



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 13, 2016 – 08:15 AM EDT

PDB ID : 5B2G  
Title : Crystal structure of human claudin-4 in complex with C-terminal fragment of Clostridium perfringens enterotoxin  
Authors : Shinoda, T.; Kimura-Someya, T.; Shirouzu, M.; Yokoyama, S.  
Deposited on : 2016-01-15  
Resolution : 3.50 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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
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)	:	rb-20027939

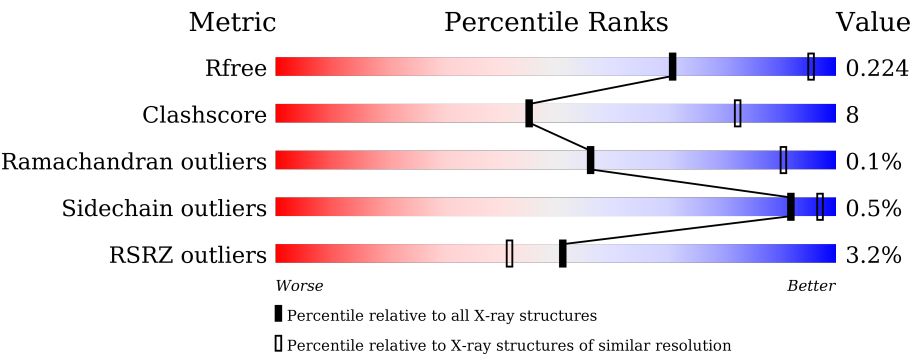
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	364	<div><div>2%</div><div><div></div><div>73%</div><div>18%</div><div>9%</div></div></div>
1	C	364	<div><div>%</div><div><div></div><div>80%</div><div>15%</div><div>• •</div></div></div>
1	E	364	<div><div>6%</div><div><div></div><div>71%</div><div>21%</div><div>8%</div></div></div>
1	G	364	<div><div>2%</div><div><div></div><div>78%</div><div>13%</div><div>• 8%</div></div></div>
2	B	142	<div><div>4%</div><div><div></div><div>70%</div><div>12%</div><div>18%</div></div></div>
2	D	142	<div><div>%</div><div><div></div><div>78%</div><div>9%</div><div>13%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	F	142	<div><div><div>4%</div><div></div><div>65%</div><div>18%</div><div>••</div><div>15%</div></div></div>
2	H	142	<div><div><div>2%</div><div></div><div>78%</div><div>9%</div><div>13%</div></div></div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14200 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 Endolysin, Claudin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	0	0
			2540	1628	438	456	18			
1	C	350	Total	C	N	O	S	0	0	0
			2657	1695	459	481	22			
1	E	335	Total	C	N	O	S	0	0	0
			2563	1642	441	460	20			
1	G	336	Total	C	N	O	S	0	0	0
			2558	1637	443	459	19			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	995	GLY	-	expression tag	UNP P00720
A	996	SER	-	expression tag	UNP P00720
A	997	SER	-	expression tag	UNP P00720
A	998	GLY	-	expression tag	UNP P00720
A	999	SER	-	expression tag	UNP P00720
A	1000	SER	-	expression tag	UNP P00720
A	1001	GLY	-	expression tag	UNP P00720
A	1012	GLY	ARG	engineered mutation	UNP P00720
A	1054	THR	CYS	engineered mutation	UNP P00720
A	1097	ALA	CYS	engineered mutation	UNP P00720
A	1137	ARG	ILE	engineered mutation	UNP P00720
A	1163	GLY	-	linker	UNP P00720
A	1347	SER	-	expression tag	UNP O14493
A	1348	GLY	-	expression tag	UNP O14493
A	1349	PRO	-	expression tag	UNP O14493
A	1350	SER	-	expression tag	UNP O14493
A	1351	SER	-	expression tag	UNP O14493
A	1352	GLY	-	expression tag	UNP O14493
A	1353	GLU	-	expression tag	UNP O14493
A	1354	ASN	-	expression tag	UNP O14493
A	1355	LEU	-	expression tag	UNP O14493

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1356	TYR	-	expression tag	UNP O14493
A	1357	PHE	-	expression tag	UNP O14493
A	1358	GLN	-	expression tag	UNP O14493
C	995	GLY	-	expression tag	UNP P00720
C	996	SER	-	expression tag	UNP P00720
C	997	SER	-	expression tag	UNP P00720
C	998	GLY	-	expression tag	UNP P00720
C	999	SER	-	expression tag	UNP P00720
C	1000	SER	-	expression tag	UNP P00720
C	1001	GLY	-	expression tag	UNP P00720
C	1012	GLY	ARG	engineered mutation	UNP P00720
C	1054	THR	CYS	engineered mutation	UNP P00720
C	1097	ALA	CYS	engineered mutation	UNP P00720
C	1137	ARG	ILE	engineered mutation	UNP P00720
C	1163	GLY	-	linker	UNP P00720
C	184	SER	-	expression tag	UNP O14493
C	185	GLY	-	expression tag	UNP O14493
C	186	PRO	-	expression tag	UNP O14493
C	187	SER	-	expression tag	UNP O14493
C	188	SER	-	expression tag	UNP O14493
C	189	GLY	-	expression tag	UNP O14493
C	190	GLU	-	expression tag	UNP O14493
C	191	ASN	-	expression tag	UNP O14493
C	192	LEU	-	expression tag	UNP O14493
C	193	TYR	-	expression tag	UNP O14493
C	194	PHE	-	expression tag	UNP O14493
C	195	GLN	-	expression tag	UNP O14493
E	995	GLY	-	expression tag	UNP P00720
E	996	SER	-	expression tag	UNP P00720
E	997	SER	-	expression tag	UNP P00720
E	998	GLY	-	expression tag	UNP P00720
E	999	SER	-	expression tag	UNP P00720
E	1000	SER	-	expression tag	UNP P00720
E	1001	GLY	-	expression tag	UNP P00720
E	1012	GLY	ARG	engineered mutation	UNP P00720
E	1054	THR	CYS	engineered mutation	UNP P00720
E	1097	ALA	CYS	engineered mutation	UNP P00720
E	1137	ARG	ILE	engineered mutation	UNP P00720
E	1163	GLY	-	linker	UNP P00720
E	1347	SER	-	expression tag	UNP O14493
E	1348	GLY	-	expression tag	UNP O14493
E	1349	PRO	-	expression tag	UNP O14493

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1350	SER	-	expression tag	UNP O14493
E	1351	SER	-	expression tag	UNP O14493
E	1352	GLY	-	expression tag	UNP O14493
E	1353	GLU	-	expression tag	UNP O14493
E	1354	ASN	-	expression tag	UNP O14493
E	1355	LEU	-	expression tag	UNP O14493
E	1356	TYR	-	expression tag	UNP O14493
E	1357	PHE	-	expression tag	UNP O14493
E	1358	GLN	-	expression tag	UNP O14493
G	995	GLY	-	expression tag	UNP P00720
G	996	SER	-	expression tag	UNP P00720
G	997	SER	-	expression tag	UNP P00720
G	998	GLY	-	expression tag	UNP P00720
G	999	SER	-	expression tag	UNP P00720
G	1000	SER	-	expression tag	UNP P00720
G	1001	GLY	-	expression tag	UNP P00720
G	1012	GLY	ARG	engineered mutation	UNP P00720
G	1054	THR	CYS	engineered mutation	UNP P00720
G	1097	ALA	CYS	engineered mutation	UNP P00720
G	1137	ARG	ILE	engineered mutation	UNP P00720
G	1163	GLY	-	linker	UNP P00720
G	1347	SER	-	expression tag	UNP O14493
G	1348	GLY	-	expression tag	UNP O14493
G	1349	PRO	-	expression tag	UNP O14493
G	1350	SER	-	expression tag	UNP O14493
G	1351	SER	-	expression tag	UNP O14493
G	1352	GLY	-	expression tag	UNP O14493
G	1353	GLU	-	expression tag	UNP O14493
G	1354	ASN	-	expression tag	UNP O14493
G	1355	LEU	-	expression tag	UNP O14493
G	1356	TYR	-	expression tag	UNP O14493
G	1357	PHE	-	expression tag	UNP O14493
G	1358	GLN	-	expression tag	UNP O14493

- Molecule 2 is a protein called Heat-labile enterotoxin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	117	Total	C	N	O	Se	0	0	0
			933	595	157	180	1			
2	D	124	Total	C	N	O	Se	0	0	0
			992	633	165	193	1			
2	F	121	Total	C	N	O	Se	0	0	0
			965	617	161	186	1			

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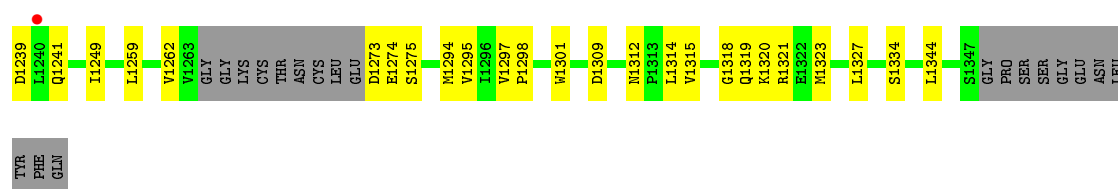
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	124	Total	C	N	O	Se	0	0	0
			992	633	165	193	1			

There are 36 discrepancies between the modelled and reference sequences:

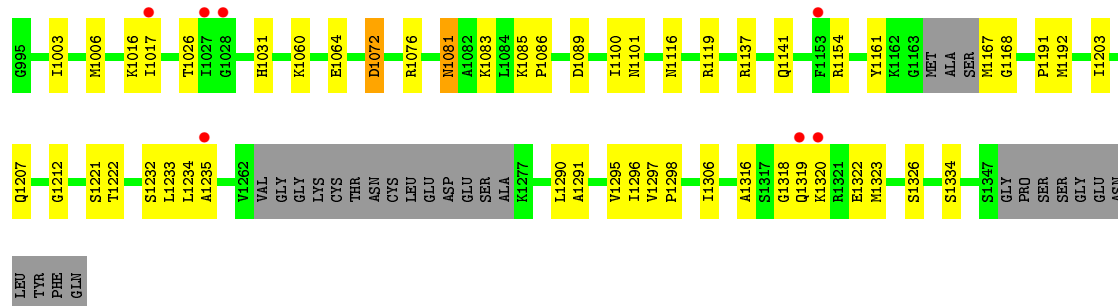
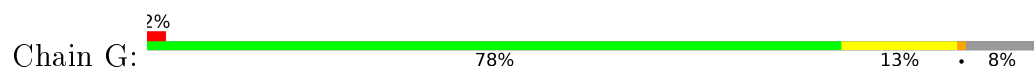
Chain	Residue	Modelled	Actual	Comment	Reference
B	178	GLY	-	expression tag	UNP P01558
B	179	SER	-	expression tag	UNP P01558
B	180	SER	-	expression tag	UNP P01558
B	181	GLY	-	expression tag	UNP P01558
B	182	SER	-	expression tag	UNP P01558
B	183	SER	-	expression tag	UNP P01558
B	184	GLY	-	expression tag	UNP P01558
B	185	ARG	-	expression tag	UNP P01558
B	186	TRP	-	expression tag	UNP P01558
D	178	GLY	-	expression tag	UNP P01558
D	179	SER	-	expression tag	UNP P01558
D	180	SER	-	expression tag	UNP P01558
D	181	GLY	-	expression tag	UNP P01558
D	182	SER	-	expression tag	UNP P01558
D	183	SER	-	expression tag	UNP P01558
D	184	GLY	-	expression tag	UNP P01558
D	185	ARG	-	expression tag	UNP P01558
D	186	TRP	-	expression tag	UNP P01558
F	178	GLY	-	expression tag	UNP P01558
F	179	SER	-	expression tag	UNP P01558
F	180	SER	-	expression tag	UNP P01558
F	181	GLY	-	expression tag	UNP P01558
F	182	SER	-	expression tag	UNP P01558
F	183	SER	-	expression tag	UNP P01558
F	184	GLY	-	expression tag	UNP P01558
F	185	ARG	-	expression tag	UNP P01558
F	186	TRP	-	expression tag	UNP P01558
H	178	GLY	-	expression tag	UNP P01558
H	179	SER	-	expression tag	UNP P01558
H	180	SER	-	expression tag	UNP P01558
H	181	GLY	-	expression tag	UNP P01558
H	182	SER	-	expression tag	UNP P01558
H	183	SER	-	expression tag	UNP P01558
H	184	GLY	-	expression tag	UNP P01558
H	185	ARG	-	expression tag	UNP P01558
H	186	TRP	-	expression tag	UNP P01558



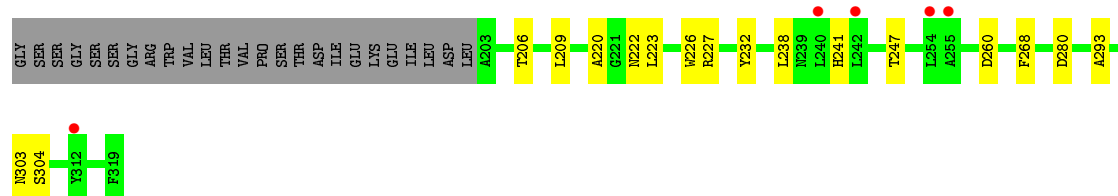




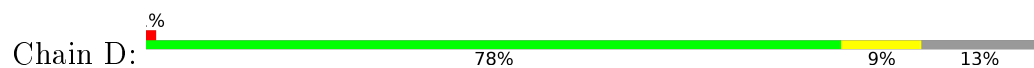
- Molecule 1: Endolysin, Claudin-4



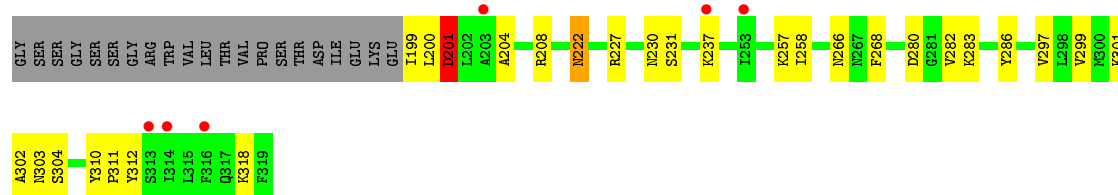
- Molecule 2: Heat-labile enterotoxin B chain



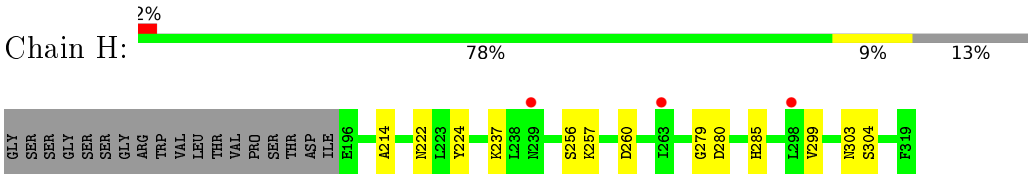
- Molecule 2: Heat-labile enterotoxin B chain



- Molecule 2: Heat-labile enterotoxin B chain



- Molecule 2: Heat-labile enterotoxin B chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.92Å 105.92Å 244.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.59 – 3.50 48.59 – 3.35	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.59-3.50) 99.8 (48.59-3.35)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.41 (at 3.33Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, $R_{free}$	0.288 , 0.309 0.216 , 0.224	Depositor DCC
$R_{free}$ test set	1688 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	129.9	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 61.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.36$ , $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.387 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14200	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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, 0.333 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	0.23	0/2581	0.44	0/3498
1	C	0.32	0/2700	0.53	3/3657 (0.1%)
1	E	0.24	0/2604	0.49	0/3528
1	G	0.22	0/2599	0.46	2/3518 (0.1%)
2	B	0.21	0/953	0.41	0/1294
2	D	0.26	0/1012	0.40	0/1373
2	F	0.29	0/985	0.45	1/1338 (0.1%)
2	H	0.22	0/1012	0.40	0/1373
All	All	0.25	0/14446	0.46	6/19579 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	156	GLN	CB-CA-C	-9.54	91.33	110.40
1	C	104	CYS	N-CA-C	-7.56	90.59	111.00
1	G	1081	ASN	CB-CA-C	5.87	122.14	110.40
2	F	201	ASP	CB-CA-C	5.78	121.96	110.40
1	G	1081	ASN	N-CA-C	-5.78	95.40	111.00
1	C	100	VAL	N-CA-C	-5.64	95.78	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	2540	0	2622	54	0
1	C	2657	0	2740	45	0
1	E	2563	0	2647	61	0
1	G	2558	0	2644	36	0
2	B	933	0	907	13	0
2	D	992	0	969	11	0
2	F	965	0	944	23	0
2	H	992	0	969	8	0
All	All	14200	0	14442	223	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 8.

All (223) 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:1213:LEU:CD1	1:A:1240:LEU:HD21	1.83	1.09
1:A:1213:LEU:HD12	1:A:1240:LEU:CD2	1.82	1.08
1:A:1213:LEU:HD12	1:A:1240:LEU:HD21	1.09	1.06
1:C:103:LYS:HG3	1:C:104:CYS:O	1.58	1.00
1:A:1213:LEU:CD1	1:A:1240:LEU:CD2	2.42	0.96
1:C:100:VAL:HG23	1:C:100:VAL:O	1.64	0.93
1:C:105:THR:O	1:E:1274:GLU:O	1.87	0.92
1:E:1318:GLY:HA2	1:E:1319:GLN:HB2	1.56	0.86
1:E:1078:ILE:HD11	1:E:1103:VAL:HG11	1.55	0.86
1:E:1052:ARG:NH2	1:E:1057:VAL:O	2.07	0.86
1:E:1007:LEU:HD13	1:E:1067:PHE:CZ	2.14	0.82
1:G:1232:SER:HB2	1:G:1233:LEU:HA	1.62	0.82
2:D:258:ILE:HD11	2:D:304:SER:HB2	1.63	0.79
1:C:69:SER:HB2	1:C:70:LEU:HA	1.66	0.76
1:C:1083:LYS:HD2	1:C:1112:ALA:HB1	1.68	0.75
1:C:100:VAL:O	1:C:100:VAL:CG2	2.35	0.73
2:B:247:THR:HG22	2:B:293:ALA:HB2	1.71	0.71
1:E:1232:SER:HB2	1:E:1233:LEU:HA	1.72	0.70
1:C:1085:LYS:HB3	1:C:1086:PRO:HD3	1.74	0.69
2:B:268:PHE:CZ	1:C:1151:THR:HG23	2.28	0.69
1:E:1007:LEU:HD13	1:E:1067:PHE:HZ	1.54	0.68
1:C:103:LYS:O	1:C:104:CYS:HB2	1.93	0.67
1:G:1232:SER:CB	1:G:1233:LEU:HA	2.23	0.67
1:A:1318:GLY:HA2	1:A:1319:GLN:HB2	1.76	0.67
1:A:1233:LEU:O	1:A:1233:LEU:HD12	1.95	0.66
1:A:1207:GLN:NE2	2:B:222:ASN:OD1	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ILE:HG22	1:C:43:SER:HB2	1.77	0.66
1:C:44:GLN:NE2	2:D:222:ASN:OD1	2.29	0.65
1:C:96:LEU:O	1:C:100:VAL:HG22	1.96	0.65
1:A:1232:SER:HB2	1:A:1233:LEU:HA	1.77	0.65
1:A:1213:LEU:HD11	1:A:1240:LEU:CD2	2.25	0.65
1:E:1219:VAL:HG22	1:E:1225:MET:HB3	1.78	0.64
1:C:69:SER:CB	1:C:70:LEU:HA	2.25	0.64
1:G:1234:LEU:HD12	1:G:1235:ALA:N	2.12	0.64
1:G:1085:LYS:NZ	1:G:1089:ASP:OD2	2.31	0.64
1:E:1006:MET:O	1:E:1009:ILE:HG22	1.97	0.63
1:E:1120:MET:O	1:E:1125:ARG:HG3	1.98	0.63
1:A:1137:ARG:NH2	2:F:201:ASP:O	2.33	0.61
1:E:1314:LEU:HD23	2:F:312:TYR:HA	1.83	0.61
1:A:1277:LYS:NZ	1:E:1238:GLN:O	2.34	0.60
2:B:206:THR:O	1:E:1140:ASN:ND2	2.34	0.60
1:C:105:THR:C	1:E:1274:GLU:O	2.40	0.60
1:E:1007:LEU:HD13	1:E:1067:PHE:CE2	2.36	0.60
1:E:1259:LEU:O	1:E:1262:VAL:HG12	2.02	0.59
1:E:1132:ASN:HA	1:E:1135:LYS:HE3	1.86	0.58
1:C:151:LEU:O	2:D:256:SER:OG	2.19	0.58
1:C:1085:LYS:NZ	1:C:1089:ASP:OD2	2.37	0.58
1:C:1060:LYS:O	1:C:1064:GLU:HG2	2.03	0.57
2:F:282:VAL:HG12	2:F:283:LYS:HG3	1.86	0.57
1:A:1002:ASN:N	1:A:1005:GLU:OE2	2.38	0.57
1:A:1087:VAL:HG21	1:A:1118:LEU:HB3	1.86	0.57
1:E:1078:ILE:CD1	1:E:1103:VAL:HG11	2.29	0.57
1:E:1232:SER:CB	1:E:1233:LEU:HA	2.32	0.57
2:F:200:LEU:HD12	2:F:204:ALA:HB2	1.87	0.57
1:C:1116:ASN:OD1	1:C:1117:SER:N	2.38	0.56
1:A:1084:LEU:HD21	1:A:1111:VAL:HB	1.87	0.56
1:E:1309:ASP:OD1	2:F:227:ARG:NH1	2.39	0.55
2:F:222:ASN:ND2	2:F:318:LYS:O	2.40	0.55
1:A:1273:ASP:OD1	2:F:208:ARG:NH2	2.39	0.55
1:E:1174:ILE:HG23	1:E:1344:LEU:HD21	1.89	0.55
1:C:156:GLN:OE1	2:D:315:LEU:HB2	2.06	0.55
1:E:1105:GLN:OE1	1:E:1145:ARG:NH1	2.39	0.55
2:B:209:LEU:HD22	2:B:226:TRP:CD1	2.42	0.54
1:A:1309:ASP:OD1	2:B:227:ARG:NH1	2.40	0.54
1:E:1155:THR:OG1	1:E:1157:THR:OG1	2.17	0.54
1:C:127:LEU:HD22	1:G:1290:LEU:HD22	1.89	0.54
2:F:237:LYS:HE2	2:F:299:VAL:HG12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1232:SER:CB	1:A:1233:LEU:HA	2.37	0.54
1:E:1196:THR:HG22	1:E:1321:ARG:HG2	1.89	0.54
1:G:1306:ILE:HG21	1:G:1323:MET:N	2.23	0.54
1:G:1318:GLY:HA2	1:G:1319:GLN:HB2	1.91	0.53
1:A:1339:LEU:O	1:A:1343:LEU:HD23	2.09	0.53
1:A:1277:LYS:NZ	1:E:1241:GLN:HB3	2.23	0.53
1:C:106:ASN:O	1:C:107:CYS:SG	2.66	0.52
1:C:133:ILE:HG13	1:C:171:SER:HB2	1.90	0.52
1:E:1318:GLY:HA2	1:E:1319:GLN:CB	2.31	0.52
1:E:1312:ASN:O	1:E:1315:VAL:HG22	2.09	0.52
2:F:199:ILE:N	2:F:201:ASP:OD1	2.43	0.52
1:G:1085:LYS:HB3	1:G:1086:PRO:HD3	1.91	0.52
2:H:237:LYS:HD3	2:H:299:VAL:CG1	2.40	0.51
1:A:1239:ASP:HB3	1:A:1301:TRP:CZ3	2.44	0.51
1:G:1296:ILE:HG13	1:G:1334:SER:HB2	1.93	0.51
1:G:1072:ASP:OD1	1:G:1076:ARG:NH2	2.44	0.51
1:A:1144:ASN:HA	1:A:1147:LYS:HE2	1.93	0.51
1:C:1069:GLN:HA	1:C:1072:ASP:OD2	2.11	0.51
2:F:301:LYS:NZ	2:F:302:ALA:O	2.34	0.51
1:E:1318:GLY:HA3	1:E:1320:LYS:N	2.26	0.51
1:A:1189:ALA:HA	1:C:19:LEU:HD22	1.93	0.50
1:A:1007:LEU:HD13	1:A:1067:PHE:CE1	2.46	0.50
1:E:1101:ASN:OD1	1:E:1145:ARG:NH2	2.42	0.50
2:D:225:ASP:OD2	2:D:227:ARG:NH1	2.41	0.50
1:A:1135:LYS:HE2	2:F:230:ASN:HB3	1.92	0.50
1:C:156:GLN:OE1	2:D:315:LEU:HD22	2.12	0.50
1:A:1296:ILE:HG13	1:A:1334:SER:HB2	1.93	0.50
2:F:237:LYS:HE2	2:F:299:VAL:CG1	2.42	0.50
1:C:146:ASP:OD1	1:C:152:VAL:HG21	2.11	0.49
2:F:303:ASN:OD1	2:F:304:SER:N	2.45	0.49
1:G:1234:LEU:HD12	1:G:1235:ALA:H	1.76	0.49
1:A:1026:THR:HG22	1:A:1027:ILE:N	2.27	0.49
1:A:1296:ILE:O	1:A:1300:SER:OG	2.26	0.48
1:G:1081:ASN:OD1	1:G:1083:LYS:HB2	2.12	0.48
2:H:303:ASN:OD1	2:H:304:SER:N	2.47	0.48
1:E:1295:VAL:HB	1:E:1334:SER:HB3	1.95	0.47
2:B:303:ASN:OD1	2:B:304:SER:N	2.48	0.47
1:A:1297:VAL:HB	1:A:1298:PRO:HD3	1.95	0.47
1:E:1249:ILE:HG22	1:E:1294:MET:HE1	1.96	0.47
1:E:1297:VAL:HB	1:E:1298:PRO:HD3	1.96	0.47
1:E:1318:GLY:HA3	1:E:1319:GLN:C	2.34	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1220:GLN:NE2	2:B:220:ALA:HB2	2.29	0.47
1:E:1078:ILE:HD11	1:E:1103:VAL:CG1	2.36	0.47
1:C:102:GLY:O	1:C:110:ASP:HB3	2.14	0.47
1:E:1020:ASP:OD1	1:E:1024:TYR:N	2.48	0.47
1:E:1144:ASN:HA	1:E:1147:LYS:HE2	1.95	0.47
1:E:1221:SER:HA	1:E:1222:THR:HA	1.52	0.47
1:G:1221:SER:HA	1:G:1222:THR:HA	1.61	0.47
1:G:1203:ILE:HD11	1:G:1319:GLN:OE1	2.15	0.47
1:C:135:PRO:O	1:C:139:THR:HG23	2.15	0.46
2:F:237:LYS:HG3	2:F:301:LYS:HB2	1.96	0.46
1:E:1318:GLY:CA	1:E:1319:GLN:C	2.84	0.46
1:G:1297:VAL:HB	1:G:1298:PRO:HD3	1.97	0.46
1:E:1009:ILE:HG21	1:E:1161:TYR:CE1	2.50	0.46
1:E:1174:ILE:CG2	1:E:1344:LEU:HD21	2.46	0.46
2:H:280:ASP:N	2:H:280:ASP:OD1	2.48	0.46
2:D:303:ASN:OD1	2:D:304:SER:N	2.49	0.46
1:E:1008:ARG:HD3	1:E:1013:LEU:HD22	1.97	0.46
1:E:1118:LEU:HD23	1:E:1121:LEU:HD21	1.98	0.46
1:G:1003:ILE:HD11	1:G:1100:ILE:CG2	2.46	0.46
1:G:1232:SER:CB	1:G:1233:LEU:CA	2.92	0.46
1:A:1318:GLY:CA	1:A:1319:GLN:HB2	2.44	0.46
1:E:1151:THR:HG21	1:E:1160:ALA:HB2	1.98	0.46
1:C:69:SER:CB	1:C:70:LEU:CA	2.94	0.45
1:E:1323:MET:HB2	1:E:1327:LEU:HD12	1.97	0.45
1:G:1003:ILE:HD11	1:G:1100:ILE:HG21	1.98	0.45
1:G:1320:LYS:NZ	1:G:1322:GLU:OE2	2.49	0.45
1:A:1232:SER:CB	1:A:1234:LEU:H	2.30	0.45
1:C:1163:GLY:O	1:C:1:MET:CB	2.64	0.45
2:F:266:ASN:HB2	2:F:297:VAL:HG23	1.98	0.45
1:A:1286:VAL:O	1:A:1289:LEU:HG	2.17	0.45
2:H:279:GLY:HA2	2:H:285:HIS:CD2	2.52	0.45
2:B:280:ASP:N	2:B:280:ASP:OD1	2.50	0.45
1:E:1232:SER:CB	1:E:1233:LEU:CA	2.95	0.45
1:C:103:LYS:O	1:C:104:CYS:CB	2.64	0.45
1:C:70:LEU:C	1:C:71:LEU:HD12	2.37	0.45
1:G:1006:MET:HG3	1:G:1161:TYR:CZ	2.52	0.45
1:E:1027:ILE:HD13	1:E:1046:LEU:HD11	1.99	0.45
1:A:1054:THR:HG23	1:A:1055:ASN:N	2.32	0.45
1:G:1016:LYS:HA	1:G:1016:LYS:HD2	1.80	0.44
1:C:156:GLN:NE2	2:D:225:ASP:OD1	2.47	0.44
2:F:280:ASP:N	2:F:280:ASP:OD1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1007:LEU:HD13	1:A:1067:PHE:HE1	1.83	0.44
1:E:1018:TYR:N	1:E:1026:THR:O	2.51	0.44
1:G:1207:GLN:NE2	2:H:222:ASN:OD1	2.50	0.44
1:E:1154:ARG:HG2	1:E:1155:THR:HG23	1.98	0.44
2:B:232:TYR:HE2	2:B:238:LEU:HD22	1.83	0.44
1:C:134:VAL:HB	1:C:135:PRO:HD3	1.99	0.44
1:G:1116:ASN:HA	1:G:1119:ARG:NH1	2.32	0.44
1:A:1209:ILE:HG12	1:A:1218:VAL:HG12	2.00	0.44
1:C:58:SER:HA	1:C:59:THR:HA	1.58	0.44
2:F:201:ASP:OD1	2:F:201:ASP:N	2.49	0.44
1:G:1060:LYS:O	1:G:1064:GLU:HG2	2.16	0.44
1:A:1221:SER:HA	1:A:1222:THR:HA	1.55	0.44
2:D:280:ASP:OD1	2:D:280:ASP:N	2.50	0.44
1:A:1085:LYS:HB3	1:A:1086:PRO:HD3	2.00	0.43
1:A:1233:LEU:HD11	1:C:103:LYS:HE2	1.99	0.43
2:D:204:ALA:HB2	2:D:237:LYS:HG3	1.99	0.43
1:E:1239:ASP:HB3	1:E:1301:TRP:CZ3	2.54	0.43
1:E:1066:LEU:HD23	1:E:1066:LEU:C	2.39	0.43
1:E:1014:ARG:HD3	1:E:1018:TYR:CD1	2.54	0.43
1:C:35:PHE:HB3	1:C:40:ILE:HG12	2.01	0.43
1:E:1222:THR:OG1	1:E:1223:GLY:N	2.51	0.43
1:G:1191:PRO:HA	1:G:1212:GLY:HA3	2.00	0.43
1:A:1232:SER:CB	1:A:1233:LEU:CA	2.97	0.43
1:A:1015:LEU:HD23	1:A:1058:ILE:O	2.19	0.43
1:A:1232:SER:HB2	1:A:1234:LEU:H	1.84	0.43
1:E:1274:GLU:HB2	1:E:1275:SER:HA	2.00	0.43
1:A:1196:THR:HG22	1:A:1321:ARG:HG2	2.01	0.42
1:A:1193:TRP:HA	1:A:1327:LEU:HD13	2.01	0.42
1:E:1026:THR:HG22	1:E:1027:ILE:N	2.34	0.42
1:G:1006:MET:SD	1:G:1101:ASN:ND2	2.92	0.42
1:G:1232:SER:HB3	1:G:1234:LEU:HG	2.01	0.42
1:G:1192:MET:O	1:G:1326:SER:OG	2.26	0.42
1:A:1024:TYR:CE1	1:A:1035:LYS:HA	2.55	0.42
1:A:1215:MET:CB	1:A:1229:VAL:HA	2.50	0.42
1:C:32:VAL:HG22	1:C:47:TRP:CE3	2.55	0.42
1:G:1137:ARG:NH1	1:G:1141:GLN:OE1	2.52	0.42
1:A:1220:GLN:NE2	2:B:220:ALA:CB	2.83	0.42
1:A:1289:LEU:HD12	1:A:1290:LEU:N	2.34	0.42
1:E:1009:ILE:HG21	1:E:1161:TYR:HE1	1.85	0.42
1:E:1262:VAL:HG13	1:E:1262:VAL:O	2.20	0.42
1:G:1167:MET:HA	1:G:1168:GLY:HA3	1.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1008:ARG:NH2	1:C:1064:GLU:OE1	2.47	0.42
1:A:1148:ARG:HB3	1:A:1161:TYR:CZ	2.55	0.41
1:C:1139:TYR:HD1	1:C:1146:ALA:CB	2.33	0.41
1:A:1135:LYS:NZ	2:F:231:SER:O	2.39	0.41
2:B:260:ASP:N	2:B:260:ASP:OD1	2.53	0.41
1:C:96:LEU:HA	1:C:99:VAL:HG22	2.02	0.41
2:D:207:GLU:OE2	2:D:232:TYR:OH	2.22	0.41
1:E:1094:VAL:HG23	1:E:1095:ARG:N	2.35	0.41
1:E:1273:ASP:OD1	1:E:1274:GLU:HA	2.20	0.41
1:A:1025:TYR:OH	1:A:1039:LEU:HG	2.20	0.41
1:E:1014:ARG:HD3	1:E:1018:TYR:HD1	1.85	0.41
1:G:1316:ALA:HB2	2:H:256:SER:HA	2.02	0.41
1:G:1026:THR:HG23	1:G:1031:HIS:C	2.41	0.41
1:A:1262:VAL:HG12	1:A:1262:VAL:O	2.19	0.41
2:F:310:TYR:HA	2:F:311:PRO:HD3	1.91	0.41
2:F:268:PHE:HB3	1:G:1154:ARG:HH12	1.85	0.41
1:G:1016:LYS:HG3	1:G:1017:ILE:H	1.85	0.41
1:C:1027:ILE:HG21	1:C:1046:LEU:CD1	2.51	0.41
1:C:1084:LEU:HD23	1:C:1099:LEU:HD21	2.03	0.41
1:C:18:TRP:O	1:C:21:VAL:HG12	2.20	0.41
1:E:1065:LYS:HA	1:E:1065:LYS:HD3	1.88	0.41
1:E:1148:ARG:O	1:E:1151:THR:HG22	2.20	0.41
2:F:286:TYR:CD1	2:F:286:TYR:N	2.88	0.41
2:H:256:SER:OG	2:H:257:LYS:N	2.53	0.41
1:C:1027:ILE:HG22	1:C:1033:LEU:HD21	2.03	0.41
2:H:214:ALA:HB1	2:H:224:TYR:CD2	2.56	0.41
2:F:200:LEU:CD1	2:F:204:ALA:HB2	2.49	0.40
1:A:1050:ILE:HD13	1:A:1052:ARG:O	2.21	0.40
1:A:1241:GLN:OE1	1:A:1244:ARG:NH1	2.55	0.40
2:F:257:LYS:HG3	2:F:258:ILE:HG23	2.02	0.40
1:G:1291:ALA:O	1:G:1295:VAL:HG23	2.22	0.40
1:A:1209:ILE:HD11	2:B:223:LEU:HD22	2.03	0.40
1:G:1318:GLY:CA	1:G:1319:GLN:HB2	2.51	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	326/364 (90%)	304 (93%)	22 (7%)	0	100	100
1	C	348/364 (96%)	313 (90%)	33 (10%)	2 (1%)	30	75
1	E	329/364 (90%)	303 (92%)	26 (8%)	0	100	100
1	G	330/364 (91%)	304 (92%)	26 (8%)	0	100	100
2	B	115/142 (81%)	109 (95%)	6 (5%)	0	100	100
2	D	122/142 (86%)	117 (96%)	5 (4%)	0	100	100
2	F	119/142 (84%)	111 (93%)	8 (7%)	0	100	100
2	H	122/142 (86%)	116 (95%)	6 (5%)	0	100	100
All	All	1811/2024 (90%)	1677 (93%)	132 (7%)	2 (0%)	56	90

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1	MET
1	C	100	VAL

### 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	268/292 (92%)	268 (100%)	0	100	100
1	C	281/292 (96%)	278 (99%)	3 (1%)	80	92
1	E	271/292 (93%)	271 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	270/292 (92%)	269 (100%)	1 (0%)	93	98
2	B	102/123 (83%)	101 (99%)	1 (1%)	82	93
2	D	109/123 (89%)	109 (100%)	0	100	100
2	F	106/123 (86%)	104 (98%)	2 (2%)	65	87
2	H	109/123 (89%)	108 (99%)	1 (1%)	84	94
All	All	1516/1660 (91%)	1508 (100%)	8 (0%)	92	97

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	241	HIS
1	C	1072	ASP
1	C	1114	PHE
1	C	4	MET
2	F	201	ASP
2	F	222	ASN
1	G	1072	ASP
2	H	260	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1220	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

There are no ligands in this entry.

## 5.7 Other polymers

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, 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	332/364 (91%)	-0.03	9 (2%) 58 47	52, 103, 148, 197	0
1	C	350/364 (96%)	-0.06	5 (1%) 78 68	46, 85, 128, 173	0
1	E	335/364 (92%)	0.17	22 (6%) 22 16	53, 119, 177, 198	0
1	G	336/364 (92%)	-0.02	7 (2%) 67 58	50, 90, 127, 148	0
2	B	116/142 (81%)	-0.13	5 (4%) 39 30	73, 110, 144, 153	0
2	D	123/142 (86%)	-0.20	1 (0%) 87 80	59, 79, 103, 116	0
2	F	120/142 (84%)	0.07	6 (5%) 32 25	80, 118, 151, 171	0
2	H	123/142 (86%)	-0.13	3 (2%) 62 52	58, 86, 113, 145	0
All	All	1835/2024 (90%)	-0.01	58 (3%) 51 42	46, 96, 155, 198	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1032	LEU	6.3
1	E	1040	ASN	6.0
1	A	1033	LEU	5.9
1	A	1058	ILE	5.9
1	G	1235	ALA	5.2
1	E	1015	LEU	4.7
1	E	1058	ILE	4.7
1	E	1041	ALA	4.2
1	E	1039	LEU	4.1
1	E	1237	PRO	4.1
1	E	1038	SER	3.8
1	A	1057	VAL	3.7
1	E	1031	HIS	3.7
1	C	1026	THR	3.7
2	F	316	PHE	3.3
1	E	1111	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	E	1112	ALA	3.1
1	E	1088	TYR	3.1
1	C	72	ALA	3.0
1	E	1079	LEU	3.0
1	C	1118	LEU	2.9
1	E	1102	MET	2.9
1	A	1153	PHE	2.9
1	G	1027	ILE	2.8
1	E	1016	LYS	2.8
1	G	1319	GLN	2.7
1	E	1240	LEU	2.7
2	B	312	TYR	2.7
1	G	1153	PHE	2.6
2	H	239	ASN	2.5
2	H	298	LEU	2.5
2	B	240	LEU	2.4
2	F	314	ILE	2.4
1	E	1060	LYS	2.4
2	B	255	ALA	2.4
1	A	1101	ASN	2.4
1	C	1017	ILE	2.3
1	E	1138	TRP	2.3
2	F	253	ILE	2.3
1	A	1102	MET	2.3
2	F	203	ALA	2.3
1	G	1320	LYS	2.3
1	E	1028	GLY	2.3
1	A	1100	ILE	2.3
2	F	237	LYS	2.3
1	G	1017	ILE	2.2
1	G	1028	GLY	2.2
1	E	1082	ALA	2.2
1	E	1030	GLY	2.2
1	E	1118	LEU	2.2
2	B	242	LEU	2.2
2	B	254	LEU	2.2
2	F	313	SER	2.2
1	E	1046	LEU	2.2
1	C	1100	ILE	2.2
2	D	263	ILE	2.1
2	H	263	ILE	2.1
1	A	1242	ALA	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.