



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:03 AM UTC

PDB ID : 3S4B / pdb_00003s4b
Title : Cellobiose phosphorylase from *Cellulomonas uda* in complex with glucose
Authors : Van Hoorebeke, A.; Stout, J.; Soetaert, W.; Van Beeumen, J.; Desmet, T.; Savvides, S.
Deposited on : 2011-05-19
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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
Mogul : 2022.3.0, CSD as543be (2022)
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

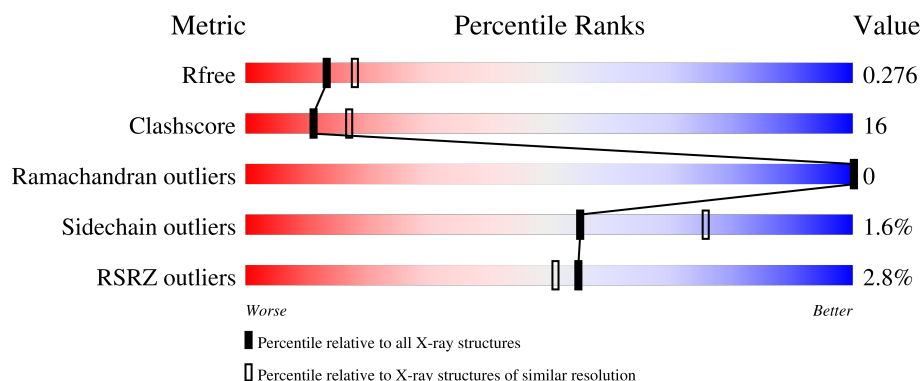
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	822	
1	B	822	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	A	823	X	-	-	-
2	GLC	A	824	X	-	-	-
2	GLC	B	823	X	-	-	-
2	GLC	B	824	X	-	-	-

2 Entry composition [i](#)

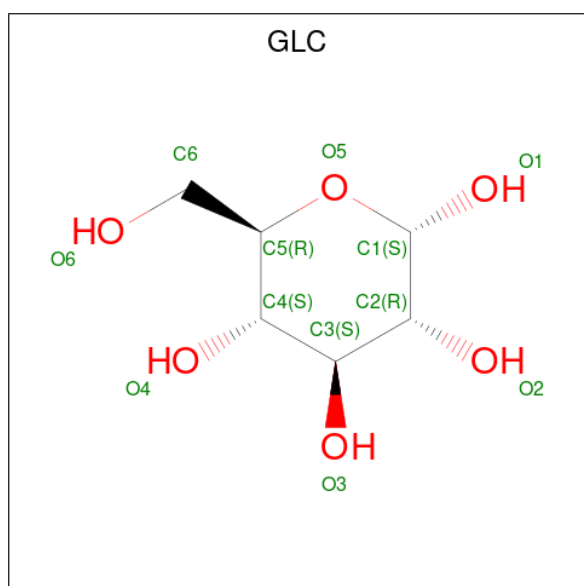
There are 3 unique types of molecules in this entry. The entry contains 25692 atoms, of which 12135 are hydrogens and 0 are deuteriums.

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 Cellobiose phosphorylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	822	Total	C	H	N	O	S	9	0	0
			12495	4063	6061	1110	1244	17			
1	B	822	Total	C	H	N	O	S	7	1	0
			12521	4070	6074	1110	1250	17			

- Molecule 2 is alpha-D-glucopyranose (CCD ID: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		

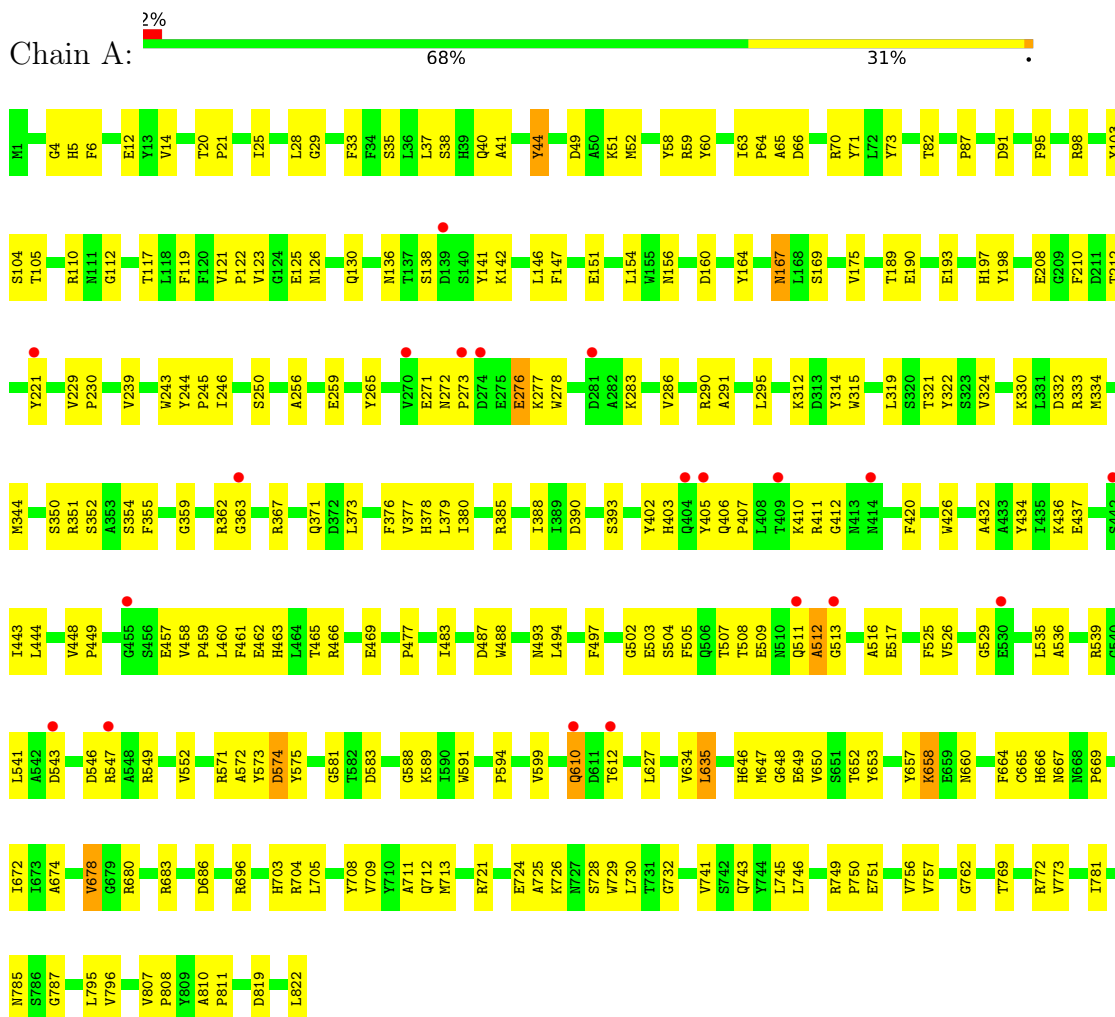
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	327	Total 327	O 327	0	0
3	B	301	Total 301	O 301	0	0

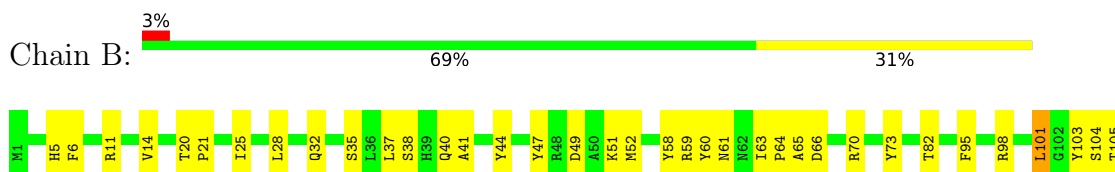
3 Residue-property plots

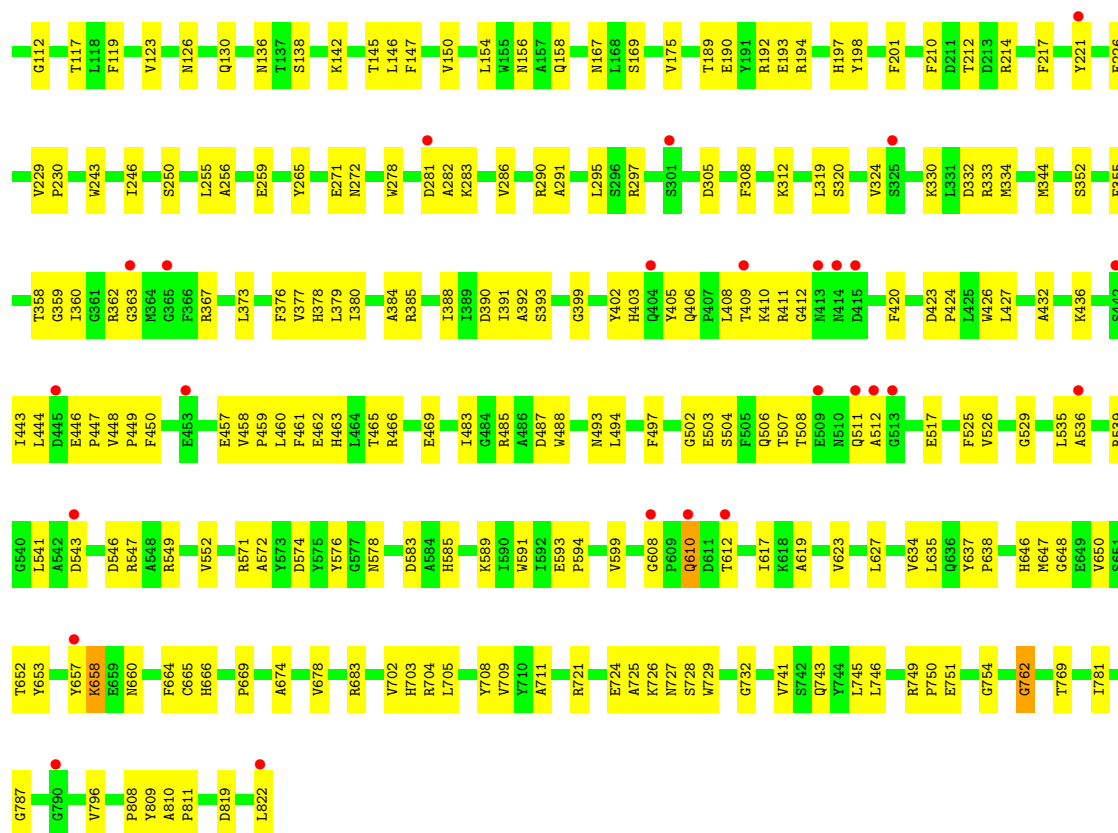
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Cellobiose phosphorylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.42Å 102.74Å 98.31Å 90.00° 96.44° 90.00°	Depositor
Resolution (Å)	19.95 – 2.40 19.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	77.3 (19.95-2.40) 81.9 (19.95-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.41Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.231 , 0.281 0.229 , 0.276	Depositor DCC
R_{free} test set	2715 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.655	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 46.7	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	0.89	EDS
Total number of atoms	25692	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.84 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.9398e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GLC

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	0.46	0/6608	0.84	10/9008 (0.1%)
1	B	0.48	0/6624	0.84	10/9029 (0.1%)
All	All	0.47	0/13232	0.84	20/18037 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	512	ALA	N-CA-C	6.98	118.97	111.36
1	A	658	LYS	CB-CA-C	-6.97	108.52	116.54
1	A	512	ALA	N-CA-C	6.87	118.84	111.36
1	B	658	LYS	CB-CA-C	-6.68	108.88	116.63
1	A	20	THR	CA-C-N	5.99	125.67	119.56
1	A	20	THR	C-N-CA	5.99	125.67	119.56
1	B	82	THR	CA-C-N	5.84	125.70	119.28
1	B	82	THR	C-N-CA	5.84	125.70	119.28
1	A	277	LYS	N-CA-C	5.64	118.19	111.71
1	A	44	TYR	N-CA-C	5.36	115.61	108.38
1	A	82	THR	CA-C-N	5.33	125.14	119.28
1	A	82	THR	C-N-CA	5.33	125.14	119.28
1	A	167	ASN	N-CA-C	5.26	119.57	113.20
1	B	762	GLY	CA-C-N	5.21	125.01	119.28
1	B	762	GLY	C-N-CA	5.21	125.01	119.28
1	B	20	THR	CA-C-N	5.15	124.81	119.56
1	B	20	THR	C-N-CA	5.15	124.81	119.56
1	A	350	SER	CB-CA-C	-5.04	110.78	116.63
1	B	608	GLY	CA-C-N	5.01	125.28	119.47
1	B	608	GLY	C-N-CA	5.01	125.28	119.47

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	6434	6061	6040	216	1
1	B	6447	6074	6052	209	0
2	A	24	0	24	0	0
2	B	24	0	24	0	0
3	A	327	0	0	12	0
3	B	301	0	0	15	0
All	All	13557	12135	12140	407	1

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 16.

All (407) 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:A:193:GLU:HG3	1:B:652:THR:HG22	1.16	1.14
1:A:193:GLU:HG3	1:B:652:THR:CG2	1.86	1.03
1:A:59:ARG:NH2	1:A:151:GLU:OE1	1.97	0.96
1:A:512:ALA:O	3:A:1083:HOH:O	1.82	0.95
1:A:504:SER:O	1:A:508:THR:HG22	1.67	0.94
1:B:504:SER:O	1:B:508:THR:HG22	1.68	0.93
1:A:741:VAL:O	1:A:746:LEU:HD13	1.66	0.93
1:B:485:ARG:HD3	1:B:506:GLN:O	1.69	0.92
1:B:278:TRP:CD1	1:B:283:LYS:HG2	2.07	0.90
1:B:574:ASP:OD2	1:B:578:ASN:HB2	1.74	0.87
1:A:278:TRP:CD1	1:A:283:LYS:HG2	2.09	0.87
1:B:385:ARG:HG3	1:B:443:ILE:HD12	1.58	0.85
1:A:741:VAL:HA	1:A:745:LEU:HD12	1.60	0.84
1:B:646:HIS:CD2	1:B:647:MET:HG3	2.15	0.82
1:B:741:VAL:HA	1:B:745:LEU:HD12	1.62	0.81
1:A:193:GLU:CG	1:B:652:THR:HG22	2.05	0.80
1:B:743:GLN:HB3	1:B:749:ARG:HB3	1.64	0.79
1:A:273:PRO:HG2	1:A:276:GLU:HG3	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:ARG:HG3	1:A:443:ILE:HD12	1.66	0.78
1:A:743:GLN:HB3	1:A:749:ARG:HB3	1.68	0.75
1:A:448:VAL:HG12	1:A:463:HIS:CE1	2.21	0.74
1:A:59:ARG:HH22	1:A:151:GLU:CD	1.93	0.74
1:A:363:GLY:HA2	1:A:406:GLN:OE1	1.87	0.73
1:B:363:GLY:HA2	1:B:406:GLN:OE1	1.87	0.73
1:B:666:HIS:CE1	1:B:732:GLY:HA3	2.25	0.72
1:B:517:GLU:O	1:B:572:ALA:HB1	1.89	0.71
1:A:571:ARG:HD3	1:A:591:TRP:CB	2.20	0.71
1:A:517:GLU:HB2	1:A:573:TYR:HB2	1.74	0.69
1:A:721:ARG:HD3	1:A:724:GLU:OE1	1.93	0.69
1:B:409:THR:HG22	1:B:411:ARG:HG2	1.75	0.69
1:B:571:ARG:HD3	1:B:591:TRP:CB	2.23	0.68
1:B:448:VAL:HG12	1:B:463:HIS:CE1	2.28	0.68
1:A:273:PRO:CG	1:A:276:GLU:HG3	2.24	0.67
1:B:721:ARG:HD3	1:B:724:GLU:OE1	1.94	0.67
1:B:444:LEU:HD11	1:B:535:LEU:HD22	1.77	0.67
1:A:193:GLU:CG	1:B:652:THR:CG2	2.68	0.67
1:B:539:ARG:HG2	1:B:541:LEU:HD13	1.76	0.66
1:A:378:HIS:CE1	1:A:379:LEU:HG	2.31	0.65
1:A:571:ARG:HD3	1:A:591:TRP:HB2	1.77	0.65
1:A:444:LEU:HD11	1:A:535:LEU:HD22	1.78	0.64
1:B:402:TYR:CG	1:B:412:GLY:HA3	2.33	0.64
1:A:539:ARG:HG2	1:A:541:LEU:HD13	1.79	0.64
1:A:465:THR:O	1:A:469:GLU:HG2	1.97	0.64
1:A:674:ALA:O	1:A:678:VAL:HG13	1.98	0.63
1:B:378:HIS:CE1	1:B:379:LEU:HG	2.33	0.63
1:A:402:TYR:CG	1:A:412:GLY:HA3	2.33	0.63
1:B:571:ARG:HD3	1:B:591:TRP:HB2	1.80	0.63
1:B:741:VAL:O	1:B:746:LEU:HG	1.98	0.63
1:B:465:THR:O	1:B:469:GLU:HG2	1.99	0.62
1:B:229:VAL:HB	1:B:230:PRO:HD3	1.82	0.61
1:A:273:PRO:HB2	1:A:276:GLU:HG3	1.81	0.61
1:B:373:LEU:HA	1:B:376:PHE:CZ	2.36	0.61
1:B:305:ASP:HA	3:B:1103:HOH:O	2.02	0.60
1:A:14:VAL:HG23	1:A:98:ARG:HG2	1.83	0.60
1:B:150:VAL:HG23	1:B:201:PHE:CZ	2.37	0.60
1:A:5:HIS:O	1:A:14:VAL:HG12	2.03	0.59
1:B:610:GLN:OE1	1:B:610:GLN:HA	2.02	0.59
1:A:373:LEU:HA	1:A:376:PHE:CZ	2.38	0.59
1:A:322:TYR:HD1	1:A:772:ARG:HD3	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:TYR:CD1	1:A:772:ARG:HD3	2.38	0.58
1:A:726:LYS:HE3	1:B:243:TRP:CE2	2.38	0.58
1:B:727:ASN:HB3	3:B:952:HOH:O	2.02	0.58
1:A:273:PRO:HG2	1:A:276:GLU:CG	2.33	0.58
1:A:437:GLU:CD	1:A:772:ARG:HH12	2.12	0.58
1:A:363:GLY:CA	1:A:406:GLN:OE1	2.51	0.58
1:A:646:HIS:CE1	1:A:647:MET:HG3	2.39	0.58
1:B:409:THR:HG22	1:B:411:ARG:CG	2.35	0.57
1:B:674:ALA:O	1:B:678:VAL:HG13	2.05	0.57
1:A:321:THR:CG2	1:A:772:ARG:NH2	2.67	0.57
1:A:796:VAL:HB	1:A:819:ASP:HB2	1.87	0.56
1:A:612:THR:HG22	1:A:612:THR:O	2.05	0.56
1:A:273:PRO:CB	1:A:276:GLU:HG3	2.36	0.56
1:A:371:GLN:OE1	1:A:666:HIS:HE1	1.87	0.56
1:A:21:PRO:HB2	1:A:704:ARG:HA	1.88	0.56
1:A:509:GLU:HB3	3:A:1093:HOH:O	2.05	0.56
1:A:635:LEU:HD13	1:A:635:LEU:O	2.06	0.56
1:A:635:LEU:HD23	1:A:667:ASN:HD21	1.71	0.56
1:A:653:TYR:CE2	1:A:658:LYS:HG3	2.41	0.56
1:B:406:GLN:HB3	3:B:1124:HOH:O	2.05	0.56
1:B:653:TYR:CE2	1:B:658:LYS:HG3	2.41	0.55
1:A:28:LEU:HB2	1:A:35:SER:HB2	1.88	0.55
1:B:28:LEU:HB2	1:B:35:SER:HB2	1.88	0.55
1:B:503:GLU:HB3	1:B:508:THR:HG21	1.89	0.55
1:B:796:VAL:HB	1:B:819:ASP:HB2	1.88	0.55
1:A:503:GLU:HB3	1:A:508:THR:HG21	1.88	0.55
1:B:150:VAL:HG21	1:B:201:PHE:CE1	2.42	0.55
1:B:11:ARG:NH2	3:B:885:HOH:O	2.40	0.55
1:B:363:GLY:CA	1:B:406:GLN:OE1	2.54	0.55
1:A:352:SER:O	1:A:359:GLY:HA2	2.06	0.55
1:B:432:ALA:O	1:B:436:LYS:HG3	2.05	0.55
1:A:610:GLN:HA	1:A:610:GLN:OE1	2.07	0.55
1:B:6:PHE:CD1	1:B:333:ARG:HD3	2.42	0.54
1:A:321:THR:HG22	1:A:772:ARG:NH2	2.21	0.54
1:B:32:GLN:HB3	1:B:47:TYR:CE1	2.43	0.54
1:B:49:ASP:HB2	1:B:408:LEU:HD13	1.89	0.54
1:A:6:PHE:CD1	1:A:333:ARG:HD3	2.42	0.54
1:B:612:THR:O	1:B:612:THR:HG22	2.07	0.54
1:A:635:LEU:HD13	1:A:635:LEU:C	2.33	0.54
1:B:373:LEU:HD23	1:B:377:VAL:HB	1.89	0.54
1:A:376:PHE:CD1	1:A:376:PHE:C	2.85	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:458:VAL:HB	1:B:459:PRO:HD2	1.90	0.54
1:A:73:TYR:HB2	1:A:147:PHE:HB2	1.89	0.54
1:B:483:ILE:HG21	1:B:494:LEU:HD12	1.90	0.54
1:A:458:VAL:HB	1:A:459:PRO:HD2	1.90	0.54
1:A:502:GLY:HA3	1:B:52:MET:HE2	1.90	0.54
1:B:344:MET:HE1	1:B:380:ILE:HD13	1.89	0.54
1:B:291:ALA:O	1:B:295:LEU:HG	2.08	0.53
1:A:680:ARG:HB3	3:A:897:HOH:O	2.08	0.53
1:B:21:PRO:HB2	1:B:704:ARG:HA	1.89	0.53
1:B:226:GLU:HB3	3:B:1073:HOH:O	2.07	0.53
1:A:49:ASP:OD1	1:A:362:ARG:HD3	2.07	0.53
1:A:487:ASP:O	1:A:488:TRP:C	2.51	0.53
1:B:49:ASP:OD1	1:B:362:ARG:HD3	2.09	0.53
1:B:373:LEU:HA	1:B:376:PHE:CE2	2.44	0.53
1:B:376:PHE:CD1	1:B:376:PHE:C	2.87	0.53
1:B:683:ARG:HG2	1:B:683:ARG:HH21	1.74	0.53
1:A:390:ASP:OD2	1:A:410:LYS:HE3	2.08	0.52
1:B:73:TYR:HB2	1:B:147:PHE:HB2	1.89	0.52
1:B:390:ASP:HB3	1:B:405:TYR:OH	2.09	0.52
1:A:665:CYS:O	1:A:669:PRO:HD3	2.08	0.52
1:A:243:TRP:CE2	1:B:726:LYS:HE3	2.44	0.52
1:A:772:ARG:HG3	1:A:773:VAL:N	2.24	0.52
1:A:448:VAL:CG1	1:A:463:HIS:CE1	2.92	0.52
1:B:390:ASP:CG	1:B:410:LYS:HE3	2.35	0.52
1:A:483:ILE:HG21	1:A:494:LEU:HD12	1.92	0.51
1:A:657:TYR:HB3	1:B:169:SER:OG	2.09	0.51
1:B:546:ASP:HA	1:B:549:ARG:CZ	2.40	0.51
1:B:324:VAL:HG13	1:B:332:ASP:OD1	2.11	0.51
1:A:525:PHE:O	1:A:529:GLY:HA3	2.11	0.51
1:A:390:ASP:HB3	1:A:405:TYR:OH	2.10	0.51
1:A:25:ILE:HG21	1:A:355:PHE:CD1	2.45	0.51
1:B:25:ILE:HA	1:B:38:SER:HA	1.93	0.51
1:A:321:THR:HG22	1:A:772:ARG:HH21	1.76	0.51
1:A:373:LEU:HA	1:A:376:PHE:CE2	2.45	0.51
1:A:546:ASP:HA	1:A:549:ARG:CZ	2.41	0.51
1:A:71:TYR:CE1	1:A:87:PRO:HD3	2.46	0.51
1:B:487:ASP:O	1:B:488:TRP:C	2.54	0.51
1:A:5:HIS:NE2	1:A:14:VAL:HG11	2.26	0.50
1:A:373:LEU:HD23	1:A:377:VAL:HB	1.93	0.50
1:A:25:ILE:HA	1:A:38:SER:HA	1.94	0.50
1:A:169:SER:OG	1:B:657:TYR:HB3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:711:ALA:HA	1:A:728:SER:HA	1.93	0.50
1:B:352:SER:O	1:B:359:GLY:HA2	2.11	0.50
1:A:728:SER:O	1:A:729:TRP:HB2	2.11	0.50
1:A:403:HIS:ND1	1:A:420:PHE:CG	2.80	0.50
1:B:403:HIS:HB2	1:B:420:PHE:CD2	2.47	0.50
1:B:403:HIS:ND1	1:B:420:PHE:CG	2.80	0.50
1:A:461:PHE:HZ	1:A:536:ALA:HB2	1.77	0.50
1:B:461:PHE:HZ	1:B:536:ALA:HB2	1.77	0.50
1:A:726:LYS:HE3	1:B:243:TRP:CZ2	2.46	0.49
1:A:757:VAL:HB	1:A:795:LEU:HD21	1.94	0.49
1:B:14:VAL:HG23	1:B:98:ARG:HG2	1.93	0.49
1:A:291:ALA:O	1:A:295:LEU:HG	2.11	0.49
1:A:49:ASP:CG	1:A:362:ARG:HD3	2.38	0.49
1:B:769:THR:HA	1:B:781:ILE:O	2.13	0.49
1:B:297:ARG:HD2	3:B:1108:HOH:O	2.12	0.49
1:B:585:HIS:O	1:B:589:LYS:NZ	2.40	0.49
1:B:525:PHE:O	1:B:529:GLY:HA3	2.13	0.49
1:B:665:CYS:O	1:B:669:PRO:HD3	2.12	0.49
1:B:574:ASP:OD2	1:B:578:ASN:CB	2.55	0.49
1:A:105:THR:HA	1:A:117:THR:O	2.13	0.48
1:A:229:VAL:HB	1:A:230:PRO:HD3	1.95	0.48
1:B:49:ASP:CG	1:B:362:ARG:HD3	2.38	0.48
1:B:410:LYS:O	1:B:411:ARG:HD3	2.14	0.48
1:B:504:SER:HB3	1:B:507:THR:OG1	2.13	0.48
1:A:189:THR:O	1:A:190:GLU:HB2	2.14	0.48
1:A:393:SER:HB2	3:A:909:HOH:O	2.13	0.48
1:A:403:HIS:HB2	1:A:420:PHE:CD2	2.48	0.48
1:B:589:LYS:HE2	3:B:1105:HOH:O	2.13	0.48
1:B:32:GLN:HB2	1:B:126:ASN:OD1	2.14	0.48
1:B:728:SER:O	1:B:729:TRP:HB2	2.13	0.48
1:A:462:GLU:O	1:A:466:ARG:HG2	2.14	0.48
1:A:516:ALA:HA	1:A:573:TYR:O	2.14	0.48
1:A:769:THR:HA	1:A:781:ILE:O	2.14	0.48
3:A:1029:HOH:O	1:B:192:ARG:HD3	2.13	0.48
1:B:709:VAL:HG11	1:B:729:TRP:CZ2	2.49	0.48
1:A:410:LYS:O	1:A:411:ARG:HD3	2.14	0.47
1:B:150:VAL:CG2	1:B:201:PHE:CZ	2.97	0.47
1:B:175:VAL:HG21	1:B:210:PHE:CE1	2.49	0.47
1:B:420:PHE:CD2	1:B:420:PHE:N	2.82	0.47
1:A:635:LEU:CD2	1:A:667:ASN:HD21	2.27	0.47
1:A:652:THR:HG22	1:B:193:GLU:HB3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:PHE:CD1	1:B:130:GLN:HG2	2.49	0.47
1:A:493:ASN:HB2	1:A:648:GLY:HA3	1.96	0.47
1:B:25:ILE:HG21	1:B:355:PHE:CD1	2.49	0.47
1:B:70:ARG:H	1:B:70:ARG:HD2	1.80	0.47
1:A:256:ALA:HB3	1:A:259:GLU:OE1	2.15	0.47
1:A:406:GLN:HB3	3:A:1004:HOH:O	2.13	0.47
1:A:437:GLU:OE1	1:A:772:ARG:NH1	2.48	0.47
1:A:477:PRO:HD2	3:A:958:HOH:O	2.13	0.47
1:A:666:HIS:O	1:A:669:PRO:HD2	2.15	0.47
1:B:130:GLN:HB2	1:B:265:TYR:HB2	1.97	0.47
1:B:711:ALA:HA	1:B:728:SER:HA	1.96	0.47
1:A:683:ARG:HG2	1:A:683:ARG:HH21	1.80	0.47
1:B:256:ALA:O	1:B:259:GLU:HB3	2.15	0.47
1:B:650:VAL:O	1:B:660:ASN:HB2	2.15	0.47
1:A:119:PHE:CD1	1:A:130:GLN:HG2	2.50	0.47
1:A:324:VAL:HG13	1:A:332:ASP:OD1	2.14	0.47
1:B:409:THR:HG22	1:B:409:THR:O	2.15	0.47
1:A:256:ALA:O	1:A:259:GLU:HB3	2.15	0.47
1:A:273:PRO:HG2	1:A:276:GLU:CD	2.40	0.47
1:A:504:SER:HB3	1:A:507:THR:OG1	2.15	0.47
1:A:635:LEU:CD2	1:A:667:ASN:ND2	2.78	0.47
1:A:25:ILE:HG21	1:A:355:PHE:CE1	2.50	0.46
1:B:402:TYR:CD2	1:B:412:GLY:HA3	2.50	0.46
1:A:52:MET:HE2	1:B:502:GLY:HA3	1.96	0.46
1:B:40:GLN:O	1:B:41:ALA:HB3	2.15	0.46
1:B:28:LEU:HD11	1:B:37:LEU:HD22	1.97	0.46
1:B:373:LEU:HD23	1:B:373:LEU:O	2.16	0.46
1:A:112:GLY:HA3	1:A:138:SER:HB3	1.97	0.46
1:A:319:LEU:N	1:A:319:LEU:HD12	2.31	0.46
1:A:574:ASP:CG	1:A:575:TYR:N	2.74	0.46
1:B:593:GLU:HB2	1:B:594:PRO:HD3	1.98	0.46
1:A:40:GLN:O	1:A:41:ALA:HB3	2.14	0.46
1:A:175:VAL:HG21	1:A:210:PHE:CE1	2.50	0.46
1:A:751:GLU:HG2	1:A:756:VAL:HG23	1.97	0.46
1:B:25:ILE:HG21	1:B:355:PHE:CE1	2.50	0.46
1:B:703:HIS:CE1	1:B:725:ALA:HB2	2.51	0.46
1:A:243:TRP:CZ2	1:B:726:LYS:HE3	2.51	0.46
1:A:448:VAL:HG21	1:A:460:LEU:HD13	1.97	0.46
1:A:666:HIS:NE2	1:A:732:GLY:HA3	2.31	0.46
1:A:709:VAL:HG11	1:A:729:TRP:CZ2	2.51	0.46
1:A:502:GLY:HA3	1:B:52:MET:CE	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:LEU:CD1	1:A:541:LEU:N	2.78	0.46
1:A:762:GLY:C	1:A:787:GLY:HA3	2.41	0.46
1:B:319:LEU:HD12	1:B:319:LEU:N	2.30	0.46
1:B:448:VAL:HG12	1:B:463:HIS:NE2	2.30	0.46
1:B:49:ASP:OD2	1:B:52:MET:HB2	2.15	0.46
1:B:666:HIS:O	1:B:669:PRO:HD2	2.16	0.46
1:B:105:THR:HA	1:B:117:THR:O	2.15	0.46
1:A:571:ARG:HG3	1:A:571:ARG:HH11	1.80	0.46
1:A:193:GLU:OE1	1:B:652:THR:HB	2.15	0.45
1:A:420:PHE:CD2	1:A:420:PHE:N	2.83	0.45
1:B:541:LEU:N	1:B:541:LEU:CD1	2.79	0.45
1:B:101:LEU:CD1	1:B:101:LEU:N	2.80	0.45
1:A:37:LEU:HD23	1:A:104:SER:HB2	1.98	0.45
1:B:409:THR:HG22	1:B:411:ARG:HB2	1.98	0.45
1:B:409:THR:O	1:B:409:THR:CG2	2.64	0.45
1:A:708:TYR:HD1	1:A:709:VAL:HG13	1.81	0.45
1:B:5:HIS:O	1:B:14:VAL:HG12	2.16	0.45
1:B:627:LEU:HD22	1:B:634:VAL:HG23	1.98	0.45
1:B:255:LEU:HD22	1:B:259:GLU:HG2	1.99	0.45
1:A:272:ASN:OD1	1:A:290:ARG:HD2	2.16	0.45
1:A:703:HIS:CE1	1:A:725:ALA:HB2	2.52	0.45
1:B:169:SER:HB2	3:B:1121:HOH:O	2.16	0.45
1:A:571:ARG:HD2	1:A:594:PRO:CG	2.47	0.45
1:B:51:LYS:O	1:B:156:ASN:HA	2.17	0.45
1:B:583:ASP:O	1:B:589:LYS:NZ	2.50	0.45
1:A:49:ASP:OD2	1:A:52:MET:HB2	2.16	0.45
1:A:402:TYR:CD2	1:A:412:GLY:HA3	2.51	0.45
1:B:65:ALA:O	1:B:66:ASP:CB	2.64	0.45
1:B:367:ARG:HB2	1:B:426:TRP:CD1	2.52	0.44
1:B:526:VAL:HA	1:B:552:VAL:HG13	1.99	0.44
1:B:683:ARG:HG2	1:B:683:ARG:NH2	2.32	0.44
1:B:702:VAL:O	1:B:704:ARG:HG2	2.17	0.44
1:B:796:VAL:HG12	3:B:860:HOH:O	2.17	0.44
1:B:448:VAL:CG1	1:B:463:HIS:CE1	2.99	0.44
1:B:599:VAL:HB	1:B:674:ALA:HB1	1.99	0.44
1:B:708:TYR:HD1	1:B:709:VAL:HG13	1.83	0.44
1:A:319:LEU:N	1:A:319:LEU:CD1	2.81	0.44
1:A:449:PRO:HA	1:A:457:GLU:HA	2.00	0.44
1:B:762:GLY:C	1:B:787:GLY:HA3	2.42	0.44
1:B:810:ALA:HB1	1:B:811:PRO:HD2	1.98	0.44
1:A:351:ARG:HA	1:A:730:LEU:HD13	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:ALA:HB3	1:B:259:GLU:OE1	2.17	0.44
1:B:112:GLY:HA3	1:B:138:SER:HB3	1.98	0.44
1:B:212:THR:HA	1:B:246:ILE:O	2.18	0.44
1:B:462:GLU:O	1:B:466:ARG:HG2	2.17	0.44
1:A:160:ASP:OD1	1:A:160:ASP:C	2.61	0.44
1:A:432:ALA:O	1:A:436:LYS:HG3	2.17	0.44
1:A:810:ALA:HB1	1:A:811:PRO:HD2	1.99	0.44
1:B:190:GLU:O	1:B:193:GLU:HG3	2.18	0.44
1:B:449:PRO:HA	1:B:457:GLU:HA	2.00	0.44
1:B:751:GLU:HB2	1:B:754:GLY:O	2.18	0.44
1:A:330:LYS:O	1:A:334:MET:HG2	2.18	0.43
1:A:449:PRO:O	1:A:463:HIS:HE1	2.00	0.43
1:A:583:ASP:O	1:A:589:LYS:NZ	2.50	0.43
1:B:450:PHE:CZ	1:B:466:ARG:HG3	2.53	0.43
1:B:390:ASP:OD2	1:B:410:LYS:HE3	2.17	0.43
1:A:28:LEU:HD11	1:A:37:LEU:HD22	2.00	0.43
1:A:65:ALA:O	1:A:66:ASP:CB	2.66	0.43
1:A:71:TYR:CD1	1:A:87:PRO:HD3	2.53	0.43
1:A:212:THR:HA	1:A:246:ILE:O	2.18	0.43
1:B:319:LEU:N	1:B:319:LEU:CD1	2.81	0.43
1:B:574:ASP:OD1	1:B:576:TYR:N	2.37	0.43
1:A:393:SER:HB3	1:A:449:PRO:HG2	1.99	0.43
1:A:571:ARG:HD3	1:A:591:TRP:HB3	1.99	0.43
1:A:712:GLN:HG2	1:A:713:MET:HB2	2.01	0.43
1:B:619:ALA:O	1:B:623:VAL:HG23	2.18	0.43
1:A:121:VAL:HA	1:A:122:PRO:HD3	1.86	0.43
1:B:272:ASN:OD1	1:B:290:ARG:HD2	2.19	0.43
1:A:70:ARG:H	1:A:70:ARG:HD2	1.84	0.43
1:A:136:ASN:ND2	1:A:142:LYS:HG2	2.33	0.43
1:A:373:LEU:HD23	1:A:373:LEU:O	2.18	0.43
1:B:60:TYR:CZ	1:B:358:THR:HA	2.54	0.43
1:B:281:ASP:CG	3:B:1128:HOH:O	2.62	0.43
1:B:448:VAL:HG21	1:B:460:LEU:HD13	1.99	0.43
1:B:198:TYR:CE2	1:B:286:VAL:HG11	2.53	0.43
1:B:721:ARG:HD3	1:B:724:GLU:CD	2.43	0.43
1:B:808:PRO:HA	3:B:1067:HOH:O	2.18	0.43
1:A:33:PHE:HB2	1:A:126:ASN:HA	2.01	0.43
1:A:344:MET:HE1	1:A:380:ILE:HD13	2.01	0.43
1:A:517:GLU:O	1:A:572:ALA:HB1	2.19	0.43
1:A:757:VAL:HB	1:A:795:LEU:CD2	2.49	0.43
1:B:44:TYR:CD2	1:B:58:TYR:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:THR:O	1:B:190:GLU:HB2	2.18	0.42
1:A:448:VAL:HG12	1:A:463:HIS:NE2	2.34	0.42
1:A:652:THR:HG22	1:B:193:GLU:CB	2.49	0.42
1:B:217:PHE:HA	1:B:229:VAL:CG2	2.50	0.42
1:B:493:ASN:HB2	1:B:648:GLY:HA3	2.00	0.42
1:B:754:GLY:HA2	1:B:809:TYR:CD2	2.54	0.42
1:A:526:VAL:HA	1:A:552:VAL:HG13	2.00	0.42
1:A:635:LEU:C	1:A:635:LEU:CD1	2.91	0.42
1:B:497:PHE:CD1	1:B:511:GLN:HG3	2.54	0.42
1:A:14:VAL:O	1:A:14:VAL:HG13	2.20	0.42
1:A:91:ASP:O	1:A:110:ARG:HD2	2.20	0.42
1:A:513:GLY:C	1:A:575:TYR:CE2	2.98	0.42
1:A:696:ARG:HB2	3:A:1141:HOH:O	2.19	0.42
1:B:571:ARG:HG3	1:B:571:ARG:HH11	1.85	0.42
1:A:40:GLN:HB2	1:A:66:ASP:OD2	2.19	0.42
1:A:51:LYS:O	1:A:156:ASN:HA	2.18	0.42
1:B:541:LEU:N	1:B:541:LEU:HD12	2.34	0.42
1:A:749:ARG:HA	1:A:750:PRO:HD3	1.90	0.42
1:A:785:ASN:HA	1:A:822:LEU:HB2	2.01	0.42
1:A:130:GLN:HB2	1:A:265:TYR:HB2	2.02	0.42
1:B:6:PHE:CE1	1:B:333:ARG:HD3	2.55	0.42
1:A:122:PRO:HB2	1:A:125:GLU:HG3	2.02	0.42
1:A:627:LEU:HD22	1:A:634:VAL:HG23	2.02	0.42
1:B:59:ARG:HH21	1:B:59:ARG:HG3	1.85	0.42
1:B:330:LYS:O	1:B:334:MET:HG2	2.20	0.42
1:B:391:ILE:HG13	1:B:392:ALA:N	2.35	0.42
1:A:373:LEU:HD23	1:A:373:LEU:C	2.45	0.42
1:A:591:TRP:CD1	1:A:591:TRP:N	2.87	0.41
1:A:807:VAL:HA	1:A:808:PRO:HD3	1.91	0.41
1:B:136:ASN:ND2	1:B:142:LYS:HG2	2.34	0.41
1:B:409:THR:HG22	1:B:411:ARG:CB	2.50	0.41
1:B:637:TYR:CD1	1:B:638:PRO:HA	2.53	0.41
1:A:60:TYR:HE1	1:A:354:SER:O	2.03	0.41
1:A:169:SER:HB2	3:A:1114:HOH:O	2.19	0.41
1:A:646:HIS:ND1	1:A:647:MET:HG3	2.34	0.41
1:B:40:GLN:HB2	1:B:66:ASP:OD2	2.20	0.41
1:B:373:LEU:HD23	1:B:373:LEU:C	2.45	0.41
1:A:95:PHE:CD2	1:A:95:PHE:C	2.99	0.41
1:A:208:GLU:HB2	3:A:1092:HOH:O	2.20	0.41
1:B:95:PHE:CD2	1:B:95:PHE:C	2.99	0.41
1:B:145:THR:HG22	1:B:147:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:GLN:HB2	1:B:360:ILE:HG23	2.02	0.41
1:B:571:ARG:HD2	1:B:594:PRO:CG	2.50	0.41
1:B:703:HIS:C	1:B:704:ARG:HG2	2.44	0.41
1:A:197:HIS:HB3	1:A:271:GLU:HG2	2.02	0.41
1:A:405:TYR:O	1:A:407:PRO:HD3	2.20	0.41
1:A:658:LYS:HE3	1:A:658:LYS:HB3	1.94	0.41
1:B:547:ARG:NH1	1:B:547:ARG:HG2	2.34	0.41
1:A:154:LEU:HD12	1:A:167:ASN:HB2	2.01	0.41
1:B:37:LEU:HD23	1:B:104:SER:HB2	2.01	0.41
1:B:446:GLU:HA	1:B:447:PRO:HD3	1.92	0.41
1:A:164:TYR:HD2	1:B:358:THR:HG22	1.85	0.41
1:A:547:ARG:NH1	1:A:547:ARG:HG2	2.35	0.41
1:B:591:TRP:CD1	1:B:591:TRP:N	2.88	0.41
1:B:705:LEU:HD13	1:B:728:SER:HB3	2.03	0.41
1:A:367:ARG:HB2	1:A:426:TRP:CD1	2.55	0.41
1:A:434:TYR:CD1	1:A:434:TYR:C	2.98	0.41
1:A:497:PHE:CD1	1:A:511:GLN:HG3	2.56	0.41
1:A:541:LEU:N	1:A:541:LEU:HD12	2.35	0.41
1:B:63:ILE:HA	1:B:64:PRO:HA	1.87	0.41
1:B:384:ALA:O	1:B:388:ILE:HG13	2.21	0.41
1:A:4:GLY:HA3	1:A:14:VAL:O	2.21	0.41
1:A:244:TYR:HA	1:A:245:PRO:HD3	1.83	0.41
1:A:705:LEU:HD11	1:A:726:LYS:HA	2.03	0.41
1:B:28:LEU:HD22	1:B:103:TYR:HA	2.03	0.41
1:B:193:GLU:CG	1:B:194:ARG:H	2.32	0.41
1:B:214:ARG:HB2	1:B:246:ILE:HG23	2.03	0.41
1:B:399:GLY:HA2	1:B:466:ARG:HB3	2.02	0.41
1:B:423:ASP:N	1:B:424:PRO:CD	2.84	0.41
1:B:539:ARG:NH1	3:B:883:HOH:O	2.52	0.41
1:B:749:ARG:HA	1:B:750:PRO:HD3	1.90	0.41
1:A:52:MET:CE	1:B:502:GLY:HA3	2.51	0.41
1:A:321:THR:HG23	3:A:1132:HOH:O	2.21	0.41
1:A:488:TRP:NE1	3:A:913:HOH:O	2.54	0.41
1:A:28:LEU:HD22	1:A:103:TYR:HA	2.02	0.40
1:A:198:TYR:CE2	1:A:286:VAL:HG11	2.56	0.40
1:A:505:PHE:CG	1:A:649:GLU:HB2	2.57	0.40
1:B:61:ASN:HA	3:B:1022:HOH:O	2.20	0.40
1:B:282:ALA:O	1:B:283:LYS:HB2	2.21	0.40
1:B:392:ALA:HB2	1:B:427:LEU:HD11	2.02	0.40
1:A:12:GLU:OE1	1:A:98:ARG:HD3	2.21	0.40
1:A:665:CYS:O	1:A:669:PRO:CD	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:TYR:CD2	1:A:58:TYR:HB3	2.56	0.40
1:A:314:TYR:HD2	1:A:315:TRP:CD1	2.39	0.40
1:A:581:GLY:HA2	1:A:588:GLY:O	2.22	0.40
1:B:308:PHE:HD2	3:B:1103:HOH:O	2.04	0.40
1:B:393:SER:HB3	1:B:449:PRO:HG2	2.02	0.40
1:A:146:LEU:O	1:A:250:SER:HA	2.21	0.40
1:A:373:LEU:HD11	1:A:388:ILE:CG1	2.51	0.40
1:A:599:VAL:HB	1:A:674:ALA:HB1	2.02	0.40
1:A:669:PRO:HA	1:A:672:ILE:HD12	2.04	0.40
1:B:146:LEU:O	1:B:250:SER:HA	2.21	0.40
1:B:154:LEU:HD12	1:B:167:ASN:HB2	2.02	0.40
1:B:197:HIS:HB3	1:B:271:GLU:HG2	2.04	0.40
1:B:511:GLN:NE2	1:B:576:TYR:OH	2.52	0.40
1:B:589:LYS:HE3	3:B:889:HOH:O	2.20	0.40
1:A:29:GLY:HA3	1:A:33:PHE:O	2.22	0.40
1:A:63:ILE:HA	1:A:64:PRO:HA	1.85	0.40
1:A:650:VAL:O	1:A:660:ASN:HB2	2.21	0.40
1:A:683:ARG:HG2	1:A:683:ARG:NH2	2.36	0.40
1:B:385:ARG:CG	1:B:443:ILE:HD12	2.40	0.40
1:B:665:CYS:O	1:B:666:HIS:C	2.65	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:TYR:OH	1:A:686:ASP:OD2[2_556]	2.17	0.03

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	820/822 (100%)	781 (95%)	39 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	821/822 (100%)	783 (95%)	38 (5%)	0	100	100
All	All	1641/1644 (100%)	1564 (95%)	77 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	666/675 (99%)	655 (98%)	11 (2%)	53	74
1	B	669/675 (99%)	658 (98%)	11 (2%)	55	76
All	All	1335/1350 (99%)	1313 (98%)	22 (2%)	55	76

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	VAL
1	A	221	TYR
1	A	239	VAL
1	A	276	GLU
1	A	312	LYS
1	A	543	ASP
1	A	574	ASP
1	A	610	GLN
1	A	635	LEU
1	A	664	PHE
1	A	678	VAL
1	B	101	LEU
1	B	123	VAL
1	B	221	TYR
1	B	312	LYS
1	B	320	SER
1	B	543	ASP
1	B	610	GLN

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Mol	Chain	Res	Type
1	B	617	ILE
1	B	635	LEU
1	B	664	PHE
1	B	822	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
1	A	177	GLN
1	A	181	HIS
1	A	292	HIS
1	A	551	HIS
1	A	646	HIS
1	A	666	HIS
1	A	760	GLN
1	B	94	HIS
1	B	158	GLN
1	B	181	HIS
1	B	292	HIS
1	B	511	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	GLC	B	824	-	12,12,12	0.57	0	17,17,17	0.54	0
2	GLC	A	824	-	12,12,12	0.58	0	17,17,17	0.61	0
2	GLC	A	823	-	12,12,12	0.61	0	17,17,17	0.64	0
2	GLC	B	823	-	12,12,12	0.72	0	17,17,17	1.40	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	B	824	-	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	A	824	-	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	A	823	-	1/1/5/5	2/2/22/22	0/1/1/1
2	GLC	B	823	-	1/1/5/5	2/2/22/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	823	GLC	O3-C3-C2	-2.84	103.67	110.38
2	B	823	GLC	O2-C2-C3	-2.70	104.01	110.38
2	B	823	GLC	O2-C2-C1	2.07	114.03	109.25

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	823	GLC	C1
2	A	824	GLC	C1
2	B	823	GLC	C1
2	B	824	GLC	C1

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	823	GLC	O5-C5-C6-O6
2	B	823	GLC	C4-C5-C6-O6
2	A	823	GLC	C4-C5-C6-O6
2	A	823	GLC	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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	822/822 (100%)	0.29	20 (2%) 59 55	8, 18, 47, 105	1 (0%)
1	B	822/822 (100%)	0.21	26 (3%) 50 46	8, 18, 48, 105	2 (0%)
All	All	1644/1644 (100%)	0.25	46 (2%) 55 51	8, 18, 48, 105	3 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	442	SER	4.4
1	A	409	THR	4.2
1	A	612	THR	4.0
1	A	414	ASN	3.7
1	B	409	THR	3.7
1	B	414	ASN	3.6
1	A	610	GLN	3.5
1	B	415	ASP	3.4
1	B	612	THR	3.3
1	A	281	ASP	3.3
1	B	608	GLY	3.1
1	A	543	ASP	2.8
1	B	442	SER	2.8
1	B	281	ASP	2.8
1	A	363	GLY	2.7
1	B	413	ASN	2.6
1	A	513	GLY	2.6
1	B	363	GLY	2.6
1	B	543	ASP	2.6
1	B	301	SER	2.5
1	A	405	TYR	2.5
1	B	610	GLN	2.4
1	B	509	GLU	2.4
1	B	221	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	536	ALA	2.4
1	B	512	ALA	2.4
1	B	365	GLY	2.3
1	A	404	GLN	2.3
1	A	221	TYR	2.3
1	A	273	PRO	2.3
1	A	511	GLN	2.2
1	B	657	TYR	2.2
1	B	822	LEU	2.2
1	B	790	GLY	2.2
1	B	453	GLU	2.1
1	A	455	GLY	2.1
1	A	547	ARG	2.0
1	B	513	GLY	2.0
1	A	139	ASP	2.0
1	B	445	ASP	2.0
1	B	404	GLN	2.0
1	B	511	GLN	2.0
1	A	270	VAL	2.0
1	A	530	GLU	2.0
1	A	274	ASP	2.0
1	B	325	SER	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	B	823	12/12	0.90	0.11	7,20,30,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLC	A	823	12/12	0.91	0.10	6,15,26,32	0
2	GLC	B	824	12/12	0.93	0.09	10,13,18,22	0
2	GLC	A	824	12/12	0.95	0.07	6,13,18,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.