



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:24 AM UTC

PDB ID : 4N0I / pdb_00004n0i
Title : Crystal structure of S. cerevisiae mitochondrial GatFAB in complex with glutamine
Authors : Araiso, Y.; Ishitani, R.; Nureki, O.
Deposited on : 2013-10-02
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
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

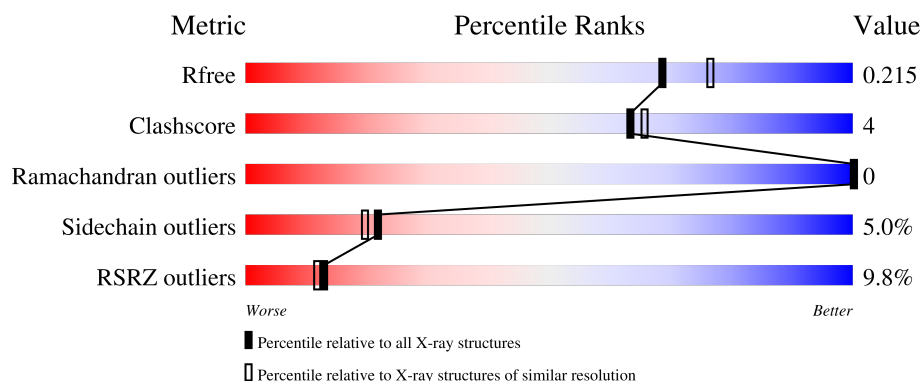
1 Overall quality at a glance

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	464	2% 86% 9% • •
2	B	325	23% 75% 12% • 12%
3	F	160	2% 70% 6% • 22%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6933 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 Glutamyl-tRNA(Gln) amidotransferase subunit A, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	5	0
			3439	2202	571	653	13			

- Molecule 2 is a protein called Glutamyl-tRNA(Gln) amidotransferase subunit B, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	285	Total	C	N	O	S	0	0	0
			2199	1397	376	420	6			

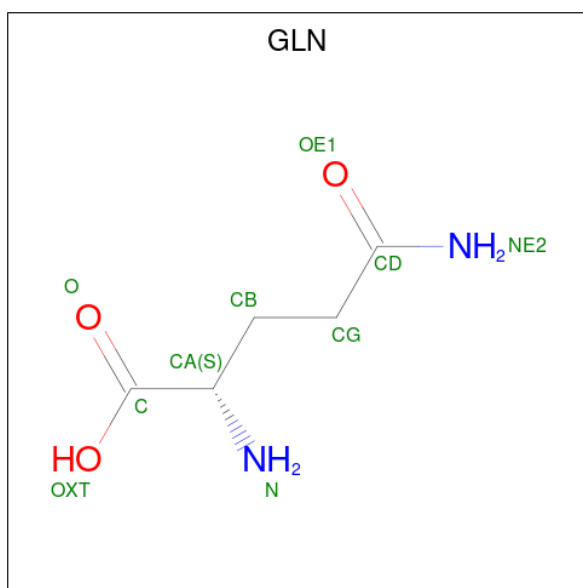
There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	330	LEU	-	expression tag	UNP P33893
B	331	GLU	-	expression tag	UNP P33893
B	332	SER	-	expression tag	UNP P33893
B	333	SER	-	expression tag	UNP P33893
B	334	GLY	-	expression tag	UNP P33893
B	335	GLU	-	expression tag	UNP P33893
B	336	ASN	-	expression tag	UNP P33893
B	337	LEU	-	expression tag	UNP P33893
B	338	TYR	-	expression tag	UNP P33893
B	339	PHE	-	expression tag	UNP P33893
B	340	GLN	-	expression tag	UNP P33893

- Molecule 3 is a protein called Glutamyl-tRNA(Gln) amidotransferase subunit F, mitochondrial.

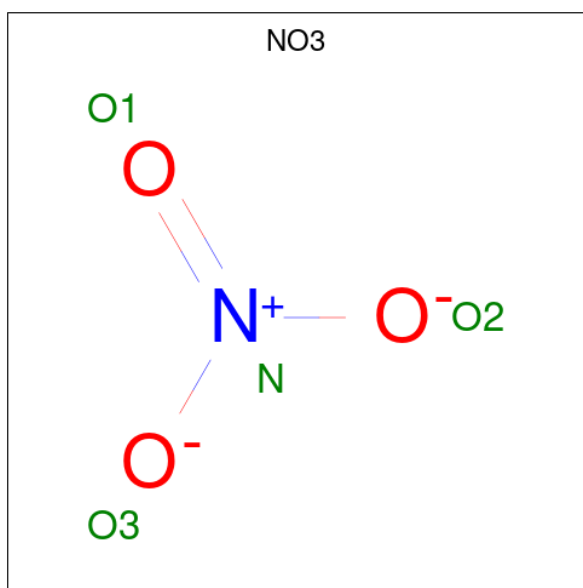
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	125	Total	C	N	O	S	0	0	0
			955	611	167	176	1			

- Molecule 4 is GLUTAMINE (CCD ID: GLN) (formula: $C_5H_{10}N_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			9	5	1	3		

- Molecule 5 is NITRATE ION (CCD ID: NO3) (formula: NO_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	N	O	0	0
			4	1	3		

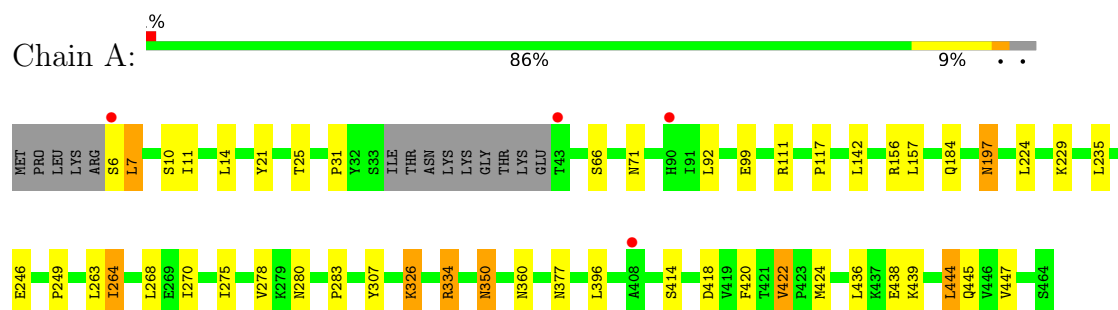
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	218	Total 218	O 218	0	0
6	B	54	Total 54	O 54	0	0
6	F	55	Total 55	O 55	0	0

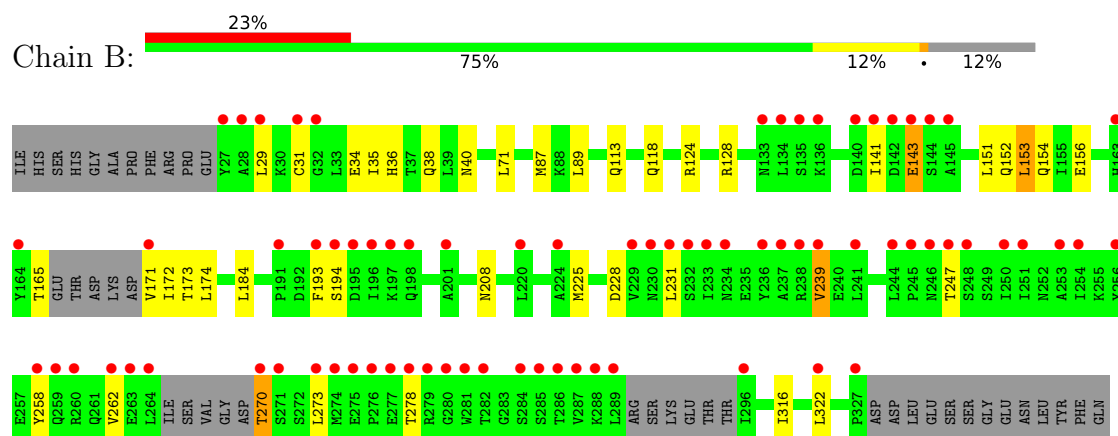
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

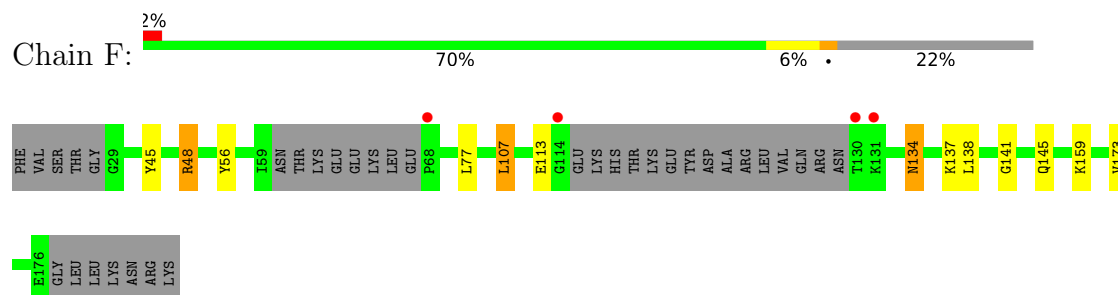
- Molecule 1: Glutamyl-tRNA(Gln) amidotransferase subunit A, mitochondrial



- Molecule 2: Glutamyl-tRNA(Gln) amidotransferase subunit B, mitochondrial



- Molecule 3: Glutamyl-tRNA(Gln) amidotransferase subunit F, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.44Å 85.43Å 194.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.46 – 2.00 26.46 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.9 (26.46-2.00) 96.9 (26.46-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.88 (at 1.99Å)	Xtriage
Refinement program	PHENIX 1.7.2 _869	Depositor
R, R_{free}	0.189 , 0.219 0.184 , 0.215	Depositor DCC
R_{free} test set	3448 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6933	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.90% 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: NO3

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	0.46	0/3518	0.80	3/4772 (0.1%)
2	B	0.43	0/2241	0.80	2/3048 (0.1%)
3	F	0.38	0/971	0.71	0/1308
All	All	0.44	0/6730	0.79	5/9128 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	422	VAL	CA-C-N	-6.19	112.18	119.05
1	A	422	VAL	C-N-CA	-6.19	112.18	119.05
2	B	124	ARG	CA-C-N	5.66	125.64	120.03
2	B	124	ARG	C-N-CA	5.66	125.64	120.03
1	A	184	GLN	N-CA-C	5.34	117.85	111.71

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	3439	0	3445	29	0
2	B	2199	0	2147	18	0
3	F	955	0	951	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	9	0	5	1	0
5	A	4	0	0	0	0
6	A	218	0	0	3	0
6	B	54	0	0	1	0
6	F	55	0	0	0	0
All	All	6933	0	6548	52	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 4.

All (52) 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:334:ARG:NH2	4:A:501:GLN:OXT	2.15	0.79
1:A:436:LEU:HD21	1:A:444:LEU:HD22	1.73	0.70
2:B:270:THR:O	2:B:270:THR:OG1	2.12	0.68
1:A:424[B]:MET:HE3	1:A:447:VAL:HG21	1.81	0.62
1:A:418:ASP:OD2	6:A:684:HOH:O	2.15	0.62
2:B:152:GLN:OE1	2:B:154:GLN:NE2	2.32	0.62
1:A:249:PRO:HB3	1:A:396:LEU:HG	1.83	0.60
1:A:377:ASN:ND2	6:A:779:HOH:O	2.33	0.60
3:F:134:ASN:ND2	3:F:137:LYS:H	2.01	0.59
1:A:283:PRO:HB2	3:F:107:LEU:HD13	1.87	0.57
1:A:263:LEU:HB3	1:A:268:LEU:HD11	1.88	0.55
1:A:229:LYS:NZ	6:A:773:HOH:O	2.39	0.54
2:B:31:CYS:HB3	2:B:193:PHE:HB2	1.89	0.54
2:B:89:LEU:HD23	2:B:153:LEU:HD21	1.92	0.52
2:B:128:ARG:NH2	6:B:429:HOH:O	2.37	0.52
1:A:307:TYR:CE2	1:A:326:LYS:HG3	2.45	0.52
2:B:113:GLN:NE2	2:B:118:GLN:HB2	2.24	0.52
2:B:29:LEU:O	2:B:194:SER:HA	2.11	0.50
1:A:7:LEU:O	1:A:11:ILE:HG12	2.12	0.50
1:A:264:ILE:HG13	1:A:270:ILE:HD12	1.93	0.49
1:A:275:ILE:O	1:A:278:VAL:HG22	2.13	0.49
2:B:113:GLN:HE22	2:B:118:GLN:HB2	1.78	0.48
1:A:438:GLU:HG2	1:A:439:LYS:HD3	1.96	0.48
1:A:156:ARG:HH11	1:A:445:GLN:HE22	1.63	0.47
1:A:66:SER:HB2	1:A:99:GLU:HG3	1.96	0.47
1:A:156:ARG:HH11	1:A:445:GLN:NE2	2.13	0.46
1:A:350:ASN:HD22	1:A:350:ASN:HA	1.58	0.46
3:F:45:TYR:HA	3:F:48:ARG:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:35:ILE:HG21	2:B:225:MET:HE3	1.97	0.45
2:B:87:MET:HE1	2:B:316:ILE:HD13	1.99	0.45
2:B:165:THR:O	2:B:165:THR:OG1	2.34	0.45
1:A:21:TYR:CE1	1:A:117:PRO:HA	2.52	0.45
1:A:197[B]:ASN:ND2	1:A:197[B]:ASN:H	2.14	0.45
2:B:34:GLU:HB2	2:B:228:ASP:HB2	2.00	0.44
3:F:134:ASN:HD22	3:F:134:ASN:C	2.26	0.43
2:B:38:GLN:HA	2:B:184:LEU:HD23	2.00	0.43
2:B:231:LEU:HD12	2:B:239:VAL:HG13	2.01	0.43
3:F:141:GLY:O	3:F:145:GLN:HG3	2.18	0.43
1:A:156:ARG:HD2	1:A:445:GLN:HE21	1.83	0.43
2:B:174:LEU:HD23	3:F:173:VAL:HG12	2.00	0.43
1:A:11:ILE:HD12	1:A:31:PRO:HG2	2.01	0.43
1:A:246:GLU:HG3	1:A:420:PHE:CZ	2.53	0.43
2:B:141:ILE:HG22	2:B:143:GLU:H	1.84	0.42
1:A:280:ASN:ND2	3:F:56:TYR:OH	2.45	0.42
1:A:360:ASN:HD21	3:F:113:GLU:H	1.67	0.42
1:A:418:ASP:HB3	1:A:422:VAL:HG13	2.02	0.42
3:F:134:ASN:HD21	3:F:137:LYS:HG2	1.84	0.42
1:A:25:THR:O	1:A:111:ARG:NH1	2.45	0.41
2:B:36:HIS:HD2	2:B:156:GLU:OE1	2.03	0.41
1:A:156:ARG:HD2	1:A:445:GLN:NE2	2.36	0.41
2:B:258:TYR:O	2:B:262:VAL:HG23	2.21	0.40
1:A:6:SER:O	1:A:10:SER:N	2.51	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	451/464 (97%)	444 (98%)	7 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	277/325 (85%)	275 (99%)	2 (1%)	0	100	100
3	F	119/160 (74%)	117 (98%)	2 (2%)	0	100	100
All	All	847/949 (89%)	836 (99%)	11 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	384/407 (94%)	368 (96%)	16 (4%)	26	25
2	B	238/292 (82%)	223 (94%)	15 (6%)	16	13
3	F	97/137 (71%)	91 (94%)	6 (6%)	16	13
All	All	719/836 (86%)	682 (95%)	37 (5%)	22	19

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	14	LEU
1	A	71	ASN
1	A	92	LEU
1	A	142	LEU
1	A	157	LEU
1	A	197[A]	ASN
1	A	197[B]	ASN
1	A	224	LEU
1	A	235	LEU
1	A	264	ILE
1	A	326	LYS
1	A	334	ARG
1	A	350	ASN
1	A	414	SER
1	A	444	LEU

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Mol	Chain	Res	Type
2	B	40	ASN
2	B	71	LEU
2	B	143	GLU
2	B	151	LEU
2	B	153	LEU
2	B	171	VAL
2	B	172	ILE
2	B	173	THR
2	B	208	ASN
2	B	239	VAL
2	B	247	THR
2	B	270	THR
2	B	273	LEU
2	B	278	THR
2	B	322	LEU
3	F	48	ARG
3	F	77	LEU
3	F	107	LEU
3	F	134	ASN
3	F	138	LEU
3	F	159	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	265	ASN
1	A	377	ASN
1	A	445	GLN
2	B	36	HIS
2	B	40	ASN
2	B	113	GLN
2	B	118	GLN
2	B	133	ASN
2	B	261	GLN
3	F	134	ASN
3	F	147	GLN

5.3.3 RNA ⓘ

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	$\# Z > 2$
5	NO3	A	502	-	1,3,3	3.42	1 (100%)	0,3,3	-	-
4	GLN	A	501	-	6,8,9	1.01	1 (16%)	6,9,11	0.96	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GLN	A	501	-	-	0/8/8/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	502	NO3	O1-N	3.42	1.41	1.24
4	A	501	GLN	OXT-C	-2.12	1.23	1.30

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	GLN	1	0

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	450/464 (96%)	-0.29	4 (0%) 81 80	12, 36, 65, 105	5 (1%)
2	B	285/325 (87%)	1.04	76 (26%) 1 1	26, 53, 166, 180	0
3	F	125/160 (78%)	0.13	4 (3%) 50 49	28, 46, 70, 85	0
All	All	860/949 (90%)	0.21	84 (9%) 13 12	12, 42, 134, 180	5 (0%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	278	THR	7.8
2	B	141	ILE	6.5
2	B	27	TYR	6.2
2	B	279	ARG	5.9
2	B	281	TRP	5.5
2	B	287	VAL	5.3
2	B	241	LEU	5.0
2	B	282	THR	5.0
2	B	259	GLN	5.0
2	B	233	ILE	4.9
2	B	231	LEU	4.7
2	B	230	ASN	4.7
2	B	260	ARG	4.5
2	B	239	VAL	4.4
2	B	264	LEU	4.4
2	B	195	ASP	4.4
2	B	247	THR	4.3
2	B	280	GLY	4.3
2	B	236	TYR	4.3
2	B	171	VAL	4.2
2	B	145	ALA	4.2
2	B	285	SER	4.2
2	B	197	LYS	4.1

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Mol	Chain	Res	Type	RSRZ
2	B	276	PRO	4.1
2	B	270	THR	4.0
2	B	29	LEU	3.9
2	B	28	ALA	3.9
2	B	244	LEU	3.9
2	B	32	GLY	3.8
2	B	322	LEU	3.8
2	B	140	ASP	3.6
2	B	196	ILE	3.6
3	F	130	THR	3.6
2	B	198	GLN	3.6
2	B	289	LEU	3.5
3	F	114	GLY	3.5
2	B	258	TYR	3.5
1	A	43	THR	3.5
2	B	284	SER	3.4
2	B	256	TYR	3.4
2	B	273	LEU	3.3
2	B	271	SER	3.3
2	B	232	SER	3.3
2	B	296	ILE	3.2
2	B	248	SER	3.2
2	B	262	VAL	3.1
2	B	229	VAL	3.1
2	B	201	ALA	3.0
2	B	234	ASN	3.0
2	B	286	THR	3.0
2	B	142	ASP	3.0
1	A	90	HIS	3.0
2	B	134	LEU	2.9
2	B	250	ILE	2.9
2	B	246	ASN	2.9
2	B	135	SER	2.9
2	B	237	ALA	2.8
2	B	253	ALA	2.8
2	B	191	PRO	2.8
2	B	288	LYS	2.8
2	B	254	ILE	2.8
2	B	164	TYR	2.7
2	B	238	ARG	2.7
2	B	263	GLU	2.6
2	B	277	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	194	SER	2.6
2	B	251	ILE	2.6
2	B	193	PHE	2.6
2	B	144	SER	2.6
3	F	68	PRO	2.5
2	B	224	ALA	2.4
2	B	136	LYS	2.4
2	B	274	MET	2.4
1	A	408	ALA	2.3
3	F	131	LYS	2.3
1	A	6	SER	2.3
2	B	31	CYS	2.3
2	B	163	HIS	2.2
2	B	220	LEU	2.2
2	B	327	PRO	2.2
2	B	245	PRO	2.1
2	B	143	GLU	2.1
2	B	275	GLU	2.1
2	B	133	ASN	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
4	GLN	A	501	9/10	0.94	0.10	33,40,47,48	0
5	NO3	A	502	4/4	0.96	0.08	36,37,38,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.