



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:58 PM UTC

PDB ID : 3P27 / pdb_00003p27
Title : Crystal structure of *S. cerevisiae* Hbs1 protein (GDP-bound form), a translational GTPase involved in RNA quality control pathways and interacting with Dom34/Pelota
Authors : van den Elzen, A.; Henri, J.; Lazar, N.; Gas, M.E.; Durand, D.; Lacroute, F.; Nicaise, M.; van Tilbeurgh, H.; Sraphin, B.; Graille, M.; Paris-Sud Yeast Structural Genomics (YSG)
Deposited on : 2010-10-01
Resolution : 2.95 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)

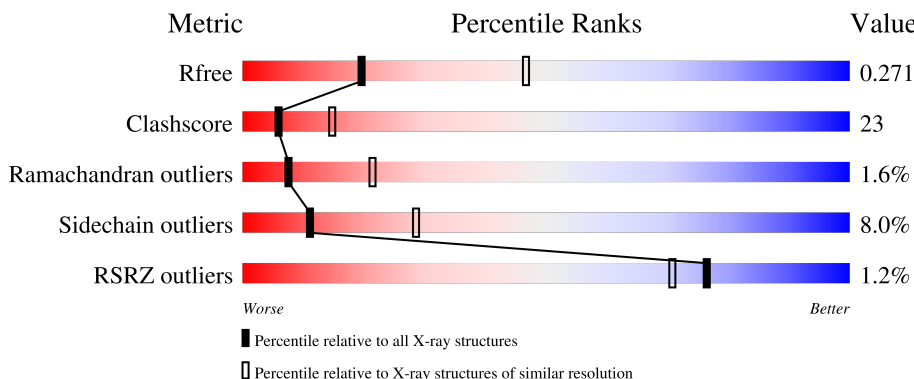
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	483	52% 34% • 10%
1	B	483	53% 34% • 8%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7043 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.

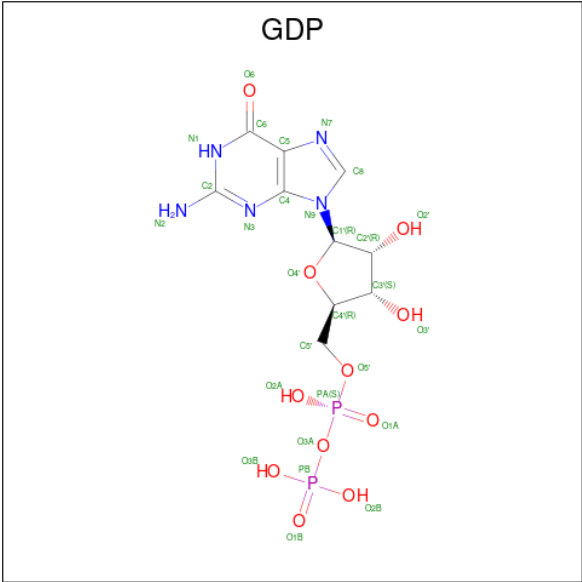
- Molecule 1 is a protein called Elongation factor 1 alpha-like protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	Se	0	0	0
			3414	2178	583	639	5	9			
1	B	444	Total	C	N	O	S	Se	0	0	0
			3509	2232	605	658	5	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	612	HIS	-	expression tag	UNP P32769
A	613	HIS	-	expression tag	UNP P32769
A	614	HIS	-	expression tag	UNP P32769
A	615	HIS	-	expression tag	UNP P32769
A	616	HIS	-	expression tag	UNP P32769
A	617	HIS	-	expression tag	UNP P32769
B	612	HIS	-	expression tag	UNP P32769
B	613	HIS	-	expression tag	UNP P32769
B	614	HIS	-	expression tag	UNP P32769
B	615	HIS	-	expression tag	UNP P32769
B	616	HIS	-	expression tag	UNP P32769
B	617	HIS	-	expression tag	UNP P32769

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

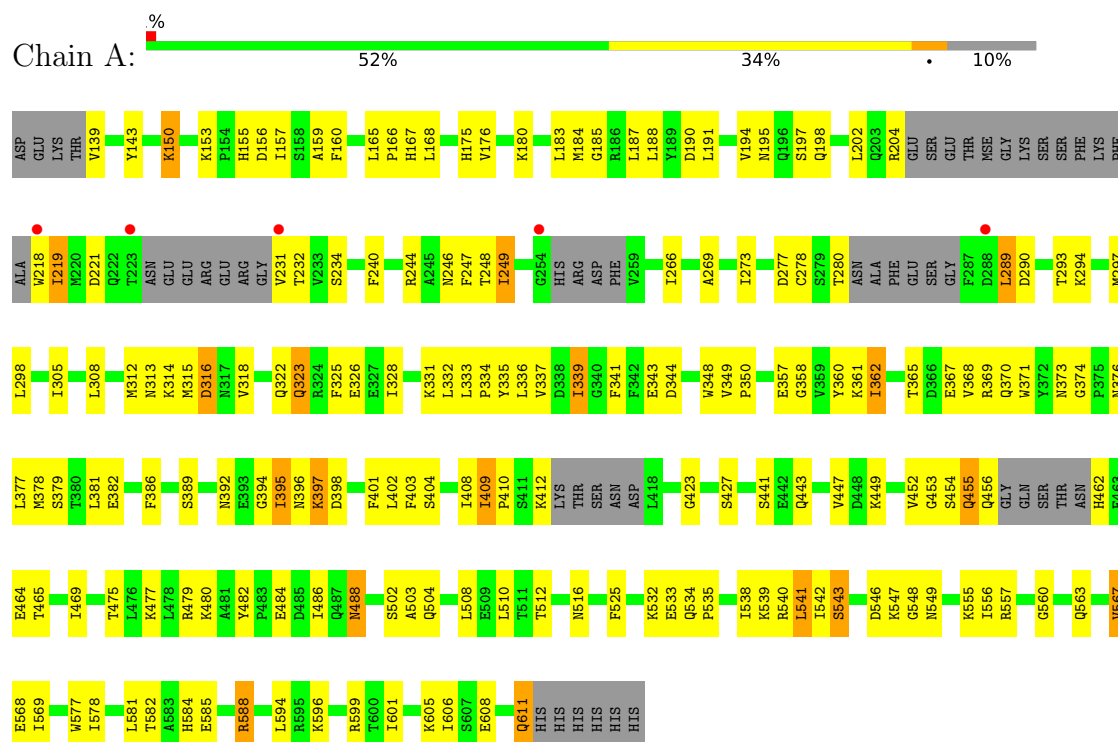
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		
3	B	33	Total	O	0	0
			33	33		

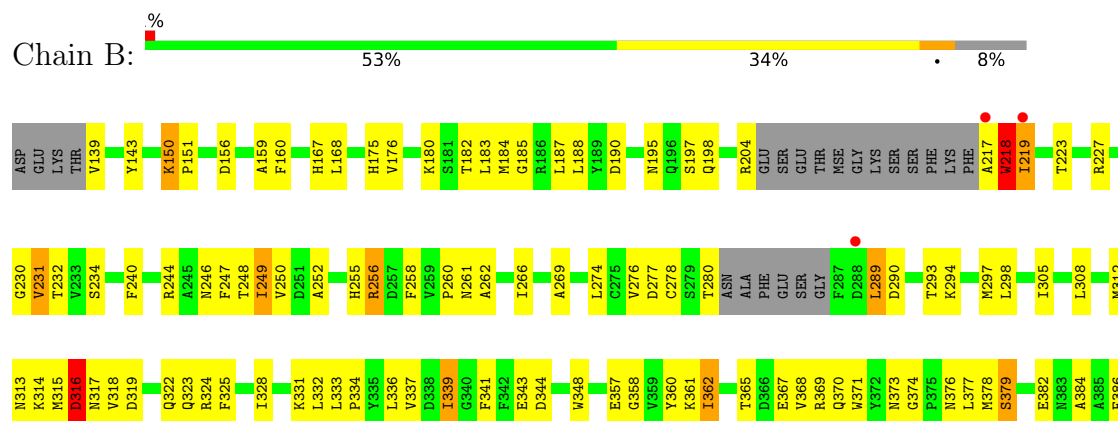
3 Residue-property plots

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: Elongation factor 1 alpha-like protein



• Molecule 1: Elongation factor 1 alpha-like protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 4 ₂ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	109.96Å 109.96Å 188.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.95 19.98 – 2.95	Depositor EDS
% Data completeness (in resolution range)	98.2 (19.98-2.95) 98.0 (19.98-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.93Å)	Xtrriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.214 , 0.278 0.207 , 0.271	Depositor DCC
R_{free} test set	1247 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	55.1	Xtrriage
Anisotropy	0.081	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7043	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.00 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0560e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0/3467	0.99	11/4667 (0.2%)
1	B	0.66	0/3566	0.98	11/4802 (0.2%)
All	All	0.65	0/7033	0.99	22/9469 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	482	TYR	CA-C-N	8.90	129.80	119.47
1	B	482	TYR	C-N-CA	8.90	129.80	119.47
1	A	482	TYR	CA-C-N	8.30	129.09	119.47
1	A	482	TYR	C-N-CA	8.30	129.09	119.47
1	B	218	TRP	N-CA-C	-6.19	104.83	112.38
1	B	150	LYS	CA-C-N	5.88	125.64	119.76
1	B	150	LYS	C-N-CA	5.88	125.64	119.76
1	A	374	GLY	CA-C-N	5.80	125.78	120.03
1	A	374	GLY	C-N-CA	5.80	125.78	120.03
1	A	397	LYS	N-CA-C	-5.70	106.33	113.28
1	A	150	LYS	CA-C-N	5.59	125.35	119.76
1	A	150	LYS	C-N-CA	5.59	125.35	119.76
1	A	409	ILE	CA-C-N	5.58	125.51	119.76
1	A	409	ILE	C-N-CA	5.58	125.51	119.76
1	A	539	LYS	N-CA-C	-5.51	105.60	112.93
1	B	374	GLY	CA-C-N	5.51	125.48	120.03
1	B	374	GLY	C-N-CA	5.51	125.48	120.03
1	B	539	LYS	N-CA-C	-5.46	105.66	112.93
1	B	316	ASP	CB-CA-C	-5.36	102.46	110.88
1	B	397	LYS	N-CA-C	-5.31	106.80	113.28
1	B	276	VAL	N-CA-C	5.21	115.46	108.17
1	A	503	ALA	N-CA-C	5.05	116.52	108.79

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	3414	0	3457	161	0
1	B	3509	0	3535	177	0
2	A	28	0	12	4	0
2	B	28	0	12	3	0
3	A	31	0	0	3	0
3	B	33	0	0	0	0
All	All	7043	0	7016	321	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 23.

All (321) 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:175:HIS:CE1	1:A:176:VAL:HG12	2.04	0.93
1:A:611:GLN:HE21	1:B:611:GLN:HE21	1.13	0.92
1:B:175:HIS:CE1	1:B:176:VAL:HG12	2.05	0.91
1:A:167:HIS:HE1	1:A:248:THR:OG1	1.56	0.88
1:B:167:HIS:HE1	1:B:248:THR:OG1	1.59	0.86
1:B:187:LEU:HD11	1:B:378:MSE:CE	2.06	0.85
1:A:297:MSE:HE1	1:A:308:LEU:HD13	1.59	0.84
1:A:512:THR:HG22	1:A:563:GLN:H	1.41	0.83
1:A:187:LEU:HD11	1:A:378:MSE:CE	2.09	0.83
1:A:187:LEU:HD11	1:A:378:MSE:HE3	1.61	0.82
1:B:452:VAL:HG22	1:B:453:GLY:H	1.42	0.82
1:B:187:LEU:HD11	1:B:378:MSE:HE3	1.61	0.81
1:B:512:THR:HG22	1:B:563:GLN:H	1.46	0.80
1:B:240:PHE:HE1	1:B:249:ILE:HD11	1.46	0.80
1:A:240:PHE:HE1	1:A:249:ILE:HD11	1.46	0.80
1:A:297:MSE:CE	1:A:308:LEU:HD13	2.12	0.79
1:B:512:THR:CG2	1:B:563:GLN:H	1.94	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:VAL:HG22	1:A:453:GLY:H	1.46	0.78
1:A:488:ASN:H	1:A:488:ASN:HD22	1.30	0.77
1:A:512:THR:CG2	1:A:563:GLN:H	1.97	0.77
1:A:358:GLY:O	1:A:376:ASN:HB2	1.85	0.76
1:B:297:MSE:CE	1:B:308:LEU:HD13	2.15	0.76
1:B:488:ASN:HD22	1:B:488:ASN:H	1.33	0.76
1:B:367:GLU:O	1:B:370:GLN:HG2	1.86	0.76
1:A:231:VAL:HG13	1:A:232:THR:H	1.52	0.75
1:B:297:MSE:HE1	1:B:308:LEU:HD13	1.68	0.75
1:B:525:PHE:HZ	1:B:569:ILE:HD12	1.51	0.74
1:B:231:VAL:HG13	1:B:232:THR:H	1.53	0.74
1:A:525:PHE:HZ	1:A:569:ILE:HD12	1.54	0.73
1:A:297:MSE:HE1	1:A:308:LEU:CD1	2.19	0.72
1:B:240:PHE:CE1	1:B:249:ILE:HD11	2.24	0.72
1:A:449:LYS:HE2	1:A:464:GLU:OE2	1.90	0.72
1:A:240:PHE:CE1	1:A:249:ILE:HD11	2.24	0.71
1:B:452:VAL:HG22	1:B:453:GLY:N	2.04	0.71
1:B:333:LEU:O	1:B:337:VAL:HG23	1.91	0.70
1:B:358:GLY:O	1:B:376:ASN:HB2	1.90	0.69
1:B:278:CYS:SG	1:B:312:MSE:HE3	2.33	0.69
1:A:452:VAL:HG22	1:A:453:GLY:N	2.08	0.69
1:B:449:LYS:HE2	1:B:464:GLU:OE2	1.92	0.69
1:A:611:GLN:HE21	1:B:611:GLN:NE2	1.88	0.69
1:B:588:ARG:HG3	1:B:588:ARG:HH11	1.56	0.69
1:A:333:LEU:O	1:A:337:VAL:HG23	1.93	0.68
1:A:367:GLU:O	1:A:370:GLN:HG2	1.93	0.68
1:B:195:ASN:HD22	1:B:197:SER:H	1.41	0.68
1:B:258:PHE:HB3	1:B:261:ASN:HD22	1.58	0.68
1:B:331:LYS:O	1:B:334:PRO:HD2	1.93	0.67
1:A:608:GLU:HG3	1:B:608:GLU:HG3	1.75	0.67
1:A:611:GLN:NE2	1:B:611:GLN:HE21	1.91	0.67
1:A:549:ASN:HB3	1:B:605:LYS:NZ	2.09	0.67
1:B:218:TRP:C	1:B:218:TRP:CD1	2.74	0.66
1:A:331:LYS:O	1:A:334:PRO:HD2	1.94	0.65
1:B:297:MSE:HE1	1:B:308:LEU:HD22	1.76	0.65
1:B:508:LEU:HD22	1:B:569:ILE:HG13	1.78	0.65
1:A:605:LYS:NZ	1:B:549:ASN:HB3	2.11	0.65
1:A:512:THR:HG21	1:A:560:GLY:O	1.96	0.65
1:B:512:THR:HG21	1:B:560:GLY:O	1.96	0.64
1:B:297:MSE:HE1	1:B:308:LEU:CD1	2.28	0.64
1:B:397:LYS:HD3	1:B:397:LYS:N	2.13	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:ASN:HD22	1:A:197:SER:H	1.45	0.63
1:A:362:ILE:HG23	3:A:6:HOH:O	1.99	0.63
1:A:508:LEU:HD22	1:A:569:ILE:HG13	1.81	0.62
1:A:542:ILE:O	1:A:543:SER:HB3	1.99	0.61
1:A:588:ARG:HG3	1:A:588:ARG:HH11	1.64	0.61
1:B:534:GLN:NE2	1:B:535:PRO:HD2	2.16	0.61
1:B:180:LYS:HB2	2:B:663:GDP:O2B	2.01	0.61
1:B:313:ASN:OD1	1:B:314:LYS:N	2.33	0.61
1:A:326:GLU:HG3	3:A:18:HOH:O	1.99	0.61
1:B:308:LEU:HD11	1:B:341:PHE:CE1	2.36	0.60
1:B:322:GLN:HB2	1:B:371:TRP:CD2	2.36	0.60
1:A:534:GLN:NE2	1:A:535:PRO:HD2	2.16	0.60
1:A:277:ASP:CG	1:A:314:LYS:HD2	2.25	0.60
1:B:240:PHE:HE1	1:B:249:ILE:CD1	2.15	0.60
1:A:175:HIS:CD2	1:A:290:ASP:OD2	2.54	0.60
1:B:277:ASP:CG	1:B:314:LYS:HD2	2.27	0.60
1:A:397:LYS:HD3	1:A:397:LYS:N	2.17	0.60
1:A:240:PHE:HE1	1:A:249:ILE:CD1	2.13	0.60
1:B:525:PHE:CZ	1:B:569:ILE:HD12	2.35	0.60
1:A:278:CYS:SG	1:A:312:MSE:HE3	2.42	0.60
1:B:175:HIS:CD2	1:B:290:ASP:OD2	2.55	0.60
1:B:316:ASP:OD1	1:B:316:ASP:N	2.31	0.60
1:A:167:HIS:CE1	1:A:248:THR:OG1	2.48	0.59
1:A:269:ALA:O	1:A:588:ARG:NH2	2.33	0.59
1:A:488:ASN:HD22	1:A:488:ASN:N	1.99	0.59
1:B:308:LEU:HD11	1:B:341:PHE:CZ	2.38	0.59
1:A:322:GLN:HB2	1:A:371:TRP:CD2	2.38	0.59
1:A:456:GLN:HG2	1:A:475:THR:OG1	2.03	0.58
1:B:185:GLY:HA2	1:B:219:ILE:HD11	1.85	0.58
1:B:297:MSE:HE1	1:B:308:LEU:CD2	2.33	0.58
1:B:452:VAL:CG2	1:B:453:GLY:H	2.16	0.58
1:A:308:LEU:HD11	1:A:341:PHE:CE1	2.39	0.58
1:B:230:GLY:O	1:B:231:VAL:C	2.47	0.58
1:B:258:PHE:HB3	1:B:261:ASN:ND2	2.19	0.58
1:B:456:GLN:HG2	1:B:475:THR:OG1	2.03	0.58
1:A:297:MSE:HE1	1:A:308:LEU:HD22	1.86	0.58
1:B:512:THR:HG22	1:B:563:GLN:N	2.16	0.58
1:B:183:LEU:HD12	1:B:183:LEU:O	2.04	0.58
1:B:218:TRP:C	1:B:218:TRP:HD1	2.12	0.57
1:B:542:ILE:O	1:B:543:SER:HB3	2.05	0.57
1:A:512:THR:HG22	1:A:563:GLN:N	2.14	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:549:ASN:HB3	1:B:605:LYS:HZ2	1.69	0.57
1:A:441:SER:OG	1:A:443:GLN:HG2	2.05	0.57
1:A:540:ARG:HB2	1:A:568:GLU:HB3	1.87	0.56
1:B:167:HIS:CE1	1:B:248:THR:OG1	2.50	0.56
1:B:512:THR:HG22	1:B:563:GLN:O	2.05	0.56
1:A:184:MSE:HE2	1:A:184:MSE:HA	1.86	0.56
1:A:316:ASP:OD2	2:A:663:GDP:N2	2.38	0.56
1:B:289:LEU:HG	1:B:294:LYS:HE3	1.87	0.56
1:B:455:GLN:O	1:B:457:GLY:N	2.39	0.56
1:A:297:MSE:CE	1:A:308:LEU:CD1	2.83	0.56
1:B:184:MSE:HA	1:B:184:MSE:HE2	1.87	0.56
1:A:289:LEU:HG	1:A:294:LYS:HE3	1.88	0.56
1:B:357:GLU:HA	1:B:362:ILE:HG13	1.87	0.56
1:A:175:HIS:CG	1:A:176:VAL:H	2.24	0.56
1:B:441:SER:OG	1:B:443:GLN:HG2	2.06	0.56
1:A:313:ASN:OD1	1:A:314:LYS:N	2.39	0.55
1:B:175:HIS:CG	1:B:176:VAL:H	2.24	0.55
1:B:488:ASN:HD22	1:B:488:ASN:N	2.01	0.55
1:A:357:GLU:HA	1:A:362:ILE:HG13	1.87	0.55
1:A:525:PHE:CZ	1:A:569:ILE:HD12	2.39	0.55
1:B:488:ASN:O	1:B:532:LYS:HE3	2.07	0.55
1:A:585:GLU:CD	1:B:540:ARG:HH22	2.14	0.54
1:A:369:ARG:HH11	1:A:373:ASN:HA	1.72	0.54
1:A:277:ASP:OD2	1:A:314:LYS:HD2	2.08	0.54
1:B:269:ALA:O	1:B:588:ARG:NH2	2.41	0.54
1:B:277:ASP:OD2	1:B:314:LYS:HD2	2.08	0.53
1:A:297:MSE:HE1	1:A:308:LEU:CD2	2.38	0.53
1:A:479:ARG:O	1:A:480:LYS:HB2	2.08	0.53
1:A:183:LEU:O	1:A:183:LEU:HD12	2.09	0.53
1:B:365:THR:HB	1:B:368:VAL:HG23	1.90	0.53
1:B:258:PHE:HD1	1:B:261:ASN:HD21	1.56	0.53
1:B:333:LEU:HD11	1:B:343:GLU:HG3	1.90	0.53
1:A:488:ASN:O	1:A:532:LYS:HE3	2.08	0.53
1:A:308:LEU:HD11	1:A:341:PHE:CZ	2.44	0.53
1:B:582:THR:HG22	1:B:606:ILE:HG22	1.91	0.53
1:A:143:TYR:C	1:A:143:TYR:CD2	2.86	0.52
1:B:256:ARG:O	1:B:256:ARG:HD3	2.10	0.52
1:B:540:ARG:HB2	1:B:568:GLU:HB3	1.89	0.52
1:A:175:HIS:CG	1:A:176:VAL:N	2.77	0.52
1:A:333:LEU:HB3	1:A:334:PRO:HD3	1.92	0.52
1:A:190:ASP:OD2	1:A:361:LYS:HB2	2.10	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:GLY:HA2	1:A:219:ILE:HD11	1.92	0.51
1:A:315:MSE:HE1	1:A:325:PHE:HB2	1.91	0.51
1:A:452:VAL:CG2	1:A:453:GLY:H	2.20	0.51
1:B:175:HIS:CG	1:B:176:VAL:N	2.78	0.51
1:B:244:ARG:HD3	1:B:382:GLU:OE2	2.09	0.51
1:A:382:GLU:O	1:A:386:PHE:HD2	1.93	0.51
1:A:180:LYS:HB2	2:A:663:GDP:O2B	2.11	0.51
1:B:452:VAL:CG2	1:B:453:GLY:N	2.74	0.51
1:B:297:MSE:CE	1:B:308:LEU:CD1	2.86	0.51
1:B:333:LEU:HB3	1:B:334:PRO:HD3	1.92	0.51
1:A:605:LYS:HZ2	1:B:549:ASN:HB3	1.76	0.51
1:A:195:ASN:HB3	1:A:198:GLN:H	1.75	0.50
1:B:588:ARG:HH11	1:B:588:ARG:CG	2.23	0.50
1:A:195:ASN:N	1:A:198:GLN:OE1	2.43	0.50
1:A:202:LEU:HD11	1:A:219:ILE:HG13	1.92	0.50
1:A:362:ILE:HG13	1:A:362:ILE:O	2.11	0.50
1:A:512:THR:HG22	1:A:563:GLN:O	2.10	0.50
1:B:219:ILE:O	1:B:223:THR:HG23	2.12	0.50
1:B:369:ARG:HH11	1:B:373:ASN:HA	1.75	0.50
1:A:333:LEU:HD11	1:A:343:GLU:HG3	1.93	0.50
1:A:502:SER:HB3	1:A:577:TRP:HB3	1.92	0.50
1:B:195:ASN:HB3	1:B:198:GLN:H	1.77	0.50
1:B:315:MSE:HE1	1:B:325:PHE:HB2	1.93	0.50
1:B:258:PHE:HD1	1:B:261:ASN:ND2	2.09	0.49
1:B:453:GLY:O	1:B:456:GLN:HB2	2.12	0.49
1:B:362:ILE:HG13	1:B:362:ILE:O	2.12	0.49
1:A:312:MSE:HE1	1:A:328:ILE:HB	1.94	0.49
1:B:588:ARG:HG3	1:B:588:ARG:NH1	2.24	0.49
1:B:143:TYR:C	1:B:143:TYR:CD2	2.90	0.49
1:B:262:ALA:O	1:B:266:ILE:HG13	2.13	0.49
1:A:333:LEU:O	1:A:333:LEU:HD12	2.13	0.48
1:B:402:LEU:HB2	1:B:427:SER:HB3	1.95	0.48
1:A:156:ASP:HB3	1:A:159:ALA:HB3	1.95	0.48
1:A:608:GLU:HA	1:B:608:GLU:HG3	1.95	0.48
1:B:538:ILE:HG21	1:B:541:LEU:HD13	1.93	0.48
1:A:582:THR:HG22	1:A:606:ILE:HG22	1.95	0.48
1:A:336:LEU:O	1:A:339:ILE:HG22	2.13	0.48
1:B:312:MSE:HE1	1:B:328:ILE:HB	1.95	0.48
1:B:392:ASN:N	1:B:392:ASN:HD22	2.09	0.48
1:B:252:ALA:H	1:B:256:ARG:HH21	1.62	0.48
1:B:502:SER:HB3	1:B:577:TRP:HB3	1.93	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:538:ILE:HG21	1:A:541:LEU:HD13	1.96	0.48
1:A:605:LYS:HZ1	1:B:549:ASN:HB3	1.77	0.48
1:B:227:ARG:HD3	1:B:457:GLY:HA3	1.95	0.48
1:B:541:LEU:HD12	1:B:567:VAL:HG22	1.95	0.48
1:A:453:GLY:O	1:A:456:GLN:HB2	2.14	0.47
1:B:449:LYS:HB2	1:B:477:LYS:HB3	1.95	0.47
1:A:394:GLY:O	1:A:395:ILE:C	2.57	0.47
1:A:452:VAL:CG2	1:A:453:GLY:N	2.77	0.47
1:B:479:ARG:O	1:B:480:LYS:HB2	2.14	0.47
1:A:167:HIS:HA	1:A:246:ASN:O	2.15	0.47
1:A:188:LEU:O	1:A:191:LEU:HB2	2.14	0.47
1:A:588:ARG:HG3	1:A:588:ARG:NH1	2.30	0.47
1:A:608:GLU:HG3	1:B:608:GLU:CG	2.43	0.47
1:B:332:LEU:HD23	1:B:348:TRP:CZ3	2.50	0.47
1:A:549:ASN:HB3	1:B:605:LYS:HZ1	1.78	0.47
1:A:332:LEU:HD23	1:A:348:TRP:CZ3	2.50	0.47
1:B:184:MSE:HB3	1:B:219:ILE:HD13	1.97	0.47
1:A:599:ARG:O	1:A:601:ILE:HG23	2.15	0.46
1:B:455:GLN:C	1:B:457:GLY:H	2.23	0.46
1:B:588:ARG:CG	1:B:588:ARG:NH1	2.76	0.46
1:A:409:ILE:HA	1:A:410:PRO:HD3	1.73	0.46
1:A:244:ARG:HD3	1:A:382:GLU:OE2	2.15	0.46
1:A:516:ASN:HD21	1:A:599:ARG:HH21	1.62	0.46
1:A:392:ASN:N	1:A:392:ASN:HD22	2.14	0.46
1:A:488:ASN:N	1:A:488:ASN:ND2	2.61	0.46
1:B:182:THR:OG1	2:B:663:GDP:H8	1.99	0.46
1:A:365:THR:HB	1:A:368:VAL:HG23	1.97	0.46
1:A:449:LYS:HB2	1:A:477:LYS:HB3	1.98	0.46
1:B:156:ASP:HB3	1:B:159:ALA:HB3	1.97	0.46
1:B:394:GLY:O	1:B:395:ILE:C	2.59	0.46
1:A:323:GLN:NE2	3:A:47:HOH:O	2.49	0.45
1:A:349:VAL:HA	1:A:350:PRO:HD3	1.86	0.45
1:A:541:LEU:HD12	1:A:567:VAL:HG22	1.99	0.45
1:A:195:ASN:HB2	1:A:198:GLN:OE1	2.16	0.45
1:A:168:LEU:O	1:A:247:PHE:HA	2.17	0.45
1:B:190:ASP:OD2	1:B:361:LYS:HB2	2.16	0.45
1:A:298:LEU:HD13	1:A:339:ILE:HG12	1.98	0.45
1:A:510:LEU:HD11	1:A:594:LEU:HD22	1.98	0.45
1:A:396:ASN:HB3	1:A:398:ASP:H	1.81	0.45
1:A:408:ILE:HD11	1:A:486:ILE:HG22	1.98	0.45
1:B:382:GLU:O	1:B:386:PHE:HD2	2.00	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:608:GLU:CG	1:B:608:GLU:HG3	2.46	0.45
1:B:298:LEU:HA	1:B:339:ILE:HD11	1.98	0.45
1:B:396:ASN:HB3	1:B:398:ASP:H	1.81	0.45
1:B:336:LEU:O	1:B:339:ILE:HG22	2.16	0.45
1:A:153:LYS:O	1:A:155:HIS:HD2	2.00	0.44
1:A:218:TRP:CD1	1:A:218:TRP:C	2.95	0.44
1:A:402:LEU:HB2	1:A:427:SER:HB3	2.00	0.44
1:A:608:GLU:HG3	1:B:608:GLU:HA	2.00	0.44
1:B:510:LEU:HD11	1:B:594:LEU:HD22	1.98	0.44
1:A:150:LYS:HE3	1:A:150:LYS:HB2	1.79	0.44
1:A:294:LYS:HG2	1:A:335:TYR:CE1	2.53	0.44
1:B:168:LEU:O	1:B:247:PHE:HA	2.16	0.44
1:A:266:ILE:HG23	1:A:305:ILE:HD12	2.00	0.43
1:B:218:TRP:HD1	1:B:218:TRP:O	2.01	0.43
1:B:312:MSE:HE1	1:B:328:ILE:CG2	2.48	0.43
1:B:217:ALA:C	1:B:219:ILE:N	2.76	0.43
1:B:317:ASN:OD1	2:B:663:GDP:N2	2.52	0.43
1:B:360:TYR:C	1:B:360:TYR:CD2	2.96	0.43
1:A:156:ASP:O	1:A:159:ALA:HB3	2.18	0.43
1:A:376:ASN:O	1:A:377:LEU:C	2.61	0.43
1:B:376:ASN:OD1	1:B:379:SER:HB2	2.18	0.43
1:B:556:ILE:HG23	1:B:556:ILE:O	2.19	0.43
1:B:333:LEU:N	1:B:334:PRO:CD	2.82	0.43
1:A:176:VAL:O	2:A:663:GDP:O1B	2.37	0.43
1:A:333:LEU:HD12	1:A:333:LEU:C	2.43	0.43
1:A:455:GLN:OE1	1:A:462:HIS:HB3	2.19	0.43
1:B:255:HIS:CB	1:B:258:PHE:HB2	2.49	0.43
1:B:258:PHE:CD1	1:B:261:ASN:ND2	2.86	0.43
1:A:369:ARG:HD3	1:A:369:ARG:HA	1.65	0.42
1:B:488:ASN:N	1:B:488:ASN:ND2	2.62	0.42
1:A:581:LEU:HD12	1:A:581:LEU:HA	1.75	0.42
1:B:167:HIS:HA	1:B:246:ASN:O	2.19	0.42
1:B:183:LEU:HD12	1:B:183:LEU:C	2.44	0.42
1:B:596:LYS:HE3	1:B:596:LYS:HB2	1.81	0.42
1:A:401:PHE:CZ	1:A:403:PHE:HB2	2.54	0.42
1:B:195:ASN:HB2	1:B:198:GLN:OE1	2.19	0.42
1:A:298:LEU:HA	1:A:339:ILE:HD11	2.01	0.42
1:B:160:PHE:CD2	1:B:160:PHE:C	2.98	0.42
1:B:455:GLN:OE1	1:B:462:HIS:HB3	2.19	0.42
1:B:156:ASP:O	1:B:159:ALA:HB3	2.19	0.42
1:B:231:VAL:HG13	1:B:232:THR:N	2.29	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:596:LYS:HG2	1:B:601:ILE:HD13	2.02	0.42
1:A:273:ILE:HD11	1:A:381:LEU:HD21	2.01	0.42
1:B:188:LEU:HD12	1:B:219:ILE:HG12	2.01	0.42
1:B:298:LEU:HA	1:B:339:ILE:CD1	2.50	0.42
1:B:516:ASN:HD21	1:B:599:ARG:HH21	1.67	0.42
1:A:298:LEU:HA	1:A:339:ILE:CD1	2.50	0.42
1:A:504:GLN:HG2	1:A:577:TRP:HH2	1.85	0.41
1:B:388:ILE:O	1:B:389:SER:C	2.62	0.41
1:B:402:LEU:HD23	1:B:402:LEU:HA	1.85	0.41
1:A:231:VAL:HG13	1:A:232:THR:N	2.28	0.41
1:A:556:ILE:O	1:A:556:ILE:HG23	2.20	0.41
1:B:188:LEU:CD1	1:B:219:ILE:HG12	2.50	0.41
1:A:160:PHE:CD2	1:A:160:PHE:C	2.98	0.41
1:A:452:VAL:HG11	1:A:465:THR:HG23	2.02	0.41
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.67	0.41
1:A:331:LYS:C	1:A:334:PRO:HD2	2.45	0.41
1:A:557:ARG:NH1	1:B:396:ASN:HD22	2.18	0.41
1:A:588:ARG:NH1	1:A:588:ARG:CG	2.82	0.41
1:B:195:ASN:N	1:B:198:GLN:OE1	2.45	0.41
1:A:404:SER:O	1:A:423:GLY:HA3	2.20	0.41
1:B:324:ARG:O	1:B:325:PHE:C	2.64	0.41
1:B:357:GLU:OE1	1:B:368:VAL:HG21	2.21	0.41
1:B:392:ASN:N	1:B:392:ASN:ND2	2.69	0.41
1:B:452:VAL:HG11	1:B:465:THR:HG23	2.03	0.41
1:B:555:LYS:HB3	1:B:557:ARG:HH21	1.86	0.41
1:A:316:ASP:C	1:A:318:VAL:H	2.27	0.41
1:A:360:TYR:C	1:A:360:TYR:CD2	2.99	0.41
1:A:596:LYS:HE3	1:A:596:LYS:HB2	1.87	0.41
1:B:150:LYS:HA	1:B:151:PRO:HD3	1.88	0.41
1:B:217:ALA:HA	1:B:219:ILE:HG22	2.02	0.41
1:B:258:PHE:CD2	1:B:260:PRO:HD2	2.56	0.41
1:B:316:ASP:C	1:B:318:VAL:H	2.29	0.41
1:B:466:ASP:C	1:B:467:VAL:HG23	2.46	0.41
1:A:175:HIS:O	1:A:180:LYS:HE2	2.21	0.41
1:B:218:TRP:CD1	1:B:218:TRP:O	2.73	0.40
1:B:266:ILE:HG23	1:B:305:ILE:HD12	2.03	0.40
1:B:388:ILE:O	1:B:390:LYS:N	2.54	0.40
1:B:409:ILE:HA	1:B:410:PRO:HD3	1.73	0.40
1:B:504:GLN:HG2	1:B:577:TRP:HH2	1.86	0.40
1:A:165:LEU:HA	1:A:166:PRO:HD3	1.91	0.40
1:A:188:LEU:HB3	1:A:194:VAL:HG23	2.02	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:MSE:O	1:A:316:ASP:C	2.61	0.40
1:A:538:ILE:HG21	1:A:541:LEU:CD1	2.51	0.40
1:B:318:VAL:O	1:B:319:ASP:C	2.64	0.40
1:B:333:LEU:O	1:B:333:LEU:HD12	2.21	0.40
1:B:437:THR:O	1:B:492:ALA:HA	2.21	0.40
1:B:599:ARG:O	1:B:601:ILE:HG23	2.21	0.40
1:A:314:LYS:HE2	2:A:663:GDP:O4'	2.21	0.40
1:A:157:ILE:HD13	1:A:157:ILE:HA	1.85	0.40
1:A:369:ARG:NH1	1:A:373:ASN:HA	2.36	0.40
1:A:555:LYS:HB3	1:A:557:ARG:HH21	1.87	0.40
1:A:585:GLU:OE1	1:B:540:ARG:NH2	2.54	0.40
1:B:175:HIS:O	1:B:180:LYS:HE2	2.21	0.40
1:B:187:LEU:HD12	1:B:187:LEU:HA	1.66	0.40
1:B:376:ASN:O	1:B:377:LEU:C	2.64	0.40
1:B:525:PHE:CE2	1:B:594:LEU:HG	2.56	0.40
1:A:312:MSE:HE1	1:A:328:ILE:CG2	2.50	0.40
1:B:384:ALA:O	1:B:387:LYS:HB2	2.21	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	419/483 (87%)	383 (91%)	30 (7%)	6 (1%)	9	24
1	B	434/483 (90%)	395 (91%)	31 (7%)	8 (2%)	6	19
All	All	853/966 (88%)	778 (91%)	61 (7%)	14 (2%)	7	21

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	231	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	546	ASP
1	B	546	ASP
1	A	455	GLN
1	B	389	SER
1	B	455	GLN
1	B	456	GLN
1	B	543	SER
1	A	395	ILE
1	A	543	SER
1	B	395	ILE
1	B	548	GLY
1	A	389	SER
1	A	548	GLY

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	383/417 (92%)	354 (92%)	29 (8%)	12	31
1	B	391/417 (94%)	358 (92%)	33 (8%)	10	26
All	All	774/834 (93%)	712 (92%)	62 (8%)	11	28

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	VAL
1	A	204	ARG
1	A	219	ILE
1	A	221	ASP
1	A	234	SER
1	A	249	ILE
1	A	280	THR
1	A	289	LEU
1	A	293	THR
1	A	316	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	323	GLN
1	A	339	ILE
1	A	344	ASP
1	A	362	ILE
1	A	379	SER
1	A	412	LYS
1	A	447	VAL
1	A	454	SER
1	A	469	ILE
1	A	484	GLU
1	A	488	ASN
1	A	533	GLU
1	A	541	LEU
1	A	547	LYS
1	A	567	VAL
1	A	578	ILE
1	A	584	HIS
1	A	588	ARG
1	A	611	GLN
1	B	139	VAL
1	B	204	ARG
1	B	218	TRP
1	B	219	ILE
1	B	234	SER
1	B	249	ILE
1	B	250	VAL
1	B	256	ARG
1	B	274	LEU
1	B	280	THR
1	B	289	LEU
1	B	293	THR
1	B	316	ASP
1	B	323	GLN
1	B	339	ILE
1	B	344	ASP
1	B	362	ILE
1	B	379	SER
1	B	422	SER
1	B	447	VAL
1	B	454	SER
1	B	469	ILE
1	B	484	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	488	ASN
1	B	511	THR
1	B	533	GLU
1	B	541	LEU
1	B	547	LYS
1	B	567	VAL
1	B	578	ILE
1	B	584	HIS
1	B	588	ARG
1	B	611	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	HIS
1	A	167	HIS
1	A	192	ASN
1	A	195	ASN
1	A	222	GLN
1	A	243	HIS
1	A	323	GLN
1	A	347	ASN
1	A	383	ASN
1	A	392	ASN
1	A	487	GLN
1	A	488	ASN
1	A	501	HIS
1	A	516	ASN
1	A	534	GLN
1	B	155	HIS
1	B	167	HIS
1	B	192	ASN
1	B	195	ASN
1	B	222	GLN
1	B	243	HIS
1	B	261	ASN
1	B	323	GLN
1	B	347	ASN
1	B	383	ASN
1	B	392	ASN
1	B	451	GLN
1	B	456	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	487	GLN
1	B	488	ASN
1	B	501	HIS
1	B	534	GLN
1	B	611	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	GDP	A	663	-	29,30,30	1.13	2 (6%)	45,47,47	1.83	8 (17%)
2	GDP	B	663	-	29,30,30	1.11	2 (6%)	45,47,47	1.87	9 (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
2	GDP	A	663	-	-	4/16/32/32	0/3/3/3
2	GDP	B	663	-	-	5/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	663	GDP	C5-C4	3.19	1.47	1.38
2	B	663	GDP	C5-C4	3.14	1.47	1.38
2	A	663	GDP	C6-N1	-2.50	1.34	1.38
2	B	663	GDP	C6-N1	-2.46	1.34	1.38

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	663	GDP	C5-C4-N3	-6.47	118.09	128.39
2	A	663	GDP	C5-C4-N3	-6.31	118.34	128.39
2	A	663	GDP	C2-N3-C4	5.31	121.45	112.30
2	B	663	GDP	C2-N3-C4	5.26	121.35	112.30
2	B	663	GDP	N9-C4-N3	4.71	135.38	125.95
2	A	663	GDP	N9-C4-N3	4.51	134.97	125.95
2	A	663	GDP	C6-C5-N7	3.55	136.75	130.29
2	B	663	GDP	C6-C5-N7	3.22	136.15	130.29
2	A	663	GDP	C4-C5-N7	-2.62	106.51	110.67
2	B	663	GDP	C4-C5-N7	-2.56	106.61	110.67
2	A	663	GDP	C5-C6-N1	2.22	118.91	113.25
2	B	663	GDP	O2A-PA-O3A	2.21	113.26	107.27
2	B	663	GDP	O6-C6-C5	-2.19	120.74	126.53
2	B	663	GDP	C5-C6-N1	2.10	118.59	113.25
2	A	663	GDP	O6-C6-C5	-2.09	121.01	126.53
2	A	663	GDP	O5'-PA-O1A	2.07	117.15	108.94
2	B	663	GDP	O3B-PB-O3A	2.05	111.52	104.64

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	663	GDP	PA-O3A-PB-O3B
2	A	663	GDP	O4'-C4'-C5'-O5'
2	B	663	GDP	O4'-C4'-C5'-O5'
2	B	663	GDP	PA-O3A-PB-O1B
2	A	663	GDP	C3'-C4'-C5'-O5'
2	A	663	GDP	PB-O3A-PA-O2A

Continued on next page...

Continued from previous page...

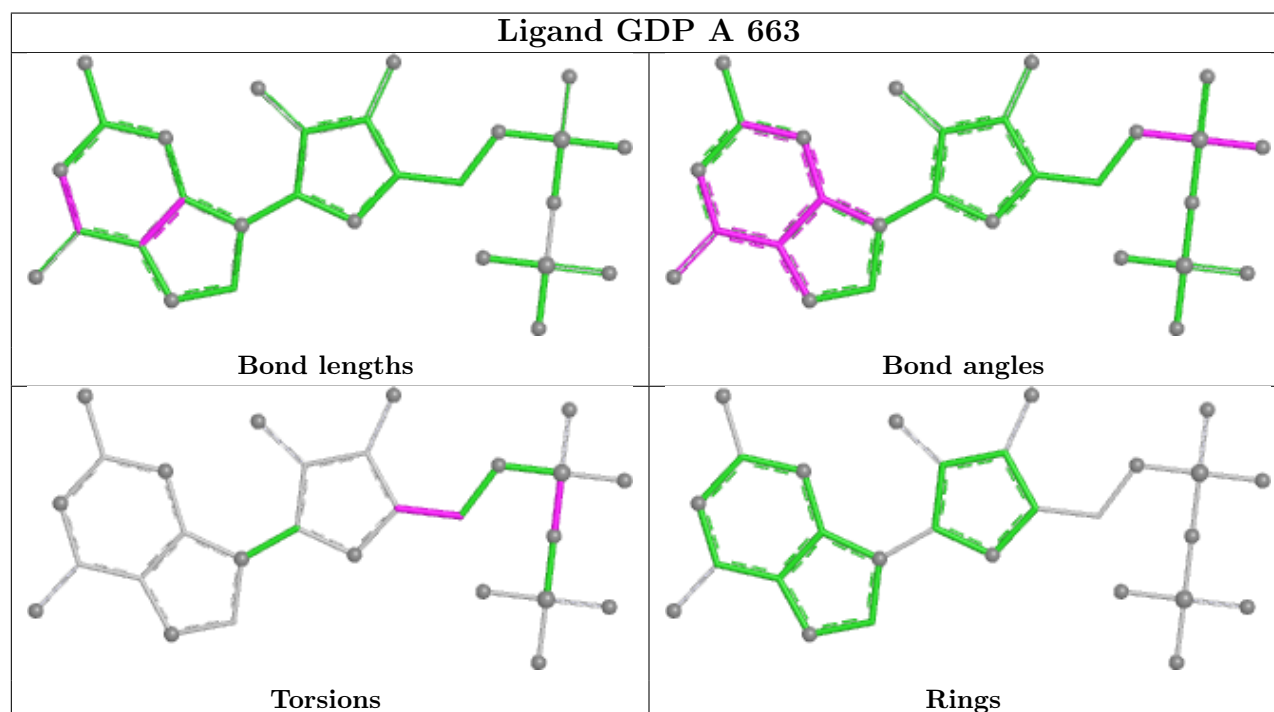
Mol	Chain	Res	Type	Atoms
2	B	663	GDP	PB-O3A-PA-O1A
2	B	663	GDP	C3'-C4'-C5'-O5'
2	A	663	GDP	PB-O3A-PA-O1A

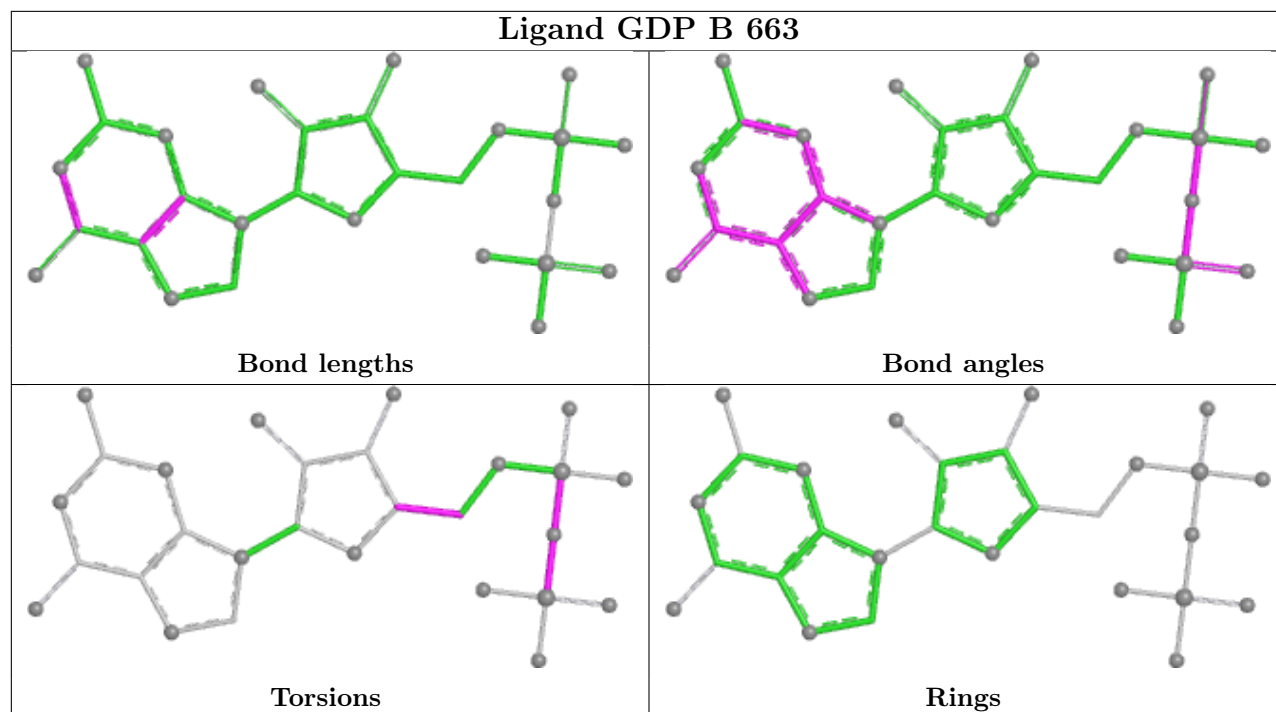
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	663	GDP	4	0
2	B	663	GDP	3	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	424/483 (87%)	-0.41	5 (1%) 76 71	23, 47, 83, 132	0
1	B	435/483 (90%)	-0.39	5 (1%) 78 73	24, 47, 91, 122	0
All	All	859/966 (88%)	-0.40	10 (1%) 76 71	23, 47, 88, 132	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	217	ALA	3.7
1	B	549	ASN	2.6
1	B	288	ASP	2.5
1	A	218	TRP	2.3
1	A	254	GLY	2.2
1	A	288	ASP	2.1
1	B	219	ILE	2.1
1	A	223	THR	2.1
1	A	231	VAL	2.1
1	B	410	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

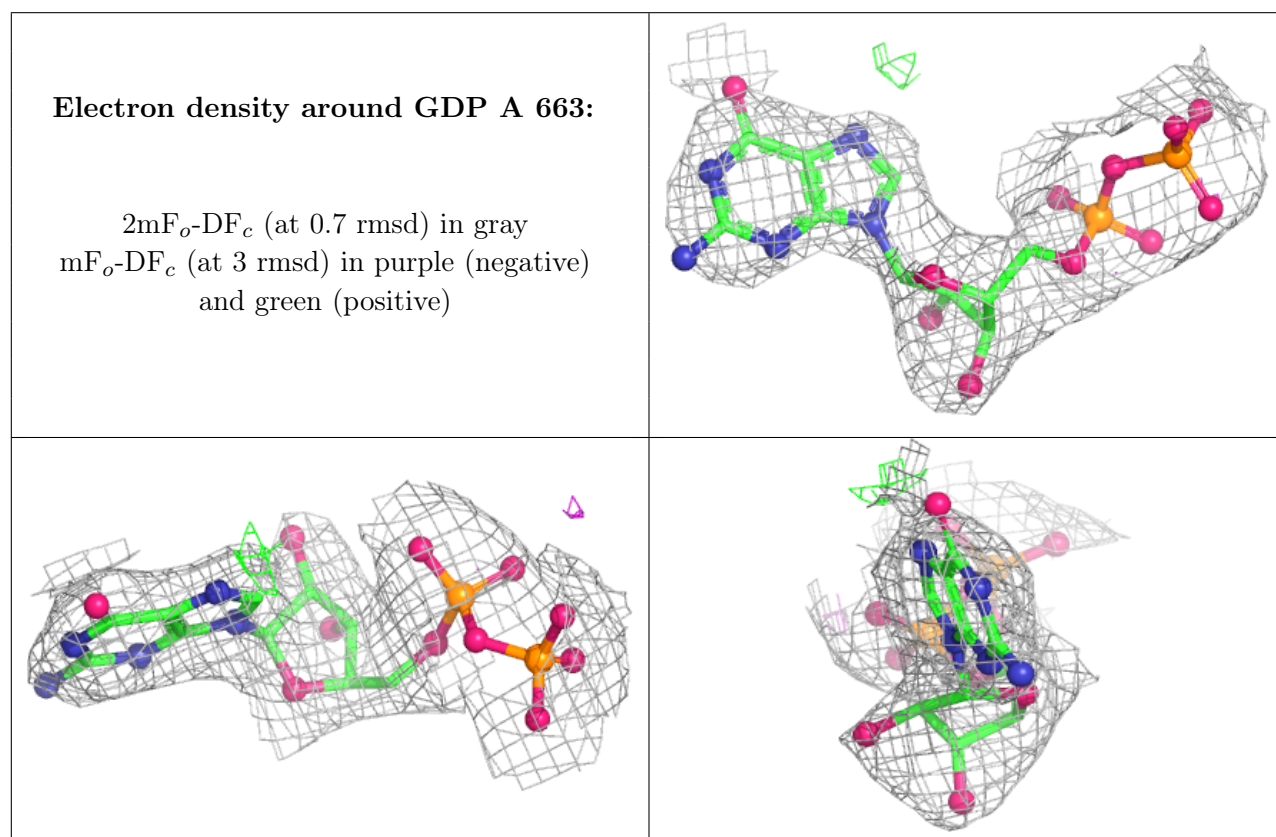
There are no oligosaccharides in this entry.

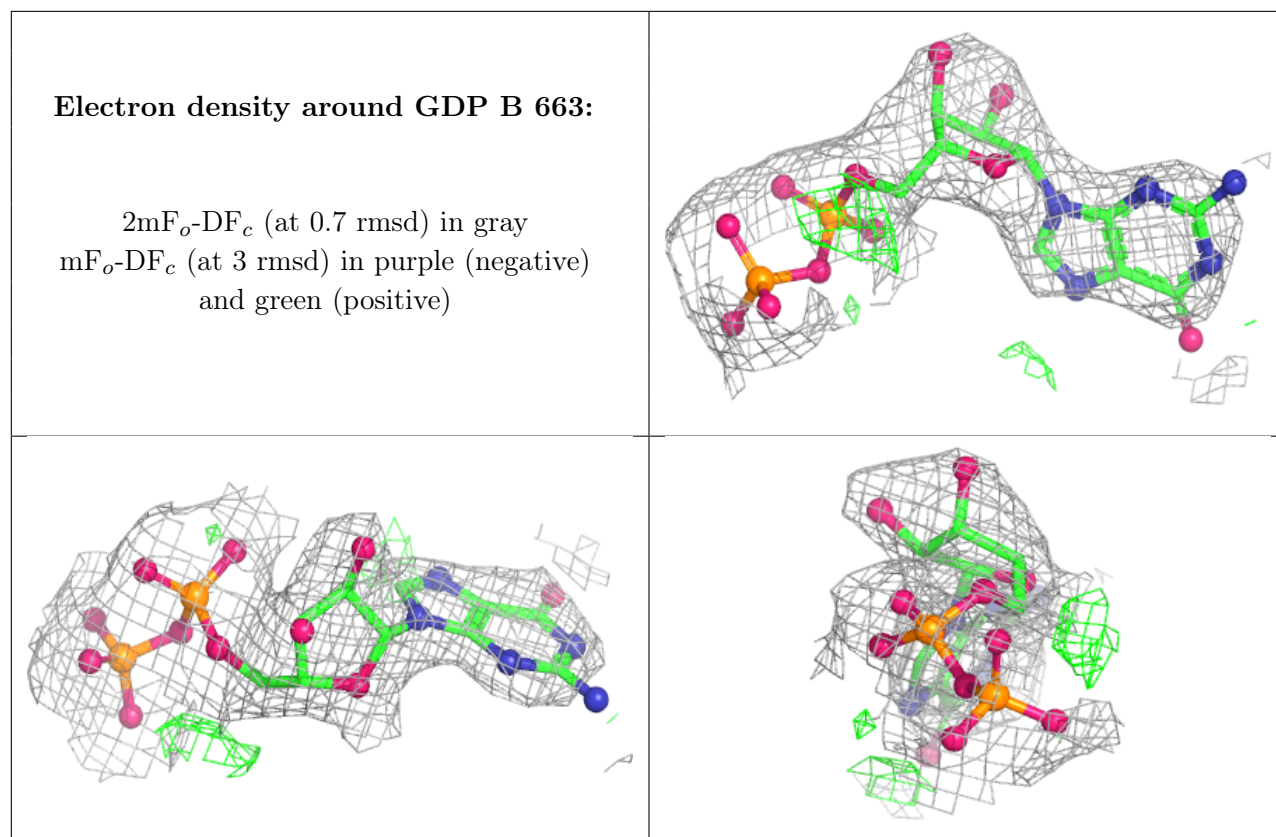
6.4 Ligands ⓘ

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	GDP	A	663	28/28	0.90	0.09	57,77,84,90	0
2	GDP	B	663	28/28	0.95	0.07	56,76,85,86	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.