



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 06:33 PM EDT

PDB ID : 5C1S / pdb_00005c1s
Title : Crystal structure of the GDP-bound fast hydrolyzing mutant (V71A/K73Q) of EhRabX3 from Entamoeba histolytica
Authors : Srivastava, V.K.; Chandra, M.; Datta, S.
Deposited on : 2015-06-15
Resolution : 3.10 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
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

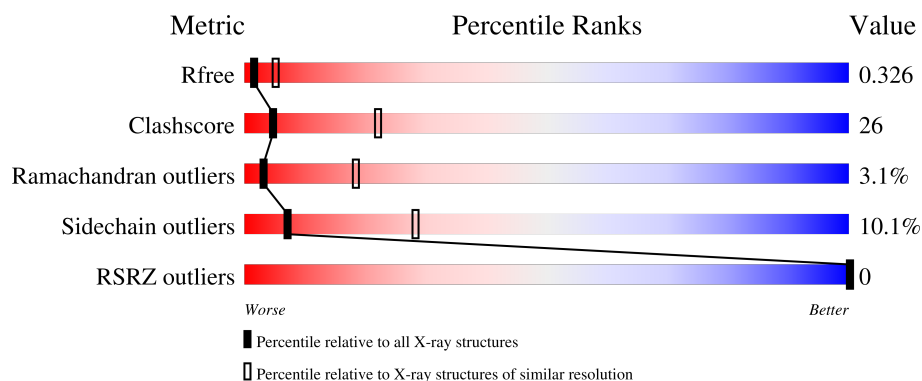
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

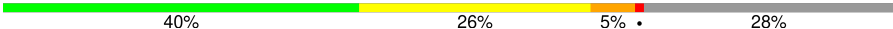
The reported resolution of this entry is 3.10 Å.

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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	391	 42% 28% 5% 26%
1	B	391	 40% 26% 5% 28%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4108 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 Small GTPase EhRabX3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	Se	0	0	0
			2033	1274	336	411	8	4			
1	B	283	Total	C	N	O	S	Se	0	0	0
			1965	1245	315	394	7	4			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-41	MSE	-	expression tag	UNP Q5NT25
A	-40	GLY	-	expression tag	UNP Q5NT25
A	-39	SER	-	expression tag	UNP Q5NT25
A	-38	SER	-	expression tag	UNP Q5NT25
A	-37	HIS	-	expression tag	UNP Q5NT25
A	-36	HIS	-	expression tag	UNP Q5NT25
A	-35	HIS	-	expression tag	UNP Q5NT25
A	-34	HIS	-	expression tag	UNP Q5NT25
A	-33	HIS	-	expression tag	UNP Q5NT25
A	-32	HIS	-	expression tag	UNP Q5NT25
A	-31	SER	-	expression tag	UNP Q5NT25
A	-30	SER	-	expression tag	UNP Q5NT25
A	-29	GLY	-	expression tag	UNP Q5NT25
A	-28	LEU	-	expression tag	UNP Q5NT25
A	-27	VAL	-	expression tag	UNP Q5NT25
A	-26	PRO	-	expression tag	UNP Q5NT25
A	-25	ARG	-	expression tag	UNP Q5NT25
A	-24	GLY	-	expression tag	UNP Q5NT25
A	-23	SER	-	expression tag	UNP Q5NT25
A	-22	HIS	-	expression tag	UNP Q5NT25
A	-21	MSE	-	expression tag	UNP Q5NT25
A	-20	ALA	-	expression tag	UNP Q5NT25
A	-19	SER	-	expression tag	UNP Q5NT25
A	-18	MSE	-	expression tag	UNP Q5NT25
A	-17	THR	-	expression tag	UNP Q5NT25

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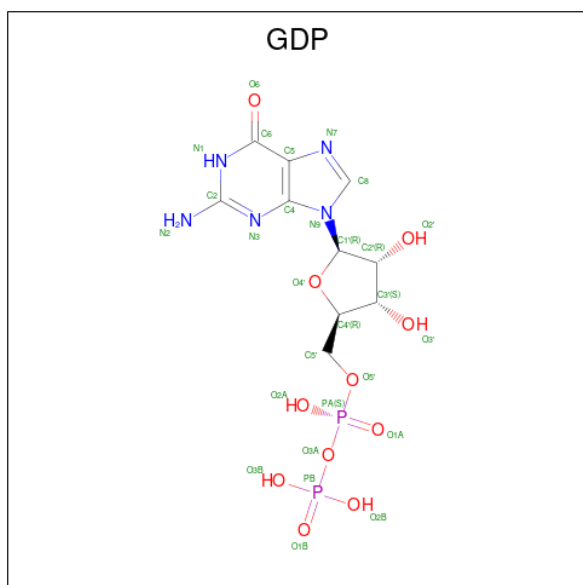
Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	GLY	-	expression tag	UNP Q5NT25
A	-15	GLY	-	expression tag	UNP Q5NT25
A	-14	GLN	-	expression tag	UNP Q5NT25
A	-13	GLN	-	expression tag	UNP Q5NT25
A	-12	MSE	-	expression tag	UNP Q5NT25
A	-11	GLY	-	expression tag	UNP Q5NT25
A	-10	ARG	-	expression tag	UNP Q5NT25
A	-9	GLY	-	expression tag	UNP Q5NT25
A	-8	SER	-	expression tag	UNP Q5NT25
A	-7	MSE	-	expression tag	UNP Q5NT25
A	-6	ALA	-	expression tag	UNP Q5NT25
A	-5	MSE	-	expression tag	UNP Q5NT25
A	-4	ILE	-	expression tag	UNP Q5NT25
A	-3	ARG	-	expression tag	UNP Q5NT25
A	-2	PRO	-	expression tag	UNP Q5NT25
A	-1	ALA	-	expression tag	UNP Q5NT25
A	0	HIS	-	expression tag	UNP Q5NT25
A	1	MSE	-	expression tag	UNP Q5NT25
A	71	ALA	VAL	engineered mutation	UNP Q5NT25
A	73	GLN	LYS	engineered mutation	UNP Q5NT25
B	-41	MSE	-	expression tag	UNP Q5NT25
B	-40	GLY	-	expression tag	UNP Q5NT25
B	-39	SER	-	expression tag	UNP Q5NT25
B	-38	SER	-	expression tag	UNP Q5NT25
B	-37	HIS	-	expression tag	UNP Q5NT25
B	-36	HIS	-	expression tag	UNP Q5NT25
B	-35	HIS	-	expression tag	UNP Q5NT25
B	-34	HIS	-	expression tag	UNP Q5NT25
B	-33	HIS	-	expression tag	UNP Q5NT25
B	-32	HIS	-	expression tag	UNP Q5NT25
B	-31	SER	-	expression tag	UNP Q5NT25
B	-30	SER	-	expression tag	UNP Q5NT25
B	-29	GLY	-	expression tag	UNP Q5NT25
B	-28	LEU	-	expression tag	UNP Q5NT25
B	-27	VAL	-	expression tag	UNP Q5NT25
B	-26	PRO	-	expression tag	UNP Q5NT25
B	-25	ARG	-	expression tag	UNP Q5NT25
B	-24	GLY	-	expression tag	UNP Q5NT25
B	-23	SER	-	expression tag	UNP Q5NT25
B	-22	HIS	-	expression tag	UNP Q5NT25
B	-21	MSE	-	expression tag	UNP Q5NT25
B	-20	ALA	-	expression tag	UNP Q5NT25

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	SER	-	expression tag	UNP Q5NT25
B	-18	MSE	-	expression tag	UNP Q5NT25
B	-17	THR	-	expression tag	UNP Q5NT25
B	-16	GLY	-	expression tag	UNP Q5NT25
B	-15	GLY	-	expression tag	UNP Q5NT25
B	-14	GLN	-	expression tag	UNP Q5NT25
B	-13	GLN	-	expression tag	UNP Q5NT25
B	-12	MSE	-	expression tag	UNP Q5NT25
B	-11	GLY	-	expression tag	UNP Q5NT25
B	-10	ARG	-	expression tag	UNP Q5NT25
B	-9	GLY	-	expression tag	UNP Q5NT25
B	-8	SER	-	expression tag	UNP Q5NT25
B	-7	MSE	-	expression tag	UNP Q5NT25
B	-6	ALA	-	expression tag	UNP Q5NT25
B	-5	MSE	-	expression tag	UNP Q5NT25
B	-4	ILE	-	expression tag	UNP Q5NT25
B	-3	ARG	-	expression tag	UNP Q5NT25
B	-2	PRO	-	expression tag	UNP Q5NT25
B	-1	ALA	-	expression tag	UNP Q5NT25
B	0	HIS	-	expression tag	UNP Q5NT25
B	1	MSE	-	expression tag	UNP Q5NT25
B	71	ALA	VAL	engineered mutation	UNP Q5NT25
B	73	GLN	LYS	engineered mutation	UNP Q5NT25

- 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
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

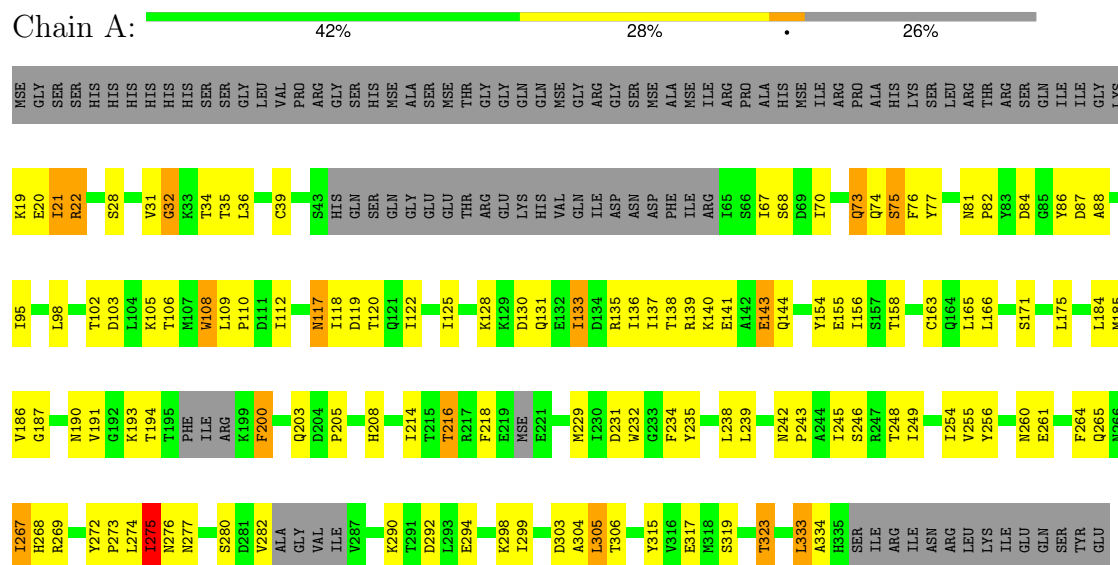
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	29	Total	O	0	0
			29	29		
3	B	25	Total	O	0	0
			25	25		

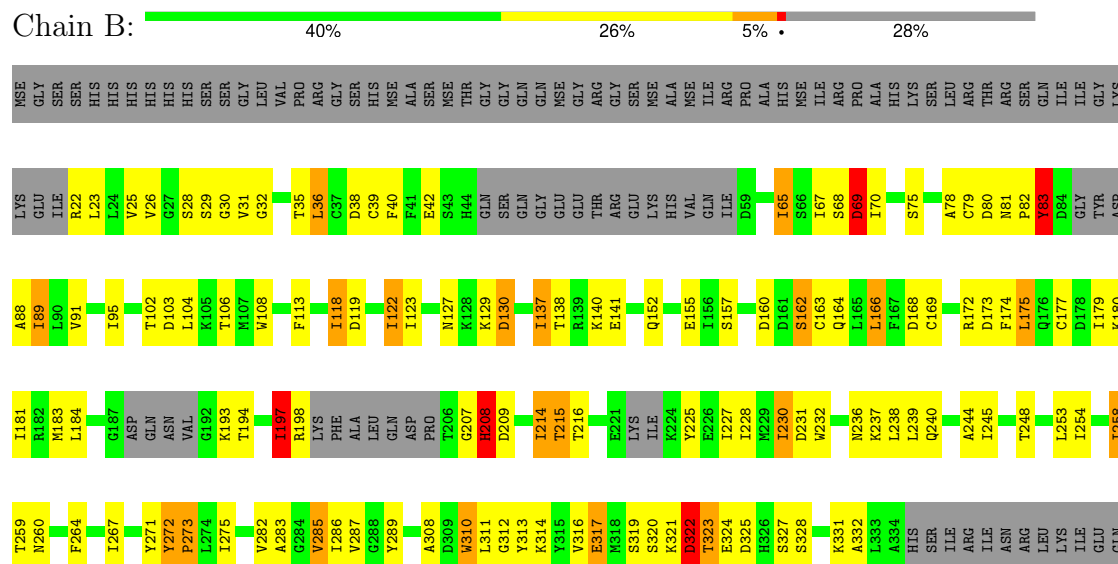
3 Residue-property plots

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: Small GTPase EhRabX3



• Molecule 1: Small GTPase EhRabX3



SER
TYR
GLU

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	63.68Å 63.68Å 324.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.62 – 3.10 38.62 – 3.10	Depositor EDS
% Data completeness (in resolution range)	97.5 (38.62-3.10) 98.8 (38.62-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.40 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.290 , 0.320 0.291 , 0.326	Depositor DCC
R_{free} test set	663 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	94.6	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 485.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.489 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4108	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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

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	0.44	1/2047 (0.0%)	1.03	14/2776 (0.5%)
1	B	0.54	1/1977 (0.1%)	1.04	11/2678 (0.4%)
All	All	0.49	2/4024 (0.0%)	1.04	25/5454 (0.5%)

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	4
1	B	0	3
All	All	0	7

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	197	ILE	CA-CB	-7.80	1.43	1.54
1	A	185	MSE	SE-CE	-5.25	1.79	1.95

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	207	GLY	N-CA-C	-8.21	103.31	115.72
1	A	205	PRO	N-CA-CB	8.00	111.48	103.48
1	B	273	PRO	N-CA-CB	7.24	110.85	103.25
1	A	299	ILE	N-CA-C	-6.91	104.72	113.22
1	B	310	TRP	N-CA-C	-6.91	104.86	113.28
1	A	82	PRO	N-CA-CB	6.80	111.05	103.44
1	A	76	PHE	N-CA-C	6.49	120.13	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	TYR	N-CA-C	-6.00	104.66	111.14
1	B	69	ASP	N-CA-C	-5.82	106.15	113.20
1	A	73	GLN	CB-CA-C	-5.74	109.94	116.54
1	B	166	LEU	N-CA-C	5.44	118.13	111.82
1	A	203	GLN	CB-CA-C	-5.42	110.34	116.63
1	B	82	PRO	N-CA-CB	5.42	108.94	103.25
1	A	290	LYS	CB-CA-C	5.38	118.47	110.62
1	B	208	HIS	CA-C-N	-5.17	117.32	126.32
1	B	208	HIS	C-N-CA	-5.17	117.32	126.32
1	A	108	TRP	N-CA-C	5.15	116.70	111.14
1	B	215	THR	CB-CA-C	-5.13	109.68	115.79
1	A	218	PHE	CB-CA-C	-5.13	109.12	115.89
1	A	200	PHE	N-CA-C	-5.10	106.84	113.16
1	A	269	ARG	N-CA-C	-5.08	106.34	112.54
1	A	274	LEU	CB-CA-C	-5.08	110.32	117.23
1	A	290	LYS	CA-CB-CG	5.07	124.24	114.10
1	A	32	GLY	N-CA-C	-5.02	105.40	112.13
1	B	237	LYS	N-CA-C	-5.00	105.54	111.69

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	119	ASP	Peptide
1	A	208	HIS	Peptide
1	A	68	SER	Peptide
1	A	75	SER	Peptide
1	B	162	SER	Peptide
1	B	258	ILE	Peptide
1	B	322	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	2033	0	1769	84	2
1	B	1965	0	1671	108	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	28	0	12	3	0
2	B	28	0	12	1	0
3	A	29	0	0	4	1
3	B	25	0	0	2	1
All	All	4108	0	3464	193	3

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 26.

All (193) 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:165:LEU:CD2	3:A:514:HOH:O	1.92	1.17
1:B:193:LYS:O	1:B:197:ILE:HG23	1.43	1.15
1:B:194:THR:O	1:B:197:ILE:HD11	1.46	1.13
1:B:194:THR:HA	1:B:197:ILE:CD1	1.78	1.12
1:B:194:THR:C	1:B:197:ILE:HD13	1.80	1.04
1:B:194:THR:C	1:B:197:ILE:CD1	2.33	1.00
1:B:194:THR:O	1:B:197:ILE:CD1	2.12	0.98
1:B:194:THR:CA	1:B:197:ILE:HD13	1.95	0.96
1:B:194:THR:CA	1:B:197:ILE:CD1	2.43	0.95
1:B:197:ILE:CG1	1:B:198:ARG:H	1.75	0.94
1:B:197:ILE:HG13	1:B:198:ARG:H	1.34	0.92
1:B:197:ILE:CG1	1:B:198:ARG:N	2.30	0.91
1:B:287:VAL:HG12	1:B:289:TYR:H	1.37	0.89
1:B:80:ASP:HA	1:B:83:TYR:HB2	1.55	0.89
1:B:194:THR:HA	1:B:197:ILE:HD13	1.54	0.86
1:B:193:LYS:O	1:B:197:ILE:CG2	2.24	0.84
1:A:158:THR:HA	1:A:163:CYS:HB2	1.62	0.82
1:A:333:LEU:HG	1:A:334:ALA:H	1.44	0.81
1:B:198:ARG:NH1	1:B:319:SER:OG	2.13	0.80
1:B:194:THR:HA	1:B:197:ILE:HD12	1.61	0.80
1:B:197:ILE:HG12	1:B:198:ARG:N	1.98	0.76
1:A:165:LEU:HD22	3:A:514:HOH:O	1.70	0.75
1:A:35:THR:HG21	2:A:401:GDP:H8	1.53	0.73
1:B:236:ASN:HA	1:B:239:LEU:HD13	1.71	0.71
1:B:283:ALA:HB1	1:B:310:TRP:N	2.06	0.71
1:B:208:HIS:O	1:B:209:ASP:CB	2.38	0.69
1:A:84:ASP:O	1:A:117:ASN:ND2	2.26	0.69
1:B:130:ASP:OD1	1:B:157:SER:OG	2.13	0.67
1:A:77:TYR:CE2	3:A:522:HOH:O	2.48	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:LEU:HD23	1:B:312:GLY:N	2.10	0.66
1:A:77:TYR:HE2	3:A:522:HOH:O	1.79	0.66
1:A:298:LYS:NZ	1:A:303:ASP:O	2.28	0.66
1:B:32:GLY:O	1:B:36:LEU:N	2.29	0.65
1:B:179:ILE:HG22	1:B:225:TYR:HA	1.78	0.65
1:B:287:VAL:HG12	1:B:289:TYR:N	2.10	0.65
1:B:258:ILE:HD11	1:B:287:VAL:HG11	1.77	0.64
1:A:235:TYR:HH	1:A:246:SER:HG	1.42	0.64
1:B:181:ILE:HB	1:B:227:ILE:HD13	1.79	0.64
1:B:31:VAL:O	1:B:127:ASN:ND2	2.30	0.63
1:B:285:VAL:HG22	1:B:286:ILE:H	1.63	0.63
1:B:324:GLU:OE2	1:B:327:SER:OG	2.16	0.63
1:A:319:SER:O	1:A:323:THR:OG1	2.14	0.62
1:A:74:GLN:O	1:A:75:SER:OG	2.16	0.62
1:B:258:ILE:O	1:B:259:THR:HG22	2.01	0.61
1:A:235:TYR:OH	1:A:246:SER:OG	2.12	0.61
1:B:194:THR:C	1:B:197:ILE:HD11	2.08	0.61
1:B:327:SER:O	1:B:331:LYS:N	2.34	0.60
1:B:332:ALA:HB1	3:B:518:HOH:O	2.01	0.60
1:A:95:ILE:HG23	1:A:137:ILE:HB	1.83	0.60
1:B:78:ALA:HA	1:B:81:ASN:HB2	1.83	0.60
1:A:333:LEU:CG	1:A:334:ALA:H	2.15	0.59
1:B:259:THR:HG23	1:B:260:ASN:H	1.68	0.59
1:A:35:THR:HG21	2:A:401:GDP:C8	2.36	0.59
1:B:129:LYS:HB2	1:B:157:SER:HB2	1.86	0.58
1:A:305:LEU:O	1:A:306:THR:OG1	2.21	0.57
1:A:117:ASN:O	1:A:120:THR:OG1	2.21	0.56
1:B:183:MSE:HE3	1:B:227:ILE:HG21	1.87	0.56
1:A:275:ILE:HG23	1:A:276:ASN:H	1.71	0.56
1:A:264:PHE:O	1:A:267:ILE:HG12	2.05	0.56
1:A:36:LEU:HA	1:A:39:CYS:HB3	1.88	0.55
1:A:280:SER:H	1:A:304:ALA:HB3	1.71	0.55
1:A:231:ASP:OD2	1:A:234:PHE:N	2.40	0.55
1:A:256:TYR:CD1	1:A:260:ASN:HB2	2.41	0.55
1:B:283:ALA:HB1	1:B:310:TRP:H	1.71	0.54
1:B:69:ASP:O	1:B:70:ILE:HG22	2.08	0.54
1:B:197:ILE:C	1:B:198:ARG:O	2.48	0.54
1:A:139:ARG:HB3	1:A:155:GLU:CD	2.32	0.54
1:B:214:ILE:HG23	1:B:215:THR:H	1.73	0.54
1:A:304:ALA:C	1:A:306:THR:H	2.14	0.53
1:B:231:ASP:OD1	1:B:232:TRP:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:ILE:O	1:B:198:ARG:C	2.48	0.52
1:B:215:THR:OG1	1:B:216:THR:N	2.29	0.52
1:A:268:HIS:O	1:A:273:PRO:HD3	2.09	0.52
1:A:186:VAL:HG13	1:A:254:ILE:HG23	1.91	0.52
1:A:333:LEU:HG	1:A:334:ALA:N	2.21	0.52
1:A:117:ASN:HD22	1:A:120:THR:HG21	1.75	0.51
1:B:238:LEU:HD13	1:B:245:ILE:HG12	1.92	0.51
1:A:102:THR:HA	1:A:105:LYS:NZ	2.25	0.51
1:B:28:SER:C	1:B:30:GLY:H	2.18	0.51
1:A:139:ARG:HD3	1:A:155:GLU:OE2	2.11	0.51
1:B:68:SER:C	1:B:70:ILE:H	2.19	0.51
1:B:39:CYS:SG	1:B:164:GLN:N	2.81	0.50
1:B:264:PHE:O	1:B:267:ILE:HG13	2.12	0.50
1:B:138:THR:HG22	1:B:141:GLU:H	1.77	0.50
1:B:113:PHE:HB3	1:B:244:ALA:HB3	1.92	0.50
1:B:197:ILE:O	1:B:198:ARG:O	2.30	0.50
1:A:298:LYS:NZ	1:A:304:ALA:O	2.26	0.49
1:B:325:ASP:O	1:B:328:SER:OG	2.28	0.49
1:A:130:ASP:OD2	1:A:131:GLN:N	2.45	0.49
2:B:401:GDP:N2	3:B:503:HOH:O	2.45	0.49
1:A:234:PHE:CE2	1:A:238:LEU:HD23	2.47	0.48
1:B:169:CYS:HA	1:B:172:ARG:HG2	1.94	0.48
1:B:104:LEU:HA	1:B:108:TRP:HE3	1.78	0.48
1:B:259:THR:HG23	1:B:260:ASN:N	2.27	0.48
1:A:276:ASN:ND2	1:A:277:ASN:O	2.46	0.48
1:A:22:ARG:HB2	1:A:86:TYR:HD2	1.79	0.48
1:A:214:ILE:HG23	1:A:229:MSE:HB2	1.95	0.48
1:B:102:THR:O	1:B:106:THR:HG23	2.14	0.48
1:B:283:ALA:HB2	1:B:308:ALA:O	2.14	0.48
1:A:186:VAL:HG21	1:A:265:GLN:HB3	1.95	0.47
1:A:261:GLU:HB2	1:A:264:PHE:CE2	2.49	0.47
1:B:123:ILE:HG12	1:B:152:GLN:HB2	1.96	0.47
1:B:316:VAL:HA	1:B:319:SER:HB3	1.95	0.47
1:A:191:VAL:O	1:A:194:THR:OG1	2.25	0.47
1:B:78:ALA:CA	1:B:81:ASN:HB2	2.45	0.47
1:A:294:GLU:HG3	1:A:305:LEU:O	2.15	0.47
1:B:258:ILE:CD1	1:B:287:VAL:HG11	2.45	0.46
1:A:238:LEU:HD11	1:A:245:ILE:HG21	1.96	0.46
1:B:23:LEU:HA	1:B:88:ALA:N	2.30	0.46
1:B:103:ASP:HA	1:B:106:THR:HG1	1.81	0.46
1:B:245:ILE:O	1:B:248:THR:OG1	2.27	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ARG:NH2	1:A:155:GLU:OE2	2.48	0.46
1:B:89:ILE:O	1:B:122:ILE:HA	2.15	0.46
1:B:138:THR:HG22	1:B:140:LYS:N	2.30	0.46
1:B:314:LYS:O	1:B:316:VAL:N	2.49	0.46
1:A:87:ASP:OD2	1:A:88:ALA:N	2.49	0.46
1:A:131:GLN:O	1:A:133:ILE:HG13	2.16	0.46
1:B:193:LYS:O	1:B:197:ILE:HD13	2.16	0.46
1:A:98:LEU:N	1:A:136:ILE:HD11	2.31	0.45
1:B:215:THR:HG1	1:B:216:THR:H	1.61	0.45
1:B:118:ILE:HG13	1:B:119:ASP:H	1.80	0.45
1:A:19:LYS:N	1:A:21:ILE:HD11	2.32	0.45
1:B:173:ASP:OD1	1:B:174:PHE:N	2.50	0.45
1:A:171:SER:O	1:A:175:LEU:N	2.41	0.45
1:B:215:THR:HB	1:B:228:ILE:HG12	1.97	0.45
1:B:78:ALA:C	1:B:81:ASN:HB2	2.42	0.45
1:A:125:ILE:HA	1:A:154:TYR:O	2.16	0.45
1:A:238:LEU:O	1:A:242:ASN:N	2.41	0.45
1:B:103:ASP:OD1	1:B:106:THR:OG1	2.35	0.45
1:B:160:ASP:OD1	1:B:162:SER:N	2.50	0.45
1:B:183:MSE:N	1:B:228:ILE:O	2.50	0.44
1:A:143:GLU:HG3	1:A:144:GLN:N	2.32	0.44
1:B:236:ASN:HA	1:B:239:LEU:CD1	2.43	0.44
1:B:163:CYS:O	1:B:166:LEU:N	2.46	0.44
1:B:168:ASP:OD1	1:B:169:CYS:N	2.50	0.44
1:A:109:LEU:HD12	1:A:110:PRO:N	2.33	0.44
1:B:163:CYS:SG	1:B:164:GLN:N	2.88	0.44
1:B:271:TYR:O	1:B:272:TYR:HD1	2.00	0.44
1:B:103:ASP:HA	1:B:106:THR:OG1	2.17	0.44
1:A:238:LEU:CD1	1:A:245:ILE:HG21	2.48	0.43
1:A:103:ASP:HA	1:A:106:THR:HG22	1.99	0.43
1:A:184:LEU:HD11	1:A:232:TRP:HA	2.00	0.43
1:B:311:LEU:C	1:B:313:TYR:N	2.77	0.43
1:B:319:SER:O	1:B:322:ASP:HB2	2.18	0.43
1:B:320:SER:O	1:B:322:ASP:N	2.51	0.43
1:A:190:ASN:O	1:A:194:THR:HG23	2.18	0.43
1:B:314:LYS:O	1:B:317:GLU:HG3	2.18	0.43
1:A:191:VAL:HG21	1:A:256:TYR:C	2.44	0.43
1:A:239:LEU:HA	1:A:242:ASN:C	2.44	0.43
1:A:256:TYR:HB3	1:A:282:VAL:CB	2.49	0.43
1:A:125:ILE:HG21	1:A:156:ILE:HD12	2.01	0.43
1:A:84:ASP:C	1:A:117:ASN:HD21	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:ARG:O	1:B:88:ALA:N	2.51	0.43
1:B:23:LEU:HG	1:B:65:ILE:HD11	2.00	0.43
1:B:26:VAL:HG23	1:B:91:VAL:HG13	2.01	0.43
1:B:184:LEU:HA	1:B:230:ILE:HG22	1.99	0.43
1:A:138:THR:OG1	1:A:141:GLU:OE2	2.37	0.43
1:B:322:ASP:O	1:B:323:THR:C	2.62	0.42
1:B:258:ILE:HD12	1:B:289:TYR:HB2	2.01	0.42
1:A:77:TYR:HE1	1:A:81:ASN:HD22	1.66	0.42
1:A:131:GLN:C	1:A:133:ILE:H	2.27	0.42
1:A:139:ARG:HB3	1:A:155:GLU:OE2	2.19	0.42
1:A:118:ILE:HG23	1:A:248:THR:HA	2.02	0.42
1:B:28:SER:HB3	1:B:108:TRP:HZ2	1.85	0.42
1:A:34:THR:OG1	1:A:35:THR:N	2.52	0.42
1:B:38:ASP:HA	1:B:42:GLU:HB2	2.01	0.42
1:A:140:LYS:HA	1:A:140:LYS:HD3	1.75	0.42
1:A:200:PHE:HB2	1:A:216:THR:OG1	2.20	0.42
1:A:77:TYR:O	1:A:77:TYR:CG	2.73	0.42
1:B:80:ASP:C	1:B:83:TYR:H	2.28	0.41
1:B:163:CYS:O	1:B:164:GLN:C	2.62	0.41
1:B:179:ILE:HG13	1:B:180:LYS:H	1.84	0.41
1:B:320:SER:O	1:B:321:LYS:C	2.63	0.41
1:B:137:ILE:HD12	1:B:137:ILE:HA	1.91	0.41
1:A:73:GLN:HB3	1:A:74:GLN:H	1.43	0.41
1:A:246:SER:HA	1:A:249:ILE:HG12	2.03	0.41
1:B:172:ARG:O	1:B:175:LEU:HG	2.20	0.41
1:A:165:LEU:HD23	1:A:165:LEU:O	2.21	0.41
1:A:242:ASN:HA	1:A:243:PRO:HD2	1.90	0.41
1:A:186:VAL:HG22	1:A:187:GLY:H	1.85	0.41
1:A:32:GLY:H	2:A:401:GDP:PB	2.43	0.41
1:A:128:LYS:C	1:A:130:ASP:H	2.27	0.41
1:B:95:ILE:HG13	1:B:137:ILE:HG22	2.02	0.41
1:B:40:PHE:CD2	1:B:67:ILE:HG21	2.56	0.41
1:A:28:SER:HA	1:A:108:TRP:HE1	1.86	0.40
1:A:156:ILE:HD11	1:A:166:LEU:HD22	2.02	0.40
1:A:268:HIS:O	1:A:272:TYR:HB2	2.21	0.40
1:B:28:SER:C	1:B:30:GLY:N	2.78	0.40
1:A:186:VAL:HG23	1:A:232:TRP:CE3	2.56	0.40
1:B:184:LEU:O	1:B:253:LEU:N	2.48	0.40
1:B:316:VAL:HA	1:B:319:SER:CB	2.51	0.40
1:A:81:ASN:OD1	1:A:84:ASP:HB2	2.21	0.40
1:A:102:THR:HA	1:A:105:LYS:HZ3	1.86	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:514:HOH:O	3:B:513:HOH:O[1_445]	1.42	0.78
1:A:193:LYS:NZ	1:B:79:CYS:O[1_445]	2.16	0.04
1:A:136:ILE:O	1:B:138:THR:OG1[6_555]	2.19	0.01

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	278/391 (71%)	218 (78%)	54 (19%)	6 (2%)	5	24
1	B	271/391 (69%)	208 (77%)	52 (19%)	11 (4%)	2	13
All	All	549/782 (70%)	426 (78%)	106 (19%)	17 (3%)	3	18

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	197	ILE
1	B	208	HIS
1	B	273	PRO
1	A	67	ILE
1	B	272	TYR
1	B	323	THR
1	A	70	ILE
1	B	285	VAL
1	B	282	VAL
1	B	322	ASP
1	A	305	LEU
1	A	317	GLU
1	B	29	SER
1	B	75	SER
1	A	275	ILE

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Mol	Chain	Res	Type
1	A	267	ILE
1	B	89	ILE

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	186/333 (56%)	170 (91%)	16 (9%)	10	34
1	B	172/333 (52%)	152 (88%)	20 (12%)	5	22
All	All	358/666 (54%)	322 (90%)	36 (10%)	7	28

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLU
1	A	21	ILE
1	A	22	ARG
1	A	31	VAL
1	A	112	ILE
1	A	117	ASN
1	A	122	ILE
1	A	133	ILE
1	A	143	GLU
1	A	216	THR
1	A	255	VAL
1	A	275	ILE
1	A	292	ASP
1	A	315	TYR
1	A	323	THR
1	A	333	LEU
1	B	25	VAL
1	B	35	THR
1	B	36	LEU
1	B	65	ILE
1	B	69	ASP

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Mol	Chain	Res	Type
1	B	83	TYR
1	B	118	ILE
1	B	122	ILE
1	B	130	ASP
1	B	137	ILE
1	B	155	GLU
1	B	175	LEU
1	B	177	CYS
1	B	197	ILE
1	B	214	ILE
1	B	230	ILE
1	B	240	GLN
1	B	254	ILE
1	B	275	ILE
1	B	317	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	A	236	ASN
1	A	277	ASN
1	B	81	ASN
1	B	260	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	280/391 (71%)	-1.28	0 100 100	64, 94, 144, 181	0
1	B	274/391 (70%)	-1.31	0 100 100	52, 95, 134, 169	0
All	All	554/782 (70%)	-1.30	0 100 100	52, 95, 141, 181	0

There are no RSRZ outliers to report.

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](#)

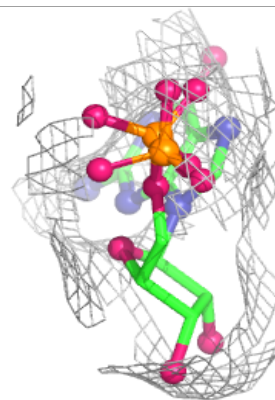
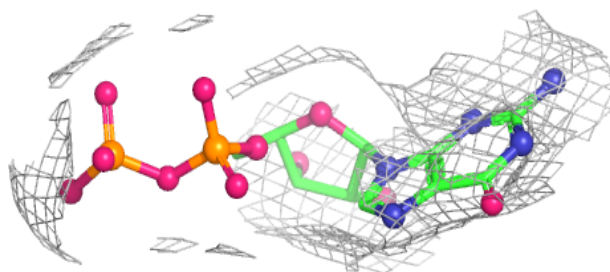
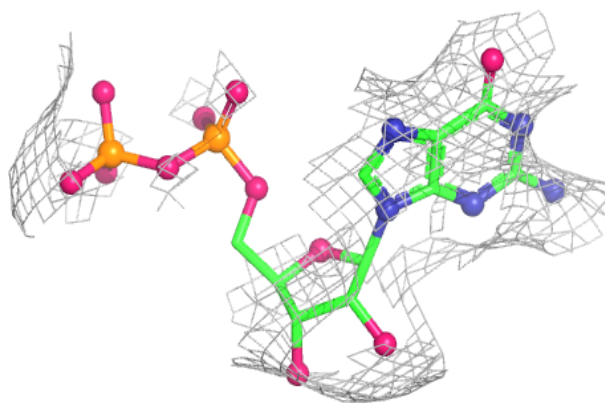
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GDP	B	401	28/28	0.98	0.04	90,127,154,158	0
2	GDP	A	401	28/28	0.99	0.03	79,105,133,136	0

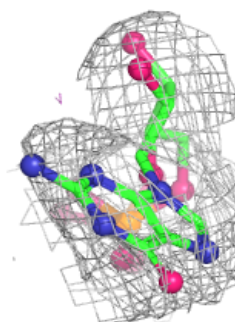
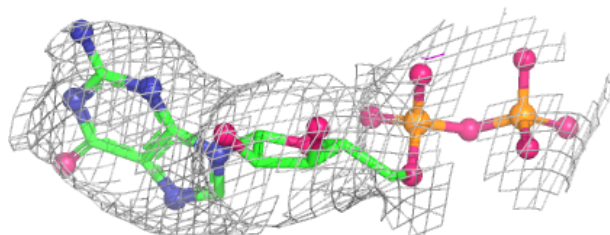
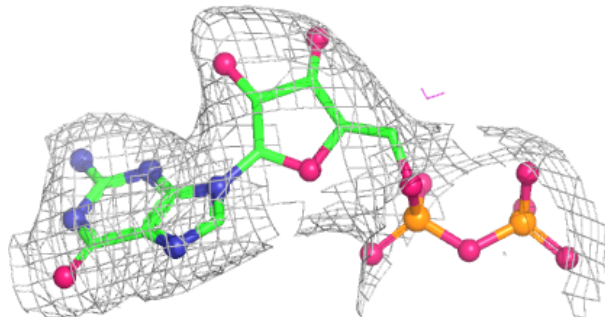
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GDP B 401:

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

**Electron density around GDP A 401:**

$2mF_o-DF_c$ (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.