



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:59 PM GMT

PDB ID : 4QL6
Title : Structure of *C. trachomatis* CT441
Authors : Kohlmann, F.; Hilgenfeld, R.; Hansen, G.
Deposited on : 2014-06-10
Resolution : 2.97 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

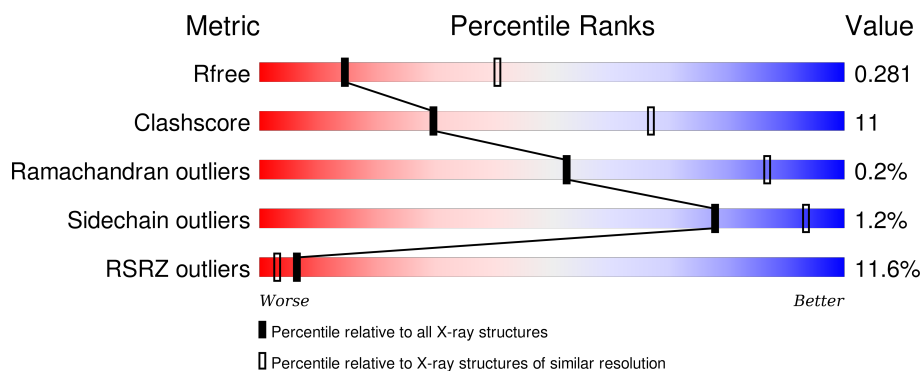
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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	634	<div> <div>8%</div> <div>59%</div> <div>18%</div> <div>23%</div> </div>
1	B	634	<div> <div>5%</div> <div>62%</div> <div>18%</div> <div>20%</div> </div>
1	C	634	<div> <div>15%</div> <div>50%</div> <div>27%</div> <div>23%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12071 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 Carboxy-terminal processing protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	0	0
			3970	2510	687	767	6			
1	B	509	Total	C	N	O	S	0	0	0
			4127	2607	715	798	7			
1	C	489	Total	C	N	O	S	0	0	0
			3974	2515	688	765	6			

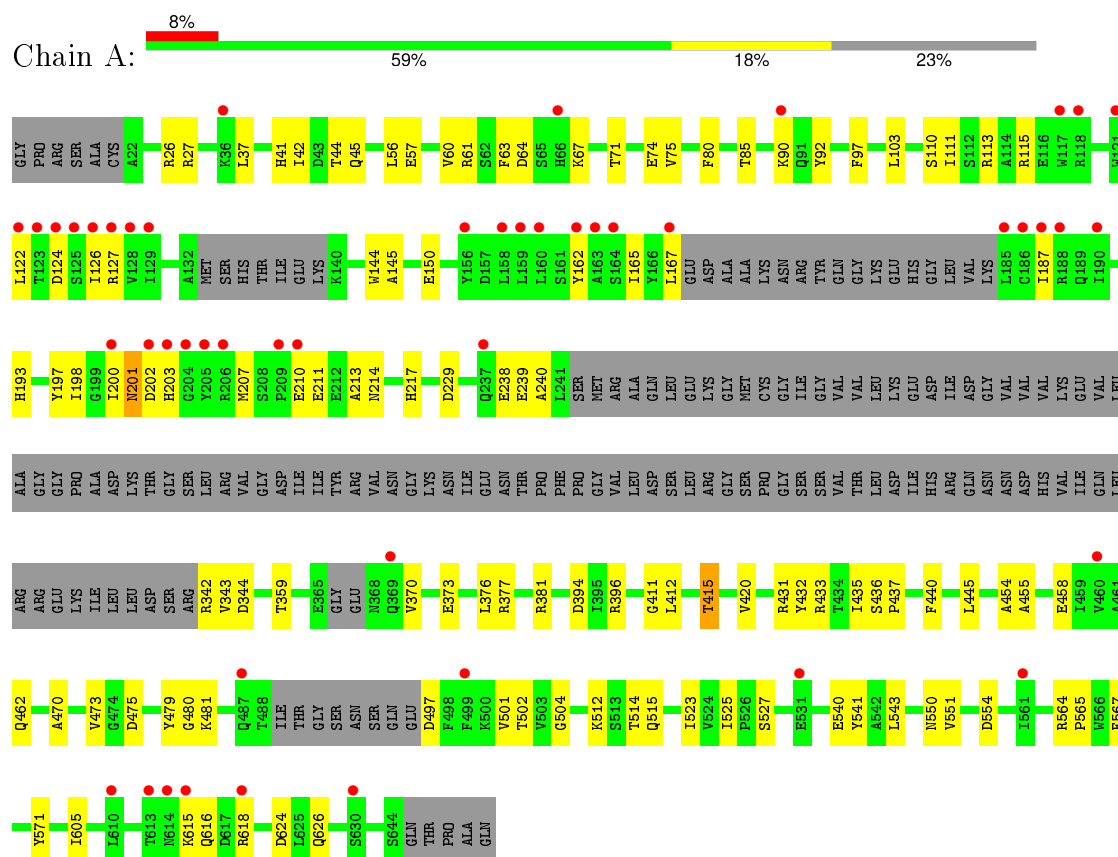
There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	-	EXPRESSION TAG	UNP B0B813
A	17	PRO	-	EXPRESSION TAG	UNP B0B813
A	18	ARG	-	EXPRESSION TAG	UNP B0B813
A	19	SER	-	EXPRESSION TAG	UNP B0B813
A	20	ALA	-	EXPRESSION TAG	UNP B0B813
A	21	CYS	-	EXPRESSION TAG	UNP B0B813
A	455	ALA	SER	ENGINEERED MUTATION	UNP B0B813
B	16	GLY	-	EXPRESSION TAG	UNP B0B813
B	17	PRO	-	EXPRESSION TAG	UNP B0B813
B	18	ARG	-	EXPRESSION TAG	UNP B0B813
B	19	SER	-	EXPRESSION TAG	UNP B0B813
B	20	ALA	-	EXPRESSION TAG	UNP B0B813
B	21	CYS	-	EXPRESSION TAG	UNP B0B813
B	455	ALA	SER	ENGINEERED MUTATION	UNP B0B813
C	16	GLY	-	EXPRESSION TAG	UNP B0B813
C	17	PRO	-	EXPRESSION TAG	UNP B0B813
C	18	ARG	-	EXPRESSION TAG	UNP B0B813
C	19	SER	-	EXPRESSION TAG	UNP B0B813
C	20	ALA	-	EXPRESSION TAG	UNP B0B813
C	21	CYS	-	EXPRESSION TAG	UNP B0B813
C	455	ALA	SER	ENGINEERED MUTATION	UNP B0B813

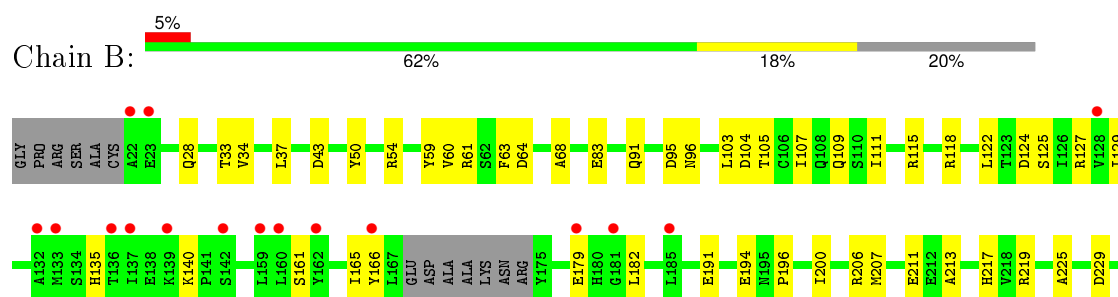
3 Residue-property plots

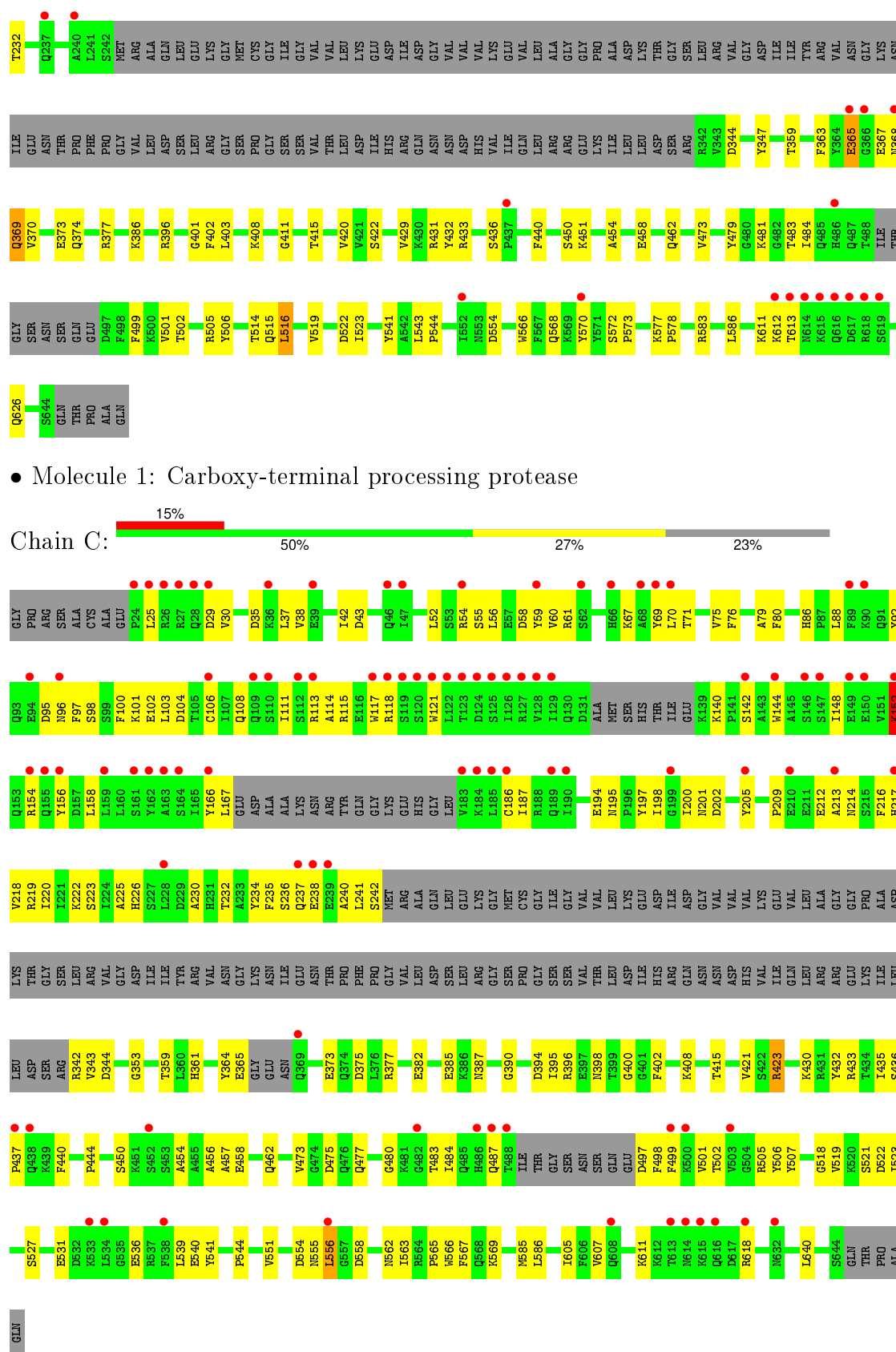
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Carboxy-terminal processing protease



• Molecule 1: Carboxy-terminal processing protease





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.73Å 183.99Å 209.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.44 – 2.97 33.44 – 2.97	Depositor EDS
% Data completeness (in resolution range)	91.5 (33.44-2.97) 91.5 (33.44-2.97)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.256 , 0.281 0.256 , 0.281	Depositor DCC
R_{free} test set	3235 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	85.9	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 47.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 64072 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12071	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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	0.26	0/4049	0.42	0/5469
1	B	0.26	0/4211	0.42	0/5687
1	C	0.26	0/4053	0.47	0/5471
All	All	0.26	0/12313	0.44	0/16627

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	3970	0	3884	67	0
1	B	4127	0	4040	73	0
1	C	3974	0	3903	130	0
All	All	12071	0	11827	268	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 11.

All (268) 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:C:67:LYS:HZ2	1:C:142:SER:H	1.23	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:235:PHE:HB2	1:C:498:PHE:HB3	1.58	0.85
1:A:111:ILE:HG22	1:A:115:ARG:HH12	1.48	0.78
1:C:42:ILE:HD11	1:C:505:ARG:HG3	1.69	0.75
1:C:423:ARG:CG	1:C:423:ARG:HH11	2.00	0.74
1:A:42:ILE:HD13	1:A:504:GLY:HA2	1.70	0.73
1:C:67:LYS:NZ	1:C:142:SER:H	1.86	0.72
1:B:129:ILE:HG13	1:B:179:GLU:HB3	1.71	0.72
1:C:423:ARG:HG3	1:C:423:ARG:HH11	1.55	0.71
1:B:363:PHE:HB2	1:B:401:GLY:HA3	1.72	0.71
1:A:41:HIS:O	1:A:45:GLN:NE2	2.24	0.70
1:C:195:ASN:HB3	1:C:200:ILE:HG13	1.75	0.69
1:C:121:TRP:HZ2	1:C:156:TYR:CD1	2.11	0.68
1:B:118:ARG:NH2	1:B:194:GLU:OE2	2.28	0.66
1:C:108:GLN:NE2	1:C:197:TYR:O	2.27	0.66
1:C:115:ARG:HD2	1:C:118:ARG:HH12	1.61	0.66
1:B:213:ALA:O	1:B:217:HIS:ND1	2.27	0.66
1:A:213:ALA:O	1:A:217:HIS:ND1	2.30	0.65
1:A:396:ARG:NH2	1:A:624:ASP:OD2	2.25	0.65
1:B:60:VAL:HG13	1:B:103:LEU:HD21	1.78	0.64
1:B:522:ASP:OD1	1:B:583:ARG:NH1	2.24	0.64
1:C:343:VAL:HG11	1:C:375:ASP:HB3	1.79	0.64
1:A:373:GLU:OE2	1:A:377:ARG:NH1	2.30	0.64
1:A:63:PHE:O	1:A:193:HIS:NE2	2.31	0.63
1:C:423:ARG:HH11	1:C:423:ARG:HB2	1.63	0.63
1:B:422:SER:HB2	1:B:429:VAL:HB	1.80	0.62
1:B:196:PRO:O	1:B:219:ARG:NH1	2.33	0.61
1:C:423:ARG:HH11	1:C:423:ARG:CB	2.14	0.60
1:C:536:GLU:HG3	1:C:539:LEU:HD12	1.82	0.60
1:A:210:GLU:O	1:A:214:ASN:N	2.33	0.60
1:B:422:SER:OG	1:B:431:ARG:NH1	2.35	0.59
1:B:568:GLN:HA	1:B:572:SER:HB2	1.84	0.59
1:B:91:GLN:HB3	1:B:96:ASN:HB3	1.83	0.59
1:A:238:GLU:OE1	1:A:238:GLU:N	2.36	0.59
1:C:556:LEU:HD13	1:C:567:PHE:HB2	1.85	0.59
1:C:225:ALA:HB1	1:C:232:THR:HG22	1.85	0.59
1:C:69:TYR:O	1:C:113:ARG:NH2	2.36	0.59
1:B:83:GLU:N	1:B:83:GLU:OE1	2.35	0.59
1:C:566:TRP:HE3	1:C:567:PHE:HD1	1.51	0.58
1:C:98:SER:HA	1:C:101:LYS:HG2	1.85	0.58
1:A:454:ALA:HA	1:A:458:GLU:HG3	1.86	0.58
1:C:423:ARG:NH1	1:C:423:ARG:HB2	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:ALA:HB2	1:B:479:TYR:HB3	1.86	0.57
1:C:67:LYS:HZ1	1:C:140:LYS:HB3	1.69	0.57
1:A:396:ARG:NH1	1:A:626:GLN:HB2	2.19	0.57
1:A:432:TYR:HE1	1:A:554:ASP:OD2	1.87	0.57
1:A:432:TYR:CE1	1:A:554:ASP:OD2	2.57	0.57
1:A:60:VAL:HG13	1:A:103:LEU:HD21	1.86	0.56
1:B:505:ARG:NH2	1:B:544:PRO:O	2.38	0.56
1:C:152:LYS:HG2	1:C:156:TYR:CE1	2.41	0.56
1:C:566:TRP:HE3	1:C:567:PHE:CD1	2.24	0.56
1:C:505:ARG:NH2	1:C:544:PRO:O	2.36	0.56
1:C:462:GLN:HB2	1:C:519:VAL:HG22	1.87	0.56
1:C:237:GLN:NE2	1:C:497:ASP:HB2	2.21	0.56
1:C:497:ASP:HB3	1:C:498:PHE:HB2	1.87	0.56
1:C:396:ARG:O	1:C:450:SER:OG	2.24	0.55
1:B:50:TYR:OH	1:B:54:ARG:NH1	2.38	0.55
1:B:577:LYS:HD2	1:B:578:PRO:HD2	1.89	0.55
1:B:566:TRP:CD2	1:C:566:TRP:HZ2	2.24	0.55
1:A:207:MET:HB2	1:A:211:GLU:HB3	1.88	0.55
1:B:111:ILE:HD13	1:B:194:GLU:HB3	1.88	0.55
1:C:230:ALA:N	1:C:541:TYR:O	2.40	0.55
1:C:483:THR:HB	1:C:502:THR:HG23	1.88	0.55
1:C:37:LEU:HD11	1:C:501:VAL:HB	1.88	0.55
1:A:567:PHE:HA	1:A:571:TYR:HB3	1.89	0.54
1:C:522:ASP:HB3	1:C:586:LEU:HD21	1.89	0.54
1:C:344:ASP:HB2	1:C:359:THR:HB	1.90	0.54
1:B:484:ILE:HG12	1:B:506:TYR:CE2	2.42	0.54
1:C:115:ARG:CD	1:C:118:ARG:HH12	2.20	0.54
1:C:118:ARG:HD2	1:C:187:ILE:HD11	1.89	0.54
1:C:71:THR:O	1:C:75:VAL:HG23	2.08	0.54
1:C:458:GLU:OE2	1:C:480:GLY:N	2.37	0.54
1:C:59:TYR:OH	1:C:104:ASP:OD2	2.25	0.54
1:C:114:ALA:O	1:C:118:ARG:HG3	2.07	0.54
1:B:225:ALA:HB1	1:B:232:THR:HG22	1.90	0.54
1:C:54:ARG:HD2	1:C:541:TYR:CE1	2.43	0.54
1:C:52:LEU:HD13	1:C:88:LEU:HB3	1.89	0.54
1:C:454:ALA:O	1:C:458:GLU:N	2.31	0.53
1:C:209:PRO:HA	1:C:212:GLU:HB3	1.89	0.53
1:B:232:THR:HG23	1:B:501:VAL:HG12	1.90	0.53
1:C:377:ARG:HA	1:C:440:PHE:HE2	1.73	0.53
1:C:433:ARG:NH1	1:C:435:ILE:HD13	2.24	0.53
1:A:71:THR:HG23	1:A:74:GLU:OE2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:TRP:CH2	1:C:156:TYR:HA	2.43	0.53
1:A:201:ASN:OD1	1:A:203:HIS:N	2.38	0.53
1:C:430:LYS:HD3	1:C:551:VAL:HG11	1.91	0.52
1:B:61:ARG:NH2	1:B:541:TYR:OH	2.43	0.52
1:C:421:VAL:O	1:C:507:TYR:N	2.43	0.52
1:C:111:ILE:HG23	1:C:194:GLU:OE2	2.10	0.52
1:C:121:TRP:HZ2	1:C:156:TYR:CE1	2.28	0.51
1:C:359:THR:HA	1:C:394:ASP:HB3	1.92	0.51
1:C:152:LYS:HG2	1:C:156:TYR:CZ	2.45	0.51
1:B:115:ARG:HH11	1:B:191:GLU:HG2	1.74	0.51
1:C:121:TRP:CZ2	1:C:156:TYR:CD1	2.98	0.51
1:C:531:GLU:OE2	1:C:611:LYS:HG2	2.11	0.51
1:C:104:ASP:OD2	1:C:216:PHE:HD1	1.93	0.51
1:B:207:MET:HB2	1:B:211:GLU:HB2	1.92	0.51
1:C:373:GLU:HG3	1:C:408:LYS:HB3	1.92	0.51
1:C:484:ILE:HG12	1:C:506:TYR:CZ	2.46	0.51
1:A:145:ALA:HB1	1:A:150:GLU:HB3	1.92	0.50
1:C:58:ASP:OD1	1:C:226:HIS:ND1	2.43	0.50
1:A:564:ARG:N	1:A:565:PRO:HD2	2.27	0.50
1:A:64:ASP:OD2	1:A:67:LYS:HA	2.12	0.50
1:C:52:LEU:HB3	1:C:88:LEU:HD23	1.94	0.50
1:C:217:HIS:O	1:C:220:ILE:HG22	2.12	0.50
1:C:432:TYR:HE2	1:C:554:ASP:OD1	1.95	0.49
1:A:497:ASP:OD1	1:A:497:ASP:N	2.45	0.49
1:C:102:GLU:O	1:C:106:CYS:N	2.41	0.49
1:A:122:LEU:HD13	1:A:187:ILE:HD13	1.93	0.49
1:B:568:GLN:O	1:B:573:PRO:HD3	2.11	0.49
1:B:566:TRP:CZ3	1:B:570:TYR:HB2	2.47	0.49
1:B:64:ASP:OD2	1:B:68:ALA:N	2.45	0.49
1:C:565:PRO:O	1:C:569:LYS:HG3	2.12	0.49
1:A:111:ILE:CG2	1:A:115:ARG:HH12	2.21	0.49
1:C:201:ASN:N	1:C:205:TYR:O	2.37	0.49
1:A:435:ILE:HG23	1:A:436:SER:O	2.11	0.49
1:A:61:ARG:NH2	1:A:541:TYR:OH	2.45	0.49
1:C:353:GLY:HA2	1:C:387:ASN:HB3	1.95	0.49
1:C:61:ARG:NH1	1:C:540:GLU:HG2	2.28	0.49
1:A:515:GLN:NE2	1:A:543:LEU:O	2.37	0.49
1:C:395:ILE:HD12	1:C:457:ALA:HB1	1.95	0.48
1:B:200:ILE:HA	1:B:206:ARG:HA	1.94	0.48
1:B:522:ASP:HB3	1:B:586:LEU:HD21	1.94	0.48
1:A:415:THR:OG1	1:A:437:PRO:O	2.30	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:ARG:NH2	1:A:370:VAL:HG11	2.28	0.48
1:C:166:TYR:HA	1:C:167:LEU:C	2.34	0.48
1:A:75:VAL:HG21	1:A:144:TRP:CD2	2.48	0.48
1:A:26:ARG:HH21	1:A:27:ARG:HH21	1.60	0.48
1:C:115:ARG:HA	1:C:118:ARG:NH1	2.28	0.48
1:C:117:TRP:CZ2	1:C:152:LYS:HG3	2.49	0.48
1:C:197:TYR:CE2	1:C:219:ARG:NH1	2.83	0.47
1:B:43:ASP:OD1	1:B:505:ARG:NH1	2.46	0.47
1:B:411:GLY:HA3	1:B:433:ARG:HE	1.79	0.47
1:C:200:ILE:HG22	1:C:201:ASN:O	2.14	0.47
1:B:368:ASN:HA	1:B:369:GLN:HA	1.50	0.47
1:A:540:GLU:OE2	1:A:540:GLU:N	2.45	0.47
1:C:37:LEU:HD13	1:C:499:PHE:CG	2.49	0.47
1:C:30:VAL:HG21	1:C:97:PHE:CZ	2.49	0.47
1:A:473:VAL:HG12	1:A:523:ILE:HB	1.97	0.47
1:B:34:VAL:O	1:B:37:LEU:HB3	2.15	0.47
1:A:165:ILE:O	1:A:167:LEU:HB2	2.14	0.47
1:B:462:GLN:HB2	1:B:519:VAL:HG22	1.96	0.47
1:B:347:TYR:OH	1:B:386:LYS:HE2	2.14	0.47
1:C:30:VAL:HG21	1:C:97:PHE:HZ	1.79	0.47
1:C:361:HIS:O	1:C:398:ASN:HA	2.15	0.47
1:C:76:PHE:CD1	1:C:80:PHE:HB2	2.50	0.47
1:B:516:LEU:HD12	1:B:516:LEU:H	1.79	0.47
1:B:450:SER:OG	1:B:451:LYS:N	2.48	0.47
1:C:556:LEU:HD13	1:C:567:PHE:CB	2.45	0.47
1:A:198:ILE:HG13	1:A:200:ILE:HG12	1.97	0.47
1:C:400:GLY:HA2	1:C:456:ALA:HB3	1.97	0.46
1:C:487:GLN:HB3	1:C:498:PHE:CZ	2.51	0.46
1:B:369:GLN:HB2	1:B:374:GLN:NE2	2.30	0.46
1:C:25:LEU:HA	1:C:29:ASP:OD2	2.15	0.46
1:B:611:LYS:O	1:B:612:LYS:HB3	2.15	0.46
1:C:423:ARG:HG3	1:C:423:ARG:NH1	2.26	0.46
1:C:98:SER:HA	1:C:101:LYS:NZ	2.30	0.46
1:C:222:LYS:HG2	1:C:234:TYR:HB3	1.97	0.46
1:B:566:TRP:CE3	1:C:566:TRP:HZ2	2.34	0.46
1:A:124:ASP:OD2	1:A:127:ARG:HB3	2.16	0.46
1:B:166:TYR:HD2	1:B:182:LEU:HD11	1.81	0.46
1:C:35:ASP:HA	1:C:38:VAL:HG22	1.97	0.46
1:C:56:LEU:HD21	1:C:79:ALA:HA	1.97	0.45
1:C:365:GLU:HB3	1:C:402:PHE:CD2	2.51	0.45
1:B:33:THR:HG22	1:B:499:PHE:HE2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:SER:O	1:C:223:SER:OG	2.24	0.45
1:C:214:ASN:O	1:C:218:VAL:N	2.37	0.45
1:B:229:ASP:OD2	1:B:502:THR:OG1	2.33	0.45
1:C:96:ASN:OD1	1:C:98:SER:N	2.36	0.45
1:B:59:TYR:OH	1:B:104:ASP:OD1	2.27	0.45
1:C:75:VAL:HG21	1:C:144:TRP:CD2	2.52	0.45
1:A:342:ARG:HB3	1:A:343:VAL:H	1.48	0.45
1:B:367:GLU:O	1:B:370:VAL:N	2.50	0.45
1:B:611:LYS:HB2	1:B:613:THR:HG22	1.98	0.44
1:A:381:ARG:HG3	1:A:440:PHE:HZ	1.82	0.44
1:A:56:LEU:HD23	1:A:85:THR:HG22	1.98	0.44
1:C:113:ARG:NH1	1:C:148:ILE:HD11	2.32	0.44
1:B:365:GLU:HB2	1:B:402:PHE:CD2	2.53	0.44
1:B:522:ASP:HB3	1:B:586:LEU:HD11	1.99	0.44
1:A:359:THR:HA	1:A:394:ASP:HB3	1.99	0.44
1:C:69:TYR:HB2	1:C:70:LEU:HD12	2.00	0.44
1:A:475:ASP:OD2	1:A:527:SER:OG	2.35	0.44
1:A:229:ASP:OD2	1:A:502:THR:HB	2.18	0.44
1:A:201:ASN:OD1	1:A:202:ASP:N	2.51	0.44
1:B:515:GLN:HA	1:B:516:LEU:HA	1.60	0.44
1:B:612:LYS:HG2	1:B:612:LYS:O	2.18	0.44
1:C:605:ILE:HD13	1:C:618:ARG:HB3	2.00	0.44
1:A:415:THR:O	1:A:435:ILE:HG22	2.18	0.44
1:C:342:ARG:HB3	1:C:364:TYR:OH	2.18	0.44
1:C:241:LEU:HD12	1:C:242:SER:N	2.32	0.44
1:A:239:GLU:HA	1:A:240:ALA:HA	1.67	0.43
1:C:201:ASN:OD1	1:C:202:ASP:N	2.51	0.43
1:C:566:TRP:CE3	1:C:567:PHE:CD1	3.06	0.43
1:A:162:TYR:O	1:A:165:ILE:HG22	2.18	0.43
1:A:420:VAL:HG13	1:A:431:ARG:HG3	2.00	0.43
1:B:373:GLU:O	1:B:377:ARG:HB2	2.17	0.43
1:A:605:ILE:HD13	1:A:618:ARG:HB3	2.00	0.43
1:A:445:LEU:HD23	1:A:470:ALA:HB2	2.00	0.43
1:B:403:LEU:HD11	1:B:420:VAL:HG11	2.01	0.43
1:C:115:ARG:NE	1:C:118:ARG:HH12	2.16	0.43
1:A:436:SER:OG	1:A:437:PRO:HD2	2.19	0.43
1:A:454:ALA:HB2	1:A:479:TYR:HB3	2.00	0.43
1:C:377:ARG:HA	1:C:440:PHE:CE2	2.54	0.43
1:C:475:ASP:OD2	1:C:527:SER:OG	2.37	0.42
1:C:585:MET:HB2	1:C:640:LEU:HD13	2.01	0.42
1:C:118:ARG:HG2	1:C:121:TRP:CZ3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:462:GLN:NE2	1:A:512:LYS:HG3	2.34	0.42
1:C:60:VAL:HG22	1:C:103:LEU:HD21	2.01	0.42
1:B:473:VAL:HG12	1:B:523:ILE:HB	2.01	0.42
1:A:376:LEU:HD23	1:A:412:LEU:HD12	2.02	0.42
1:B:122:LEU:O	1:B:125:SER:OG	2.22	0.42
1:C:236:SER:O	1:C:240:ALA:N	2.53	0.42
1:A:480:GLY:HA2	1:A:514:THR:O	2.19	0.42
1:A:63:PHE:HB2	1:A:197:TYR:CZ	2.55	0.42
1:B:28:GLN:HA	1:B:95:ASP:OD2	2.18	0.42
1:B:105:THR:O	1:B:109:GLN:HG3	2.19	0.42
1:B:115:ARG:CD	1:B:118:ARG:HH12	2.33	0.42
1:C:562:ASN:OD1	1:C:563:ILE:N	2.52	0.42
1:C:154:ARG:O	1:C:158:LEU:N	2.53	0.42
1:B:481:LYS:O	1:B:514:THR:OG1	2.37	0.42
1:C:436:SER:OG	1:C:437:PRO:HD2	2.20	0.42
1:A:455:ALA:HB2	1:A:481:LYS:HD3	2.02	0.42
1:C:195:ASN:HA	1:C:198:ILE:HD13	2.02	0.42
1:C:118:ARG:HA	1:C:121:TRP:CE3	2.55	0.42
1:B:483:THR:HB	1:B:502:THR:HG23	2.02	0.42
1:A:90:LYS:HD3	1:A:90:LYS:HA	1.88	0.42
1:A:525:ILE:HG22	1:A:626:GLN:HG3	2.03	0.41
1:C:555:ASN:C	1:C:556:LEU:HG	2.41	0.41
1:C:61:ARG:NH2	1:C:541:TYR:OH	2.42	0.41
1:B:408:LYS:HD3	1:B:433:ARG:HH22	1.85	0.41
1:C:477:GLN:HG3	1:C:518:GLY:H	1.85	0.41
1:A:92:TYR:HD1	1:A:97:PHE:HE1	1.68	0.41
1:B:140:LYS:HG2	1:B:140:LYS:H	1.64	0.41
1:B:124:ASP:OD2	1:B:127:ARG:HB3	2.21	0.41
1:C:432:TYR:OH	1:C:558:ASP:OD1	2.36	0.41
1:C:56:LEU:HA	1:C:100:PHE:HE1	1.86	0.41
1:C:382:GLU:OE2	1:C:385:GLU:OE2	2.39	0.41
1:A:37:LEU:HD12	1:A:501:VAL:HG22	2.03	0.41
1:C:415:THR:O	1:C:435:ILE:HG22	2.21	0.41
1:C:475:ASP:O	1:C:521:SER:OG	2.25	0.41
1:C:390:GLY:HA2	1:C:444:PRO:HG2	2.02	0.41
1:B:432:TYR:CE1	1:B:554:ASP:OD2	2.74	0.41
1:B:63:PHE:CZ	1:B:107:ILE:HG23	2.56	0.41
1:C:226:HIS:HD2	1:C:232:THR:O	2.03	0.41
1:B:454:ALA:HB2	1:B:479:TYR:CB	2.51	0.41
1:C:238:GLU:OE1	1:C:238:GLU:N	2.43	0.41
1:A:57:GLU:HG2	1:A:61:ARG:HE	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:ASN:OD1	1:A:551:VAL:N	2.54	0.41
1:A:411:GLY:HA3	1:A:433:ARG:HG2	2.03	0.41
1:B:161:SER:O	1:B:165:ILE:HG12	2.21	0.41
1:C:121:TRP:CZ2	1:C:156:TYR:CG	3.09	0.40
1:B:543:LEU:HA	1:B:544:PRO:HD3	1.93	0.40
1:A:75:VAL:HG21	1:A:144:TRP:CG	2.56	0.40
1:B:63:PHE:CE2	1:B:107:ILE:HG23	2.57	0.40
1:B:344:ASP:HB2	1:B:359:THR:HB	2.04	0.40
1:C:121:TRP:HZ2	1:C:156:TYR:CG	2.39	0.40
1:B:454:ALA:HA	1:B:458:GLU:HG3	2.03	0.40
1:C:88:LEU:O	1:C:92:TYR:N	2.46	0.40
1:B:37:LEU:HD13	1:B:499:PHE:CG	2.56	0.40
1:B:377:ARG:HG2	1:B:440:PHE:CZ	2.56	0.40
1:A:110:SER:HA	1:A:113:ARG:HB3	2.03	0.40
1:B:396:ARG:NH1	1:B:626:GLN:HB2	2.36	0.40
1:C:42:ILE:HG23	1:C:43:ASP:N	2.37	0.40
1:A:41:HIS:HB3	1:A:44:THR:O	2.22	0.40
1:C:607:VAL:O	1:C:611:LYS:HG3	2.22	0.40
1:A:344:ASP:HB2	1:A:359:THR:HB	2.03	0.40
1:C:209:PRO:O	1:C:213:ALA:N	2.55	0.40
1:C:473:VAL:HG12	1:C:523:ILE:HB	2.03	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	
1	A	477/634 (75%)	467 (98%)	9 (2%)	1 (0%)	52	87
1	B	501/634 (79%)	490 (98%)	11 (2%)	0	100	100
1	C	477/634 (75%)	463 (97%)	12 (2%)	2 (0%)	39	79
All	All	1455/1902 (76%)	1420 (98%)	32 (2%)	3 (0%)	52	87

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	PHE
1	C	152	LYS
1	C	86	HIS

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	
1	A	439/560 (78%)	434 (99%)	5 (1%)	80	94
1	B	456/560 (81%)	450 (99%)	6 (1%)	76	93
1	C	441/560 (79%)	436 (99%)	5 (1%)	80	94
All	All	1336/1680 (80%)	1320 (99%)	16 (1%)	78	93

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	ILE
1	A	201	ASN
1	A	415	THR
1	A	615	LYS
1	A	616	GLN
1	B	135	HIS
1	B	365	GLU
1	B	369	GLN
1	B	415	THR
1	B	436	SER
1	B	516	LEU
1	C	95	ASP
1	C	152	LYS
1	C	186	CYS
1	C	423	ARG
1	C	556	LEU

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

sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	HIS
1	A	588	GLN
1	B	374	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 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	489/634 (77%)	0.60	48 (9%) 10 5	50, 83, 156, 199	0
1	B	509/634 (80%)	0.42	33 (6%) 22 11	56, 76, 107, 126	0
1	C	489/634 (77%)	1.03	92 (18%) 2 1	58, 108, 205, 238	0
All	All	1487/1902 (78%)	0.68	173 (11%) 6 3	50, 86, 180, 238	0

All (173) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	184	LYS	12.4
1	A	129	ILE	11.2
1	C	123	THR	10.8
1	A	162	TYR	9.4
1	C	119	SER	9.0
1	C	120	SER	8.6
1	C	183	VAL	8.1
1	C	121	TRP	7.7
1	A	128	VAL	7.2
1	C	117	TRP	7.1
1	B	613	THR	6.4
1	B	614	ASN	6.3
1	C	110	SER	6.3
1	B	618	ARG	6.1
1	A	163	ALA	6.0
1	A	164	SER	5.6
1	B	142	SER	5.6
1	A	126	ILE	5.4
1	C	500	LYS	5.4
1	C	437	PRO	5.4
1	C	109	GLN	5.4
1	A	122	LEU	5.2
1	B	437	PRO	5.1

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Mol	Chain	Res	Type	RSRZ
1	C	486	HIS	5.0
1	A	185	LEU	5.0
1	A	187	ILE	5.0
1	B	133	MET	4.7
1	C	217	HIS	4.7
1	C	28	GLN	4.6
1	C	499	PHE	4.5
1	A	186	CYS	4.5
1	C	166	TYR	4.4
1	C	126	ILE	4.4
1	C	185	LEU	4.3
1	C	614	ASN	4.2
1	B	132	ALA	4.2
1	C	69	TYR	4.1
1	A	36	LYS	4.1
1	C	149	GLU	4.1
1	A	160	LEU	4.1
1	C	613	THR	4.1
1	B	166	TYR	4.0
1	A	66	HIS	4.0
1	C	46	GLN	4.0
1	A	613	THR	4.0
1	C	62	SER	3.9
1	B	615	LYS	3.9
1	B	616	GLN	3.9
1	A	159	LEU	3.9
1	C	128	VAL	3.9
1	C	162	TYR	3.8
1	C	113	ARG	3.8
1	A	209	PRO	3.8
1	C	142	SER	3.7
1	C	129	ILE	3.7
1	B	368	ASN	3.7
1	C	438	GLN	3.7
1	C	144	TRP	3.6
1	B	619	SER	3.6
1	A	90	LYS	3.6
1	C	616	GLN	3.6
1	B	612	LYS	3.5
1	B	137	ILE	3.5
1	C	112	SER	3.5
1	B	136	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	238	GLU	3.5
1	C	156	TYR	3.4
1	C	147	SER	3.4
1	C	89	PHE	3.4
1	C	39	GLU	3.3
1	B	22	ALA	3.3
1	C	190	ILE	3.3
1	C	106	CYS	3.2
1	B	617	ASP	3.2
1	A	206	ARG	3.2
1	B	128	VAL	3.2
1	C	124	ASP	3.2
1	C	25	LEU	3.2
1	C	205	TYR	3.1
1	C	163	ALA	3.1
1	C	186	CYS	3.1
1	A	210	GLU	3.1
1	A	156	TYR	3.1
1	C	26	ARG	3.0
1	C	533	LYS	3.0
1	A	118	ARG	3.0
1	B	139	LYS	3.0
1	C	96	ASN	3.0
1	A	203	HIS	3.0
1	B	570	TYR	3.0
1	C	27	ARG	3.0
1	C	189	GLN	3.0
1	C	118	ARG	3.0
1	A	200	ILE	2.9
1	C	199	GLY	2.9
1	C	239	GLU	2.9
1	C	213	ALA	2.9
1	C	164	SER	2.9
1	A	614	ASN	2.9
1	B	240	ALA	2.9
1	B	237	GLN	2.9
1	C	24	PRO	2.8
1	A	561	ILE	2.8
1	C	29	ASP	2.8
1	B	23	GLU	2.8
1	B	179	GLU	2.8
1	C	487	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	482	GLY	2.7
1	C	127	ARG	2.7
1	C	210	GLU	2.7
1	A	158	LEU	2.7
1	C	59	TYR	2.7
1	A	123	THR	2.7
1	C	122	LEU	2.7
1	A	237	GLN	2.7
1	C	161	SER	2.7
1	C	154	ARG	2.7
1	A	125	SER	2.7
1	C	632	ASN	2.6
1	C	538	PHE	2.6
1	C	228	LEU	2.6
1	A	205	TYR	2.6
1	A	127	ARG	2.6
1	C	150	GLU	2.6
1	A	610	LEU	2.6
1	C	70	LEU	2.6
1	A	615	LYS	2.5
1	B	552	ILE	2.5
1	C	452	SER	2.5
1	A	188	ARG	2.5
1	C	488	THR	2.5
1	A	460	VAL	2.5
1	A	487	GLN	2.5
1	C	68	ALA	2.5
1	B	366	GLY	2.5
1	C	125	SER	2.5
1	C	534	LEU	2.4
1	A	121	TRP	2.4
1	B	162	TYR	2.4
1	C	615	LYS	2.4
1	C	152	LYS	2.3
1	A	369	GLN	2.3
1	C	608	GLN	2.3
1	A	202	ASP	2.3
1	C	618	ARG	2.3
1	A	117	TRP	2.3
1	A	531	GLU	2.3
1	B	486	HIS	2.2
1	B	185	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	190	ILE	2.2
1	C	94	GLU	2.2
1	A	204	GLY	2.2
1	C	556	LEU	2.1
1	C	369	GLN	2.1
1	C	66	HIS	2.1
1	A	499	PHE	2.1
1	B	181	GLY	2.1
1	C	155	GLN	2.1
1	C	237	GLN	2.1
1	C	36	LYS	2.1
1	A	167	LEU	2.1
1	C	503	VAL	2.1
1	A	618	ARG	2.1
1	A	630	SER	2.1
1	B	160	LEU	2.1
1	C	159	LEU	2.1
1	C	54	ARG	2.1
1	C	47	ILE	2.0
1	B	159	LEU	2.0
1	C	146	SER	2.0
1	C	90	LYS	2.0
1	A	124	ASP	2.0
1	B	365	GLU	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.