



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 20, 2023 – 12:12 PM JST

PDB ID : 7CJY  
Title : Drimenol synthase from *Persicaria hydropiper*  
Authors : Xiao, W.; Zhou, H.; Zhou, M.T.; Liu, L.; Xiang, Z.  
Deposited on : 2020-07-14  
Resolution : 2.20 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

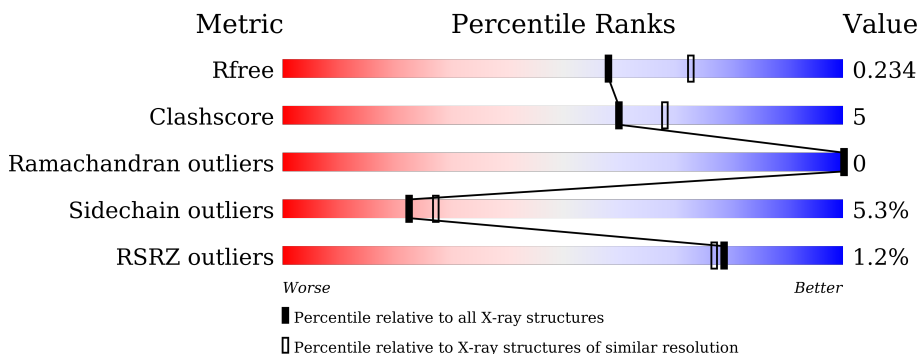
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	542	 2% 85% 14% .
1	B	542	 % 85% 13% .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8908 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 (-)-drimenol synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	540	4424	2847	746	806	25	0	2	0
1	B	540	4427	2848	750	804	25	0	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	GLY	-	expression tag	UNP W0FFD7
B	18	GLY	-	expression tag	UNP W0FFD7

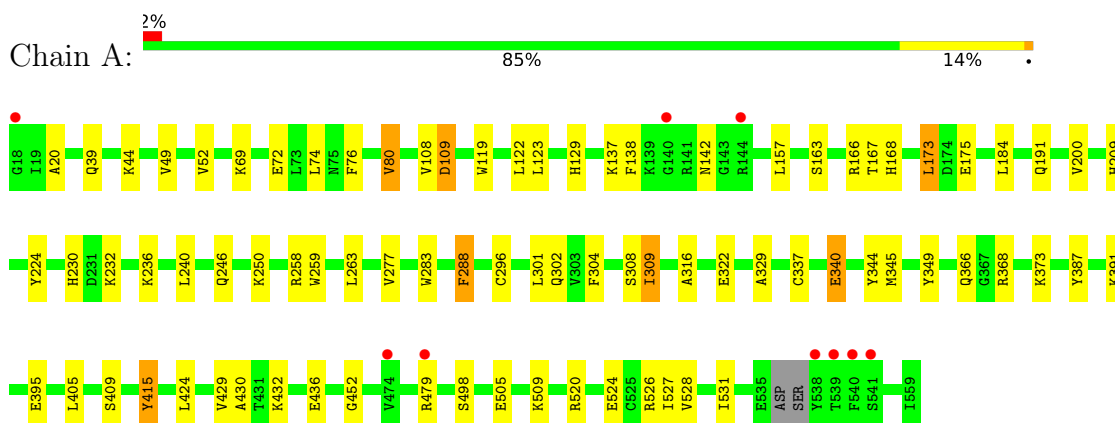
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	26	Total	O	0	0
			26	26		
2	B	31	Total	O	0	0
			31	31		

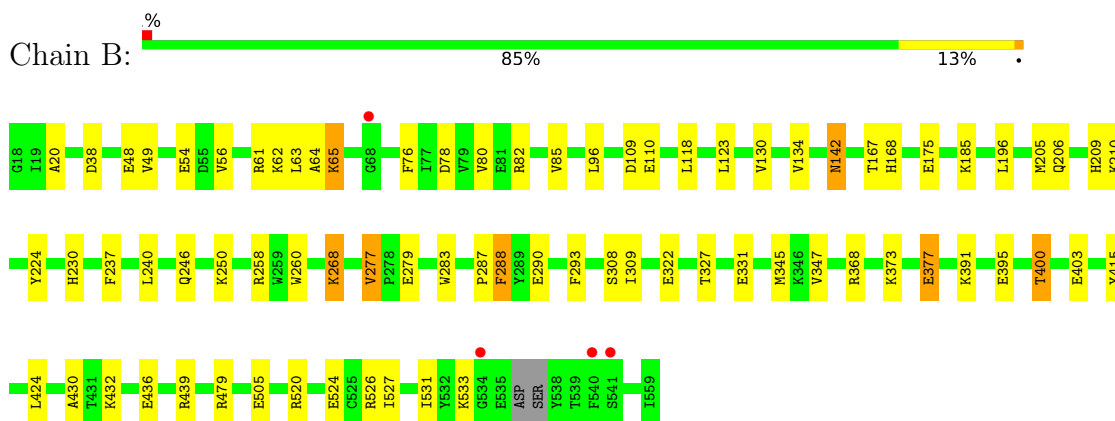
### 3 Residue-property plots [i](#)

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: (-)-drimenol synthase



- Molecule 1: (-)-drimenol synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	232.53Å 60.09Å 106.22Å 90.00° 111.88° 90.00°	Depositor
Resolution (Å)	28.74 – 2.20 98.57 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.3 (28.74-2.20) 99.1 (98.57-2.15)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.53 (at 2.14Å)	Xtrriage
Refinement program	PHENIX 1.17.1	Depositor
R, $R_{free}$	0.195 , 0.234 0.195 , 0.234	Depositor DCC
$R_{free}$ test set	3690 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.7	Xtrriage
Anisotropy	0.414	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 53.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8908	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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	0.42	0/4549	0.56	0/6149
1	B	0.43	0/4557	0.58	0/6159
All	All	0.42	0/9106	0.57	0/12308

There are no bond length outliers.

There are no bond angle outliers.

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	4424	0	4327	45	0
1	B	4427	0	4337	40	0
2	A	26	0	0	1	0
2	B	31	0	0	1	0
All	All	8908	0	8664	81	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 (81) 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:167:THR:HG22	1:A:168:HIS:H	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ALA:HB1	1:B:308:SER:HB2	1.73	0.70
1:A:49:VAL:HG22	1:A:240:LEU:HB3	1.77	0.67
1:B:400:THR:HG22	1:B:403:GLU:H	1.60	0.66
1:B:167:THR:HG22	1:B:168:HIS:H	1.61	0.65
1:B:49:VAL:HG22	1:B:240:LEU:HB3	1.79	0.64
1:A:308:SER:HB2	1:B:20:ALA:HB1	1.81	0.62
1:A:76:PHE:O	1:A:80:VAL:HG13	2.00	0.60
1:B:142:ASN:OD1	1:B:142:ASN:N	2.37	0.57
1:A:69:LYS:HB3	1:A:72:GLU:HG3	1.87	0.56
1:B:283[A]:TRP:CE3	1:B:531:ILE:HD13	2.41	0.55
1:B:61:ARG:O	1:B:65:LYS:HG2	2.07	0.54
1:B:368:ARG:NH1	1:B:430:ALA:O	2.40	0.54
1:A:167:THR:HG22	1:A:168:HIS:N	2.22	0.54
1:A:69:LYS:HB3	1:A:72:GLU:CG	2.38	0.53
1:A:167:THR:CG2	1:A:505:GLU:HB2	2.39	0.53
1:B:78:ASP:OD2	1:B:82:ARG:NH2	2.39	0.53
1:B:123:LEU:HD23	1:B:130:VAL:HG11	1.91	0.52
1:A:109:ASP:OD1	1:A:109:ASP:N	2.41	0.52
1:B:80:VAL:HG23	1:B:85:VAL:HG23	1.91	0.51
1:A:283[B]:TRP:CE3	1:A:531:ILE:HD13	2.45	0.51
1:B:210:LYS:NZ	2:B:602:HOH:O	2.44	0.50
1:B:527:ILE:HG13	1:B:531:ILE:HD12	1.94	0.49
1:B:224:TYR:CE2	1:B:230:HIS:HB2	2.48	0.49
1:A:387:TYR:CE2	1:A:391:LYS:HE2	2.48	0.49
1:A:108:VAL:HG21	1:A:119:TRP:CE2	2.47	0.48
1:A:329:ALA:O	1:A:349:TYR:OH	2.26	0.48
1:B:283[B]:TRP:HZ3	1:B:415:TYR:HH	1.58	0.48
1:B:209:HIS:O	1:B:526:ARG:HD3	2.14	0.47
1:B:377[A]:GLU:HG2	1:B:439:ARG:NH2	2.30	0.47
1:B:76:PHE:O	1:B:80:VAL:HG12	2.15	0.47
1:A:283[A]:TRP:HZ3	1:A:415:TYR:HH	1.57	0.47
1:B:432:LYS:O	1:B:436:GLU:HG3	2.16	0.46
1:B:287:PRO:HB2	1:B:288:PHE:CE1	2.51	0.46
1:A:209:HIS:O	1:A:526:ARG:HD3	2.16	0.46
1:A:288:PHE:CZ	1:A:296:CYS:HB3	2.51	0.46
1:B:290:GLU:HG2	1:B:293:PHE:CD2	2.51	0.46
1:A:191:GLN:HB3	2:A:606:HOH:O	2.16	0.45
1:B:246:GLN:O	1:B:250:LYS:HG3	2.17	0.45
1:B:327:THR:O	1:B:331:GLU:HG3	2.17	0.45
1:A:344:TYR:CE1	1:B:268:LYS:HG2	2.51	0.45
1:A:432:LYS:O	1:A:436:GLU:HG3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:ALA:HB1	1:B:268:LYS:O	2.17	0.44
1:A:337:CYS:HA	1:A:340:GLU:HG3	1.98	0.44
1:B:167:THR:CG2	1:B:505:GLU:HB2	2.47	0.44
1:A:137:LYS:HE3	1:A:138:PHE:CZ	2.52	0.44
1:A:232:LYS:HD3	1:A:232:LYS:HA	1.87	0.43
1:A:246:GLN:O	1:A:250:LYS:HG3	2.17	0.43
1:B:391:LYS:HE3	1:B:395:GLU:OE2	2.17	0.43
1:B:322:GLU:HB3	1:B:345:MET:HE3	2.00	0.43
1:A:259:TRP:CH2	1:A:302:GLN:HG2	2.53	0.43
1:A:527:ILE:HG13	1:A:531:ILE:HD12	1.99	0.43
1:B:64:ALA:HB2	1:B:96:LEU:HD21	2.00	0.43
1:A:163:SER:O	1:A:166:ARG:HD3	2.18	0.43
1:B:288:PHE:CD1	1:B:288:PHE:N	2.86	0.43
1:A:405:LEU:HA	1:A:409:SER:OG	2.19	0.43
1:A:301:LEU:O	1:A:304:PHE:HB3	2.18	0.42
1:A:415:TYR:HD1	1:A:415:TYR:HA	1.68	0.42
1:A:184:LEU:HD22	1:A:200:VAL:HG13	2.01	0.42
1:B:415:TYR:CZ	1:B:524:GLU:HB3	2.55	0.42
1:B:205:MET:HG2	1:B:533:LYS:HD3	2.01	0.42
1:A:309:ILE:HD12	1:A:309:ILE:HA	1.84	0.42
1:A:74:LEU:HD23	1:A:74:LEU:HA	1.87	0.42
1:B:56:VAL:HG21	1:B:237:PHE:CD2	2.55	0.42
1:B:78:ASP:HB2	1:B:118:LEU:HD11	2.01	0.42
1:A:52:VAL:HG11	1:A:236:LYS:HE3	2.02	0.42
1:A:167:THR:HG21	1:A:505:GLU:HB2	2.02	0.42
1:A:322:GLU:HB3	1:A:345:MET:HE3	2.02	0.42
1:A:391:LYS:O	1:A:395:GLU:HG2	2.20	0.41
1:A:415:TYR:CZ	1:A:524:GLU:HB3	2.55	0.41
1:A:368:ARG:NH2	1:A:430:ALA:O	2.47	0.41
1:A:122:LEU:HD23	1:A:122:LEU:HA	1.87	0.41
1:A:288:PHE:CE2	1:A:296:CYS:HB3	2.55	0.41
1:B:415:TYR:HD1	1:B:415:TYR:HA	1.62	0.41
1:A:224:TYR:CZ	1:A:230:HIS:HB2	2.56	0.41
1:A:163:SER:HA	1:A:173:LEU:HG	2.02	0.41
1:B:167:THR:HG22	1:B:168:HIS:N	2.34	0.41
1:B:206:GLN:OE1	1:B:533:LYS:HE3	2.21	0.41
1:B:373:LYS:HE2	1:B:373:LYS:HB3	1.64	0.41
1:B:260:TRP:CD1	1:B:277:VAL:HG22	2.56	0.40
1:A:452:GLY:HA2	1:A:528:VAL:HG11	2.03	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	538/542 (99%)	524 (97%)	14 (3%)	0	100	100
1	B	540/542 (100%)	529 (98%)	11 (2%)	0	100	100
All	All	1078/1084 (99%)	1053 (98%)	25 (2%)	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	478/481 (99%)	452 (95%)	26 (5%)	22	26
1	B	478/481 (99%)	451 (94%)	27 (6%)	21	25
All	All	956/962 (99%)	903 (94%)	53 (6%)	22	26

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	44	LYS
1	A	80	VAL
1	A	109	ASP
1	A	123	LEU
1	A	129	HIS
1	A	142	ASN
1	A	157	LEU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	173	LEU
1	A	175[A]	GLU
1	A	175[B]	GLU
1	A	258	ARG
1	A	263	LEU
1	A	277	VAL
1	A	288	PHE
1	A	309	ILE
1	A	340	GLU
1	A	366	GLN
1	A	373	LYS
1	A	415	TYR
1	A	424	LEU
1	A	429	VAL
1	A	479	ARG
1	A	498	SER
1	A	509	LYS
1	A	520	ARG
1	B	38	ASP
1	B	48	GLU
1	B	54	GLU
1	B	62	LYS
1	B	63	LEU
1	B	65	LYS
1	B	109	ASP
1	B	110	GLU
1	B	134	VAL
1	B	142	ASN
1	B	175	GLU
1	B	185	LYS
1	B	196	LEU
1	B	258[A]	ARG
1	B	258[B]	ARG
1	B	268	LYS
1	B	277	VAL
1	B	279	GLU
1	B	288	PHE
1	B	309	ILE
1	B	347	VAL
1	B	377[A]	GLU
1	B	377[B]	GLU
1	B	400	THR

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Mol	Chain	Res	Type
1	B	424	LEU
1	B	479	ARG
1	B	520	ARG

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

Mol	Chain	Res	Type
1	A	98	GLN
1	A	292	GLN
1	A	381	GLN
1	B	219	HIS
1	B	551	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	540/542 (99%)	0.17	9 (1%) 70 68	48, 69, 96, 142	0
1	B	540/542 (99%)	0.14	4 (0%) 87 86	47, 68, 100, 139	0
All	All	1080/1084 (99%)	0.16	13 (1%) 79 77	47, 68, 99, 142	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	540	PHE	5.4
1	A	540	PHE	3.5
1	B	541	SER	3.2
1	A	18	GLY	2.9
1	A	539	THR	2.8
1	A	541	SER	2.7
1	A	538	TYR	2.4
1	B	534	GLY	2.4
1	B	68	GLY	2.2
1	A	474	VAL	2.1
1	A	144	ARG	2.1
1	A	140	GLY	2.0
1	A	479	ARG	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.