



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:05 PM GMT

PDB ID : 5AFR
Title : N-terminal fragment of dynein heavy chain
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Deposited on : 2015-01-23
Resolution : 5.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

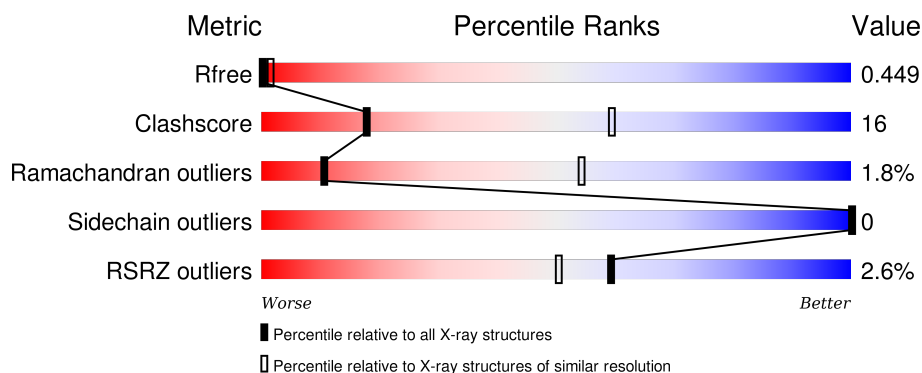
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.00 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}	91344	1119 (6.22-3.60)
Clashscore	102246	1019 (6.22-3.66)
Ramachandran outliers	100387	1158 (6.22-3.60)
Sidechain outliers	100360	1136 (6.22-3.60)
RSRZ outliers	91569	1122 (6.22-3.60)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	557	<div> <div>2%</div> <div>53% 6% 40%</div> </div>
1	B	557	<div> <div>%</div> <div>51% 8% 41%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3318 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.

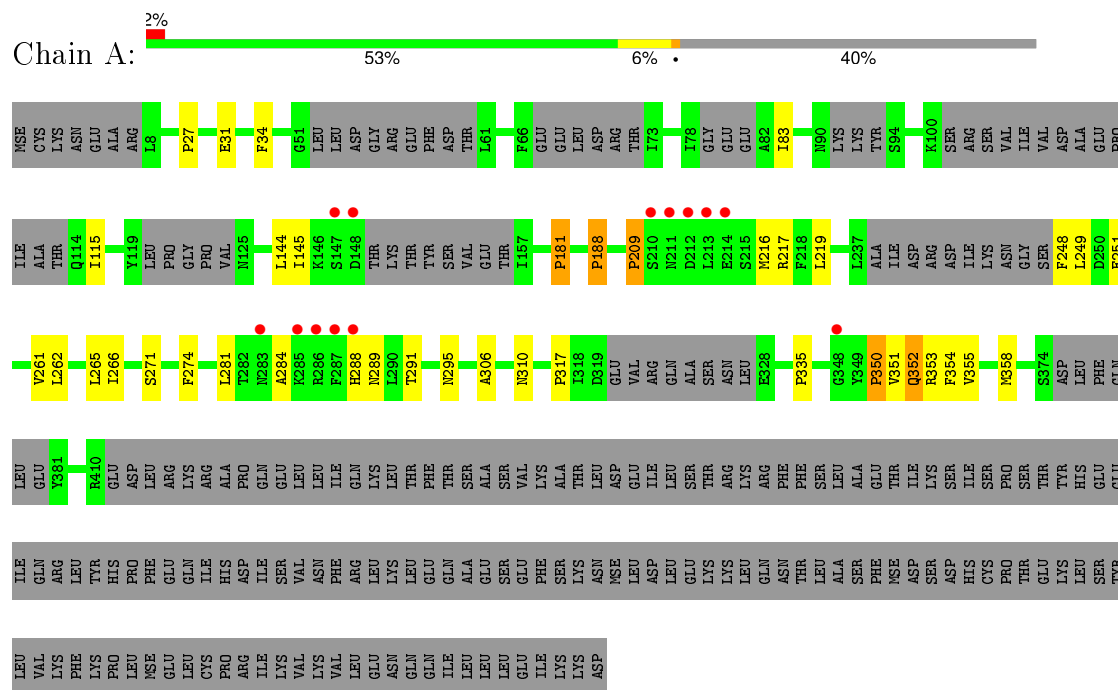
- Molecule 1 is a protein called DYNEIN HEAVY CHAIN, CYTOPLASMIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	Se	0	0	0
			1664	995	332	332	5			
1	B	330	Total	C	N	O	Se	0	0	0
			1654	989	330	330	5			

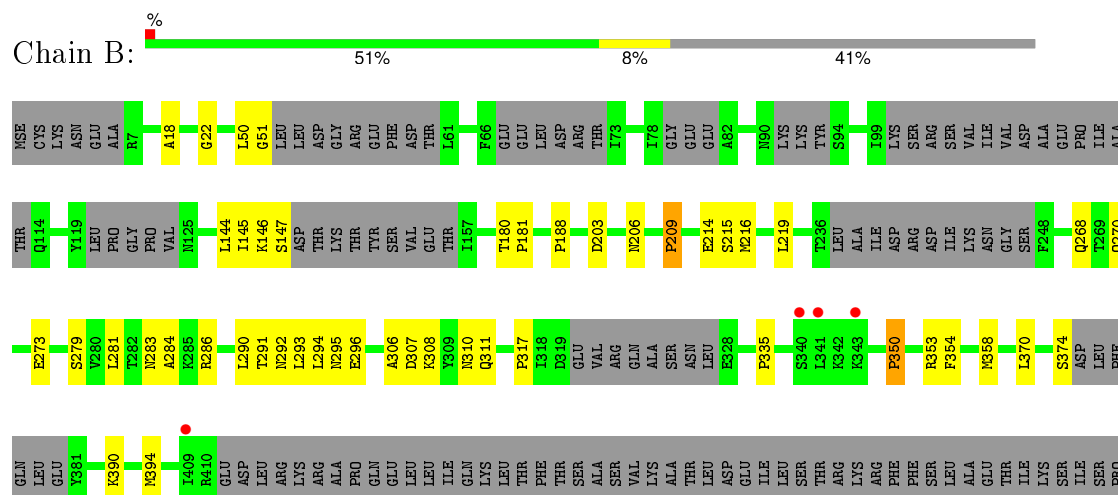
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: DYNEIN HEAVY CHAIN, CYTOPLASMIC



• Molecule 1: DYNEIN HEAVY CHAIN, CYTOPLASMIC



SER	THR	TYR	HIS	GLU	ILE	GLN	ARG	LEU	TYR	PRO	PHE	GLU	GLN	ILE	HIS	ASP	ILE	SER	VAL	ASN	PHE	ARG	LYS	LEU	LEU	GLU	GLN	ALA	GLU	SER	GLU	PHE	SER	LYS	ASN	ASP	LEU	LEU	GLU	LYS	LYS	GLN	ASN	THR	ALA	LEU	SER	PHE	ASP	ASP	CYS	PRO
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THR	GLU	LYS	LEU	SER	TYR	LEU	VAL	LYS	PHE	LYS	PRO	LEU	MSE	GLU	LEU	CYS	PRO	ARG	ILE	LYS	VAL	LYS	VAL	LEU	GLU	ASN	GLN	GLN	ILE	LEU	LEU	GLU	ILE	LYS	LYS	ASP
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.90Å 148.86Å 179.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	114.62 – 5.00 51.71 – 5.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (114.62-5.00) 99.9 (51.71-5.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.79 (at 5.10Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.393 , 0.437 0.395 , 0.449	Depositor DCC
R_{free} test set	414 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	230.0	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.63 , 2691.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	1 of 8801 reflections (0.011%)	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	3318	wwPDB-VP
Average B, all atoms (Å ²)	201.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.56	0/1648	0.67	6/2269 (0.3%)
1	B	0.59	0/1638	0.71	5/2255 (0.2%)
All	All	0.57	0/3286	0.69	11/4524 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	181	PRO	N-CA-CB	7.14	111.86	103.30
1	B	188	PRO	N-CA-CB	7.03	111.73	103.30
1	A	335	PRO	N-CA-CB	6.80	111.46	103.30
1	B	350	PRO	N-CA-CB	6.68	111.32	103.30
1	B	335	PRO	N-CA-CB	6.56	111.17	103.30
1	A	209	PRO	N-CA-CB	6.49	111.09	103.30
1	A	188	PRO	N-CA-CB	5.94	110.42	103.30
1	A	27	PRO	N-CA-CB	5.89	110.37	103.30
1	A	350	PRO	N-CA-CB	5.29	109.65	103.30
1	A	181	PRO	N-CA-CB	5.24	109.59	103.30
1	B	209	PRO	N-CA-CB	5.07	109.38	103.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1664	0	752	35	0
1	B	1654	0	748	41	0
All	All	3318	0	1500	76	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 16.

All (76) 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:214:GLU:O	1:B:216:MSE:HA	1.07	1.19
1:B:18:ALA:O	1:B:22:GLY:N	1.81	1.12
1:B:214:GLU:O	1:B:216:MSE:CA	2.03	1.06
1:A:31:GLU:O	1:A:34:PHE:N	1.88	1.05
1:B:215:SER:CB	1:B:219:LEU:H	1.75	0.99
1:B:370:LEU:O	1:B:374:SER:N	1.96	0.97
1:A:271:SER:O	1:A:274:PHE:N	1.97	0.96
1:A:281:LEU:O	1:A:284:ALA:CA	2.16	0.93
1:A:281:LEU:O	1:A:284:ALA:HB3	1.70	0.92
1:B:306:ALA:O	1:B:310:ASN:N	2.03	0.91
1:B:214:GLU:C	1:B:216:MSE:HA	1.93	0.90
1:A:281:LEU:O	1:A:284:ALA:N	2.07	0.86
1:B:281:LEU:O	1:B:286:ARG:CB	2.26	0.84
1:A:281:LEU:O	1:A:284:ALA:CB	2.26	0.83
1:B:292:ASN:O	1:B:296:GLU:N	2.10	0.83
1:A:248:PHE:O	1:A:251:GLU:N	2.13	0.81
1:B:215:SER:CB	1:B:219:LEU:N	2.47	0.77
1:B:215:SER:CB	1:B:219:LEU:CB	2.63	0.77
1:B:215:SER:CB	1:B:216:MSE:O	2.34	0.76
1:A:216:MSE:O	1:A:219:LEU:N	2.18	0.75
1:B:18:ALA:O	1:B:22:GLY:CA	2.34	0.75
1:B:307:ASP:O	1:B:311:GLN:N	2.15	0.75
1:A:271:SER:O	1:A:274:PHE:CA	2.35	0.74
1:A:31:GLU:O	1:A:34:PHE:CA	2.39	0.70
1:A:271:SER:O	1:A:274:PHE:HA	1.92	0.69
1:B:281:LEU:O	1:B:284:ALA:O	2.12	0.68
1:B:291:THR:O	1:B:294:LEU:N	2.27	0.66
1:A:31:GLU:CB	1:A:34:PHE:CB	2.73	0.66
1:B:281:LEU:C	1:B:286:ARG:CB	2.64	0.66
1:B:268:GLN:O	1:B:273:GLU:CB	2.45	0.65
1:B:215:SER:CB	1:B:216:MSE:C	2.66	0.64
1:B:144:LEU:O	1:B:146:LYS:N	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ASP:O	1:B:206:ASN:N	2.31	0.63
1:A:262:LEU:O	1:A:266:ILE:N	2.33	0.61
1:A:31:GLU:O	1:A:34:PHE:CB	2.50	0.60
1:B:291:THR:O	1:B:295:ASN:N	2.29	0.60
1:B:144:LEU:O	1:B:145:ILE:C	2.40	0.59
1:B:203:ASP:O	1:B:206:ASN:CA	2.51	0.59
1:A:291:THR:O	1:A:295:ASN:N	2.24	0.58
1:A:281:LEU:O	1:A:284:ALA:O	2.22	0.58
1:A:306:ALA:O	1:A:310:ASN:N	2.27	0.57
1:A:351:VAL:O	1:A:353:ARG:N	2.38	0.57
1:B:50:LEU:O	1:B:51:GLY:O	2.22	0.57
1:A:281:LEU:O	1:A:284:ALA:C	2.42	0.56
1:A:281:LEU:HA	1:A:284:ALA:HB3	1.86	0.56
1:A:351:VAL:C	1:A:353:ARG:N	2.58	0.56
1:A:281:LEU:CA	1:A:284:ALA:HB3	2.36	0.56
1:A:281:LEU:C	1:A:284:ALA:HB3	2.26	0.56
1:B:18:ALA:O	1:B:22:GLY:HA3	2.04	0.55
1:A:281:LEU:C	1:A:284:ALA:H	2.10	0.54
1:A:351:VAL:O	1:A:352:GLN:C	2.45	0.54
1:B:279:SER:O	1:B:283:ASN:N	2.41	0.54
1:A:271:SER:C	1:A:274:PHE:H	2.06	0.52
1:A:248:PHE:O	1:A:249:LEU:C	2.48	0.51
1:A:216:MSE:O	1:A:217:ARG:C	2.52	0.48
1:B:144:LEU:O	1:B:147:SER:N	2.43	0.48
1:A:281:LEU:HA	1:A:284:ALA:CB	2.44	0.48
1:B:281:LEU:O	1:B:286:ARG:CA	2.62	0.48
1:A:288:HIS:O	1:A:289:ASN:C	2.53	0.47
1:B:215:SER:CB	1:B:216:MSE:CA	2.92	0.47
1:B:291:THR:O	1:B:292:ASN:C	2.53	0.47
1:B:290:LEU:O	1:B:293:LEU:N	2.38	0.47
1:A:352:GLN:O	1:A:355:VAL:N	2.49	0.46
1:B:144:LEU:C	1:B:146:LYS:N	2.68	0.45
1:B:350:PRO:CB	1:B:353:ARG:CB	2.95	0.45
1:A:354:PHE:O	1:A:358:MSE:HG2	2.17	0.45
1:B:270:GLN:O	1:B:273:GLU:CB	2.65	0.44
1:B:308:LYS:O	1:B:311:GLN:O	2.35	0.43
1:B:291:THR:C	1:B:293:LEU:N	2.70	0.43
1:A:144:LEU:O	1:A:145:ILE:C	2.57	0.42
1:B:354:PHE:O	1:B:358:MSE:HG2	2.19	0.42
1:A:261:VAL:O	1:A:265:LEU:N	2.48	0.41
1:B:390:LYS:O	1:B:394:MSE:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:ASN:O	1:B:296:GLU:CB	2.68	0.41
1:A:351:VAL:C	1:A:353:ARG:H	2.24	0.41
1:B:215:SER:CB	1:B:219:LEU:CA	3.00	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	310/557 (56%)	291 (94%)	11 (4%)	8 (3%)	7	46
1	B	308/557 (55%)	294 (96%)	11 (4%)	3 (1%)	19	65
All	All	618/1114 (56%)	585 (95%)	22 (4%)	11 (2%)	11	54

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	83	ILE
1	A	181	PRO
1	A	317	PRO
1	B	317	PRO
1	A	209	PRO
1	A	350	PRO
1	A	352	GLN
1	B	209	PRO
1	A	115	ILE
1	A	188	PRO
1	B	180	THR

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	5/502 (1%)	5 (100%)	0	100	100
1	B	5/502 (1%)	5 (100%)	0	100	100
All	All	10/1004 (1%)	10 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

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 ⓘ

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	327/557 (58%)	-0.30	13 (3%) 42 35	33, 179, 457, 500	0
1	B	325/557 (58%)	-0.32	4 (1%) 81 74	21, 186, 417, 500	0
All	All	652/1114 (58%)	-0.31	17 (2%) 59 50	21, 184, 438, 500	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	LEU	7.0
1	A	148	ASP	6.3
1	A	212	ASP	5.9
1	A	214	GLU	5.5
1	B	340	SER	4.2
1	A	211	ASN	3.8
1	A	286	ARG	3.7
1	A	285	LYS	3.5
1	A	210	SER	3.3
1	A	147	SER	3.0
1	A	287	PHE	2.7
1	B	341	LEU	2.7
1	A	348	GLY	2.7
1	B	343	LYS	2.5
1	B	409	ILE	2.4
1	A	283	ASN	2.3
1	A	288	HIS	2.1

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 [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.