



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:33 AM UTC

PDB ID : 5OL8 / pdb_00005ol8
Title : Structure of human mitochondrial transcription elongation factor (TEFM) C-terminal domain
Authors : Hillen, H.S.; Parshin, A.V.; Agaronyan, K.; Morozov, Y.; Graber, J.J.; Chernev, A.; Schwinghammer, K.; Urlaub, H.; Anikin, M.; Cramer, P.; Temiakov, D.
Deposited on : 2017-07-27
Resolution : 1.90 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

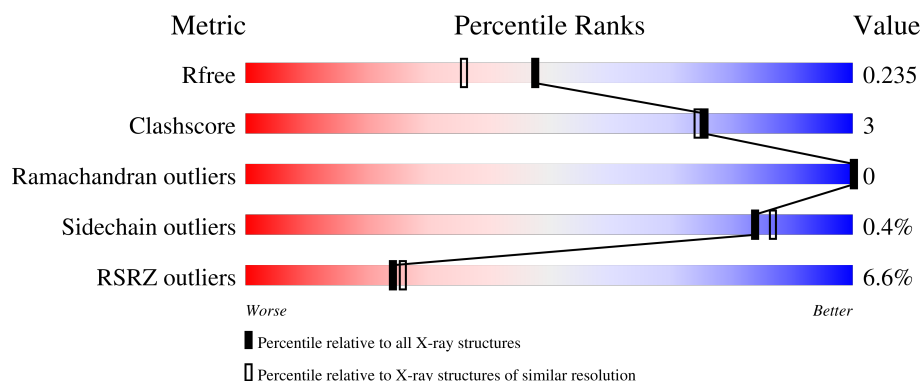
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	319	5% 58% 6% 36%
1	B	319	3% 59% 5% 36%
1	C	319	4% 59% 5% 36%
1	D	319	5% 61% . 36%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6877 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 Transcription elongation factor, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1681	1089	284	301	7			
1	B	205	Total	C	N	O	S	0	0	0
			1681	1089	284	301	7			
1	C	205	Total	C	N	O	S	0	0	0
			1681	1089	284	301	7			
1	D	205	Total	C	N	O	S	0	0	0
			1681	1089	284	301	7			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	MET	-	initiating methionine	UNP Q96QE5
A	361	LEU	-	expression tag	UNP Q96QE5
A	362	GLU	-	expression tag	UNP Q96QE5
A	363	HIS	-	expression tag	UNP Q96QE5
A	364	HIS	-	expression tag	UNP Q96QE5
A	365	HIS	-	expression tag	UNP Q96QE5
A	366	HIS	-	expression tag	UNP Q96QE5
A	367	HIS	-	expression tag	UNP Q96QE5
A	368	HIS	-	expression tag	UNP Q96QE5
B	50	MET	-	initiating methionine	UNP Q96QE5
B	361	LEU	-	expression tag	UNP Q96QE5
B	362	GLU	-	expression tag	UNP Q96QE5
B	363	HIS	-	expression tag	UNP Q96QE5
B	364	HIS	-	expression tag	UNP Q96QE5
B	365	HIS	-	expression tag	UNP Q96QE5
B	366	HIS	-	expression tag	UNP Q96QE5
B	367	HIS	-	expression tag	UNP Q96QE5
B	368	HIS	-	expression tag	UNP Q96QE5
C	50	MET	-	initiating methionine	UNP Q96QE5
C	361	LEU	-	expression tag	UNP Q96QE5
C	362	GLU	-	expression tag	UNP Q96QE5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	363	HIS	-	expression tag	UNP Q96QE5
C	364	HIS	-	expression tag	UNP Q96QE5
C	365	HIS	-	expression tag	UNP Q96QE5
C	366	HIS	-	expression tag	UNP Q96QE5
C	367	HIS	-	expression tag	UNP Q96QE5
C	368	HIS	-	expression tag	UNP Q96QE5
D	50	MET	-	initiating methionine	UNP Q96QE5
D	361	LEU	-	expression tag	UNP Q96QE5
D	362	GLU	-	expression tag	UNP Q96QE5
D	363	HIS	-	expression tag	UNP Q96QE5
D	364	HIS	-	expression tag	UNP Q96QE5
D	365	HIS	-	expression tag	UNP Q96QE5
D	366	HIS	-	expression tag	UNP Q96QE5
D	367	HIS	-	expression tag	UNP Q96QE5
D	368	HIS	-	expression tag	UNP Q96QE5

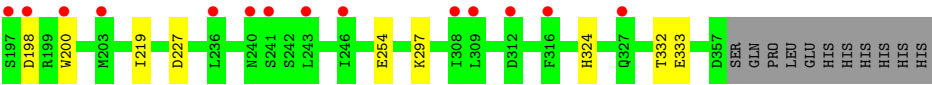
- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



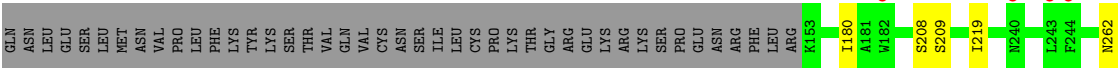
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total 38	O 38	0	0
3	B	29	Total 29	O 29	0	0
3	C	34	Total 34	O 34	0	0
3	D	28	Total 28	O 28	0	0



● Molecule 1: Transcription elongation factor, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.48Å 112.54Å 88.84Å 90.00° 110.13° 90.00°	Depositor
Resolution (Å)	46.23 – 1.90 46.23 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.23-1.90) 98.2 (46.23-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 1.90Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.195 , 0.228 0.202 , 0.235	Depositor DCC
R_{free} test set	3509 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6877	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% 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: GOL

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	0.13	0/1717	0.32	0/2314
1	B	0.14	0/1717	0.31	0/2314
1	C	0.14	0/1717	0.32	0/2314
1	D	0.13	0/1717	0.32	0/2314
All	All	0.13	0/6868	0.32	0/9256

There are no bond length outliers.

There are no bond angle outliers.

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	1681	0	1708	15	0
1	B	1681	0	1708	9	0
1	C	1681	0	1708	11	0
1	D	1681	0	1708	6	0
2	A	12	0	16	2	0
2	D	12	0	16	3	0
3	A	38	0	0	5	0
3	B	29	0	0	0	0
3	C	34	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	28	0	0	1	0
All	All	6877	0	6864	41	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 3.

All (41) 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:B:332:THR:HB	1:B:337:VAL:HG21	1.60	0.81
1:C:324:HIS:ND1	3:C:402:HOH:O	2.22	0.73
1:A:153:LYS:N	3:A:503:HOH:O	2.26	0.67
1:D:208:SER:HA	2:D:402:GOL:H32	1.78	0.65
1:A:254:GLU:OE2	3:A:501:HOH:O	2.15	0.63
1:C:179:ARG:HG2	1:C:198:ASP:HA	1.83	0.61
1:C:167:VAL:HG23	1:C:227:ASP:HB2	1.86	0.57
1:B:165:LYS:NZ	1:B:357:ASP:OD1	2.37	0.57
1:B:297:LYS:NZ	1:B:333:GLU:HG3	2.21	0.56
1:A:295:SER:CB	1:A:336:ARG:HH12	2.20	0.55
1:A:320:ASP:OD1	1:A:321:LYS:HG2	2.07	0.53
1:A:308:ILE:HG22	1:A:310:LYS:H	1.72	0.53
1:A:254:GLU:OE2	2:A:401:GOL:H11	2.09	0.53
1:D:297:LYS:HE3	1:D:333:GLU:HG3	1.93	0.50
1:A:336:ARG:NH1	1:A:338:GLU:OE2	2.44	0.50
1:B:180:ILE:HG21	1:B:219:ILE:HD13	1.94	0.49
1:A:180:ILE:HG21	1:A:219:ILE:HD13	1.94	0.49
1:A:262:ASN:HB3	1:A:265:PHE:HB2	1.95	0.49
1:C:180:ILE:HG21	1:C:219:ILE:HD13	1.95	0.48
1:C:161:ARG:HG2	1:C:165:LYS:HE3	1.95	0.48
1:D:262:ASN:HB3	1:D:265:PHE:HB2	1.93	0.48
1:D:180:ILE:HG21	1:D:219:ILE:HD13	1.96	0.48
1:B:161:ARG:O	1:B:165:LYS:HG2	2.14	0.47
1:C:175:PHE:HB3	1:C:180:ILE:HG13	1.96	0.47
1:B:187:ARG:HA	1:B:351:TYR:CE1	2.50	0.47
1:B:172:SER:O	1:B:182:TRP:HA	2.16	0.46
1:A:163:ARG:HG3	3:A:523:HOH:O	2.16	0.45
1:C:332:THR:OG1	3:C:401:HOH:O	2.20	0.45
1:A:199:ARG:HG3	1:A:219:ILE:HD11	2.00	0.43
1:C:297:LYS:NZ	1:C:333:GLU:O	2.51	0.43
1:A:209:SER:H	2:A:402:GOL:H12	1.84	0.42
1:A:302:GLN:O	1:A:306:ASP:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:LYS:NZ	3:D:507:HOH:O	2.53	0.42
1:B:171:ILE:HD11	1:B:224:PRO:HG2	2.01	0.42
1:C:200:TRP:C	1:C:200:TRP:CD1	2.98	0.41
1:D:209:SER:N	2:D:402:GOL:H11	2.35	0.41
1:B:253:MET:HE3	1:B:253:MET:HB3	1.92	0.41
1:C:254:GLU:OE2	2:D:402:GOL:O3	2.20	0.41
1:C:172:SER:O	1:C:182:TRP:HA	2.20	0.41
1:A:263:LYS:NZ	3:A:506:HOH:O	2.53	0.41
1:A:153:LYS:O	3:A:502:HOH:O	2.22	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	203/319 (64%)	200 (98%)	3 (2%)	0	100	100
1	B	203/319 (64%)	198 (98%)	5 (2%)	0	100	100
1	C	203/319 (64%)	197 (97%)	6 (3%)	0	100	100
1	D	203/319 (64%)	197 (97%)	6 (3%)	0	100	100
All	All	812/1276 (64%)	792 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	186/292 (64%)	185 (100%)	1 (0%)	81	84
1	B	186/292 (64%)	185 (100%)	1 (0%)	81	84
1	C	186/292 (64%)	186 (100%)	0	100	100
1	D	186/292 (64%)	185 (100%)	1 (0%)	81	84
All	All	744/1168 (64%)	741 (100%)	3 (0%)	84	87

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	240	ASN
1	B	232	GLU
1	D	323	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	240	ASN
1	A	324	HIS
1	A	346	GLN
1	B	346	GLN
1	C	195	GLN
1	C	272	GLN
1	C	346	GLN
1	D	251	HIS
1	D	346	GLN

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

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	401	-	5,5,5	0.36	0	5,5,5	0.59	0
2	GOL	D	402	-	5,5,5	0.39	0	5,5,5	0.32	0
2	GOL	A	402	-	5,5,5	0.38	0	5,5,5	0.35	0
2	GOL	D	401	-	5,5,5	0.40	0	5,5,5	0.27	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
2	GOL	A	401	-	-	4/4/4/4	-
2	GOL	D	402	-	-	2/4/4/4	-
2	GOL	A	402	-	-	2/4/4/4	-
2	GOL	D	401	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	GOL	C1-C2-C3-O3
2	A	402	GOL	O1-C1-C2-C3
2	D	401	GOL	O1-C1-C2-C3
2	D	402	GOL	O1-C1-C2-C3
2	A	402	GOL	O1-C1-C2-O2
2	D	402	GOL	O1-C1-C2-O2
2	A	401	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	401	GOL	O2-C2-C3-O3
2	D	401	GOL	O1-C1-C2-O2
2	A	401	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	GOL	1	0
2	D	402	GOL	3	0
2	A	402	GOL	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	205/319 (64%)	0.71	15 (7%) 21 22	39, 64, 104, 139	0
1	B	205/319 (64%)	0.61	10 (4%) 35 37	41, 63, 105, 118	0
1	C	205/319 (64%)	0.64	14 (6%) 23 24	41, 61, 99, 125	0
1	D	205/319 (64%)	0.77	15 (7%) 21 22	42, 65, 105, 143	0
All	All	820/1276 (64%)	0.68	54 (6%) 24 26	39, 63, 104, 143	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	308	ILE	5.8
1	A	308	ILE	5.5
1	A	309	LEU	5.0
1	D	309	LEU	5.0
1	C	309	LEU	3.6
1	D	316	PHE	3.5
1	B	334	LEU	3.2
1	A	244	PHE	3.1
1	A	240	ASN	3.0
1	B	306	ASP	2.9
1	A	310	LYS	2.9
1	D	323	VAL	2.9
1	D	334	LEU	2.8
1	D	240	ASN	2.8
1	B	309	LEU	2.8
1	D	330	LEU	2.7
1	B	308	ILE	2.7
1	B	244	PHE	2.7
1	C	241	SER	2.6
1	A	337	VAL	2.6
1	B	323	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	316	PHE	2.6
1	A	206	ILE	2.5
1	C	327	GLN	2.5
1	C	308	ILE	2.5
1	C	243	LEU	2.4
1	D	243	LEU	2.4
1	C	312	ASP	2.4
1	A	153	LYS	2.4
1	A	178	ARG	2.3
1	C	240	ASN	2.3
1	C	203	MET	2.3
1	B	238	ILE	2.2
1	D	327	GLN	2.2
1	A	200	TRP	2.2
1	C	200	TRP	2.2
1	C	246	ILE	2.2
1	D	319	SER	2.2
1	A	238	ILE	2.2
1	B	337	VAL	2.2
1	C	198	ASP	2.1
1	D	357	ASP	2.1
1	B	241	SER	2.1
1	C	197	SER	2.1
1	D	182	TRP	2.1
1	C	236	LEU	2.1
1	D	244	PHE	2.1
1	D	307	SER	2.1
1	A	334	LEU	2.1
1	A	291	ASP	2.1
1	A	329	PHE	2.0
1	C	316	PHE	2.0
1	B	167	VAL	2.0
1	D	322	ILE	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
2	GOL	A	402	6/6	0.79	0.18	65,74,77,86	0
2	GOL	D	401	6/6	0.84	0.20	82,93,94,97	0
2	GOL	D	402	6/6	0.84	0.15	70,78,80,82	0
2	GOL	A	401	6/6	0.85	0.14	71,73,75,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.