Full wwPDB X-ray Structure Validation Report (i)

May 16, 2020 – 10:05 am BST

PDB ID	:	5CHE
Title	:	Crystal structure of Arabidopsis glutamyl-tRNA reductase in complex with its regulatory proteins
Authors	:	Fang, Y.; Liu, L.
Deposited on	:	2015-07-10
Resolution	:	3.20 Å(reported)

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

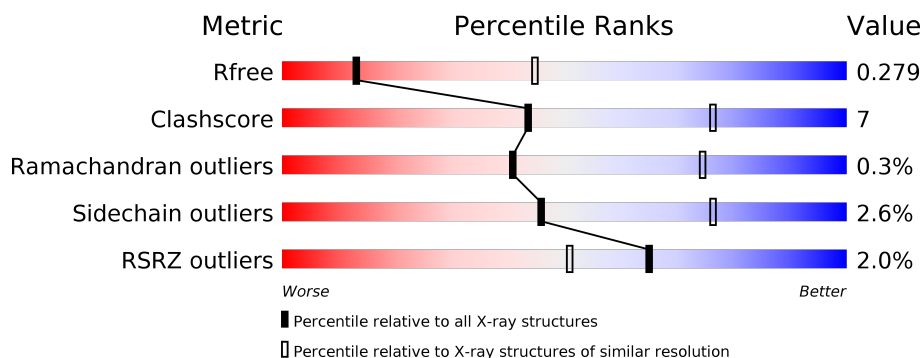|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	Percentile Ranks	Value
Rfree		0.279
Clashscore		7
Ramachandran outliers		0.3%
Sidechain outliers		2.6%
RSRZ outliers		2.0%

Worse   *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 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 <=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	472	76% 14% • 10%
1	B	472	70% 19% • 10%
2	C	310	66% 14% • 19%
2	D	310	66% 14% • 19%
3	E	159	60% 14% 25%
3	F	159	54% 19% 26%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12371 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 reductase 1, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total 3292	C 2055	N 591	O 621	S 25	0	1	0
1	B	423	Total 3243	C 2029	N 586	O 603	S 25	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	SER	-	expression tag	UNP P42804
B	72	SER	-	expression tag	UNP P42804

- Molecule 2 is a protein called Glutamyl-tRNA reductase-binding protein, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	250	Total 1970	C 1246	N 335	O 376	S 13	0	0	0
2	D	250	Total 1992	C 1260	N 339	O 380	S 13	0	1	0

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	8	MET	-	initiating methionine	UNP Q9LU39
C	9	GLY	-	expression tag	UNP Q9LU39
C	10	SER	-	expression tag	UNP Q9LU39
C	11	SER	-	expression tag	UNP Q9LU39
C	12	HIS	-	expression tag	UNP Q9LU39
C	13	HIS	-	expression tag	UNP Q9LU39
C	14	HIS	-	expression tag	UNP Q9LU39
C	15	HIS	-	expression tag	UNP Q9LU39
C	16	HIS	-	expression tag	UNP Q9LU39
C	17	HIS	-	expression tag	UNP Q9LU39

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Chain	Residue	Modelled	Actual	Comment	Reference
C	18	SER	-	expression tag	UNP Q9LU39
C	19	SER	-	expression tag	UNP Q9LU39
C	20	GLY	-	expression tag	UNP Q9LU39
C	21	LEU	-	expression tag	UNP Q9LU39
C	22	VAL	-	expression tag	UNP Q9LU39
C	23	PRO	-	expression tag	UNP Q9LU39
C	24	ARG	-	expression tag	UNP Q9LU39
C	25	GLY	-	expression tag	UNP Q9LU39
C	26	SER	-	expression tag	UNP Q9LU39
C	27	HIS	-	expression tag	UNP Q9LU39
C	28	MET	-	expression tag	UNP Q9LU39
C	29	ALA	-	expression tag	UNP Q9LU39
C	30	SER	-	expression tag	UNP Q9LU39
C	31	MET	-	expression tag	UNP Q9LU39
C	32	THR	-	expression tag	UNP Q9LU39
C	33	GLY	-	expression tag	UNP Q9LU39
C	34	GLY	-	expression tag	UNP Q9LU39
C	35	GLN	-	expression tag	UNP Q9LU39
C	36	GLN	-	expression tag	UNP Q9LU39
C	37	MET	-	expression tag	UNP Q9LU39
C	38	GLY	-	expression tag	UNP Q9LU39
C	39	ARG	-	expression tag	UNP Q9LU39
C	40	GLY	-	expression tag	UNP Q9LU39
C	41	SER	-	expression tag	UNP Q9LU39
D	8	MET	-	initiating methionine	UNP Q9LU39
D	9	GLY	-	expression tag	UNP Q9LU39
D	10	SER	-	expression tag	UNP Q9LU39
D	11	SER	-	expression tag	UNP Q9LU39
D	12	HIS	-	expression tag	UNP Q9LU39
D	13	HIS	-	expression tag	UNP Q9LU39
D	14	HIS	-	expression tag	UNP Q9LU39
D	15	HIS	-	expression tag	UNP Q9LU39
D	16	HIS	-	expression tag	UNP Q9LU39
D	17	HIS	-	expression tag	UNP Q9LU39
D	18	SER	-	expression tag	UNP Q9LU39
D	19	SER	-	expression tag	UNP Q9LU39
D	20	GLY	-	expression tag	UNP Q9LU39
D	21	LEU	-	expression tag	UNP Q9LU39
D	22	VAL	-	expression tag	UNP Q9LU39
D	23	PRO	-	expression tag	UNP Q9LU39
D	24	ARG	-	expression tag	UNP Q9LU39
D	25	GLY	-	expression tag	UNP Q9LU39

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Chain	Residue	Modelled	Actual	Comment	Reference
D	26	SER	-	expression tag	UNP Q9LU39
D	27	HIS	-	expression tag	UNP Q9LU39
D	28	MET	-	expression tag	UNP Q9LU39
D	29	ALA	-	expression tag	UNP Q9LU39
D	30	SER	-	expression tag	UNP Q9LU39
D	31	MET	-	expression tag	UNP Q9LU39
D	32	THR	-	expression tag	UNP Q9LU39
D	33	GLY	-	expression tag	UNP Q9LU39
D	34	GLY	-	expression tag	UNP Q9LU39
D	35	GLN	-	expression tag	UNP Q9LU39
D	36	GLN	-	expression tag	UNP Q9LU39
D	37	MET	-	expression tag	UNP Q9LU39
D	38	GLY	-	expression tag	UNP Q9LU39
D	39	ARG	-	expression tag	UNP Q9LU39
D	40	GLY	-	expression tag	UNP Q9LU39
D	41	SER	-	expression tag	UNP Q9LU39

- Molecule 3 is a protein called Protein FLUORESCENT IN BLUE LIGHT, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	119	Total 951	C 599	N 160	O 190	S 2	0	1	0
3	F	117	Total 917	C 579	N 157	O 179	S 2	0	0	0

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	159	MET	-	initiating methionine	UNP Q940U6
E	160	LYS	-	expression tag	UNP Q940U6
E	161	TYR	-	expression tag	UNP Q940U6
E	162	LEU	-	expression tag	UNP Q940U6
E	163	LEU	-	expression tag	UNP Q940U6
E	164	PRO	-	expression tag	UNP Q940U6
E	165	THR	-	expression tag	UNP Q940U6
E	166	ALA	-	expression tag	UNP Q940U6
E	167	ALA	-	expression tag	UNP Q940U6
E	168	ALA	-	expression tag	UNP Q940U6
E	169	GLY	-	expression tag	UNP Q940U6
E	170	LEU	-	expression tag	UNP Q940U6
E	171	LEU	-	expression tag	UNP Q940U6
E	172	LEU	-	expression tag	UNP Q940U6

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Chain	Residue	Modelled	Actual	Comment	Reference
E	173	LEU	-	expression tag	UNP Q940U6
E	174	ALA	-	expression tag	UNP Q940U6
E	175	ALA	-	expression tag	UNP Q940U6
E	176	GLN	-	expression tag	UNP Q940U6
E	177	PRO	-	expression tag	UNP Q940U6
E	178	ALA	-	expression tag	UNP Q940U6
E	179	MET	-	expression tag	UNP Q940U6
E	180	ALA	-	expression tag	UNP Q940U6
E	181	MET	-	expression tag	UNP Q940U6
E	182	ASP	-	expression tag	UNP Q940U6
E	183	ILE	-	expression tag	UNP Q940U6
E	184	GLY	-	expression tag	UNP Q940U6
E	185	ILE	-	expression tag	UNP Q940U6
E	186	ASN	-	expression tag	UNP Q940U6
E	187	SER	-	expression tag	UNP Q940U6
E	188	ASP	-	expression tag	UNP Q940U6
E	189	PRO	-	expression tag	UNP Q940U6
E	190	HIS	-	expression tag	UNP Q940U6
E	191	HIS	-	expression tag	UNP Q940U6
E	192	HIS	-	expression tag	UNP Q940U6
E	193	HIS	-	expression tag	UNP Q940U6
E	194	HIS	-	expression tag	UNP Q940U6
E	195	HIS	-	expression tag	UNP Q940U6
F	159	MET	-	initiating methionine	UNP Q940U6
F	160	LYS	-	expression tag	UNP Q940U6
F	161	TYR	-	expression tag	UNP Q940U6
F	162	LEU	-	expression tag	UNP Q940U6
F	163	LEU	-	expression tag	UNP Q940U6
F	164	PRO	-	expression tag	UNP Q940U6
F	165	THR	-	expression tag	UNP Q940U6
F	166	ALA	-	expression tag	UNP Q940U6
F	167	ALA	-	expression tag	UNP Q940U6
F	168	ALA	-	expression tag	UNP Q940U6
F	169	GLY	-	expression tag	UNP Q940U6
F	170	LEU	-	expression tag	UNP Q940U6
F	171	LEU	-	expression tag	UNP Q940U6
F	172	LEU	-	expression tag	UNP Q940U6
F	173	LEU	-	expression tag	UNP Q940U6
F	174	ALA	-	expression tag	UNP Q940U6
F	175	ALA	-	expression tag	UNP Q940U6
F	176	GLN	-	expression tag	UNP Q940U6
F	177	PRO	-	expression tag	UNP Q940U6

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Chain	Residue	Modelled	Actual	Comment	Reference
F	178	ALA	-	expression tag	UNP Q940U6
F	179	MET	-	expression tag	UNP Q940U6
F	180	ALA	-	expression tag	UNP Q940U6
F	181	MET	-	expression tag	UNP Q940U6
F	182	ASP	-	expression tag	UNP Q940U6
F	183	ILE	-	expression tag	UNP Q940U6
F	184	GLY	-	expression tag	UNP Q940U6
F	185	ILE	-	expression tag	UNP Q940U6
F	186	ASN	-	expression tag	UNP Q940U6
F	187	SER	-	expression tag	UNP Q940U6
F	188	ASP	-	expression tag	UNP Q940U6
F	189	PRO	-	expression tag	UNP Q940U6
F	190	HIS	-	expression tag	UNP Q940U6
F	191	HIS	-	expression tag	UNP Q940U6
F	192	HIS	-	expression tag	UNP Q940U6
F	193	HIS	-	expression tag	UNP Q940U6
F	194	HIS	-	expression tag	UNP Q940U6
F	195	HIS	-	expression tag	UNP Q940U6

- Molecule 4 is water.

| Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total O 2 2	0	0
4	B	2	Total O 2 2	0	0
4	C	2	Total O 2 2	0	0

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 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.
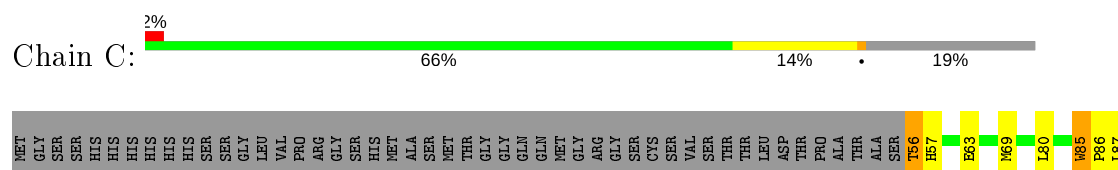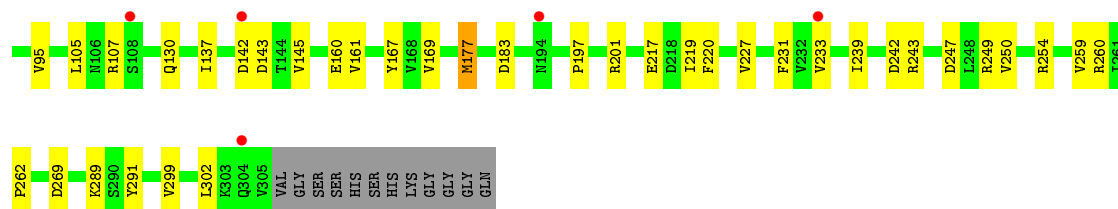
- Molecule 1: Glutamyl-tRNA reductase 1, chloroplastic
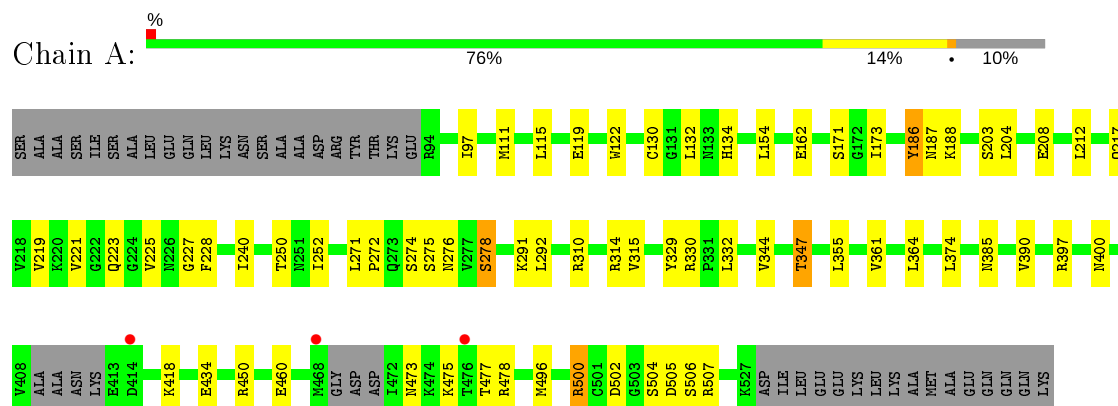
Chain A:

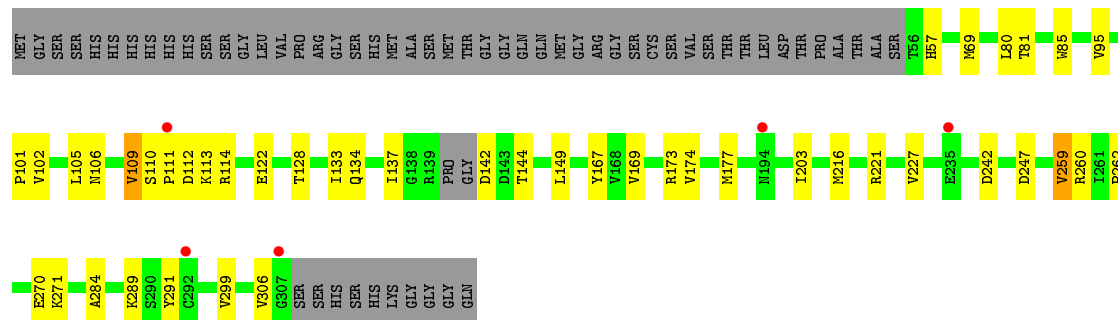- Molecule 1: Glutamyl-tRNA reductase 1, chloroplastic

Chain B:
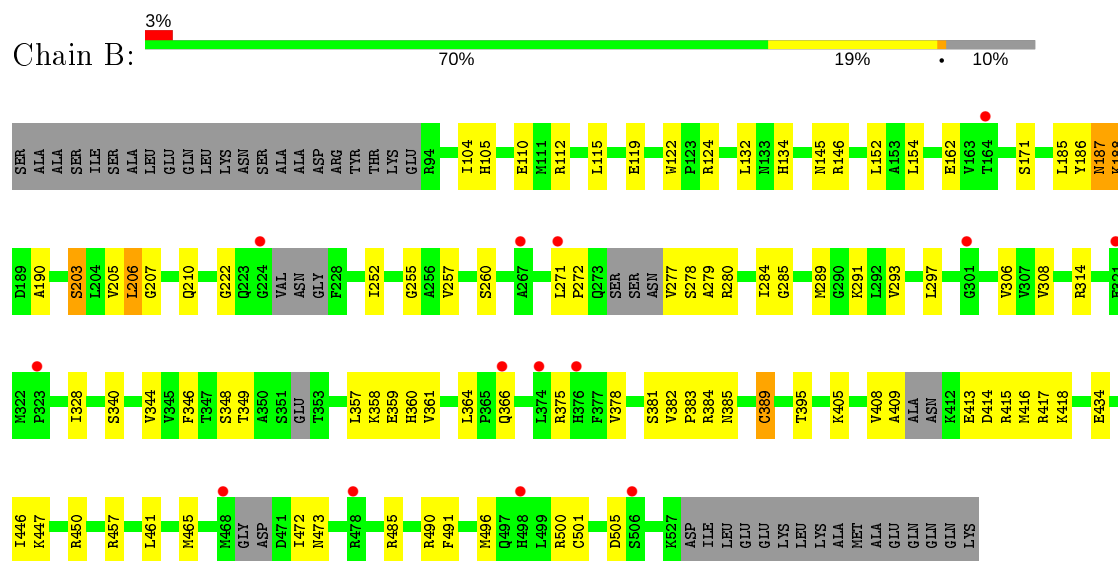
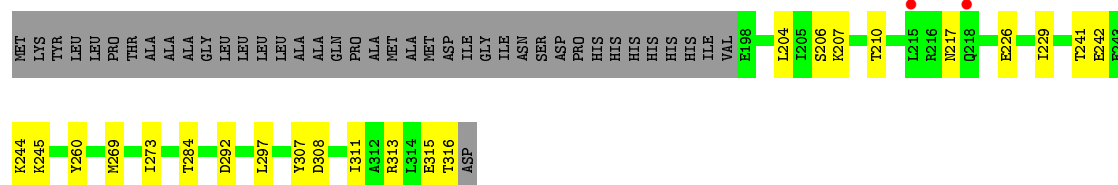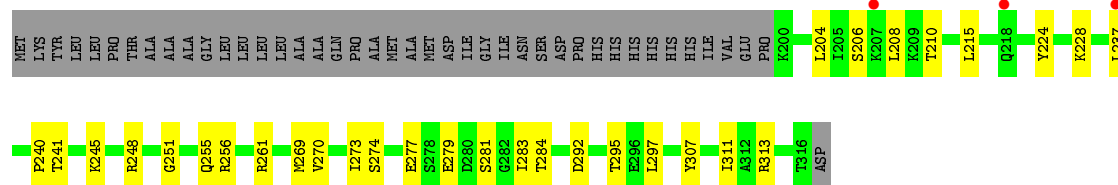- Molecule 2: Glutamyl-tRNA reductase-binding protein, chloroplastic

Chain C:

• Molecule 2: Glutamyl-tRNA reductase-binding protein, chloroplastic

Chain D:

• Molecule 3: Protein FLUORESCENT IN BLUE LIGHT, chloroplastic

Chain E:

• Molecule 3: Protein FLUORESCENT IN BLUE LIGHT, chloroplastic

Chain F:

4   Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	216.99Å 53.21Å 203.76Å 90.00° 108.36° 90.00°	Depositor
Resolution (Å)	39.27 – 3.20 48.23 – 3.20	Depositor EDS
% Data completeness (in resolution range)	87.2 (39.27-3.20) 87.4 (48.23-3.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.00 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.223 , 0.278 0.225 , 0.279	Depositor DCC
R_{free} test set	1636 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	58.7	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 5.9	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	12371	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1] Intensities estimated from amplitudes.
[2] Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.32	0/3338	0.52	3/4501 (0.1%)
1	B	0.33	0/3283	0.57	6/4423 (0.1%)
2	C	0.29	0/2013	0.45	0/2731
2	D	0.31	0/2034	0.46	1/2755 (0.0%)
3	E	0.30	0/967	0.47	0/1294
3	F	0.26	0/929	0.44	0/1243
All	All	0.31	0/12564	0.50	10/16947 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	187	ASN	CB-CA-C	9.99	130.39	110.40
1	A	502	ASP	N-CA-C	-8.89	86.98	111.00
1	B	188	LYS	N-CA-CB	-8.62	95.08	110.60
1	A	186	TYR	N-CA-C	7.29	130.67	111.00
1	B	186	TYR	CB-CA-C	-5.86	98.67	110.40
1	B	222	GLY	N-CA-C	5.57	127.03	113.10
1	A	228	PHE	N-CA-C	-5.34	96.59	111.00
1	B	188	LYS	N-CA-C	-5.14	97.12	111.00
1	B	280	ARG	N-CA-C	5.04	124.60	111.00
2	D	306	VAL	N-CA-C	5.02	124.55	111.00

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	3292	0	3363	37	0
1	B	3243	0	3309	54	0
2	C	1970	0	1928	27	0
2	D	1992	0	1955	24	0
3	E	951	0	959	16	0
3	F	917	0	926	18	0
4	A	2	0	0	1	0
4	B	2	0	0	0	0
4	C	2	0	0	1	0
All	All	12371	0	12440	166	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 7.

All (166) 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:252:ILE:HG22	1:B:418:LYS:HB3	1.61	0.80
1:A:271:LEU:HD11	1:A:276:ASN:HA	1.65	0.78
2:C:80:LEU:HD21	2:D:80:LEU:HD21	1.66	0.78
2:D:227:VAL:HG12	2:D:299:VAL:HB	1.67	0.77
3:F:215:LEU:O	3:F:256:ARG:NH1	2.21	0.73
1:A:504:SER:OG	1:A:505:ASP:N	2.22	0.73
1:A:475:LYS:HG3	1:A:478:ARG:HH21	1.54	0.72
1:A:450:ARG:NH1	3:E:292:ASP:OD2	2.24	0.71
1:A:496:MET:HB3	3:E:311:ILE:HD12	1.71	0.70
1:B:450:ARG:NH1	3:F:292:ASP:OD2	2.26	0.68
1:A:203:SER:HB3	1:A:204:LEU:HD23	1.75	0.68
2:C:183:ASP:OD1	2:D:106:ASN:ND2	2.26	0.68
3:F:204:LEU:HD12	3:F:237:LEU:HD12	1.76	0.67
1:B:291:LYS:HE3	1:B:314:ARG:HH12	1.60	0.67
2:D:69:MET:HB3	2:D:95:VAL:HG21	1.77	0.67
1:B:306:VAL:HG22	1:B:328:ILE:HD11	1.75	0.67
2:C:227:VAL:HG12	2:C:299:VAL:HB	1.76	0.66
2:C:69:MET:HB3	2:C:95:VAL:HG21	1.77	0.65
1:A:310:ARG:HG2	1:A:332:LEU:HD21	1.78	0.65
1:A:115:LEU:HD13	1:A:171:SER:HB3	1.80	0.64
1:A:130:CYS:O	4:A:601:HOH:O	2.15	0.64
2:D:242:ASP:OD1	2:D:260:ARG:NH2	2.30	0.62
1:B:496:MET:HB3	3:F:311:ILE:HD12	1.82	0.62
1:B:382:VAL:HA	1:B:383:PRO:C	2.20	0.61
3:F:261:ARG:HH21	3:F:297:LEU:HD13	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:457:ARG:HD2	1:B:485:ARG:HG2	1.82	0.61
1:B:252:ILE:HA	1:B:418:LYS:HD3	1.82	0.61
3:E:284:THR:HG22	3:E:313:ARG:HE	1.66	0.60
2:D:101:PRO:HG2	2:D:169:VAL:HB	1.83	0.60
2:C:239:ILE:HD11	2:C:249:ARG:HB2	1.82	0.60
1:B:278:SER:OG	1:B:279:ALA:N	2.33	0.60
3:F:279:GLU:HG3	3:F:281:SER:H	1.67	0.59
1:B:349:THR:OG1	1:B:385:ASN:ND2	2.35	0.59
1:B:366:GLN:HG2	1:B:395:THR:HG21	1.84	0.58
1:B:188:LYS:HD3	1:B:434:GLU:HG3	1.86	0.57
1:A:188:LYS:HE2	1:A:434:GLU:HG2	1.86	0.57
1:B:145:ASN:OD1	1:B:146:ARG:NH1	2.38	0.57
3:F:248:ARG:HG3	3:F:270:VAL:HG11	1.86	0.57
2:D:109:VAL:HG13	2:D:167:TYR:HE1	1.70	0.57
1:A:460:GLU:OE2	1:B:490:ARG:NH1	2.33	0.56
3:E:307:TYR:O	3:E:311:ILE:HG12	2.06	0.56
3:E:292:ASP:OD1	3:E:307:TYR:OH	2.18	0.55
3:E:260:TYR:HB2	3:E:297:LEU:HD21	1.88	0.55
1:B:277:VAL:HA	1:B:278:SER:HB3	1.87	0.54
2:C:269:ASP:OD1	2:C:269:ASP:N	2.40	0.54
1:B:124:ARG:HB2	1:B:124:ARG:NH1	2.24	0.53
1:B:115:LEU:HD13	1:B:171:SER:HB3	1.89	0.53
2:D:203:ILE:HD13	2:D:270:GLU:HB2	1.90	0.53
2:C:160:GLU:HG3	2:C:161:VAL:H	1.72	0.53
1:B:408:VAL:HG12	1:B:409:ALA:H	1.74	0.53
1:B:134:HIS:ND1	1:B:162:GLU:OE2	2.42	0.52
1:A:275:SER:HB3	1:A:278:SER:HB3	1.92	0.52
1:B:340:SER:HA	1:B:375:ARG:HH12	1.75	0.52
1:A:97:ILE:O	1:A:186:TYR:O	2.27	0.52
1:B:381:SER:O	1:B:384:ARG:HA	2.09	0.52
2:C:231:PHE:HD1	2:C:254:ARG:HD2	1.75	0.51
2:D:57:HIS:CG	2:D:262:PRO:HB3	2.46	0.51
1:B:185:LEU:HD11	2:D:271:LYS:HG3	1.92	0.51
1:A:500:ARG:HG2	3:E:308:ASP:OD1	2.10	0.50
1:B:446:ILE:HG23	1:B:496:MET:SD	2.50	0.50
2:D:122:GLU:HA	2:D:128:THR:HG22	1.94	0.50
2:D:247:ASP:OD2	2:D:260:ARG:NH1	2.43	0.50
2:C:85:TRP:CD1	2:C:86:PRO:HD2	2.46	0.50
1:B:252:ILE:CG2	1:B:418:LYS:HB3	2.36	0.50
2:C:231:PHE:HB3	2:C:254:ARG:HE	1.77	0.50
1:B:358:LYS:HE2	1:B:389:CYS:SG	2.51	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:259:VAL:HG12	2:C:302:LEU:HD22	1.94	0.49
2:D:80:LEU:HD12	2:D:112:ASP:OD2	2.13	0.49
2:C:231:PHE:CD1	2:C:254:ARG:HD2	2.47	0.49
3:E:241:THR:HG22	3:E:245:LYS:HE3	1.95	0.49
1:B:361:VAL:HG13	1:B:364:LEU:HD12	1.94	0.48
2:D:102:VAL:HG21	2:D:149:LEU:HD22	1.94	0.48
1:A:252:ILE:HA	1:A:418:LYS:HD3	1.96	0.48
2:C:137:ILE:HG12	2:C:169:VAL:HG22	1.96	0.48
2:D:105:LEU:HB2	2:D:167:TYR:CE2	2.49	0.48
2:C:233:VAL:HG13	2:C:250:VAL:HG13	1.95	0.48
1:B:413:GLU:HG3	1:B:414:ASP:H	1.79	0.48
3:F:269:MET:O	3:F:273:ILE:HG13	2.13	0.47
1:A:315:VAL:HG21	1:A:329:TYR:HD1	1.79	0.47
3:E:315:GLU:N	3:E:316:THR:HA	2.28	0.47
2:C:130:GLN:HG3	2:C:177:MET:HG3	1.97	0.47
1:B:409:ALA:HB2	1:B:415:ARG:NH2	2.30	0.47
3:F:251:GLY:O	3:F:255:GLN:HG3	2.15	0.47
3:E:206:SER:O	3:E:210:THR:OG1	2.26	0.47
3:F:284:THR:OG1	3:F:313:ARG:NE	2.40	0.46
2:C:143:ASP:OD1	2:C:145:VAL:HG22	2.15	0.46
1:A:134:HIS:HA	1:A:154:LEU:HD23	1.97	0.46
1:B:124:ARG:HB2	1:B:124:ARG:HH11	1.80	0.46
1:B:361:VAL:HA	1:B:364:LEU:HG	1.96	0.46
1:B:132:LEU:HD13	1:B:162:GLU:HB3	1.97	0.46
3:F:224:TYR:CE2	3:F:228:LYS:HD2	2.50	0.46
1:B:285:GLY:HA3	1:B:348:SER:O	2.16	0.46
1:B:289:MET:O	1:B:293:VAL:HG23	2.16	0.46
1:A:374:LEU:HD13	1:A:397:ARG:HE	1.80	0.45
1:B:408:VAL:HG12	1:B:409:ALA:N	2.31	0.45
1:B:105:HIS:HE2	1:B:203:SER:HB3	1.81	0.45
1:A:332:LEU:HD22	1:A:332:LEU:H	1.81	0.45
1:A:208:GLU:HG3	1:A:292:LEU:HD21	1.98	0.45
1:B:472:ILE:HA	1:B:473:ASN:HA	1.79	0.45
3:E:207:LYS:HZ1	3:E:229:ILE:HD12	1.81	0.45
3:F:206:SER:O	3:F:210:THR:OG1	2.20	0.45
1:A:347:THR:OG1	1:A:385:ASN:HB2	2.17	0.45
3:F:307:TYR:O	3:F:311:ILE:HG12	2.17	0.45
2:C:289:LYS:HB3	2:C:291:TYR:CE2	2.51	0.45
2:D:221:ARG:HG2	2:D:284:ALA:HB1	1.99	0.45
3:E:269:MET:O	3:E:273:ILE:HG13	2.17	0.45
3:E:207:LYS:NZ	3:E:226:GLU:OE1	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:VAL:O	1:A:223:GLN:HG2	2.17	0.44
1:A:225:VAL:HG12	1:A:227:GLY:H	1.82	0.44
1:A:329:TYR:O	1:A:330:ARG:HD2	2.18	0.44
1:B:284:ILE:HG13	1:B:308:VAL:HB	1.99	0.44
2:C:56:THR:HG22	2:C:57:HIS:H	1.82	0.44
3:F:274:SER:HB3	3:F:279:GLU:O	2.18	0.44
1:B:409:ALA:HB2	1:B:415:ARG:HH22	1.83	0.44
2:D:110:SER:HB2	2:D:111:PRO:HD2	1.99	0.44
1:A:217:GLN:O	1:A:221:VAL:HG23	2.18	0.44
1:B:187:ASN:O	1:B:190:ALA:HB3	2.18	0.44
2:D:109:VAL:HG13	2:D:167:TYR:CE1	2.49	0.44
3:F:241:THR:HG22	3:F:245:LYS:HE3	2.00	0.44
1:A:212:LEU:HD11	1:A:240:ILE:HG23	2.00	0.43
1:A:473:ASN:HB2	1:A:477:THR:OG1	2.18	0.43
1:A:111:MET:HG3	1:A:173:ILE:HD13	2.00	0.43
1:B:104:ILE:HA	1:B:112:ARG:NH1	2.34	0.43
3:E:204:LEU:HD21	3:E:242:GLU:HB3	2.00	0.43
1:A:132:LEU:HD13	1:A:162:GLU:HB3	2.01	0.43
1:A:119:GLU:HA	1:A:122:TRP:CD2	2.54	0.43
1:B:346:PHE:HD1	1:B:378:VAL:HB	1.83	0.43
1:A:355:LEU:N	1:A:385:ASN:O	2.38	0.43
1:A:505:ASP:O	1:A:507:ARG:N	2.52	0.43
2:C:217:GLU:OE1	4:C:401:HOH:O	2.21	0.42
2:C:63:GLU:OE1	2:C:243:ARG:NH2	2.36	0.42
2:C:197:PRO:HB2	2:C:201:ARG:NH1	2.35	0.42
2:D:289:LYS:HB3	2:D:291:TYR:CE2	2.54	0.42
3:F:240:PRO:HB3	3:F:277:GLU:HG3	2.01	0.42
2:C:105:LEU:HG	2:C:107:ARG:H	1.83	0.42
2:D:134:GLN:HE22	2:D:173:ARG:HH21	1.66	0.42
3:F:283:ILE:HB	3:F:313:ARG:NH2	2.35	0.42
1:B:293:VAL:O	1:B:297:LEU:HB2	2.19	0.42
1:B:105:HIS:HA	1:B:405:LYS:NZ	2.35	0.42
1:B:447:LYS:HD3	3:F:295:THR:HB	2.00	0.42
2:D:142:ASP:OD2	2:D:144:THR:OG1	2.26	0.42
2:D:133:ILE:HG12	2:D:174:VAL:HG22	2.00	0.42
1:B:207:GLY:O	1:B:210:GLN:HG3	2.20	0.42
1:A:361:VAL:HA	1:A:364:LEU:HG	2.01	0.41
1:B:271:LEU:HA	1:B:272:PRO:HD2	1.77	0.41
2:C:57:HIS:HB3	2:C:262:PRO:HB3	2.01	0.41
1:B:346:PHE:CD1	1:B:378:VAL:HB	2.54	0.41
1:A:275:SER:HB3	1:A:278:SER:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:GLU:HA	1:B:122:TRP:CE2	2.56	0.41
3:E:241:THR:HA	3:E:244:LYS:HE2	2.02	0.41
1:B:206:LEU:HB3	1:B:210:GLN:HB2	2.02	0.41
1:B:382:VAL:CA	1:B:383:PRO:C	2.89	0.41
2:C:247:ASP:OD1	2:C:260:ARG:HD3	2.20	0.41
1:A:272:PRO:C	1:A:274:SER:H	2.24	0.40
1:B:500:ARG:HD2	1:B:500:ARG:HA	1.97	0.40
1:A:272:PRO:O	1:A:274:SER:N	2.54	0.40
1:A:291:LYS:CE	1:A:314:ARG:HH22	2.33	0.40
1:B:461:LEU:O	1:B:465:MET:HG2	2.22	0.40
2:C:137:ILE:HG23	2:C:167:TYR:HB3	2.03	0.40
2:D:227:VAL:HG11	2:D:259:VAL:HG11	2.02	0.40
1:B:357:LEU:HD23	1:B:360:HIS:CE1	2.55	0.40
2:C:242:ASP:OD1	2:C:260:ARG:NH2	2.38	0.40
1:B:134:HIS:HA	1:B:154:LEU:HD23	2.02	0.40
2:C:219:ILE:HD13	2:C:219:ILE:HA	1.96	0.40
2:D:113:LYS:O	2:D:137:ILE:HG22	2.21	0.40
3:E:284:THR:HG22	3:E:313:ARG:NE	2.33	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	422/472 (89%)	406 (96%)	14 (3%)	2 (0%)	29	67
1	B	411/472 (87%)	392 (95%)	16 (4%)	3 (1%)	22	61
2	C	248/310 (80%)	241 (97%)	7 (3%)	0	100	100
2	D	247/310 (80%)	241 (98%)	6 (2%)	0	100	100
3	E	118/159 (74%)	115 (98%)	3 (2%)	0	100	100
3	F	115/159 (72%)	114 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1561/1882 (83%)	1509 (97%)	47 (3%)	5 (0%)	41 74

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	505	ASP
1	A	506	SER
1	B	203	SER
1	A	187	ASN
1	B	255	GLY

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	363/400 (91%)	356 (98%)	7 (2%)	57 81
1	B	352/400 (88%)	339 (96%)	13 (4%)	34 68
2	C	220/269 (82%)	214 (97%)	6 (3%)	44 75
2	D	224/269 (83%)	217 (97%)	7 (3%)	40 72
3	E	99/129 (77%)	98 (99%)	1 (1%)	76 90
3	F	93/129 (72%)	92 (99%)	1 (1%)	73 88
All	All	1351/1596 (85%)	1316 (97%)	35 (3%)	46 76

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	250	THR
1	A	278	SER
1	A	344	VAL
1	A	347	THR
1	A	390	VAL
1	A	400	ASN
1	A	500	ARG

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Mol	Chain	Res	Type
1	B	110	GLU
1	B	152	LEU
1	B	205	VAL
1	B	206	LEU
1	B	257	VAL
1	B	260	SER
1	B	344	VAL
1	B	359	GLU
1	B	389	CYS
1	B	416	MET
1	B	417	ARG
1	B	491	PHE
1	B	501	CYS
2	C	56	THR
2	C	85	TRP
2	C	87	LEU
2	C	142	ASP
2	C	177	MET
2	C	220	PHE
2	D	81	THR
2	D	85	TRP
2	D	109	VAL
2	D	114	ARG
2	D	177	MET
2	D	216	MET
2	D	259	VAL
3	E	217	ASN
3	F	208	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	385	ASN
2	D	120	GLN
2	D	134	GLN

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 (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	427/472 (90%)	-0.12	3 (0%)	87	81	2, 12, 45, 90	0
1	B	423/472 (89%)	0.18	14 (3%)	46	30	5, 25, 63, 85	0
2	C	250/310 (80%)	0.15	5 (2%)	65	51	4, 21, 59, 83	0
2	D	250/310 (80%)	0.06	5 (2%)	65	51	6, 24, 65, 87	0
3	E	119/159 (74%)	-0.04	2 (1%)	70	57	6, 22, 47, 58	0
3	F	117/159 (73%)	0.15	3 (2%)	56	40	17, 48, 71, 83	0
All	All	1586/1882 (84%)	0.06	32 (2%)	65	51	2, 21, 64, 90	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	218	GLN	4.2
1	A	476	THR	4.0
2	D	194	ASN	3.4
1	A	414	ASP	3.4
3	E	218	GLN	3.2
2	D	307	GLY	2.9
1	B	506	SER	2.8
2	C	194	ASN	2.8
2	C	142	ASP	2.8
1	B	321	GLU	2.8
2	C	233	VAL	2.7
1	B	366	GLN	2.6
2	D	292	CYS	2.5
1	B	267	ALA	2.5
1	A	468	MET	2.4
1	B	468	MET	2.4
1	B	374	LEU	2.4
2	C	108	SER	2.4
1	B	323	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	D	235	GLU	2.3
1	B	301	GLY	2.3
2	C	304	GLN	2.3
3	E	215	LEU	2.2
1	B	376	HIS	2.2
3	F	237	LEU	2.2
1	B	164	THR	2.2
1	B	478	ARG	2.1
1	B	498	HIS	2.1
1	B	224	GLY	2.1
3	F	207	LYS	2.1
2	D	111	PRO	2.0
1	B	271	LEU	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 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.