



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:18 AM UTC

PDB ID : 2X2E / pdb_00002x2e
Title : Dynamin GTPase dimer, long axis form
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Deposited on : 2010-01-12
Resolution : 2.00 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

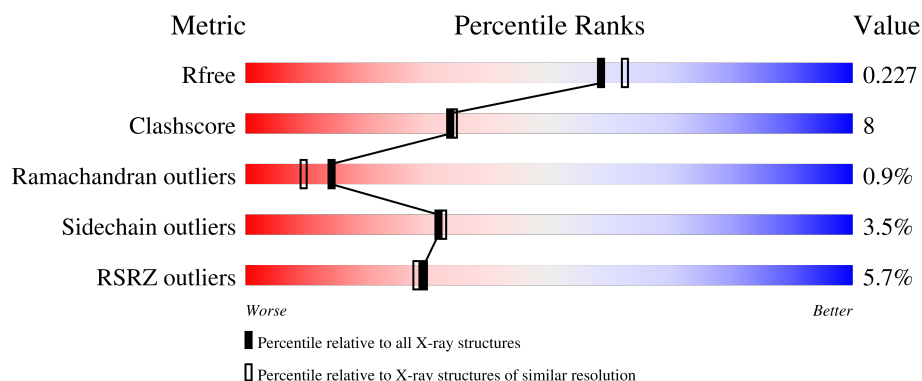
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	353	4% 83% 13% ..
1	D	353	7% 74% 19% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6039 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 DYNAMIN-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	Se	0	0	0
			2675	1682	474	509	2	8			
1	D	341	Total	C	N	O	S	Se	0	0	0
			2675	1682	474	509	2	8			

There are 28 discrepancies between the modelled and reference sequences:

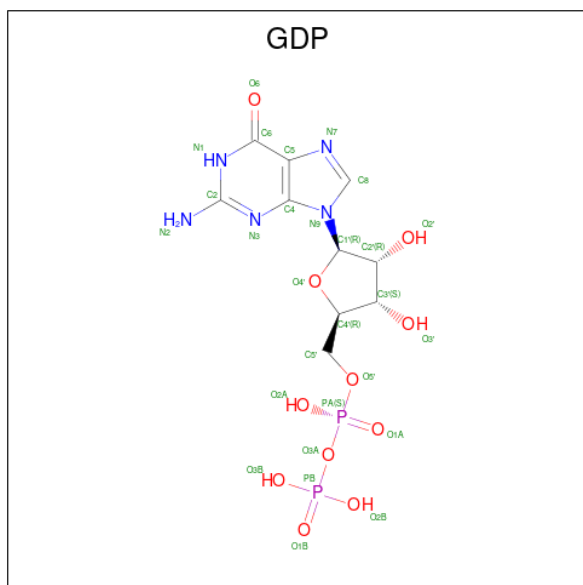
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP Q05193
A	2	PRO	-	cloning artifact	UNP Q05193
A	3	GLU	-	cloning artifact	UNP Q05193
A	4	PHE	-	cloning artifact	UNP Q05193
A	5	SER	-	cloning artifact	UNP Q05193
A	321	LYS	-	linker	UNP Q05193
A	322	HIS	-	linker	UNP Q05193
A	323	GLY	-	linker	UNP Q05193
A	324	THR	-	linker	UNP Q05193
A	325	ASP	-	linker	UNP Q05193
A	326	SER	-	linker	UNP Q05193
A	327	ARG	-	linker	UNP Q05193
A	328	VAL	-	linker	UNP Q05193
A	744	ASN	ASP	variant	UNP Q05193
D	1	GLY	-	cloning artifact	UNP Q05193
D	2	PRO	-	cloning artifact	UNP Q05193
D	3	GLU	-	cloning artifact	UNP Q05193
D	4	PHE	-	cloning artifact	UNP Q05193
D	5	SER	-	cloning artifact	UNP Q05193
D	321	LYS	-	linker	UNP Q05193
D	322	HIS	-	linker	UNP Q05193
D	323	GLY	-	linker	UNP Q05193
D	324	THR	-	linker	UNP Q05193
D	325	ASP	-	linker	UNP Q05193
D	326	SER	-	linker	UNP Q05193

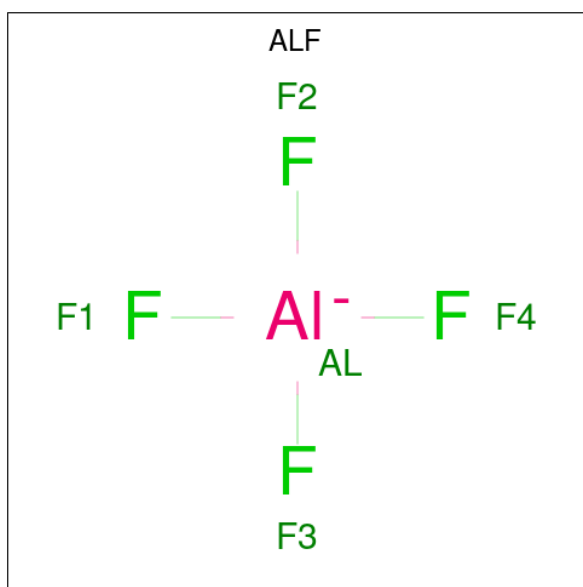
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Chain	Residue	Modelled	Actual	Comment	Reference
D	327	ARG	-	linker	UNP Q05193
D	328	VAL	-	linker	UNP Q05193
D	744	ASN	ASP	variant	UNP Q05193

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Al	F	0	0
			5	1	4		
3	D	1	Total	Al	F	0	0
			5	1	4		

- 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	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		
5	D	1	Total	Na	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	332	Total	O	0	0
			332	332		

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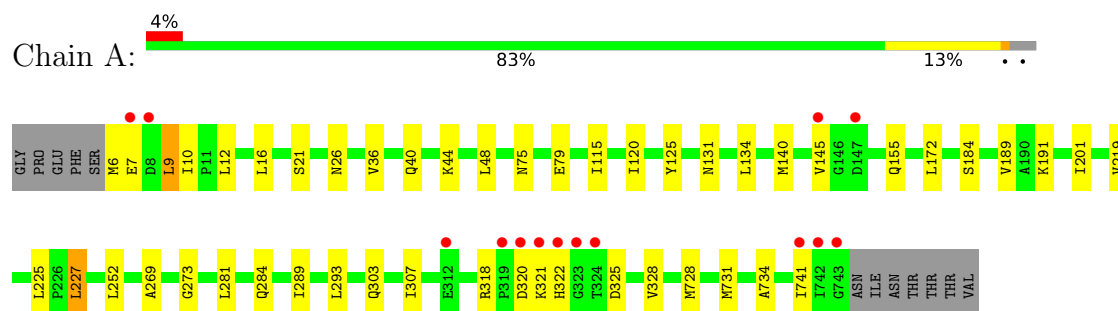
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	287	Total 287	O 287	0	0

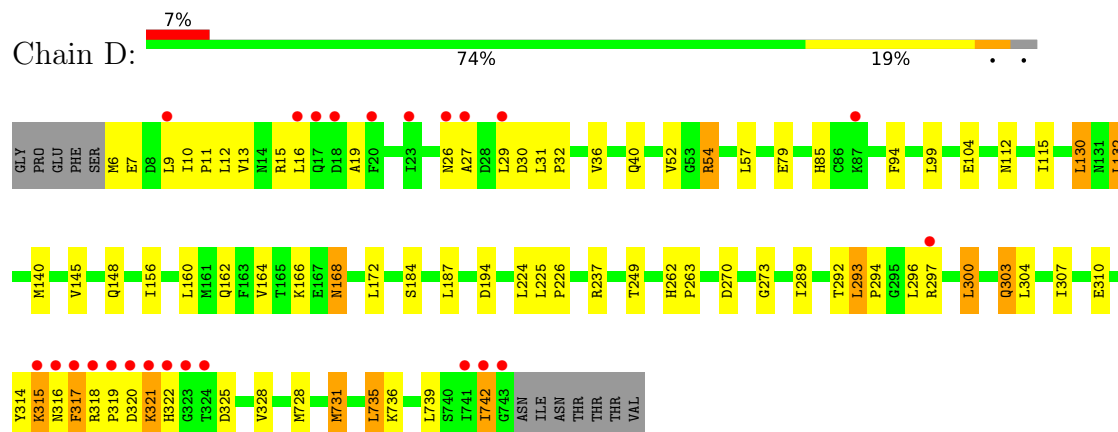
3 Residue-property plots

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: DYNAMIN-1



• Molecule 1: DYNAMIN-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	43.61Å 81.02Å 181.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 30.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.00) 99.8 (30.00-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.19 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.185 , 0.230 0.189 , 0.227	Depositor DCC
R_{free} test set	899 reflections (2.02%)	wwPDB-VP
Wilson B-factor (Å ²)	13.2	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.9	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	6039	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.44% 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

5.1 Standard geometry

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

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	0.68	2/2710 (0.1%)	0.98	7/3654 (0.2%)
1	D	0.63	2/2710 (0.1%)	0.98	11/3654 (0.3%)
All	All	0.65	4/5420 (0.1%)	0.98	18/7308 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	728	MSE	SE-CE	-9.04	1.68	1.95
1	D	728	MSE	SE-CE	-7.89	1.71	1.95
1	D	731	MSE	SE-CE	-6.54	1.75	1.95
1	A	731	MSE	SE-CE	-5.60	1.78	1.95

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	321	LYS	N-CA-C	-11.10	99.05	112.59
1	A	9	LEU	N-CA-C	-10.41	100.58	113.28
1	A	7	GLU	N-CA-C	-8.41	101.15	112.26
1	A	184	SER	N-CA-C	8.15	120.88	109.15
1	A	120	ILE	N-CA-C	-7.60	97.50	108.36
1	D	184	SER	N-CA-C	7.27	119.62	109.15
1	A	269	ALA	N-CA-C	6.66	119.39	111.33
1	D	27	ALA	N-CA-C	6.26	118.38	110.24
1	D	320	ASP	N-CA-C	6.15	119.17	110.50
1	D	226	PRO	N-CA-C	5.92	120.37	111.14
1	D	194	ASP	CA-C-N	5.68	125.80	119.32
1	D	194	ASP	C-N-CA	5.68	125.80	119.32
1	A	273	GLY	N-CA-C	5.41	118.35	112.29
1	D	164	VAL	N-CA-C	5.18	119.73	113.00
1	D	273	GLY	N-CA-C	5.14	118.05	112.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	321	LYS	N-CA-C	5.10	121.66	110.80
1	D	317	PHE	N-CA-C	5.07	116.97	107.99
1	D	237	ARG	N-CA-C	5.03	117.53	110.23

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	2675	0	2724	33	0
1	D	2675	0	2724	53	0
2	A	28	0	12	0	0
2	D	28	0	12	0	0
3	A	5	0	0	1	0
3	D	5	0	0	1	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	332	0	0	3	0
6	D	287	0	0	9	0
All	All	6039	0	5472	84	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 8.

All (84) 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:201:ILE:HD12	1:A:281:LEU:HD23	1.48	0.95
1:D:13:VAL:HG11	1:D:31:LEU:HD11	1.64	0.80
1:A:21:SER:HA	1:A:26:ASN:HD21	1.50	0.76
1:D:315:LYS:HE3	1:D:318:ARG:HD3	1.69	0.73
1:A:201:ILE:HD11	1:A:284:GLN:HG3	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:ASN:HB2	6:D:2124:HOH:O	1.88	0.71
1:D:32:PRO:HG3	1:D:289:ILE:HG12	1.76	0.68
1:D:317:PHE:CZ	1:D:319:PRO:HB2	2.33	0.64
1:A:191:LYS:HE3	1:A:225:LEU:HD21	1.78	0.64
1:D:297:ARG:NH2	6:D:2273:HOH:O	2.30	0.64
1:D:731:MSE:HG2	6:D:2006:HOH:O	2.00	0.62
1:A:201:ILE:HD12	1:A:281:LEU:CD2	2.27	0.61
1:D:318:ARG:HD2	6:D:2280:HOH:O	2.00	0.61
1:A:325:ASP:O	1:A:328:VAL:HG12	2.00	0.61
1:D:19:ALA:HB1	1:D:731:MSE:HG3	1.83	0.61
1:A:10:ILE:HG13	1:A:131:ASN:HB3	1.83	0.60
1:A:9:LEU:H	1:A:9:LEU:HD23	1.66	0.60
1:D:739:LEU:HA	1:D:742:ILE:HG22	1.83	0.59
1:D:249:THR:HB	6:D:2221:HOH:O	2.04	0.57
1:D:300:LEU:HB3	1:D:739:LEU:HD21	1.86	0.57
1:D:16:LEU:HD21	1:D:735:LEU:HD13	1.85	0.57
1:D:6:MSE:SE	1:D:9:LEU:HD21	2.54	0.56
1:A:140:MSE:SE	1:A:189:VAL:HG11	2.54	0.56
1:A:320:ASP:HB3	1:A:322:HIS:H	1.70	0.55
1:D:13:VAL:CG1	1:D:31:LEU:HD11	2.35	0.55
1:D:11:PRO:O	1:D:15:ARG:HG2	2.07	0.55
1:D:292:THR:HG23	6:D:2271:HOH:O	2.07	0.54
1:D:145:VAL:H	1:D:148:GLN:NE2	2.08	0.52
1:D:325:ASP:O	1:D:328:VAL:HG12	2.09	0.52
1:A:10:ILE:HD12	6:A:2137:HOH:O	2.10	0.51
1:D:270:ASP:HB2	6:D:2254:HOH:O	2.10	0.51
1:A:48:LEU:HD22	1:A:134:LEU:HD13	1.92	0.50
1:D:10:ILE:N	1:D:11:PRO:CD	2.74	0.50
1:A:189:VAL:HG13	6:A:2182:HOH:O	2.11	0.50
1:D:30:ASP:HB3	1:D:168:ASN:OD1	2.12	0.50
1:A:318:ARG:C	1:A:320:ASP:N	2.69	0.50
1:D:130:LEU:HB3	1:D:132:LEU:HD22	1.93	0.49
1:A:145:VAL:HG21	1:D:224:LEU:CD1	2.43	0.49
1:D:168:ASN:HB2	6:D:2015:HOH:O	2.13	0.48
1:D:6:MSE:HA	1:D:9:LEU:CD2	2.44	0.47
1:A:303:GLN:O	1:A:307:ILE:HG12	2.15	0.47
1:D:13:VAL:HG11	1:D:31:LEU:CD1	2.38	0.47
1:A:172:LEU:HD13	1:A:281:LEU:HD21	1.97	0.46
1:A:79:GLU:HA	1:A:125:TYR:O	2.16	0.46
1:D:40:GLN:HG3	3:D:1745:ALF:F3	2.05	0.46
1:D:162:GLN:O	1:D:166:LYS:HE3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:54:ARG:HA	1:D:54:ARG:HD2	1.73	0.46
1:D:315:LYS:HE3	1:D:318:ARG:HH11	1.80	0.46
1:A:16:LEU:HA	1:A:734:ALA:HB1	1.97	0.46
1:A:12:LEU:HG	1:A:741:ILE:HG13	1.98	0.46
1:D:6:MSE:HA	1:D:9:LEU:HD23	1.98	0.46
1:D:187:LEU:HD13	1:D:225:LEU:HD23	1.97	0.46
1:A:9:LEU:H	1:A:9:LEU:CD2	2.29	0.45
1:D:742:ILE:HG23	1:D:742:ILE:O	2.15	0.45
1:A:145:VAL:HG21	1:D:224:LEU:HD11	1.98	0.45
1:A:75:ASN:H	1:A:131:ASN:ND2	2.15	0.45
1:D:85:HIS:CE1	1:D:104:GLU:HG3	2.51	0.45
1:D:739:LEU:HD13	1:D:742:ILE:HG21	1.99	0.44
1:A:6:MSE:SE	1:A:289:ILE:HD12	2.67	0.44
1:A:115:ILE:HD11	1:A:155:GLN:HE21	1.83	0.43
1:A:189:VAL:HG12	6:A:2192:HOH:O	2.19	0.43
1:D:303:GLN:O	1:D:307:ILE:HG12	2.18	0.43
1:A:227:LEU:HD12	1:A:227:LEU:HA	1.90	0.43
1:A:6:MSE:HE1	1:A:131:ASN:O	2.19	0.42
1:A:115:ILE:HD12	1:A:115:ILE:C	2.44	0.42
1:D:12:LEU:HD13	1:D:12:LEU:HA	1.90	0.42
1:D:318:ARG:N	1:D:319:PRO:CD	2.81	0.42
1:A:36:VAL:HG21	1:A:44:LYS:HB3	2.01	0.42
1:D:300:LEU:HB3	1:D:739:LEU:CD2	2.50	0.42
1:D:29:LEU:HD12	6:D:2011:HOH:O	2.19	0.42
1:D:262:HIS:HA	1:D:263:PRO:HD3	1.85	0.41
1:D:314:TYR:O	1:D:315:LYS:HB2	2.20	0.41
1:A:140:MSE:SE	1:A:189:VAL:CG1	3.18	0.41
1:D:36:VAL:HG23	1:D:36:VAL:O	2.20	0.41
1:D:293:LEU:N	1:D:294:PRO:CD	2.84	0.41
1:D:79:GLU:HG2	1:D:94:PHE:CE2	2.54	0.41
1:A:40:GLN:HG3	3:A:1745:ALF:F3	2.09	0.41
1:D:6:MSE:HE2	1:D:289:ILE:HG23	2.02	0.41
1:A:318:ARG:C	1:A:320:ASP:H	2.27	0.41
1:D:13:VAL:HA	1:D:16:LEU:HD12	2.02	0.41
1:D:140:MSE:HG3	1:D:160:LEU:HD11	2.01	0.41
1:D:156:ILE:O	1:D:160:LEU:HG	2.21	0.41
1:D:115:ILE:C	1:D:115:ILE:HD12	2.46	0.41
1:D:19:ALA:CB	1:D:731:MSE:HG3	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	339/353 (96%)	330 (97%)	9 (3%)	0	100	100
1	D	339/353 (96%)	326 (96%)	7 (2%)	6 (2%)	6	3
All	All	678/706 (96%)	656 (97%)	16 (2%)	6 (1%)	14	9

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	315	LYS
1	D	316	ASN
1	D	321	LYS
1	D	26	ASN
1	D	322	HIS
1	D	742	ILE

5.3.2 Protein sidechains [i](#)

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	298/301 (99%)	294 (99%)	4 (1%)	61	68
1	D	298/301 (99%)	281 (94%)	17 (6%)	18	15
All	All	596/602 (99%)	575 (96%)	21 (4%)	32	32

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	VAL
1	A	227	LEU
1	A	252	LEU
1	A	293	LEU
1	D	7	GLU
1	D	52	VAL
1	D	54	ARG
1	D	57	LEU
1	D	99	LEU
1	D	130	LEU
1	D	132	LEU
1	D	168	ASN
1	D	172	LEU
1	D	293	LEU
1	D	296	LEU
1	D	300	LEU
1	D	303	GLN
1	D	304	LEU
1	D	310	GLU
1	D	735	LEU
1	D	736	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	26	ASN
1	A	33	GLN
1	A	75	ASN
1	A	128	HIS
1	A	131	ASN
1	A	155	GLN
1	A	183	ASN
1	A	278	GLN
1	A	287	ASN
1	A	298	ASN
1	A	301	GLN
1	D	26	ASN
1	D	112	ASN
1	D	121	ASN
1	D	148	GLN
1	D	155	GLN
1	D	162	GLN

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Mol	Chain	Res	Type
1	D	196	GLN
1	D	733	HIS

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	ALF	A	1745	5,6,2,4	4,4,4	1.50	0	-		
3	ALF	D	1745	5,6,2,4	4,4,4	1.51	0	-		
2	GDP	A	1744	3,5,4	29,30,30	0.70	1 (3%)	45,47,47	0.93	3 (6%)
2	GDP	D	1744	3,5,4	29,30,30	0.61	0	45,47,47	0.94	3 (6%)

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	1744	3,5,4	-	0/16/32/32	0/3/3/3
2	GDP	D	1744	3,5,4	-	0/16/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1744	GDP	PB-O2B	-2.20	1.46	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1744	GDP	O2A-PA-O3A	3.07	115.58	107.27
2	A	1744	GDP	O2A-PA-O3A	2.84	114.96	107.27
2	A	1744	GDP	O3A-PA-O1A	-2.34	103.65	110.70
2	D	1744	GDP	O3A-PA-O1A	-2.30	103.78	110.70
2	A	1744	GDP	O2B-PB-O3A	-2.09	97.61	104.64
2	D	1744	GDP	O2B-PB-O1B	2.08	118.95	110.83

There are no chirality outliers.

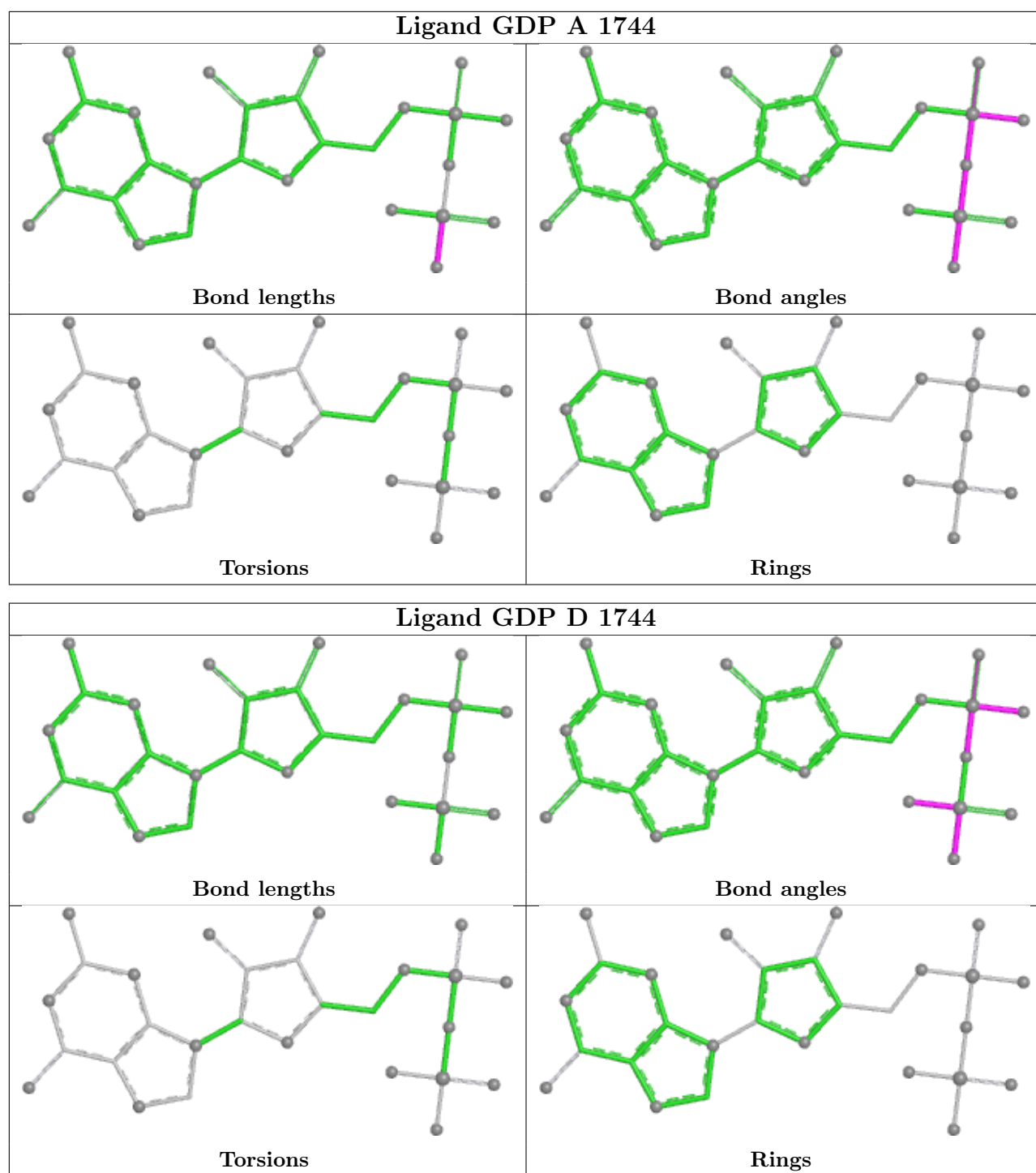
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1745	ALF	1	0
3	D	1745	ALF	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	333/353 (94%)	-0.27	14 (4%) 40 39	3, 12, 35, 61	0
1	D	333/353 (94%)	0.04	24 (7%) 21 20	3, 12, 48, 67	0
All	All	666/706 (94%)	-0.11	38 (5%) 29 28	3, 12, 45, 67	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	319	PRO	6.7
1	D	317	PHE	5.4
1	D	316	ASN	5.1
1	A	322	HIS	4.7
1	A	323	GLY	4.6
1	D	318	ARG	4.4
1	D	315	LYS	4.1
1	A	321	LYS	4.0
1	D	27	ALA	3.7
1	D	323	GLY	3.7
1	D	20	PHE	3.6
1	D	321	LYS	3.6
1	D	322	HIS	3.5
1	D	320	ASP	3.5
1	D	324	THR	3.5
1	A	7	GLU	3.3
1	A	741	ILE	3.2
1	D	741	ILE	3.1
1	D	742	ILE	3.1
1	D	23	ILE	3.0
1	A	8	ASP	3.0
1	A	319	PRO	2.9
1	D	9	LEU	2.8
1	A	324	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	17	GLN	2.7
1	D	26	ASN	2.6
1	D	87	LYS	2.5
1	D	743	GLY	2.5
1	A	743	GLY	2.3
1	A	147	ASP	2.3
1	A	320	ASP	2.2
1	A	145	VAL	2.2
1	D	16	LEU	2.2
1	D	18	ASP	2.1
1	D	29	LEU	2.1
1	A	742	ILE	2.1
1	A	312	GLU	2.0
1	D	297	ARG	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 [i](#)

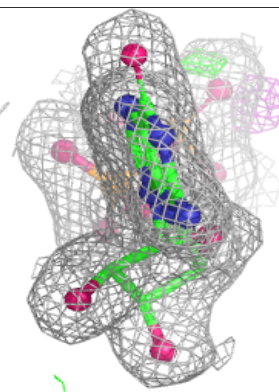
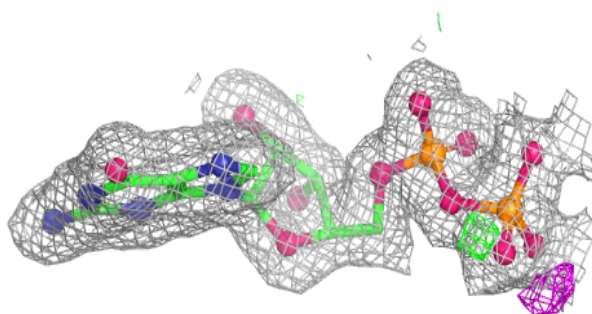
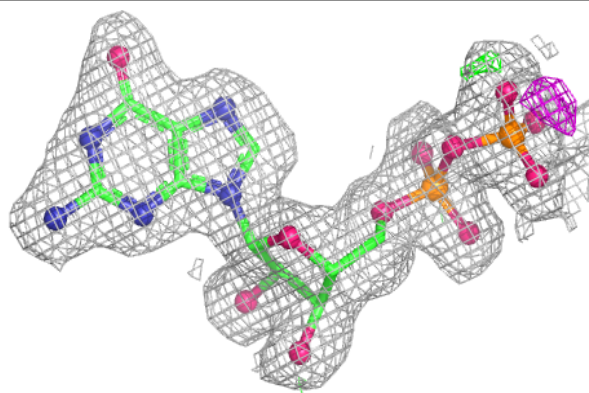
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NA	A	1747	1/1	0.93	0.05	8,8,8,8	0
5	NA	D	1747	1/1	0.95	0.04	9,9,9,9	0
3	ALF	A	1745	5/5	0.96	0.07	14,14,15,15	0
3	ALF	D	1745	5/5	0.96	0.06	11,12,12,13	0
4	MG	D	1746	1/1	0.98	0.02	4,4,4,4	0
2	GDP	A	1744	28/28	0.98	0.04	3,4,5,5	0
2	GDP	D	1744	28/28	0.98	0.04	4,5,7,7	0
4	MG	A	1746	1/1	0.99	0.03	5,5,5,5	0

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

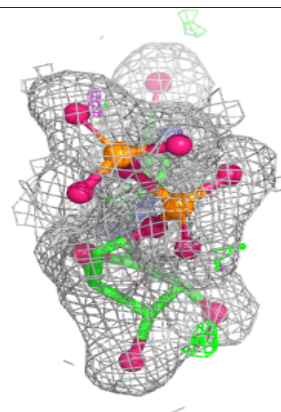
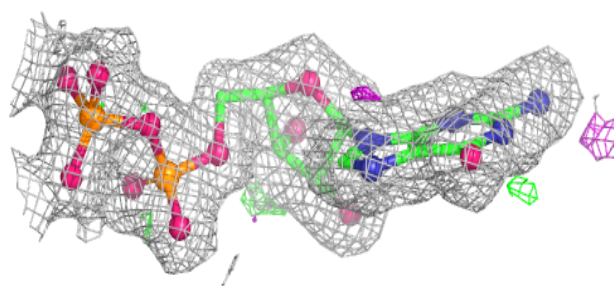
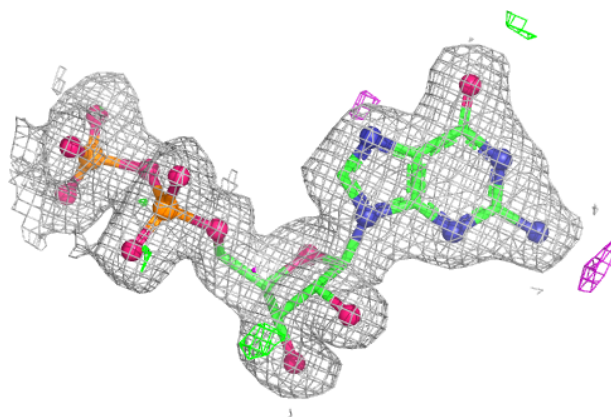
Electron density around GDP A 1744:

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



Electron density around GDP D 1744:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.