



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:59 AM UTC

PDB ID : 1UWI / pdb_00001uwi
Title : CRYSTAL STRUCTURE OF MUTATED BETA-GLYCOSIDASE FROM
SULFOLOBUS SOLFATARICUS, WORKING AT MODERATE TEMPER-
ATURE
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Deposited on : 2004-02-05
Resolution : 2.55 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

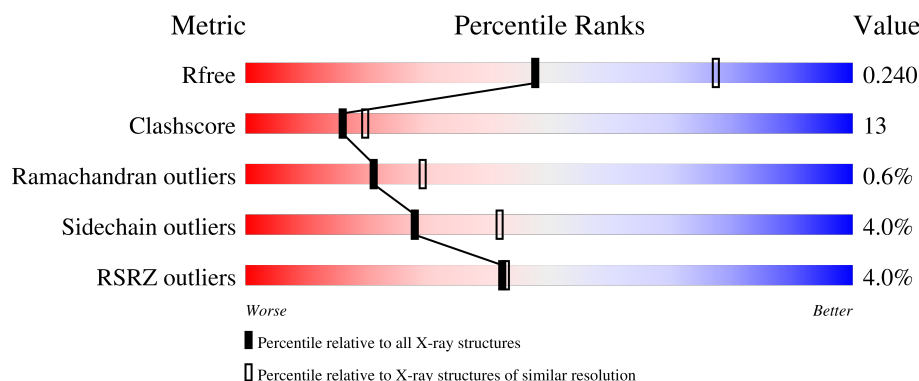
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	489	3% 75% 22% .
1	B	489	4% 73% 25% .
1	C	489	4% 76% 21% .
1	D	489	4% 76% 20% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16494 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 BETA-GALACTOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	0	0
			4017	2573	693	739	12			
1	B	489	Total	C	N	O	S	0	0	0
			4017	2573	693	739	12			
1	C	489	Total	C	N	O	S	0	0	0
			4017	2573	693	739	12			
1	D	489	Total	C	N	O	S	0	0	0
			4017	2573	693	739	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	LEU	PRO	engineered mutation	UNP P22498
A	68	ASN	ASP	engineered mutation	UNP P22498
A	331	GLY	GLU	engineered mutation	UNP P22498
A	459	SER	ALA	engineered mutation	UNP P22498
B	29	LEU	PRO	engineered mutation	UNP P22498
B	68	ASN	ASP	engineered mutation	UNP P22498
B	331	GLY	GLU	engineered mutation	UNP P22498
B	459	SER	ALA	engineered mutation	UNP P22498
C	29	LEU	PRO	engineered mutation	UNP P22498
C	68	ASN	ASP	engineered mutation	UNP P22498
C	331	GLY	GLU	engineered mutation	UNP P22498
C	459	SER	ALA	engineered mutation	UNP P22498
D	29	LEU	PRO	engineered mutation	UNP P22498
D	68	ASN	ASP	engineered mutation	UNP P22498
D	331	GLY	GLU	engineered mutation	UNP P22498
D	459	SER	ALA	engineered mutation	UNP P22498

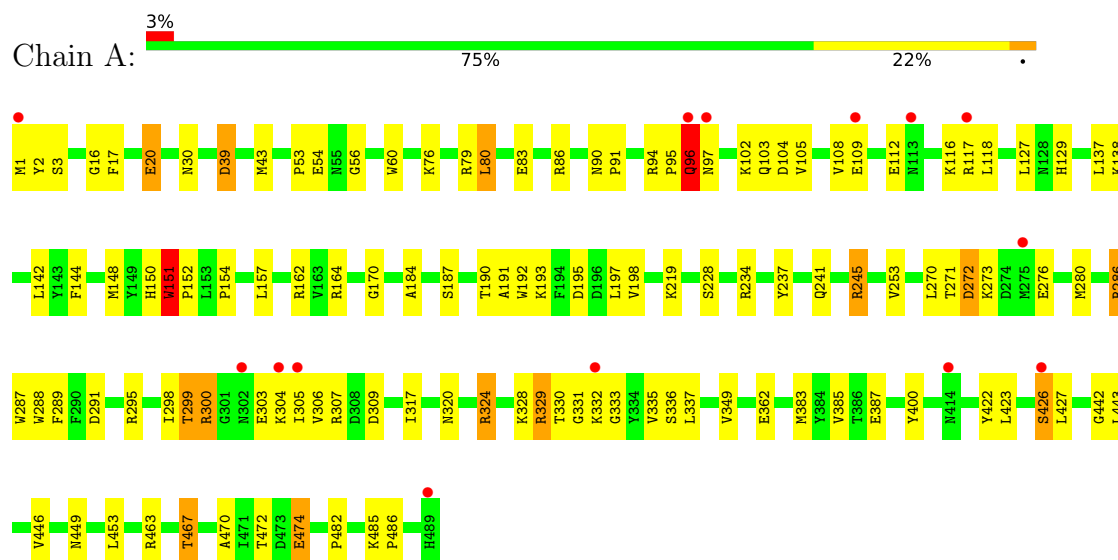
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	128	Total 128	O 128	0	0
2	B	102	Total 102	O 102	0	0
2	C	81	Total 81	O 81	0	0
2	D	115	Total 115	O 115	0	0

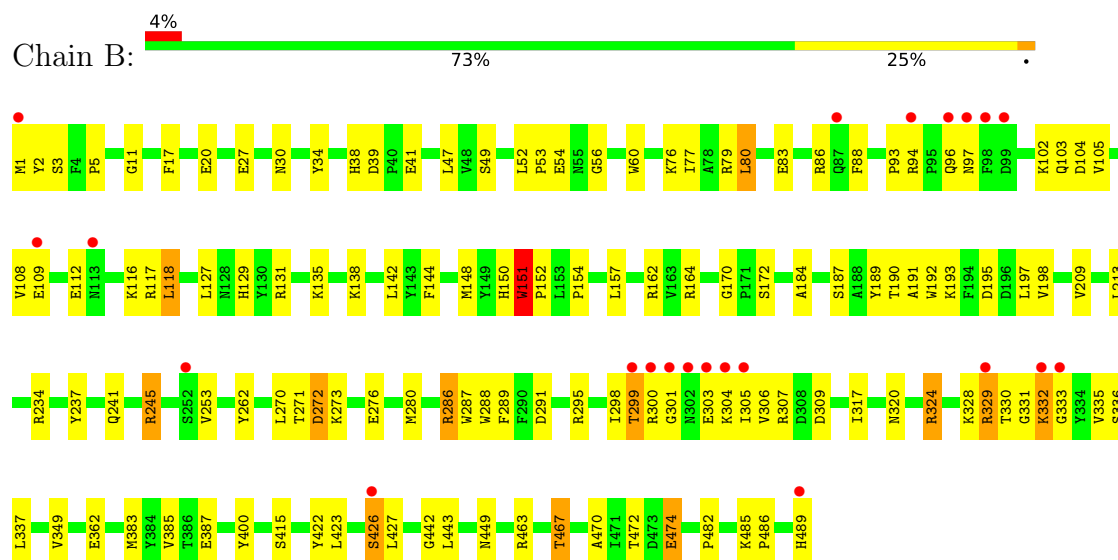
3 Residue-property plots

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.

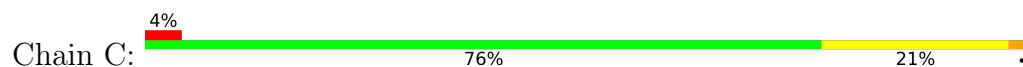
• Molecule 1: BETA-GALACTOSIDASE

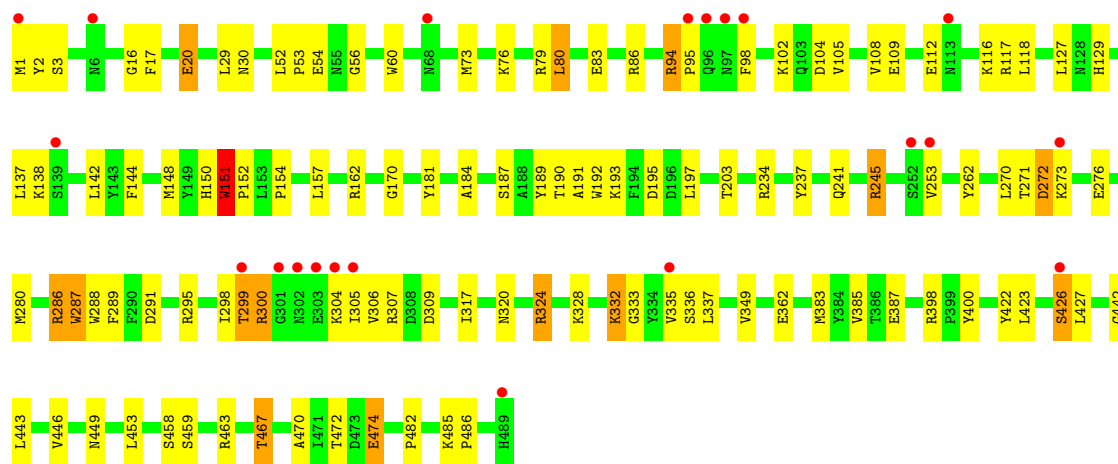


• Molecule 1: BETA-GALACTOSIDASE

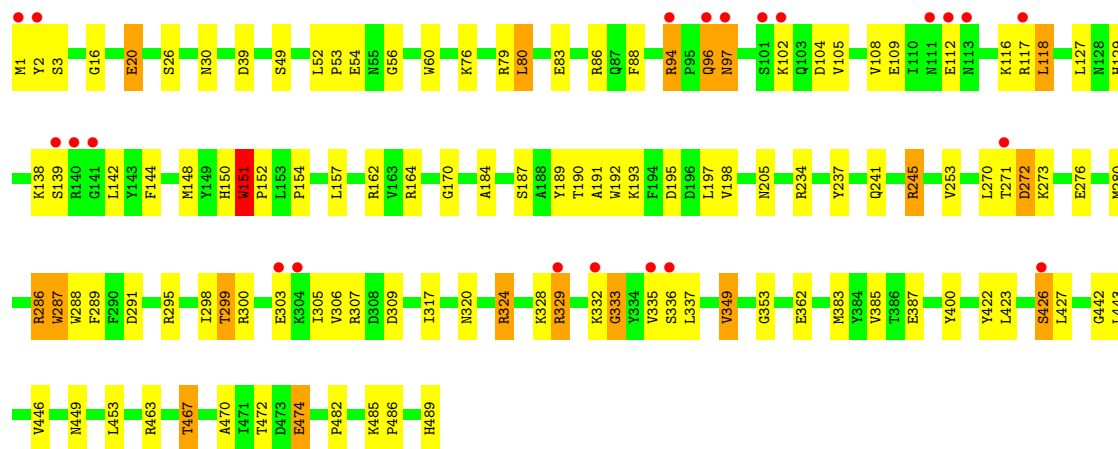
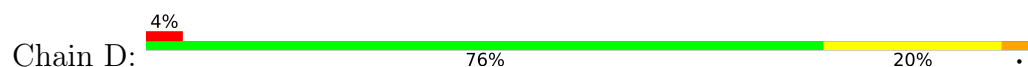


• Molecule 1: BETA-GALACTOSIDASE





• Molecule 1: BETA-GALACTOSIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.08Å 132.43Å 219.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.44 – 2.55 28.44 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.2 (28.44-2.55) 99.1 (28.44-2.55)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.241 0.209 , 0.240	Depositor DCC
R_{free} test set	5638 reflections (7.07%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.213	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16494	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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	0.40	0/4145	0.93	14/5629 (0.2%)
1	B	0.41	0/4145	0.92	15/5629 (0.3%)
1	C	0.40	0/4145	0.92	11/5629 (0.2%)
1	D	0.40	0/4145	0.92	13/5629 (0.2%)
All	All	0.40	0/16580	0.92	53/22516 (0.2%)

There are no bond length outliers.

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	443	LEU	N-CA-C	-8.83	101.10	112.23
1	A	443	LEU	N-CA-C	-8.70	101.27	112.23
1	B	443	LEU	N-CA-C	-8.69	101.28	112.23
1	D	443	LEU	N-CA-C	-8.48	101.54	112.23
1	A	442	GLY	N-CA-C	7.90	122.33	112.14
1	D	442	GLY	N-CA-C	7.67	122.04	112.14
1	B	442	GLY	N-CA-C	7.67	122.03	112.14
1	C	442	GLY	N-CA-C	7.55	121.88	112.14
1	C	426	SER	N-CA-C	6.93	125.56	110.80
1	A	426	SER	N-CA-C	6.83	125.36	110.80
1	B	426	SER	N-CA-C	6.76	125.21	110.80
1	D	426	SER	N-CA-C	6.71	125.10	110.80
1	B	301	GLY	N-CA-C	-6.56	105.64	111.67
1	C	20	GLU	N-CA-C	6.53	118.08	110.97
1	C	56	GLY	CA-C-N	6.52	126.45	119.28
1	C	56	GLY	C-N-CA	6.52	126.45	119.28
1	A	20	GLU	N-CA-C	6.51	118.07	110.97
1	D	56	GLY	CA-C-N	6.46	126.39	119.28
1	D	56	GLY	C-N-CA	6.46	126.39	119.28
1	B	56	GLY	CA-C-N	6.46	126.39	119.28
1	B	56	GLY	C-N-CA	6.46	126.39	119.28
1	A	96	GLN	N-CA-C	-6.38	102.29	110.53
1	B	20	GLU	N-CA-C	6.34	117.88	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	20	GLU	N-CA-C	6.34	117.88	110.97
1	A	56	GLY	CA-C-N	6.24	126.14	119.28
1	A	56	GLY	C-N-CA	6.24	126.14	119.28
1	D	144	PHE	N-CA-C	5.98	118.16	108.41
1	B	30	ASN	N-CA-C	5.95	120.81	112.72
1	B	144	PHE	N-CA-C	5.90	118.01	108.34
1	B	385	VAL	N-CA-C	-5.87	98.03	106.55
1	A	385	VAL	N-CA-C	-5.85	98.07	106.55
1	C	30	ASN	N-CA-C	5.79	120.60	112.72
1	D	30	ASN	N-CA-C	5.77	120.57	112.72
1	A	30	ASN	N-CA-C	5.71	120.49	112.72
1	D	385	VAL	N-CA-C	-5.64	98.37	106.55
1	B	49	SER	N-CA-C	5.61	117.39	111.28
1	A	144	PHE	N-CA-C	5.53	117.41	108.34
1	C	144	PHE	N-CA-C	5.35	117.12	108.34
1	D	49	SER	N-CA-C	5.30	117.80	111.71
1	C	385	VAL	N-CA-C	-5.29	98.89	106.55
1	A	198	VAL	N-CA-C	5.23	116.32	109.21
1	B	198	VAL	N-CA-C	5.23	116.32	109.21
1	C	151	TRP	N-CA-C	5.20	121.31	109.81
1	B	151	TRP	N-CA-C	5.19	121.28	109.81
1	D	151	TRP	N-CA-C	5.17	121.23	109.81
1	D	287	TRP	N-CA-C	5.15	117.56	111.33
1	D	198	VAL	N-CA-C	5.13	116.19	109.21
1	C	287	TRP	N-CA-C	5.12	117.26	111.11
1	B	172	SER	N-CA-C	5.08	118.52	112.72
1	A	151	TRP	N-CA-C	5.04	120.95	109.81
1	B	415	SER	N-CA-C	-5.03	107.14	113.28
1	A	39	ASP	CA-C-N	5.02	124.62	119.05
1	A	39	ASP	C-N-CA	5.02	124.62	119.05

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	4017	0	3802	113	0
1	B	4017	0	3802	105	0
1	C	4017	0	3802	97	0
1	D	4017	0	3802	101	0
2	A	128	0	0	2	0
2	B	102	0	0	5	0
2	C	81	0	0	2	0
2	D	115	0	0	4	0
All	All	16494	0	15208	414	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 13.

All (414) 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:148:MET:HE2	1:A:187:SER:HA	1.34	1.09
1:C:148:MET:HE2	1:C:187:SER:HA	1.35	1.07
1:B:148:MET:HE2	1:B:187:SER:HA	1.34	1.05
1:D:148:MET:HE2	1:D:187:SER:HA	1.36	1.03
1:D:148:MET:HE1	1:D:190:THR:HB	1.54	0.88
1:B:148:MET:HE1	1:B:190:THR:HB	1.56	0.88
1:A:148:MET:HE1	1:A:190:THR:HB	1.57	0.85
1:B:305:ILE:HG22	1:B:306:VAL:H	1.43	0.83
1:C:148:MET:HE1	1:C:190:THR:HB	1.59	0.83
1:A:305:ILE:HG22	1:A:306:VAL:H	1.40	0.83
1:C:305:ILE:HG22	1:C:306:VAL:H	1.44	0.83
1:A:192:TRP:HA	1:A:253:VAL:HG11	1.62	0.82
1:D:305:ILE:HG22	1:D:306:VAL:H	1.43	0.82
1:C:192:TRP:HA	1:C:253:VAL:HG11	1.62	0.81
1:D:192:TRP:HA	1:D:253:VAL:HG11	1.61	0.81
1:C:104:ASP:HA	1:C:241:GLN:HE22	1.44	0.81
1:B:192:TRP:HA	1:B:253:VAL:HG11	1.64	0.79
1:B:104:ASP:HA	1:B:241:GLN:HE22	1.46	0.78
1:A:104:ASP:HA	1:A:241:GLN:HE22	1.47	0.78
1:D:104:ASP:HA	1:D:241:GLN:HE22	1.47	0.78
1:B:329:ARG:HD3	1:B:330:THR:N	1.99	0.78
1:C:463:ARG:HD2	2:C:2073:HOH:O	1.85	0.76
1:D:127:LEU:HD21	1:D:193:LYS:HD3	1.69	0.75
1:A:127:LEU:HD21	1:A:193:LYS:HD3	1.68	0.74
1:B:485:LYS:HG2	2:B:2099:HOH:O	1.88	0.74
1:B:127:LEU:HD21	1:B:193:LYS:HD3	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:LYS:HG3	1:A:333:GLY:H	1.53	0.73
1:B:241:GLN:HE21	1:B:307:ARG:HH22	1.37	0.72
1:D:241:GLN:HE21	1:D:307:ARG:HH22	1.34	0.72
1:B:76:LYS:HA	1:B:142:LEU:HD23	1.72	0.72
1:D:148:MET:HE1	1:D:190:THR:CB	2.20	0.72
1:A:241:GLN:HE21	1:A:307:ARG:HH22	1.38	0.71
1:C:127:LEU:HD21	1:C:193:LYS:HD3	1.69	0.71
1:C:485:LYS:HG2	2:C:2077:HOH:O	1.90	0.71
1:B:305:ILE:HG22	1:B:306:VAL:N	2.06	0.70
1:C:76:LYS:HA	1:C:142:LEU:HD23	1.72	0.70
1:A:463:ARG:O	1:A:467:THR:HB	1.92	0.70
1:B:148:MET:HE1	1:B:190:THR:CB	2.21	0.70
1:A:305:ILE:HG22	1:A:306:VAL:N	2.05	0.70
1:A:94:ARG:NH1	1:A:94:ARG:HB3	2.07	0.70
1:D:305:ILE:HG22	1:D:306:VAL:N	2.07	0.70
1:A:299:THR:HB	1:A:304:LYS:HE3	1.74	0.69
1:A:76:LYS:HA	1:A:142:LEU:HD23	1.74	0.69
1:A:148:MET:HE1	1:A:190:THR:CB	2.23	0.68
1:C:305:ILE:HG22	1:C:306:VAL:N	2.08	0.68
1:C:463:ARG:O	1:C:467:THR:HB	1.94	0.68
1:D:76:LYS:HA	1:D:142:LEU:HD23	1.75	0.68
1:D:329:ARG:HG3	1:D:329:ARG:HH11	1.57	0.68
1:B:463:ARG:O	1:B:467:THR:HB	1.94	0.67
1:C:241:GLN:HE21	1:C:307:ARG:HH22	1.40	0.67
1:A:328:LYS:HD2	1:A:337:LEU:HD21	1.77	0.67
1:B:96:GLN:HG3	1:B:97:ASN:H	1.60	0.67
1:C:148:MET:HE1	1:C:190:THR:CB	2.23	0.67
1:D:463:ARG:O	1:D:467:THR:HB	1.95	0.67
1:B:328:LYS:HD2	1:B:337:LEU:HD21	1.76	0.66
1:A:280:MET:HE1	1:A:333:GLY:HA2	1.78	0.65
1:B:105:VAL:H	1:B:241:GLN:NE2	1.95	0.64
1:A:324:ARG:HD3	1:A:362:GLU:OE1	1.97	0.64
1:C:105:VAL:H	1:C:241:GLN:NE2	1.96	0.64
1:D:324:ARG:HD3	1:D:362:GLU:OE1	1.97	0.64
1:A:105:VAL:H	1:A:241:GLN:NE2	1.95	0.64
1:D:241:GLN:NE2	1:D:307:ARG:HH12	1.95	0.63
1:D:192:TRP:CA	1:D:253:VAL:HG11	2.29	0.63
1:B:271:THR:HG22	1:B:272:ASP:N	2.14	0.63
1:A:271:THR:HG22	1:A:272:ASP:N	2.14	0.63
1:D:328:LYS:HD2	1:D:337:LEU:HD21	1.80	0.62
1:B:324:ARG:HD3	1:B:362:GLU:OE1	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271:THR:HG22	1:C:272:ASP:N	2.13	0.62
1:C:328:LYS:HD2	1:C:337:LEU:HD21	1.79	0.62
1:D:96:GLN:O	1:D:97:ASN:OD1	2.17	0.62
1:B:286:ARG:NH1	2:B:2057:HOH:O	2.32	0.62
1:D:105:VAL:H	1:D:241:GLN:NE2	1.97	0.62
1:C:324:ARG:HD3	1:C:362:GLU:OE1	1.99	0.62
1:B:96:GLN:CG	1:B:97:ASN:H	2.13	0.62
1:C:280:MET:HE1	1:C:333:GLY:HA2	1.82	0.62
1:A:241:GLN:NE2	1:A:307:ARG:HH12	1.96	0.61
1:B:241:GLN:NE2	1:B:307:ARG:HH12	1.98	0.61
1:C:151:TRP:HB2	1:C:152:PRO:HD3	1.83	0.61
1:D:271:THR:HG22	1:D:272:ASP:N	2.15	0.60
1:A:241:GLN:HE21	1:A:307:ARG:HH12	1.49	0.60
1:D:109:GLU:O	1:D:109:GLU:HG3	2.01	0.60
1:A:79:ARG:C	1:A:80:LEU:HD12	2.25	0.60
1:C:79:ARG:C	1:C:80:LEU:HD12	2.27	0.60
1:C:280:MET:HE1	1:C:332:LYS:O	2.02	0.60
1:B:79:ARG:C	1:B:80:LEU:HD12	2.27	0.60
1:B:94:ARG:HB3	1:B:94:ARG:NH1	2.17	0.60
1:C:241:GLN:NE2	1:C:307:ARG:HH12	1.99	0.60
1:A:109:GLU:O	1:A:109:GLU:HG3	2.00	0.59
1:B:109:GLU:HG3	1:B:109:GLU:O	2.01	0.59
1:C:109:GLU:O	1:C:109:GLU:HG3	2.02	0.59
1:A:151:TRP:HB2	1:A:152:PRO:HD3	1.84	0.59
1:A:286:ARG:HD2	1:A:287:TRP:CZ3	2.37	0.59
1:B:192:TRP:CA	1:B:253:VAL:HG11	2.31	0.59
1:C:192:TRP:CA	1:C:253:VAL:HG11	2.30	0.59
1:B:241:GLN:HE21	1:B:307:ARG:HH12	1.51	0.59
1:B:286:ARG:HD2	1:B:287:TRP:CZ3	2.38	0.58
1:D:79:ARG:C	1:D:80:LEU:HD12	2.28	0.58
1:C:271:THR:HG21	1:C:273:LYS:NZ	2.19	0.58
1:C:286:ARG:HD2	1:C:287:TRP:CZ3	2.38	0.58
1:D:329:ARG:HD3	1:D:333:GLY:O	2.04	0.58
1:D:117:ARG:NH1	1:D:117:ARG:HB2	2.19	0.58
1:B:271:THR:HG21	1:B:273:LYS:NZ	2.19	0.58
1:A:117:ARG:HB2	1:A:117:ARG:NH1	2.19	0.58
1:C:3:SER:HA	1:C:470:ALA:HB2	1.87	0.57
1:D:241:GLN:HE21	1:D:307:ARG:NH2	2.01	0.57
1:A:192:TRP:CA	1:A:253:VAL:HG11	2.30	0.57
1:A:237:TYR:CE1	1:A:300:ARG:HB3	2.38	0.57
1:D:151:TRP:HB2	1:D:152:PRO:HD3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:241:GLN:HE21	1:D:307:ARG:HH12	1.51	0.57
1:A:3:SER:HA	1:A:470:ALA:HB2	1.86	0.57
1:A:271:THR:HG21	1:A:273:LYS:NZ	2.20	0.57
1:C:117:ARG:HB2	1:C:117:ARG:NH1	2.20	0.57
1:C:94:ARG:HB3	1:C:94:ARG:NH1	2.19	0.57
1:C:241:GLN:HE21	1:C:307:ARG:HH12	1.51	0.57
1:C:485:LYS:HB2	1:C:486:PRO:HD3	1.86	0.57
1:D:3:SER:HA	1:D:470:ALA:HB2	1.85	0.57
1:B:151:TRP:HB2	1:B:152:PRO:HD3	1.85	0.57
1:B:241:GLN:HE21	1:B:307:ARG:NH2	2.03	0.57
1:A:485:LYS:HB2	1:A:486:PRO:HD3	1.87	0.56
1:D:271:THR:HG21	1:D:273:LYS:HZ3	1.71	0.56
1:B:3:SER:HA	1:B:470:ALA:HB2	1.87	0.56
1:B:271:THR:HG21	1:B:273:LYS:HZ3	1.69	0.56
1:B:485:LYS:HB2	1:B:486:PRO:HD3	1.87	0.56
1:B:280:MET:HE1	1:B:333:GLY:HA2	1.88	0.56
1:D:271:THR:HG21	1:D:273:LYS:NZ	2.20	0.56
1:D:286:ARG:HD2	1:D:287:TRP:CZ3	2.41	0.56
1:A:320:ASN:CG	1:A:387:GLU:HB2	2.31	0.56
1:B:270:LEU:HD22	1:B:337:LEU:HD11	1.86	0.56
1:D:150:HIS:O	1:D:151:TRP:HB2	2.05	0.55
1:A:303:GLU:HB3	1:A:305:ILE:HD11	1.87	0.55
1:C:320:ASN:CG	1:C:387:GLU:HB2	2.31	0.55
1:D:286:ARG:HG3	1:D:286:ARG:HH11	1.71	0.55
1:B:117:ARG:NH1	1:B:117:ARG:HB2	2.21	0.55
1:D:320:ASN:CG	1:D:387:GLU:HB2	2.31	0.54
1:A:330:THR:O	1:A:332:LYS:N	2.40	0.54
1:D:270:LEU:HD22	1:D:337:LEU:CD1	2.37	0.54
1:B:270:LEU:HD22	1:B:337:LEU:CD1	2.37	0.54
1:B:286:ARG:HG3	1:B:286:ARG:HH11	1.73	0.54
1:B:320:ASN:CG	1:B:387:GLU:HB2	2.33	0.54
1:A:270:LEU:HD22	1:A:337:LEU:CD1	2.38	0.54
1:B:329:ARG:HD3	1:B:330:THR:H	1.73	0.54
1:A:103:GLN:OE1	1:A:300:ARG:HD3	2.08	0.54
1:A:150:HIS:O	1:A:151:TRP:HB2	2.08	0.54
1:A:270:LEU:HD22	1:A:337:LEU:HD11	1.89	0.54
1:C:241:GLN:HE21	1:C:307:ARG:NH2	2.06	0.54
1:C:270:LEU:HD22	1:C:337:LEU:HD11	1.89	0.53
1:D:485:LYS:HB2	1:D:486:PRO:HD3	1.89	0.53
1:C:270:LEU:HD22	1:C:337:LEU:CD1	2.38	0.53
1:D:270:LEU:HD22	1:D:337:LEU:HD11	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:305:ILE:CG2	1:D:306:VAL:H	2.20	0.53
1:B:237:TYR:CE1	1:B:300:ARG:HB3	2.44	0.52
1:C:271:THR:HG21	1:C:273:LYS:HZ3	1.74	0.52
1:A:329:ARG:HG3	1:A:329:ARG:HH11	1.73	0.52
1:B:489:HIS:HB3	2:B:2101:HOH:O	2.08	0.52
1:A:300:ARG:HD2	1:A:300:ARG:O	2.09	0.52
1:A:286:ARG:HG3	1:A:286:ARG:HH11	1.74	0.52
1:D:1:MET:HB3	2:D:2002:HOH:O	2.09	0.52
1:B:317:ILE:O	1:B:383:MET:HB2	2.10	0.52
1:B:400:TYR:HB2	1:B:482:PRO:HG3	1.92	0.52
1:A:463:ARG:NH2	2:A:2119:HOH:O	2.43	0.52
1:A:241:GLN:HE21	1:A:307:ARG:NH2	2.04	0.51
1:A:400:TYR:HB2	1:A:482:PRO:HG3	1.92	0.51
1:B:150:HIS:O	1:B:151:TRP:HB2	2.10	0.51
1:A:80:LEU:HD12	1:A:80:LEU:N	2.25	0.51
1:D:192:TRP:HA	1:D:253:VAL:CG1	2.39	0.51
1:B:328:LYS:CD	1:B:337:LEU:HD21	2.41	0.51
1:C:150:HIS:O	1:C:151:TRP:HB2	2.10	0.51
1:A:184:ALA:HB2	1:A:245:ARG:HB3	1.93	0.51
1:D:328:LYS:CD	1:D:337:LEU:HD21	2.41	0.51
1:D:329:ARG:HG3	1:D:329:ARG:NH1	2.26	0.51
1:C:400:TYR:HB2	1:C:482:PRO:HG3	1.92	0.50
1:C:328:LYS:CD	1:C:337:LEU:HD21	2.41	0.50
1:D:303:GLU:N	1:D:303:GLU:CD	2.69	0.50
1:A:271:THR:HG21	1:A:273:LYS:HZ3	1.76	0.50
1:B:299:THR:HB	1:B:304:LYS:HE3	1.92	0.50
1:B:305:ILE:CG2	1:B:306:VAL:H	2.19	0.50
1:C:423:LEU:HD12	1:C:423:LEU:N	2.27	0.50
1:D:400:TYR:HB2	1:D:482:PRO:HG3	1.93	0.50
1:A:305:ILE:CG2	1:A:306:VAL:H	2.18	0.50
1:A:328:LYS:CD	1:A:337:LEU:HD21	2.40	0.50
1:A:335:VAL:HG12	1:A:336:SER:O	2.12	0.50
1:D:317:ILE:O	1:D:383:MET:HB2	2.12	0.50
1:B:103:GLN:OE1	1:B:300:ARG:HD3	2.11	0.50
1:C:80:LEU:HD12	1:C:80:LEU:N	2.26	0.50
1:C:299:THR:HB	1:C:304:LYS:HE3	1.94	0.50
1:C:184:ALA:HB2	1:C:245:ARG:HB3	1.94	0.50
1:A:162:ARG:NH1	1:A:170:GLY:N	2.60	0.49
1:D:94:ARG:NH1	1:D:94:ARG:HB3	2.26	0.49
1:C:138:LYS:HD2	1:C:197:LEU:O	2.12	0.49
1:C:286:ARG:HG3	1:C:286:ARG:HH11	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:LYS:HD3	1:B:47:LEU:HD23	1.94	0.49
1:A:96:GLN:H	1:A:96:GLN:CD	2.17	0.49
1:A:192:TRP:HA	1:A:253:VAL:CG1	2.40	0.49
1:B:53:PRO:HD2	1:B:54:GLU:OE2	2.13	0.49
1:A:329:ARG:HD3	1:A:333:GLY:O	2.13	0.49
1:C:317:ILE:O	1:C:383:MET:HB2	2.12	0.49
1:A:138:LYS:HD2	1:A:197:LEU:O	2.12	0.48
1:B:138:LYS:HD2	1:B:197:LEU:O	2.13	0.48
1:D:162:ARG:NH1	1:D:170:GLY:N	2.61	0.48
1:D:53:PRO:HD2	1:D:54:GLU:OE2	2.13	0.48
1:A:474:GLU:CD	1:A:474:GLU:H	2.21	0.48
1:A:53:PRO:HD2	1:A:54:GLU:OE2	2.14	0.48
1:A:317:ILE:O	1:A:383:MET:HB2	2.13	0.48
1:C:474:GLU:H	1:C:474:GLU:CD	2.20	0.48
1:D:138:LYS:HD2	1:D:197:LEU:O	2.14	0.48
1:A:154:PRO:HG2	1:A:157:LEU:HD12	1.94	0.48
1:A:105:VAL:H	1:A:241:GLN:HE22	1.60	0.48
1:B:80:LEU:HD12	1:B:80:LEU:N	2.28	0.48
1:D:80:LEU:HD12	1:D:80:LEU:N	2.29	0.48
1:D:154:PRO:HG2	1:D:157:LEU:HD12	1.96	0.48
1:A:108:VAL:HG23	1:A:245:ARG:HG2	1.96	0.48
1:B:184:ALA:HB2	1:B:245:ARG:HB3	1.96	0.48
1:C:162:ARG:NH1	1:C:170:GLY:N	2.61	0.48
1:C:53:PRO:HD2	1:C:54:GLU:OE2	2.13	0.47
1:D:489:HIS:HB3	2:D:2115:HOH:O	2.14	0.47
1:A:94:ARG:HG2	1:A:95:PRO:CD	2.44	0.47
1:C:335:VAL:HG12	1:C:336:SER:O	2.14	0.47
1:D:83:GLU:HB2	1:D:86:ARG:HG3	1.96	0.47
1:A:96:GLN:OE1	1:A:96:GLN:N	2.42	0.47
1:A:305:ILE:CG2	1:A:306:VAL:N	2.75	0.47
1:C:422:TYR:C	1:C:423:LEU:HD12	2.39	0.47
1:A:423:LEU:N	1:A:423:LEU:HD12	2.30	0.47
1:B:474:GLU:CD	1:B:474:GLU:H	2.21	0.47
1:C:150:HIS:O	1:C:152:PRO:HD3	2.15	0.47
1:B:271:THR:CG2	1:B:272:ASP:N	2.77	0.47
1:D:184:ALA:HB2	1:D:245:ARG:HB3	1.96	0.47
1:B:105:VAL:H	1:B:241:GLN:HE22	1.60	0.47
1:C:154:PRO:HG2	1:C:157:LEU:HD12	1.97	0.47
1:C:192:TRP:HA	1:C:253:VAL:CG1	2.40	0.47
1:D:108:VAL:HG23	1:D:245:ARG:HG2	1.96	0.47
1:D:422:TYR:C	1:D:423:LEU:HD12	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:ARG:HB3	1:C:94:ARG:CZ	2.45	0.47
1:D:150:HIS:O	1:D:152:PRO:HD3	2.15	0.47
1:A:94:ARG:HB3	1:A:94:ARG:HH11	1.80	0.47
1:A:271:THR:CG2	1:A:272:ASP:N	2.77	0.47
1:B:422:TYR:C	1:B:423:LEU:HD12	2.40	0.47
1:C:271:THR:CG2	1:C:272:ASP:N	2.76	0.47
1:D:241:GLN:HE21	1:D:307:ARG:NH1	2.12	0.47
1:A:422:TYR:C	1:A:423:LEU:HD12	2.40	0.47
1:B:162:ARG:NH1	1:B:170:GLY:N	2.62	0.47
1:A:241:GLN:HE21	1:A:307:ARG:NH1	2.11	0.46
1:C:241:GLN:HE21	1:C:307:ARG:NH1	2.14	0.46
1:A:94:ARG:HG2	1:A:95:PRO:HD2	1.97	0.46
1:B:192:TRP:HA	1:B:253:VAL:CG1	2.42	0.46
1:B:335:VAL:HG12	1:B:336:SER:O	2.14	0.46
1:D:96:GLN:OE1	1:D:96:GLN:N	2.48	0.46
1:D:298:ILE:HG12	1:D:299:THR:N	2.30	0.46
1:D:474:GLU:CD	1:D:474:GLU:H	2.21	0.46
1:A:96:GLN:O	1:A:97:ASN:OD1	2.34	0.46
1:C:105:VAL:H	1:C:241:GLN:HE22	1.63	0.46
1:D:191:ALA:O	1:D:195:ASP:HB2	2.16	0.46
1:A:60:TRP:O	1:A:129:HIS:HE1	1.98	0.46
1:C:95:PRO:HB2	1:C:98:PHE:HB2	1.98	0.46
1:D:60:TRP:O	1:D:129:HIS:HE1	1.97	0.46
1:B:108:VAL:HG23	1:B:245:ARG:HG2	1.96	0.46
1:A:329:ARG:HG3	1:A:329:ARG:NH1	2.31	0.46
1:C:83:GLU:HB2	1:C:86:ARG:HG3	1.96	0.46
1:C:253:VAL:O	1:C:253:VAL:HG12	2.16	0.46
1:D:423:LEU:HD12	1:D:423:LEU:N	2.30	0.46
1:B:288:TRP:CG	1:B:289:PHE:N	2.84	0.46
1:A:237:TYR:CZ	1:A:300:ARG:HB3	2.51	0.46
1:B:154:PRO:HG2	1:B:157:LEU:HD12	1.98	0.46
1:B:241:GLN:HE21	1:B:307:ARG:NH1	2.12	0.46
1:B:276:GLU:O	1:B:280:MET:HG3	2.16	0.46
1:D:271:THR:CG2	1:D:272:ASP:N	2.78	0.46
1:B:60:TRP:O	1:B:129:HIS:HE1	1.99	0.45
1:B:83:GLU:HB2	1:B:86:ARG:HG3	1.98	0.45
1:C:60:TRP:O	1:C:129:HIS:HE1	1.98	0.45
1:A:150:HIS:O	1:A:152:PRO:HD3	2.16	0.45
1:A:276:GLU:O	1:A:280:MET:HG3	2.16	0.45
1:B:307:ARG:HB3	1:B:309:ASP:OD1	2.16	0.45
1:D:237:TYR:CE1	1:D:300:ARG:HB3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ARG:HH11	1:A:300:ARG:HG2	1.81	0.45
1:C:98:PHE:CZ	1:C:181:TYR:HB2	2.51	0.45
1:D:300:ARG:HG2	1:D:300:ARG:HH11	1.82	0.45
1:A:94:ARG:HB3	1:A:94:ARG:CZ	2.47	0.45
1:A:138:LYS:HE2	1:A:138:LYS:HA	1.98	0.45
1:A:298:ILE:HG12	1:A:299:THR:N	2.31	0.45
1:B:138:LYS:HA	1:B:138:LYS:HE2	1.99	0.45
1:B:298:ILE:HG12	1:B:299:THR:N	2.32	0.45
1:D:105:VAL:H	1:D:241:GLN:HE22	1.62	0.45
1:B:150:HIS:O	1:B:152:PRO:HD3	2.17	0.45
1:B:423:LEU:HD12	1:B:423:LEU:N	2.31	0.45
1:B:329:ARG:HG3	1:B:329:ARG:HH11	1.82	0.45
1:C:237:TYR:CE1	1:C:300:ARG:HB3	2.52	0.45
1:C:298:ILE:HG12	1:C:299:THR:N	2.32	0.45
1:D:335:VAL:HG12	1:D:336:SER:O	2.16	0.45
1:B:191:ALA:O	1:B:195:ASP:HB2	2.16	0.45
1:D:446:VAL:HG22	1:D:453:LEU:CD2	2.47	0.45
1:D:288:TRP:CG	1:D:289:PHE:N	2.84	0.44
1:D:291:ASP:O	1:D:295:ARG:HB2	2.17	0.44
1:C:108:VAL:HG23	1:C:245:ARG:HG2	1.98	0.44
1:C:191:ALA:O	1:C:195:ASP:HB2	2.17	0.44
1:A:298:ILE:C	1:A:304:LYS:HE2	2.42	0.44
1:D:298:ILE:HG12	1:D:299:THR:H	1.83	0.44
1:A:162:ARG:HH12	1:A:170:GLY:N	2.16	0.44
1:C:300:ARG:HD2	1:C:300:ARG:O	2.17	0.44
1:D:276:GLU:O	1:D:280:MET:HG3	2.17	0.44
1:A:83:GLU:HB2	1:A:86:ARG:HG3	1.99	0.44
1:A:138:LYS:HE3	1:A:142:LEU:O	2.18	0.44
1:A:191:ALA:O	1:A:195:ASP:HB2	2.17	0.44
1:D:16:GLY:HA2	1:D:20:GLU:HG3	2.00	0.44
1:A:104:ASP:CA	1:A:241:GLN:HE22	2.24	0.44
1:B:329:ARG:HD3	1:B:329:ARG:C	2.41	0.44
1:C:245:ARG:HA	1:C:245:ARG:NE	2.33	0.44
1:C:3:SER:HA	1:C:470:ALA:CB	2.48	0.44
1:D:1:MET:HG2	1:D:2:TYR:N	2.33	0.44
1:D:138:LYS:HE3	1:D:142:LEU:O	2.18	0.44
1:D:112:GLU:O	1:D:116:LYS:HG2	2.18	0.43
1:D:150:HIS:O	1:D:151:TRP:CB	2.66	0.43
1:A:3:SER:HA	1:A:470:ALA:CB	2.48	0.43
1:D:245:ARG:HA	1:D:245:ARG:NE	2.33	0.43
1:B:39:ASP:OD2	1:B:164:ARG:NH2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:GLN:CG	1:B:97:ASN:N	2.79	0.43
1:C:104:ASP:CA	1:C:241:GLN:HE22	2.24	0.43
1:C:288:TRP:CG	1:C:289:PHE:N	2.86	0.43
1:C:307:ARG:HB3	1:C:309:ASP:OD1	2.18	0.43
1:D:3:SER:HA	1:D:470:ALA:CB	2.47	0.43
1:A:117:ARG:HB2	1:A:117:ARG:HH11	1.84	0.43
1:B:253:VAL:HG12	1:B:253:VAL:O	2.17	0.43
1:C:298:ILE:HG12	1:C:299:THR:H	1.84	0.43
1:C:1:MET:HG2	1:C:2:TYR:N	2.33	0.43
1:C:276:GLU:O	1:C:280:MET:HG3	2.19	0.43
1:A:332:LYS:HG3	1:A:333:GLY:N	2.27	0.43
1:A:446:VAL:HG22	1:A:453:LEU:CD2	2.49	0.43
1:C:332:LYS:HE3	1:C:332:LYS:HA	2.01	0.43
1:D:307:ARG:HB3	1:D:309:ASP:OD1	2.19	0.43
1:A:112:GLU:O	1:A:116:LYS:HG2	2.19	0.43
1:A:288:TRP:CG	1:A:289:PHE:N	2.86	0.43
1:A:300:ARG:HG2	1:A:300:ARG:NH1	2.33	0.43
1:C:305:ILE:CG2	1:C:306:VAL:H	2.21	0.43
1:C:112:GLU:O	1:C:116:LYS:HG2	2.19	0.42
1:C:138:LYS:HE3	1:C:142:LEU:O	2.19	0.42
1:C:291:ASP:O	1:C:295:ARG:HB2	2.17	0.42
1:A:298:ILE:HG12	1:A:299:THR:H	1.84	0.42
1:C:137:LEU:HD12	1:C:142:LEU:HD12	2.01	0.42
1:A:43:MET:HE1	2:A:2023:HOH:O	2.18	0.42
1:A:304:LYS:HD3	1:A:305:ILE:N	2.35	0.42
1:B:138:LYS:HE3	1:B:142:LEU:O	2.19	0.42
1:C:117:ARG:HB2	1:C:117:ARG:HH11	1.84	0.42
1:A:253:VAL:HG12	1:A:253:VAL:O	2.18	0.42
1:C:1:MET:HE2	1:C:472:THR:HA	2.02	0.42
1:D:39:ASP:OD2	1:D:164:ARG:NH2	2.52	0.42
1:A:150:HIS:O	1:A:151:TRP:CB	2.67	0.42
1:A:245:ARG:NE	1:A:245:ARG:HA	2.34	0.42
1:B:1:MET:HE2	1:B:472:THR:HA	2.02	0.42
1:C:446:VAL:HG22	1:C:453:LEU:CD2	2.49	0.42
1:D:138:LYS:HE2	1:D:138:LYS:HA	2.01	0.42
1:B:298:ILE:HG12	1:B:299:THR:H	1.84	0.42
1:D:117:ARG:HB2	1:D:117:ARG:HH11	1.83	0.42
1:D:162:ARG:HH12	1:D:170:GLY:N	2.18	0.42
1:B:234:ARG:HG3	1:B:234:ARG:HH11	1.84	0.42
1:C:94:ARG:HG2	1:C:95:PRO:HD2	2.02	0.42
1:C:138:LYS:HE2	1:C:138:LYS:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:GLU:OE1	1:A:305:ILE:HD11	2.20	0.42
1:B:76:LYS:CA	1:B:142:LEU:HD23	2.47	0.42
1:B:112:GLU:O	1:B:116:LYS:HG2	2.19	0.42
1:B:245:ARG:HA	1:B:245:ARG:NE	2.35	0.42
1:C:16:GLY:HA2	1:C:20:GLU:HG3	2.01	0.42
1:C:262:TYR:O	1:C:320:ASN:HB2	2.20	0.42
1:D:241:GLN:NE2	1:D:307:ARG:HH22	2.10	0.41
1:D:253:VAL:O	1:D:253:VAL:HG12	2.18	0.41
1:A:1:MET:HG2	1:A:2:TYR:N	2.34	0.41
1:A:16:GLY:HA2	1:A:20:GLU:HG3	2.02	0.41
1:D:234:ARG:HH11	1:D:234:ARG:HG3	1.86	0.41
1:D:349:VAL:CG1	1:D:353:GLY:HA2	2.50	0.41
1:A:241:GLN:NE2	1:A:307:ARG:HH22	2.12	0.41
1:D:79:ARG:NH2	1:D:205:ASN:OD1	2.51	0.41
1:D:139:SER:HA	2:D:2041:HOH:O	2.20	0.41
1:A:137:LEU:HD12	1:A:142:LEU:HD12	2.03	0.41
1:A:307:ARG:HB3	1:A:309:ASP:OD1	2.20	0.41
1:B:11:GLY:HA3	1:B:77:ILE:O	2.20	0.41
1:B:34:TYR:CZ	1:B:38:HIS:CE1	3.09	0.41
1:B:463:ARG:HD2	2:B:2094:HOH:O	2.20	0.41
1:B:1:MET:HG2	1:B:2:TYR:N	2.36	0.41
1:C:162:ARG:HH12	1:C:170:GLY:N	2.18	0.41
1:C:189:TYR:CZ	1:C:193:LYS:HD2	2.56	0.41
1:D:300:ARG:HG2	1:D:300:ARG:NH1	2.35	0.41
1:A:228:SER:HA	1:B:41:GLU:OE2	2.20	0.41
1:A:291:ASP:O	1:A:295:ARG:HB2	2.20	0.41
1:B:3:SER:HA	1:B:470:ALA:CB	2.48	0.41
1:B:27:GLU:H	1:B:27:GLU:CD	2.29	0.41
1:D:60:TRP:O	1:D:129:HIS:CE1	2.73	0.41
1:D:349:VAL:HG13	1:D:353:GLY:HA2	2.03	0.41
1:A:39:ASP:OD2	1:A:164:ARG:NH2	2.54	0.41
1:A:234:ARG:HG3	1:A:234:ARG:HH11	1.86	0.41
1:B:5:PRO:HD3	2:B:2002:HOH:O	2.19	0.41
1:D:463:ARG:HD2	2:D:2109:HOH:O	2.21	0.41
1:A:1:MET:HE2	1:A:472:THR:HA	2.02	0.41
1:A:17:PHE:HB2	1:A:152:PRO:HG2	2.03	0.41
1:D:1:MET:HE2	1:D:472:THR:HA	2.03	0.41
1:D:88:PHE:CE1	1:D:118:LEU:HD22	2.56	0.41
1:D:151:TRP:N	1:D:151:TRP:CD1	2.89	0.41
1:B:17:PHE:HB2	1:B:152:PRO:HG2	2.03	0.41
1:B:88:PHE:CE1	1:B:118:LEU:HD22	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:ASP:O	1:B:295:ARG:HB2	2.21	0.41
1:C:52:LEU:HA	1:C:53:PRO:HD3	1.92	0.41
1:D:189:TYR:CZ	1:D:193:LYS:HD2	2.56	0.41
1:B:209:VAL:HG13	1:B:213:LEU:HD12	2.03	0.40
1:C:17:PHE:HB2	1:C:152:PRO:HG2	2.03	0.40
1:C:148:MET:HB2	1:C:203:THR:O	2.21	0.40
1:C:234:ARG:HG3	1:C:234:ARG:HH11	1.85	0.40
1:D:26:SER:O	1:D:86:ARG:NH2	2.51	0.40
1:B:52:LEU:HB3	1:B:54:GLU:OE2	2.20	0.40
1:A:90:ASN:HA	1:A:91:PRO:HD3	1.93	0.40
1:B:131:ARG:O	1:B:135:LYS:HG3	2.21	0.40
1:B:189:TYR:CZ	1:B:193:LYS:HD2	2.56	0.40
1:B:93:PRO:O	1:B:94:ARG:C	2.64	0.40
1:B:162:ARG:HH12	1:B:170:GLY:N	2.20	0.40
1:B:262:TYR:O	1:B:320:ASN:HB2	2.22	0.40
1:B:271:THR:CG2	1:B:273:LYS:HG2	2.51	0.40
1:C:398:ARG:HD2	1:C:458:SER:OG	2.21	0.40
1:D:52:LEU:HA	1:D:53:PRO:HD3	1.91	0.40
1:C:73:MET:HE2	1:C:459:SER:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	487/489 (100%)	463 (95%)	21 (4%)	3 (1%)	21	29
1	B	487/489 (100%)	463 (95%)	20 (4%)	4 (1%)	16	22
1	C	487/489 (100%)	467 (96%)	18 (4%)	2 (0%)	30	39
1	D	487/489 (100%)	465 (96%)	19 (4%)	3 (1%)	21	29
All	All	1948/1956 (100%)	1858 (95%)	78 (4%)	12 (1%)	21	29

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	TRP
1	A	331	GLY
1	B	151	TRP
1	C	151	TRP
1	D	151	TRP
1	B	331	GLY
1	B	332	LYS
1	A	426	SER
1	B	426	SER
1	C	426	SER
1	D	426	SER
1	D	333	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	421/421 (100%)	405 (96%)	16 (4%)	29	44
1	B	421/421 (100%)	405 (96%)	16 (4%)	29	44
1	C	421/421 (100%)	404 (96%)	17 (4%)	28	42
1	D	421/421 (100%)	403 (96%)	18 (4%)	26	39
All	All	1684/1684 (100%)	1617 (96%)	67 (4%)	28	42

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LEU
1	A	96	GLN
1	A	102	LYS
1	A	118	LEU
1	A	245	ARG
1	A	272	ASP
1	A	286	ARG
1	A	299	THR

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Mol	Chain	Res	Type
1	A	300	ARG
1	A	324	ARG
1	A	329	ARG
1	A	349	VAL
1	A	427	LEU
1	A	449	ASN
1	A	467	THR
1	A	474	GLU
1	B	80	LEU
1	B	102	LYS
1	B	118	LEU
1	B	245	ARG
1	B	272	ASP
1	B	286	ARG
1	B	299	THR
1	B	303	GLU
1	B	324	ARG
1	B	329	ARG
1	B	332	LYS
1	B	349	VAL
1	B	427	LEU
1	B	449	ASN
1	B	467	THR
1	B	474	GLU
1	C	29	LEU
1	C	80	LEU
1	C	94	ARG
1	C	102	LYS
1	C	118	LEU
1	C	245	ARG
1	C	272	ASP
1	C	286	ARG
1	C	299	THR
1	C	300	ARG
1	C	324	ARG
1	C	332	LYS
1	C	349	VAL
1	C	427	LEU
1	C	449	ASN
1	C	467	THR
1	C	474	GLU
1	D	80	LEU

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Mol	Chain	Res	Type
1	D	94	ARG
1	D	96	GLN
1	D	97	ASN
1	D	102	LYS
1	D	118	LEU
1	D	245	ARG
1	D	272	ASP
1	D	286	ARG
1	D	299	THR
1	D	324	ARG
1	D	329	ARG
1	D	332	LYS
1	D	349	VAL
1	D	427	LEU
1	D	449	ASN
1	D	467	THR
1	D	474	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	128	ASN
1	A	129	HIS
1	A	241	GLN
1	A	302	ASN
1	A	342	HIS
1	A	410	HIS
1	A	430	ASN
1	A	479	ASN
1	B	55	ASN
1	B	96	GLN
1	B	128	ASN
1	B	129	HIS
1	B	147	ASN
1	B	241	GLN
1	B	268	GLN
1	B	342	HIS
1	B	410	HIS
1	B	430	ASN
1	B	479	ASN
1	C	55	ASN

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Mol	Chain	Res	Type
1	C	128	ASN
1	C	129	HIS
1	C	147	ASN
1	C	241	GLN
1	C	342	HIS
1	C	410	HIS
1	C	430	ASN
1	C	479	ASN
1	D	55	ASN
1	D	62	ASN
1	D	97	ASN
1	D	128	ASN
1	D	129	HIS
1	D	147	ASN
1	D	241	GLN
1	D	264	ASN
1	D	302	ASN
1	D	342	HIS
1	D	410	HIS
1	D	430	ASN
1	D	477	HIS
1	D	479	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	489/489 (100%)	-0.00	14 (2%) 53 55	19, 28, 54, 76	0
1	B	489/489 (100%)	0.01	22 (4%) 38 38	19, 28, 56, 78	0
1	C	489/489 (100%)	0.10	21 (4%) 40 40	19, 29, 56, 84	0
1	D	489/489 (100%)	0.03	22 (4%) 38 38	18, 28, 53, 72	0
All	All	1956/1956 (100%)	0.03	79 (4%) 42 43	18, 29, 55, 84	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	141	GLY	6.2
1	C	299	THR	5.1
1	C	304	LYS	4.8
1	D	113	ASN	4.1
1	C	68	ASN	4.0
1	D	139	SER	4.0
1	D	1	MET	3.8
1	C	302	ASN	3.7
1	B	303	GLU	3.6
1	A	97	ASN	3.4
1	A	304	LYS	3.3
1	B	301	GLY	3.3
1	D	304	LYS	3.2
1	D	332	LYS	3.2
1	B	304	LYS	3.1
1	C	96	GLN	3.1
1	B	97	ASN	3.0
1	C	489	HIS	3.0
1	D	101	SER	3.0
1	C	139	SER	2.9
1	A	113	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	140	ARG	2.9
1	A	414	ASN	2.8
1	C	426	SER	2.8
1	A	305	ILE	2.7
1	A	96	GLN	2.7
1	A	489	HIS	2.7
1	B	87	GLN	2.7
1	B	96	GLN	2.7
1	B	332	LYS	2.7
1	B	252	SER	2.7
1	D	303	GLU	2.6
1	C	98	PHE	2.6
1	D	112	GLU	2.6
1	D	97	ASN	2.6
1	A	117	ARG	2.5
1	B	113	ASN	2.5
1	D	96	GLN	2.5
1	C	252	SER	2.5
1	D	111	ASN	2.5
1	D	426	SER	2.5
1	D	102	LYS	2.5
1	A	1	MET	2.5
1	C	301	GLY	2.5
1	A	426	SER	2.5
1	D	117	ARG	2.4
1	C	6	ASN	2.4
1	C	113	ASN	2.4
1	C	1	MET	2.4
1	D	94	ARG	2.4
1	B	333	GLY	2.4
1	B	489	HIS	2.4
1	B	329	ARG	2.3
1	B	305	ILE	2.3
1	B	302	ASN	2.3
1	B	1	MET	2.3
1	D	335	VAL	2.3
1	D	329	ARG	2.3
1	C	305	ILE	2.2
1	B	94	ARG	2.2
1	C	253	VAL	2.2
1	D	2	TYR	2.2
1	B	300	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	303	GLU	2.2
1	C	95	PRO	2.2
1	A	109	GLU	2.2
1	C	335	VAL	2.1
1	A	332	LYS	2.1
1	B	299	THR	2.1
1	B	109	GLU	2.1
1	A	275	MET	2.1
1	A	302	ASN	2.1
1	C	273	LYS	2.1
1	B	98	PHE	2.1
1	B	99	ASP	2.1
1	D	336	SER	2.1
1	C	97	ASN	2.0
1	B	426	SER	2.0
1	D	271	THR	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.