



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 16, 2026 – 10:25 AM UTC

PDB ID : 3M6D / pdb_00003m6d
Title : The crystal structure of the d307a mutant of glycoside Hydrolase (family 31) from ruminococcus obeum atcc 29174
Authors : Tan, K.; Tesar, C.; Freeman, L.; Babnigg, G.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2010-03-15
Resolution : 2.90 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

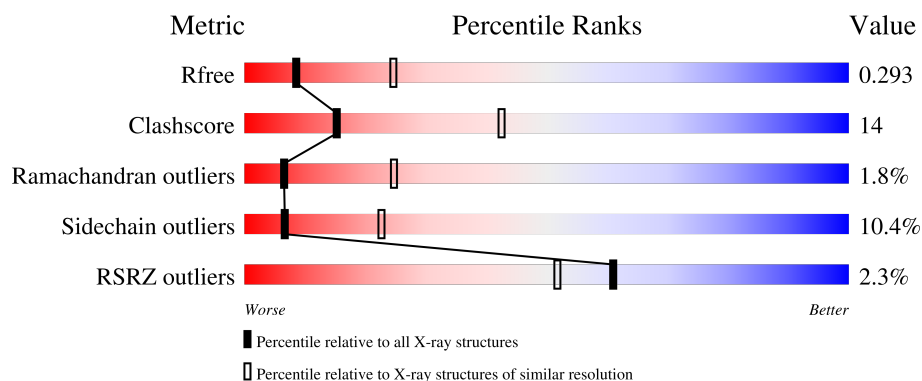
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	666	2% 62% 32% 5%
1	B	666	2% 65% 29% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10735 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	655	Total	C	N	O	S	0	0	0
			5357	3442	878	1003	34			
1	B	658	Total	C	N	O	S	0	0	0
			5378	3453	881	1009	35			

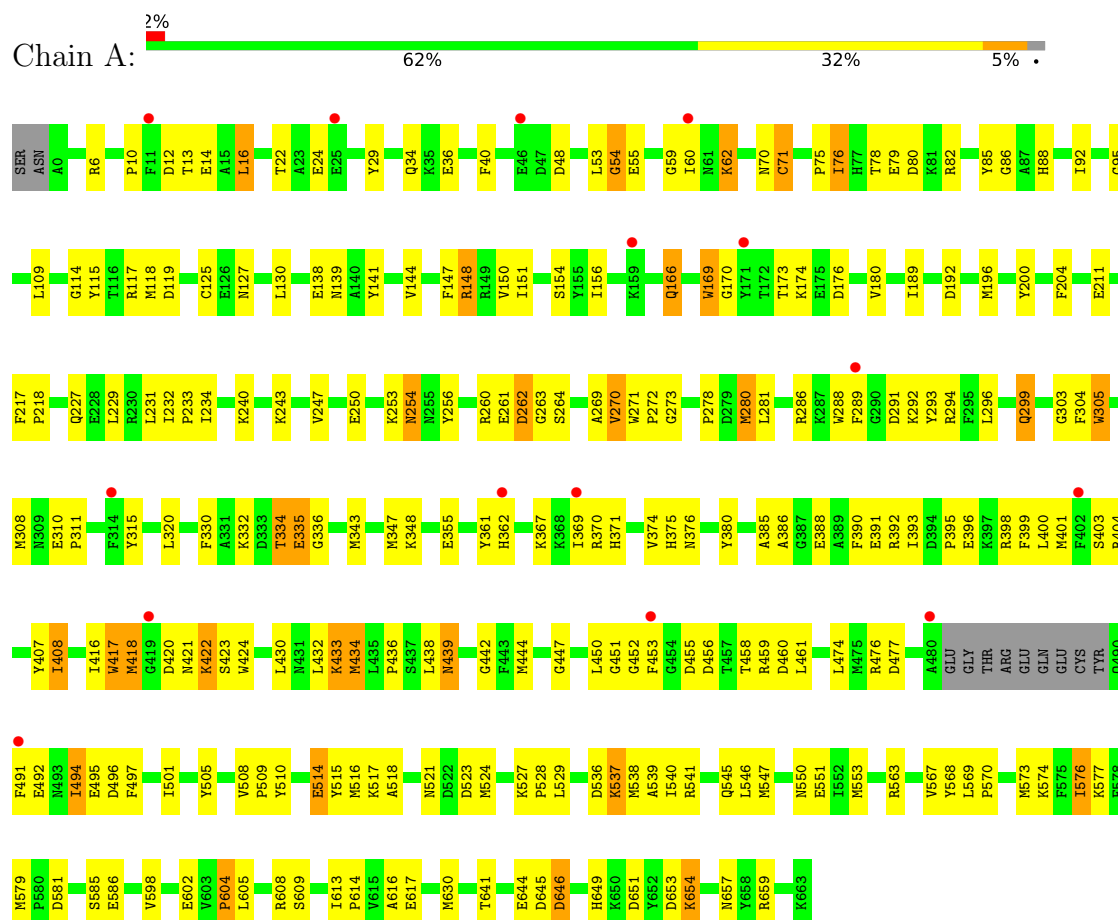
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A5ZY13
A	-1	ASN	-	expression tag	UNP A5ZY13
A	0	ALA	-	expression tag	UNP A5ZY13
A	307	ALA	ASP	engineered mutation	UNP A5ZY13
B	-2	SER	-	expression tag	UNP A5ZY13
B	-1	ASN	-	expression tag	UNP A5ZY13
B	0	ALA	-	expression tag	UNP A5ZY13
B	307	ALA	ASP	engineered mutation	UNP A5ZY13

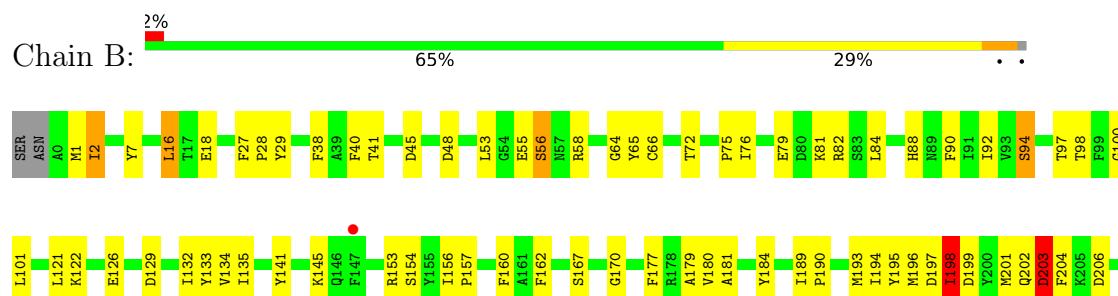
3 Residue-property plots

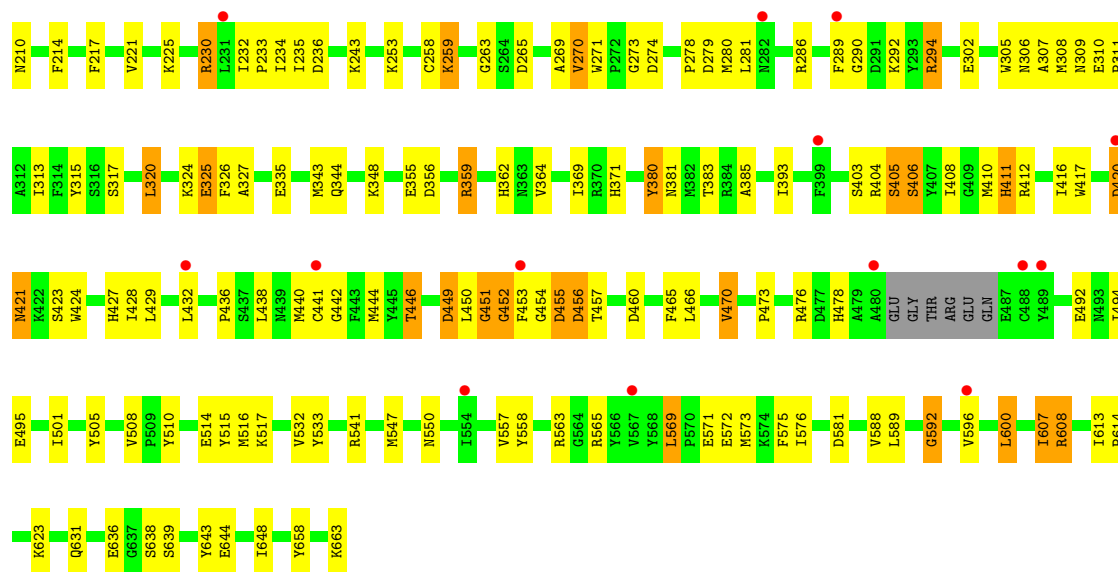
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Uncharacterized protein



• Molecule 1: Uncharacterized protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.86Å 124.03Å 87.83Å 90.00° 108.15° 90.00°	Depositor
Resolution (Å)	49.78 – 2.90 49.78 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.78-2.90) 99.0 (49.78-2.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.218 , 0.297 0.208 , 0.293	Depositor DCC
R_{free} test set	1481 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	54.9	Xtriage
Anisotropy	0.488	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10735	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% 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 ⓘ

5.1 Standard geometry ⓘ

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	0.77	1/5498 (0.0%)	1.00	4/7412 (0.1%)
1	B	0.78	0/5519	1.02	5/7441 (0.1%)
All	All	0.77	1/11017 (0.0%)	1.01	9/14853 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	613	ILE	CA-C	5.23	1.58	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	569	LEU	CA-C-N	7.07	126.77	119.56
1	A	569	LEU	C-N-CA	7.07	126.77	119.56
1	B	420	ASP	N-CA-C	6.83	117.59	108.38
1	A	59	GLY	N-CA-C	5.84	121.53	112.60
1	A	60	ILE	N-CA-C	-5.55	105.12	110.72
1	B	455	ASP	N-CA-C	-5.39	105.15	112.26
1	B	508	VAL	N-CA-CB	5.27	113.82	110.50
1	B	592	GLY	N-CA-C	5.14	118.10	110.60
1	B	532	VAL	N-CA-C	-5.00	107.85	111.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	5357	0	5136	164	0
1	B	5378	0	5149	127	0
All	All	10735	0	10285	285	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 14.

All (285) 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:450:LEU:O	1:B:451:GLY:O	1.64	1.16
1:A:343:MET:HE2	1:A:347:MET:HE2	1.38	1.02
1:A:514:GLU:OE2	1:A:517:LYS:HE3	1.59	1.01
1:A:270:VAL:HG12	1:A:271:TRP:H	1.29	0.94
1:A:76:ILE:HG23	1:A:422:LYS:HZ2	1.35	0.91
1:B:455:ASP:C	1:B:457:THR:H	1.81	0.88
1:B:429:LEU:HD22	1:B:558:TYR:CE1	2.11	0.86
1:A:420:ASP:O	1:A:451:GLY:HA3	1.76	0.85
1:B:230:ARG:HG3	1:B:230:ARG:HH11	1.40	0.85
1:B:270:VAL:HG12	1:B:271:TRP:H	1.39	0.85
1:A:75:PRO:O	1:A:420:ASP:HB2	1.78	0.84
1:A:514:GLU:OE2	1:A:514:GLU:HA	1.84	0.77
1:B:327:ALA:HB2	1:B:343:MET:HE1	1.67	0.77
1:B:234:ILE:HD11	1:B:307:ALA:CB	2.16	0.76
1:A:55:GLU:HG3	1:A:434:MET:HE3	1.67	0.76
1:A:491:PHE:O	1:A:497:PHE:HE2	1.70	0.74
1:A:16:LEU:H	1:A:16:LEU:HD12	1.54	0.73
1:A:510:TYR:CE1	1:A:514:GLU:HG2	2.26	0.70
1:A:362:HIS:NE2	1:A:371:HIS:HD2	1.89	0.70
1:A:459:ARG:HD2	1:A:492:GLU:HG3	1.74	0.70
1:B:455:ASP:C	1:B:457:THR:N	2.46	0.70
1:A:545:GLN:NE2	1:A:553:MET:HE3	2.06	0.70
1:A:343:MET:HE2	1:A:347:MET:CE	2.21	0.69
1:B:450:LEU:C	1:B:451:GLY:O	2.35	0.68
1:A:436:PRO:HG3	1:A:546:LEU:HD22	1.76	0.68
1:A:296:LEU:CD1	1:A:304:PHE:HE1	2.08	0.67
1:B:2:ILE:HG23	1:B:135:ILE:HG12	1.75	0.67
1:A:296:LEU:HA	1:A:299:GLN:HG3	1.78	0.66
1:A:545:GLN:NE2	1:A:553:MET:CE	2.58	0.65
1:B:575:PHE:HB2	1:B:589:LEU:HD12	1.78	0.65
1:B:270:VAL:HG12	1:B:271:TRP:N	2.12	0.65
1:A:447:GLY:HA3	1:A:474:LEU:HB3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:LEU:HD22	1:B:133:TYR:CE1	2.32	0.65
1:B:201:MET:HB2	1:B:204:PHE:HA	1.78	0.65
1:B:455:ASP:O	1:B:457:THR:N	2.30	0.64
1:A:391:GLU:O	1:A:395:PRO:HG3	1.97	0.64
1:B:452:GLY:O	1:B:454:GLY:N	2.30	0.64
1:B:225:LYS:HE2	1:B:302:GLU:OE2	1.97	0.63
1:A:497:PHE:O	1:A:501:ILE:HG12	1.98	0.63
1:A:491:PHE:O	1:A:497:PHE:CE2	2.52	0.62
1:A:303:GLY:HA3	1:A:400:LEU:HB3	1.81	0.62
1:A:330:PHE:O	1:B:324:LYS:HE2	2.00	0.62
1:B:234:ILE:HD11	1:B:307:ALA:HB2	1.80	0.61
1:A:269:ALA:HA	1:A:273:GLY:O	1.99	0.61
1:B:278:PRO:HB2	1:B:280:MET:HE2	1.82	0.61
1:B:643:TYR:C	1:B:643:TYR:CD2	2.77	0.61
1:A:494:ILE:HA	1:A:497:PHE:HD2	1.66	0.61
1:A:169:TRP:HD1	1:A:170:GLY:N	1.98	0.60
1:A:296:LEU:HD12	1:A:304:PHE:HE1	1.65	0.60
1:A:250:GLU:O	1:A:254:ASN:ND2	2.34	0.60
1:B:234:ILE:HD11	1:B:307:ALA:HB3	1.82	0.60
1:A:436:PRO:HG3	1:A:546:LEU:CD2	2.31	0.60
1:B:58:ARG:NH1	1:B:79:GLU:O	2.32	0.60
1:A:270:VAL:HG12	1:A:271:TRP:N	2.08	0.60
1:A:6:ARG:HG3	1:A:10:PRO:HG3	1.84	0.59
1:B:309:ASN:ND2	1:B:403:SER:OG	2.35	0.59
1:A:551:GLU:O	1:A:608:ARG:HG3	2.02	0.59
1:B:259:LYS:HG2	1:B:263:GLY:O	2.01	0.59
1:B:29:TYR:O	1:B:40:PHE:HE1	1.86	0.58
1:A:76:ILE:HG23	1:A:422:LYS:NZ	2.16	0.58
1:A:286:ARG:HG2	1:A:385:ALA:HB2	1.86	0.58
1:B:269:ALA:HA	1:B:273:GLY:O	2.03	0.58
1:A:509:PRO:HB2	1:A:630:MET:HE1	1.85	0.58
1:A:169:TRP:CD1	1:A:169:TRP:C	2.81	0.58
1:A:148:ARG:HB3	1:A:148:ARG:HH21	1.68	0.58
1:A:29:TYR:O	1:A:40:PHE:HE1	1.87	0.57
1:A:398:ARG:NH1	1:A:645:ASP:OD2	2.37	0.57
1:B:449:ASP:OD2	1:B:478:HIS:HB3	2.03	0.57
1:B:88:HIS:NE2	1:B:406:SER:HB2	2.20	0.57
1:A:494:ILE:HA	1:A:497:PHE:CD2	2.39	0.57
1:B:320:LEU:O	1:B:324:LYS:HG3	2.05	0.56
1:B:157:PRO:HG2	1:B:162:PHE:HE2	1.69	0.56
1:B:230:ARG:HH11	1:B:230:ARG:CG	2.15	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:ASN:HD21	1:B:404:ARG:H	1.53	0.56
1:A:310:GLU:N	1:A:311:PRO:HA	2.21	0.56
1:A:538:MET:HG2	1:A:568:TYR:CZ	2.41	0.55
1:A:605:LEU:HD12	1:A:605:LEU:C	2.32	0.55
1:A:169:TRP:HH2	1:A:271:TRP:CH2	2.25	0.55
1:A:48:ASP:OD1	1:A:95:GLY:HA3	2.07	0.55
1:A:217:PHE:HB3	1:A:218:PRO:HD3	1.90	0.54
1:B:45:ASP:HB3	1:B:48:ASP:OD1	2.07	0.54
1:A:396:GLU:O	1:A:649:HIS:HA	2.07	0.54
1:A:424:TRP:HB3	1:B:79:GLU:OE2	2.08	0.54
1:B:210:ASN:O	1:B:214:PHE:HB2	2.07	0.54
1:B:100:GLY:O	1:B:101:LEU:HD23	2.08	0.54
1:A:459:ARG:NH2	1:A:496:ASP:OD2	2.41	0.53
1:B:557:VAL:HG22	1:B:565:ARG:HG2	1.89	0.53
1:A:453:PHE:HB3	1:A:456:ASP:OD2	2.08	0.53
1:B:636:GLU:CD	1:B:663:LYS:HE3	2.34	0.53
1:B:362:HIS:NE2	1:B:371:HIS:HD2	2.05	0.53
1:A:218:PRO:HG3	1:A:299:GLN:HB3	1.90	0.53
1:B:177:PHE:HD1	1:B:194:ILE:HG21	1.74	0.52
1:B:444:MET:O	1:B:473:PRO:HG2	2.09	0.52
1:B:573:MET:HG2	1:B:608:ARG:HA	1.90	0.52
1:A:651:ASP:OD1	1:A:654:LYS:HG3	2.09	0.52
1:A:293:TYR:CE1	1:A:386:ALA:HB2	2.45	0.52
1:A:505:TYR:HA	1:A:508:VAL:HG23	1.92	0.52
1:A:646:ASP:OD1	1:A:646:ASP:C	2.52	0.52
1:B:64:GLY:O	1:B:65:TYR:HB2	2.10	0.52
1:B:94:SER:HA	1:B:98:THR:HG23	1.91	0.51
1:B:380:TYR:C	1:B:380:TYR:CD1	2.88	0.51
1:A:88:HIS:HB3	1:A:408:ILE:HD11	1.92	0.51
1:B:356:ASP:OD1	1:B:359:ARG:NH2	2.44	0.51
1:B:38:PHE:CZ	1:B:132:ILE:HD11	2.45	0.51
1:A:271:TRP:HB2	1:A:272:PRO:HD3	1.92	0.51
1:A:343:MET:CE	1:A:347:MET:HE2	2.26	0.51
1:B:411:HIS:HB2	1:B:442:GLY:O	2.11	0.51
1:A:117:ARG:NH1	1:A:119:ASP:OD2	2.44	0.51
1:A:109:LEU:HD12	1:A:125:CYS:HB3	1.92	0.50
1:B:309:ASN:OD1	1:B:405:SER:HB3	2.11	0.50
1:A:156:ILE:HD13	1:A:516:MET:HE1	1.93	0.50
1:A:510:TYR:CE1	1:A:514:GLU:CG	2.94	0.50
1:B:286:ARG:HG2	1:B:385:ALA:HB2	1.94	0.50
1:B:190:PRO:HB2	1:B:505:TYR:CE1	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:LYS:HE2	1:B:265:ASP:OD1	2.11	0.50
1:B:167:SER:O	1:B:478:HIS:HA	2.11	0.50
1:A:539:ALA:O	1:A:541:ARG:N	2.45	0.49
1:A:16:LEU:HA	1:A:151:ILE:HA	1.95	0.49
1:B:280:MET:O	1:B:381:ASN:HB3	2.13	0.49
1:A:371:HIS:HE1	1:A:375:HIS:ND1	2.10	0.49
1:B:55:GLU:HB2	1:B:438:LEU:HD21	1.95	0.49
1:A:227:GLN:O	1:A:229:LEU:HG	2.13	0.49
1:B:510:TYR:HB2	1:B:614:PRO:HD2	1.94	0.49
1:A:204:PHE:HB2	1:A:272:PRO:HG3	1.96	0.48
1:B:90:PHE:CZ	1:B:92:ILE:HD11	2.48	0.48
1:A:574:LYS:CE	1:A:586:GLU:OE2	2.61	0.48
1:B:75:PRO:O	1:B:421:ASN:HB3	2.13	0.48
1:A:536:ASP:HB3	1:A:539:ALA:HB3	1.95	0.48
1:A:147:PHE:CE2	1:A:151:ILE:HG21	2.50	0.47
1:A:86:GLY:HA3	1:A:88:HIS:CE1	2.50	0.47
1:A:538:MET:HG2	1:A:568:TYR:CE1	2.49	0.47
1:A:70:ASN:O	1:A:71:CYS:C	2.58	0.47
1:A:654:LYS:O	1:A:657:ASN:HB2	2.14	0.47
1:B:141:TYR:CZ	1:B:145:LYS:HD3	2.49	0.47
1:A:247:VAL:HG22	1:A:292:LYS:HE2	1.97	0.47
1:A:547:MET:HE1	1:A:608:ARG:HD3	1.96	0.47
1:A:278:PRO:O	1:A:280:MET:HG2	2.15	0.47
1:A:148:ARG:NH1	1:A:444:MET:SD	2.88	0.46
1:A:296:LEU:HD13	1:A:304:PHE:HE1	1.78	0.46
1:A:447:GLY:CA	1:A:474:LEU:HB3	2.45	0.46
1:B:410:MET:O	1:B:412:ARG:N	2.48	0.46
1:B:423:SER:OG	1:B:452:GLY:HA2	2.14	0.46
1:B:658:TYR:CD2	1:B:658:TYR:N	2.83	0.46
1:A:444:MET:HE1	1:A:518:ALA:HB3	1.97	0.46
1:A:574:LYS:HB3	1:A:576:ILE:HD13	1.97	0.46
1:B:289:PHE:O	1:B:290:GLY:C	2.57	0.46
1:A:78:THR:HB	1:B:424:TRP:CZ2	2.50	0.46
1:A:82:ARG:NH2	1:B:456:ASP:OD1	2.39	0.46
1:A:88:HIS:HA	1:A:408:ILE:HD12	1.98	0.46
1:A:545:GLN:HE21	1:A:553:MET:HE3	1.77	0.46
1:A:53:LEU:O	1:A:54:GLY:C	2.58	0.46
1:A:234:ILE:HD12	1:A:305:TRP:CZ2	2.50	0.46
1:A:189:ILE:HD11	1:A:494:ILE:HG12	1.96	0.46
1:A:510:TYR:HB2	1:A:614:PRO:HG2	1.98	0.46
1:A:568:TYR:O	1:A:570:PRO:HD3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ILE:HD11	1:A:144:VAL:HG22	1.97	0.46
1:B:196:MET:HB2	1:B:233:PRO:HA	1.98	0.46
1:B:235:ILE:HB	1:B:306:ASN:ND2	2.31	0.46
1:B:438:LEU:HD13	1:B:446:THR:HG21	1.98	0.46
1:A:196:MET:SD	1:A:231:LEU:HD13	2.57	0.45
1:B:428:ILE:HG22	1:B:558:TYR:HB3	1.97	0.45
1:A:78:THR:OG1	1:A:80:ASP:OD1	2.32	0.45
1:A:362:HIS:NE2	1:A:371:HIS:CD2	2.78	0.45
1:A:422:LYS:HD2	1:A:424:TRP:CZ2	2.51	0.45
1:A:570:PRO:HD2	1:A:573:MET:HE3	1.98	0.45
1:B:184:TYR:CD2	1:B:189:ILE:HB	2.50	0.45
1:A:114:GLY:HA2	1:A:117:ARG:O	2.16	0.45
1:B:162:PHE:HZ	1:B:515:TYR:CG	2.34	0.45
1:A:390:PHE:CE2	1:A:399:PHE:HB2	2.51	0.45
1:B:450:LEU:HD13	1:B:465:PHE:CE1	2.52	0.45
1:A:196:MET:HB2	1:A:233:PRO:HA	1.99	0.45
1:A:232:ILE:HD12	1:A:400:LEU:HD23	1.99	0.45
1:B:7:TYR:O	1:B:129:ASP:HA	2.17	0.45
1:B:121:LEU:HD12	1:B:122:LYS:N	2.31	0.45
1:A:644:GLU:OE2	1:A:659:ARG:NH2	2.49	0.45
1:B:230:ARG:HG3	1:B:230:ARG:NH1	2.19	0.45
1:A:16:LEU:HB3	1:A:150:VAL:O	2.17	0.45
1:A:148:ARG:NH1	1:A:523:ASP:O	2.49	0.45
1:A:166:GLN:NE2	1:A:477:ASP:OD2	2.50	0.45
1:B:533:TYR:CE2	1:B:547:MET:HE2	2.52	0.45
1:A:85:TYR:CE1	1:A:404:ARG:HD2	2.51	0.45
1:A:515:TYR:C	1:A:515:TYR:CD2	2.95	0.45
1:B:278:PRO:HB2	1:B:280:MET:CE	2.46	0.45
1:B:436:PRO:O	1:B:440:MET:HG3	2.17	0.45
1:B:515:TYR:C	1:B:515:TYR:CD2	2.95	0.45
1:B:271:TRP:HE1	1:B:308:MET:HE3	1.81	0.44
1:A:169:TRP:CD1	1:A:170:GLY:N	2.84	0.44
1:B:417:TRP:CD1	1:B:476:ARG:HH21	2.36	0.44
1:A:291:ASP:O	1:A:294:ARG:HB2	2.16	0.44
1:A:423:SER:HB3	1:A:461:LEU:HD21	2.00	0.44
1:B:197:ASP:O	1:B:199:ASP:N	2.50	0.44
1:B:466:LEU:O	1:B:470:VAL:HG23	2.16	0.44
1:B:281:LEU:HB3	1:B:364:VAL:HG22	1.98	0.44
1:A:577:LYS:HB2	1:A:585:SER:OG	2.18	0.44
1:A:537:LYS:HD2	1:A:537:LYS:HA	1.53	0.44
1:A:196:MET:HE2	1:A:200:TYR:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:PHE:O	1:A:292:LYS:HB2	2.18	0.44
1:B:160:PHE:CD1	1:B:160:PHE:C	2.96	0.44
1:A:62:LYS:HA	1:A:62:LYS:CE	2.48	0.43
1:A:361:TYR:CE1	1:A:370:ARG:HG3	2.53	0.43
1:A:154:SER:OG	1:A:442:GLY:O	2.31	0.43
1:B:514:GLU:OE2	1:B:517:LYS:NZ	2.35	0.43
1:A:281:LEU:CD1	1:A:374:VAL:HG21	2.49	0.43
1:A:348:LYS:HD2	1:B:344:GLN:NE2	2.33	0.43
1:A:421:ASN:O	1:A:422:LYS:HG3	2.17	0.43
1:B:429:LEU:HD12	1:B:432:LEU:HD23	1.99	0.43
1:A:176:ASP:O	1:A:180:VAL:HG23	2.18	0.43
1:B:53:LEU:HA	1:B:88:HIS:O	2.17	0.43
1:A:260:ARG:NH1	1:A:264:SER:OG	2.51	0.43
1:A:517:LYS:O	1:A:521:ASN:HB2	2.18	0.43
1:B:575:PHE:HB2	1:B:589:LEU:CD1	2.47	0.43
1:B:195:TYR:CD2	1:B:232:ILE:HB	2.53	0.43
1:B:217:PHE:O	1:B:221:VAL:HG23	2.19	0.43
1:A:390:PHE:CD2	1:A:399:PHE:HB2	2.54	0.43
1:B:294:ARG:NH2	1:B:393:ILE:HG12	2.33	0.43
1:B:428:ILE:CG2	1:B:558:TYR:HB3	2.49	0.43
1:A:417:TRP:NE1	1:A:476:ARG:NH2	2.65	0.43
1:B:230:ARG:CG	1:B:230:ARG:NH1	2.79	0.43
1:B:310:GLU:N	1:B:311:PRO:HA	2.34	0.43
1:A:130:LEU:N	1:A:130:LEU:HD23	2.34	0.43
1:A:204:PHE:CB	1:A:272:PRO:HG3	2.49	0.43
1:A:296:LEU:HD13	1:A:304:PHE:CE1	2.53	0.43
1:B:156:ILE:HG21	1:B:516:MET:HE1	2.00	0.43
1:A:169:TRP:CH2	1:A:271:TRP:CH2	3.06	0.42
1:A:447:GLY:HA3	1:A:474:LEU:HD23	2.00	0.42
1:B:280:MET:O	1:B:281:LEU:HD12	2.19	0.42
1:B:313:ILE:HG22	1:B:315:TYR:N	2.33	0.42
1:A:458:THR:OG1	1:A:460:ASP:OD1	2.34	0.42
1:B:92:ILE:N	1:B:92:ILE:HD12	2.33	0.42
1:A:53:LEU:HA	1:A:88:HIS:O	2.19	0.42
1:A:433:LYS:O	1:A:436:PRO:HD2	2.19	0.42
1:A:439:ASN:HB3	1:A:524:MET:HE3	2.01	0.42
1:B:607:ILE:HD12	1:B:613:ILE:HG12	2.01	0.42
1:A:574:LYS:HE2	1:A:586:GLU:OE2	2.18	0.42
1:B:232:ILE:O	1:B:305:TRP:NE1	2.52	0.42
1:B:327:ALA:CB	1:B:343:MET:HE1	2.44	0.42
1:B:571:GLU:O	1:B:572:GLU:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:MET:HE3	1:A:343:MET:HA	2.02	0.42
1:A:422:LYS:HD2	1:A:424:TRP:HZ2	1.84	0.42
1:B:179:ALA:O	1:B:180:VAL:C	2.63	0.42
1:B:193:MET:HG2	1:B:232:ILE:HD11	2.00	0.42
1:A:13:THR:O	1:A:380:TYR:HE2	2.03	0.42
1:B:383:THR:HG21	1:B:410:MET:HE3	2.02	0.42
1:A:256:TYR:CD2	1:A:288:TRP:CD2	3.08	0.42
1:A:617:GLU:OE1	1:A:617:GLU:HA	2.20	0.42
1:B:101:LEU:HD22	1:B:134:VAL:HG22	2.02	0.42
1:B:325:GLU:HG3	1:B:326:PHE:N	2.34	0.42
1:B:643:TYR:CD2	1:B:644:GLU:N	2.88	0.42
1:A:139:ASN:OD1	1:A:141:TYR:HB3	2.19	0.41
1:B:270:VAL:CG1	1:B:271:TRP:N	2.82	0.41
1:B:569:LEU:O	1:B:592:GLY:N	2.52	0.41
1:A:315:TYR:CE1	1:A:320:LEU:HD22	2.55	0.41
1:B:206:ASP:OD2	1:B:236:ASP:O	2.38	0.41
1:A:260:ARG:HB2	1:A:262:ASP:OD1	2.21	0.41
1:A:376:ASN:O	1:A:407:TYR:HB2	2.20	0.41
1:A:436:PRO:HB3	1:A:528:PRO:HG3	2.01	0.41
1:A:651:ASP:O	1:A:651:ASP:CG	2.63	0.41
1:A:523:ASP:OD1	1:A:524:MET:N	2.51	0.41
1:A:261:GLU:C	1:A:263:GLY:N	2.77	0.41
1:A:563:ARG:HD2	1:B:541:ARG:HD2	2.01	0.41
1:A:514:GLU:OE2	1:A:517:LYS:CE	2.49	0.41
1:B:202:GLN:O	1:B:203:ASP:C	2.64	0.41
1:A:418:MET:H	1:A:418:MET:HG2	1.74	0.41
1:A:421:ASN:OD1	1:A:421:ASN:N	2.53	0.41
1:B:27:PHE:HA	1:B:28:PRO:HD3	1.94	0.41
1:B:189:ILE:HA	1:B:190:PRO:HD3	1.98	0.41
1:B:460:ASP:HB2	1:B:600:LEU:HD21	2.03	0.41
1:A:55:GLU:HB2	1:A:438:LEU:HD21	2.03	0.41
1:A:334:THR:C	1:A:336:GLY:N	2.79	0.41
1:A:348:LYS:HB3	1:A:348:LYS:HE2	1.82	0.41
1:B:153:ARG:O	1:B:412:ARG:HG2	2.20	0.41
1:B:306:ASN:HB3	1:B:309:ASN:HD22	1.86	0.41
1:B:380:TYR:HA	1:B:410:MET:SD	2.61	0.41
1:A:200:TYR:CZ	1:A:233:PRO:HB2	2.56	0.40
1:A:417:TRP:CD1	1:A:476:ARG:NH2	2.89	0.40
1:A:598:VAL:HG11	1:A:604:PRO:HG3	2.03	0.40
1:B:56:SER:CA	1:B:84:LEU:HD12	2.51	0.40
1:A:527:LYS:NZ	1:A:550:ASN:OD1	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:596:VAL:O	1:B:596:VAL:HG23	2.21	0.40
1:A:616:ALA:HB2	1:A:630:MET:HE3	2.04	0.40
1:B:258:CYS:SG	1:B:279:ASP:HA	2.61	0.40
1:B:408:ILE:HD11	1:B:441:CYS:HB3	2.04	0.40
1:A:22:THR:HG22	1:A:24:GLU:OE2	2.22	0.40
1:A:579:MET:HB3	1:A:581:ASP:OD1	2.21	0.40
1:B:198:ILE:HG22	1:B:201:MET:HE2	2.04	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	651/666 (98%)	591 (91%)	49 (8%)	11 (2%)	7	26
1	B	654/666 (98%)	588 (90%)	53 (8%)	13 (2%)	6	22
All	All	1305/1332 (98%)	1179 (90%)	102 (8%)	24 (2%)	6	25

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	451	GLY
1	B	453	PHE
1	B	638	SER
1	A	54	GLY
1	A	540	ILE
1	B	198	ILE
1	B	456	ASP
1	A	138	GLU
1	A	253	LYS
1	B	411	HIS
1	A	71	CYS

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Mol	Chain	Res	Type
1	A	262	ASP
1	A	335	GLU
1	A	653	ASP
1	B	600	LEU
1	A	452	GLY
1	B	181	ALA
1	B	203	ASP
1	B	170	GLY
1	B	270	VAL
1	A	604	PRO
1	B	470	VAL
1	A	270	VAL
1	B	452	GLY

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	563/574 (98%)	503 (89%)	60 (11%)	6	22
1	B	565/574 (98%)	508 (90%)	57 (10%)	7	24
All	All	1128/1148 (98%)	1011 (90%)	117 (10%)	7	23

All (117) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASP
1	A	14	GLU
1	A	16	LEU
1	A	34	GLN
1	A	36	GLU
1	A	62	LYS
1	A	76	ILE
1	A	79	GLU
1	A	115	TYR
1	A	118	MET

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Mol	Chain	Res	Type
1	A	127	ASN
1	A	148	ARG
1	A	166	GLN
1	A	169	TRP
1	A	173	THR
1	A	174	LYS
1	A	192	ASP
1	A	211	GLU
1	A	240	LYS
1	A	243	LYS
1	A	254	ASN
1	A	280	MET
1	A	299	GLN
1	A	305	TRP
1	A	308	MET
1	A	332	LYS
1	A	334	THR
1	A	335	GLU
1	A	355	GLU
1	A	367	LYS
1	A	369	ILE
1	A	388	GLU
1	A	392	ARG
1	A	393	ILE
1	A	401	MET
1	A	403	SER
1	A	408	ILE
1	A	416	ILE
1	A	417	TRP
1	A	418	MET
1	A	422	LYS
1	A	430	LEU
1	A	432	LEU
1	A	433	LYS
1	A	434	MET
1	A	439	ASN
1	A	450	LEU
1	A	455	ASP
1	A	494	ILE
1	A	495	GLU
1	A	514	GLU
1	A	529	LEU

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Mol	Chain	Res	Type
1	A	537	LYS
1	A	567	VAL
1	A	576	ILE
1	A	602	GLU
1	A	609	SER
1	A	641	THR
1	A	646	ASP
1	A	654	LYS
1	B	1	MET
1	B	2	ILE
1	B	16	LEU
1	B	18	GLU
1	B	41	THR
1	B	56	SER
1	B	66	CYS
1	B	72	THR
1	B	76	ILE
1	B	81	LYS
1	B	82	ARG
1	B	94	SER
1	B	97	THR
1	B	126	GLU
1	B	154	SER
1	B	198	ILE
1	B	203	ASP
1	B	230	ARG
1	B	243	LYS
1	B	253	LYS
1	B	259	LYS
1	B	274	ASP
1	B	292	LYS
1	B	294	ARG
1	B	317	SER
1	B	320	LEU
1	B	325	GLU
1	B	335	GLU
1	B	348	LYS
1	B	355	GLU
1	B	359	ARG
1	B	369	ILE
1	B	380	TYR
1	B	405	SER

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Mol	Chain	Res	Type
1	B	406	SER
1	B	416	ILE
1	B	420	ASP
1	B	421	ASN
1	B	427	HIS
1	B	446	THR
1	B	449	ASP
1	B	492	GLU
1	B	494	ILE
1	B	495	GLU
1	B	501	ILE
1	B	550	ASN
1	B	563	ARG
1	B	569	LEU
1	B	576	ILE
1	B	581	ASP
1	B	588	VAL
1	B	607	ILE
1	B	608	ARG
1	B	623	LYS
1	B	631	GLN
1	B	639	SER
1	B	648	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	89	ASN
1	A	166	GLN
1	A	187	ASN
1	A	254	ASN
1	A	255	ASN
1	A	306	ASN
1	A	344	GLN
1	A	371	HIS
1	A	427	HIS
1	A	493	ASN
1	B	70	ASN
1	B	77	HIS
1	B	139	ASN
1	B	146	GLN

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Mol	Chain	Res	Type
1	B	306	ASN
1	B	344	GLN
1	B	371	HIS
1	B	381	ASN
1	B	421	ASN
1	B	431	ASN
1	B	493	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	655/666 (98%)	-0.17	15 (2%)	61 52	33, 54, 69, 104	4 (0%)
1	B	658/666 (98%)	-0.24	15 (2%)	61 52	34, 53, 70, 92	1 (0%)
All	All	1313/1332 (98%)	-0.20	30 (2%)	61 52	33, 54, 69, 104	5 (0%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	402	PHE	5.0
1	B	488	CYS	4.5
1	A	362	HIS	4.1
1	B	596	VAL	4.0
1	A	60	ILE	3.4
1	B	480	ALA	3.4
1	B	147	PHE	3.3
1	B	289	PHE	3.2
1	A	289	PHE	3.1
1	A	369	ILE	3.0
1	A	46	GLU	2.9
1	B	489	TYR	2.9
1	A	480	ALA	2.8
1	B	441	CYS	2.7
1	B	399	PHE	2.6
1	A	419	GLY	2.5
1	A	159	LYS	2.5
1	B	432	LEU	2.5
1	A	491	PHE	2.3
1	A	25	GLU	2.2
1	B	420	ASP	2.2
1	B	231	LEU	2.2
1	A	314	PHE	2.2
1	B	453	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	453	PHE	2.1
1	B	282	ASN	2.1
1	A	11	PHE	2.0
1	A	171	TYR	2.0
1	B	554	ILE	2.0
1	B	567	VAL	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.