



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 12:35 AM UTC

PDB ID : 2QN6 / pdb_00002qn6
Title : Structure of an archaeal heterotrimeric initiation factor 2 reveals a nucleotide state between the GTP and the GDP states
Authors : Mechulam, Y.; Yatime, L.; Blanquet, S.; Schmitt, E.
Deposited on : 2007-07-18
Resolution : 2.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

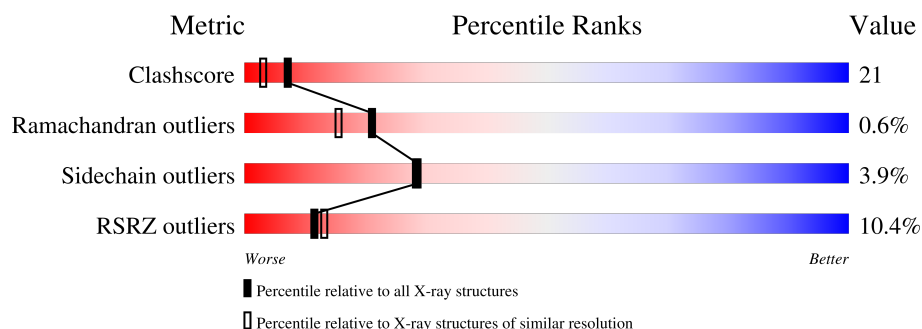
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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))
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	414	10% 59% 34% • 5%
2	B	93	2% 76% 18% • •
3	C	18	39% 50% 50%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4177 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 Translation initiation factor 2 gamma subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			3044	1953	516	565	10			

- Molecule 2 is a protein called Translation initiation factor 2 alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	90	Total	C	N	O	S	0	0	0
			698	442	117	138	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	174	MET	-	expression tag	UNP Q97Z79

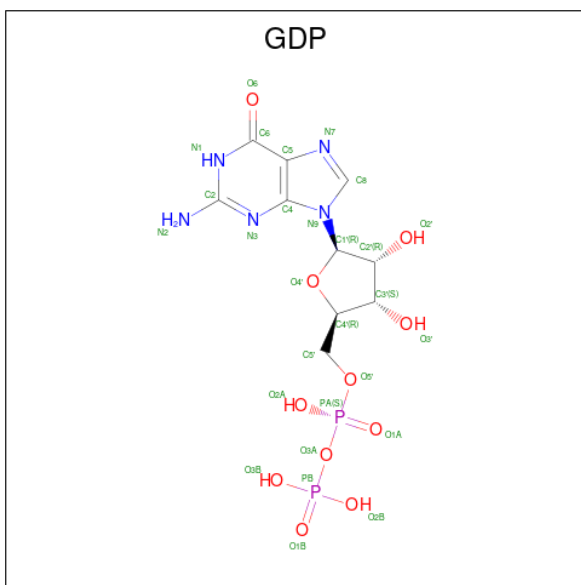
- Molecule 3 is a protein called Translation initiation factor 2 beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	18	Total	C	N	O	S	0	0	0
			153	97	23	32	1			

- 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		

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



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

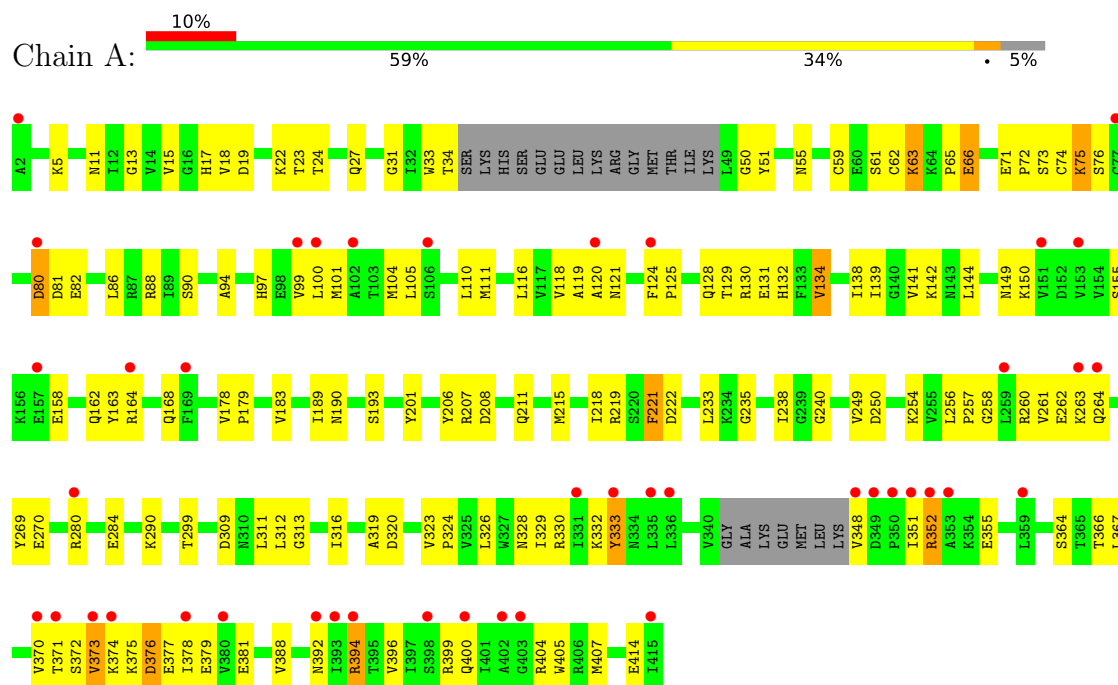
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	201	Total O 201 201	0	0
6	B	50	Total O 50 50	0	0
6	C	2	Total O 2 2	0	0

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: Translation initiation factor 2 gamma subunit



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	122.12Å 51.85Å 98.03Å 90.00° 94.42° 90.00°	Depositor
Resolution (Å)	100.00 – 2.15 97.74 – 2.15	Depositor EDS
% Data completeness (in resolution range)	79.2 (100.00-2.15) 88.1 (97.74-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.14Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.264 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.607	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	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	0.94	EDS
Total number of atoms	4177	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GDP, MG

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	0.42	0/3100	0.94	11/4204 (0.3%)
2	B	0.49	0/703	0.92	1/949 (0.1%)
3	C	0.39	0/155	0.84	0/205
All	All	0.43	0/3958	0.94	12/5358 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	332	LYS	N-CA-C	-6.66	99.80	109.59
1	A	222	ASP	N-CA-C	-5.91	97.69	108.02
1	A	333	TYR	N-CA-C	5.55	118.30	109.81
1	A	31	GLY	N-CA-C	-5.46	106.67	115.08
1	A	142	LYS	N-CA-C	5.42	117.27	111.36
1	A	71	GLU	CA-C-N	5.31	126.48	119.84
1	A	71	GLU	C-N-CA	5.31	126.48	119.84
1	A	280	ARG	N-CA-C	5.29	118.04	109.06
1	A	206	TYR	N-CA-C	-5.23	102.55	110.28
1	A	62	CYS	N-CA-C	5.12	116.54	110.19
1	A	326	LEU	N-CA-C	5.01	117.57	109.40
2	B	188	ASN	N-CA-C	-5.01	106.59	113.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3044	0	3141	147	0
2	B	698	0	736	16	0
3	C	153	0	151	10	0
4	A	1	0	0	0	0
5	A	28	0	12	1	0
6	A	201	0	0	10	0
6	B	50	0	0	2	0
6	C	2	0	0	1	0
All	All	4177	0	4040	164	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 21.

All (164) 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:394:ARG:HB2	1:A:394:ARG:HH11	1.18	1.06
1:A:218:ILE:HG12	1:A:240:GLY:HA2	1.53	0.89
1:A:394:ARG:HH11	1:A:394:ARG:CB	1.84	0.89
1:A:63:LYS:HD2	1:A:63:LYS:H	1.43	0.84
1:A:59:CYS:HG	1:A:74:CYS:HG	1.23	0.83
1:A:13:GLY:HA3	1:A:111:MET:HE3	1.62	0.81
1:A:139:ILE:HB	1:A:394:ARG:HH21	1.49	0.77
2:B:184:THR:HG22	6:B:557:HOH:O	1.85	0.77
1:A:75:LYS:HD2	1:A:76:SER:N	2.03	0.73
1:A:139:ILE:HG13	1:A:141:VAL:HG23	1.69	0.72
1:A:15:VAL:C	1:A:22:LYS:HD3	2.14	0.72
1:A:261:VAL:CG1	1:A:263:LYS:HE2	2.19	0.72
1:A:139:ILE:CB	1:A:394:ARG:HH21	2.01	0.71
1:A:63:LYS:H	1:A:63:LYS:CD	2.02	0.71
1:A:215:MET:HE3	1:A:316:ILE:HD12	1.73	0.71
1:A:189:ILE:HD12	3:C:18:LEU:HD23	1.73	0.70
1:A:104:MET:HB2	1:A:139:ILE:HD13	1.73	0.70
1:A:139:ILE:CG2	1:A:394:ARG:HH21	2.05	0.69
1:A:101:MET:HE1	1:A:132:HIS:HD2	1.57	0.68
1:A:351:ILE:HG22	6:A:697:HOH:O	1.91	0.68
1:A:218:ILE:O	1:A:312:LEU:HD23	1.93	0.68
1:A:94:ALA:HB3	1:A:100:LEU:HG	1.76	0.68
1:A:34:THR:HG23	1:A:50:GLY:HA3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:GLN:HG3	1:A:33:TRP:CE2	2.29	0.67
1:A:392:ASN:HD21	1:A:414:GLU:HG2	1.59	0.67
2:B:177:VAL:HG21	2:B:236:PRO:HG3	1.76	0.66
1:A:249:VAL:O	1:A:250:ASP:HB2	1.96	0.65
1:A:131:GLU:O	1:A:134:VAL:HG22	1.96	0.64
1:A:97:HIS:CE1	1:A:128:GLN:HG2	2.32	0.64
1:A:190:ASN:ND2	3:C:13:ARG:HH11	1.96	0.62
1:A:254:LYS:HE3	1:A:319:ALA:O	1.99	0.62
1:A:352:ARG:O	1:A:355:GLU:HB2	1.98	0.62
1:A:73:SER:OG	1:A:75:LYS:HG3	2.00	0.62
1:A:17:HIS:CD2	1:A:128:GLN:HB2	2.35	0.62
1:A:211:GLN:HG2	6:A:524:HOH:O	2.00	0.61
1:A:11:ASN:HD22	1:A:110:LEU:HB3	1.65	0.61
1:A:260:ARG:NH2	1:A:262:GLU:OE1	2.34	0.61
1:A:263:LYS:O	1:A:264:GLN:HB2	2.01	0.60
1:A:394:ARG:CB	1:A:394:ARG:NH1	2.61	0.59
1:A:351:ILE:HD11	1:A:370:VAL:HG21	1.85	0.59
1:A:34:THR:CG2	1:A:50:GLY:HA3	2.34	0.58
1:A:219:ARG:HE	1:A:309:ASP:CG	2.10	0.58
1:A:11:ASN:ND2	1:A:110:LEU:HB3	2.19	0.57
1:A:233:LEU:HD23	2:B:193:VAL:CG2	2.35	0.57
1:A:179:PRO:HG3	1:A:201:TYR:CD2	2.40	0.57
1:A:189:ILE:CD1	3:C:18:LEU:HD23	2.35	0.57
1:A:193:SER:HB3	3:C:10:MET:HE2	1.86	0.57
1:A:375:LYS:HA	6:A:752:HOH:O	2.04	0.57
1:A:97:HIS:ND1	1:A:128:GLN:HG2	2.20	0.56
2:B:178:LYS:O	2:B:179:MET:HE2	2.06	0.56
1:A:75:LYS:HD2	1:A:75:LYS:C	2.31	0.56
1:A:5:LYS:NZ	1:A:284:GLU:OE1	2.38	0.55
2:B:219:ILE:HG12	2:B:229:VAL:HG22	1.88	0.55
1:A:324:PRO:HB2	1:A:388:VAL:CG2	2.37	0.55
1:A:120:ALA:HA	1:A:162:GLN:OE1	2.07	0.55
1:A:155:SER:OG	1:A:158:GLU:HG3	2.07	0.54
1:A:94:ALA:H	1:A:100:LEU:HD21	1.72	0.54
1:A:371:THR:OG1	1:A:379:GLU:HB2	2.08	0.54
1:A:352:ARG:NH1	1:A:355:GLU:CD	2.67	0.53
1:A:320:ASP:HB3	6:A:595:HOH:O	2.07	0.53
1:A:374:LYS:HE2	6:A:649:HOH:O	2.08	0.53
1:A:262:GLU:O	1:A:263:LYS:HD3	2.09	0.53
1:A:101:MET:HB3	1:A:407:MET:HE1	1.92	0.52
1:A:66:GLU:HG3	3:C:13:ARG:NH2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:HIS:CG	1:A:128:GLN:HG2	2.45	0.52
1:A:105:LEU:HB3	1:A:396:VAL:HG22	1.91	0.52
1:A:330:ARG:NH1	1:A:379:GLU:OE2	2.41	0.52
2:B:194:GLU:HG3	6:B:666:HOH:O	2.09	0.51
1:A:163:TYR:CE2	3:C:4:GLU:HG3	2.45	0.51
1:A:333:TYR:CE2	1:A:378:ILE:HG23	2.46	0.51
2:B:182:LEU:O	2:B:264:VAL:HG22	2.11	0.51
1:A:313:GLY:O	1:A:364:SER:HB3	2.11	0.51
1:A:55:ASN:ND2	1:A:88:ARG:HH11	2.09	0.51
1:A:18:VAL:O	1:A:19:ASP:HB2	2.11	0.50
1:A:15:VAL:O	1:A:22:LYS:HD3	2.12	0.50
1:A:348:VAL:HG13	1:A:348:VAL:O	2.11	0.50
1:A:388:VAL:HG23	1:A:388:VAL:O	2.11	0.50
1:A:392:ASN:ND2	1:A:414:GLU:HG2	2.25	0.50
1:A:104:MET:HB2	1:A:139:ILE:CD1	2.42	0.50
1:A:235:GLY:HA3	1:A:299:THR:O	2.11	0.50
1:A:376:ASP:O	1:A:376:ASP:CG	2.55	0.50
3:C:4:GLU:O	3:C:8:VAL:HG23	2.12	0.49
1:A:101:MET:CE	1:A:132:HIS:HD2	2.23	0.49
1:A:139:ILE:HG22	1:A:394:ARG:HH21	1.75	0.49
1:A:65:PRO:HG2	1:A:193:SER:OG	2.11	0.49
1:A:328:ASN:OD1	1:A:381:GLU:HA	2.12	0.49
2:B:182:LEU:HG	2:B:264:VAL:HG21	1.95	0.49
1:A:94:ALA:HB3	1:A:100:LEU:CD2	2.43	0.49
1:A:261:VAL:HG11	1:A:263:LYS:HE2	1.93	0.49
1:A:183:VAL:HG12	1:A:190:ASN:HB3	1.95	0.49
2:B:242:ALA:O	2:B:246:ILE:HG13	2.13	0.49
1:A:189:ILE:O	1:A:190:ASN:HB2	2.13	0.49
1:A:374:LYS:HG3	1:A:375:LYS:H	1.78	0.49
1:A:101:MET:O	1:A:105:LEU:HB2	2.13	0.48
1:A:164:ARG:O	1:A:168:GLN:HB2	2.12	0.48
1:A:99:VAL:HG23	1:A:100:LEU:N	2.28	0.48
1:A:63:LYS:CD	1:A:63:LYS:N	2.72	0.48
1:A:116:LEU:C	1:A:116:LEU:HD23	2.39	0.48
1:A:139:ILE:CG2	1:A:394:ARG:NH2	2.76	0.48
1:A:351:ILE:HD13	1:A:378:ILE:HG21	1.94	0.48
1:A:63:LYS:H	1:A:63:LYS:CE	2.27	0.47
1:A:17:HIS:HD2	1:A:128:GLN:H	1.62	0.47
1:A:260:ARG:HD2	1:A:269:TYR:OH	2.14	0.47
1:A:193:SER:HB3	3:C:10:MET:CE	2.44	0.47
1:A:27:GLN:HG3	1:A:33:TRP:CD2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:ILE:HA	6:A:690:HOH:O	2.14	0.47
1:A:59:CYS:CB	1:A:74:CYS:HG	2.29	0.46
1:A:15:VAL:HG11	1:A:104:MET:HE1	1.97	0.46
1:A:366:THR:HG22	1:A:367:LEU:O	2.15	0.46
1:A:134:VAL:O	1:A:138:ILE:HG13	2.16	0.46
1:A:324:PRO:HD2	1:A:388:VAL:O	2.15	0.46
1:A:23:THR:HG23	1:A:34:THR:HB	1.97	0.46
1:A:94:ALA:HB3	1:A:100:LEU:CG	2.42	0.46
1:A:370:VAL:HG11	1:A:373:VAL:HG22	1.98	0.46
1:A:330:ARG:HH11	1:A:330:ARG:HG3	1.80	0.46
2:B:180:SER:HA	2:B:229:VAL:O	2.17	0.45
1:A:139:ILE:HG22	1:A:394:ARG:NH2	2.31	0.45
1:A:144:LEU:O	1:A:178:VAL:HG13	2.16	0.45
1:A:101:MET:HE1	1:A:132:HIS:CD2	2.46	0.45
1:A:94:ALA:N	1:A:100:LEU:HD21	2.32	0.44
1:A:63:LYS:HG2	1:A:66:GLU:OE2	2.17	0.44
1:A:81:ASP:OD1	1:A:82:GLU:N	2.51	0.44
1:A:218:ILE:O	1:A:219:ARG:HD2	2.16	0.44
1:A:221:PHE:N	1:A:221:PHE:CD1	2.85	0.44
1:A:17:HIS:CD2	1:A:128:GLN:H	2.35	0.44
1:A:130:ARG:O	1:A:134:VAL:HG13	2.17	0.44
1:A:139:ILE:HB	1:A:394:ARG:NH2	2.25	0.44
1:A:238:ILE:HD11	1:A:311:LEU:HD12	1.99	0.44
1:A:330:ARG:NH1	1:A:379:GLU:CD	2.75	0.44
1:A:86:LEU:HB3	6:A:576:HOH:O	2.17	0.44
1:A:290:LYS:HD2	6:A:582:HOH:O	2.17	0.43
1:A:375:LYS:C	1:A:377:GLU:N	2.74	0.43
1:A:256:LEU:HA	1:A:257:PRO:C	2.42	0.43
1:A:116:LEU:HD22	1:A:118:VAL:HG23	2.00	0.43
1:A:233:LEU:HD23	2:B:193:VAL:HG23	2.01	0.43
3:C:14:LEU:C	3:C:14:LEU:HD23	2.43	0.43
1:A:352:ARG:NH1	1:A:355:GLU:OE1	2.52	0.43
1:A:400:GLN:HA	1:A:405:TRP:HA	2.00	0.43
1:A:262:GLU:C	1:A:263:LYS:HD3	2.44	0.43
2:B:182:LEU:HB3	2:B:264:VAL:CG2	2.49	0.43
1:A:17:HIS:HB2	1:A:129:THR:OG1	2.19	0.42
1:A:208:ASP:HB3	1:A:211:GLN:HE21	1.84	0.42
2:B:177:VAL:HG12	2:B:179:MET:HE3	2.01	0.42
3:C:9:GLU:HG3	6:C:713:HOH:O	2.18	0.42
1:A:149:ASN:ND2	1:A:150:LYS:HG2	2.34	0.42
1:A:51:TYR:CE2	1:A:90:SER:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:VAL:HG21	6:A:692:HOH:O	2.19	0.41
1:A:312:LEU:C	1:A:312:LEU:HD13	2.45	0.41
1:A:80:ASP:HA	6:A:654:HOH:O	2.19	0.41
1:A:351:ILE:O	1:A:351:ILE:HG23	2.21	0.41
2:B:237:LYS:O	2:B:241:GLU:HG3	2.21	0.41
1:A:258:GLY:HA3	1:A:270:GLU:O	2.21	0.41
1:A:323:VAL:HA	1:A:324:PRO:HD3	1.89	0.41
1:A:352:ARG:HG2	1:A:352:ARG:HH11	1.85	0.41
2:B:182:LEU:CG	2:B:264:VAL:HG21	2.51	0.41
1:A:124:PHE:HA	1:A:125:PRO:HA	1.84	0.41
1:A:179:PRO:HG3	1:A:201:TYR:CE2	2.55	0.41
1:A:330:ARG:HD3	1:A:377:GLU:OE2	2.21	0.41
1:A:372:SER:OG	1:A:379:GLU:HG2	2.20	0.40
1:A:24:THR:OG1	5:A:417:GDP:O1A	2.36	0.40
1:A:116:LEU:HB2	1:A:144:LEU:HD11	2.03	0.40
1:A:162:GLN:O	1:A:163:TYR:C	2.65	0.40
1:A:119:ALA:C	1:A:121:ASN:N	2.78	0.40
2:B:208:GLU:CD	2:B:208:GLU:H	2.28	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	387/414 (94%)	365 (94%)	19 (5%)	3 (1%)	16	10
2	B	88/93 (95%)	86 (98%)	2 (2%)	0	100	100
3	C	16/18 (89%)	16 (100%)	0	0	100	100
All	All	491/525 (94%)	467 (95%)	21 (4%)	3 (1%)	21	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	PRO
1	A	66	GLU
1	A	373	VAL

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	337/356 (95%)	325 (96%)	12 (4%)	31	31
2	B	80/84 (95%)	77 (96%)	3 (4%)	29	29
3	C	18/18 (100%)	16 (89%)	2 (11%)	6	3
All	All	435/458 (95%)	418 (96%)	17 (4%)	28	28

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	SER
1	A	63	LYS
1	A	75	LYS
1	A	80	ASP
1	A	134	VAL
1	A	207	ARG
1	A	221	PHE
1	A	352	ARG
1	A	376	ASP
1	A	394	ARG
1	A	399	ARG
1	A	404	ARG
2	B	193	VAL
2	B	208	GLU
2	B	215	LEU
3	C	2	SER
3	C	6	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	27	GLN
1	A	55	ASN
1	A	165	GLN
1	A	187	HIS
1	A	190	ASN
1	A	211	GLN
1	A	392	ASN
2	B	206	ASN
2	B	209	GLN
2	B	244	ASN
2	B	249	ASN
2	B	258	ASN

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](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	GDP	A	417	4	29,30,30	1.46	2 (6%)	45,47,47	1.64	8 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GDP	A	417	4	-	2/16/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	417	GDP	PA-O3A	5.85	1.65	1.59
5	A	417	GDP	C6-N1	2.32	1.43	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	417	GDP	O2A-PA-O3A	5.54	122.24	107.27
5	A	417	GDP	O3A-PA-O1A	-4.65	96.72	110.70
5	A	417	GDP	O3B-PB-O2B	-2.94	96.77	107.80
5	A	417	GDP	O3B-PB-O3A	2.82	114.08	104.64
5	A	417	GDP	O3'-C3'-C4'	-2.45	104.05	111.08
5	A	417	GDP	C1'-N9-C8	2.26	133.16	126.73
5	A	417	GDP	O4'-C4'-C5'	2.15	116.22	109.33
5	A	417	GDP	O2B-PB-O1B	2.11	119.07	110.83

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	417	GDP	PB-O3A-PA-O2A
5	A	417	GDP	PB-O3A-PA-O1A

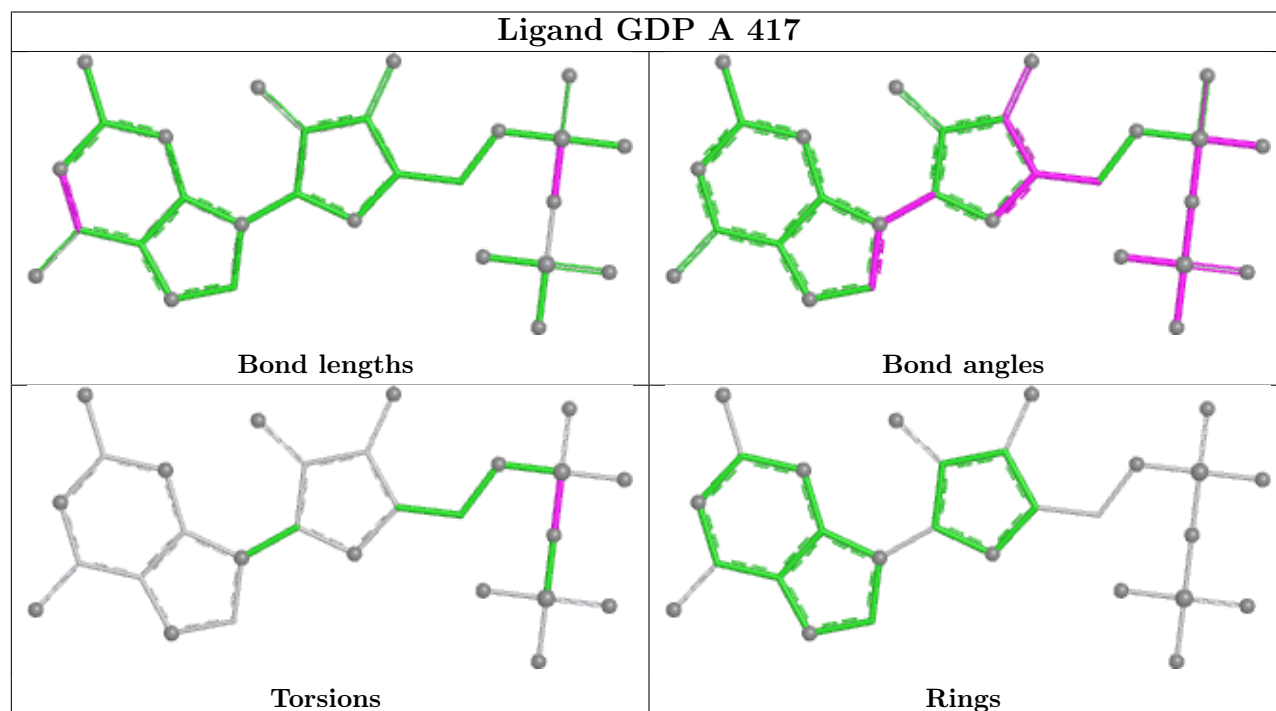
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	417	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	393/414 (94%)	0.78	43 (10%) 10 12	22, 46, 69, 82	0
2	B	90/93 (96%)	0.29	2 (2%) 62 66	26, 36, 50, 67	0
3	C	18/18 (100%)	1.55	7 (38%) 1 1	51, 64, 76, 78	0
All	All	501/525 (95%)	0.72	52 (10%) 11 13	22, 44, 69, 82	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	373	VAL	5.0
1	A	402	ALA	3.7
1	A	336	LEU	3.6
1	A	351	ILE	3.6
2	B	177	VAL	3.5
3	C	12	ASP	3.5
1	A	164	ARG	3.4
1	A	99	VAL	3.4
1	A	393	ILE	3.3
1	A	100	LEU	3.3
1	A	348	VAL	3.1
1	A	264	GLN	3.0
1	A	353	ALA	2.9
1	A	403	GLY	2.9
1	A	392	ASN	2.8
1	A	151	VAL	2.8
1	A	335	LEU	2.7
1	A	374	LYS	2.6
1	A	77	CYS	2.5
1	A	352	ARG	2.5
1	A	102	ALA	2.5
1	A	169	PHE	2.5
1	A	2	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	120	ALA	2.4
1	A	415	ILE	2.4
1	A	80	ASP	2.4
1	A	124	PHE	2.4
1	A	157	GLU	2.3
2	B	265	LYS	2.3
3	C	2	SER	2.3
1	A	359	LEU	2.3
1	A	349	ASP	2.3
1	A	378	ILE	2.3
1	A	371	THR	2.3
3	C	5	LYS	2.3
1	A	331	ILE	2.3
3	C	8	VAL	2.3
1	A	394	ARG	2.3
3	C	14	LEU	2.2
1	A	370	VAL	2.2
1	A	153	VAL	2.2
1	A	106	SER	2.2
1	A	280	ARG	2.1
3	C	19	PRO	2.1
1	A	398	SER	2.1
1	A	350	PRO	2.1
1	A	263	LYS	2.1
3	C	18	LEU	2.1
1	A	380	VAL	2.1
1	A	259	LEU	2.0
1	A	400	GLN	2.0
1	A	333	TYR	2.0

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

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

6.3 Carbohydrates ⓘ

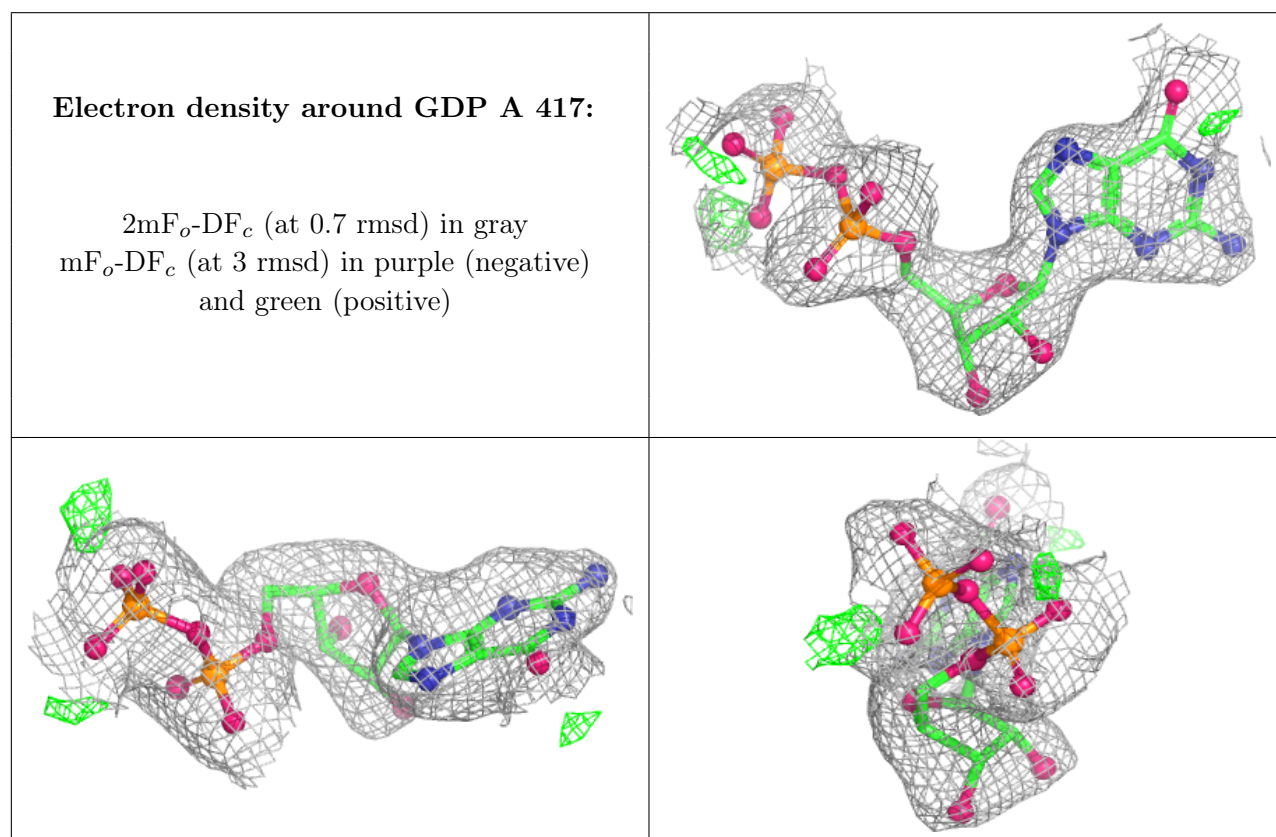
There are no oligosaccharides in this entry.

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
4	MG	A	416	1/1	0.92	0.14	51,51,51,51	0
5	GDP	A	417	28/28	0.92	0.11	49,60,64,65	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.