



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:09 PM GMT

PDB ID : 3WA0  
Title : Crystal structure of merlin complexed with DCAF1/VprBP  
Authors : Mori, T.; Gotoh, S.; Shirakawa, M.; Hakoshima, T.  
Deposited on : 2013-04-20  
Resolution : 2.31 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](mailto: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.

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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

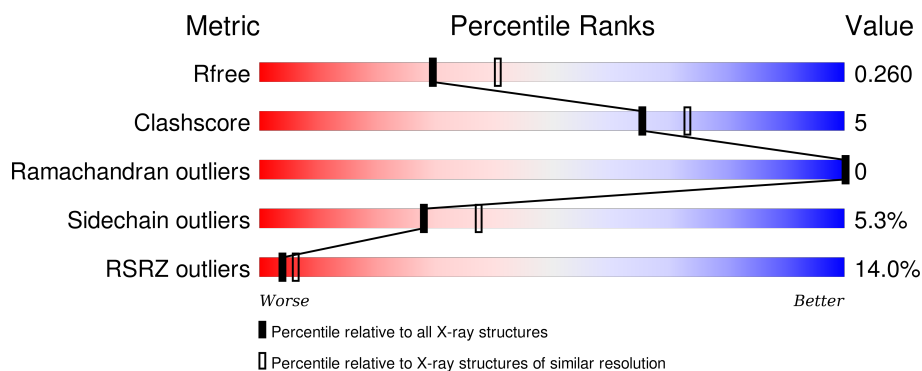
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.31 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	4425 (2.34-2.30)
Clashscore	102246	5057 (2.34-2.30)
Ramachandran outliers	100387	5008 (2.34-2.30)
Sidechain outliers	100360	5007 (2.34-2.30)
RSRZ outliers	91569	4432 (2.34-2.30)

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  $\geq 3$ , 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	301	<div> <div>15%</div> <div>81%</div> <div>12%</div> <div>• •</div> </div>
1	B	301	<div> <div>14%</div> <div>81%</div> <div>13%</div> <div>• •</div> </div>
1	C	301	<div> <div>9%</div> <div>83%</div> <div>14%</div> <div>• •</div> </div>
1	D	301	<div> <div>15%</div> <div>83%</div> <div>15%</div> <div>• •</div> </div>
1	E	301	<div> <div>12%</div> <div>85%</div> <div>11%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	301	<div><div></div><div>15%</div><div>82%</div><div>12%</div><div></div><div></div></div>
2	G	96	<div><div></div><div>3%</div><div>7%</div><div></div><div>89%</div><div></div></div>
2	H	96	<div><div></div><div>3%</div><div>8%</div><div></div><div>89%</div><div></div></div>
3	I	7	<div><div></div><div></div><div>100%</div><div></div></div>
4	J	8	<div><div></div><div></div><div>88%</div><div>13%</div><div></div></div>
5	K	6	<div><div></div><div></div><div>100%</div><div></div></div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15373 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 Merlin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2409	1559	402	435	13			
1	B	290	Total	C	N	O	S	0	0	0
			2412	1562	400	437	13			
1	C	293	Total	C	N	O	S	0	0	0
			2435	1575	406	441	13			
1	D	295	Total	C	N	O	S	0	0	0
			2457	1592	409	443	13			
1	E	292	Total	C	N	O	S	0	0	0
			2429	1572	404	440	13			
1	F	291	Total	C	N	O	S	0	0	0
			2427	1572	403	439	13			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	GLY	-	EXPRESSION TAG	UNP P46662
A	15	PRO	-	EXPRESSION TAG	UNP P46662
A	16	LEU	-	EXPRESSION TAG	UNP P46662
A	17	GLY	-	EXPRESSION TAG	UNP P46662
A	18	SER	-	EXPRESSION TAG	UNP P46662
B	14	GLY	-	EXPRESSION TAG	UNP P46662
B	15	PRO	-	EXPRESSION TAG	UNP P46662
B	16	LEU	-	EXPRESSION TAG	UNP P46662
B	17	GLY	-	EXPRESSION TAG	UNP P46662
B	18	SER	-	EXPRESSION TAG	UNP P46662
C	14	GLY	-	EXPRESSION TAG	UNP P46662
C	15	PRO	-	EXPRESSION TAG	UNP P46662
C	16	LEU	-	EXPRESSION TAG	UNP P46662
C	17	GLY	-	EXPRESSION TAG	UNP P46662
C	18	SER	-	EXPRESSION TAG	UNP P46662
D	14	GLY	-	EXPRESSION TAG	UNP P46662
D	15	PRO	-	EXPRESSION TAG	UNP P46662

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Chain	Residue	Modelled	Actual	Comment	Reference
D	16	LEU	-	EXPRESSION TAG	UNP P46662
D	17	GLY	-	EXPRESSION TAG	UNP P46662
D	18	SER	-	EXPRESSION TAG	UNP P46662
E	14	GLY	-	EXPRESSION TAG	UNP P46662
E	15	PRO	-	EXPRESSION TAG	UNP P46662
E	16	LEU	-	EXPRESSION TAG	UNP P46662
E	17	GLY	-	EXPRESSION TAG	UNP P46662
E	18	SER	-	EXPRESSION TAG	UNP P46662
F	14	GLY	-	EXPRESSION TAG	UNP P46662
F	15	PRO	-	EXPRESSION TAG	UNP P46662
F	16	LEU	-	EXPRESSION TAG	UNP P46662
F	17	GLY	-	EXPRESSION TAG	UNP P46662
F	18	SER	-	EXPRESSION TAG	UNP P46662

- Molecule 2 is a protein called Protein VPRBP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	11	Total	C	N	O	0	0	0
			80	50	11	19			
2	H	11	Total	C	N	O	0	0	0
			76	48	12	16			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	1411	GLY	-	EXPRESSION TAG	UNP Q9Y4B6
G	1412	PRO	-	EXPRESSION TAG	UNP Q9Y4B6
G	1413	LEU	-	EXPRESSION TAG	UNP Q9Y4B6
G	1414	GLY	-	EXPRESSION TAG	UNP Q9Y4B6
G	1415	SER	-	EXPRESSION TAG	UNP Q9Y4B6
G	1416	TYR	-	EXPRESSION TAG	UNP Q9Y4B6
H	1411	GLY	-	EXPRESSION TAG	UNP Q9Y4B6
H	1412	PRO	-	EXPRESSION TAG	UNP Q9Y4B6
H	1413	LEU	-	EXPRESSION TAG	UNP Q9Y4B6
H	1414	GLY	-	EXPRESSION TAG	UNP Q9Y4B6
H	1415	SER	-	EXPRESSION TAG	UNP Q9Y4B6
H	1416	TYR	-	EXPRESSION TAG	UNP Q9Y4B6

- Molecule 3 is a protein called Protein VPRBP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	I	7	Total	C	N	O	0	0	0
			35	21	7	7			

- Molecule 4 is a protein called Protein VPRBP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	J	8	Total	C	N	O	0	0	0
			40	24	8	8			

- Molecule 5 is a protein called Protein VPRBP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	K	6	Total	C	N	O	0	0	0
			30	18	6	6			

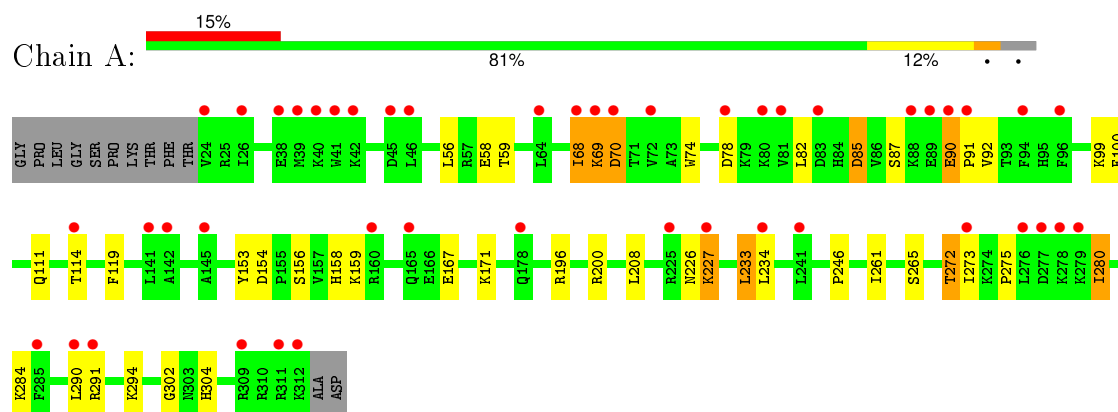
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	70	Total	O	0	0
			70	70		
6	B	85	Total	O	0	0
			85	85		
6	C	102	Total	O	0	0
			102	102		
6	D	94	Total	O	0	0
			94	94		
6	E	86	Total	O	0	0
			86	86		
6	F	105	Total	O	0	0
			105	105		
6	H	1	Total	O	0	0
			1	1		

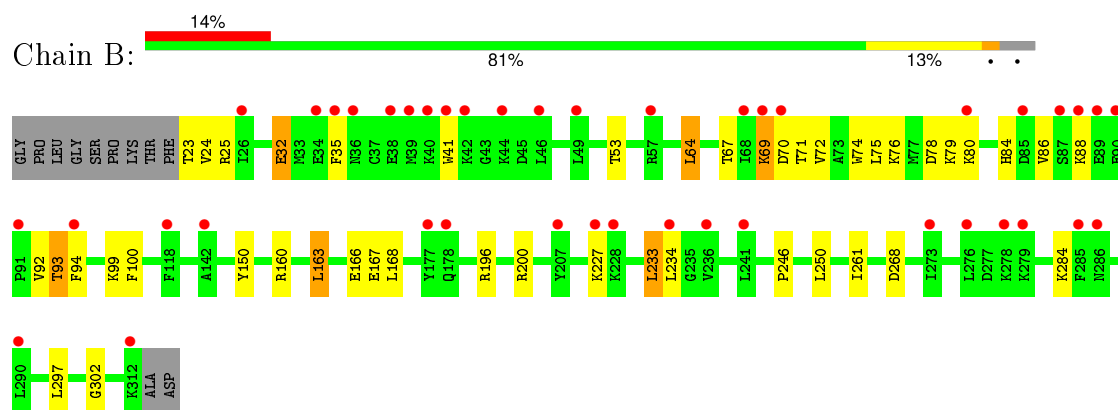
### 3 Residue-property plots [i](#)

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.

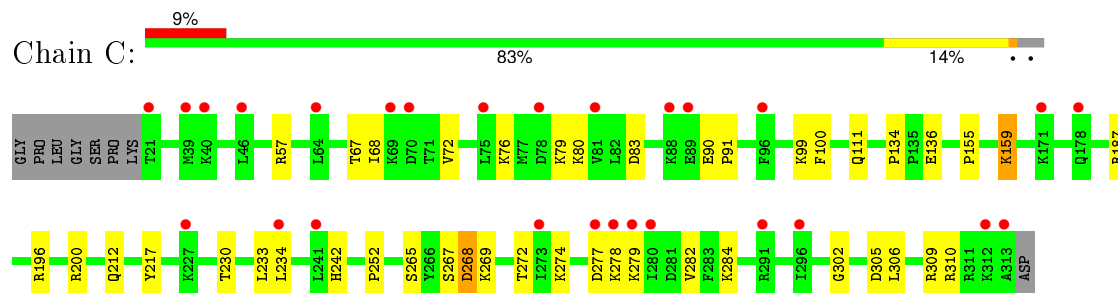
#### • Molecule 1: Merlin



#### • Molecule 1: Merlin

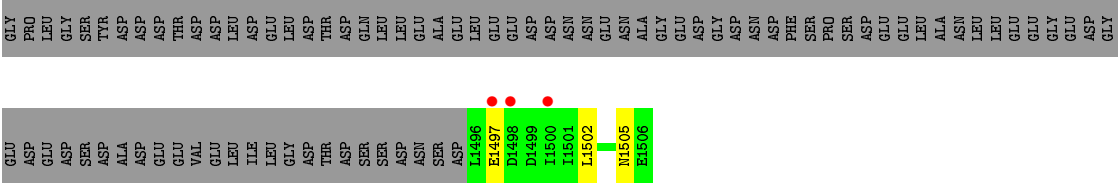


#### • Molecule 1: Merlin



#### • Molecule 1: Merlin





• Molecule 3: Protein VPRBP

Chain I: 100%

There are no outlier residues recorded for this chain.

• Molecule 4: Protein VPRBP

Chain J: 88% 13%



• Molecule 5: Protein VPRBP

Chain K: 100%

There are no outlier residues recorded for this chain.

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.36Å 135.76Å 238.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.07 – 2.31 20.07 – 2.31	Depositor EDS
% Data completeness (in resolution range)	98.8 (20.07-2.31) 99.0 (20.07-2.31)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.71 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.235 , 0.264 0.230 , 0.260	Depositor DCC
$R_{free}$ test set	6149 reflections (5.37%)	DCC
Wilson B-factor (Å <sup>2</sup> )	52.2	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 55.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 120576 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	15373	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.35	0/2467	0.52	0/3328
1	B	0.35	0/2470	0.51	0/3334
1	C	0.38	0/2494	0.51	0/3366
1	D	0.36	0/2517	0.53	0/3396
1	E	0.36	0/2488	0.50	0/3359
1	F	0.36	0/2486	0.51	0/3356
2	G	0.46	0/79	0.65	0/106
2	H	0.47	0/75	1.03	1/100 (1.0%)
All	All	0.36	0/15076	0.52	1/20345 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1497	GLU	CB-CA-C	-5.61	99.18	110.40

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	2409	0	2402	36	0
1	B	2412	0	2405	34	0
1	C	2435	0	2419	24	0
1	D	2457	0	2458	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2429	0	2412	15	0
1	F	2427	0	2419	21	0
2	G	80	0	70	5	0
2	H	76	0	68	2	0
3	I	35	0	9	0	0
4	J	40	0	10	1	0
5	K	30	0	9	0	0
6	A	70	0	0	1	0
6	B	85	0	0	0	0
6	C	102	0	0	1	0
6	D	94	0	0	0	0
6	E	86	0	0	2	0
6	F	105	0	0	0	0
6	H	1	0	0	0	0
All	All	15373	0	14681	147	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 5.

All (147) 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:25:ARG:HB3	1:B:93:THR:CG2	1.69	1.23
1:A:68:ILE:HD13	1:A:291:ARG:NH1	1.59	1.17
1:B:25:ARG:CB	1:B:93:THR:HG23	1.78	1.12
1:B:25:ARG:HD3	1:B:93:THR:HG21	1.34	1.10
1:A:68:ILE:CD1	1:A:291:ARG:NH1	2.28	0.95
1:A:68:ILE:HD13	1:A:291:ARG:HH12	1.31	0.93
1:A:304:HIS:CD2	2:G:1485:LEU:HD22	2.03	0.92
1:A:68:ILE:CD1	1:A:291:ARG:HH11	1.82	0.90
1:B:25:ARG:HD3	1:B:93:THR:CG2	2.02	0.88
1:F:23:THR:HB	1:F:90:GLU:OE2	1.78	0.82
1:B:25:ARG:HB3	1:B:93:THR:HG23	0.86	0.78
1:A:304:HIS:CG	2:G:1485:LEU:HD22	2.19	0.78
1:A:68:ILE:HD12	1:A:291:ARG:HH11	1.49	0.76
1:C:80:LYS:HG3	1:C:83:ASP:HB2	1.72	0.71
1:E:79:LYS:HE2	1:E:83:ASP:HB3	1.72	0.71
1:F:68:ILE:HD11	1:F:291:ARG:HE	1.57	0.69
1:D:227:LYS:HB2	1:D:284:LYS:HE3	1.74	0.69
1:A:87:SER:O	1:A:92:VAL:HG22	1.94	0.67
1:B:78:ASP:HB2	2:G:1487:ASP:O	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:22:PHE:CE2	1:D:92:VAL:HG21	2.30	0.66
1:A:265:SER:HB2	1:A:272:THR:HG23	1.78	0.65
1:B:25:ARG:CB	1:B:93:THR:CG2	2.54	0.64
1:D:20:LYS:HG2	1:D:39:MET:HE3	1.81	0.63
1:A:261:ILE:HD12	1:A:273:ILE:HG23	1.81	0.63
1:F:25:ARG:N	1:F:92:VAL:O	2.27	0.62
1:B:163:LEU:HG	1:B:168:LEU:HD11	1.82	0.62
1:B:69:LYS:O	1:B:70:ASP:HB2	2.00	0.61
1:F:85:ASP:OD1	1:F:85:ASP:N	2.34	0.60
1:B:25:ARG:CD	1:B:93:THR:HG21	2.23	0.60
1:B:25:ARG:CD	1:B:93:THR:CG2	2.77	0.60
1:F:42:LYS:HA	1:F:80:LYS:HA	1.84	0.59
1:A:227:LYS:HD2	1:A:284:LYS:HB2	1.82	0.59
1:D:22:PHE:HE2	1:D:92:VAL:HG21	1.68	0.58
1:B:261:ILE:O	2:H:1505:ASN:ND2	2.38	0.57
1:C:265:SER:OG	1:C:272:THR:OG1	2.22	0.57
1:E:172:ARG:NH1	6:E:473:HOH:O	2.37	0.56
1:B:79:LYS:HG2	1:B:80:LYS:O	2.06	0.56
1:E:68:ILE:HD12	1:E:291:ARG:NH1	2.20	0.56
1:A:85:ASP:OD1	1:A:85:ASP:N	2.39	0.56
1:C:67:THR:HG22	1:C:72:VAL:HG12	1.89	0.55
1:F:88:LYS:O	1:F:89:GLU:C	2.42	0.55
1:F:22:PHE:HD2	1:F:39:MET:HE2	1.72	0.55
1:F:76:LYS:NZ	1:F:78:ASP:HB2	2.21	0.55
1:A:154:ASP:H	1:A:158:HIS:CE1	2.25	0.54
1:B:93:THR:O	1:B:94:PHE:CD1	2.60	0.54
1:B:25:ARG:CG	1:B:93:THR:CG2	2.85	0.54
1:A:171:LYS:NZ	6:A:466:HOH:O	2.40	0.54
1:A:56:LEU:HD13	1:A:114:THR:HG22	1.88	0.54
1:A:69:LYS:O	1:A:70:ASP:HB2	2.09	0.53
1:A:304:HIS:CG	2:G:1485:LEU:CD2	2.92	0.53
1:A:111:GLN:HB2	1:A:114:THR:HG23	1.92	0.51
1:C:90:GLU:HG3	1:C:91:PRO:HA	1.91	0.51
1:C:99:LYS:NZ	6:C:439:HOH:O	2.36	0.51
1:B:24:VAL:HG12	1:B:92:VAL:HB	1.93	0.50
1:A:226:ASN:HB2	1:A:227:LYS:HZ1	1.75	0.50
1:A:233:LEU:HD22	1:A:246:PRO:HA	1.93	0.50
1:B:76:LYS:HD2	2:G:1488:THR:HA	1.92	0.50
1:F:234:LEU:HD21	1:F:241:LEU:HD22	1.93	0.50
1:F:100:PHE:CZ	1:F:302:GLY:HA3	2.46	0.50
1:A:58:GLU:OE1	1:A:114:THR:HG21	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:LYS:HG2	1:B:88:LYS:O	2.12	0.49
1:B:100:PHE:CZ	1:B:302:GLY:HA3	2.48	0.48
1:E:233:LEU:HD22	1:E:246:PRO:HA	1.95	0.48
1:D:42:LYS:HE3	1:D:44:LYS:HB2	1.94	0.48
1:A:78:ASP:OD1	1:A:78:ASP:N	2.38	0.48
1:B:233:LEU:HD22	1:B:246:PRO:HA	1.96	0.48
1:C:155:PRO:HG2	1:F:196:ARG:HG2	1.95	0.48
1:D:179:MET:HB3	1:D:183:MET:HB2	1.96	0.47
1:B:25:ARG:NH2	1:B:32:GLU:OE2	2.44	0.47
1:A:226:ASN:HB2	1:A:227:LYS:NZ	2.30	0.47
1:D:25:ARG:NH2	1:D:32:GLU:OE2	2.34	0.47
1:B:35:PHE:CE1	1:B:53:THR:HG21	2.49	0.47
1:B:74:TRP:CD1	1:B:99:LYS:HE3	2.49	0.47
1:F:68:ILE:HD11	1:F:291:ARG:NE	2.27	0.47
1:E:26:ILE:HB	1:E:33:MET:HB2	1.97	0.47
1:A:167:GLU:HG3	1:A:167:GLU:O	2.16	0.46
1:E:74:TRP:CD1	1:E:99:LYS:HE3	2.50	0.46
1:B:64:LEU:HA	1:B:64:LEU:HD12	1.84	0.46
1:A:294:LYS:HD3	1:A:294:LYS:HA	1.62	0.46
1:C:272:THR:HG22	1:C:284:LYS:HG2	1.99	0.45
1:D:269:LYS:NZ	1:D:290:LEU:HD11	2.32	0.45
1:C:274:LYS:HZ1	1:C:278:LYS:HD2	1.81	0.45
1:A:272:THR:HB	1:A:284:LYS:HG2	1.98	0.45
1:B:64:LEU:HD23	1:B:75:LEU:HD22	1.98	0.45
1:B:71:THR:HG22	1:B:72:VAL:O	2.17	0.45
1:C:79:LYS:HZ3	1:C:83:ASP:HB3	1.82	0.45
1:F:46:LEU:O	1:F:50:VAL:HG23	2.16	0.45
1:A:119:PHE:HA	1:A:208:LEU:HD21	1.99	0.45
1:D:88:LYS:HG3	1:D:88:LYS:O	2.16	0.45
1:D:68:ILE:HD12	1:D:85:ASP:OD2	2.16	0.45
1:C:76:LYS:HB2	1:C:79:LYS:HB3	1.99	0.45
1:C:159:LYS:HB3	1:C:159:LYS:HE2	1.59	0.45
1:F:87:SER:OG	1:F:89:GLU:HG2	2.17	0.45
1:B:297:LEU:HD23	2:H:1502:LEU:HD11	1.98	0.45
1:E:224:ILE:HG12	1:E:285:PHE:CD1	2.51	0.44
1:D:263:ASN:OD1	1:D:264:ILE:N	2.38	0.44
1:D:275:PRO:HD2	1:D:282:VAL:HG12	1.99	0.44
1:B:84:HIS:HB2	1:B:86:VAL:HG12	1.98	0.44
1:D:100:PHE:CZ	1:D:302:GLY:HA3	2.53	0.44
1:C:134:PRO:HB2	1:E:194:GLU:CD	2.38	0.44
1:D:212:GLN:HA	1:D:217:TYR:CG	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HG	1:B:168:LEU:CD1	2.47	0.44
1:F:227:LYS:HA	1:F:227:LYS:HD3	1.70	0.44
1:C:268:ASP:HB3	1:C:269:LYS:H	1.56	0.44
1:A:100:PHE:CZ	1:A:302:GLY:HA3	2.52	0.44
1:C:277:ASP:OD1	1:C:279:LYS:HG2	2.18	0.44
1:B:79:LYS:HG2	1:B:80:LYS:N	2.33	0.43
1:C:136:GLU:OE1	1:C:187:ARG:NH2	2.52	0.43
1:D:22:PHE:CE1	1:D:82:LEU:HD11	2.53	0.43
1:C:305:ASP:O	1:C:309:ARG:HG3	2.18	0.43
1:E:253:LYS:HE2	1:E:253:LYS:HB3	1.82	0.43
1:A:74:TRP:CD1	1:A:99:LYS:HE3	2.54	0.43
1:E:263:ASN:HB2	4:J:6:UNK:O	2.19	0.43
1:A:56:LEU:CD1	1:A:114:THR:HG22	2.49	0.43
1:C:90:GLU:HA	1:C:91:PRO:HA	1.80	0.43
1:A:290:LEU:HD12	1:A:290:LEU:H	1.84	0.43
1:A:275:PRO:HG2	1:A:280:ILE:HG12	2.00	0.42
1:B:76:LYS:HB2	1:B:79:LYS:HB3	2.01	0.42
1:B:160:ARG:HE	1:B:160:ARG:HB2	1.58	0.42
1:C:57:ARG:N	1:C:111:GLN:OE1	2.42	0.42
1:A:153:TYR:HE2	1:E:155:PRO:HG2	1.84	0.42
1:A:156:SER:O	1:A:159:LYS:HE2	2.20	0.42
1:C:100:PHE:CZ	1:C:302:GLY:HA3	2.54	0.42
1:F:167:GLU:O	1:F:167:GLU:HG3	2.19	0.41
1:D:265:SER:HB2	1:D:272:THR:OG1	2.19	0.41
1:B:150:TYR:OH	1:B:166:GLU:OE2	2.31	0.41
1:D:67:THR:O	1:D:67:THR:HG23	2.19	0.41
1:A:90:GLU:HA	1:A:91:PRO:C	2.39	0.41
1:F:265:SER:HB2	1:F:272:THR:HB	2.01	0.41
1:D:44:LYS:HA	1:D:77:MET:HE1	2.03	0.41
1:F:200:ARG:HD3	1:F:204:GLU:OE2	2.20	0.41
1:D:141:LEU:HB3	1:D:207:TYR:CE2	2.55	0.41
1:C:306:LEU:O	1:C:310:ARG:HG3	2.21	0.41
1:F:270:GLU:HB2	1:F:286:ASN:ND2	2.35	0.41
1:B:227:LYS:HD3	1:B:284:LYS:HB2	2.01	0.41
1:E:182:GLU:OE1	1:E:182:GLU:N	2.51	0.41
1:D:292:VAL:O	1:D:296:ILE:HG13	2.20	0.41
1:A:74:TRP:CG	1:A:99:LYS:HE3	2.56	0.41
1:C:242:HIS:HB3	1:C:252:PRO:HB3	2.02	0.41
1:C:277:ASP:HB2	1:D:279:LYS:HD3	2.02	0.40
1:C:187:ARG:NH1	6:E:443:HOH:O	2.54	0.40
1:E:75:LEU:HD12	1:E:75:LEU:HA	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:57:ARG:N	1:F:111:GLN:OE1	2.39	0.40
1:F:305:ASP:O	1:F:309:ARG:HG3	2.21	0.40
1:E:88:LYS:HG2	1:E:88:LYS:O	2.20	0.40
1:C:212:GLN:HA	1:C:217:TYR:CG	2.56	0.40
1:E:111:GLN:HB3	1:E:113:ILE:HG22	2.04	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	287/301 (95%)	283 (99%)	4 (1%)	0	100	100
1	B	288/301 (96%)	287 (100%)	1 (0%)	0	100	100
1	C	291/301 (97%)	289 (99%)	2 (1%)	0	100	100
1	D	293/301 (97%)	290 (99%)	3 (1%)	0	100	100
1	E	290/301 (96%)	288 (99%)	2 (1%)	0	100	100
1	F	289/301 (96%)	285 (99%)	4 (1%)	0	100	100
2	G	9/96 (9%)	9 (100%)	0	0	100	100
2	H	9/96 (9%)	9 (100%)	0	0	100	100
All	All	1756/1998 (88%)	1740 (99%)	16 (1%)	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	261/272 (96%)	247 (95%)	14 (5%)	27	36
1	B	262/272 (96%)	247 (94%)	15 (6%)	25	34
1	C	263/272 (97%)	253 (96%)	10 (4%)	40	54
1	D	267/272 (98%)	254 (95%)	13 (5%)	31	41
1	E	263/272 (97%)	246 (94%)	17 (6%)	21	27
1	F	264/272 (97%)	250 (95%)	14 (5%)	28	37
2	G	8/85 (9%)	7 (88%)	1 (12%)	6	5
2	H	7/85 (8%)	7 (100%)	0	100	100
All	All	1595/1802 (88%)	1511 (95%)	84 (5%)	28	37

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	THR
1	A	68	ILE
1	A	69	LYS
1	A	70	ASP
1	A	82	LEU
1	A	85	ASP
1	A	90	GLU
1	A	196	ARG
1	A	200	ARG
1	A	227	LYS
1	A	233	LEU
1	A	234	LEU
1	A	272	THR
1	A	280	ILE
1	B	23	THR
1	B	32	GLU
1	B	41	TRP
1	B	64	LEU
1	B	67	THR
1	B	69	LYS
1	B	93	THR
1	B	163	LEU
1	B	167	GLU
1	B	196	ARG

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Mol	Chain	Res	Type
1	B	200	ARG
1	B	233	LEU
1	B	234	LEU
1	B	250	LEU
1	B	268	ASP
1	C	68	ILE
1	C	159	LYS
1	C	196	ARG
1	C	200	ARG
1	C	230	THR
1	C	233	LEU
1	C	234	LEU
1	C	267	SER
1	C	268	ASP
1	C	282	VAL
1	D	21	THR
1	D	44	LYS
1	D	69	LYS
1	D	70	ASP
1	D	72	VAL
1	D	78	ASP
1	D	90	GLU
1	D	107	GLU
1	D	196	ARG
1	D	234	LEU
1	D	241	LEU
1	D	279	LYS
1	D	308	MET
1	E	23	THR
1	E	24	VAL
1	E	38	GLU
1	E	68	ILE
1	E	69	LYS
1	E	78	ASP
1	E	85	ASP
1	E	90	GLU
1	E	107	GLU
1	E	176	LEU
1	E	200	ARG
1	E	224	ILE
1	E	225	ARG
1	E	233	LEU

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Mol	Chain	Res	Type
1	E	234	LEU
1	E	250	LEU
1	E	254	ILE
1	F	41	TRP
1	F	42	LYS
1	F	68	ILE
1	F	69	LYS
1	F	85	ASP
1	F	176	LEU
1	F	196	ARG
1	F	200	ARG
1	F	233	LEU
1	F	234	LEU
1	F	250	LEU
1	F	268	ASP
1	F	288	SER
1	F	290	LEU
2	G	1482	GLU

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

Mol	Chain	Res	Type
1	C	298	GLN
1	F	65	GLN
1	F	263	ASN
1	F	286	ASN
1	F	298	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	289/301 (96%)	0.80	46 (15%) 3 4	33, 59, 104, 130	0
1	B	290/301 (96%)	0.81	42 (14%) 3 6	29, 57, 109, 125	0
1	C	293/301 (97%)	0.59	27 (9%) 11 16	29, 54, 91, 108	0
1	D	295/301 (98%)	0.74	44 (14%) 3 5	30, 62, 101, 129	0
1	E	292/301 (97%)	0.63	37 (12%) 5 8	29, 57, 92, 110	0
1	F	291/301 (96%)	0.75	46 (15%) 3 5	27, 54, 104, 115	0
2	G	11/96 (11%)	1.08	3 (27%) 1 1	56, 68, 92, 94	0
2	H	11/96 (11%)	1.34	3 (27%) 1 1	65, 68, 100, 102	0
3	I	0/7	-	-	-	-
4	J	0/8	-	-	-	-
5	K	0/6	-	-	-	-
All	All	1772/2019 (87%)	0.73	248 (13%) 4 6	27, 57, 101, 130	0

All (248) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	LYS	10.9
1	C	313	ALA	7.3
1	D	281	ASP	6.4
1	E	279	LYS	6.4
1	D	280	ILE	6.2
1	D	277	ASP	5.8
1	D	284	LYS	5.6
1	F	279	LYS	5.6
1	A	278	LYS	5.4
1	E	278	LYS	5.3
1	B	40	LYS	5.3
1	A	83	ASP	5.3
1	B	38	GLU	5.2

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Mol	Chain	Res	Type	RSRZ
1	D	279	LYS	5.0
1	D	313	ALA	4.9
1	A	291	ARG	4.9
1	F	280	ILE	4.8
1	D	229	GLY	4.7
1	D	40	LYS	4.7
1	E	276	LEU	4.7
1	B	70	ASP	4.6
1	F	291	ARG	4.6
1	B	42	LYS	4.6
1	A	64	LEU	4.6
1	B	312	LYS	4.5
1	F	229	GLY	4.5
1	A	41	TRP	4.5
1	A	279	LYS	4.5
1	A	42	LYS	4.4
1	A	88	LYS	4.4
1	A	312	LYS	4.4
1	F	268	ASP	4.4
1	C	88	LYS	4.3
1	F	40	LYS	4.3
1	B	178	GLN	4.2
1	C	241	LEU	4.2
1	F	46	LEU	4.2
1	F	21	THR	4.2
1	A	46	LEU	4.2
1	D	234	LEU	4.2
1	D	19	PRO	4.2
1	C	39	MET	4.2
2	H	1498	ASP	4.2
1	D	241	LEU	4.1
1	D	227	LYS	4.1
1	F	290	LEU	4.1
1	B	290	LEU	4.1
1	B	46	LEU	4.0
1	D	288	SER	4.0
1	F	88	LYS	4.0
1	B	227	LYS	4.0
1	A	39	MET	4.0
1	F	90	GLU	3.9
1	F	64	LEU	3.9
1	B	90	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	34	GLU	3.8
1	B	39	MET	3.8
1	E	277	ASP	3.8
1	C	89	GLU	3.7
1	E	46	LEU	3.7
1	E	291	ARG	3.7
1	C	40	LYS	3.7
1	F	87	SER	3.7
1	F	228	LYS	3.7
1	D	21	THR	3.7
1	F	78	ASP	3.6
1	E	89	GLU	3.6
1	C	78	ASP	3.6
1	E	78	ASP	3.6
1	D	46	LEU	3.6
1	B	68	ILE	3.5
1	F	234	LEU	3.5
1	D	274	LYS	3.5
1	B	41	TRP	3.5
1	D	75	LEU	3.5
1	A	69	LYS	3.5
1	C	69	LYS	3.5
1	D	291	ARG	3.5
2	G	1487	ASP	3.4
1	C	234	LEU	3.4
1	A	90	GLU	3.4
1	D	278	LYS	3.4
1	F	75	LEU	3.4
1	E	288	SER	3.4
1	A	277	ASP	3.4
1	E	227	LYS	3.4
1	D	64	LEU	3.4
1	A	68	ILE	3.4
1	E	269	LYS	3.3
1	F	41	TRP	3.3
1	F	296	ILE	3.3
1	E	40	LYS	3.2
1	B	26	ILE	3.2
1	D	236	VAL	3.2
1	B	228	LYS	3.2
1	A	234	LEU	3.2
1	E	274	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	78	ASP	3.2
1	A	80	LYS	3.2
1	E	90	GLU	3.1
1	E	64	LEU	3.1
1	B	80	LYS	3.1
1	C	64	LEU	3.1
1	E	236	VAL	3.1
1	E	228	LYS	3.1
1	A	89	GLU	3.1
1	B	69	LYS	3.1
1	F	69	LYS	3.0
1	F	80	LYS	3.0
1	D	285	PHE	3.0
1	D	178	GLN	3.0
1	D	20	LYS	2.9
1	D	312	LYS	2.9
1	F	91	PRO	2.9
1	F	22	PHE	2.9
1	A	38	GLU	2.9
1	D	289	LYS	2.9
1	A	78	ASP	2.9
1	B	94	PHE	2.9
1	B	234	LEU	2.9
1	E	39	MET	2.9
1	E	69	LYS	2.9
1	C	277	ASP	2.9
1	B	279	LYS	2.9
1	E	88	LYS	2.9
1	F	227	LYS	2.9
1	A	309	ARG	2.8
1	B	236	VAL	2.8
1	E	24	VAL	2.8
1	B	44	LYS	2.8
1	A	70	ASP	2.8
1	A	276	LEU	2.8
1	D	225	ARG	2.8
1	C	46	LEU	2.8
1	F	70	ASP	2.8
1	B	57	ARG	2.8
1	D	68	ILE	2.7
1	D	94	PHE	2.7
1	A	227	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	34	GLU	2.7
1	B	49	LEU	2.7
1	B	88	LYS	2.7
1	F	94	PHE	2.7
1	F	276	LEU	2.7
1	D	273	ILE	2.7
1	F	26	ILE	2.7
1	F	269	LYS	2.7
1	A	165	GLN	2.7
1	B	273	ILE	2.6
1	B	89	GLU	2.6
1	B	241	LEU	2.6
1	C	21	THR	2.6
1	B	36	ASN	2.6
1	D	268	ASP	2.6
1	E	70	ASP	2.6
1	A	290	LEU	2.6
1	A	273	ILE	2.6
1	F	285	PHE	2.6
1	F	241	LEU	2.6
1	F	207	TYR	2.6
1	A	311	ARG	2.5
1	E	234	LEU	2.5
1	D	253	LYS	2.5
1	B	177	TYR	2.5
1	F	281	ASP	2.5
1	B	91	PRO	2.5
1	F	83	ASP	2.5
1	A	26	ILE	2.5
1	A	141	LEU	2.5
1	B	87	SER	2.5
1	E	26	ILE	2.5
1	A	178	GLN	2.5
1	C	291	ARG	2.5
1	E	265	SER	2.5
1	C	273	ILE	2.5
1	C	178	GLN	2.4
1	E	83	ASP	2.4
1	E	75	LEU	2.4
1	F	235	GLY	2.4
1	B	278	LYS	2.4
1	C	278	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	279	LYS	2.4
1	C	312	LYS	2.4
1	D	290	LEU	2.4
1	E	241	LEU	2.4
1	D	286	ASN	2.4
1	E	235	GLY	2.4
1	E	289	LYS	2.4
1	F	224	ILE	2.4
1	A	285	PHE	2.3
1	B	276	LEU	2.3
1	E	81	VAL	2.3
2	H	1497	GLU	2.3
1	A	142	ALA	2.3
1	A	241	LEU	2.3
1	D	230	THR	2.3
1	F	57	ARG	2.3
1	A	96	PHE	2.3
1	B	118	PHE	2.3
1	E	285	PHE	2.3
1	C	75	LEU	2.3
1	A	225	ARG	2.3
1	A	94	PHE	2.2
1	D	269	LYS	2.2
1	E	230	THR	2.2
1	F	89	GLU	2.2
2	H	1500	ILE	2.2
1	D	228	LYS	2.2
1	A	91	PRO	2.2
1	A	160	ARG	2.2
1	C	280	ILE	2.2
1	E	160	ARG	2.2
1	D	81	VAL	2.2
1	F	278	LYS	2.2
1	A	114	THR	2.2
1	C	296	ILE	2.2
1	C	70	ASP	2.2
1	D	69	LYS	2.1
1	A	72	VAL	2.1
2	G	1486	GLY	2.1
1	B	207	TYR	2.1
1	B	142	ALA	2.1
1	D	85	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	280	ILE	2.1
1	B	285	PHE	2.1
1	D	266	TYR	2.1
1	A	24	VAL	2.1
1	F	277	ASP	2.1
1	F	122	VAL	2.1
1	C	96	PHE	2.1
1	F	79	LYS	2.1
1	B	286	ASN	2.1
2	G	1479	GLU	2.1
1	C	227	LYS	2.1
1	B	35	PHE	2.1
1	E	94	PHE	2.1
1	F	38	GLU	2.1
1	F	208	LEU	2.1
1	A	145	ALA	2.1
1	D	208	LEU	2.0
1	A	45	ASP	2.0
1	D	80	LYS	2.0
1	C	81	VAL	2.0
1	B	85	ASP	2.0
1	E	165	GLN	2.0
1	F	240	GLY	2.0
1	C	171	LYS	2.0
1	A	81	VAL	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 carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.