



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 11:32 AM EDT

PDB ID : 7JQY
Title : Crystal structure of Cfl1-D123S from Burkholderia cenocepacia
Authors : Taher, N.M.; Madden, D.R.
Deposited on : 2020-08-11
Resolution : 2.15 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

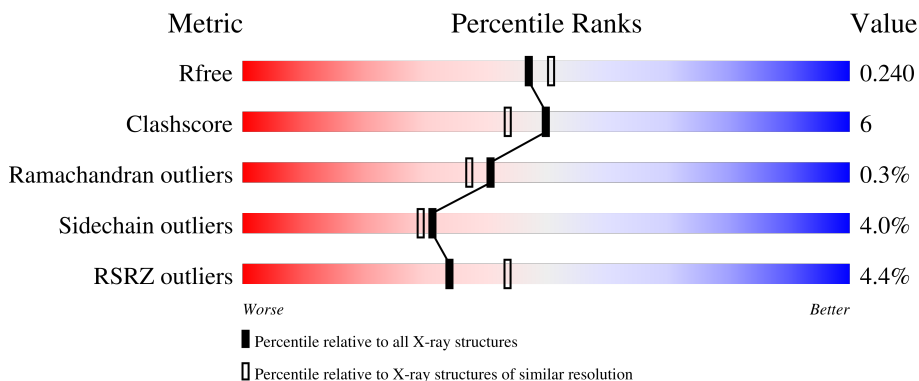
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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.



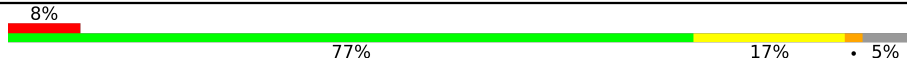


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	309	
1	B	309	
1	C	309	
1	D	309	
1	E	309	

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Mol	Chain	Length	Quality of chain
1	F	309	
1	G	309	
1	H	309	

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 19289 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 Cif-like 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	293	2308	1479	411	408	10	0	0	0
1	B	294	2317	1484	412	411	10	0	0	0
1	C	293	2310	1479	411	410	10	0	0	0
1	D	294	2317	1484	412	411	10	0	0	0
1	E	294	2317	1484	412	411	10	0	0	0
1	F	293	2308	1479	411	408	10	0	0	0
1	G	293	2308	1479	411	408	10	0	0	0
1	H	294	2317	1484	412	411	10	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	SER	ASP	engineered mutation	UNP B4EJL9
B	123	SER	ASP	engineered mutation	UNP B4EJL9
C	123	SER	ASP	engineered mutation	UNP B4EJL9
D	123	SER	ASP	engineered mutation	UNP B4EJL9
E	123	SER	ASP	engineered mutation	UNP B4EJL9
F	123	SER	ASP	engineered mutation	UNP B4EJL9
G	123	SER	ASP	engineered mutation	UNP B4EJL9
H	123	SER	ASP	engineered mutation	UNP B4EJL9

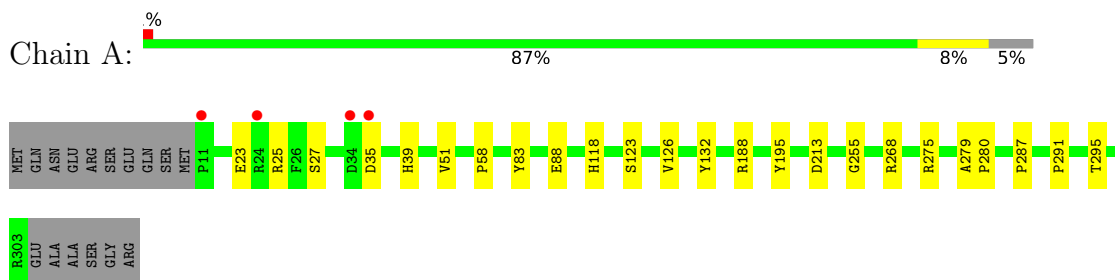
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	183	Total 183	O 183	0	0
2	B	135	Total 135	O 135	0	0
2	C	118	Total 118	O 118	0	0
2	D	94	Total 94	O 94	0	0
2	E	87	Total 87	O 87	0	0
2	F	57	Total 57	O 57	0	0
2	G	64	Total 64	O 64	0	0
2	H	49	Total 49	O 49	0	0

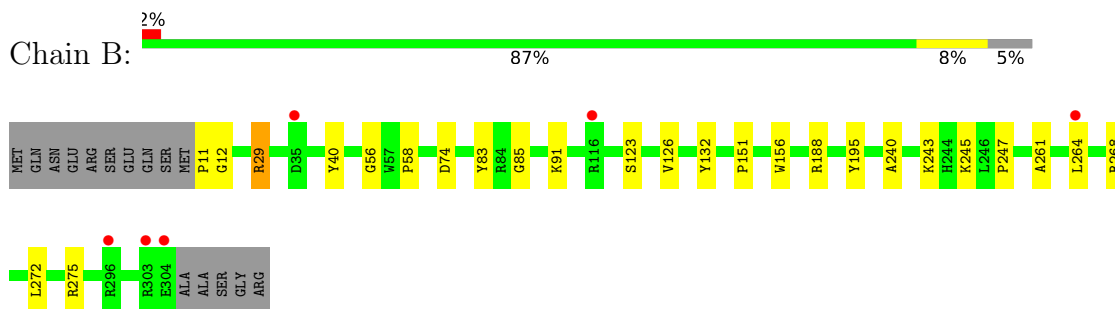
3 Residue-property plots [i](#)

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.

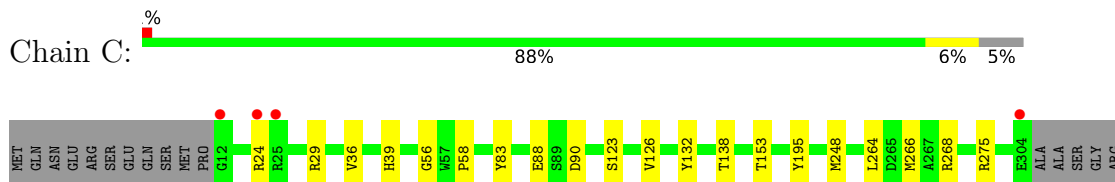
- Molecule 1: Cif-like 1



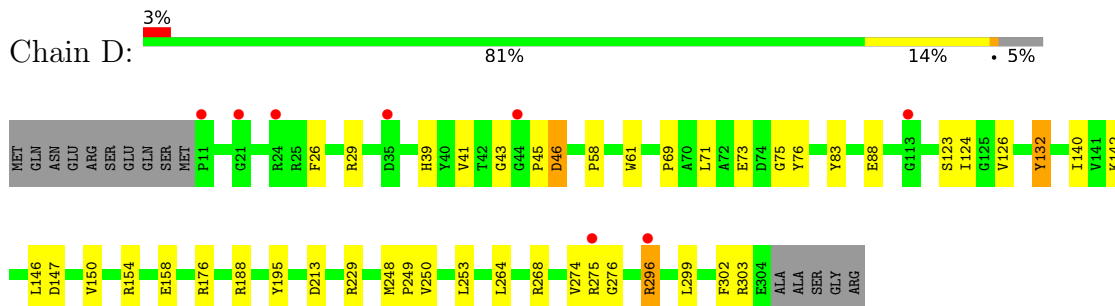
- Molecule 1: Cif-like 1



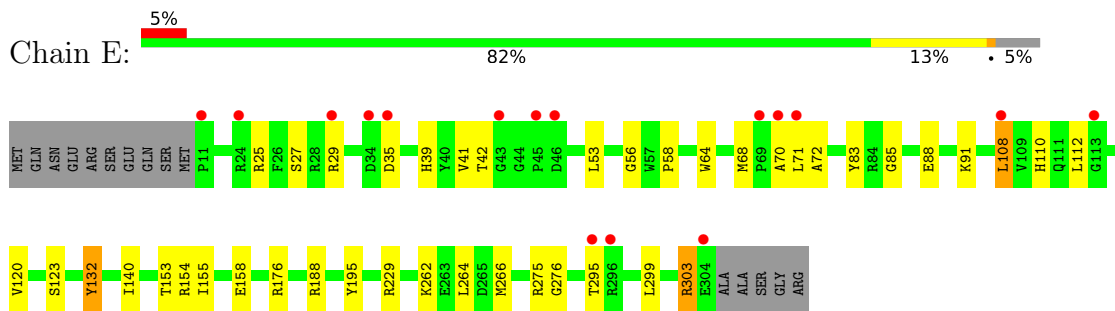
- Molecule 1: Cif-like 1



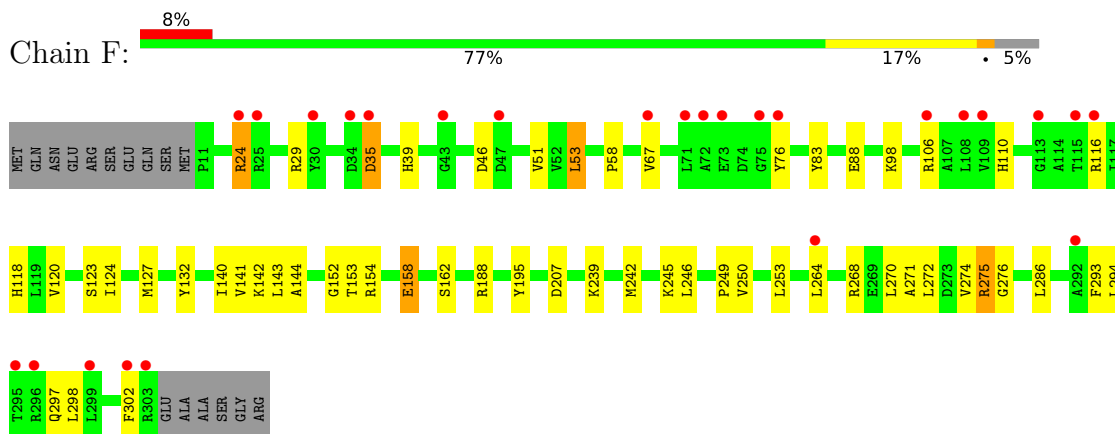
- Molecule 1: Cif-like 1



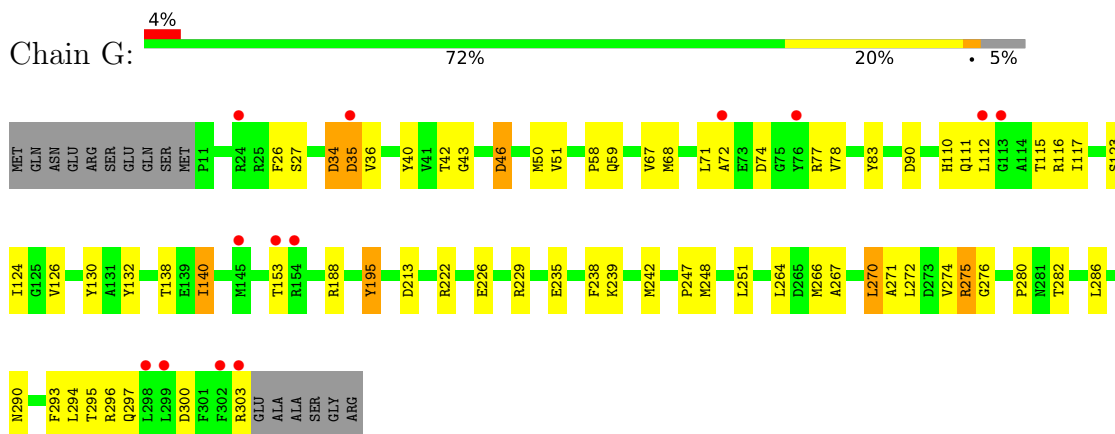
• Molecule 1: Cif-like 1



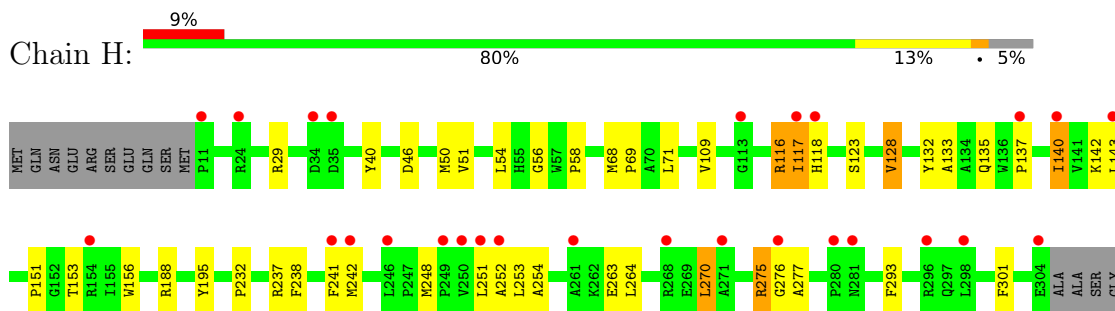
• Molecule 1: Cif-like 1



• Molecule 1: Cif-like 1



• Molecule 1: Cif-like 1



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	195.00Å 98.40Å 170.00Å 90.00° 118.60° 90.00°	Depositor
Resolution (Å)	44.71 – 2.15 48.47 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.71-2.15) 99.5 (48.47-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.16Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.209 , 0.240 0.209 , 0.240	Depositor DCC
R_{free} test set	7620 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtrriage
Anisotropy	0.586	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19289	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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, 0.333 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	0.38	0/2374	0.55	0/3225
1	B	0.41	0/2383	0.55	0/3237
1	C	0.41	0/2375	0.55	0/3226
1	D	0.55	0/2383	0.62	0/3237
1	E	0.53	0/2383	0.62	0/3237
1	F	0.61	0/2374	0.65	0/3225
1	G	0.62	0/2374	0.75	0/3225
1	H	0.60	0/2383	0.63	0/3237
All	All	0.52	0/19029	0.62	0/25849

There are no bond length outliers.

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	2308	0	2265	13	0
1	B	2317	0	2271	12	0
1	C	2310	0	2263	10	0
1	D	2317	0	2271	30	0
1	E	2317	0	2271	22	0
1	F	2308	0	2265	36	0
1	G	2308	0	2265	57	0
1	H	2317	0	2271	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	183	0	0	3	0
2	B	135	0	0	2	0
2	C	118	0	0	1	0
2	D	94	0	0	2	0
2	E	87	0	0	0	0
2	F	57	0	0	2	0
2	G	64	0	0	0	0
2	H	49	0	0	1	0
All	All	19289	0	18142	212	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 (212) 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:G:50:MET:HB3	1:G:117:ILE:HG22	1.47	0.96
1:G:50:MET:CE	1:G:117:ILE:HG21	1.98	0.94
1:H:116:ARG:HG2	1:H:116:ARG:HH21	1.35	0.91
1:G:50:MET:HE3	1:G:117:ILE:HG21	1.57	0.85
1:D:142:LYS:HG2	1:D:249:PRO:HB2	1.59	0.84
1:F:275:ARG:HH11	1:F:275:ARG:HB3	1.54	0.73
1:G:50:MET:HB3	1:G:117:ILE:CG2	2.19	0.72
1:G:51:VAL:HG13	1:G:51:VAL:O	1.90	0.71
1:G:271:ALA:O	1:G:274:VAL:CG2	2.39	0.70
1:H:264:LEU:HD21	1:H:276:GLY:HA3	1.73	0.70
1:H:116:ARG:HH21	1:H:116:ARG:CG	2.03	0.69
1:C:138:THR:HG21	1:G:300:ASP:HB3	1.72	0.69
1:F:35:ASP:OD2	1:F:35:ASP:N	2.23	0.69
1:F:132:TYR:CD1	1:F:140:ILE:HD11	2.28	0.69
1:G:50:MET:HE2	1:G:117:ILE:HG21	1.72	0.69
1:D:69:PRO:O	1:D:73:GLU:HG3	1.93	0.68
1:E:53:LEU:HD23	1:E:120:VAL:HB	1.75	0.67
1:H:140:ILE:HD11	1:H:143:LEU:HB2	1.78	0.66
1:D:250:VAL:HB	1:D:274:VAL:HG12	1.78	0.66
1:E:153:THR:HG21	1:E:266:MET:HG2	1.79	0.65
1:E:64:TRP:O	1:E:68:MET:HG3	1.97	0.65
1:G:35:ASP:HB2	1:G:36:VAL:HG23	1.77	0.65
1:H:140:ILE:CD1	1:H:143:LEU:HB2	2.27	0.65
1:H:140:ILE:HD11	1:H:142:LYS:C	2.18	0.64
1:H:251:LEU:HD12	1:H:275:ARG:HG2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:264:LEU:CD2	1:H:276:GLY:HA3	2.29	0.63
1:G:271:ALA:O	1:G:274:VAL:HG23	2.00	0.62
1:F:294:LEU:O	1:F:298:LEU:HG	1.99	0.62
1:G:50:MET:HE2	1:G:117:ILE:CG2	2.29	0.62
1:E:25:ARG:NH1	1:E:72:ALA:HB3	2.14	0.62
1:G:46:ASP:OD2	1:G:46:ASP:N	2.32	0.61
1:F:253:LEU:HD22	1:F:293:PHE:HE1	1.65	0.61
1:G:72:ALA:HB2	1:G:78:VAL:HG23	1.82	0.61
1:D:248:MET:HB2	1:D:249:PRO:CD	2.29	0.61
1:H:50:MET:HB3	1:H:117:ILE:HG22	1.83	0.61
1:H:153:THR:HG22	1:H:242:MET:HE3	1.82	0.61
1:D:46:ASP:N	1:D:46:ASP:OD1	2.33	0.60
1:G:275:ARG:HB3	1:G:275:ARG:CZ	2.31	0.60
1:D:264:LEU:HD13	1:D:276:GLY:HA3	1.85	0.59
1:G:188:ARG:HH21	1:G:213:ASP:CG	2.06	0.59
1:G:282:THR:HG21	1:G:290:ASN:HB2	1.85	0.59
1:F:53:LEU:HD22	1:F:53:LEU:N	2.18	0.58
1:H:254:ALA:HB2	1:H:263:GLU:HG3	1.84	0.58
1:C:123:SER:O	1:C:126:VAL:HG12	2.04	0.58
1:C:138:THR:HG21	1:G:300:ASP:CB	2.34	0.58
1:A:35:ASP:OD2	1:A:35:ASP:N	2.37	0.57
1:C:153:THR:HG21	1:C:266:MET:HG2	1.87	0.57
1:G:115:THR:O	1:G:115:THR:OG1	2.22	0.57
1:F:268:ARG:HG2	1:F:274:VAL:HG21	1.85	0.57
1:D:268:ARG:HG2	1:D:274:VAL:HG21	1.86	0.56
1:D:248:MET:HB2	1:D:249:PRO:HD2	1.86	0.56
1:G:267:ALA:HA	1:G:270:LEU:HD12	1.88	0.56
1:G:35:ASP:N	1:G:35:ASP:OD1	2.38	0.55
1:G:68:MET:HG2	1:G:78:VAL:HG11	1.88	0.55
1:H:238:PHE:O	1:H:242:MET:HG3	2.06	0.55
1:H:140:ILE:HD13	1:H:248:MET:HG3	1.89	0.55
1:H:56:GLY:HA3	1:H:123:SER:HB3	1.88	0.55
1:C:56:GLY:HA3	1:C:123:SER:HB3	1.89	0.54
1:F:246:LEU:HB2	1:F:271:ALA:HB2	1.89	0.54
1:G:153:THR:HG21	1:G:266:MET:HG2	1.90	0.54
1:H:54:LEU:HD13	1:H:128:VAL:CG2	2.37	0.54
1:D:154:ARG:O	1:D:158:GLU:HG3	2.08	0.54
1:D:296:ARG:HA	1:D:296:ARG:HE	1.73	0.53
1:G:27:SER:N	1:G:42:THR:O	2.35	0.53
1:H:253:LEU:HD22	1:H:277:ALA:HB3	1.91	0.53
1:G:117:ILE:C	1:G:117:ILE:HD12	2.29	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:296:ARG:HH21	1:D:299:LEU:HB2	1.73	0.53
1:F:153:THR:HG22	1:F:242:MET:HE1	1.89	0.53
1:C:138:THR:HG21	1:G:300:ASP:CG	2.30	0.52
1:F:132:TYR:CE1	1:F:140:ILE:HD11	2.45	0.52
1:A:27:SER:OG	1:B:11:PRO:HA	2.09	0.52
1:G:293:PHE:CE1	1:G:297:GLN:HG3	2.45	0.52
1:E:27:SER:O	1:E:41:VAL:HA	2.09	0.52
1:E:25:ARG:HH11	1:E:72:ALA:HB3	1.74	0.52
1:A:23:GLU:O	1:B:12:GLY:HA2	2.09	0.52
1:D:264:LEU:O	1:D:268:ARG:HG3	2.09	0.51
1:A:268:ARG:NE	2:A:403:HOH:O	2.29	0.51
1:E:70:ALA:CB	1:E:295:THR:HG21	2.40	0.51
1:H:252:ALA:O	1:H:253:LEU:HD23	2.11	0.51
1:B:245:LYS:HE3	1:B:272:LEU:HD23	1.93	0.51
1:B:56:GLY:HA3	1:B:123:SER:HB3	1.93	0.51
1:E:154:ARG:O	1:E:158:GLU:HG3	2.11	0.51
1:F:120:VAL:HG22	1:F:144:ALA:HB3	1.92	0.51
1:G:188:ARG:NH2	1:G:213:ASP:OD2	2.44	0.50
1:G:238:PHE:O	1:G:242:MET:HG2	2.12	0.50
1:H:109:VAL:HG11	1:H:117:ILE:HD13	1.94	0.50
1:F:253:LEU:CD2	1:F:293:PHE:HE1	2.24	0.50
1:H:142:LYS:HB3	1:H:301:PHE:CE2	2.47	0.50
1:B:261:ALA:O	1:B:264:LEU:HB2	2.13	0.49
1:F:275:ARG:HB3	1:F:275:ARG:NH1	2.23	0.49
1:F:293:PHE:C	1:F:293:PHE:CD1	2.85	0.49
1:G:51:VAL:O	1:G:51:VAL:CG1	2.60	0.49
1:D:126:VAL:HG13	1:D:150:VAL:HG12	1.94	0.49
1:D:123:SER:O	1:D:126:VAL:HG12	2.12	0.49
1:F:297:GLN:NE2	2:F:405:HOH:O	2.35	0.49
1:G:123:SER:O	1:G:126:VAL:HG12	2.12	0.49
1:G:251:LEU:HD13	1:G:275:ARG:HD3	1.93	0.49
1:D:41:VAL:HG21	1:D:61:TRP:CE2	2.48	0.49
1:D:274:VAL:HG22	2:D:402:HOH:O	2.13	0.48
1:E:70:ALA:HB3	1:E:295:THR:HG21	1.94	0.48
1:G:280:PRO:O	1:G:282:THR:HG23	2.13	0.48
1:D:248:MET:CB	1:D:249:PRO:CD	2.91	0.48
1:E:29:ARG:HD2	1:E:42:THR:CG2	2.44	0.48
1:G:72:ALA:HB2	1:G:78:VAL:CG2	2.43	0.48
1:F:264:LEU:HD13	1:F:276:GLY:HA3	1.96	0.47
1:G:248:MET:H	1:G:248:MET:HE3	1.78	0.47
1:D:71:LEU:O	1:D:76:TYR:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:29:ARG:HD2	1:E:42:THR:HG21	1.97	0.47
1:E:39:HIS:CG	1:E:88:GLU:HB2	2.49	0.47
1:G:40:TYR:OH	1:G:112:LEU:HD21	2.15	0.47
1:A:291:PRO:O	1:A:295:THR:HG23	2.15	0.47
1:E:85:GLY:HA2	1:E:91:LYS:HG3	1.96	0.47
1:E:155:ILE:HD12	1:E:262:LYS:HG2	1.96	0.47
1:F:207:ASP:OD2	2:F:401:HOH:O	2.21	0.46
1:G:235:GLU:O	1:G:239:LYS:HG3	2.14	0.46
1:A:287:PRO:HD2	2:A:478:HOH:O	2.15	0.46
1:D:45:PRO:HG2	1:D:75:GLY:HA2	1.97	0.46
1:D:123:SER:OG	1:D:124:ILE:N	2.48	0.46
1:F:106:ARG:HA	1:F:132:TYR:OH	2.16	0.46
1:G:117:ILE:HD12	1:G:117:ILE:O	2.16	0.46
1:D:76:TYR:OH	1:D:302:PHE:HB3	2.14	0.46
1:H:151:PRO:HA	1:H:156:TRP:CG	2.51	0.46
1:H:237:ARG:NH2	1:H:241:PHE:HZ	2.14	0.46
1:A:23:GLU:HB2	2:A:472:HOH:O	2.16	0.46
1:E:299:LEU:O	1:E:303:ARG:HD3	2.16	0.46
1:G:123:SER:OG	1:G:124:ILE:N	2.48	0.46
1:A:188:ARG:HH21	1:A:213:ASP:CG	2.20	0.45
1:E:56:GLY:HA3	1:E:123:SER:HB3	1.98	0.45
1:G:50:MET:HA	1:G:77:ARG:HB3	1.97	0.45
1:H:71:LEU:HD23	1:H:71:LEU:HA	1.67	0.45
1:F:154:ARG:O	1:F:158:GLU:HB2	2.17	0.45
1:C:264:LEU:HD21	1:C:268:ARG:HH21	1.82	0.45
1:A:123:SER:O	1:A:126:VAL:HG12	2.16	0.45
1:B:268:ARG:NH2	2:B:403:HOH:O	2.33	0.45
1:H:51:VAL:HA	1:H:118:HIS:O	2.17	0.45
1:F:51:VAL:HG23	1:F:76:TYR:HB3	1.99	0.45
1:G:264:LEU:HD13	1:G:276:GLY:HA3	1.99	0.45
1:G:275:ARG:HB3	1:G:275:ARG:NH2	2.32	0.45
1:H:133:ALA:O	1:H:137:PRO:HG3	2.17	0.45
1:F:264:LEU:O	1:F:268:ARG:HG3	2.17	0.44
1:G:140:ILE:HG22	1:G:248:MET:SD	2.57	0.44
1:H:140:ILE:HD11	1:H:143:LEU:CB	2.47	0.44
1:H:264:LEU:CG	1:H:276:GLY:HA3	2.47	0.44
1:E:264:LEU:HD13	1:E:276:GLY:HA3	2.00	0.44
1:A:39:HIS:CG	1:A:88:GLU:HB2	2.52	0.44
1:B:123:SER:O	1:B:126:VAL:HG12	2.18	0.44
1:H:135:GLN:OE1	1:H:237:ARG:NH2	2.47	0.44
1:A:255:GLY:HA2	1:A:279:ALA:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:147:ASP:OD1	2:D:401:HOH:O	2.21	0.44
1:D:188:ARG:HH21	1:D:213:ASP:CG	2.21	0.44
1:E:176:ARG:HA	1:E:229:ARG:CZ	2.48	0.44
1:C:36:VAL:HG23	1:C:90:ASP:HB2	2.00	0.44
1:D:296:ARG:HE	1:D:296:ARG:CA	2.30	0.44
1:F:123:SER:OG	1:F:124:ILE:N	2.49	0.44
1:F:143:LEU:O	1:F:250:VAL:HA	2.18	0.44
1:F:242:MET:HE2	1:F:270:LEU:HD11	2.00	0.44
1:D:39:HIS:CG	1:D:88:GLU:HB2	2.53	0.43
1:F:294:LEU:HD11	1:F:298:LEU:HD21	2.00	0.43
1:G:222:ARG:O	1:G:226:GLU:HG2	2.18	0.43
1:G:272:LEU:HA	1:G:272:LEU:HD23	1.56	0.43
1:D:132:TYR:CD1	1:D:140:ILE:HD11	2.53	0.43
1:H:29:ARG:HB2	1:H:40:TYR:CZ	2.53	0.43
1:G:67:VAL:HG21	1:G:294:LEU:HD23	2.01	0.43
1:H:253:LEU:HD13	1:H:293:PHE:HE1	1.83	0.43
1:H:40:TYR:OH	2:H:401:HOH:O	2.21	0.43
1:H:116:ARG:HG2	1:H:116:ARG:NH2	2.13	0.43
1:F:245:LYS:HD2	1:F:272:LEU:HG	2.00	0.43
1:F:249:PRO:HB3	1:F:275:ARG:NH2	2.34	0.43
1:C:29:ARG:NE	2:C:406:HOH:O	2.46	0.43
1:E:108:LEU:HD11	1:E:112:LEU:HD11	2.01	0.43
1:G:117:ILE:C	1:G:117:ILE:CD1	2.87	0.43
1:C:39:HIS:CG	1:C:88:GLU:HB2	2.53	0.42
1:B:240:ALA:O	1:B:243:LYS:HG2	2.18	0.42
1:G:247:PRO:HD2	1:G:248:MET:HE3	2.01	0.42
1:G:26:PHE:CD1	1:G:43:GLY:HA3	2.55	0.42
1:D:188:ARG:NH2	1:D:213:ASP:OD2	2.50	0.42
1:G:34:ASP:OD2	1:G:34:ASP:N	2.51	0.42
1:E:132:TYR:CD1	1:E:140:ILE:HD11	2.54	0.42
1:H:242:MET:HG2	1:H:270:LEU:HD23	2.01	0.42
1:G:71:LEU:HD23	1:G:71:LEU:HA	1.70	0.42
1:G:226:GLU:HA	1:G:229:ARG:HD2	2.02	0.42
1:F:39:HIS:CG	1:F:88:GLU:HB2	2.54	0.42
1:F:152:GLY:O	1:F:239:LYS:HE3	2.19	0.42
1:D:176:ARG:HA	1:D:229:ARG:CZ	2.50	0.42
1:G:36:VAL:HG13	1:G:90:ASP:HB2	2.00	0.42
1:G:59:GLN:NE2	1:G:195:TYR:OH	2.38	0.42
1:F:118:HIS:HB3	1:F:302:PHE:CZ	2.55	0.41
1:E:71:LEU:HD23	1:E:71:LEU:HA	1.72	0.41
1:F:158:GLU:O	1:F:162:SER:OG	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:PRO:HA	1:B:156:TRP:CG	2.55	0.41
1:B:85:GLY:HA2	1:B:91:LYS:HG3	2.03	0.41
1:D:26:PHE:CD1	1:D:43:GLY:HA3	2.56	0.41
1:G:67:VAL:HA	1:G:295:THR:CG2	2.51	0.41
1:B:247:PRO:O	2:B:401:HOH:O	2.21	0.41
1:E:108:LEU:CD1	1:E:112:LEU:HD11	2.51	0.41
1:F:98:LYS:HE2	1:F:127:MET:HE2	2.03	0.41
1:F:141:VAL:HG12	1:F:142:LYS:HG3	2.02	0.41
1:F:24:ARG:H	1:F:24:ARG:HG2	1.49	0.41
1:G:130:TYR:CE1	1:G:270:LEU:HD22	2.56	0.41
1:H:140:ILE:HD12	1:H:143:LEU:HB2	2.00	0.41
1:H:264:LEU:HD21	1:H:276:GLY:CA	2.46	0.41
1:D:146:LEU:HD22	1:D:253:LEU:HB2	2.02	0.40
1:F:245:LYS:HD3	1:F:271:ALA:C	2.42	0.40
1:H:68:MET:N	1:H:69:PRO:CD	2.84	0.40
1:A:51:VAL:HG22	1:A:118:HIS:HB2	2.02	0.40
1:B:29:ARG:HB3	1:B:40:TYR:CZ	2.56	0.40
1:F:67:VAL:HG21	1:F:294:LEU:HD23	2.03	0.40
1:G:138:THR:O	1:G:138:THR:HG22	2.22	0.40
1:H:116:ARG:CG	1:H:116:ARG:NH2	2.72	0.40
1:A:279:ALA:HA	1:A:280:PRO:HD3	1.98	0.40
1:G:303:ARG:HD2	1:G:303:ARG:HA	1.81	0.40
1:H:140:ILE:HD11	1:H:142:LYS:O	2.20	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	291/309 (94%)	282 (97%)	8 (3%)	1 (0%)	41 37
1	B	292/309 (94%)	283 (97%)	8 (3%)	1 (0%)	41 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	291/309 (94%)	282 (97%)	8 (3%)	1 (0%)	41	37
1	D	292/309 (94%)	284 (97%)	7 (2%)	1 (0%)	41	37
1	E	292/309 (94%)	283 (97%)	8 (3%)	1 (0%)	41	37
1	F	291/309 (94%)	281 (97%)	9 (3%)	1 (0%)	41	37
1	G	291/309 (94%)	278 (96%)	12 (4%)	1 (0%)	41	37
1	H	292/309 (94%)	285 (98%)	6 (2%)	1 (0%)	41	37
All	All	2332/2472 (94%)	2258 (97%)	66 (3%)	8 (0%)	41	37

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	58	PRO
1	D	58	PRO
1	E	58	PRO
1	F	58	PRO
1	H	58	PRO
1	G	58	PRO
1	A	58	PRO
1	B	58	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	228/241 (95%)	223 (98%)	5 (2%)	52	55
1	B	229/241 (95%)	222 (97%)	7 (3%)	40	39
1	C	228/241 (95%)	222 (97%)	6 (3%)	46	47
1	D	229/241 (95%)	221 (96%)	8 (4%)	36	34
1	E	229/241 (95%)	220 (96%)	9 (4%)	32	30
1	F	228/241 (95%)	215 (94%)	13 (6%)	20	16
1	G	228/241 (95%)	213 (93%)	15 (7%)	16	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	229/241 (95%)	218 (95%)	11 (5%)	25	22
All	All	1828/1928 (95%)	1754 (96%)	74 (4%)	31	29

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	83	TYR
1	A	132	TYR
1	A	195	TYR
1	A	275	ARG
1	B	29	ARG
1	B	74	ASP
1	B	83	TYR
1	B	132	TYR
1	B	188	ARG
1	B	195	TYR
1	B	275	ARG
1	C	24	ARG
1	C	83	TYR
1	C	132	TYR
1	C	195	TYR
1	C	248	MET
1	C	275	ARG
1	D	29	ARG
1	D	46	ASP
1	D	83	TYR
1	D	132	TYR
1	D	195	TYR
1	D	275	ARG
1	D	296	ARG
1	D	303	ARG
1	E	35	ASP
1	E	83	TYR
1	E	108	LEU
1	E	110	HIS
1	E	132	TYR
1	E	188	ARG
1	E	195	TYR
1	E	275	ARG
1	E	303	ARG
1	F	24	ARG

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Mol	Chain	Res	Type
1	F	29	ARG
1	F	35	ASP
1	F	46	ASP
1	F	53	LEU
1	F	83	TYR
1	F	110	HIS
1	F	116	ARG
1	F	158	GLU
1	F	188	ARG
1	F	195	TYR
1	F	275	ARG
1	F	286	LEU
1	G	34	ASP
1	G	35	ASP
1	G	46	ASP
1	G	74	ASP
1	G	83	TYR
1	G	110	HIS
1	G	111	GLN
1	G	116	ARG
1	G	132	TYR
1	G	140	ILE
1	G	195	TYR
1	G	270	LEU
1	G	275	ARG
1	G	286	LEU
1	G	296	ARG
1	H	46	ASP
1	H	116	ARG
1	H	117	ILE
1	H	128	VAL
1	H	132	TYR
1	H	140	ILE
1	H	188	ARG
1	H	195	TYR
1	H	232	PRO
1	H	270	LEU
1	H	275	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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, 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	293/309 (94%)	0.18	4 (1%) 75 80	30, 38, 52, 68	0
1	B	294/309 (95%)	-0.05	6 (2%) 65 72	29, 43, 63, 77	0
1	C	293/309 (94%)	-0.27	4 (1%) 75 80	32, 43, 59, 79	0
1	D	294/309 (95%)	0.03	8 (2%) 54 63	36, 57, 80, 89	0
1	E	294/309 (95%)	0.24	16 (5%) 25 34	36, 58, 80, 93	0
1	F	293/309 (94%)	0.47	26 (8%) 9 14	35, 67, 88, 98	0
1	G	293/309 (94%)	0.59	13 (4%) 34 43	34, 78, 104, 112	0
1	H	294/309 (95%)	0.49	27 (9%) 9 13	38, 72, 96, 110	0
All	All	2348/2472 (94%)	0.21	104 (4%) 34 43	29, 54, 93, 112	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	35	ASP	5.0
1	F	73	GLU	4.2
1	E	35	ASP	4.2
1	B	35	ASP	4.1
1	H	241	PHE	4.0
1	E	24	ARG	4.0
1	E	296	ARG	3.9
1	F	25	ARG	3.8
1	F	47	ASP	3.8
1	G	113	GLY	3.8
1	F	34	ASP	3.6
1	C	304	GLU	3.5
1	F	113	GLY	3.5
1	H	276	GLY	3.5
1	H	296	ARG	3.4
1	H	11	PRO	3.4

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Mol	Chain	Res	Type	RSRZ
1	H	261	ALA	3.3
1	F	299	LEU	3.3
1	H	271	ALA	3.3
1	F	303	ARG	3.2
1	F	72	ALA	3.2
1	H	117	ILE	3.2
1	B	304	GLU	3.2
1	H	113	GLY	3.1
1	F	302	PHE	3.1
1	H	34	ASP	3.1
1	F	295	THR	3.1
1	H	154	ARG	3.1
1	H	298	LEU	3.1
1	H	242	MET	3.1
1	A	35	ASP	3.1
1	D	35	ASP	3.1
1	D	113	GLY	3.0
1	F	24	ARG	3.0
1	B	264	LEU	3.0
1	E	45	PRO	3.0
1	E	46	ASP	3.0
1	H	140	ILE	2.9
1	H	24	ARG	2.9
1	E	34	ASP	2.9
1	F	264	LEU	2.9
1	G	24	ARG	2.8
1	E	29	ARG	2.8
1	D	24	ARG	2.8
1	A	24	ARG	2.8
1	F	115	THR	2.8
1	D	21	GLY	2.7
1	B	116	ARG	2.7
1	H	246	LEU	2.7
1	E	304	GLU	2.7
1	F	292	ALA	2.7
1	G	303	ARG	2.7
1	D	44	GLY	2.6
1	E	113	GLY	2.6
1	F	43	GLY	2.6
1	H	281	ASN	2.6
1	F	67	VAL	2.5
1	E	43	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	71	LEU	2.5
1	E	11	PRO	2.5
1	F	71	LEU	2.4
1	H	304	GLU	2.4
1	G	35	ASP	2.4
1	F	108	LEU	2.4
1	H	250	VAL	2.4
1	G	153	THR	2.4
1	D	296	ARG	2.4
1	H	280	PRO	2.4
1	E	70	ALA	2.4
1	B	303	ARG	2.4
1	H	118	HIS	2.3
1	H	137	PRO	2.3
1	H	251	LEU	2.3
1	G	154	ARG	2.3
1	E	69	PRO	2.3
1	F	30	TYR	2.3
1	G	299	LEU	2.3
1	A	11	PRO	2.3
1	H	268	ARG	2.2
1	H	249	PRO	2.2
1	E	295	THR	2.2
1	F	116	ARG	2.2
1	F	35	ASP	2.2
1	C	12	GLY	2.2
1	G	298	LEU	2.2
1	F	296	ARG	2.2
1	E	108	LEU	2.2
1	G	302	PHE	2.2
1	B	296	ARG	2.2
1	F	106	ARG	2.1
1	G	112	LEU	2.1
1	H	143	LEU	2.1
1	G	76	TYR	2.1
1	G	72	ALA	2.1
1	H	252	ALA	2.1
1	F	109	VAL	2.1
1	F	76	TYR	2.1
1	D	11	PRO	2.1
1	C	25	ARG	2.1
1	C	24	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	145	MET	2.1
1	A	34	ASP	2.1
1	F	75	GLY	2.0
1	D	275	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.