	C1'-N9-C8	2.54	133.96	126.73
3	A	500	GDP	O6-C6-C5	-2.54	119.84	126.53
3	A	500	GDP	O3B-PB-O3A	2.48	112.95	104.64
3	B	500	GDP	O3B-PB-O3A	2.42	112.76	104.64
3	B	500	GDP	N1-C2-N3	-2.30	119.12	123.32
3	A	500	GDP	N1-C2-N3	-2.24	119.22	123.32
3	B	500	GDP	O4'-C4'-C5'	2.01	115.79	109.33

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	500	GDP	PA-O3A-PB-O3B

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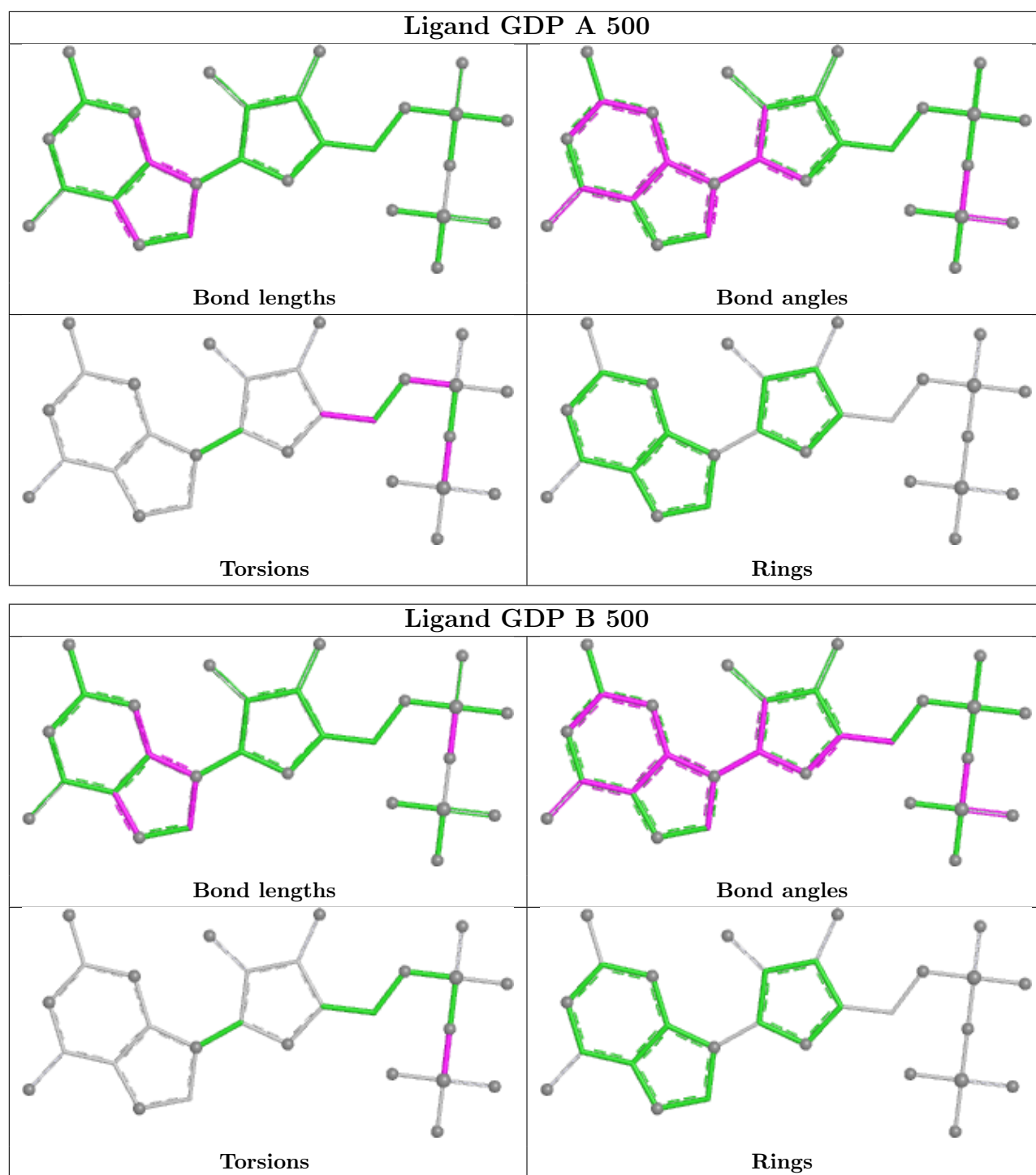
Mol	Chain	Res	Type	Atoms
3	A	500	GDP	C5'-O5'-PA-O2A
3	B	500	GDP	PA-O3A-PB-O3B
3	A	500	GDP	C5'-O5'-PA-O3A
3	A	500	GDP	C5'-O5'-PA-O1A
3	B	500	GDP	PA-O3A-PB-O1B
3	A	500	GDP	PA-O3A-PB-O2B
3	B	500	GDP	PA-O3A-PB-O2B
3	A	500	GDP	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	GDP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	386/394 (97%)	0.21	11 (2%) 55 56	21, 48, 80, 101	0
1	B	386/394 (97%)	0.28	14 (3%) 46 46	26, 48, 81, 99	0
2	C	3/12 (25%)	0.36	0 100 100	33, 33, 38, 54	0
2	D	3/12 (25%)	-0.07	0 100 100	33, 33, 40, 52	0
All	All	778/812 (95%)	0.25	25 (3%) 50 51	21, 48, 80, 101	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	ILE	3.1
1	A	39	TYR	2.9
1	A	47	ASP	2.8
1	B	42	ALA	2.8
1	B	38	THR	2.7
1	B	182	ALA	2.7
1	B	59	GLY	2.7
1	A	8	THR	2.6
1	A	42	ALA	2.6
1	B	47	ASP	2.5
1	A	149	VAL	2.5
1	A	72	PRO	2.4
1	B	10	PRO	2.4
1	B	140	VAL	2.3
1	B	44	ARG	2.3
1	A	11	HIS	2.3
1	A	146	LEU	2.3
1	B	8	THR	2.2
1	A	184	TRP	2.2
1	B	181	ASP	2.2
1	B	194	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	140	VAL	2.1
1	B	111	PRO	2.1
1	B	145	LEU	2.1
1	A	143	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	BB7	D	6	9/10	0.89	0.12	50,54,66,66	0
2	BB8	D	8	11/13	0.90	0.12	35,37,40,42	0
2	9BB	C	12	17/17	0.91	0.14	38,63,79,80	0
2	9BB	D	12	17/17	0.92	0.12	37,65,79,81	0
2	BB7	C	6	9/10	0.93	0.12	49,55,62,63	0
2	BB8	C	8	11/13	0.93	0.09	27,30,32,33	0
2	BB6	C	4	7/8	0.93	0.10	40,43,49,51	0
2	MH6	D	11	4/7	0.95	0.07	33,34,36,37	0
2	MEN	C	3	8/10	0.95	0.08	30,34,38,40	0
2	BB9	D	2	6/7	0.95	0.08	32,33,35,36	0
2	BB9	C	2	6/7	0.96	0.08	35,36,38,41	0
2	MH6	C	11	4/7	0.96	0.07	32,33,35,40	0
2	BB6	D	4	7/8	0.97	0.07	33,34,40,41	0
2	MEN	D	3	8/10	0.97	0.06	25,30,35,35	0
2	BB9	D	10	5/7	0.98	0.06	28,28,29,29	0
2	BB9	D	9	5/7	0.98	0.05	30,32,34,37	0
2	BB9	C	10	5/7	0.99	0.05	25,26,28,29	0
2	BB9	C	9	5/7	0.99	0.04	23,25,27,27	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

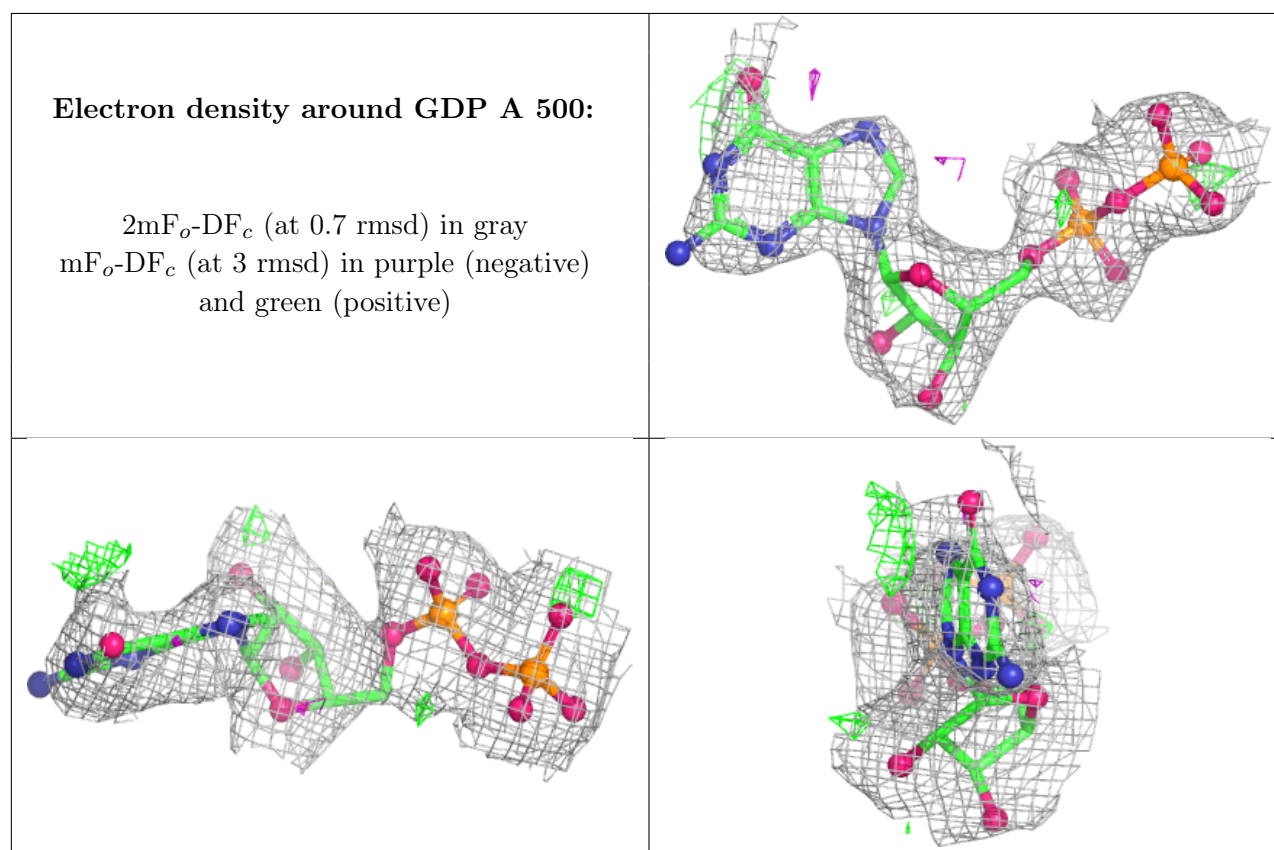
6.4 Ligands [i](#)

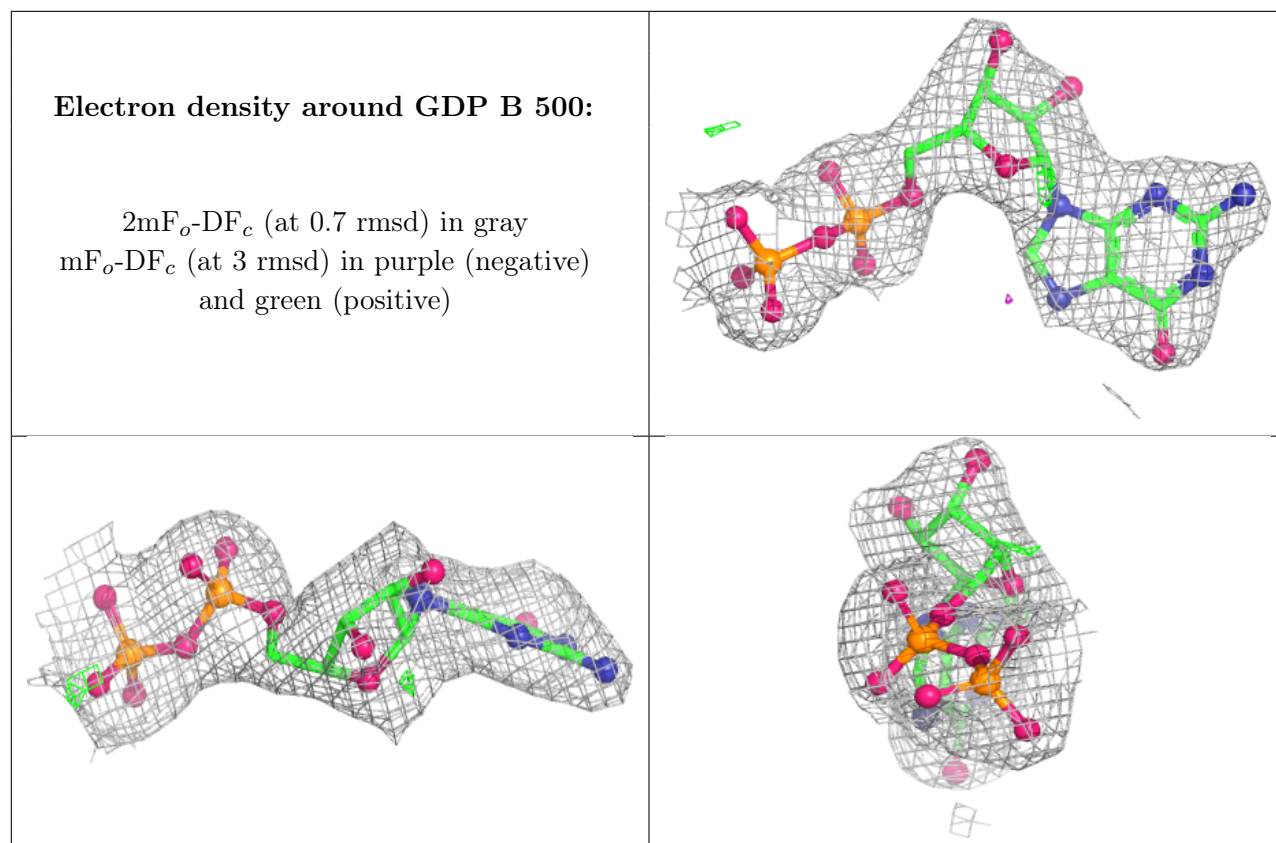
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GDP	A	500	28/28	0.90	0.13	51,75,81,81	0
3	GDP	B	500	28/28	0.93	0.10	39,60,63,64	0
4	MG	A	394	1/1	0.96	0.07	41,41,41,41	0
4	MG	B	394	1/1	0.99	0.08	46,46,46,46	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.