Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 04:45 AM EDT

PDB ID	:	6NKO
Title	:	Crystal structure of ForH
Authors	:	Zheng, J.; Irani, S.; Zhang, Y.
Deposited on	:	2019-01-07
Resolution	:	2.40 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 (i) 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 (1)) were used in the production of this report:

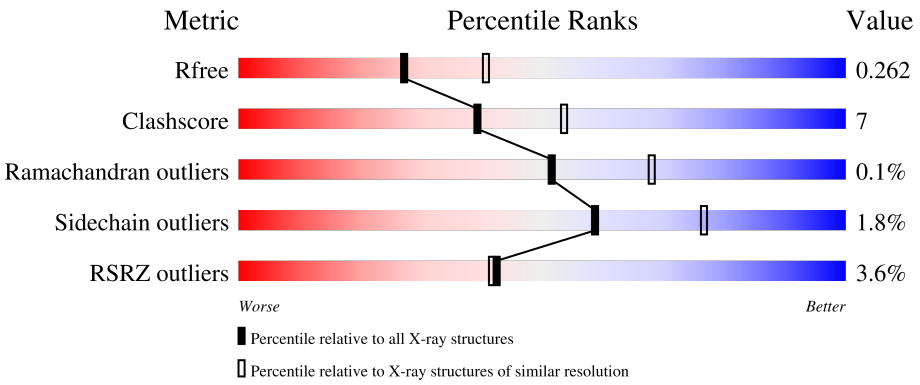
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	216	5% 72% 19% 9%
1	B	216	6% 73% 15% • 11%
1	C	216	. % 73% 15% 12%
1	D	216	. % 77% 11% 12%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5940 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 ForH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	196	1480	933	263	279	5	0	0	0
1	B	192	1453	919	256	273	5	0	0	0
1	C	190	1439	908	254	272	5	0	0	0
1	D	190	1442	911	254	272	5	0	0	0

- Molecule 2 is water.

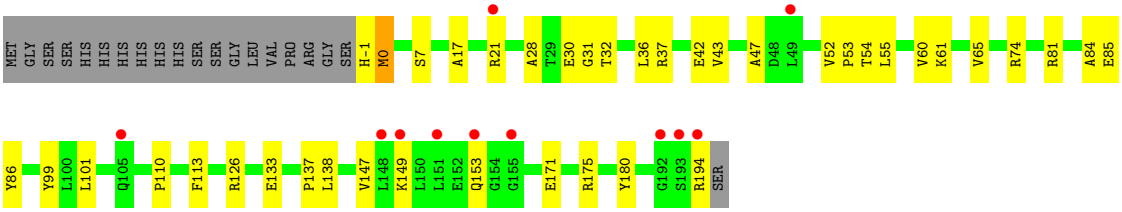
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	26	26	26	0	0
2	B	22	22	22	0	0
2	C	38	38	38	0	0
2	D	40	40	40	0	0

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.

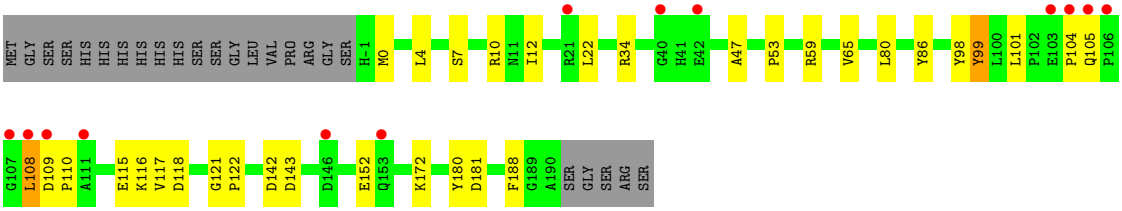
• Molecule 1: ForH

Chain A:



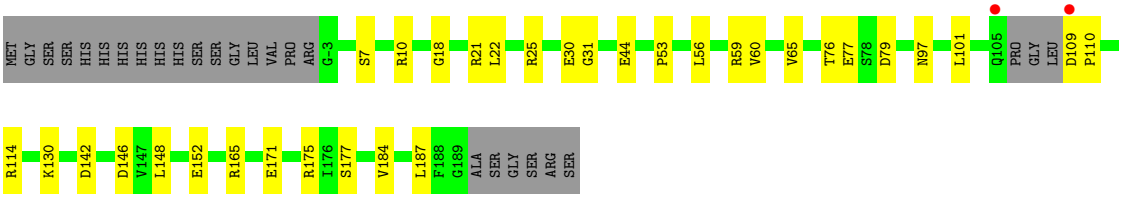
• Molecule 1: ForH

Chain B:



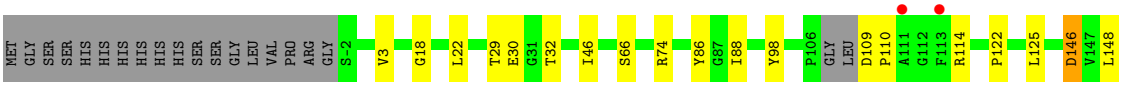
• Molecule 1: ForH

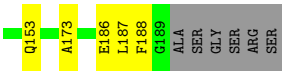
Chain C:



• Molecule 1: ForH

Chain D:




Q153
A173
E186
L187
F188
G189
ALA
SER
GLY
SER
ARG
SER

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	57.22Å 111.40Å 163.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.93 – 2.40 48.93 – 2.40	Depositor EDS
% Data completeness (in resolution range)	89.7 (48.93-2.40) 89.7 (48.93-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.63 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.212 , 0.262 0.211 , 0.262	Depositor DCC
R_{free} test set	1864 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	44.9	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5940	wwPDB-VP
Average B, all atoms (Å^2)	50.0	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 25.77 % 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 2.9785e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.30	0/1506	0.47	0/2042
1	B	0.26	0/1479	0.48	1/2007 (0.0%)
1	C	0.26	0/1463	0.43	0/1982
1	D	0.26	0/1467	0.43	0/1989
All	All	0.27	0/5915	0.45	1/8020 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	108	LEU	C-N-CA	5.17	134.61	121.70

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	A	1480	0	1499	32	0
1	B	1453	0	1473	26	0
1	C	1439	0	1454	23	0
1	D	1442	0	1458	17	0
2	A	26	0	0	0	0
2	B	22	0	0	0	0
2	C	38	0	0	1	0
2	D	40	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5940	0	5884	86	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 (86) 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:105:GLN:H	1:B:108:LEU:HD22	1.05	1.08
1:B:105:GLN:N	1:B:108:LEU:HD22	1.84	0.91
1:C:22:LEU:HD21	1:C:152:GLU:HG3	1.63	0.78
1:D:146:ASP:OD2	2:D:201:HOH:O	2.02	0.76
1:B:105:GLN:O	1:B:108:LEU:CD2	2.38	0.72
1:C:10:ARG:NH2	1:C:142:ASP:OD1	2.27	0.67
1:A:52:VAL:CG1	1:A:60:VAL:HG23	2.25	0.66
1:B:105:GLN:O	1:B:108:LEU:HD23	1.94	0.66
1:C:18:GLY:HA3	1:C:148:LEU:HD21	1.78	0.66
1:A:175:ARG:NH1	1:D:186:GLU:OE2	2.28	0.66
1:B:109:ASP:OD2	1:B:110:PRO:HD2	1.96	0.65
1:A:52:VAL:HG22	1:A:55:LEU:HD12	1.78	0.65
1:B:10:ARG:NH2	1:B:142:ASP:OD1	2.29	0.64
1:B:104:PRO:HA	1:B:108:LEU:HD13	1.79	0.64
1:A:53:PRO:O	1:A:54:THR:HG22	1.99	0.63
1:A:47:ALA:C	1:A:53:PRO:HG3	2.21	0.60
1:B:104:PRO:HA	1:B:108:LEU:CD1	2.31	0.60
1:B:22:LEU:HD11	1:B:152:GLU:HG3	1.83	0.60
1:D:66:SER:HA	1:D:88:ILE:HD12	1.84	0.58
1:B:105:GLN:O	1:B:108:LEU:HD22	2.03	0.57
1:A:149:LYS:O	1:A:153:GLN:HG2	2.05	0.57
1:D:29:THR:HG23	1:D:46:ILE:HD12	1.88	0.56
1:C:21:ARG:NH1	2:C:206:HOH:O	2.34	0.56
1:D:3:VAL:HG13	1:D:46:ILE:HD11	1.87	0.56
1:C:7:SER:N	1:C:97:ASN:OD1	2.36	0.55
1:C:77:GLU:H	1:C:77:GLU:CD	2.09	0.55
1:C:76:THR:HG23	1:C:79:ASP:H	1.73	0.54
1:C:101:LEU:HD23	1:C:187:LEU:HD22	1.88	0.54
1:C:109:ASP:N	1:C:109:ASP:OD1	2.40	0.54
1:A:86:TYR:HA	1:B:65:VAL:HB	1.89	0.54
1:B:47:ALA:HB1	1:B:53:PRO:HB3	1.90	0.54
1:A:52:VAL:O	1:A:52:VAL:HG13	2.08	0.53
1:D:18:GLY:HA3	1:D:148:LEU:HD21	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ARG:O	1:A:85:GLU:HG2	2.09	0.53
1:D:30:GLU:HG3	1:D:32:THR:H	1.73	0.52
1:C:53:PRO:HG2	1:C:60:VAL:HG12	1.93	0.51
1:B:101:LEU:HD11	1:B:180:TYR:CE1	2.46	0.50
1:C:146:ASP:OD1	1:C:165:ARG:NH2	2.45	0.50
1:A:-1:HIS:HB3	1:A:0:MET:SD	2.52	0.49
1:D:74:ARG:NH2	2:D:206:HOH:O	2.43	0.49
1:A:28:ALA:HB1	1:A:32:THR:HB	1.94	0.49
1:A:110:PRO:HA	1:A:113:PHE:CE2	2.48	0.49
1:B:4:LEU:HG	1:B:12:ILE:HD11	1.93	0.49
1:A:138:LEU:HD21	1:A:147:VAL:HG21	1.94	0.49
1:A:30:GLU:HG2	1:A:31:GLY:N	2.29	0.48
1:B:143:ASP:CG	1:B:172:LYS:HZ2	2.16	0.48
1:A:36:LEU:HD13	1:A:43:VAL:HG11	1.95	0.48
1:D:125:LEU:HG	1:D:173:ALA:HB1	1.96	0.48
1:A:17:ALA:O	1:A:21:ARG:HG3	2.14	0.48
1:C:30:GLU:CG	1:C:31:GLY:H	2.26	0.48
1:B:80:LEU:HB3	1:D:86:TYR:HE1	1.78	0.48
1:A:171:GLU:HB2	1:D:188:PHE:HB3	1.96	0.47
1:B:65:VAL:HG22	1:C:60:VAL:HG11	1.96	0.47
1:B:99:TYR:CE1	1:B:116:LYS:HD3	2.49	0.47
1:A:52:VAL:HG11	1:A:60:VAL:HG23	1.96	0.47
1:A:133:GLU:OE2	1:D:114:ARG:NH2	2.32	0.47
1:A:52:VAL:HG12	1:A:61:LYS:HA	1.96	0.46
1:A:52:VAL:CG1	1:A:52:VAL:O	2.63	0.46
1:C:25:ARG:NH1	1:C:44:GLU:OE1	2.49	0.46
1:D:18:GLY:O	1:D:22:LEU:HD13	2.15	0.46
1:B:59:ARG:NH2	1:B:115:GLU:OE2	2.48	0.46
1:B:99:TYR:CZ	1:B:116:LYS:HD3	2.51	0.45
1:A:47:ALA:HB1	1:A:53:PRO:HB3	1.98	0.45
1:C:30:GLU:HG2	1:C:31:GLY:H	1.81	0.45
1:B:181:ASP:OD2	1:C:177:SER:OG	2.30	0.44
1:A:65:VAL:HB	1:B:86:TYR:HA	2.00	0.44
1:A:101:LEU:HD11	1:A:180:TYR:CE1	2.52	0.44
1:A:101:LEU:HD11	1:A:180:TYR:HE1	1.81	0.44
1:A:52:VAL:CG2	1:A:55:LEU:HD12	2.46	0.43
1:A:42:GLU:OE2	1:A:42:GLU:N	2.38	0.43
1:C:101:LEU:HD22	1:C:184:VAL:HA	1.99	0.43
1:A:52:VAL:HG13	1:A:60:VAL:HG23	1.99	0.43
1:C:110:PRO:O	1:C:114:ARG:HG3	2.18	0.43
1:B:117:VAL:O	1:C:130:LYS:NZ	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:TYR:CE1	1:B:122:PRO:HG3	2.54	0.42
1:C:56:LEU:HB3	1:C:59:ARG:HH12	1.84	0.42
1:A:-1:HIS:HB3	1:A:0:MET:CE	2.49	0.42
1:B:118:ASP:OD2	1:B:121:GLY:N	2.47	0.42
1:D:187:LEU:HA	1:D:187:LEU:HD12	1.81	0.42
1:A:84:ALA:HB1	1:C:65:VAL:HG21	2.01	0.42
1:D:153:GLN:OE1	1:D:153:GLN:N	2.53	0.41
1:B:188:PHE:HB3	1:C:171:GLU:HB2	2.03	0.41
1:A:7:SER:HB3	1:A:99:TYR:CD1	2.56	0.41
1:A:126:ARG:NH2	1:D:122:PRO:HG2	2.36	0.41
1:D:109:ASP:HB3	1:D:110:PRO:HD3	2.03	0.41
1:C:165:ARG:HD3	1:C:165:ARG:O	2.22	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	194/216 (90%)	185 (95%)	8 (4%)	1 (0%)	29 41
1	B	190/216 (88%)	184 (97%)	6 (3%)	0	100 100
1	C	186/216 (86%)	178 (96%)	8 (4%)	0	100 100
1	D	186/216 (86%)	177 (95%)	9 (5%)	0	100 100
All	All	756/864 (88%)	724 (96%)	31 (4%)	1 (0%)	51 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	PRO

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	152/169 (90%)	148 (97%)	4 (3%)	46	66
1	B	149/169 (88%)	145 (97%)	4 (3%)	44	65
1	C	148/169 (88%)	147 (99%)	1 (1%)	84	92
1	D	149/169 (88%)	147 (99%)	2 (1%)	69	84
All	All	598/676 (88%)	587 (98%)	11 (2%)	59	76

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	MET
1	A	37	ARG
1	A	74	ARG
1	A	194	ARG
1	B	0	MET
1	B	7	SER
1	B	34	ARG
1	B	99	TYR
1	C	175	ARG
1	D	98	TYR
1	D	146	ASP

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

Mol	Chain	Res	Type
1	B	105	GLN
1	B	153	GLN
1	D	96	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 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, 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	196/216 (90%)	0.15	11 (5%)	24	23	31, 47, 70, 100	0
1	B	192/216 (88%)	0.29	13 (6%)	17	15	34, 51, 79, 101	0
1	C	190/216 (87%)	-0.07	2 (1%)	80	79	30, 44, 70, 95	0
1	D	190/216 (87%)	-0.13	2 (1%)	80	79	31, 45, 72, 92	0
All	All	768/864 (88%)	0.06	28 (3%)	42	42	30, 47, 73, 101	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	105	GLN	5.6
1	A	194	ARG	5.3
1	B	104	PRO	4.7
1	B	108	LEU	4.7
1	B	106	PRO	4.6
1	B	107	GLY	3.8
1	A	193	SER	3.4
1	A	151	LEU	3.2
1	B	105	GLN	3.2
1	B	146	ASP	3.1
1	B	111	ALA	3.1
1	C	109	ASP	3.0
1	A	192	GLY	3.0
1	B	40	GLY	3.0
1	A	155	GLY	3.0
1	A	49	LEU	2.8
1	B	103	GLU	2.6
1	B	153	GLN	2.4
1	B	109	ASP	2.3
1	A	149	LYS	2.3
1	B	42	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	148	LEU	2.2
1	D	113	PHE	2.2
1	A	153	GLN	2.2
1	A	105	GLN	2.1
1	B	21	ARG	2.1
1	D	111	ALA	2.0
1	A	21	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 (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.