



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:30 AM GMT

PDB ID : 3B39  
Title : Structure of the DnaG primase catalytic domain bound to ssDNA  
Authors : Corn, J.E.; Pelton, J.G.; Berger, J.M.  
Deposited on : 2007-10-19  
Resolution : 2.35 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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

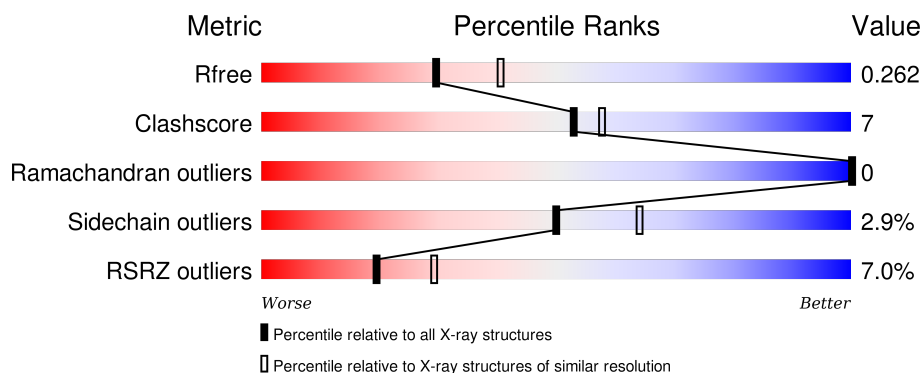
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	C	15	<div> <div>7%</div> <div>20% 13% 7% 60%</div> </div>
1	D	15	<div> <div>53% 47%</div> </div>
2	A	322	<div> <div>6%</div> <div>81% 16% ..</div> </div>
2	B	322	<div> <div>8%</div> <div>79% 17% .</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5354 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 DNA chain called DNA (5'-D(\*DCP\*DAP\*DAP\*DAP\*DGP\*DCP\*DCP\*DAP\*DAP\*DAP\*DAP\*DGP\*DGP\*DAP\*DC)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	6	Total	C	N	O	P	0	0	0
			108	49	23	30	6			
1	D	8	Total	C	N	O	P	0	0	0
			132	59	28	37	8			

- Molecule 2 is a protein called DNA primase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	317	Total	C	N	O	S	0	0	0
			2511	1577	451	471	12			
2	B	310	Total	C	N	O	S	0	0	0
			2457	1542	444	459	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	108	SER	-	EXPRESSION TAG	UNP P0ABS5
A	109	ASN	-	EXPRESSION TAG	UNP P0ABS5
A	110	ALA	-	EXPRESSION TAG	UNP P0ABS5
A	320	CYS	ARG	ENGINEERED	UNP P0ABS5
B	108	SER	-	EXPRESSION TAG	UNP P0ABS5
B	109	ASN	-	EXPRESSION TAG	UNP P0ABS5
B	110	ALA	-	EXPRESSION TAG	UNP P0ABS5
B	320	CYS	ARG	ENGINEERED	UNP P0ABS5

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	62	Total	O	0	0
			62	62		

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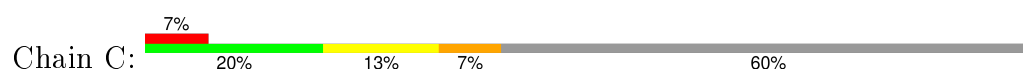
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	67	Total 67	O 67	0	0
3	C	8	Total 8	O 8	0	0
3	D	9	Total 9	O 9	0	0

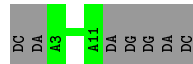
### 3 Residue-property plots [i](#)

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.

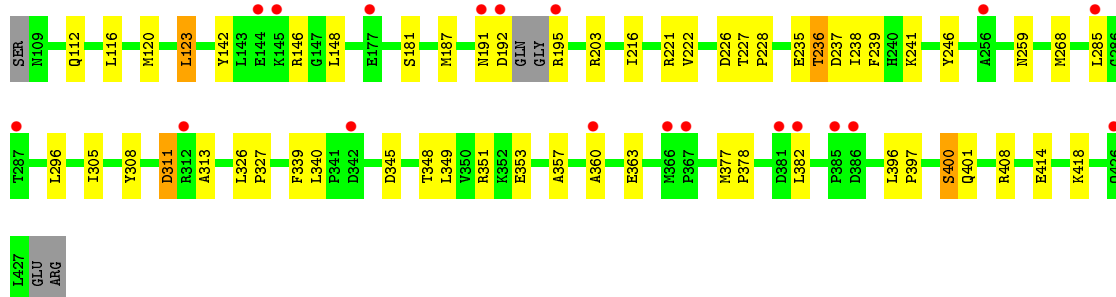
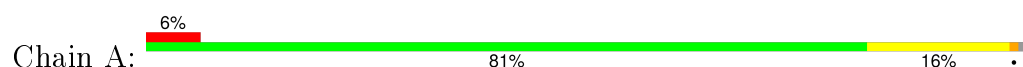
- Molecule 1: DNA (5'-D(\*DCP\*DAP\*DAP\*DAP\*DGP\*DCP\*DCP\*DAP\*DAP\*DAP\*DAP\*DGP\*DGP\*DAP\*DC)-3')



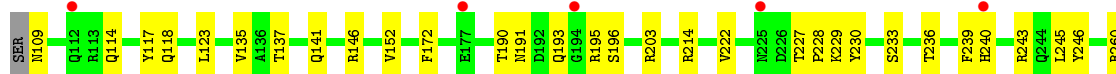
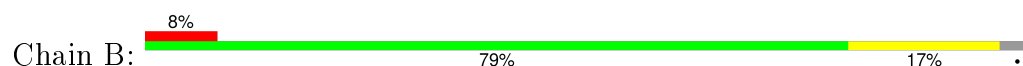
- Molecule 1: DNA (5'-D(\*DCP\*DAP\*DAP\*DAP\*DGP\*DCP\*DCP\*DAP\*DAP\*DAP\*DAP\*DGP\*DGP\*DAP\*DC)-3')



- Molecule 2: DNA primase



- Molecule 2: DNA primase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	136.83 Å   136.83 Å   71.07 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	40.00 – 2.35 37.89 – 2.35	Depositor EDS
% Data completeness (in resolution range)	94.6 (40.00-2.35) 94.6 (37.89-2.35)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.67 (at 2.34 Å)	Xtriage
Refinement program	REFMAC 5.3.0021	Depositor
R, $R_{free}$	0.212 , 0.254 0.216 , 0.262	Depositor DCC
$R_{free}$ test set	1519 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.5	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 52.2	EDS
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 30064 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5354	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	C	0.80	0/121	1.49	1/185 (0.5%)
1	D	0.71	0/147	1.15	0/222
2	A	0.54	0/2560	0.62	0/3463
2	B	0.54	0/2507	0.62	0/3392
All	All	0.55	0/5335	0.68	1/7262 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3	DA	P-O3'-C3'	-5.81	112.73	119.70

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	C	108	0	55	4	0
1	D	132	0	66	0	0
2	A	2511	0	2470	35	0
2	B	2457	0	2414	31	0
3	A	62	0	0	0	0
3	B	67	0	0	0	0
3	C	8	0	0	0	0
3	D	9	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5354	0	5005	67	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 7.

All (67) 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)
2:B:310:GLY:O	2:B:315:ARG:NH1	1.91	1.02
2:A:236:THR:HG22	2:A:238:ILE:H	1.27	0.99
2:A:246:TYR:HB2	2:A:285:LEU:HD21	1.71	0.72
2:A:345:ASP:OD1	2:A:348:THR:HG23	1.94	0.67
2:A:236:THR:HG22	2:A:238:ILE:N	2.07	0.67
2:A:396:LEU:O	2:A:400:SER:OG	2.15	0.65
2:A:222:VAL:HG21	2:A:226:ASP:O	2.00	0.62
2:A:340:LEU:HD22	2:A:349:LEU:HD12	1.82	0.62
2:A:236:THR:CG2	2:A:237:ASP:N	2.64	0.61
2:B:114:GLN:HE21	2:B:118:GLN:HE21	1.48	0.60
2:A:360:ALA:O	2:A:363:GLU:HG2	2.03	0.57
2:A:353:GLU:HB3	2:A:357:ALA:HB3	1.86	0.56
2:A:305:ILE:HD12	2:A:305:ILE:N	2.21	0.55
1:C:6:DC:H2''	2:B:109:ASN:HB2	1.91	0.53
2:B:190:THR:OG1	2:B:196:SER:OG	2.26	0.53
2:B:236:THR:O	2:B:239:PHE:O	2.26	0.53
1:C:2:DA:H2'	1:C:3:DA:C8	2.45	0.52
2:B:117:TYR:CE1	2:B:214:ARG:HD2	2.44	0.52
1:C:6:DC:C2'	2:B:109:ASN:HB2	2.41	0.51
2:B:372:LEU:CD2	2:B:411:LEU:HD22	2.40	0.50
1:C:2:DA:H2''	1:C:3:DA:O5'	2.11	0.50
2:A:112:GLN:O	2:A:116:LEU:HD13	2.11	0.50
2:B:114:GLN:HE21	2:B:118:GLN:NE2	2.09	0.49
2:A:414:GLU:O	2:A:418:LYS:HG2	2.13	0.49
2:B:341:PRO:HG2	2:B:344:GLU:HG3	1.95	0.49
2:A:191:ASN:O	2:A:192:ASP:HB2	2.14	0.47
2:A:116:LEU:O	2:A:120:MET:HG2	2.14	0.47
2:B:349:LEU:O	2:B:353:GLU:HG2	2.16	0.45
2:B:372:LEU:HD22	2:B:411:LEU:HD22	1.98	0.45
2:B:260:ARG:NH1	2:B:363:GLU:OE2	2.48	0.44
2:B:191:ASN:HB3	2:B:193:GLN:H	1.82	0.44
2:B:229:LYS:CG	2:B:230:TYR:CD2	3.00	0.44
2:B:245:LEU:HD23	2:B:284:SER:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:146:ARG:O	2:A:351:ARG:NH2	2.50	0.44
2:A:408:ARG:HH12	2:B:408:ARG:HH22	1.64	0.44
2:B:137:THR:O	2:B:141:GLN:HG3	2.17	0.44
2:B:377:MET:HB2	2:B:378:PRO:HD3	1.98	0.44
2:A:191:ASN:HB3	2:A:195:ARG:HA	1.99	0.44
2:A:246:TYR:CB	2:A:285:LEU:HD21	2.45	0.44
2:A:326:LEU:HB2	2:A:327:PRO:HD3	2.00	0.44
2:B:229:LYS:HG2	2:B:230:TYR:N	2.33	0.43
2:A:142:TYR:OH	2:A:221:ARG:NH1	2.51	0.43
2:B:340:LEU:HD13	2:B:345:ASP:HA	2.01	0.43
2:A:203:ARG:NH1	2:A:222:VAL:HG13	2.34	0.43
2:A:142:TYR:OH	2:A:221:ARG:CZ	2.66	0.43
2:B:203:ARG:NH1	2:B:222:VAL:HG13	2.33	0.43
2:A:227:THR:OG1	2:A:228:PRO:HA	2.19	0.43
2:A:148:LEU:HD11	2:A:268:MET:HE1	1.99	0.43
2:B:227:THR:HA	2:B:228:PRO:C	2.39	0.43
2:A:236:THR:HG22	2:A:237:ASP:N	2.33	0.42
2:B:246:TYR:HB2	2:B:285:LEU:HD21	2.01	0.42
2:A:235:GLU:OE2	2:A:241:LYS:HG3	2.20	0.42
2:A:123:LEU:HD23	2:A:187:MET:HG3	2.02	0.42
2:A:192:ASP:H	2:A:195:ARG:HA	1.84	0.42
2:A:216:ILE:O	2:A:239:PHE:HB2	2.19	0.42
2:A:377:MET:N	2:A:378:PRO:HD2	2.34	0.42
2:A:311:ASP:OD2	2:A:313:ALA:HB3	2.19	0.42
2:A:308:TYR:O	2:A:339:PHE:HA	2.19	0.42
2:B:369:SER:OG	2:B:411:LEU:HD23	2.20	0.41
2:B:240:HIS:HB3	2:B:243:ARG:HH11	1.85	0.41
2:B:146:ARG:NH2	2:B:347:ASP:OD2	2.51	0.41
2:A:123:LEU:CD2	2:A:187:MET:HG3	2.51	0.41
2:B:123:LEU:HD23	2:B:172:PHE:CE1	2.56	0.41
2:B:191:ASN:HB2	2:B:195:ARG:H	1.86	0.40
2:A:397:PRO:O	2:A:401:GLN:HG3	2.21	0.40
2:B:382:LEU:HD21	2:B:391:LEU:HD22	2.04	0.40
2:B:392:SER:O	2:B:396:LEU:HD12	2.22	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	
2	A	313/322 (97%)	306 (98%)	7 (2%)	0	100	100
2	B	308/322 (96%)	298 (97%)	10 (3%)	0	100	100
All	All	621/644 (96%)	604 (97%)	17 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	
2	A	265/269 (98%)	257 (97%)	8 (3%)	48	62
2	B	258/269 (96%)	251 (97%)	7 (3%)	52	67
All	All	523/538 (97%)	508 (97%)	15 (3%)	50	64

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	123	LEU
2	A	181	SER
2	A	236	THR
2	A	259	ASN
2	A	296	LEU
2	A	311	ASP
2	A	382	LEU
2	A	400	SER

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Mol	Chain	Res	Type
2	B	135	VAL
2	B	152	VAL
2	B	233	SER
2	B	284	SER
2	B	296	LEU
2	B	309	ASP
2	B	326	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	118	GLN
2	B	178	ASN
2	B	191	ASN

### 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 carbohydrates 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	C	6/15 (40%)	0.21	1 (16%) 2 4	65, 68, 76, 84	0
1	D	8/15 (53%)	0.53	0 100 100	69, 75, 93, 97	0
2	A	317/322 (98%)	0.32	19 (5%) 25 38	38, 50, 64, 78	0
2	B	310/322 (96%)	0.33	25 (8%) 15 23	35, 47, 71, 78	0
All	All	641/674 (95%)	0.33	45 (7%) 19 29	35, 49, 70, 97	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	381	ASP	5.3
2	B	382	LEU	5.1
2	B	387	GLY	5.0
2	B	391	LEU	5.0
2	B	225	ASN	4.4
2	B	389	ALA	4.4
2	B	385	PRO	4.4
2	A	287	THR	4.1
2	A	191	ASN	3.9
2	B	388	ARG	3.9
2	B	384	THR	3.7
2	A	342	ASP	3.6
2	A	386	ASP	3.6
2	B	415	LEU	3.5
2	B	112	GLN	3.5
2	B	381	ASP	3.4
2	B	380	VAL	3.3
2	A	195	ARG	3.3
2	A	385	PRO	3.3
2	A	145	LYS	3.2
2	A	177	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
2	A	285	LEU	3.0
2	B	285	LEU	2.8
2	A	192	ASP	2.8
2	B	390	ARG	2.8
2	B	379	GLN	2.7
2	A	144	GLU	2.7
2	B	411	LEU	2.6
2	A	367	PRO	2.5
2	A	382	LEU	2.5
2	B	392	SER	2.5
2	B	386	ASP	2.3
2	B	194	GLY	2.3
2	A	426	GLN	2.3
2	B	177	GLU	2.2
2	A	256	ALA	2.2
2	B	312	ARG	2.2
2	A	312	ARG	2.1
2	A	360	ALA	2.1
2	B	356	GLU	2.1
1	C	2	DA	2.1
2	B	383	SER	2.1
2	A	366	MET	2.0
2	B	410	TYR	2.0
2	B	240	HIS	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.