



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:22 PM GMT

PDB ID : 4I36
Title : Crystal Structure of the Bacillus stearothermophilus Phosphofructokinase Mutant D12A
Authors : Mosser, R.; Reddy, M.; Bruning, J.B.; Sacchettini, J.C.; Reinhart, G.D.
Deposited on : 2012-11-25
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

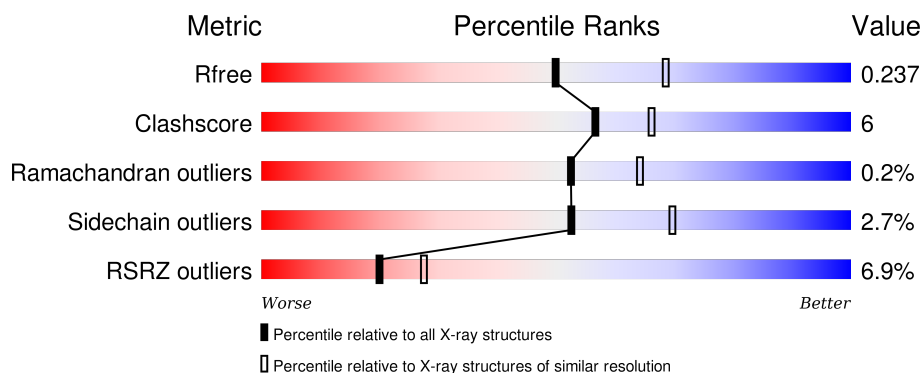
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 ≥ 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	A	319	<div> <div>6%</div> <div>85%</div> <div>14%</div> </div>
1	B	319	<div> <div>9%</div> <div>86%</div> <div>13%</div> </div>
1	C	319	<div> <div>10%</div> <div>87%</div> <div>12%</div> </div>
1	D	319	<div> <div>3%</div> <div>83%</div> <div>15%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9846 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 6-phosphofructokinase.

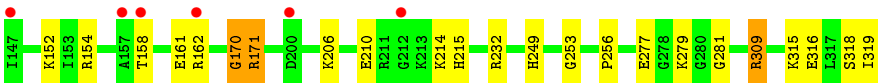
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	2	0
			2370	1479	430	452	9			
1	B	319	Total	C	N	O	S	0	3	0
			2388	1495	427	457	9			
1	C	319	Total	C	N	O	S	0	2	0
			2380	1493	427	451	9			
1	D	319	Total	C	N	O	S	0	1	0
			2379	1487	431	452	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
B	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
C	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
D	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	76	Total	O	0	0
			76	76		
2	B	72	Total	O	0	0
			72	72		
2	C	75	Total	O	0	0
			75	75		
2	D	106	Total	O	0	0
			106	106		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.05Å 112.66Å 129.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.89 – 2.30 38.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.89-2.30) 99.9 (38.89-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.201 , 0.254 0.185 , 0.237	Depositor DCC
R_{free} test set	3205 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	43.6	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	2 of 63132 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9846	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	A	0.28	0/2412	0.45	1/3258 (0.0%)
1	B	0.27	0/2433	0.45	0/3284
1	C	0.26	0/2422	0.45	0/3267
1	D	0.29	0/2417	0.48	1/3262 (0.0%)
All	All	0.28	0/9684	0.46	2/13071 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	7	LEU	CA-CB-CG	6.13	129.40	115.30
1	A	7	LEU	CA-CB-CG	5.58	128.12	115.30

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	2370	0	2358	39	0
1	B	2388	0	2380	28	0
1	C	2380	0	2388	26	0
1	D	2379	0	2377	36	0
2	A	76	0	0	1	0
2	B	72	0	0	2	0
2	C	75	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	106	0	0	4	0
All	All	9846	0	9503	116	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 6.

All (116) 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:D:28:ILE:HD13	1:D:54:VAL:HG13	1.57	0.86
1:D:102:GLY:HA2	1:D:130:ILE:HD11	1.61	0.81
1:C:3:ARG:HD2	1:C:35:TYR:CE1	2.20	0.77
1:D:3:ARG:HG2	1:D:33:GLU:HB2	1.65	0.77
1:D:277:GLU:HG3	1:D:279:LYS:HD3	1.68	0.74
1:D:79:GLU:O	1:D:83:LYS:HG2	1.86	0.74
1:D:38:TYR:O	1:D:43:GLY:HA3	1.89	0.71
1:B:97:LEU:HB3	1:B:119[B]:CYS:SG	2.30	0.71
1:C:38:TYR:O	1:C:43:GLY:HA3	1.90	0.71
1:A:188:THR:HB	1:A:218:ILE:HD12	1.76	0.67
1:D:113:THR:HG21	1:D:281:GLY:N	2.12	0.66
1:B:7:LEU:HB3	1:B:37:VAL:HB	1.79	0.65
1:D:3:ARG:HD3	1:D:35:TYR:CE1	2.32	0.65
1:B:187:GLU:HA	1:B:318:SER:HB2	1.79	0.65
1:A:38:TYR:O	1:A:43:GLY:HA3	1.98	0.64
1:A:28:ILE:HD13	1:A:54:VAL:HG13	1.79	0.63
1:A:60:ILE:HA	1:A:63:ARG:HG3	1.81	0.62
1:C:226:SER:O	1:C:230:PHE:HD2	1.82	0.61
1:A:113:THR:HG21	1:A:281:GLY:H	1.66	0.61
1:C:7:LEU:HB3	1:C:37:VAL:HB	1.83	0.60
1:C:102:GLY:HA2	1:C:130:ILE:HD11	1.82	0.60
1:B:158:THR:HG22	1:B:159:SER:H	1.66	0.59
1:C:59:ASP:HB2	1:D:214:LYS:HE3	1.85	0.59
1:D:315:LYS:HD2	2:D:492:HOH:O	2.02	0.59
1:A:113:THR:HG21	1:A:281:GLY:N	2.18	0.59
1:A:97:LEU:HB3	1:A:119[A]:CYS:SG	2.43	0.58
1:B:102:GLY:HA2	1:B:130:ILE:HD11	1.85	0.58
1:B:38:TYR:O	1:B:43:GLY:HA3	2.02	0.58
1:D:206:LYS:O	1:D:210:GLU:HG2	2.03	0.58
1:B:44:LEU:HD11	1:B:89:LEU:HG	1.86	0.58
1:B:128:ASN:OD1	1:B:135:PHE:HA	2.04	0.57
1:C:158:THR:O	1:C:160:HIS:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:GLY:HA2	1:A:84:LYS:HE3	1.86	0.56
1:B:158:THR:HG22	1:B:159:SER:N	2.21	0.56
1:B:33:GLU:OE2	1:B:51:LYS:HE3	2.06	0.56
1:C:44:LEU:HD11	1:C:89:LEU:HG	1.88	0.56
1:A:158:THR:HG22	1:D:70:THR:HG23	1.88	0.54
1:D:7:LEU:C	1:D:7:LEU:HD23	2.28	0.54
1:D:82:GLN:O	1:D:86:ILE:HG13	2.08	0.54
1:D:7:LEU:HG	1:D:40:GLY:HA2	1.90	0.53
1:D:113:THR:HG21	1:D:281:GLY:H	1.71	0.53
1:A:206:LYS:O	1:A:210:GLU:HG2	2.08	0.53
1:A:160:HIS:HD2	2:D:500:HOH:O	1.91	0.53
1:D:112:LEU:HD13	1:D:119[B]:CYS:SG	2.49	0.53
1:C:158:THR:O	1:C:161:GLU:N	2.42	0.53
1:A:210:GLU:OE1	1:A:210:GLU:HA	2.10	0.52
1:B:250:VAL:HG23	1:C:152:LYS:HB2	1.92	0.52
1:A:160:HIS:CD2	2:D:500:HOH:O	2.63	0.51
1:C:97:LEU:HB3	1:C:119[A]:CYS:SG	2.50	0.51
1:D:154:ARG:O	1:D:215:HIS:CE1	2.64	0.51
1:A:7:LEU:HB3	1:A:37:VAL:HB	1.92	0.51
1:A:162:ARG:O	1:A:162:ARG:HG2	2.11	0.51
1:C:89:LEU:HD13	1:C:117:PHE:CE1	2.47	0.50
1:B:110:LYS:O	1:B:113:THR:HB	2.13	0.49
1:D:158:THR:HB	1:D:161:GLU:HB2	1.94	0.49
1:A:113:THR:HG22	1:A:119[A]:CYS:H	1.78	0.49
1:B:250:VAL:HG22	1:C:152:LYS:HD3	1.95	0.49
1:A:128:ASN:OD1	1:A:135:PHE:HA	2.13	0.49
1:A:113:THR:HG22	1:A:119[B]:CYS:H	1.79	0.48
1:D:232:ARG:NH2	2:D:481:HOH:O	2.47	0.48
1:D:128:ASN:OD1	1:D:135:PHE:HA	2.14	0.48
1:A:113:THR:CG2	1:A:119[A]:CYS:H	2.27	0.48
1:C:7:LEU:CD1	1:C:97:LEU:HD21	2.43	0.48
1:A:160:HIS:O	1:A:241:GLU:HB2	2.13	0.48
1:A:113:THR:HA	1:A:117:PHE:O	2.14	0.47
1:D:113:THR:HG21	1:D:281:GLY:CA	2.44	0.47
1:A:113:THR:CG2	1:A:119[B]:CYS:H	2.27	0.47
1:D:15:GLY:N	1:D:256:PRO:HG3	2.30	0.47
1:C:311:TYR:CE2	1:C:315:LYS:HE3	2.50	0.47
1:C:3:ARG:HD2	1:C:35:TYR:CD1	2.50	0.47
1:A:188:THR:HB	1:A:218:ILE:CD1	2.41	0.46
1:D:318:SER:O	1:D:319:ILE:C	2.53	0.46
2:A:442:HOH:O	1:D:158:THR:HG23	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:GLU:HA	1:C:318:SER:HB2	1.97	0.46
1:D:309:ARG:NH2	1:D:316:GLU:OE2	2.48	0.46
1:B:63:ARG:NH1	1:D:63:ARG:HD2	2.30	0.45
1:C:277:GLU:OE2	1:C:279:LYS:HE2	2.16	0.45
1:A:8:THR:HA	1:A:100:ILE:O	2.16	0.45
1:C:257:THR:O	1:C:261:ARG:HG3	2.16	0.45
1:B:169:MET:HE1	1:B:249:HIS:HA	1.99	0.45
1:B:250:VAL:HG22	1:C:152:LYS:CD	2.47	0.44
1:C:163:THR:HG21	1:C:205:LEU:HD11	1.99	0.44
1:A:317:LEU:O	1:B:25:ARG:HD2	2.18	0.44
1:C:201:VAL:O	1:C:204:ARG:HB2	2.17	0.44
1:B:198:MET:O	1:B:202:ILE:HG12	2.17	0.44
1:A:162:ARG:HG3	1:A:214:LYS:O	2.17	0.43
1:A:261:ARG:CZ	1:B:147:ILE:HD12	2.49	0.43
1:D:130:ILE:HA	1:D:131:PRO:HD3	1.85	0.43
1:B:130:ILE:HA	1:B:131:PRO:HD3	1.89	0.43
1:A:277:GLU:OE1	1:A:279:LYS:HE3	2.17	0.43
1:B:8:THR:HA	1:B:100:ILE:O	2.18	0.43
1:D:117:PHE:HA	1:D:118:PRO:HD3	1.88	0.43
1:A:113:THR:HG22	1:A:118:PRO:HA	2.01	0.43
1:A:160:HIS:ND1	1:D:72:ARG:NH2	2.67	0.43
1:C:205:LEU:HD22	1:C:240:PHE:CD2	2.54	0.42
1:A:144:ASN:HA	1:A:147:ILE:HG12	2.01	0.42
1:C:226:SER:O	1:C:230:PHE:CD2	2.68	0.42
1:A:152:LYS:O	1:D:253:GLY:HA3	2.20	0.42
1:B:153:ILE:HD12	1:B:166:ILE:HD11	2.00	0.42
1:B:252:ARG:NH1	2:B:461:HOH:O	2.34	0.42
1:A:95:GLU:O	1:A:118:PRO:HD2	2.20	0.42
1:D:90:LYS:HB2	1:D:90:LYS:HE3	1.71	0.42
1:B:7:LEU:C	1:B:7:LEU:CD2	2.88	0.42
1:B:82:GLN:O	1:B:86:ILE:HG13	2.20	0.42
1:A:243:ARG:HD2	1:D:249:HIS:CE1	2.55	0.42
1:B:103:ASP:OD1	1:B:103:ASP:N	2.52	0.42
1:B:24:VAL:HA	1:B:34:VAL:HG21	2.00	0.42
1:A:83:LYS:O	1:A:87:GLU:HG2	2.20	0.41
1:A:160:HIS:ND1	1:D:72:ARG:CZ	2.84	0.41
1:D:170:GLY:O	1:D:171:ARG:CB	2.68	0.41
1:A:257:THR:O	1:A:261:ARG:HG3	2.20	0.41
1:C:117:PHE:HA	1:C:118:PRO:HD3	1.79	0.41
1:A:318:SER:HB3	2:B:467:HOH:O	2.20	0.41
1:A:253:GLY:HA3	1:D:152:LYS:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ASN:OD1	1:C:135:PHE:HA	2.21	0.40
1:B:274:LEU:HD22	1:B:279:LYS:HD2	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	318/319 (100%)	311 (98%)	7 (2%)	0	100	100
1	B	320/319 (100%)	312 (98%)	8 (2%)	0	100	100
1	C	319/319 (100%)	308 (97%)	10 (3%)	1 (0%)	46	57
1	D	318/319 (100%)	310 (98%)	6 (2%)	2 (1%)	30	36
All	All	1275/1276 (100%)	1241 (97%)	31 (2%)	3 (0%)	52	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	159	SER
1	D	171	ARG
1	D	170	GLY

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	241/246 (98%)	233 (97%)	8 (3%)	45	61
1	B	243/246 (99%)	237 (98%)	6 (2%)	55	73
1	C	242/246 (98%)	238 (98%)	4 (2%)	68	83
1	D	242/246 (98%)	234 (97%)	8 (3%)	45	61
All	All	968/984 (98%)	942 (97%)	26 (3%)	52	70

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	54	VAL
1	A	63	ARG
1	A	66	THR
1	A	79	GLU
1	A	86	ILE
1	A	215	HIS
1	A	309	ARG
1	B	7	LEU
1	B	66	THR
1	B	70	THR
1	B	103	ASP
1	B	161	GLU
1	B	215	HIS
1	C	7	LEU
1	C	66	THR
1	C	215	HIS
1	C	309	ARG
1	D	7	LEU
1	D	54	VAL
1	D	59	ASP
1	D	66	THR
1	D	70	THR
1	D	103	ASP
1	D	162	ARG
1	D	309	ARG

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

Mol	Chain	Res	Type
1	B	209	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 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, 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	318/319 (99%)	0.29	19 (5%) 25 33	24, 45, 76, 95	0
1	B	319/319 (100%)	0.49	28 (8%) 12 18	26, 45, 81, 102	0
1	C	319/319 (100%)	0.58	33 (10%) 9 13	27, 50, 84, 100	0
1	D	319/319 (100%)	0.25	8 (2%) 61 70	25, 41, 67, 88	0
All	All	1275/1276 (99%)	0.40	88 (6%) 20 27	24, 45, 79, 102	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	259[A]	PHE	6.7
1	B	160	HIS	4.7
1	B	259[A]	PHE	4.6
1	B	79	GLU	4.2
1	A	113	THR	4.1
1	C	230	PHE	4.1
1	C	202	ILE	4.0
1	B	113	THR	3.9
1	B	80	GLU	3.8
1	C	72	ARG	3.8
1	D	72	ARG	3.8
1	B	86	ILE	3.7
1	B	74	PRO	3.5
1	A	301	ALA	3.5
1	A	209	HIS	3.5
1	C	200	ASP	3.3
1	D	158	THR	3.3
1	D	74	PRO	3.2
1	A	205	LEU	3.2
1	C	238	THR	3.2
1	C	71	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	215	HIS	3.1
1	A	160	HIS	3.1
1	B	78	THR	3.1
1	A	86	ILE	3.1
1	D	200	ASP	3.1
1	A	199	ASN	3.1
1	C	195	ASP	3.0
1	B	240	PHE	2.9
1	B	319	ILE	2.9
1	B	232	ARG	2.9
1	B	158	THR	2.8
1	A	116	GLY	2.8
1	D	157	ALA	2.8
1	B	92	HIS	2.8
1	C	159	SER	2.7
1	C	223	GLY	2.7
1	C	301	ALA	2.7
1	C	158	THR	2.7
1	A	211	ARG	2.7
1	D	162	ARG	2.7
1	B	203	ALA	2.6
1	A	197	ASP	2.6
1	C	196	TYR	2.6
1	B	87	GLU	2.6
1	A	201	VAL	2.6
1	C	297	ALA	2.6
1	B	89	LEU	2.6
1	A	117	PHE	2.5
1	A	237	ALA	2.5
1	B	76	PHE	2.5
1	C	197	ASP	2.5
1	A	200	ASP	2.5
1	C	3	ARG	2.5
1	C	240	PHE	2.5
1	C	92	HIS	2.4
1	C	221	ALA	2.4
1	C	207	ARG	2.4
1	C	211	ARG	2.4
1	B	112	LEU	2.4
1	D	147	ILE	2.4
1	D	212	GLY	2.4
1	B	91	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	159	SER	2.4
1	C	224	VAL	2.4
1	B	81	GLY	2.3
1	A	162	ARG	2.3
1	A	240	PHE	2.2
1	B	205	LEU	2.2
1	C	93	GLY	2.2
1	A	210	GLU	2.2
1	B	200	ASP	2.2
1	C	194	ALA	2.2
1	C	172	HIS	2.2
1	C	110	LYS	2.2
1	B	93	GLY	2.2
1	C	215	HIS	2.2
1	B	46	ALA	2.2
1	B	199	ASN	2.1
1	C	199	ASN	2.1
1	B	84	LYS	2.1
1	C	226	SER	2.1
1	C	49	ILE	2.1
1	C	143	LEU	2.0
1	C	319	ILE	2.0
1	C	48	ASN	2.0
1	B	85	GLY	2.0
1	B	234	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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.