



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 07:27 AM EDT

PDB ID : 6NTE  
Title : Crystal Structure of Synechocystis Dihydroxyacid Dehydratase (DHAD)  
Authors : MacTavish, B.; Bruner, S.  
Deposited on : 2019-01-29  
Resolution : 2.33 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 : 2.36.2  
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) : 2.36.2

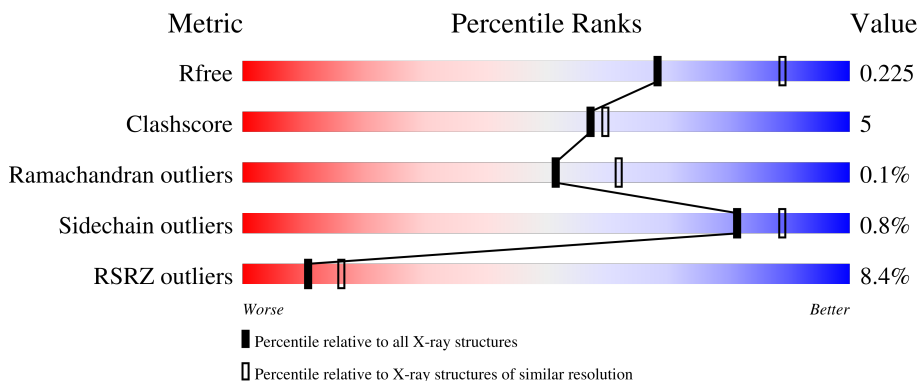
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.33 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	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	561	 9% 77% 10% 13%
1	B	561	 6% 77% 9% 14%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7279 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 Dihydroxy-acid dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	486	3515	2217	600	676	22	0	0	0
1	B	485	3508	2212	599	675	22	0	0	0

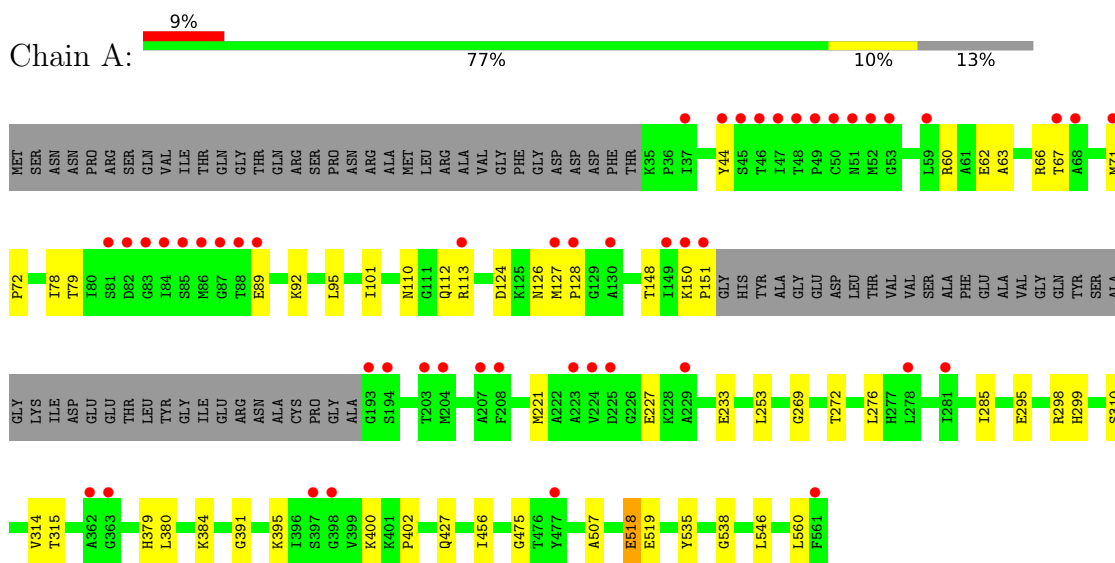
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	124	124	124	0	0
2	B	132	132	132	0	0

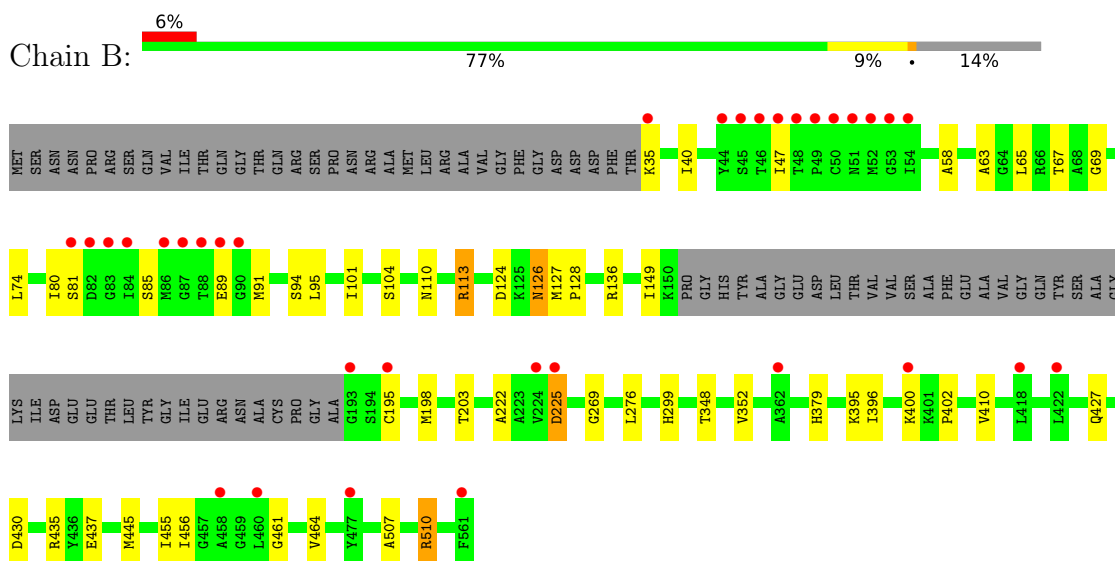
### 3 Residue-property plots [i](#)

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: Dihydroxy-acid dehydratase



- Molecule 1: Dihydroxy-acid dehydratase



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.79Å 74.26Å 80.68Å 90.00° 90.15° 90.00°	Depositor
Resolution (Å)	37.13 – 2.33 37.13 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.3 (37.13-2.33) 99.3 (37.13-2.33)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.79 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.172 , 0.225 0.172 , 0.225	Depositor DCC
$R_{free}$ test set	2153 reflections (5.35%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.6	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 54.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.021 for h,-k,-l 0.015 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7279	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% 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	0.45	0/3569	0.64	3/4843 (0.1%)
1	B	0.44	0/3561	0.59	1/4831 (0.0%)
All	All	0.44	0/7130	0.62	4/9674 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	113	ARG	NE-CZ-NH2	-11.92	114.34	120.30
1	A	427	GLN	CA-CB-CG	6.61	127.95	113.40
1	A	113	ARG	CB-CG-CD	5.84	126.78	111.60
1	B	400	LYS	CB-CG-CD	-5.53	97.22	111.60

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	3515	0	3559	38	0
1	B	3508	0	3552	34	0
2	A	124	0	0	1	0
2	B	132	0	0	0	0
All	All	7279	0	7111	69	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 (69) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:MET:HG3	1:B:128:PRO:HD3	1.64	0.80
1:B:198:MET:HE3	1:B:203:THR:HA	1.66	0.78
1:A:518:GLU:OE1	2:A:601:HOH:O	2.02	0.77
1:A:60:ARG:NH1	1:A:233:GLU:OE1	2.21	0.74
1:B:222:ALA:HB3	1:B:225:ASP:OD1	1.92	0.70
1:A:519:GLU:O	1:A:519:GLU:HG3	1.94	0.68
1:A:66:ARG:HH22	1:A:71:MET:HE1	1.58	0.68
1:A:44:TYR:HA	1:A:79:THR:HG21	1.79	0.64
1:A:66:ARG:HH22	1:A:71:MET:CE	2.11	0.63
1:A:379:HIS:HD2	1:A:395:LYS:HE3	1.63	0.62
1:A:44:TYR:HA	1:A:79:THR:CG2	2.32	0.60
1:A:150:LYS:HG2	1:A:151:PRO:HD2	1.82	0.60
1:A:295:GLU:OE2	1:A:299:HIS:NE2	2.37	0.58
1:B:101:ILE:HD11	1:B:126:ASN:ND2	2.18	0.58
1:A:402:PRO:HB2	1:A:507:ALA:HB3	1.87	0.56
1:B:101:ILE:CD1	1:B:126:ASN:ND2	2.70	0.55
1:B:101:ILE:CD1	1:B:126:ASN:HD22	2.20	0.55
1:B:80:ILE:HD11	1:B:94:SER:HA	1.90	0.54
1:A:66:ARG:NH2	1:A:71:MET:CE	2.71	0.53
1:B:198:MET:HE3	1:B:203:THR:CA	2.37	0.53
1:B:455:ILE:HD12	1:B:464:VAL:HG21	1.91	0.53
1:B:35:LYS:HD3	1:B:69:GLY:O	2.10	0.52
1:A:89:GLU:OE1	1:A:535:TYR:OH	2.30	0.49
1:B:379:HIS:HD2	1:B:395:LYS:HE3	1.76	0.49
1:B:427:GLN:HG2	1:B:430:ASP:OD2	2.11	0.49
1:B:299:HIS:CE1	1:B:510:ARG:NH1	2.80	0.49
1:A:127:MET:N	1:A:128:PRO:HD2	2.28	0.49
1:A:269:GLY:HA2	1:A:379:HIS:CE1	2.48	0.49
1:A:538:GLY:HA2	1:B:136:ARG:HD3	1.94	0.49
1:B:63:ALA:O	1:B:67:THR:HG23	2.13	0.49
1:A:148:THR:HG21	1:A:315:THR:HG21	1.95	0.48
1:A:95:LEU:HD11	1:A:276:LEU:HD21	1.96	0.48
1:A:78:ILE:HG21	1:B:104:SER:HB3	1.97	0.47
1:B:456:ILE:HD13	1:B:461:GLY:HA3	1.96	0.47
1:B:269:GLY:HA2	1:B:379:HIS:CE1	2.50	0.47
1:B:124:ASP:O	1:B:127:MET:HG2	2.15	0.47
1:B:95:LEU:HD11	1:B:276:LEU:HD21	1.96	0.46
1:A:253:LEU:HB2	1:A:285:ILE:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:MET:SD	1:A:227:GLU:HG3	2.56	0.46
1:B:402:PRO:HB2	1:B:507:ALA:HB3	1.98	0.46
1:A:151:PRO:HG3	1:A:314:VAL:HG21	1.98	0.46
1:A:101:ILE:HD12	1:A:126:ASN:OD1	2.16	0.45
1:A:63:ALA:O	1:A:67:THR:HG23	2.17	0.45
1:A:272:THR:HG22	1:A:380:LEU:HD22	1.99	0.45
1:B:40:ILE:HD11	1:B:65:LEU:HD12	1.98	0.45
1:B:110:ASN:O	1:B:113:ARG:NH1	2.50	0.45
1:B:89:GLU:O	1:B:89:GLU:HG2	2.18	0.44
1:B:198:MET:CE	1:B:203:THR:HA	2.43	0.44
1:B:348:THR:O	1:B:352:VAL:HG13	2.17	0.44
1:A:124:ASP:O	1:A:127:MET:HG2	2.18	0.43
1:A:456:ILE:HD11	1:A:475:GLY:O	2.19	0.42
1:A:298:ARG:HD3	1:A:380:LEU:O	2.20	0.42
1:A:89:GLU:OE2	1:A:92:LYS:HD2	2.20	0.42
1:B:85:SER:O	1:B:91:MET:HG2	2.20	0.41
1:B:149:ILE:HG23	1:B:195:CYS:HB2	2.01	0.41
1:B:437:GLU:O	1:B:445:MET:HA	2.20	0.41
1:A:384:LYS:HG2	1:A:391:GLY:HA2	2.01	0.41
1:A:95:LEU:HD23	1:A:95:LEU:HA	1.87	0.41
1:A:150:LYS:HG2	1:A:151:PRO:CD	2.49	0.41
1:B:95:LEU:HD23	1:B:95:LEU:HA	1.82	0.41
1:B:410:VAL:HG11	1:B:435:ARG:CZ	2.50	0.41
1:A:110:ASN:HD22	1:A:110:ASN:HA	1.71	0.41
1:A:62:GLU:HG3	1:A:72:PRO:CG	2.52	0.40
1:A:400:LYS:HA	1:A:400:LYS:HD3	1.79	0.40
1:A:546:LEU:HD22	1:A:560:LEU:HD22	2.03	0.40
1:B:396:ILE:H	1:B:396:ILE:HG13	1.73	0.40
1:A:112:GLN:OE1	1:B:81:SER:N	2.45	0.40
1:A:150:LYS:NZ	1:A:151:PRO:HG2	2.35	0.40
1:B:58:ALA:HB1	1:B:74:LEU:HD21	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	482/561 (86%)	467 (97%)	15 (3%)	0	100	100
1	B	481/561 (86%)	466 (97%)	14 (3%)	1 (0%)	47	55
All	All	963/1122 (86%)	933 (97%)	29 (3%)	1 (0%)	51	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	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	366/435 (84%)	364 (100%)	2 (0%)	88	93
1	B	365/435 (84%)	361 (99%)	4 (1%)	73	83
All	All	731/870 (84%)	725 (99%)	6 (1%)	81	89

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	310	SER
1	A	518	GLU
1	B	113	ARG
1	B	126	ASN
1	B	225	ASP
1	B	510	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	527	GLN

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Mol	Chain	Res	Type
1	B	55	ASN
1	B	126	ASN
1	B	376	GLN
1	B	427	GLN

### 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

### 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	486/561 (86%)	0.39	49 (10%) <b>7</b> <b>11</b>	21, 33, 69, 117	0
1	B	485/561 (86%)	0.20	33 (6%) <b>17</b> <b>25</b>	21, 32, 61, 117	0
All	All	971/1122 (86%)	0.30	82 (8%) <b>11</b> <b>16</b>	21, 32, 68, 117	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	PRO	11.5
1	B	49	PRO	9.7
1	B	83	GLY	9.5
1	A	52	MET	9.4
1	B	50	CYS	9.2
1	A	151	PRO	8.2
1	A	50	CYS	7.9
1	B	52	MET	7.6
1	B	82	ASP	6.7
1	A	48	THR	6.6
1	A	47	ILE	6.1
1	B	48	THR	5.7
1	A	53	GLY	5.4
1	A	86	MET	5.3
1	A	82	ASP	5.2
1	A	51	ASN	4.8
1	A	224	VAL	4.8
1	A	88	THR	4.7
1	A	45	SER	4.6
1	B	81	SER	4.4
1	B	51	ASN	4.4
1	A	150	LYS	4.3
1	B	84	ILE	4.3
1	A	87	GLY	4.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	85	SER	4.0
1	A	149	ILE	3.9
1	B	86	MET	3.9
1	B	44	TYR	3.9
1	B	46	THR	3.8
1	A	194	SER	3.6
1	A	84	ILE	3.6
1	A	561	PHE	3.5
1	A	363	GLY	3.5
1	A	398	GLY	3.4
1	A	362	ALA	3.4
1	B	224	VAL	3.3
1	A	89	GLU	3.3
1	A	46	THR	3.3
1	B	45	SER	3.2
1	A	193	GLY	3.2
1	B	47	ILE	3.2
1	B	193	GLY	3.0
1	A	225	ASP	3.0
1	B	35	LYS	2.9
1	A	71	MET	2.8
1	A	81	SER	2.8
1	B	477	TYR	2.8
1	B	53	GLY	2.7
1	B	90	GLY	2.7
1	B	418	LEU	2.7
1	A	68	ALA	2.7
1	A	127	MET	2.7
1	A	207	ALA	2.6
1	B	362	ALA	2.6
1	A	37	ILE	2.5
1	B	88	THR	2.5
1	B	458	ALA	2.4
1	A	281	ILE	2.4
1	A	59	LEU	2.4
1	A	278	LEU	2.4
1	A	67	THR	2.4
1	A	204	MET	2.4
1	B	460	LEU	2.4
1	A	477	TYR	2.3
1	A	44	TYR	2.2
1	A	113	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	89	GLU	2.2
1	A	83	GLY	2.2
1	A	223	ALA	2.2
1	B	225	ASP	2.2
1	B	400	LYS	2.1
1	A	397	SER	2.1
1	A	229	ALA	2.1
1	B	87	GLY	2.1
1	B	195	CYS	2.1
1	A	130	ALA	2.1
1	A	208	PHE	2.1
1	B	561	PHE	2.0
1	B	422	LEU	2.0
1	A	203	THR	2.0
1	A	128	PRO	2.0
1	B	54	ILE	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.