Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2024 – 02:52 PM JST

PDB ID	:	8WV2
Title	:	Crystal structure of urethanase from Candida parapsilosis and structure-based Engineering to improve the catalytic activity and stability
Authors	:	Zhao, T.; Wu, H.
Deposited on	:	2023-10-22
Resolution	:	2.66 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:

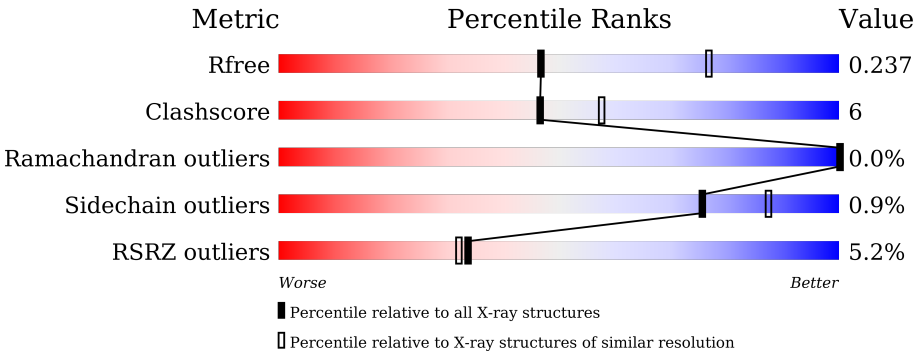
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.66 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}	164625	1003 (2.66-2.66)
Clashscore	180529	1063 (2.66-2.66)
Ramachandran outliers	177936	1052 (2.66-2.66)
Sidechain outliers	177891	1052 (2.66-2.66)
RSRZ outliers	164620	1003 (2.66-2.66)

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 <=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	551	89% 11%
1	B	551	89% 11%
1	C	551	88% 12%
1	D	551	30% 79% 21%
1	E	551	89% 11%
1	F	551	86% 13%

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Mol	Chain	Length	Quality of chain
1	G	551	2% 87% 13%
1	H	551	3% 88% 12%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 35789 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 Amidase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	B	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	C	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	E	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	F	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	G	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	H	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			
1	D	551	Total	C	N	O	S	0	0	0
			4354	2795	720	822	17			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	161	Total	O	0	0
			161	161		
2	B	117	Total	O	0	0
			117	117		
2	C	100	Total	O	0	0
			100	100		
2	E	148	Total	O	0	0
			148	148		
2	F	139	Total	O	0	0
			139	139		
2	G	117	Total	O	0	0
			117	117		

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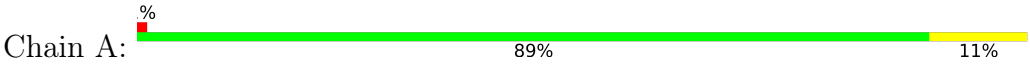
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	H	104	Total O 104 104	0	0
2	D	71	Total O 71 71	0	0

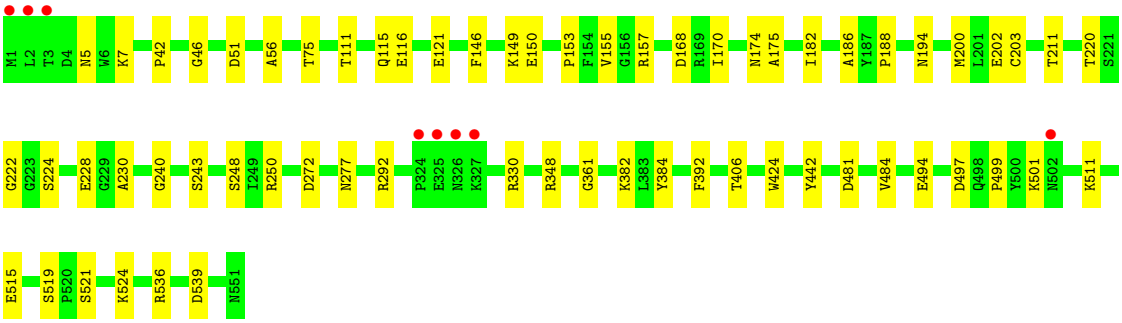
3 Residue-property plots (i)

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: Amidase family protein

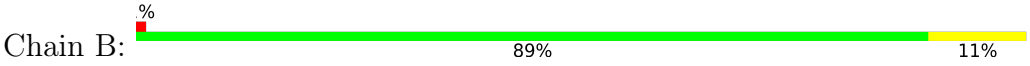
Chain A:

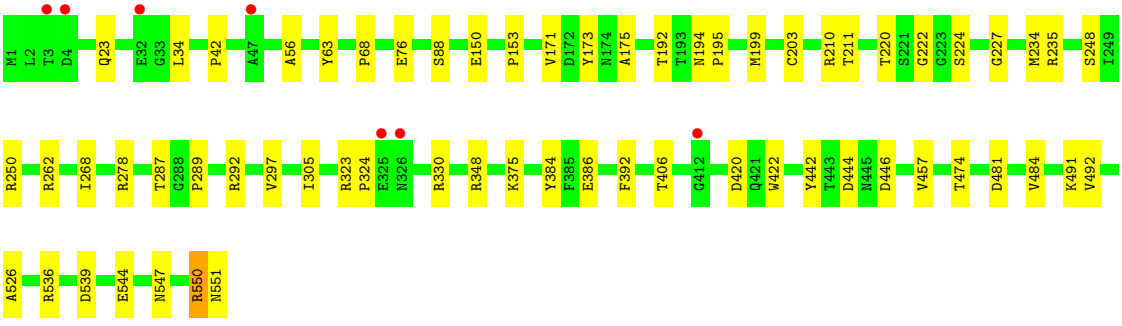




• Molecule 1: Amidase family protein

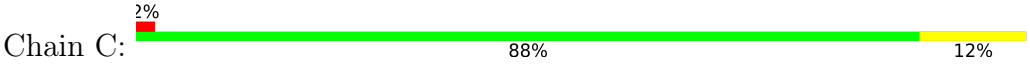
Chain B:

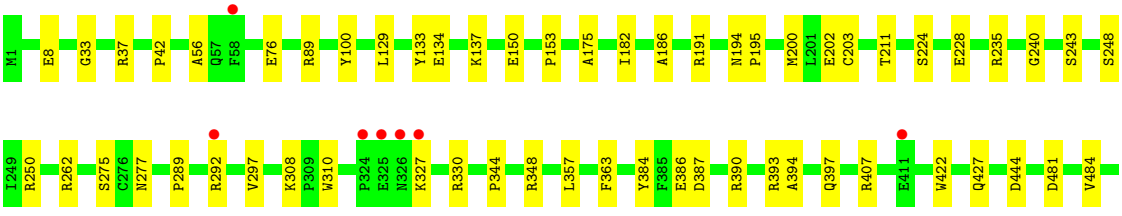


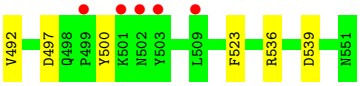


• Molecule 1: Amidase family protein

Chain C:

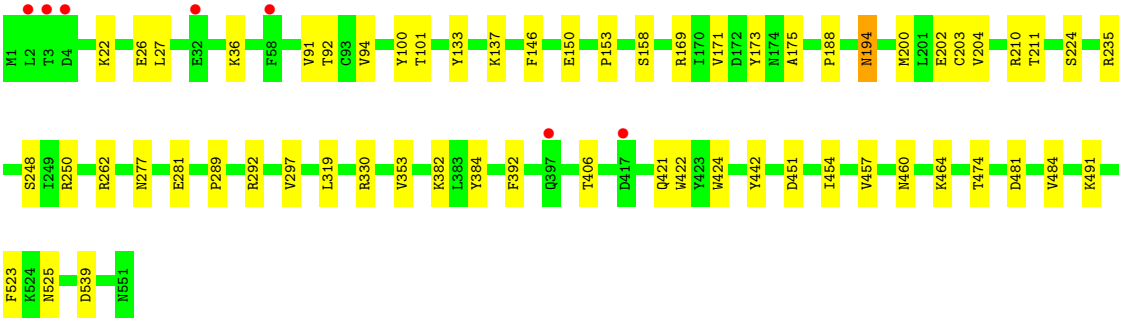






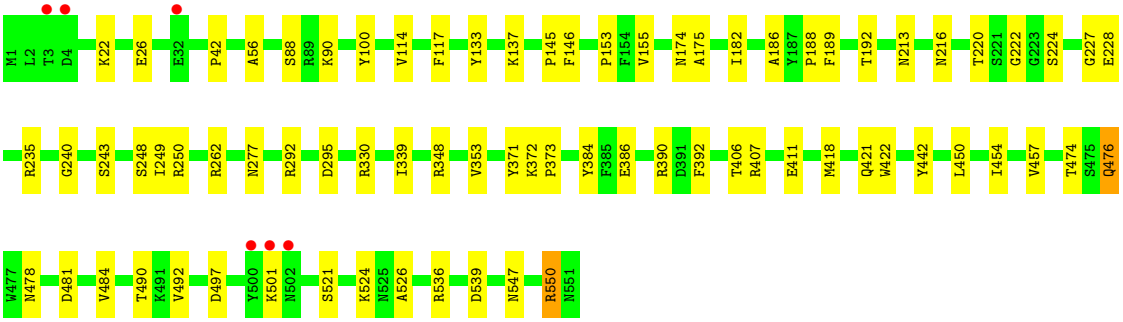
• Molecule 1: Amidase family protein

Chain E:



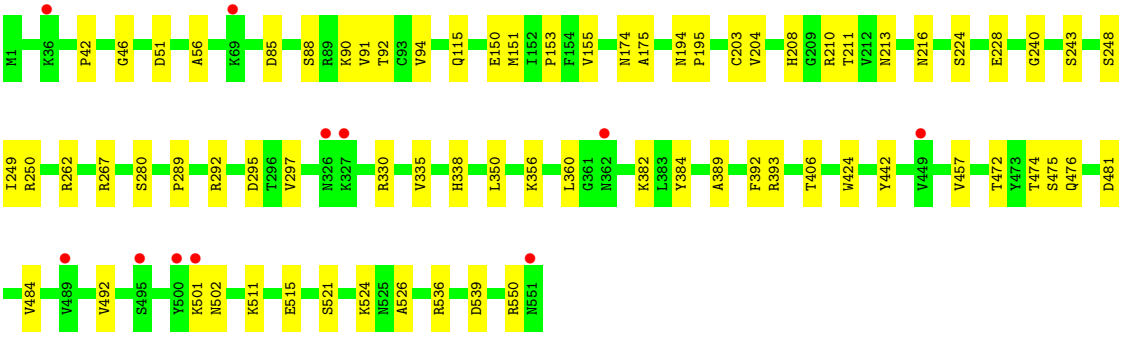
• Molecule 1: Amidase family protein

Chain F:



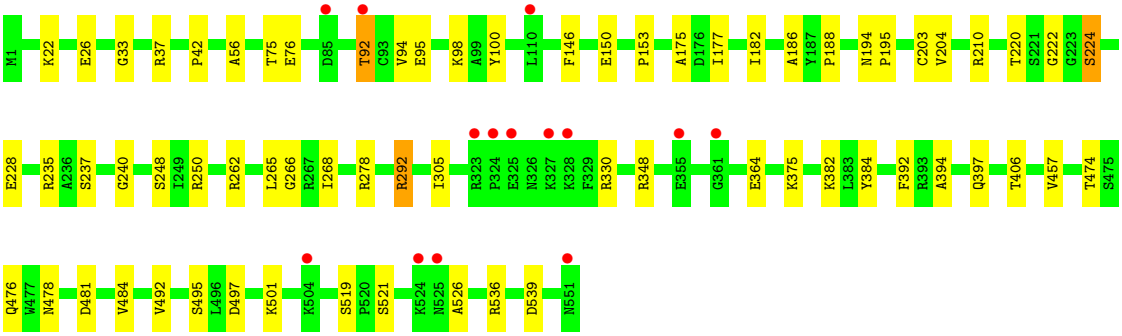
• Molecule 1: Amidase family protein

Chain G:



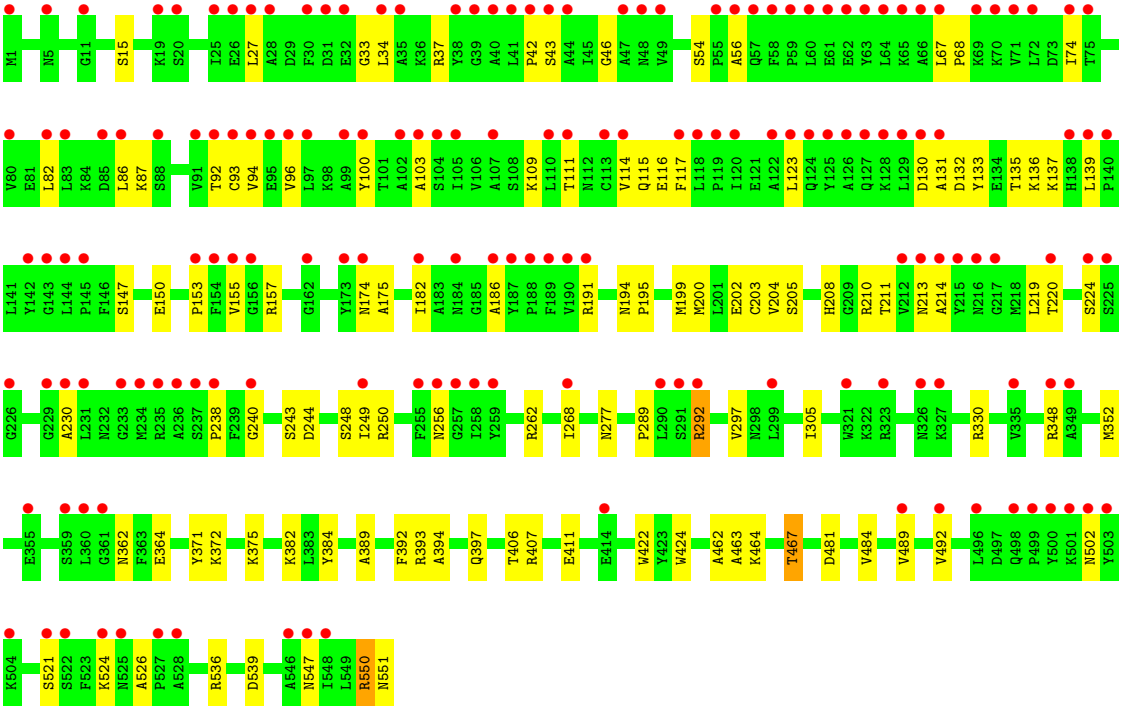
• Molecule 1: Amidase family protein

Chain H:



• Molecule 1: Amidase family protein

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	63.42Å 110.16Å 180.58Å 75.84° 83.62° 73.95°	Depositor
Resolution (Å)	48.91 – 2.66 48.91 – 2.66	Depositor EDS
% Data completeness (in resolution range)	98.3 (48.91-2.66) 98.3 (48.91-2.66)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.70 (at 2.65Å)	Xtriage
Refinement program	PHENIX (1.16_3549: ???)	Depositor
R, R_{free}	0.204 , 0.237 0.204 , 0.237	Depositor DCC
R_{free} test set	6428 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	29.9	Xtriage
Anisotropy	1.043	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 41.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	35789	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.33	0/4460	0.52	0/6061
1	B	0.35	0/4460	0.53	0/6061
1	C	0.31	0/4460	0.52	0/6061
1	D	0.36	0/4460	0.55	0/6061
1	E	0.33	0/4460	0.51	0/6061
1	F	0.33	0/4460	0.51	0/6061
1	G	0.36	0/4460	0.53	0/6061
1	H	0.34	0/4460	0.52	0/6061
All	All	0.34	0/35680	0.52	0/48488

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

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	4354	0	4331	46	0
1	B	4354	0	4331	40	0
1	C	4354	0	4331	48	0
1	D	4354	0	4331	84	0
1	E	4354	0	4331	42	0
1	F	4354	0	4331	51	0
1	G	4354	0	4331	41	0
1	H	4354	0	4331	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	161	0	0	5	0
2	B	117	0	0	9	0
2	C	100	0	0	12	0
2	D	71	0	0	6	0
2	E	148	0	0	5	0
2	F	139	0	0	6	0
2	G	117	0	0	4	0
2	H	104	0	0	6	0
All	All	35789	0	34648	389	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 (389) 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:E:171:VAL:HG13	1:E:173:TYR:CE2	2.00	0.96
1:A:111:THR:HG21	1:A:230:ALA:HB1	1.62	0.81
1:F:235:ARG:HG2	1:F:292:ARG:HD3	1.64	0.80
1:E:171:VAL:HG12	1:E:173:TYR:H	1.49	0.77
1:E:171:VAL:HG13	1:E:173:TYR:CD2	2.21	0.74
1:G:330:ARG:NH2	1:G:442:TYR:O	2.20	0.74
1:A:153:PRO:HD2	1:A:175:ALA:HB2	1.70	0.74
1:B:150:GLU:HB3	1:B:194:ASN:ND2	2.03	0.74
1:G:338:HIS:NE2	2:G:602:HOH:O	2.21	0.74
1:E:235:ARG:HG2	1:E:292:ARG:HD3	1.70	0.73
1:G:150:GLU:HB3	1:G:194:ASN:ND2	2.03	0.73
1:C:224:SER:HB3	1:C:248:SER:HB3	1.70	0.72
1:G:153:PRO:HD2	1:G:175:ALA:HB2	1.71	0.72
1:F:153:PRO:HD2	1:F:175:ALA:HB2	1.71	0.71
1:E:330:ARG:NH2	1:E:442:TYR:O	2.24	0.70
1:D:362:ASN:ND2	2:D:601:HOH:O	2.23	0.70
1:D:150:GLU:HB3	1:D:194:ASN:ND2	2.07	0.70
1:E:171:VAL:CG1	1:E:173:TYR:CD2	2.75	0.69
1:G:267:ARG:NH1	2:G:604:HOH:O	2.23	0.69
1:H:330:ARG:NH1	1:H:364:GLU:OE2	2.25	0.69
1:D:132:ASP:HB3	1:D:135:THR:OG1	1.92	0.69
1:D:199:MET:SD	2:D:669:HOH:O	2.51	0.69
1:C:292:ARG:HD2	2:C:635:HOH:O	1.93	0.69
1:E:171:VAL:CG1	1:E:173:TYR:CE2	2.75	0.69
1:G:228:GLU:HG2	2:G:614:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:GLU:HB3	1:C:194:ASN:ND2	2.09	0.68
1:C:153:PRO:HD2	1:C:175:ALA:HB2	1.74	0.68
1:D:33:GLY:O	1:D:37:ARG:HG3	1.93	0.68
1:D:92:THR:HG23	1:D:130:ASP:CG	2.14	0.68
1:C:150:GLU:HB3	1:C:194:ASN:HD21	1.57	0.68
1:D:362:ASN:ND2	2:D:603:HOH:O	2.26	0.68
1:B:150:GLU:HB3	1:B:194:ASN:HD21	1.58	0.68
1:A:330:ARG:NH2	1:A:442:TYR:O	2.26	0.68
1:C:481:ASP:OD2	1:C:536:ARG:NH1	2.27	0.68
1:A:75:THR:O	1:A:292:ARG:NH2	2.25	0.67
1:F:100:TYR:OH	1:F:292:ARG:NH2	2.28	0.67
1:D:86:LEU:HD21	1:D:96:VAL:HG21	1.77	0.66
1:C:100:TYR:OH	1:C:292:ARG:NH2	2.29	0.66
1:F:292:ARG:NH1	2:F:605:HOH:O	2.28	0.66
1:B:153:PRO:HD2	1:B:175:ALA:HB2	1.78	0.66
1:D:74:ILE:HG23	1:D:82:LEU:HD21	1.78	0.65
1:B:330:ARG:NH2	1:B:442:TYR:O	2.29	0.65
1:F:277:ASN:ND2	2:F:602:HOH:O	2.23	0.65
1:E:100:TYR:OH	1:E:292:ARG:NH2	2.29	0.64
1:D:213:ASN:ND2	1:D:256:ASN:OD1	2.30	0.64
1:C:492:VAL:HG21	1:C:523:PHE:O	1.96	0.64
1:H:153:PRO:HD2	1:H:175:ALA:HB2	1.80	0.64
1:D:46:GLY:N	2:D:610:HOH:O	2.31	0.63
1:G:295:ASP:OD1	1:G:550:ARG:NH2	2.30	0.63
1:E:153:PRO:HD2	1:E:175:ALA:HB2	1.81	0.62
1:D:371:TYR:OH	1:D:372:LYS:NZ	2.33	0.62
1:C:134:GLU:OE1	1:C:137:LYS:NZ	2.26	0.62
1:H:150:GLU:HB3	1:H:194:ASN:ND2	2.15	0.62
1:F:547:ASN:OD1	1:F:550:ARG:NH1	2.32	0.61
1:A:224:SER:OG	1:A:248:SER:HB3	2.01	0.61
1:G:92:THR:HG22	1:G:94:VAL:H	1.65	0.61
1:D:375:LYS:NZ	2:D:612:HOH:O	2.33	0.61
1:D:407:ARG:O	1:D:411:GLU:HG2	2.00	0.61
1:G:150:GLU:HB3	1:G:194:ASN:HD21	1.66	0.60
1:G:224:SER:H	1:G:248:SER:HB3	1.66	0.60
1:D:150:GLU:HB3	1:D:194:ASN:HD21	1.66	0.60
1:G:501:LYS:HG2	1:G:502:ASN:N	2.17	0.60
1:C:235:ARG:HG2	1:C:292:ARG:HD3	1.84	0.60
1:E:92:THR:HG22	1:E:94:VAL:H	1.67	0.59
1:B:323:ARG:HG3	1:B:324:PRO:HD2	1.84	0.59
1:B:250:ARG:HD3	1:B:484:VAL:HB	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:150:GLU:HB3	1:E:194:ASN:HD21	1.68	0.58
1:D:289:PRO:HG2	1:D:297:VAL:HG13	1.86	0.58
1:H:262:ARG:HE	1:H:478:ASN:ND2	2.02	0.58
1:H:268:ILE:HD11	1:H:305:ILE:HG13	1.86	0.58
1:D:136:LYS:HB3	1:D:139:LEU:HD12	1.85	0.57
1:H:75:THR:O	1:H:292:ARG:NH2	2.37	0.57
1:B:220:THR:HG22	1:B:222:GLY:H	1.69	0.57
1:D:204:VAL:HG22	1:D:210:ARG:HG3	1.87	0.57
1:D:250:ARG:HD3	1:D:484:VAL:HB	1.87	0.57
1:D:348:ARG:NH1	2:D:609:HOH:O	2.31	0.57
1:D:330:ARG:HG2	1:D:364:GLU:HB3	1.87	0.56
1:A:494:GLU:OE2	1:A:524:LYS:HE2	2.04	0.56
1:H:92:THR:CG2	1:H:95:GLU:H	2.17	0.56
1:D:492:VAL:HG23	1:D:526:ALA:O	2.05	0.56
1:D:116:GLU:OE1	1:D:191:ARG:NH1	2.37	0.56
1:E:250:ARG:HD3	1:E:484:VAL:HB	1.87	0.56
1:E:200:MET:CE	1:E:202:GLU:HB2	2.36	0.56
1:H:204:VAL:HG22	1:H:210:ARG:HG3	1.87	0.56
1:H:224:SER:H	1:H:248:SER:HB3	1.71	0.56
1:D:153:PRO:HD2	1:D:175:ALA:HB2	1.87	0.55
1:E:150:GLU:HB3	1:E:194:ASN:ND2	2.22	0.55
1:F:330:ARG:NH2	1:F:442:TYR:O	2.39	0.55
1:A:499:PRO:HB3	2:A:644:HOH:O	2.06	0.55
1:C:33:GLY:O	1:C:37:ARG:HG3	2.06	0.54
1:A:224:SER:H	1:A:248:SER:HB3	1.71	0.54
1:A:42:PRO:HG2	1:A:56:ALA:HB2	1.89	0.54
1:G:382:LYS:HD2	1:G:424:TRP:NE1	2.22	0.54
1:H:33:GLY:O	1:H:37:ARG:HG3	2.07	0.54
1:D:195:PRO:HA	1:D:205:SER:HB3	1.90	0.54
1:C:89:ARG:CD	2:C:695:HOH:O	2.55	0.54
1:D:87:LYS:HA	1:D:137:LYS:HG3	1.89	0.54
1:C:89:ARG:HD2	2:C:695:HOH:O	2.08	0.54
1:A:149:LYS:NZ	1:A:224:SER:OG	2.36	0.54
1:D:34:LEU:HD22	1:D:68:PRO:HD3	1.90	0.54
1:A:250:ARG:HD3	1:A:484:VAL:HB	1.90	0.54
1:B:171:VAL:HG12	1:B:173:TYR:H	1.72	0.54
1:G:250:ARG:HD3	1:G:484:VAL:HB	1.90	0.53
1:E:460:ASN:ND2	2:E:623:HOH:O	2.39	0.53
1:D:203:CYS:HB3	1:D:211:THR:HB	1.89	0.53
1:D:213:ASN:HD22	1:D:214:ALA:H	1.54	0.53
1:A:511:LYS:HE2	1:A:515:GLU:OE2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:ILE:HD11	1:B:305:ILE:HG13	1.91	0.53
1:F:22:LYS:O	1:F:26:GLU:HG3	2.09	0.53
1:D:42:PRO:HG2	1:D:56:ALA:HB2	1.89	0.53
1:F:348:ARG:NH1	1:F:497:ASP:OD2	2.42	0.53
1:D:224:SER:H	1:D:248:SER:HB3	1.74	0.53
1:G:42:PRO:HG2	1:G:56:ALA:HB2	1.90	0.53
1:G:492:VAL:HG23	1:G:526:ALA:O	2.08	0.53
1:A:519:SER:OG	1:A:521:SER:OG	2.27	0.53
1:F:224:SER:HB3	1:F:248:SER:HB3	1.90	0.53
1:B:420:ASP:OD2	1:C:327:LYS:NZ	2.36	0.53
1:C:191:ARG:HG2	1:C:191:ARG:HH11	1.73	0.52
1:H:250:ARG:HD3	1:H:484:VAL:HB	1.91	0.52
1:E:200:MET:HE2	1:E:202:GLU:HB2	1.92	0.52
1:F:348:ARG:NH2	1:F:490:THR:OG1	2.42	0.52
1:F:547:ASN:O	1:F:550:ARG:HG3	2.10	0.52
1:D:463:ALA:HB1	1:D:467:THR:HG22	1.91	0.52
1:F:407:ARG:O	1:F:411:GLU:HG2	2.08	0.52
1:D:93:CYS:N	1:D:130:ASP:OD1	2.29	0.52
1:D:230:ALA:HA	1:D:256:ASN:ND2	2.25	0.52
1:F:250:ARG:HD3	1:F:484:VAL:HB	1.91	0.52
1:H:492:VAL:HG23	1:H:526:ALA:O	2.09	0.52
1:G:195:PRO:HB3	1:G:203:CYS:HA	1.91	0.52
1:A:200:MET:CE	1:A:202:GLU:HB2	2.40	0.52
1:B:481:ASP:OD2	1:B:536:ARG:NH1	2.43	0.52
1:C:308:LYS:HG2	1:C:310:TRP:CZ2	2.45	0.51
1:H:476:GLN:HG3	1:H:476:GLN:O	2.09	0.51
1:D:463:ALA:HB1	1:D:467:THR:CG2	2.40	0.51
1:B:457:VAL:O	1:B:474:THR:HG23	2.10	0.51
1:C:203:CYS:HB3	1:C:211:THR:HB	1.91	0.51
1:F:42:PRO:HG2	1:F:56:ALA:HB2	1.93	0.51
1:F:492:VAL:HG23	1:F:526:ALA:O	2.09	0.51
1:B:192:THR:HB	1:B:227:GLY:HA3	1.93	0.51
1:E:36:LYS:NZ	2:E:613:HOH:O	2.28	0.51
1:B:550:ARG:O	1:B:551:ASN:HB2	2.11	0.51
1:A:220:THR:HG23	1:A:222:GLY:H	1.76	0.51
1:B:287:THR:HG22	2:B:646:HOH:O	2.09	0.51
1:E:133:TYR:O	1:E:137:LYS:HB2	2.11	0.51
1:A:203:CYS:HB3	1:A:211:THR:HB	1.93	0.51
1:D:87:LYS:HG3	1:D:137:LYS:HG2	1.93	0.50
1:H:348:ARG:NH1	1:H:497:ASP:OD1	2.44	0.50
1:C:289:PRO:HG2	1:C:297:VAL:HG13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:67:LEU:HD11	1:D:103:ALA:HA	1.92	0.50
1:B:171:VAL:HG13	1:B:173:TYR:CD2	2.47	0.50
1:E:262:ARG:NH2	1:E:481:ASP:HA	2.27	0.50
1:D:115:GLN:OE1	1:D:208:HIS:HB3	2.12	0.50
1:C:89:ARG:HG3	2:C:695:HOH:O	2.11	0.50
1:C:200:MET:HG2	1:C:202:GLU:H	1.75	0.50
1:E:203:CYS:HB3	1:E:211:THR:HB	1.92	0.50
1:A:228:GLU:OE1	1:A:240:GLY:HA3	2.12	0.50
1:B:76:GLU:OE1	1:B:235:ARG:NH1	2.45	0.50
1:E:224:SER:H	1:E:248:SER:HB3	1.76	0.49
1:D:27:LEU:HD11	1:D:123:LEU:HB2	1.92	0.49
1:H:42:PRO:HG2	1:H:56:ALA:HB2	1.94	0.49
1:D:155:VAL:HG23	1:D:174:ASN:OD1	2.12	0.49
1:D:224:SER:OG	1:D:248:SER:HB3	2.12	0.49
1:C:387:ASP:O	1:C:390:ARG:HD3	2.11	0.49
1:A:200:MET:HE3	1:A:202:GLU:HB2	1.93	0.49
1:F:224:SER:H	1:F:248:SER:HB3	1.77	0.49
1:H:237:SER:O	1:H:292:ARG:NH1	2.45	0.49
1:E:422:TRP:CZ3	1:F:277:ASN:HA	2.48	0.49
1:A:111:THR:CG2	1:A:230:ALA:HB1	2.39	0.49
1:C:262:ARG:NH1	1:C:481:ASP:HA	2.27	0.48
1:G:521:SER:O	1:G:524:LYS:HG3	2.13	0.48
1:F:182:ILE:HA	1:F:186:ALA:O	2.12	0.48
1:F:295:ASP:OD1	1:F:550:ARG:NH2	2.47	0.48
1:G:228:GLU:OE1	1:G:240:GLY:HA3	2.13	0.48
1:G:472:THR:HA	1:G:475:SER:OG	2.13	0.48
1:H:22:LYS:NZ	1:H:26:GLU:OE2	2.36	0.48
1:E:146:PHE:O	1:E:188:PRO:HA	2.14	0.48
1:E:451:ASP:O	2:E:601:HOH:O	2.20	0.48
1:H:195:PRO:HB3	1:H:203:CYS:HA	1.95	0.48
1:B:392:PHE:CZ	1:B:406:THR:HG21	2.49	0.48
1:C:277:ASN:HA	1:D:422:TRP:CZ3	2.48	0.48
1:F:386:GLU:OE2	1:F:421:GLN:NE2	2.43	0.48
1:H:497:ASP:OD1	2:H:601:HOH:O	2.20	0.48
1:B:278:ARG:HD2	2:B:603:HOH:O	2.13	0.48
1:D:92:THR:HG23	1:D:130:ASP:OD1	2.13	0.48
1:A:220:THR:HG23	1:A:222:GLY:N	2.29	0.48
1:A:150:GLU:HB3	1:A:194:ASN:HD21	1.78	0.48
1:F:220:THR:OG1	2:F:601:HOH:O	2.20	0.48
1:F:521:SER:O	1:F:524:LYS:HG3	2.14	0.47
1:A:392:PHE:CZ	1:A:406:THR:HG21	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:CYS:HB3	1:B:211:THR:HB	1.96	0.47
1:D:268:ILE:HD11	1:D:305:ILE:HG13	1.95	0.47
1:H:519:SER:OG	1:H:521:SER:OG	2.29	0.47
1:D:87:LYS:CG	1:D:137:LYS:HG2	2.44	0.47
1:E:27:LEU:HD23	1:E:101:THR:HG21	1.96	0.47
1:E:171:VAL:HG12	1:E:173:TYR:N	2.25	0.47
1:F:250:ARG:NH1	2:F:617:HOH:O	2.44	0.47
1:B:34:LEU:HD13	1:B:68:PRO:HD3	1.95	0.47
1:B:88:SER:O	2:B:601:HOH:O	2.20	0.47
1:D:262:ARG:NH2	1:D:481:ASP:HA	2.29	0.47
1:B:63:TYR:HB2	1:B:234:MET:HE1	1.96	0.47
1:B:375:LYS:NZ	2:B:622:HOH:O	2.47	0.47
1:D:382:LYS:HG3	1:D:424:TRP:CE2	2.50	0.47
1:D:521:SER:O	1:D:524:LYS:HG3	2.15	0.47
1:F:353:VAL:HG11	1:F:454:ILE:HD13	1.97	0.46
1:H:481:ASP:OD2	1:H:536:ARG:NH1	2.48	0.46
1:B:224:SER:H	1:B:248:SER:HB3	1.80	0.46
1:H:228:GLU:OE1	1:H:240:GLY:HA3	2.15	0.46
1:E:277:ASN:HA	1:F:422:TRP:CZ3	2.51	0.46
1:A:348:ARG:NH1	1:A:497:ASP:OD2	2.47	0.46
1:G:389:ALA:O	1:G:393:ARG:HG3	2.16	0.46
1:H:92:THR:HG23	1:H:95:GLU:H	1.79	0.46
1:D:243:SER:O	1:D:249:ILE:HG13	2.15	0.46
1:F:145:PRO:HB2	1:F:189:PHE:CD2	2.51	0.46
1:G:88:SER:OG	1:G:90:LYS:HG3	2.15	0.46
1:H:278:ARG:NH2	2:H:623:HOH:O	2.42	0.46
1:F:371:TYR:CZ	1:F:372:LYS:HE3	2.51	0.46
1:G:335:VAL:HA	1:G:350:LEU:HD13	1.98	0.46
1:D:133:TYR:CZ	1:D:137:LYS:HD3	2.50	0.46
1:D:220:THR:HG22	1:D:463:ALA:O	2.16	0.46
1:A:224:SER:H	1:A:248:SER:CB	2.28	0.46
1:G:151:MET:HE1	1:G:243:SER:HB3	1.98	0.46
1:H:182:ILE:HA	1:H:186:ALA:O	2.16	0.46
1:A:150:GLU:HB3	1:A:194:ASN:ND2	2.30	0.46
1:A:155:VAL:HG23	2:A:648:HOH:O	2.16	0.46
1:C:386:GLU:HG2	2:C:641:HOH:O	2.16	0.45
1:G:203:CYS:HB3	1:G:211:THR:HB	1.98	0.45
1:D:244:ASP:HB2	1:D:249:ILE:HB	1.98	0.45
1:A:277:ASN:HA	1:B:422:TRP:CZ3	2.51	0.45
1:F:224:SER:H	1:F:248:SER:CB	2.29	0.45
1:H:501:LYS:HA	1:H:501:LYS:HD2	1.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:SER:HB2	1:A:524:LYS:HE3	1.97	0.45
1:C:42:PRO:HG2	1:C:56:ALA:HB2	1.97	0.45
1:C:89:ARG:CG	2:C:695:HOH:O	2.65	0.45
1:D:131:ALA:C	1:D:133:TYR:H	2.17	0.45
1:D:219:LEU:HD13	1:D:462:ALA:O	2.17	0.45
1:E:319:LEU:O	2:E:602:HOH:O	2.21	0.45
1:F:262:ARG:NH2	1:F:481:ASP:HA	2.32	0.45
1:G:155:VAL:HG23	1:G:174:ASN:OD1	2.17	0.45
1:C:8:GLU:OE1	2:C:601:HOH:O	2.21	0.45
1:C:195:PRO:HB3	1:C:203:CYS:HA	1.97	0.45
1:C:330:ARG:NH2	1:C:444:ASP:HB3	2.32	0.45
1:H:92:THR:HG22	1:H:95:GLU:HB2	1.99	0.45
1:A:157:ARG:HA	1:A:157:ARG:HD3	1.65	0.45
1:C:224:SER:H	1:C:248:SER:CB	2.29	0.45
1:H:457:VAL:O	1:H:474:THR:HG23	2.16	0.45
1:B:195:PRO:HB3	1:B:203:CYS:HA	1.99	0.45
1:B:42:PRO:HG2	1:B:56:ALA:HB2	1.98	0.45
1:C:250:ARG:HD3	1:C:484:VAL:HB	1.98	0.45
1:F:228:GLU:OE1	1:F:240:GLY:HA3	2.17	0.45
1:G:204:VAL:HG22	1:G:210:ARG:HG3	1.98	0.45
1:G:280:SER:O	2:G:601:HOH:O	2.20	0.45
1:A:348:ARG:NH1	1:A:497:ASP:OD1	2.50	0.44
1:D:147:SER:O	1:D:240:GLY:HA2	2.17	0.44
1:D:213:ASN:ND2	1:D:214:ALA:H	2.15	0.44
1:C:393:ARG:NH1	2:C:623:HOH:O	2.49	0.44
1:G:457:VAL:O	1:G:474:THR:HG23	2.17	0.44
1:H:392:PHE:CZ	1:H:406:THR:HG21	2.52	0.44
1:H:394:ALA:O	1:H:397:GLN:HG3	2.17	0.44
1:B:492:VAL:HG23	1:B:526:ALA:O	2.17	0.44
1:E:353:VAL:HG11	1:E:454:ILE:HD13	1.99	0.44
1:E:457:VAL:O	1:E:474:THR:HG23	2.18	0.44
1:G:262:ARG:NH1	1:G:481:ASP:HA	2.32	0.44
1:G:289:PRO:HG2	1:G:297:VAL:HG13	2.00	0.44
1:H:94:VAL:HG12	1:H:98:LYS:HE2	1.99	0.44
1:H:224:SER:H	1:H:248:SER:CB	2.31	0.44
1:D:86:LEU:HD21	1:D:96:VAL:CG2	2.44	0.44
1:A:182:ILE:HA	1:A:186:ALA:O	2.18	0.44
1:B:210:ARG:NH2	2:B:628:HOH:O	2.51	0.44
1:B:224:SER:H	1:B:248:SER:CB	2.31	0.44
1:D:92:THR:CG2	1:D:130:ASP:CG	2.84	0.44
1:A:382:LYS:HG3	1:A:424:TRP:CE2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:511:LYS:HE2	1:G:515:GLU:OE2	2.17	0.44
1:D:481:ASP:OD2	1:D:536:ARG:NH1	2.50	0.44
1:A:155:VAL:HG13	1:A:174:ASN:OD1	2.17	0.44
1:A:243:SER:O	1:A:248:SER:HB2	2.17	0.44
1:A:250:ARG:NH1	2:A:631:HOH:O	2.50	0.44
1:E:92:THR:HG22	1:E:94:VAL:N	2.31	0.44
1:G:476:GLN:O	1:G:476:GLN:HG3	2.17	0.44
1:H:265:LEU:O	2:H:602:HOH:O	2.20	0.44
1:D:194:ASN:OD1	1:D:194:ASN:N	2.51	0.44
1:A:116:GLU:OE2	1:A:157:ARG:NH2	2.45	0.43
1:H:382:LYS:HA	2:H:631:HOH:O	2.18	0.43
1:G:356:LYS:O	1:G:360:LEU:HD22	2.17	0.43
1:A:481:ASP:OD2	1:A:536:ARG:NH1	2.51	0.43
1:C:394:ALA:O	1:C:397:GLN:HG3	2.17	0.43
1:F:155:VAL:HG23	1:F:174:ASN:OD1	2.18	0.43
1:G:481:ASP:OD2	1:G:536:ARG:NH1	2.51	0.43
1:H:92:THR:HG22	1:H:95:GLU:H	1.83	0.43
1:D:182:ILE:HA	1:D:186:ALA:O	2.18	0.43
1:F:392:PHE:CZ	1:F:406:THR:HG21	2.53	0.43
1:D:213:ASN:ND2	1:D:214:ALA:N	2.66	0.43
1:F:192:THR:HB	1:F:227:GLY:HA3	2.01	0.43
1:F:243:SER:O	1:F:249:ILE:HG13	2.19	0.43
1:E:158:SER:HA	1:E:169:ARG:O	2.18	0.43
1:H:76:GLU:OE1	1:H:235:ARG:HD2	2.19	0.43
1:D:521:SER:HB2	1:D:524:LYS:HE3	2.01	0.43
1:B:348:ARG:NH2	1:B:491:LYS:O	2.40	0.43
1:C:133:TYR:O	1:C:137:LYS:HB3	2.18	0.43
1:D:92:THR:HG22	1:D:94:VAL:H	1.83	0.43
1:B:330:ARG:NH1	1:B:444:ASP:HB3	2.33	0.43
1:E:224:SER:OG	1:E:248:SER:HB3	2.19	0.43
1:F:481:ASP:OD2	1:F:536:ARG:NH1	2.52	0.43
1:D:389:ALA:O	1:D:393:ARG:HG3	2.19	0.43
1:C:275:SER:HB2	2:C:627:HOH:O	2.19	0.42
1:F:220:THR:HG23	1:F:222:GLY:N	2.33	0.42
1:F:450:LEU:HD23	1:F:450:LEU:HA	1.94	0.42
1:H:100:TYR:OH	1:H:292:ARG:NH2	2.47	0.42
1:D:200:MET:HE2	1:D:202:GLU:HB2	2.00	0.42
1:C:240:GLY:O	1:C:289:PRO:HA	2.19	0.42
1:H:146:PHE:O	1:H:188:PRO:HA	2.19	0.42
1:H:519:SER:HG	1:H:521:SER:HG	1.65	0.42
1:D:43:SER:HB2	1:D:109:LYS:HE3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:VAL:HG11	1:D:117:PHE:CZ	2.55	0.42
1:C:182:ILE:HA	1:C:186:ALA:O	2.20	0.42
1:C:308:LYS:NZ	2:C:618:HOH:O	2.42	0.42
1:C:348:ARG:NH1	1:C:497:ASP:OD1	2.52	0.42
1:D:550:ARG:O	1:D:551:ASN:HB2	2.19	0.42
1:C:243:SER:O	1:C:248:SER:HB2	2.20	0.42
1:G:115:GLN:OE1	1:G:208:HIS:HB3	2.20	0.42
1:G:392:PHE:CZ	1:G:406:THR:HG21	2.53	0.42
1:D:392:PHE:CZ	1:D:406:THR:HG21	2.55	0.42
1:F:220:THR:HG23	1:F:222:GLY:H	1.85	0.42
1:D:132:ASP:HB3	1:D:135:THR:HG1	1.83	0.42
1:D:464:LYS:O	1:D:467:THR:HG22	2.20	0.42
1:B:536:ARG:NH2	2:B:630:HOH:O	2.52	0.42
1:E:392:PHE:CZ	1:E:406:THR:HG21	2.55	0.42
1:G:85:ASP:HB3	1:G:91:VAL:HG23	2.00	0.42
1:D:34:LEU:HD22	1:D:68:PRO:CD	2.50	0.42
1:B:23:GLN:O	2:B:602:HOH:O	2.21	0.42
1:C:422:TRP:CZ3	1:D:277:ASN:HA	2.55	0.42
1:F:390:ARG:HD2	2:F:610:HOH:O	2.20	0.42
1:F:457:VAL:O	1:F:474:THR:HG23	2.20	0.42
1:G:243:SER:O	1:G:249:ILE:HG13	2.19	0.42
1:B:199:MET:O	2:B:604:HOH:O	2.22	0.42
1:C:129:LEU:HG	2:C:636:HOH:O	2.18	0.42
1:E:289:PRO:HG2	1:E:297:VAL:HG13	2.01	0.42
1:F:114:VAL:HG11	1:F:117:PHE:CZ	2.55	0.42
1:A:146:PHE:O	1:A:188:PRO:HA	2.19	0.41
1:A:272:ASP:OD1	2:A:601:HOH:O	2.21	0.41
1:E:491:LYS:NZ	1:E:525:ASN:HD22	2.18	0.41
1:G:213:ASN:HB3	1:G:216:ASN:O	2.19	0.41
1:C:228:GLU:OE1	1:C:240:GLY:HA3	2.19	0.41
1:C:357:LEU:HD22	1:C:363:PHE:CD1	2.56	0.41
1:E:22:LYS:O	1:E:26:GLU:HG3	2.20	0.41
1:F:262:ARG:HE	1:F:478:ASN:ND2	2.19	0.41
1:D:200:MET:CE	1:D:202:GLU:HB2	2.51	0.41
1:A:115:GLN:HG2	1:A:150:GLU:HG3	2.03	0.41
1:E:464:LYS:HE3	1:E:523:PHE:CZ	2.55	0.41
1:D:394:ALA:O	1:D:397:GLN:HG3	2.20	0.41
1:A:5:ASN:ND2	1:A:7:LYS:HB2	2.36	0.41
1:G:46:GLY:HA3	1:G:51:ASP:O	2.20	0.41
1:D:352:MET:HE2	1:D:489:VAL:HG12	2.02	0.41
1:H:495:SER:O	2:H:603:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:292:ARG:H	1:D:292:ARG:HG3	1.45	0.41
1:D:547:ASN:O	1:D:550:ARG:HG3	2.21	0.41
1:C:427:GLN:HA	2:C:602:HOH:O	2.20	0.41
1:C:224:SER:H	1:C:248:SER:HB3	1.84	0.41
1:E:382:LYS:HG3	1:E:424:TRP:CE2	2.55	0.41
1:F:339:ILE:HD11	1:F:373:PRO:HG2	2.03	0.41
1:F:501:LYS:HA	1:F:501:LYS:HD2	1.90	0.41
1:D:213:ASN:HD22	1:D:214:ALA:N	2.18	0.41
1:A:168:ASP:C	1:A:170:ILE:HD12	2.41	0.41
1:A:361:GLY:O	2:A:602:HOH:O	2.22	0.41
1:A:501:LYS:HA	1:A:501:LYS:HD2	1.82	0.41
1:B:289:PRO:HG2	1:B:297:VAL:HG13	2.02	0.41
1:E:421:GLN:HB2	1:F:418:MET:HE3	2.02	0.41
1:F:88:SER:OG	1:F:90:LYS:HG3	2.20	0.41
1:G:224:SER:H	1:G:248:SER:CB	2.33	0.41
1:H:177:ILE:HD12	1:H:268:ILE:CG2	2.51	0.41
1:H:262:ARG:NH2	1:H:481:ASP:HA	2.36	0.41
1:E:281:GLU:OE1	2:E:604:HOH:O	2.22	0.41
1:F:213:ASN:HB3	1:F:216:ASN:O	2.21	0.40
1:H:220:THR:HG23	1:H:222:GLY:H	1.86	0.40
1:D:137:LYS:HB3	1:D:137:LYS:HE2	1.39	0.40
1:A:46:GLY:HA3	1:A:51:ASP:O	2.21	0.40
1:A:121:GLU:OE2	1:A:157:ARG:NH1	2.54	0.40
1:B:262:ARG:NH1	1:B:481:ASP:HA	2.36	0.40
1:E:204:VAL:HG22	1:E:210:ARG:HG3	2.03	0.40
1:H:266:GLY:N	2:H:611:HOH:O	2.28	0.40
1:D:100:TYR:CE2	1:D:238:PRO:HG3	2.57	0.40
1:F:146:PHE:O	1:F:188:PRO:HA	2.21	0.40
1:C:344:PRO:HG3	1:C:500:TYR:CE1	2.56	0.40
1:F:133:TYR:O	1:F:137:LYS:HB2	2.21	0.40
1:B:386:GLU:HG2	2:B:620:HOH:O	2.22	0.40
1:B:544:GLU:O	1:B:547:ASN:HB3	2.21	0.40
1:C:76:GLU:OE2	1:C:235:ARG:NH1	2.52	0.40
1:F:476:GLN:NE2	2:F:625:HOH:O	2.49	0.40
1:D:27:LEU:CD1	1:D:123:LEU:HB2	2.51	0.40
1:D:157:ARG:HB3	1:D:191:ARG:NH1	2.36	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	549/551 (100%)	535 (97%)	14 (3%)	0	100	100
1	B	549/551 (100%)	533 (97%)	16 (3%)	0	100	100
1	C	549/551 (100%)	532 (97%)	17 (3%)	0	100	100
1	D	549/551 (100%)	527 (96%)	21 (4%)	1 (0%)	44	61
1	E	549/551 (100%)	532 (97%)	17 (3%)	0	100	100
1	F	549/551 (100%)	532 (97%)	17 (3%)	0	100	100
1	G	549/551 (100%)	533 (97%)	16 (3%)	0	100	100
1	H	549/551 (100%)	533 (97%)	16 (3%)	0	100	100
All	All	4392/4408 (100%)	4257 (97%)	134 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	502	ASN

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	476/476 (100%)	474 (100%)	2 (0%)	89	95
1	B	476/476 (100%)	471 (99%)	5 (1%)	70	84
1	C	476/476 (100%)	473 (99%)	3 (1%)	84	92
1	D	476/476 (100%)	468 (98%)	8 (2%)	56	75

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	E	476/476 (100%)	472 (99%)	4 (1%)	79 89
1	F	476/476 (100%)	472 (99%)	4 (1%)	79 89
1	G	476/476 (100%)	473 (99%)	3 (1%)	84 92
1	H	476/476 (100%)	470 (99%)	6 (1%)	65 80
All	All	3808/3808 (100%)	3773 (99%)	35 (1%)	75 87

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	384	TYR
1	A	539	ASP
1	B	292	ARG
1	B	384	TYR
1	B	446	ASP
1	B	539	ASP
1	B	550	ARG
1	C	384	TYR
1	C	407	ARG
1	C	539	ASP
1	E	91	VAL
1	E	194	ASN
1	E	384	TYR
1	E	539	ASP
1	F	384	TYR
1	F	476	GLN
1	F	539	ASP
1	F	550	ARG
1	G	292	ARG
1	G	384	TYR
1	G	539	ASP
1	H	92	THR
1	H	224	SER
1	H	292	ARG
1	H	375	LYS
1	H	384	TYR
1	H	539	ASP
1	D	15	SER
1	D	54	SER
1	D	111	THR
1	D	292	ARG
1	D	384	TYR

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Mol	Chain	Res	Type
1	D	467	THR
1	D	539	ASP
1	D	550	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	194	ASN
1	B	194	ASN
1	C	57	GLN
1	C	476	GLN
1	C	516	GLN
1	E	525	ASN
1	G	525	ASN
1	H	476	GLN
1	D	213	ASN

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)

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, 95^{th} 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	551/551 (100%)	-0.29	8 (1%) 71 69	10, 26, 50, 109	0
1	B	551/551 (100%)	-0.13	7 (1%) 74 72	10, 30, 57, 110	0
1	C	551/551 (100%)	0.20	12 (2%) 62 60	15, 37, 71, 128	0
1	D	551/551 (100%)	1.43	165 (29%) 1 1	22, 56, 93, 119	0
1	E	551/551 (100%)	-0.14	7 (1%) 74 72	13, 30, 54, 101	0
1	F	551/551 (100%)	-0.19	6 (1%) 77 75	10, 30, 54, 104	0
1	G	551/551 (100%)	0.12	11 (1%) 64 62	17, 35, 64, 106	0
1	H	551/551 (100%)	0.34	14 (2%) 58 56	18, 40, 68, 107	0
All	All	4408/4408 (100%)	0.17	230 (5%) 34 32	10, 34, 72, 128	0

All (230) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	71	VAL	5.5
1	D	236	ALA	5.0
1	D	39	GLY	4.8
1	D	80	VAL	4.6
1	D	113	CYS	4.4
1	D	234	MET	4.4
1	D	25	ILE	4.1
1	E	2	LEU	4.0
1	D	83	LEU	4.0
1	D	292	ARG	4.0
1	D	257	GLY	3.8
1	D	38	TYR	3.8
1	G	551	ASN	3.8
1	D	67	LEU	3.8
1	D	96	VAL	3.8
1	D	69	LYS	3.7

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Mol	Chain	Res	Type	RSRZ
1	F	3	THR	3.7
1	D	42	PRO	3.7
1	C	326	ASN	3.7
1	B	3	THR	3.7
1	D	27	LEU	3.7
1	D	119	PRO	3.6
1	D	348	ARG	3.6
1	D	126	ALA	3.6
1	D	525	ASN	3.5
1	D	125	TYR	3.5
1	D	64	LEU	3.5
1	D	86	LEU	3.5
1	D	326	ASN	3.5
1	D	123	LEU	3.4
1	D	56	ALA	3.4
1	D	237	SER	3.4
1	D	213	ASN	3.4
1	D	47	ALA	3.3
1	D	114	VAL	3.3
1	D	355	GLU	3.3
1	D	130	ASP	3.3
1	D	60	LEU	3.3
1	D	28	ALA	3.3
1	D	129	LEU	3.3
1	D	189	PHE	3.3
1	D	102	ALA	3.3
1	D	143	GLY	3.3
1	D	186	ALA	3.2
1	D	154	PHE	3.2
1	E	58	PHE	3.2
1	D	111	THR	3.2
1	D	145	PRO	3.2
1	D	120	ILE	3.2
1	D	124	GLN	3.2
1	F	502	ASN	3.2
1	D	92	THR	3.2
1	D	30	PHE	3.1
1	D	75	THR	3.1
1	D	142	TYR	3.1
1	D	49	VAL	3.1
1	D	156	GLY	3.1
1	D	94	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	63	TYR	3.1
1	D	215	TYR	3.1
1	F	4	ASP	3.0
1	D	91	VAL	3.0
1	D	256	ASN	3.0
1	G	69	LYS	3.0
1	D	131	ALA	3.0
1	D	155	VAL	3.0
1	D	190	VAL	2.9
1	D	327	LYS	2.9
1	D	184	ASN	2.9
1	D	240	GLY	2.9
1	D	187	TYR	2.9
1	D	58	PHE	2.9
1	D	188	PRO	2.9
1	D	72	LEU	2.9
1	D	290	LEU	2.9
1	D	66	ALA	2.9
1	D	299	LEU	2.9
1	H	327	LYS	2.9
1	D	44	ALA	2.9
1	D	105	ILE	2.8
1	D	233	GLY	2.8
1	C	502	ASN	2.8
1	D	547	ASN	2.8
1	D	95	GLU	2.8
1	D	65	LYS	2.8
1	D	85	ASP	2.8
1	D	323	ARG	2.8
1	H	324	PRO	2.7
1	D	153	PRO	2.7
1	D	212	VAL	2.7
1	D	498	GLN	2.7
1	C	327	LYS	2.7
1	D	238	PRO	2.7
1	D	521	SER	2.7
1	D	99	ALA	2.7
1	D	214	ALA	2.7
1	A	2	LEU	2.7
1	D	110	LEU	2.7
1	G	500	TYR	2.7
1	D	127	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	82	LEU	2.7
1	D	93	CYS	2.7
1	D	43	SER	2.7
1	E	4	ASP	2.6
1	G	495	SER	2.6
1	D	11	GLY	2.6
1	G	327	LYS	2.6
1	B	32	GLU	2.6
1	D	1	MET	2.6
1	D	48	ASN	2.6
1	D	496	LEU	2.6
1	D	144	LEU	2.6
1	D	231	LEU	2.6
1	D	220	THR	2.6
1	D	226	GLY	2.6
1	D	35	ALA	2.5
1	D	546	ALA	2.5
1	E	417	ASP	2.5
1	D	489	VAL	2.5
1	F	501	LYS	2.5
1	H	504	LYS	2.5
1	D	41	LEU	2.5
1	D	174	ASN	2.5
1	A	3	THR	2.5
1	D	229	GLY	2.5
1	D	122	ALA	2.5
1	C	411	GLU	2.4
1	D	140	PRO	2.4
1	D	117	PHE	2.4
1	D	107	ALA	2.4
1	D	361	GLY	2.4
1	D	522	SER	2.4
1	G	449	VAL	2.4
1	D	70	LYS	2.4
1	D	230	ALA	2.4
1	H	551	ASN	2.4
1	D	217	GLY	2.4
1	D	34	LEU	2.4
1	D	118	LEU	2.4
1	D	235	ARG	2.4
1	A	1	MET	2.4
1	D	55	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	138	HIS	2.4
1	D	88	SER	2.4
1	C	503	TYR	2.4
1	D	414	GLU	2.3
1	C	499	PRO	2.3
1	C	58	PHE	2.3
1	D	224	SER	2.3
1	D	97	LEU	2.3
1	H	524	LYS	2.3
1	D	191	ARG	2.3
1	D	335	VAL	2.3
1	D	258	ILE	2.3
1	H	361	GLY	2.3
1	D	139	LEU	2.3
1	D	259	TYR	2.3
1	C	501	LYS	2.3
1	H	328	LYS	2.3
1	D	26	GLU	2.3
1	D	32	GLU	2.3
1	D	500	TYR	2.3
1	A	324	PRO	2.3
1	D	527	PRO	2.3
1	E	397	GLN	2.3
1	D	225	SER	2.3
1	H	323	ARG	2.2
1	G	362	ASN	2.2
1	D	268	ILE	2.2
1	C	324	PRO	2.2
1	E	3	THR	2.2
1	G	326	ASN	2.2
1	A	327	LYS	2.2
1	B	412	GLY	2.2
1	H	110	LEU	2.2
1	D	162	GLY	2.2
1	D	104	SER	2.2
1	A	325	GLU	2.2
1	H	325	GLU	2.2
1	D	62	GLU	2.2
1	B	326	ASN	2.2
1	H	525	ASN	2.2
1	D	31	ASP	2.2
1	D	19	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	291	SER	2.2
1	A	502	ASN	2.2
1	D	100	TYR	2.2
1	D	57	GLN	2.2
1	D	74	ILE	2.2
1	D	255	PHE	2.2
1	B	47	ALA	2.2
1	D	103	ALA	2.2
1	D	499	PRO	2.1
1	D	128	LYS	2.1
1	D	249	ILE	2.1
1	D	524	LYS	2.1
1	F	32	GLU	2.1
1	A	326	ASN	2.1
1	D	173	TYR	2.1
1	C	509	LEU	2.1
1	C	325	GLU	2.1
1	D	502	ASN	2.1
1	E	32	GLU	2.1
1	D	59	PRO	2.1
1	D	61	GLU	2.1
1	D	321	TRP	2.1
1	G	501	LYS	2.1
1	D	360	LEU	2.1
1	F	500	TYR	2.1
1	D	40	ALA	2.1
1	B	325	GLU	2.0
1	H	355	GLU	2.0
1	G	36	LYS	2.0
1	H	92	THR	2.0
1	D	182	ILE	2.0
1	D	548	ILE	2.0
1	D	492	VAL	2.0
1	C	292	ARG	2.0
1	D	5	ASN	2.0
1	H	85	ASP	2.0
1	D	501	LYS	2.0
1	D	504	LYS	2.0
1	G	489	VAL	2.0
1	D	20	SER	2.0
1	D	359	SER	2.0
1	D	216	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	4	ASP	2.0
1	D	349	ALA	2.0
1	D	503	TYR	2.0
1	D	528	ALA	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.