



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:49 AM UTC

PDB ID : 7JH5 / pdb_00007jh5
Title : Co-LOCKR: de novo designed protein switch
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Deposited on : 2020-07-20
Resolution : 2.10 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

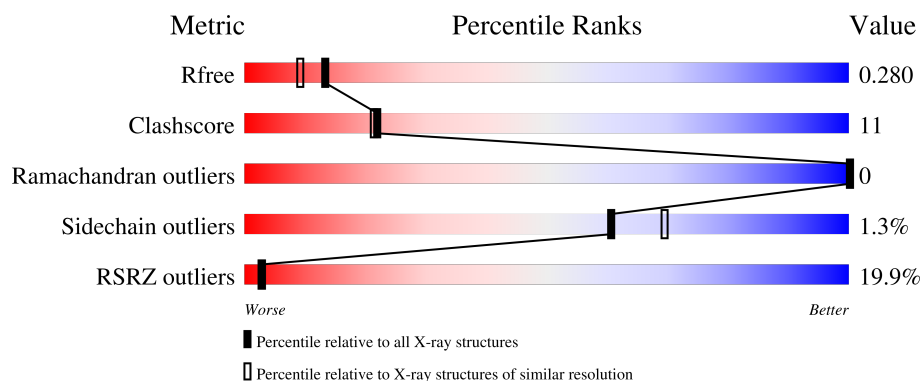
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 ≥ 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	291	21% 75% 23% ..
1	B	291	17% 79% 17% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8770 atoms, of which 4303 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 Co-LOCKR: de novo designed protein switch.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	H	N	O	0	3	0
			4379	1360	2201	400	418			
1	B	279	Total	C	H	N	O	0	4	0
			4201	1316	2102	381	402			

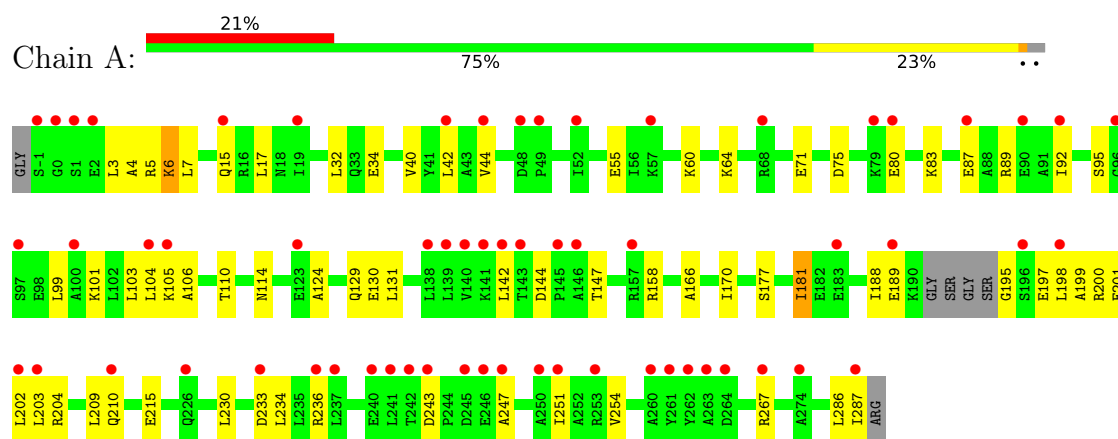
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	87	Total	O	0	0
			87	87		
2	B	103	Total	O	0	0
			103	103		

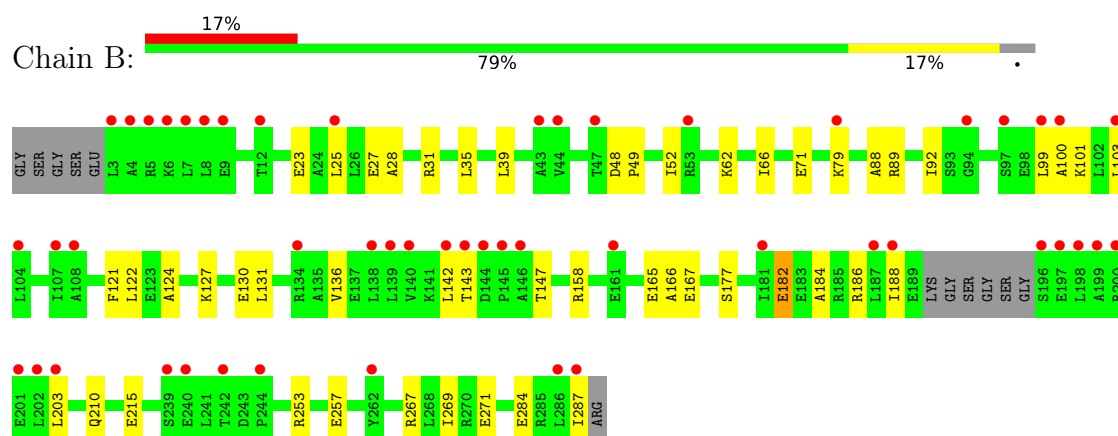
3 Residue-property plots [i](#)

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: Co-LOCKR: de novo designed protein switch



- Molecule 1: Co-LOCKR: de novo designed protein switch



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.84Å 52.91Å 75.59Å 90.00° 102.64° 90.00°	Depositor
Resolution (Å)	42.38 – 2.10 42.38 – 2.10	Depositor EDS
% Data completeness (in resolution range)	85.2 (42.38-2.10) 84.7 (42.38-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.10Å)	Xtriage
Refinement program	PHENIX dev 3026	Depositor
R, R_{free}	0.257 , 0.288 0.254 , 0.280	Depositor DCC
R_{free} test set	2000 reflections (7.23%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8770	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 59.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7780e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.25	1/2185 (0.0%)	0.36	0/2944
1	B	0.22	1/2113 (0.0%)	0.32	0/2852
All	All	0.23	2/4298 (0.0%)	0.34	0/5796

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	181	ILE	C-O	-8.06	1.15	1.24
1	B	182	GLU	C-O	-5.78	1.17	1.24

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	2178	2201	2194	47	1
1	B	2099	2102	2084	46	0
2	A	87	0	0	10	0
2	B	103	0	0	11	1
All	All	4467	4303	4278	92	2

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 11.

All (92) 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:234:LEU:HD13	1:A:254:VAL:HG21	1.60	0.82
1:B:39:LEU:HG	1:B:136:VAL:HG11	1.62	0.81
1:B:103:LEU:HD11	1:B:188:ILE:HD11	1.66	0.78
1:B:99:LEU:HD11	1:B:188:ILE:HG12	1.66	0.78
1:A:5:ARG:CG	1:A:287:ILE:HG21	2.13	0.78
1:B:167:GLU:OE1	2:B:301:HOH:O	2.01	0.77
1:B:127:LYS:NZ	1:B:165[B]:GLU:OE1	2.20	0.75
1:A:215[A]:GLU:OE1	2:A:301:HOH:O	2.06	0.73
1:B:130:GLU:OE2	2:B:302:HOH:O	2.07	0.73
1:B:215:GLU:OE2	2:B:303:HOH:O	2.08	0.72
1:A:197:GLU:OE2	2:A:302:HOH:O	2.08	0.72
1:A:3:LEU:HG	1:A:92:ILE:HD13	1.72	0.72
1:A:71:GLU:OE1	2:A:304:HOH:O	2.10	0.70
1:A:144:ASP:OD2	2:A:303:HOH:O	2.09	0.70
1:B:27:GLU:OE2	1:B:31:ARG:NE	2.25	0.70
1:A:42:LEU:HD13	1:A:55:GLU:HB3	1.74	0.69
1:A:130:GLU:OE1	2:A:305:HOH:O	2.10	0.68
1:A:3:LEU:HD21	1:A:95:SER:OG	1.94	0.68
1:B:66:ILE:O	2:B:304:HOH:O	2.11	0.67
1:B:167:GLU:OE2	2:B:305:HOH:O	2.13	0.66
1:A:15[A]:GLN:HG2	1:A:209:LEU:HD22	1.77	0.65
1:B:167:GLU:OE1	2:B:306:HOH:O	2.15	0.65
1:A:87:GLU:O	2:A:306:HOH:O	2.15	0.64
1:B:131:LEU:HD13	1:B:158[A]:ARG:HG2	1.80	0.64
1:B:103:LEU:HD11	1:B:188:ILE:CG1	2.28	0.63
1:B:103:LEU:HD11	1:B:188:ILE:CD1	2.29	0.61
1:B:188:ILE:HD12	1:B:203:LEU:HD13	1.82	0.61
1:B:142:LEU:HD13	1:B:147:THR:HB	1.83	0.60
1:B:35:LEU:HD21	1:B:62:LYS:HE2	1.84	0.60
1:B:103:LEU:CD1	1:B:188:ILE:HD11	2.32	0.59
1:A:267:ARG:NH1	2:A:309:HOH:O	2.33	0.58
1:A:188:ILE:HG13	1:A:203:LEU:HD12	1.85	0.58
1:A:4:ALA:HB1	1:A:202:LEU:HD11	1.87	0.56
1:B:71:GLU:HG3	1:B:122:LEU:HD21	1.86	0.56
1:B:49:PRO:HA	1:B:52:ILE:HG22	1.88	0.55
1:B:131:LEU:HD13	1:B:158[B]:ARG:HG2	1.89	0.54
1:A:177:SER:OG	1:A:210:GLN:HG2	2.08	0.53
1:B:271:GLU:OE1	2:B:307:HOH:O	2.19	0.53
1:A:4:ALA:HB1	1:A:202:LEU:CD1	2.39	0.53
1:B:143:THR:HG22	1:B:143:THR:O	2.08	0.53
1:A:99:LEU:HD11	1:A:188:ILE:CD1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:LEU:HD13	1:A:80:GLU:HG2	1.91	0.52
1:A:32:LEU:HD22	1:A:129:GLN:HG2	1.90	0.52
1:B:284:GLU:HA	1:B:287:ILE:HD12	1.93	0.51
1:A:230:LEU:HB3	1:A:254:VAL:HG13	1.92	0.51
1:A:189:GLU:OE1	1:A:200:ARG:NH2	2.44	0.50
1:A:234:LEU:HD13	1:A:254:VAL:CG2	2.37	0.50
1:A:201:GLU:HG3	1:A:286:LEU:HD13	1.93	0.50
1:B:89:ARG:HA	1:B:92:ILE:HG22	1.94	0.50
1:A:17:LEU:HD22	1:A:80:GLU:HG3	1.93	0.49
1:A:124:ALA:HB2	1:A:166:ALA:HB2	1.94	0.49
1:A:195:GLY:O	1:A:200:ARG:NH1	2.46	0.49
1:A:60:LYS:O	1:A:64:LYS:HG2	2.13	0.49
1:A:131:LEU:HD13	1:A:158:ARG:HG2	1.94	0.48
1:A:142:LEU:HD13	1:A:147:THR:HB	1.95	0.48
1:A:101:LYS:O	1:A:105:LYS:HG2	2.13	0.48
1:A:188:ILE:HD12	1:A:199:ALA:HB1	1.95	0.48
1:B:253[B]:ARG:NH1	2:B:316:HOH:O	2.47	0.47
1:A:233:ASP:OD1	1:A:236:ARG:NH2	2.47	0.47
1:B:124:ALA:HB2	1:B:166:ALA:HB2	1.97	0.47
1:A:243:ASP:N	2:A:314:HOH:O	2.43	0.46
1:A:5:ARG:CG	1:A:287:ILE:HD13	2.46	0.46
1:A:34:GLU:OE2	1:B:186:ARG:NH1	2.43	0.46
1:A:195:GLY:N	2:A:317:HOH:O	2.48	0.46
1:A:89:ARG:HG2	1:A:104:LEU:HD21	1.98	0.46
1:B:25:LEU:HD13	1:B:121:PHE:HD2	1.81	0.45
1:A:114:ASN:HB3	2:A:355:HOH:O	2.17	0.45
1:B:23:GLU:HG3	1:B:269:ILE:HG21	1.99	0.45
1:A:106:ALA:O	1:A:110:THR:HG23	2.16	0.45
1:B:253[B]:ARG:NH1	2:B:318:HOH:O	2.49	0.45
1:B:177:SER:OG	1:B:210:GLN:HG2	2.16	0.45
1:A:7:LEU:HD23	1:A:103:LEU:HB3	1.98	0.45
1:B:253[A]:ARG:NH1	1:B:257:GLU:OE2	2.50	0.45
1:A:83:LYS:O	1:A:87:GLU:HG3	2.17	0.43
1:B:28:ALA:HB2	2:B:341:HOH:O	2.17	0.43
1:A:6:LYS:HA	1:A:6:LYS:HD2	1.88	0.43
1:B:79:LYS:HE2	1:B:79:LYS:N	2.34	0.43
1:B:35:LEU:HD11	1:B:62:LYS:HB3	2.01	0.42
1:A:198:LEU:O	1:A:202:LEU:HG	2.20	0.42
1:A:40:VAL:HG12	1:A:251:ILE:HD11	2.01	0.42
1:B:25:LEU:HB3	1:B:121:PHE:HE2	1.84	0.42
1:B:89:ARG:HA	1:B:92:ILE:CG2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:ALA:HB3	1:B:203:LEU:HD11	2.01	0.42
1:A:44:VAL:HG11	1:A:247:ALA:HB3	2.01	0.41
1:B:267:ARG:NE	2:B:309:HOH:O	2.28	0.41
1:B:203:LEU:HD23	1:B:203:LEU:C	2.45	0.41
1:A:166:ALA:O	1:A:170:ILE:HG12	2.21	0.41
1:B:92:ILE:HD11	1:B:100:ALA:C	2.46	0.41
1:B:143:THR:O	1:B:143:THR:CG2	2.68	0.41
1:B:88:ALA:O	1:B:92:ILE:HG22	2.21	0.41
1:B:92:ILE:HD11	1:B:100:ALA:CB	2.51	0.41
1:B:103:LEU:CG	1:B:188:ILE:HD11	2.51	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:374:HOH:O	2:B:389:HOH:O[2_646]	2.10	0.10
1:A:75:ASP:OD1	1:A:204:ARG:HH22[2_555]	1.55	0.05

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	284/291 (98%)	281 (99%)	3 (1%)	0	100	100
1	B	279/291 (96%)	277 (99%)	2 (1%)	0	100	100
All	All	563/582 (97%)	558 (99%)	5 (1%)	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	201/237 (85%)	199 (99%)	2 (1%)	68	76
1	B	188/237 (79%)	185 (98%)	3 (2%)	55	64
All	All	389/474 (82%)	384 (99%)	5 (1%)	61	69

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	181	ILE
1	B	48	ASP
1	B	101	LYS
1	B	182	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	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 oligosaccharides 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, 95th 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 ²)	Q<0.9
1	A	285/291 (97%)	1.31	62 (21%)	2 2	14, 35, 56, 87	3 (1%)
1	B	279/291 (95%)	1.19	50 (17%)	3 4	13, 31, 60, 78	3 (1%)
All	All	564/582 (96%)	1.25	112 (19%)	3 3	13, 33, 57, 87	6 (1%)

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	196	SER	6.8
1	B	197	GLU	6.7
1	B	100	ALA	5.9
1	B	198	LEU	5.7
1	A	1	SER	5.5
1	B	203	LEU	5.0
1	A	240	GLU	4.9
1	B	199	ALA	4.8
1	A	96	GLY	4.7
1	A	262	TYR	4.6
1	B	200	ARG	4.6
1	B	3	LEU	4.2
1	A	44	VAL	4.2
1	A	97	SER	4.1
1	A	196	SER	4.1
1	B	5	ARG	4.0
1	B	4	ALA	3.9
1	B	240	GLU	3.7
1	B	140	VAL	3.7
1	A	92	ILE	3.7
1	B	103	LEU	3.6
1	A	274	ALA	3.5
1	B	144	ASP	3.5
1	B	99	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	145	PRO	3.4
1	A	-1	SER	3.4
1	A	141	LYS	3.4
1	B	287	ILE	3.4
1	A	139	LEU	3.4
1	B	201	GLU	3.3
1	B	108	ALA	3.3
1	B	107	ILE	3.2
1	A	267	ARG	3.2
1	A	142	LEU	3.1
1	A	241	LEU	3.1
1	A	264	ASP	3.1
1	B	8	LEU	3.1
1	A	0	GLY	3.0
1	A	226	GLN	3.0
1	A	261	TYR	3.0
1	A	263	ALA	3.0
1	B	146	ALA	3.0
1	A	42	LEU	3.0
1	B	138	LEU	3.0
1	A	2	GLU	3.0
1	B	104	LEU	2.9
1	A	104	LEU	2.9
1	A	251	ILE	2.8
1	A	123	GLU	2.8
1	B	202	LEU	2.8
1	B	187	LEU	2.7
1	A	233	ASP	2.7
1	A	105	LYS	2.7
1	A	138	LEU	2.7
1	A	68	ARG	2.7
1	A	100	ALA	2.6
1	B	25	LEU	2.6
1	A	49	PRO	2.6
1	A	202	LEU	2.6
1	B	139	LEU	2.6
1	A	287	ILE	2.6
1	A	242	THR	2.6
1	B	43	ALA	2.6
1	B	97	SER	2.6
1	B	47	THR	2.5
1	B	188	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	260	ALA	2.5
1	B	142	LEU	2.5
1	A	210	GLN	2.5
1	A	246	GLU	2.5
1	A	198	LEU	2.5
1	A	203	LEU	2.5
1	A	237	LEU	2.5
1	A	48	ASP	2.4
1	A	52	ILE	2.4
1	B	262	TYR	2.4
1	B	12	THR	2.4
1	A	19	ILE	2.4
1	B	6	LYS	2.4
1	A	247	ALA	2.3
1	A	243	ASP	2.3
1	A	245	ASP	2.3
1	B	242	THR	2.3
1	A	87	GLU	2.3
1	A	183	GLU	2.3
1	A	79	LYS	2.3
1	A	143	THR	2.3
1	B	94	GLY	2.3
1	A	80	GLU	2.3
1	A	189	GLU	2.3
1	B	239	SER	2.3
1	B	44	VAL	2.3
1	A	157	ARG	2.3
1	A	145	PRO	2.2
1	B	181	ILE	2.2
1	B	161	GLU	2.2
1	A	57	LYS	2.2
1	B	143	THR	2.2
1	A	140	VAL	2.2
1	A	253	ARG	2.2
1	A	90	GLU	2.1
1	A	250	ALA	2.1
1	B	7	LEU	2.1
1	B	286	LEU	2.1
1	B	53	ARG	2.1
1	B	134	ARG	2.1
1	B	244	PRO	2.1
1	B	79	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	15[A]	GLN	2.0
1	A	236	ARG	2.0
1	B	9	GLU	2.0
1	A	146	ALA	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 oligosaccharides 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.