



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:02 PM GMT

PDB ID : 2DDK  
Title : Crystal structure of human myo-inositol monophosphatase 2 (IMPA2) (orthorhombic form)  
Authors : Arai, R.; Ito, K.; Kamo-Uchikubo, T.; Bessho, Y.; Ohba, H.; Ohnishi, T.; Yoshikawa, T.; Shirouzu, M.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2006-01-30  
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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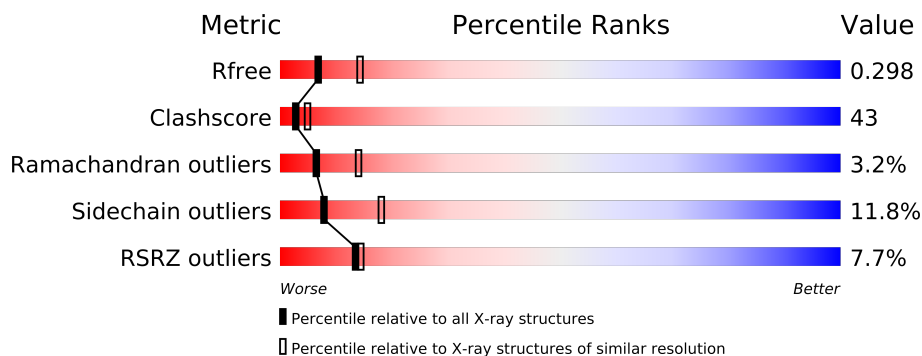
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.70 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))
$R_{free}$	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	299	
1	B	299	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3913 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Inositol monophosphatase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1970	1243	354	361	12			
1	B	249	Total	C	N	O	S	0	0	0
			1921	1216	344	349	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	CLONING ARTIFACT	UNP O14732
A	-9	GLY	-	CLONING ARTIFACT	UNP O14732
A	-8	SER	-	CLONING ARTIFACT	UNP O14732
A	-7	HIS	-	CLONING ARTIFACT	UNP O14732
A	-6	MET	-	CLONING ARTIFACT	UNP O14732
A	-5	GLU	-	CLONING ARTIFACT	UNP O14732
A	-4	LEU	-	CLONING ARTIFACT	UNP O14732
A	-3	PRO	-	CLONING ARTIFACT	UNP O14732
A	-2	GLY	-	CLONING ARTIFACT	UNP O14732
A	-1	SER	-	CLONING ARTIFACT	UNP O14732
A	0	SER	-	CLONING ARTIFACT	UNP O14732
B	-10	GLY	-	CLONING ARTIFACT	UNP O14732
B	-9	GLY	-	CLONING ARTIFACT	UNP O14732
B	-8	SER	-	CLONING ARTIFACT	UNP O14732
B	-7	HIS	-	CLONING ARTIFACT	UNP O14732
B	-6	MET	-	CLONING ARTIFACT	UNP O14732
B	-5	GLU	-	CLONING ARTIFACT	UNP O14732
B	-4	LEU	-	CLONING ARTIFACT	UNP O14732
B	-3	PRO	-	CLONING ARTIFACT	UNP O14732
B	-2	GLY	-	CLONING ARTIFACT	UNP O14732
B	-1	SER	-	CLONING ARTIFACT	UNP O14732
B	0	SER	-	CLONING ARTIFACT	UNP O14732

- Molecule 2 is water.

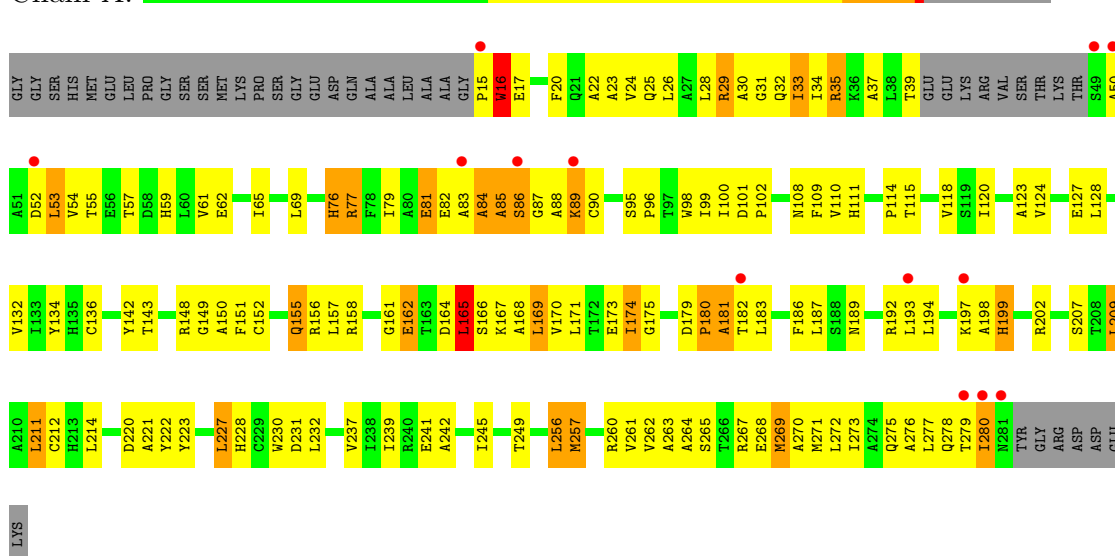
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total 9	O 9	0	0
2	B	13	Total 13	O 13	0	0

### 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 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.

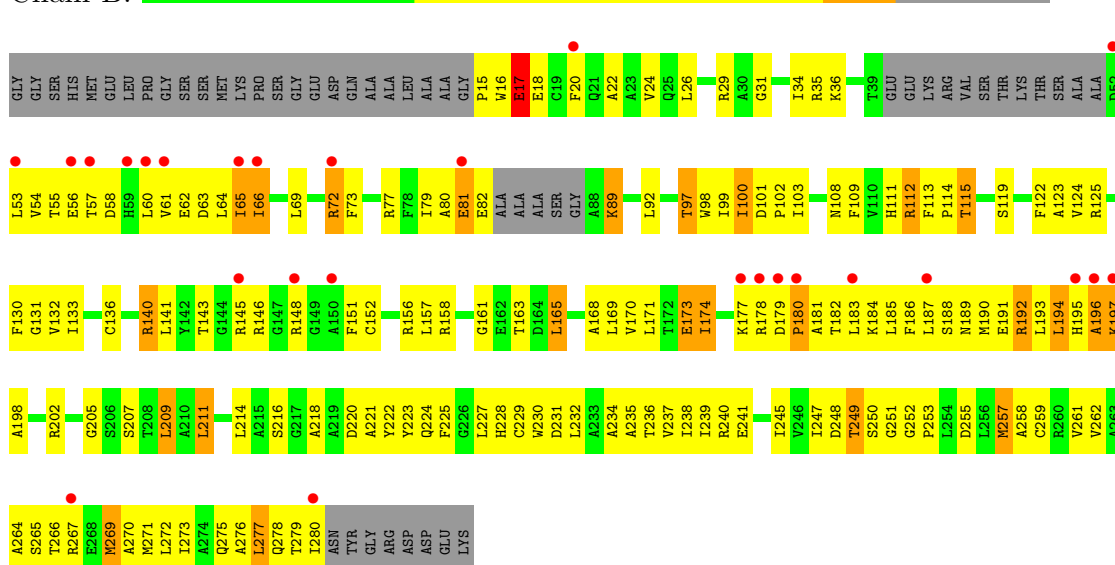
#### • Molecule 1: Inositol monophosphatase 2

Chain A:



#### • Molecule 1: Inositol monophosphatase 2

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.46Å 95.64Å 107.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.82 – 2.70 47.82 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.8 (47.82-2.70) 94.6 (47.82-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.260 , 0.301 0.257 , 0.298	Depositor DCC
$R_{free}$ test set	1788 reflections (9.87%)	DCC
Wilson B-factor (Å <sup>2</sup> )	63.3	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 46.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 18986 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3913	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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	0.43	0/2005	0.74	2/2717 (0.1%)
1	B	0.41	0/1955	0.72	0/2647
All	All	0.42	0/3960	0.73	2/5364 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	TRP	N-CA-C	7.62	131.56	111.00
1	A	118	VAL	N-CA-C	-5.18	97.01	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1970	0	1982	163	0
1	B	1921	0	1937	183	0
2	A	9	0	0	1	0
2	B	13	0	0	0	0
All	All	3913	0	3919	338	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 43.

All (338) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:192:ARG:HH21	1:B:192:ARG:HB2	1.17	1.07
1:B:34:ILE:HD11	1:B:136:CYS:SG	1.96	1.06
1:B:192:ARG:HH11	1:B:278:GLN:HB2	1.21	1.04
1:B:192:ARG:HB2	1:B:192:ARG:NH2	1.74	1.02
1:B:171:LEU:HD23	1:B:202:ARG:HB2	1.39	1.02
1:A:84:ALA:HB1	1:A:89:LYS:CE	1.90	1.01
1:B:81:GLU:HG2	1:B:230:TRP:CE2	1.96	1.00
1:A:15:PRO:HD2	1:A:124:VAL:HG21	1.44	1.00
1:A:53:LEU:H	1:A:53:LEU:HD23	1.24	0.99
1:A:171:LEU:HD22	1:A:202:ARG:HB2	1.44	0.97
1:A:33:ILE:HD11	1:A:61:VAL:HG22	1.46	0.97
1:B:228:HIS:HD1	1:B:230:TRP:HZ2	1.14	0.95
1:B:189:ASN:HD22	1:B:192:ARG:HH12	1.12	0.95
1:A:84:ALA:HB1	1:A:89:LYS:CD	1.97	0.95
1:B:250:SER:O	1:B:252:GLY:N	2.04	0.91
1:B:189:ASN:HA	1:B:192:ARG:HH22	1.37	0.90
1:A:87:GLY:O	1:A:89:LYS:HE3	1.72	0.89
1:B:214:LEU:HD23	1:B:239:ILE:HD13	1.55	0.89
1:B:192:ARG:NH1	1:B:278:GLN:HB2	1.87	0.88
1:B:193:LEU:HD22	1:B:223:TYR:HE2	1.38	0.88
1:B:22:ALA:O	1:B:26:LEU:HD23	1.78	0.83
1:B:165:LEU:HA	1:B:269:MET:HE1	1.60	0.82
1:B:53:LEU:C	1:B:55:THR:H	1.83	0.80
1:B:80:ALA:O	1:B:82:GLU:N	2.15	0.80
1:A:83:ALA:O	1:A:85:ALA:N	2.15	0.80
1:A:88:ALA:C	1:A:89:LYS:HG3	2.00	0.80
1:A:35:ARG:HD2	1:A:136:CYS:O	1.82	0.80
1:A:84:ALA:HB1	1:A:89:LYS:HE2	1.64	0.78
1:A:100:ILE:CD1	1:A:120:ILE:HG23	2.15	0.77
1:B:24:VAL:HG13	1:B:141:LEU:HD21	1.65	0.76
1:B:192:ARG:HH21	1:B:192:ARG:CB	1.99	0.76
1:A:170:VAL:HG11	1:A:193:LEU:HD21	1.68	0.75
1:A:193:LEU:HD22	1:A:223:TYR:HE2	1.50	0.75
1:B:236:THR:O	1:B:240:ARG:HG3	1.86	0.75
1:B:130:PHE:C	1:B:237:VAL:HG11	2.07	0.75
1:B:133:ILE:HD13	1:B:238:ILE:HD12	1.67	0.75
1:B:245:ILE:HD11	1:B:270:ALA:CB	2.17	0.74
1:B:190:MET:O	1:B:194:LEU:HB2	1.87	0.74
1:B:189:ASN:ND2	1:B:192:ARG:HH12	1.86	0.73
1:A:278:GLN:O	1:A:279:THR:HG23	1.88	0.73
1:B:132:VAL:HG22	1:B:143:THR:HG22	1.70	0.73
1:B:133:ILE:CD1	1:B:238:ILE:HD12	2.19	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:81:GLU:OE2	1:A:101:ASP:HA	1.87	0.73
1:A:148:ARG:HH21	1:A:148:ARG:HG3	1.53	0.73
1:A:84:ALA:CA	1:A:89:LYS:HG2	2.19	0.72
1:B:181:ALA:C	1:B:183:LEU:H	1.93	0.72
1:B:179:ASP:OD1	1:B:182:THR:HG23	1.88	0.72
1:A:84:ALA:CB	1:A:89:LYS:HG2	2.21	0.71
1:A:15:PRO:CD	1:A:124:VAL:HG21	2.22	0.70
1:B:271:MET:O	1:B:275:GLN:HG3	1.91	0.70
1:B:99:ILE:HG22	1:B:230:TRP:HA	1.73	0.69
1:B:57:THR:O	1:B:61:VAL:HG22	1.91	0.69
1:A:84:ALA:HA	1:A:89:LYS:HG2	1.74	0.69
1:A:166:SER:HA	1:A:197:LYS:O	1.92	0.69
1:A:84:ALA:HB1	1:A:89:LYS:CG	2.22	0.69
1:B:15:PRO:HA	1:B:122:PHE:CE1	2.29	0.68
1:A:87:GLY:O	1:A:89:LYS:CE	2.42	0.68
1:A:192:ARG:HD3	1:A:276:ALA:O	1.93	0.68
1:A:101:ASP:OD2	1:A:231:ASP:OD1	2.12	0.68
1:A:84:ALA:HB1	1:A:89:LYS:HG2	1.76	0.68
1:A:89:LYS:O	1:A:257:MET:HE1	1.95	0.67
1:A:149:GLY:CA	1:A:241:GLU:HG3	2.23	0.67
1:B:130:PHE:HA	1:B:237:VAL:HG21	1.76	0.67
1:B:196:ALA:O	1:B:197:LYS:HB2	1.95	0.67
1:A:152:CYS:CB	1:A:157:LEU:HD11	2.25	0.66
1:A:151:PHE:CD1	1:A:156:ARG:HA	2.30	0.66
1:A:168:ALA:HA	1:A:220:ASP:OD2	1.94	0.66
1:A:174:ILE:HD13	1:A:175:GLY:N	2.11	0.65
1:B:183:LEU:HD11	1:B:187:LEU:HD23	1.78	0.65
1:A:264:ALA:HB1	1:A:269:MET:HB3	1.77	0.65
1:B:189:ASN:CA	1:B:192:ARG:HH22	2.07	0.65
1:B:174:ILE:HG23	1:B:186:PHE:CD1	2.32	0.65
1:B:193:LEU:HD22	1:B:223:TYR:CE2	2.28	0.65
1:A:193:LEU:HD22	1:A:223:TYR:CE2	2.31	0.64
1:B:228:HIS:ND1	1:B:230:TRP:CZ2	2.64	0.64
1:B:195:HIS:C	1:B:197:LYS:H	1.98	0.64
1:A:22:ALA:O	1:A:26:LEU:HD23	1.97	0.64
1:A:57:THR:O	1:A:61:VAL:HG23	1.98	0.64
1:A:165:LEU:O	1:A:198:ALA:HA	1.97	0.64
1:A:99:ILE:HG22	1:A:230:TRP:HA	1.79	0.64
1:A:33:ILE:HD11	1:A:61:VAL:CG2	2.25	0.64
1:A:54:VAL:HG12	1:A:109:PHE:CD2	2.34	0.63
1:A:89:LYS:O	1:A:90:CYS:HB2	1.98	0.63
1:B:247:ILE:HB	1:B:252:GLY:O	1.98	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:131:GLY:N	1:B:237:VAL:HG11	2.14	0.63
1:B:245:ILE:HD11	1:B:270:ALA:HB3	1.80	0.63
1:B:29:ARG:HH21	1:B:64:LEU:HD21	1.64	0.63
1:B:53:LEU:C	1:B:55:THR:N	2.52	0.63
1:B:189:ASN:HA	1:B:192:ARG:NH2	2.11	0.62
1:B:278:GLN:HG2	1:B:280:ILE:HG13	1.81	0.62
1:B:100:ILE:C	1:B:100:ILE:HD12	2.20	0.61
1:B:232:LEU:H	1:B:232:LEU:HD23	1.63	0.61
1:A:151:PHE:HA	1:A:157:LEU:HD13	1.81	0.61
1:A:245:ILE:HD11	1:A:267:ARG:HA	1.82	0.61
1:B:89:LYS:HE3	1:B:257:MET:HE3	1.82	0.61
1:B:62:GLU:O	1:B:65:ILE:HG22	1.99	0.61
1:A:158:ARG:HD3	1:A:242:ALA:O	2.01	0.61
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.82	0.61
1:B:69:LEU:HD23	1:B:98:TRP:CE2	2.36	0.60
1:B:181:ALA:C	1:B:183:LEU:N	2.54	0.60
1:B:185:LEU:HD13	1:B:280:ILE:O	2.01	0.60
1:B:174:ILE:HD12	1:B:174:ILE:H	1.67	0.60
1:B:64:LEU:HD22	1:B:64:LEU:O	2.01	0.60
1:A:81:GLU:OE2	1:A:102:PRO:HD2	2.02	0.60
1:B:245:ILE:CD1	1:B:267:ARG:HA	2.31	0.60
1:B:89:LYS:HG3	1:B:257:MET:CE	2.31	0.60
1:A:261:VAL:HG22	1:A:262:VAL:N	2.16	0.60
1:A:50:ALA:HB3	1:A:52:ASP:OD1	2.01	0.60
1:A:50:ALA:C	1:A:52:ASP:H	2.03	0.60
1:B:188:SER:O	1:B:192:ARG:NH2	2.34	0.60
1:B:189:ASN:HD22	1:B:192:ARG:NH1	1.93	0.59
1:B:261:VAL:HG22	1:B:262:VAL:N	2.17	0.59
1:B:81:GLU:HG2	1:B:230:TRP:NE1	2.17	0.59
1:A:54:VAL:HG11	1:A:110:VAL:HG22	1.84	0.59
1:A:245:ILE:CD1	1:A:267:ARG:HA	2.32	0.59
1:A:268:GLU:O	1:A:272:LEU:HD13	2.03	0.58
1:B:145:ARG:HH22	1:B:148:ARG:HH21	1.51	0.58
1:B:195:HIS:C	1:B:197:LYS:N	2.55	0.58
1:A:214:LEU:C	1:A:214:LEU:HD23	2.24	0.58
1:B:174:ILE:HG23	1:B:186:PHE:CE1	2.39	0.58
1:B:115:THR:HG22	1:B:209:LEU:HD23	1.86	0.58
1:B:225:PHE:CZ	1:B:249:THR:HG21	2.39	0.57
1:B:165:LEU:HG	1:B:269:MET:HE1	1.86	0.57
1:A:271:MET:O	1:A:275:GLN:HB2	2.03	0.57
1:A:76:HIS:HD2	1:A:98:TRP:HE1	1.51	0.57
1:A:30:ALA:O	1:A:33:ILE:HD13	2.05	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:165:LEU:HA	1:B:269:MET:CE	2.33	0.57
1:A:161:GLY:H	1:A:265:SER:HB2	1.68	0.57
1:A:15:PRO:HD2	1:A:124:VAL:CG2	2.24	0.57
1:B:227:LEU:HD23	1:B:227:LEU:H	1.70	0.57
1:A:108:ASN:OD1	1:B:202:ARG:NH1	2.32	0.57
1:B:101:ASP:OD2	1:B:231:ASP:OD1	2.23	0.57
1:B:53:LEU:HA	1:B:55:THR:HG23	1.86	0.57
1:B:89:LYS:NZ	1:B:259:CYS:SG	2.76	0.57
1:B:170:VAL:HG11	1:B:193:LEU:HD21	1.87	0.57
1:B:245:ILE:HD11	1:B:270:ALA:HB2	1.86	0.57
1:A:54:VAL:CG2	1:A:55:THR:N	2.68	0.57
1:A:89:LYS:O	1:A:257:MET:CE	2.53	0.56
1:A:84:ALA:O	1:A:89:LYS:HE2	2.05	0.56
1:A:35:ARG:CD	1:A:136:CYS:O	2.52	0.56
1:A:148:ARG:HG3	1:A:148:ARG:NH2	2.20	0.56
1:B:111:HIS:C	1:B:112:ARG:HE	2.08	0.56
1:A:79:ILE:HD12	1:A:99:ILE:HD12	1.87	0.56
1:B:229:CYS:HA	1:B:232:LEU:CD2	2.35	0.56
1:B:171:LEU:HD23	1:B:202:ARG:CB	2.25	0.56
1:B:163:THR:HG22	1:B:266:THR:OG1	2.06	0.56
1:A:84:ALA:O	1:A:86:SER:N	2.39	0.56
1:A:33:ILE:CD1	1:A:61:VAL:HG22	2.27	0.56
1:B:64:LEU:C	1:B:64:LEU:HD13	2.26	0.56
1:B:92:LEU:HD12	1:B:97:THR:HG21	1.87	0.56
1:A:76:HIS:CD2	1:A:98:TRP:HE1	2.24	0.56
1:A:152:CYS:HB3	1:A:157:LEU:HD11	1.87	0.56
1:B:65:ILE:C	1:B:65:ILE:HD13	2.26	0.56
1:B:81:GLU:HG2	1:B:230:TRP:CZ2	2.40	0.56
1:A:54:VAL:HG23	1:A:55:THR:N	2.21	0.55
1:B:183:LEU:CD1	1:B:187:LEU:HD23	2.34	0.55
1:B:77:ARG:HH21	1:B:77:ARG:HG2	1.70	0.55
1:A:83:ALA:C	1:A:85:ALA:H	2.09	0.55
1:A:170:VAL:C	1:A:171:LEU:HD23	2.27	0.55
1:A:100:ILE:HD12	1:A:120:ILE:HG23	1.86	0.55
1:A:53:LEU:CD2	1:A:53:LEU:H	2.01	0.55
1:B:72:ARG:HB3	1:B:73:PHE:CE1	2.42	0.55
1:B:228:HIS:HB2	1:B:230:TRP:CE2	2.42	0.54
1:B:157:LEU:HD22	1:B:216:SER:HA	1.89	0.54
1:B:161:GLY:H	1:B:265:SER:HB2	1.70	0.54
1:B:272:LEU:O	1:B:275:GLN:HB2	2.07	0.54
1:B:185:LEU:HD13	1:B:280:ILE:C	2.26	0.54
1:B:69:LEU:HD23	1:B:98:TRP:CD2	2.43	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:166:SER:O	1:A:199:HIS:HB2	2.07	0.54
1:A:54:VAL:HG12	1:A:109:PHE:CE2	2.43	0.54
1:A:183:LEU:HD11	1:B:194:LEU:HD22	1.89	0.54
1:B:165:LEU:HD13	1:B:196:ALA:HB1	1.90	0.54
1:B:186:PHE:HD2	1:B:187:LEU:HD22	1.71	0.54
1:B:54:VAL:HG11	1:B:109:PHE:HE1	1.73	0.54
1:B:177:LYS:C	1:B:178:ARG:HD2	2.28	0.54
1:B:62:GLU:O	1:B:63:ASP:C	2.45	0.54
1:A:183:LEU:O	1:A:187:LEU:HD23	2.08	0.54
1:B:192:ARG:HH11	1:B:278:GLN:CB	2.08	0.53
1:B:228:HIS:ND1	1:B:230:TRP:HZ2	1.94	0.53
1:B:211:LEU:O	1:B:214:LEU:HB3	2.08	0.53
1:A:221:ALA:HA	1:A:263:ALA:O	2.09	0.53
1:B:66:ILE:O	1:B:66:ILE:HD13	2.08	0.53
1:A:186:PHE:HD1	1:A:187:LEU:HD22	1.73	0.53
1:A:155:GLN:NE2	1:A:156:ARG:O	2.42	0.53
1:A:76:HIS:HD2	1:A:98:TRP:NE1	2.06	0.53
1:B:99:ILE:CG2	1:B:230:TRP:HA	2.36	0.53
1:A:30:ALA:HB1	1:A:61:VAL:HG13	1.90	0.53
1:B:89:LYS:HE3	1:B:257:MET:CE	2.39	0.53
1:B:99:ILE:N	1:B:99:ILE:HD12	2.23	0.53
1:B:195:HIS:O	1:B:197:LYS:N	2.42	0.53
1:B:97:THR:HG23	1:B:123:ALA:HB3	1.91	0.53
1:A:199:HIS:HD2	1:B:111:HIS:HE1	1.57	0.53
1:B:111:HIS:O	1:B:112:ARG:HB2	2.08	0.53
1:A:111:HIS:O	1:B:218:ALA:HA	2.09	0.53
1:A:173:GLU:HG3	2:A:290:HOH:O	2.09	0.52
1:A:179:ASP:OD1	1:A:181:ALA:HB3	2.08	0.52
1:B:56:GLU:O	1:B:60:LEU:HG	2.10	0.52
1:B:108:ASN:HB3	1:B:113:PHE:O	2.08	0.52
1:B:183:LEU:HD11	1:B:187:LEU:CD2	2.39	0.52
1:A:150:ALA:HB2	1:A:241:GLU:HB3	1.92	0.52
1:B:178:ARG:N	1:B:178:ARG:HD2	2.25	0.52
1:A:115:THR:HG23	1:A:209:LEU:HD23	1.92	0.51
1:A:249:THR:OG1	1:A:260:ARG:HD2	2.10	0.51
1:A:99:ILE:CG2	1:A:230:TRP:HA	2.40	0.51
1:A:192:ARG:HD2	1:A:277:LEU:HA	1.93	0.51
1:A:83:ALA:C	1:A:85:ALA:N	2.62	0.51
1:A:29:ARG:O	1:A:32:GLN:HB2	2.11	0.51
1:A:149:GLY:C	1:A:241:GLU:HG3	2.31	0.51
1:A:151:PHE:HD1	1:A:156:ARG:HA	1.73	0.51
1:B:54:VAL:HG11	1:B:109:PHE:CE1	2.45	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:80:ALA:CB	1:B:100:ILE:HG13	2.41	0.51
1:A:84:ALA:C	1:A:86:SER:N	2.65	0.50
1:A:169:LEU:HB2	1:B:111:HIS:CD2	2.46	0.50
1:A:84:ALA:CB	1:A:89:LYS:CG	2.87	0.50
1:B:92:LEU:CD1	1:B:97:THR:HG21	2.42	0.50
1:A:33:ILE:HD13	1:A:34:ILE:N	2.27	0.50
1:B:29:ARG:NH2	1:B:64:LEU:HD21	2.26	0.50
1:A:16:TRP:HD1	1:A:17:GLU:N	2.10	0.50
1:B:257:MET:HG3	1:B:257:MET:O	2.12	0.49
1:A:142:TYR:CE2	1:A:212:CYS:HB3	2.47	0.49
1:B:184:LYS:O	1:B:188:SER:OG	2.26	0.49
1:B:245:ILE:HD12	1:B:267:ARG:HD2	1.94	0.49
1:A:30:ALA:HA	1:A:33:ILE:HD12	1.95	0.49
1:B:16:TRP:HD1	1:B:17:GLU:H	1.53	0.49
1:A:85:ALA:C	1:A:87:GLY:H	2.16	0.49
1:B:80:ALA:HB2	1:B:100:ILE:HG13	1.94	0.49
1:B:61:VAL:O	1:B:65:ILE:HG22	2.13	0.48
1:B:156:ARG:NH1	1:B:158:ARG:NH1	2.62	0.48
1:B:229:CYS:HA	1:B:232:LEU:HD21	1.96	0.48
1:A:62:GLU:OE2	1:A:82:GLU:CB	2.62	0.48
1:B:58:ASP:OD2	1:B:102:PRO:HG3	2.14	0.48
1:B:192:ARG:HB3	1:B:276:ALA:HB1	1.96	0.47
1:B:146:ARG:HH21	1:B:146:ARG:HG3	1.79	0.47
1:A:84:ALA:O	1:A:85:ALA:C	2.50	0.47
1:B:62:GLU:O	1:B:65:ILE:N	2.45	0.47
1:B:272:LEU:HD23	1:B:275:GLN:NE2	2.29	0.47
1:A:132:VAL:HG13	1:A:143:THR:HG22	1.96	0.47
1:B:31:GLY:O	1:B:35:ARG:HB2	2.15	0.47
1:B:229:CYS:HA	1:B:232:LEU:HD23	1.96	0.47
1:A:179:ASP:O	1:A:181:ALA:N	2.47	0.47
1:B:165:LEU:HD13	1:B:196:ALA:CB	2.44	0.47
1:A:31:GLY:HA3	1:A:134:TYR:CD1	2.49	0.47
1:B:20:PHE:CE2	1:B:145:ARG:HD2	2.50	0.47
1:A:194:LEU:HD22	1:B:178:ARG:NH2	2.30	0.47
1:B:225:PHE:HE1	1:B:277:LEU:HD11	1.79	0.47
1:A:194:LEU:HD22	1:B:178:ARG:HH21	1.79	0.47
1:A:280:ILE:N	1:A:280:ILE:HD13	2.30	0.47
1:A:34:ILE:HG23	1:A:109:PHE:HE1	1.80	0.46
1:B:225:PHE:CE1	1:B:277:LEU:HD11	2.50	0.46
1:A:228:HIS:HB3	1:A:230:TRP:CE2	2.50	0.46
1:A:214:LEU:HD11	1:A:263:ALA:C	2.36	0.46
1:A:245:ILE:HD11	1:A:270:ALA:CB	2.46	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:26:LEU:HD12	1:B:65:ILE:HA	1.97	0.46
1:A:261:VAL:CG2	1:A:262:VAL:N	2.78	0.46
1:B:151:PHE:CD1	1:B:156:ARG:HA	2.51	0.46
1:B:168:ALA:HA	1:B:220:ASP:OD1	2.15	0.46
1:B:15:PRO:HA	1:B:122:PHE:CZ	2.50	0.46
1:B:184:LYS:HD3	1:B:184:LYS:HA	1.64	0.46
1:A:50:ALA:C	1:A:52:ASP:N	2.69	0.46
1:A:214:LEU:HD11	1:A:263:ALA:O	2.16	0.45
1:B:156:ARG:NH1	1:B:158:ARG:HH12	2.15	0.45
1:B:278:GLN:C	1:B:280:ILE:N	2.70	0.45
1:A:123:ALA:HA	1:A:127:GLU:O	2.17	0.45
1:A:84:ALA:HB1	1:A:89:LYS:HD3	1.92	0.44
1:B:100:ILE:HD12	1:B:101:ASP:N	2.32	0.44
1:A:69:LEU:HD13	1:A:98:TRP:CD2	2.52	0.44
1:B:278:GLN:O	1:B:280:ILE:N	2.50	0.44
1:B:89:LYS:HG3	1:B:257:MET:HE1	1.99	0.44
1:B:102:PRO:HG2	1:B:103:ILE:H	1.82	0.44
1:A:227:LEU:CD1	1:A:227:LEU:O	2.65	0.44
1:A:220:ASP:O	1:A:264:ALA:HA	2.18	0.44
1:A:54:VAL:HG12	1:A:109:PHE:HD2	1.83	0.44
1:B:278:GLN:C	1:B:280:ILE:H	2.21	0.43
1:B:145:ARG:O	1:B:241:GLU:HG3	2.18	0.43
1:A:179:ASP:O	1:A:180:PRO:C	2.56	0.43
1:A:162:GLU:OE2	1:A:167:LYS:HB2	2.18	0.43
1:B:224:GLN:HB3	1:B:227:LEU:HD22	2.00	0.43
1:A:171:LEU:O	1:A:222:TYR:HA	2.18	0.43
1:B:211:LEU:HB3	1:B:235:ALA:HB1	2.00	0.43
1:A:100:ILE:HG22	1:A:101:ASP:N	2.33	0.43
1:A:81:GLU:H	1:A:81:GLU:HG2	1.10	0.43
1:A:85:ALA:O	1:A:87:GLY:N	2.52	0.43
1:A:29:ARG:HA	1:A:32:GLN:NE2	2.33	0.43
1:A:207:SER:O	1:A:211:LEU:HD22	2.19	0.43
1:A:111:HIS:CE1	1:B:169:LEU:HB2	2.54	0.43
1:B:255:ASP:O	1:B:258:ALA:HB3	2.19	0.43
1:A:193:LEU:HD11	1:A:273:ILE:HD12	2.01	0.43
1:A:20:PHE:O	1:A:23:ALA:HB3	2.19	0.43
1:A:55:THR:HG22	1:A:59:HIS:CE1	2.53	0.43
1:B:80:ALA:O	1:B:81:GLU:C	2.58	0.42
1:A:211:LEU:HG	1:A:239:ILE:HD11	2.00	0.42
1:B:248:ASP:HB2	1:B:253:PRO:O	2.19	0.42
1:B:119:SER:OG	1:B:234:ALA:HB3	2.19	0.42
1:A:88:ALA:C	1:A:89:LYS:CG	2.76	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:149:GLY:HA2	1:A:241:GLU:HG3	2.00	0.42
1:A:232:LEU:HD11	1:A:256:LEU:HD12	2.02	0.42
1:B:140:ARG:HB3	1:B:152:CYS:SG	2.60	0.42
1:B:131:GLY:O	1:B:143:THR:HA	2.20	0.42
1:B:111:HIS:O	1:B:112:ARG:CB	2.66	0.42
1:B:124:VAL:O	1:B:125:ARG:HB2	2.19	0.42
1:A:24:VAL:O	1:A:25:GLN:C	2.57	0.42
1:B:132:VAL:HG22	1:B:143:THR:CG2	2.46	0.42
1:B:224:GLN:HG2	1:B:225:PHE:N	2.35	0.42
1:B:173:GLU:OE2	1:B:205:GLY:N	2.50	0.42
1:A:189:ASN:ND2	1:A:277:LEU:CB	2.83	0.42
1:A:269:MET:O	1:A:273:ILE:HG12	2.20	0.42
1:A:279:THR:HB	1:A:280:ILE:HD13	2.02	0.42
1:A:227:LEU:CD1	1:A:232:LEU:HD23	2.50	0.41
1:A:128:LEU:HB3	1:A:237:VAL:CG2	2.50	0.41
1:B:197:LYS:O	1:B:198:ALA:C	2.59	0.41
1:A:189:ASN:HD22	1:A:277:LEU:HD22	1.85	0.41
1:B:273:ILE:O	1:B:277:LEU:HB2	2.19	0.41
1:B:100:ILE:CD1	1:B:100:ILE:C	2.86	0.41
1:A:164:ASP:C	1:A:166:SER:H	2.23	0.41
1:A:37:ALA:C	1:A:39:THR:H	2.22	0.41
1:B:79:ILE:O	1:B:99:ILE:HA	2.21	0.41
1:A:249:THR:HA	1:A:262:VAL:HG23	2.03	0.41
1:A:84:ALA:CB	1:A:89:LYS:CD	2.85	0.41
1:A:232:LEU:C	1:A:232:LEU:HD12	2.41	0.41
1:A:223:TYR:CD1	1:A:223:TYR:C	2.93	0.41
1:B:222:TYR:OH	1:B:224:GLN:NE2	2.54	0.41
1:B:228:HIS:CB	1:B:230:TRP:CE2	3.04	0.41
1:B:207:SER:O	1:B:211:LEU:HD22	2.21	0.41
1:A:77:ARG:HA	1:A:77:ARG:NE	2.35	0.41
1:A:29:ARG:O	1:A:33:ILE:HG23	2.21	0.40
1:A:152:CYS:HB2	1:A:157:LEU:HD11	1.98	0.40
1:B:232:LEU:N	1:B:232:LEU:HD23	2.33	0.40
1:A:214:LEU:CD1	1:A:263:ALA:O	2.69	0.40
1:B:62:GLU:HA	1:B:65:ILE:HG22	2.03	0.40
1:A:69:LEU:HD22	1:A:98:TRP:CZ2	2.56	0.40
1:A:269:MET:HE2	1:A:269:MET:HB3	1.87	0.40
1:B:227:LEU:O	1:B:259:CYS:HA	2.22	0.40
1:B:221:ALA:HA	1:B:264:ALA:HA	2.03	0.40
1:B:65:ILE:O	1:B:69:LEU:HD13	2.21	0.40
1:A:26:LEU:HD12	1:A:65:ILE:HA	2.03	0.40
1:B:225:PHE:HE1	1:B:277:LEU:CD1	2.34	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:108:ASN:O	1:B:109:PHE:C	2.58	0.40
1:A:227:LEU:HD13	1:A:232:LEU:HD23	2.03	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	254/299 (85%)	216 (85%)	30 (12%)	8 (3%)	7	14
1	B	243/299 (81%)	198 (82%)	37 (15%)	8 (3%)	6	13
All	All	497/598 (83%)	414 (83%)	67 (14%)	16 (3%)	6	14

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	ALA
1	B	81	GLU
1	B	180	PRO
1	B	251	GLY
1	A	85	ALA
1	A	86	SER
1	A	181	ALA
1	A	199	HIS
1	B	17	GLU
1	B	114	PRO
1	A	114	PRO
1	B	279	THR
1	A	165	LEU
1	B	196	ALA
1	A	180	PRO
1	B	197	LYS



### 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	205/236 (87%)	182 (89%)	23 (11%)	9	20
1	B	202/236 (86%)	177 (88%)	25 (12%)	7	16
All	All	407/472 (86%)	359 (88%)	48 (12%)	8	18

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	TRP
1	A	28	LEU
1	A	29	ARG
1	A	33	ILE
1	A	35	ARG
1	A	53	LEU
1	A	76	HIS
1	A	77	ARG
1	A	81	GLU
1	A	89	LYS
1	A	155	GLN
1	A	162	GLU
1	A	165	LEU
1	A	169	LEU
1	A	174	ILE
1	A	182	THR
1	A	209	LEU
1	A	211	LEU
1	A	227	LEU
1	A	256	LEU
1	A	257	MET
1	A	269	MET
1	A	280	ILE
1	B	17	GLU
1	B	18	GLU
1	B	36	LYS
1	B	65	ILE
1	B	66	ILE
1	B	72	ARG

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Mol	Chain	Res	Type
1	B	89	LYS
1	B	97	THR
1	B	100	ILE
1	B	112	ARG
1	B	115	THR
1	B	140	ARG
1	B	165	LEU
1	B	173	GLU
1	B	174	ILE
1	B	180	PRO
1	B	191	GLU
1	B	192	ARG
1	B	194	LEU
1	B	209	LEU
1	B	211	LEU
1	B	249	THR
1	B	257	MET
1	B	269	MET
1	B	277	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	59	HIS
1	A	76	HIS
1	A	135	HIS
1	A	155	GLN
1	A	189	ASN
1	A	199	HIS
1	B	32	GLN
1	B	135	HIS
1	B	189	ASN
1	B	199	HIS
1	B	224	GLN
1	B	275	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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	258/299 (86%)	0.30	13 (5%) 28 30	27, 54, 90, 112	0
1	B	249/299 (83%)	0.62	26 (10%) 7 7	31, 65, 100, 110	0
All	All	507/598 (84%)	0.46	39 (7%) 13 14	27, 60, 98, 112	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	279	THR	7.6
1	A	280	ILE	7.2
1	A	281	ASN	4.9
1	B	66	ILE	4.8
1	B	180	PRO	4.5
1	B	280	ILE	4.4
1	B	60	LEU	4.3
1	B	61	VAL	4.2
1	B	52	ASP	4.2
1	B	183	LEU	4.0
1	B	20	PHE	3.8
1	A	50	ALA	3.7
1	B	267	ARG	3.4
1	B	150	ALA	3.3
1	A	89	LYS	3.1
1	A	49	SER	3.0
1	A	15	PRO	2.9
1	B	57	THR	2.8
1	B	65	ILE	2.7
1	A	86	SER	2.7
1	B	187	LEU	2.7
1	A	83	ALA	2.6
1	B	197	LYS	2.4
1	A	193	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	177	LYS	2.4
1	B	81	GLU	2.4
1	B	195	HIS	2.4
1	B	196	ALA	2.4
1	B	72	ARG	2.3
1	A	52	ASP	2.3
1	A	197	LYS	2.3
1	B	53	LEU	2.2
1	B	148	ARG	2.2
1	B	145	ARG	2.2
1	B	56	GLU	2.2
1	A	182	THR	2.1
1	B	179	ASP	2.1
1	B	59	HIS	2.1
1	B	178	ARG	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 ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.