



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 20, 2023 – 03:08 PM JST

PDB ID : 7CS9  
Title : AtPrR1 in apo form  
Authors : Shao, K.; Zhang, P.  
Deposited on : 2020-08-14  
Resolution : 2.80 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

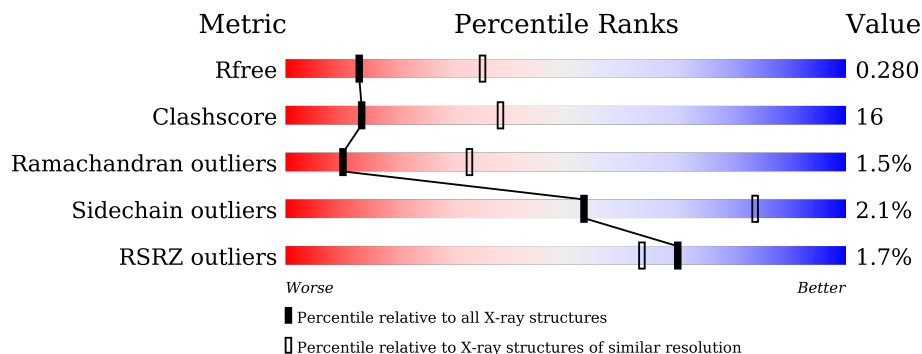
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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	317	 72% 22% . .
1	B	317	 75% 18% . .
1	C	317	 55% 38% . .
1	D	317	 67% 26% . .

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9564 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.

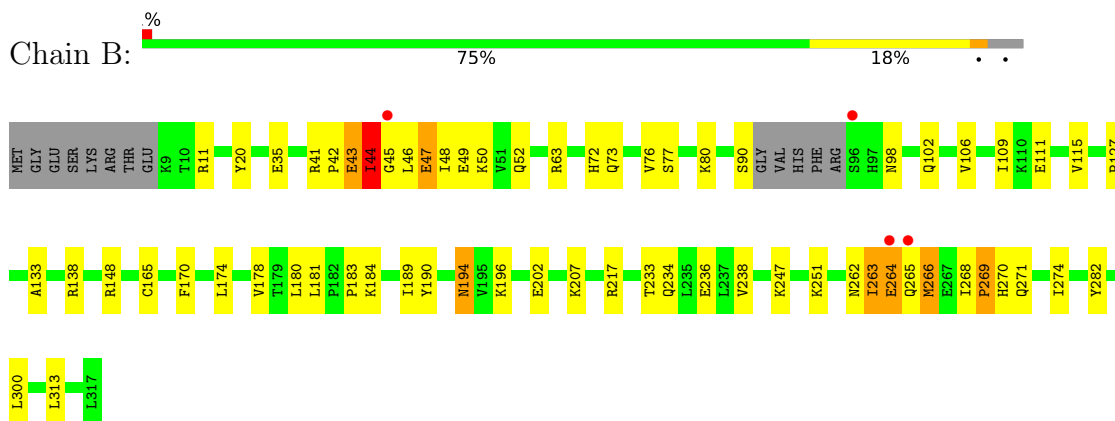
- Molecule 1 is a protein called Pinorensinol reductase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	304	2390	1528	404	447	11	0	0	0
1	C	304	2395	1533	404	447	11	0	0	0
1	A	303	2384	1525	403	445	11	0	0	0
1	D	304	2395	1533	404	447	11	0	0	0

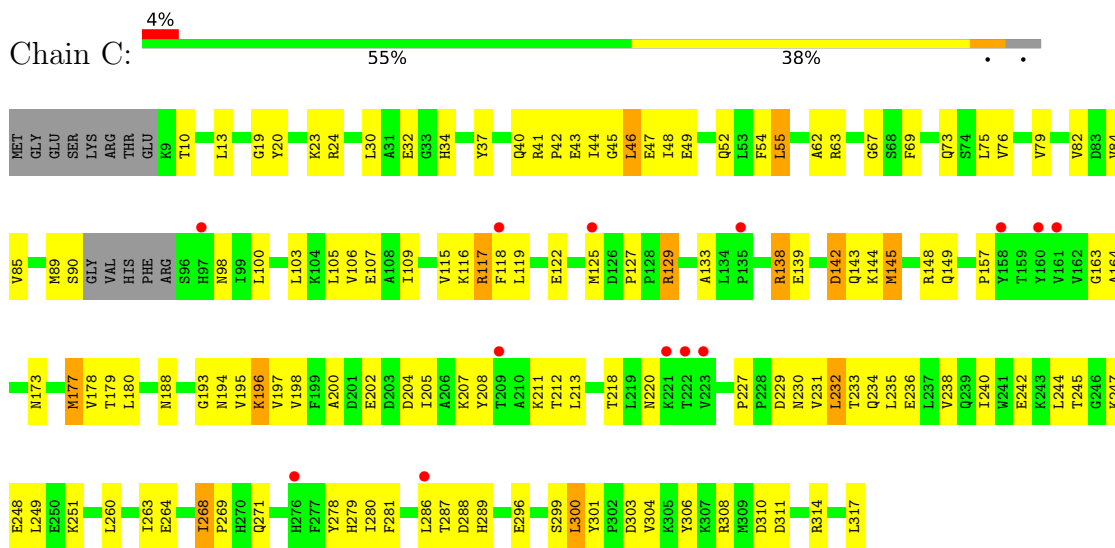
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

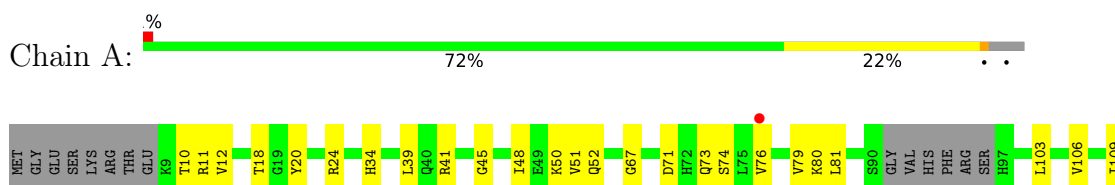
- Molecule 1: Pinoresinol reductase 1

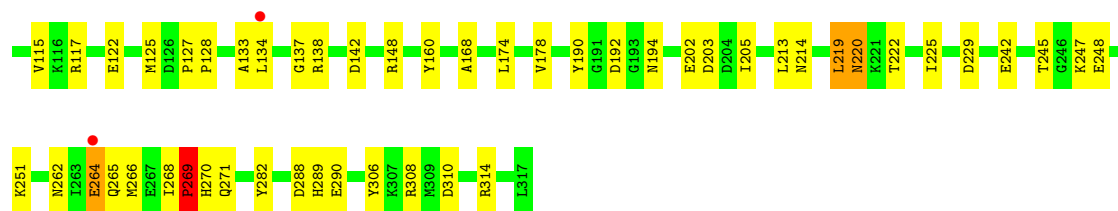


- Molecule 1: Pinoresinol reductase 1



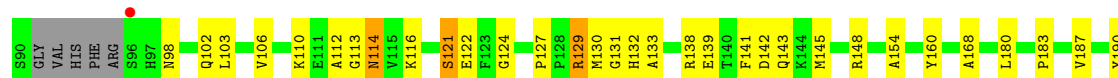
- Molecule 1: Pinoresinol reductase 1





● Molecule 1: Pinoreosinol reductase 1

Chain D: 67% 26%



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.73Å 134.92Å 142.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.27 – 2.80 39.27 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.27-2.80) 91.8 (39.27-2.80)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.80 (at 2.81Å)	Xtrriage
Refinement program	PHENIX 1.10_2155, PHENIX 1.10_2155	Depositor
R, $R_{free}$	0.221 , 0.280 0.221 , 0.280	Depositor DCC
$R_{free}$ test set	2003 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	54.5	Xtrriage
Anisotropy	0.623	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 36.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.022 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	9564	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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	0.61	1/2431 (0.0%)	0.67	1/3288 (0.0%)
1	B	0.55	0/2437	0.71	1/3296 (0.0%)
1	C	0.63	3/2442 (0.1%)	0.80	5/3303 (0.2%)
1	D	0.49	0/2442	0.68	3/3303 (0.1%)
All	All	0.57	4/9752 (0.0%)	0.72	10/13190 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	5
1	D	0	1
All	All	0	8

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	269	PRO	N-CD	-17.11	1.23	1.47
1	C	117	ARG	NE-CZ	-7.50	1.23	1.33
1	C	117	ARG	CZ-NH1	-6.92	1.24	1.33
1	C	196	LYS	CB-CG	6.90	1.71	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	232	LEU	CB-CG-CD1	-10.77	92.70	111.00
1	D	129	ARG	NE-CZ-NH1	-8.59	116.01	120.30
1	A	134	LEU	CA-CB-CG	7.10	131.62	115.30
1	B	45	GLY	N-CA-C	-6.99	95.62	113.10
1	C	300	LEU	CA-CB-CG	6.02	129.15	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	55	LEU	CA-CB-CG	5.71	128.42	115.30
1	C	196	LYS	CB-CG-CD	5.68	126.37	111.60
1	D	30	LEU	CB-CG-CD2	-5.44	101.75	111.00
1	D	129	ARG	NE-CZ-NH2	5.03	122.81	120.30
1	C	142	ASP	CB-CG-OD1	-5.01	113.79	118.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	264	GLU	Peptide
1	A	269	PRO	Peptide
1	B	263	ILE	Peptide
1	B	264	GLU	Peptide
1	B	269	PRO	Peptide
1	B	43	GLU	Peptide
1	B	44	ILE	Peptide
1	D	265	GLN	Peptide

## 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	2384	0	2390	62	0
1	B	2390	0	2395	44	0
1	C	2395	0	2409	135	0
1	D	2395	0	2411	74	0
All	All	9564	0	9605	306	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (306) 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:C:196:LYS:HZ2	1:C:233:THR:HG22	1.28	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:VAL:HG11	1:C:213:LEU:HD21	1.47	0.97
1:D:130:MET:O	1:D:138:ARG:NH2	2.02	0.93
1:D:208:TYR:HE2	1:D:304:VAL:HG11	1.32	0.93
1:C:242:GLU:HA	1:C:245:THR:HG22	1.51	0.92
1:B:184:LYS:O	1:B:247:LYS:NZ	2.03	0.91
1:C:196:LYS:NZ	1:C:233:THR:HG22	1.85	0.90
1:A:242:GLU:OE2	1:A:251:LYS:NZ	2.04	0.89
1:B:35:GLU:OE1	1:B:63:ARG:NH2	2.06	0.88
1:D:208:TYR:CE2	1:D:304:VAL:HG11	2.11	0.86
1:C:145:MET:HG3	1:C:148:ARG:HD3	1.56	0.85
1:D:110:LYS:HG3	1:D:154:ALA:HB1	1.58	0.84
1:B:77:SER:HA	1:B:80:LYS:HD2	1.59	0.84
1:C:32:GLU:HG2	1:C:34:HIS:HE1	1.42	0.84
1:A:34:HIS:CE1	1:A:213:LEU:HD21	2.13	0.83
1:C:69:PHE:HA	1:C:75:LEU:HD11	1.62	0.82
1:B:73:GLN:NE2	1:B:111:GLU:OE2	2.12	0.82
1:A:117:ARG:HH21	1:A:219:LEU:HB2	1.45	0.82
1:C:163:GLY:O	1:C:198:VAL:HG21	1.81	0.79
1:C:229:ASP:HB2	1:C:308:ARG:HD2	1.64	0.78
1:C:204:ASP:OD2	1:C:207:LYS:NZ	2.17	0.78
1:D:133:ALA:O	1:D:138:ARG:NH1	2.17	0.78
1:C:196:LYS:HD2	1:C:233:THR:CB	2.14	0.78
1:C:48:ILE:HD12	1:C:49:GLU:H	1.49	0.77
1:C:245:THR:HG23	1:C:247:LYS:H	1.48	0.76
1:C:117:ARG:NH2	1:C:213:LEU:O	2.20	0.75
1:C:268:ILE:HA	1:C:271:GLN:HG3	1.67	0.75
1:A:229:ASP:OD2	1:A:308:ARG:NH1	2.20	0.75
1:C:263:ILE:O	1:C:271:GLN:NE2	2.20	0.74
1:C:48:ILE:HD12	1:C:49:GLU:N	2.02	0.74
1:C:48:ILE:CD1	1:C:49:GLU:HG2	2.17	0.73
1:A:109:ILE:HG23	1:A:115:VAL:HG21	1.70	0.73
1:A:18:THR:HG21	1:A:50:LYS:HD3	1.71	0.73
1:A:52:GLN:HE22	1:D:24:ARG:HH22	1.35	0.72
1:A:73:GLN:HA	1:A:76:VAL:HG22	1.72	0.71
1:A:268:ILE:O	1:A:270:HIS:N	2.23	0.71
1:C:233:THR:HG23	1:C:235:LEU:H	1.55	0.70
1:C:227:PRO:HB2	1:C:230:ASN:HB2	1.73	0.70
1:B:194:ASN:O	1:B:194:ASN:ND2	2.24	0.70
1:C:129:ARG:HH12	1:C:289:HIS:CG	2.11	0.69
1:C:249:LEU:O	1:C:251:LYS:NZ	2.26	0.69
1:C:138:ARG:NH2	1:C:142:ASP:OD1	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:129:ARG:HH12	1:D:289:HIS:CG	2.11	0.68
1:A:245:THR:HG21	1:A:247:LYS:HE2	1.74	0.68
1:C:79:VAL:HG22	1:C:85:VAL:HG11	1.76	0.67
1:C:90:SER:O	1:C:98:ASN:ND2	2.28	0.67
1:C:196:LYS:HD2	1:C:233:THR:N	2.09	0.67
1:C:196:LYS:HZ1	1:C:236:GLU:CG	2.08	0.67
1:C:264:GLU:HA	1:C:271:GLN:HE22	1.60	0.67
1:A:117:ARG:NH1	1:A:213:LEU:O	2.28	0.66
1:D:11:ARG:NH2	1:D:83:ASP:OD2	2.28	0.66
1:D:139:GLU:O	1:D:143:GLN:HG3	1.94	0.66
1:B:268:ILE:O	1:B:270:HIS:N	2.20	0.66
1:C:196:LYS:HD2	1:C:233:THR:HB	1.76	0.66
1:B:52:GLN:NE2	1:C:20:TYR:OH	2.29	0.65
1:D:29:CYS:HB3	1:D:36:THR:HG21	1.76	0.65
1:C:196:LYS:HD2	1:C:233:THR:H	1.62	0.65
1:D:20:TYR:HE2	1:D:202:GLU:HG2	1.62	0.64
1:C:32:GLU:HG2	1:C:34:HIS:CE1	2.30	0.64
1:D:20:TYR:CE2	1:D:202:GLU:HG2	2.33	0.64
1:B:44:ILE:HA	1:B:47:GLU:H	1.62	0.64
1:C:127:PRO:HB2	1:C:145:MET:HE2	1.80	0.63
1:C:196:LYS:HD2	1:C:233:THR:HG22	1.79	0.63
1:A:265:GLN:HE21	1:A:271:GLN:HE22	1.46	0.63
1:A:103:LEU:O	1:A:106:VAL:HG12	1.99	0.63
1:C:196:LYS:HZ1	1:C:236:GLU:HG2	1.64	0.62
1:A:122:GLU:OE1	1:A:148:ARG:NH2	2.23	0.62
1:D:127:PRO:HB2	1:D:145:MET:HE3	1.82	0.62
1:C:139:GLU:O	1:C:143:GLN:HG3	1.99	0.62
1:D:145:MET:HE1	1:D:148:ARG:HD2	1.82	0.61
1:C:232:LEU:HD11	1:C:236:GLU:HB2	1.82	0.61
1:C:311:ASP:OD1	1:C:314:ARG:NH2	2.33	0.61
1:C:84:VAL:HG11	1:C:213:LEU:CD2	2.27	0.61
1:C:196:LYS:HD2	1:C:233:THR:CG2	2.30	0.61
1:C:200:ALA:HB3	1:C:205:ILE:HD11	1.81	0.61
1:A:20:TYR:CE2	1:A:202:GLU:HG2	2.36	0.61
1:C:85:VAL:HG23	1:C:118:PHE:CD1	2.36	0.60
1:A:20:TYR:HE2	1:A:202:GLU:HG2	1.66	0.60
1:B:234:GLN:O	1:B:238:VAL:HG23	2.01	0.60
1:A:34:HIS:HE1	1:A:213:LEU:HD21	1.64	0.60
1:C:164:ALA:HA	1:C:198:VAL:CG2	2.30	0.60
1:C:244:LEU:HD11	1:C:310:ASP:OD2	2.01	0.60
1:C:41:ARG:HG3	1:C:42:PRO:HD2	1.81	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:GLN:HG3	1:C:44:ILE:CD1	2.31	0.60
1:C:164:ALA:HA	1:C:198:VAL:HG23	1.84	0.60
1:C:188:ASN:O	1:C:234:GLN:NE2	2.34	0.60
1:C:195:VAL:O	1:C:233:THR:OG1	2.17	0.60
1:D:102:GLN:O	1:D:106:VAL:HG13	2.02	0.59
1:D:129:ARG:NH1	1:D:289:HIS:CG	2.71	0.59
1:A:168:ALA:HB1	1:D:48:ILE:HD11	1.85	0.58
1:A:11:ARG:NH2	1:A:81:LEU:O	2.32	0.58
1:D:14:VAL:HG11	1:D:26:VAL:HG23	1.83	0.58
1:B:189:ILE:HD11	1:B:251:LYS:HB3	1.86	0.58
1:A:268:ILE:HG22	1:A:269:PRO:HD2	1.85	0.58
1:C:103:LEU:O	1:C:106:VAL:HG22	2.05	0.57
1:A:52:GLN:NE2	1:D:24:ARG:HH22	2.02	0.57
1:B:133:ALA:HB3	1:B:138:ARG:HG3	1.86	0.57
1:C:37:TYR:CZ	1:C:63:ARG:HD3	2.40	0.57
1:C:260:LEU:HD11	1:C:278:TYR:CE1	2.39	0.57
1:C:13:LEU:HB3	1:C:85:VAL:HG12	1.86	0.56
1:C:40:GLN:HG3	1:C:44:ILE:HD11	1.86	0.56
1:C:13:LEU:N	1:C:82:VAL:HG21	2.20	0.56
1:C:303:ASP:OD1	1:C:303:ASP:N	2.38	0.56
1:C:13:LEU:H	1:C:82:VAL:HG21	1.70	0.56
1:C:84:VAL:CG1	1:C:213:LEU:HD21	2.30	0.56
1:B:233:THR:OG1	1:B:236:GLU:HG3	2.05	0.56
1:C:208:TYR:OH	1:C:304:VAL:HG21	2.06	0.56
1:D:235:LEU:O	1:D:239:GLN:HG3	2.05	0.56
1:B:165:CYS:HB3	1:B:170:PHE:CD1	2.40	0.56
1:C:37:TYR:CE2	1:C:63:ARG:HD3	2.41	0.55
1:C:105:LEU:O	1:C:109:ILE:HG13	2.06	0.55
1:C:196:LYS:CD	1:C:233:THR:H	2.18	0.55
1:A:205:ILE:HG23	1:A:225:ILE:CD1	2.35	0.55
1:C:194:ASN:HA	1:C:233:THR:HG21	1.88	0.55
1:B:41:ARG:O	1:B:43:GLU:HG3	2.05	0.55
1:C:100:LEU:HD21	1:C:143:GLN:HB3	1.89	0.55
1:D:138:ARG:HG2	1:D:142:ASP:OD2	2.07	0.55
1:D:263:ILE:HA	1:D:266:MET:HG3	1.89	0.54
1:B:262:ASN:O	1:B:266:MET:HB2	2.07	0.54
1:C:85:VAL:HG23	1:C:118:PHE:HD1	1.72	0.54
1:C:196:LYS:NZ	1:C:236:GLU:HG2	2.21	0.54
1:C:242:GLU:OE2	1:C:248:GLU:HA	2.07	0.54
1:C:300:LEU:HB2	1:C:301:TYR:CE2	2.43	0.54
1:C:127:PRO:HD2	1:C:148:ARG:NH1	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:GLY:O	1:C:23:LYS:HG3	2.07	0.54
1:B:20:TYR:HE2	1:B:202:GLU:HG2	1.72	0.54
1:D:103:LEU:O	1:D:106:VAL:HG22	2.08	0.54
1:A:262:ASN:HA	1:A:264:GLU:OE1	2.08	0.53
1:C:127:PRO:HD2	1:C:148:ARG:HH12	1.73	0.53
1:D:73:GLN:HA	1:D:76:VAL:HG22	1.90	0.53
1:A:219:LEU:HD23	1:A:220:ASN:OD1	2.08	0.53
1:A:290:GLU:OE1	1:A:290:GLU:N	2.41	0.53
1:C:48:ILE:HD13	1:C:49:GLU:HG2	1.88	0.53
1:A:265:GLN:NE2	1:A:271:GLN:OE1	2.42	0.53
1:A:270:HIS:HD2	1:A:270:HIS:O	1.91	0.53
1:C:227:PRO:HB3	1:C:306:TYR:CE2	2.44	0.53
1:B:41:ARG:O	1:B:43:GLU:N	2.43	0.52
1:C:196:LYS:HG3	1:C:232:LEU:HA	1.90	0.52
1:C:306:TYR:CE1	1:C:308:ARG:NH2	2.67	0.52
1:A:242:GLU:HG2	1:A:248:GLU:HA	1.91	0.52
1:A:268:ILE:C	1:A:270:HIS:H	2.13	0.52
1:C:208:TYR:CE1	1:C:304:VAL:HG21	2.44	0.52
1:A:160:TYR:HB2	1:A:222:THR:HG23	1.91	0.52
1:D:110:LYS:HE2	1:D:154:ALA:HA	1.91	0.52
1:A:192:ASP:HB2	1:A:194:ASN:ND2	2.25	0.52
1:C:129:ARG:HH22	1:C:289:HIS:CG	2.28	0.51
1:A:39:LEU:HD11	1:A:67:GLY:HA3	1.91	0.51
1:B:102:GLN:O	1:B:106:VAL:HG23	2.09	0.51
1:C:173:ASN:HB3	1:C:180:LEU:HA	1.92	0.51
1:D:48:ILE:O	1:D:52:GLN:HG3	2.11	0.51
1:C:145:MET:O	1:C:149:GLN:HG3	2.09	0.51
1:B:265:GLN:NE2	1:B:271:GLN:HE22	2.09	0.51
1:D:270:HIS:CD2	1:D:274:ILE:HG13	2.45	0.51
1:C:196:LYS:CD	1:C:233:THR:N	2.74	0.51
1:D:127:PRO:HD2	1:D:148:ARG:NH2	2.26	0.51
1:B:271:GLN:O	1:B:274:ILE:HG22	2.11	0.51
1:B:127:PRO:HD2	1:B:148:ARG:NH2	2.26	0.51
1:C:212:THR:HG23	1:C:218:THR:OG1	2.11	0.50
1:C:125:MET:SD	1:C:280:ILE:HD11	2.51	0.50
1:D:29:CYS:HB3	1:D:36:THR:CG2	2.41	0.50
1:D:212:THR:HA	1:D:301:TYR:OH	2.12	0.50
1:A:268:ILE:CG2	1:A:269:PRO:HD2	2.41	0.50
1:D:32:GLU:OE1	1:D:211:LYS:HE3	2.11	0.50
1:C:45:GLY:C	1:C:46:LEU:HD12	2.32	0.49
1:D:41:ARG:HH21	1:D:43:GLU:CD	2.15	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:ILE:C	1:C:46:LEU:H	2.15	0.49
1:C:177:MET:O	1:C:179:THR:N	2.45	0.49
1:D:69:PHE:HA	1:D:75:LEU:HD11	1.94	0.49
1:C:208:TYR:CZ	1:C:304:VAL:HG21	2.47	0.49
1:D:148:ARG:HD3	1:D:160:TYR:CE1	2.47	0.49
1:D:183:PRO:HG3	1:D:187:VAL:HG22	1.95	0.49
1:C:82:VAL:HG22	1:C:84:VAL:H	1.77	0.49
1:C:129:ARG:HH22	1:C:289:HIS:CB	2.26	0.49
1:C:157:PRO:HA	1:C:220:ASN:HD21	1.77	0.49
1:A:265:GLN:HE21	1:A:271:GLN:NE2	2.08	0.49
1:D:122:GLU:O	1:D:124:GLY:N	2.46	0.49
1:C:314:ARG:O	1:C:317:LEU:HD12	2.13	0.48
1:A:190:TYR:HB3	1:A:282:TYR:CZ	2.48	0.48
1:B:76:VAL:HG12	1:B:80:LYS:HZ2	1.78	0.48
1:A:117:ARG:NH1	1:A:213:LEU:C	2.67	0.48
1:B:20:TYR:CE2	1:B:202:GLU:HG2	2.48	0.48
1:B:44:ILE:CB	1:B:47:GLU:H	2.26	0.48
1:C:144:LYS:O	1:C:148:ARG:HG3	2.13	0.48
1:C:79:VAL:HG11	1:C:109:ILE:HA	1.96	0.47
1:C:133:ALA:HB2	1:C:279:HIS:CE1	2.50	0.47
1:B:73:GLN:HA	1:B:76:VAL:HG23	1.95	0.47
1:C:238:VAL:O	1:C:242:GLU:HG2	2.14	0.47
1:A:202:GLU:CD	1:A:202:GLU:H	2.18	0.47
1:A:20:TYR:O	1:A:24:ARG:HD2	2.14	0.47
1:A:133:ALA:O	1:A:138:ARG:NH2	2.47	0.47
1:C:227:PRO:HB3	1:C:306:TYR:CZ	2.50	0.47
1:A:12:VAL:HG21	1:A:213:LEU:HD11	1.96	0.47
1:A:20:TYR:HE1	1:D:52:GLN:NE2	2.13	0.47
1:A:48:ILE:HD11	1:D:168:ALA:HB1	1.96	0.47
1:D:13:LEU:HA	1:D:37:TYR:O	2.15	0.47
1:C:20:TYR:HE2	1:C:202:GLU:HG2	1.80	0.47
1:C:122:GLU:OE2	1:C:148:ARG:HD2	2.15	0.47
1:C:193:GLY:O	1:C:233:THR:HG21	2.15	0.47
1:D:37:TYR:CD2	1:D:63:ARG:HB2	2.50	0.47
1:C:40:GLN:HB2	1:C:54:PHE:CE2	2.50	0.47
1:C:240:ILE:O	1:C:244:LEU:HD13	2.15	0.47
1:C:10:THR:HB	1:C:34:HIS:HD2	1.80	0.46
1:D:145:MET:CE	1:D:148:ARG:HD2	2.45	0.46
1:C:43:GLU:O	1:C:46:LEU:HD13	2.16	0.46
1:C:287:THR:O	1:C:289:HIS:N	2.40	0.46
1:B:196:LYS:HG2	1:B:233:THR:HG22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:HIS:O	1:B:76:VAL:HG23	2.16	0.46
1:C:196:LYS:O	1:C:287:THR:OG1	2.27	0.46
1:D:270:HIS:HD2	1:D:274:ILE:HG13	1.81	0.46
1:B:266:MET:HE1	1:D:270:HIS:NE2	2.31	0.46
1:C:40:GLN:HB2	1:C:54:PHE:HE2	1.80	0.46
1:B:266:MET:HE2	1:B:266:MET:HB3	1.90	0.45
1:C:301:TYR:O	1:C:304:VAL:HG22	2.15	0.45
1:C:296:GLU:O	1:C:300:LEU:HD12	2.16	0.45
1:D:44:ILE:HD12	1:D:44:ILE:H	1.82	0.45
1:D:294:ASP:OD2	1:D:294:ASP:N	2.44	0.45
1:C:196:LYS:HG3	1:C:232:LEU:CA	2.47	0.45
1:A:41:ARG:HA	1:A:67:GLY:O	2.17	0.45
1:A:71:ASP:OD2	1:A:74:SER:HB3	2.17	0.45
1:A:127:PRO:N	1:A:128:PRO:HD2	2.32	0.45
1:A:310:ASP:O	1:A:314:ARG:HG3	2.17	0.45
1:D:190:TYR:HB3	1:D:282:TYR:CE1	2.52	0.45
1:B:44:ILE:C	1:B:46:LEU:H	2.19	0.45
1:A:10:THR:HG21	1:A:214:ASN:HB3	1.99	0.45
1:D:266:MET:HB2	1:D:271:GLN:HG2	1.99	0.45
1:B:48:ILE:HG23	1:B:49:GLU:N	2.32	0.45
1:A:270:HIS:O	1:A:270:HIS:CD2	2.70	0.45
1:C:198:VAL:HG22	1:C:286:LEU:HD13	1.99	0.44
1:C:211:LYS:HB2	1:C:211:LYS:HE2	1.73	0.44
1:D:89:MET:N	1:D:121:SER:OG	2.44	0.44
1:C:116:LYS:O	1:C:157:PRO:HG2	2.17	0.44
1:D:132:HIS:N	1:D:138:ARG:HH22	2.16	0.44
1:C:196:LYS:HG3	1:C:232:LEU:C	2.38	0.44
1:C:314:ARG:HA	1:C:317:LEU:CD1	2.47	0.44
1:D:268:ILE:HB	1:D:269:PRO:HD3	1.99	0.44
1:A:142:ASP:OD2	1:A:142:ASP:N	2.50	0.44
1:D:264:GLU:HG3	1:D:265:GLN:OE1	2.18	0.44
1:C:69:PHE:HZ	1:C:89:MET:HE1	1.82	0.44
1:D:39:LEU:HD11	1:D:67:GLY:HA3	1.98	0.44
1:D:301:TYR:O	1:D:304:VAL:HG12	2.17	0.44
1:B:189:ILE:CD1	1:B:251:LYS:HB3	2.47	0.44
1:C:212:THR:HG1	1:C:301:TYR:HH	1.55	0.44
1:D:98:ASN:N	1:D:98:ASN:OD1	2.51	0.44
1:C:287:THR:C	1:C:289:HIS:H	2.20	0.44
1:B:174:LEU:HD12	1:B:183:PRO:HD2	2.00	0.44
1:D:79:VAL:HG22	1:D:85:VAL:HG21	1.99	0.44
1:B:44:ILE:C	1:B:46:LEU:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:ILE:CA	1:B:47:GLU:H	2.28	0.43
1:C:196:LYS:HD2	1:C:233:THR:CA	2.49	0.43
1:A:127:PRO:HD2	1:A:148:ARG:NH2	2.33	0.43
1:D:131:GLY:C	1:D:138:ARG:HH22	2.20	0.43
1:D:187:VAL:HG21	1:D:249:LEU:HD13	1.99	0.43
1:D:197:VAL:HG21	1:D:281:PHE:CE2	2.53	0.43
1:C:48:ILE:HD12	1:C:49:GLU:HG2	1.97	0.43
1:C:198:VAL:HG12	1:C:231:VAL:HG22	2.01	0.43
1:C:202:GLU:HA	1:C:205:ILE:HD12	2.01	0.43
1:A:264:GLU:HG2	1:A:265:GLN:H	1.83	0.43
1:B:313:LEU:HD23	1:B:313:LEU:HA	1.81	0.43
1:D:297:ALA:O	1:D:301:TYR:HD1	2.02	0.43
1:A:73:GLN:HA	1:A:76:VAL:CG2	2.46	0.43
1:D:266:MET:O	1:D:267:GLU:HB2	2.18	0.43
1:C:118:PHE:C	1:C:119:LEU:HD12	2.40	0.43
1:C:234:GLN:HB2	1:C:281:PHE:HE1	1.83	0.43
1:D:304:VAL:O	1:D:304:VAL:HG13	2.18	0.43
1:A:79:VAL:HG11	1:A:109:ILE:HA	2.02	0.42
1:D:114:ASN:O	1:D:114:ASN:ND2	2.45	0.42
1:B:190:TYR:HB3	1:B:282:TYR:CE2	2.54	0.42
1:D:23:LYS:HD3	1:D:53:LEU:HD22	2.01	0.42
1:D:233:THR:HG23	1:D:236:GLU:OE1	2.19	0.42
1:A:125:MET:O	1:A:127:PRO:HD3	2.20	0.42
1:D:127:PRO:HB2	1:D:145:MET:CE	2.48	0.42
1:B:11:ARG:HA	1:B:35:GLU:HB2	2.01	0.42
1:B:263:ILE:HG13	1:B:271:GLN:HG2	2.01	0.42
1:C:98:ASN:OD1	1:C:98:ASN:N	2.52	0.42
1:C:129:ARG:NH2	1:C:289:HIS:ND1	2.67	0.42
1:B:217:ARG:NH1	1:B:300:LEU:HD22	2.35	0.42
1:C:89:MET:HB3	1:C:98:ASN:HB2	2.01	0.42
1:A:52:GLN:HE22	1:D:24:ARG:NH2	2.10	0.42
1:C:73:GLN:O	1:C:76:VAL:HB	2.19	0.42
1:C:234:GLN:O	1:C:238:VAL:HG23	2.20	0.42
1:C:232:LEU:HA	1:C:232:LEU:HD12	1.62	0.42
1:D:112:ALA:C	1:D:114:ASN:H	2.23	0.42
1:D:187:VAL:CG2	1:D:249:LEU:HD13	2.50	0.42
1:D:168:ALA:O	1:D:180:LEU:HD12	2.20	0.41
1:B:90:SER:O	1:B:98:ASN:ND2	2.53	0.41
1:C:196:LYS:HA	1:C:233:THR:HA	2.02	0.41
1:C:79:VAL:HG13	1:C:115:VAL:HG21	2.02	0.41
1:C:30:LEU:HD11	1:C:62:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ARG:HG2	1:A:203:ASP:OD1	2.19	0.41
1:D:73:GLN:CA	1:D:76:VAL:HG22	2.51	0.41
1:A:45:GLY:HA2	1:A:51:VAL:CG2	2.50	0.41
1:A:266:MET:O	1:A:266:MET:HG2	2.20	0.41
1:C:197:VAL:HG11	1:C:280:ILE:HG22	2.01	0.41
1:A:168:ALA:CB	1:D:48:ILE:HD11	2.50	0.41
1:A:288:ASP:OD1	1:A:289:HIS:ND1	2.33	0.41
1:C:196:LYS:HE3	1:C:232:LEU:HD12	2.03	0.41
1:D:13:LEU:HD12	1:D:37:TYR:HB2	2.03	0.40
1:D:300:LEU:HB2	1:D:301:TYR:CE1	2.56	0.40
1:B:47:GLU:OE2	1:B:50:LYS:NZ	2.46	0.40
1:C:304:VAL:O	1:C:304:VAL:HG23	2.21	0.40
1:A:306:TYR:CZ	1:A:308:ARG:NH1	2.89	0.40
1:B:180:LEU:O	1:B:181:LEU:HD23	2.22	0.40
1:C:41:ARG:HA	1:C:67:GLY:O	2.21	0.40
1:A:80:LYS:HE2	1:A:80:LYS:HB2	1.81	0.40
1:B:109:ILE:HG23	1:B:115:VAL:HG21	2.03	0.40
1:D:130:MET:HG3	1:D:141:PHE:CE1	2.57	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	299/317 (94%)	279 (93%)	16 (5%)	4 (1%)	12 36
1	B	300/317 (95%)	286 (95%)	9 (3%)	5 (2%)	9 29
1	C	300/317 (95%)	283 (94%)	11 (4%)	6 (2%)	7 24
1	D	300/317 (95%)	283 (94%)	14 (5%)	3 (1%)	15 44
All	All	1199/1268 (95%)	1131 (94%)	50 (4%)	18 (2%)	10 33

All (18) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	B	44	ILE
1	B	178	VAL
1	C	178	VAL
1	C	288	ASP
1	A	178	VAL
1	D	267	GLU
1	B	264	GLU
1	B	269	PRO
1	C	47	GLU
1	A	174	LEU
1	A	269	PRO
1	B	42	PRO
1	C	177	MET
1	C	268	ILE
1	A	137	GLY
1	D	113	GLY
1	C	269	PRO
1	D	268	ILE

### 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	256/270 (95%)	254 (99%)	2 (1%)	81	94
1	B	257/270 (95%)	253 (98%)	4 (2%)	62	88
1	C	259/270 (96%)	250 (96%)	9 (4%)	36	70
1	D	259/270 (96%)	252 (97%)	7 (3%)	44	78
All	All	1031/1080 (96%)	1009 (98%)	22 (2%)	53	84

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	47	GLU
1	B	194	ASN
1	B	207	LYS
1	B	266	MET

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Mol	Chain	Res	Type
1	C	24	ARG
1	C	46	LEU
1	C	52	GLN
1	C	55	LEU
1	C	107	GLU
1	C	129	ARG
1	C	138	ARG
1	C	145	MET
1	C	299	SER
1	A	219	LEU
1	A	220	ASN
1	D	9	LYS
1	D	74	SER
1	D	114	ASN
1	D	116	LYS
1	D	121	SER
1	D	212	THR
1	D	264	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	52	GLN
1	B	194	ASN
1	B	265	GLN
1	C	34	HIS
1	C	271	GLN
1	A	34	HIS
1	A	52	GLN
1	A	143	GLN
1	A	265	GLN
1	A	270	HIS
1	A	271	GLN
1	D	52	GLN

### 5.3.3 RNA

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 monosaccharides 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	303/317 (95%)	-0.11	3 (0%) 82 77	42, 60, 86, 110	0
1	B	304/317 (95%)	0.03	4 (1%) 77 72	30, 54, 81, 122	0
1	C	304/317 (95%)	0.36	13 (4%) 35 25	53, 82, 106, 143	0
1	D	304/317 (95%)	-0.08	1 (0%) 94 93	38, 60, 83, 136	0
All	All	1215/1268 (95%)	0.05	21 (1%) 70 63	30, 64, 96, 143	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	222	THR	3.9
1	B	45	GLY	3.7
1	B	265	GLN	3.6
1	C	221	LYS	3.6
1	D	96	SER	3.4
1	C	135	PRO	3.4
1	B	96	SER	3.1
1	B	264	GLU	3.0
1	C	276	HIS	2.7
1	C	97	HIS	2.7
1	C	161	VAL	2.6
1	C	160	TYR	2.5
1	C	158	TYR	2.4
1	A	76	VAL	2.3
1	A	264	GLU	2.3
1	C	118	PHE	2.1
1	A	134	LEU	2.1
1	C	223	VAL	2.1
1	C	209	THR	2.0
1	C	125	MET	2.0
1	C	286	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 monosaccharides 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.