



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 02:29 AM UTC

PDB ID : 1PMK / pdb_00001pmk
Title : KRINGLE-KRINGLE INTERACTIONS IN MULTIMER KRINGLE STRUCTURES
Authors : Padmanabhan, K.; Tulinsky, A.
Deposited on : 1994-04-25
Resolution : 2.25 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

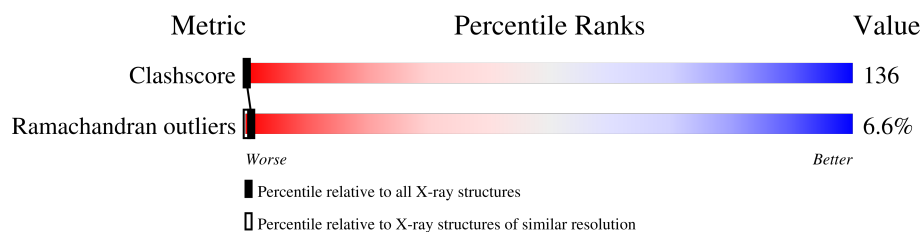
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	88	
1	B	88	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1362 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 PLASMINOGEN KRINGLE 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	78	Total	C	N	O	S	0	0	0
			621	380	113	120	8			
1	B	78	Total	C	N	O	S	0	0	0
			621	380	113	120	8			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	54	Total	O	0	0
			54	54		
2	B	66	Total	O	0	0
			66	66		

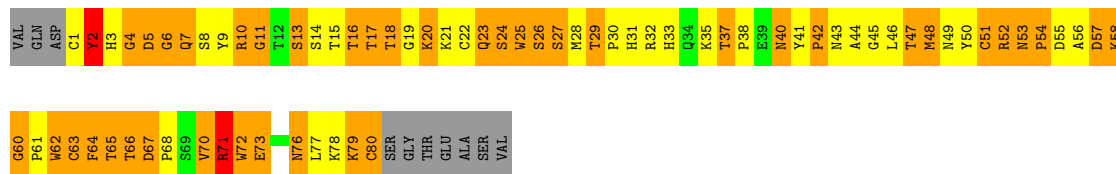
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

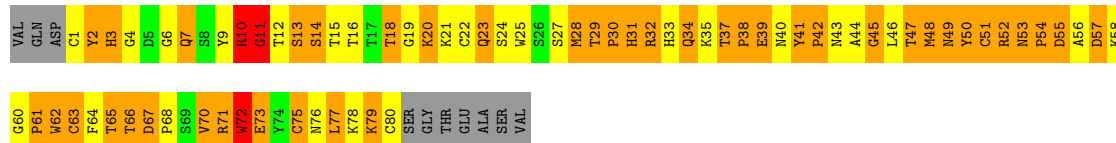
• Molecule 1: PLASMINOGEN KRINGLE 4

Chain A:  7% 33% 47% 11%



• Molecule 1: PLASMINOGEN KRINGLE 4

Chain B:  7% 31% 48% 11%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	32.73Å 49.09Å 46.15Å 90.00° 100.60° 90.00°	Depositor
Resolution (Å)	7.00 – 2.25	Depositor
% Data completeness (in resolution range)	(Not available) (7.00-2.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.165 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1362	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

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	1.28	1/641 (0.2%)	2.96	83/870 (9.5%)
1	B	1.27	0/641	3.04	90/870 (10.3%)
All	All	1.27	1/1282 (0.1%)	3.00	173/1740 (9.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	37	THR	CA-CB	5.21	1.61	1.53

All (173) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	39	GLU	CB-CG-CD	14.22	136.78	112.60
1	B	43	ASN	CA-CB-CG	-14.00	98.60	112.60
1	A	71	ARG	NE-CZ-NH1	13.12	134.62	121.50
1	A	53	ASN	O-C-N	12.56	130.91	121.12
1	B	34	GLN	CB-CG-CD	11.29	131.79	112.60
1	B	66	THR	N-CA-CB	11.26	127.67	110.44
1	A	23	GLN	O-C-N	11.25	135.68	122.85
1	A	25	TRP	CA-CB-CG	11.11	134.70	113.60
1	B	42	PRO	O-C-N	10.74	134.59	122.24
1	B	55	ASP	O-C-N	10.64	133.91	122.20
1	A	57	ASP	CA-CB-CG	-10.62	101.98	112.60
1	B	75	CYS	N-CA-CB	-9.80	96.34	110.45
1	B	61	PRO	N-CA-C	-9.73	94.67	111.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	VAL	N-CA-CB	-9.07	101.17	112.15
1	B	55	ASP	CA-C-O	-9.06	110.83	120.89
1	A	52	ARG	NE-CZ-NH2	-8.76	111.32	119.20
1	A	57	ASP	CA-C-N	8.74	131.81	120.44
1	A	57	ASP	C-N-CA	8.74	131.81	120.44
1	B	75	CYS	CB-CA-C	8.68	132.95	113.33
1	B	43	ASN	CA-C-O	-8.66	111.36	120.55
1	B	79	LYS	CA-C-N	8.58	137.14	121.70
1	B	79	LYS	C-N-CA	8.58	137.14	121.70
1	A	40	ASN	OD1-CG-ND2	8.48	131.08	122.60
1	B	62	TRP	N-CA-CB	-8.30	95.93	110.87
1	B	31	HIS	N-CA-CB	8.20	123.60	110.65
1	B	43	ASN	O-C-N	8.10	130.70	122.12
1	A	14	SER	N-CA-CB	-8.05	98.71	110.70
1	B	13	SER	N-CA-C	-8.04	97.16	109.95
1	B	65	THR	N-CA-CB	-8.02	98.89	110.36
1	B	29	THR	N-CA-CB	8.02	125.23	110.08
1	B	43	ASN	OD1-CG-ND2	7.96	130.56	122.60
1	A	43	ASN	CA-CB-CG	-7.83	104.77	112.60
1	B	2	TYR	CA-C-N	7.71	133.45	121.19
1	B	2	TYR	C-N-CA	7.71	133.45	121.19
1	B	28	MET	O-C-N	7.67	130.59	122.23
1	B	31	HIS	CA-C-N	7.66	133.60	121.66
1	B	31	HIS	C-N-CA	7.66	133.60	121.66
1	A	7	GLN	CA-C-N	7.64	133.49	120.72
1	A	7	GLN	C-N-CA	7.64	133.49	120.72
1	A	67	ASP	CB-CA-C	7.59	117.69	110.17
1	A	72	TRP	CA-C-O	-7.59	112.97	121.40
1	A	4	GLY	CA-C-O	-7.59	110.78	121.40
1	A	50	TYR	N-CA-C	7.55	120.75	110.55
1	B	29	THR	CA-CB-CG2	7.55	123.34	110.50
1	A	27	SER	O-C-N	7.41	131.43	122.68
1	A	25	TRP	CB-CA-C	7.36	126.33	110.21
1	B	51	CYS	O-C-N	7.35	131.85	122.89
1	A	62	TRP	CA-C-O	7.34	130.34	121.68
1	A	23	GLN	CA-C-O	-7.32	113.17	121.81
1	A	11	GLY	O-C-N	7.19	131.68	122.91
1	B	29	THR	O-C-N	7.16	127.35	120.55
1	B	23	GLN	N-CA-C	-7.10	100.97	110.55
1	A	6	GLY	CA-C-O	7.10	128.18	119.88
1	A	65	THR	CB-CA-C	7.09	121.64	109.51
1	B	13	SER	N-CA-CB	7.06	122.15	110.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	ASN	CA-C-O	-7.01	114.31	121.32
1	B	49	ASN	CA-C-O	-7.00	113.08	121.34
1	A	7	GLN	N-CA-C	-7.00	104.21	112.89
1	B	66	THR	N-CA-C	-6.96	104.53	113.16
1	A	5	ASP	CA-CB-CG	6.96	119.56	112.60
1	A	3	HIS	O-C-N	6.90	129.95	122.64
1	A	71	ARG	NH1-CZ-NH2	-6.88	110.36	119.30
1	B	18	THR	N-CA-C	-6.88	105.51	114.04
1	B	55	ASP	N-CA-CB	6.86	121.47	109.18
1	B	49	ASN	CA-C-N	-6.80	110.91	122.67
1	B	49	ASN	C-N-CA	-6.80	110.91	122.67
1	B	34	GLN	CA-C-O	6.79	128.28	120.00
1	A	52	ARG	CD-NE-CZ	6.70	133.78	124.40
1	B	7	GLN	CG-CD-NE2	6.70	126.44	116.40
1	B	18	THR	CB-CA-C	6.66	121.00	109.15
1	A	71	ARG	N-CA-CB	6.63	121.70	110.49
1	A	71	ARG	CD-NE-CZ	6.59	133.63	124.40
1	A	67	ASP	CA-C-O	6.58	124.89	119.36
1	A	73	GLU	CA-CB-CG	6.56	127.22	114.10
1	A	25	TRP	N-CA-C	-6.52	105.04	112.87
1	B	72	TRP	CA-C-O	6.44	129.71	120.51
1	B	10	ARG	CD-NE-CZ	-6.43	115.39	124.40
1	A	71	ARG	CG-CD-NE	6.41	126.11	112.00
1	B	51	CYS	CA-CB-SG	-6.40	99.67	114.40
1	B	43	ASN	N-CA-C	6.32	118.17	111.28
1	B	31	HIS	CA-C-O	6.32	127.03	120.40
1	B	63	CYS	O-C-N	6.31	130.36	123.10
1	B	3	HIS	CB-CA-C	-6.31	100.37	110.84
1	B	53	ASN	N-CA-CB	6.29	120.46	111.21
1	A	8	SER	CA-C-O	6.28	127.10	119.38
1	B	37	THR	CA-CB-CG2	6.26	121.15	110.50
1	B	50	TYR	CB-CA-C	6.24	121.25	110.45
1	A	72	TRP	O-C-N	6.17	130.99	122.78
1	A	20	LYS	N-CA-C	-6.13	102.44	110.53
1	A	18	THR	O-C-N	6.12	129.07	121.84
1	B	53	ASN	O-C-N	6.09	126.85	121.37
1	B	56	ALA	N-CA-C	-6.08	103.10	110.88
1	A	65	THR	CA-C-N	6.06	134.06	121.94
1	A	65	THR	C-N-CA	6.06	134.06	121.94
1	B	56	ALA	O-C-N	6.04	129.99	121.83
1	B	2	TYR	CB-CA-C	6.04	122.44	110.42
1	B	66	THR	CA-CB-CG2	6.00	120.70	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	67	ASP	O-C-N	6.00	128.22	121.32
1	B	38	PRO	O-C-N	5.97	129.40	122.17
1	A	29	THR	O-C-N	5.96	125.78	120.48
1	A	79	LYS	CA-C-O	5.94	128.77	121.99
1	A	7	GLN	CB-CG-CD	5.92	122.66	112.60
1	A	79	LYS	N-CA-CB	5.92	118.48	110.38
1	A	23	GLN	OE1-CD-NE2	-5.91	116.69	122.60
1	B	7	GLN	CA-C-N	5.90	129.28	120.31
1	B	7	GLN	C-N-CA	5.90	129.28	120.31
1	A	13	SER	CA-C-N	5.90	132.23	122.73
1	A	13	SER	C-N-CA	5.90	132.23	122.73
1	A	63	CYS	CA-C-O	5.90	127.54	121.23
1	A	23	GLN	N-CA-C	-5.89	102.60	110.55
1	B	41	TYR	CA-C-N	5.85	125.99	119.32
1	B	41	TYR	C-N-CA	5.85	125.99	119.32
1	B	14	SER	CA-CB-OG	5.84	122.78	111.10
1	B	11	GLY	CA-C-O	5.83	130.72	120.57
1	A	24	SER	CA-C-O	5.83	127.52	120.81
1	A	80	CYS	CA-C-O	-5.82	110.91	120.80
1	B	57	ASP	CA-C-N	5.75	128.24	120.65
1	B	57	ASP	C-N-CA	5.75	128.24	120.65
1	A	64	PHE	CA-CB-CG	-5.74	108.06	113.80
1	A	16	THR	N-CA-CB	-5.70	101.75	110.42
1	A	66	THR	N-CA-C	-5.67	106.34	113.20
1	A	51	CYS	CB-CA-C	5.66	118.73	109.90
1	A	64	PHE	CA-C-O	-5.65	114.27	120.43
1	A	29	THR	CA-CB-CG2	5.58	119.99	110.50
1	A	42	PRO	O-C-N	5.58	128.47	122.17
1	A	29	THR	CA-CB-OG1	-5.56	101.26	109.60
1	A	10	ARG	CD-NE-CZ	5.56	132.18	124.40
1	A	64	PHE	CA-C-N	-5.55	112.57	122.07
1	A	64	PHE	C-N-CA	-5.55	112.57	122.07
1	A	29	THR	N-CA-CB	5.55	118.57	110.30
1	B	49	ASN	OD1-CG-ND2	5.47	128.07	122.60
1	A	51	CYS	CA-C-O	-5.46	114.69	120.92
1	B	23	GLN	CA-C-N	-5.46	113.62	121.42
1	B	23	GLN	C-N-CA	-5.46	113.62	121.42
1	B	71	ARG	NE-CZ-NH2	-5.43	114.31	119.20
1	A	72	TRP	N-CA-C	-5.42	101.21	108.86
1	A	26	SER	CA-CB-OG	5.41	121.92	111.10
1	A	3	HIS	N-CA-CB	5.39	121.88	110.29
1	B	73	GLU	O-C-N	5.39	129.30	123.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	18	THR	CA-CB-OG1	-5.37	101.55	109.60
1	B	34	GLN	OE1-CD-NE2	-5.33	117.27	122.60
1	B	20	LYS	O-C-N	-5.28	115.28	122.36
1	B	13	SER	CB-CA-C	5.28	119.04	109.37
1	B	70	VAL	O-C-N	5.27	128.66	123.18
1	B	79	LYS	N-CA-C	5.26	117.65	110.55
1	B	55	ASP	N-CA-C	-5.25	101.80	109.63
1	A	71	ARG	O-C-N	5.25	129.58	122.59
1	A	52	ARG	N-CA-C	5.23	115.45	108.38
1	B	43	ASN	CB-CG-OD1	-5.23	110.34	120.80
1	A	76	ASN	OD1-CG-ND2	5.22	127.82	122.60
1	A	58	LYS	N-CA-C	5.21	116.65	111.07
1	A	60	GLY	O-C-N	5.20	126.97	121.77
1	B	47	THR	CA-C-O	5.20	127.95	120.51
1	B	54	PRO	N-CA-C	-5.20	107.27	113.98
1	B	57	ASP	N-CA-C	-5.17	103.85	110.53
1	A	19	GLY	CA-C-O	5.16	124.84	119.06
1	B	34	GLN	CB-CA-C	5.16	119.86	110.11
1	B	52	ARG	NE-CZ-NH2	5.16	123.84	119.20
1	A	2	TYR	CA-CB-CG	5.15	123.17	113.90
1	A	8	SER	CA-CB-OG	5.15	121.39	111.10
1	B	76	ASN	CA-C-N	5.14	130.94	121.85
1	B	76	ASN	C-N-CA	5.14	130.94	121.85
1	A	76	ASN	N-CA-C	5.13	117.64	107.98
1	B	77	LEU	CA-C-O	-5.13	115.24	120.99
1	A	54	PRO	O-C-N	5.13	129.51	122.38
1	A	16	THR	O-C-N	5.12	128.67	122.68
1	B	52	ARG	NE-CZ-NH1	-5.11	116.39	121.50
1	B	72	TRP	CA-C-N	5.07	130.18	121.87
1	B	72	TRP	C-N-CA	5.07	130.18	121.87
1	B	32	ARG	O-C-N	5.05	129.59	122.77
1	A	64	PHE	O-C-N	5.05	128.97	123.27
1	B	63	CYS	CB-CA-C	-5.03	99.67	111.09
1	A	47	THR	CA-C-O	5.02	127.87	121.55

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	10	ARG	Sidechain

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	621	0	567	171	0
1	B	621	0	567	153	0
2	A	54	0	0	35	0
2	B	66	0	0	26	0
All	All	1362	0	1134	322	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 136.

All (322) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:MET:HB2	1:B:50:TYR:CE1	1.79	1.17
1:A:51:CYS:O	1:A:52:ARG:HD3	1.46	1.14
1:B:28:MET:HE3	1:B:32:ARG:HG2	1.23	1.12
1:A:53:ASN:HD21	1:A:57:ASP:CB	1.64	1.11
1:B:2:TYR:OH	1:B:60:GLY:HA2	1.56	1.06
1:B:30:PRO:HB3	2:B:260:HOH:O	1.55	1.05
1:A:7:GLN:OE1	1:A:56:ALA:HA	1.59	1.02
1:A:30:PRO:HA	2:A:264:HOH:O	1.60	1.02
1:B:65:THR:HG22	1:B:67:ASP:H	1.23	1.01
1:A:31:HIS:CE1	1:A:71:ARG:HA	1.95	1.00
1:A:2:TYR:HB3	1:A:78:LYS:O	1.62	1.00
1:A:29:THR:HG22	1:A:30:PRO:HD3	1.45	0.98
1:A:37:THR:OG1	1:A:40:ASN:ND2	1.98	0.97
1:A:28:MET:HG2	1:A:32:ARG:HA	1.46	0.96
1:A:44:ALA:HB3	1:A:46:LEU:HD13	1.46	0.96
1:A:73:GLU:HB3	2:A:282:HOH:O	1.63	0.96
1:A:78:LYS:HD3	1:A:79:LYS:O	1.65	0.94
1:A:53:ASN:OD1	1:A:55:ASP:O	1.84	0.94
1:A:21:LYS:O	1:A:66:THR:HG22	1.67	0.94
1:B:65:THR:HB	1:B:70:VAL:O	1.65	0.94
1:A:16:THR:CG2	1:A:20:LYS:HB2	1.99	0.93
1:B:70:VAL:HG11	1:B:73:GLU:OE1	1.69	0.93
1:A:53:ASN:HD21	1:A:57:ASP:HB2	1.34	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HB3	2:A:279:HOH:O	1.71	0.89
1:A:62:TRP:HB2	1:A:72:TRP:CZ3	2.07	0.89
1:A:20:LYS:HG3	1:A:73:GLU:OE2	1.73	0.89
1:A:63:CYS:O	1:A:72:TRP:HB2	1.73	0.88
1:B:2:TYR:OH	1:B:60:GLY:CA	2.22	0.88
1:A:16:THR:HG21	1:A:20:LYS:HB2	1.56	0.87
1:A:5:ASP:O	1:A:60:GLY:HA2	1.73	0.87
1:B:63:CYS:O	1:B:72:TRP:HB2	1.73	0.87
1:B:1:CYS:SG	1:B:2:TYR:N	2.48	0.86
1:B:18:THR:OG1	1:B:73:GLU:OE2	1.92	0.86
1:A:53:ASN:HD21	1:A:57:ASP:HB3	1.38	0.86
1:A:20:LYS:HD2	1:A:70:VAL:HG11	1.59	0.85
1:B:48:MET:HB2	1:B:50:TYR:HE1	1.40	0.85
1:A:28:MET:HE1	2:A:313:HOH:O	1.77	0.83
1:A:61:PRO:HG2	1:A:77:LEU:HD11	1.59	0.83
1:B:60:GLY:O	1:B:62:TRP:HD1	1.62	0.82
1:B:41:TYR:HB2	1:B:46:LEU:HD22	1.59	0.82
1:A:25:TRP:C	2:A:304:HOH:O	2.23	0.81
1:B:44:ALA:HB1	2:B:295:HOH:O	1.79	0.81
1:A:35:LYS:HZ1	1:A:64:PHE:HZ	1.26	0.80
1:A:20:LYS:CD	1:A:65:THR:HG21	2.12	0.80
1:A:54:PRO:O	2:A:298:HOH:O	1.97	0.80
1:A:48:MET:HE2	1:B:48:MET:HE1	1.64	0.79
1:A:53:ASN:ND2	1:A:57:ASP:HB3	1.97	0.79
1:A:29:THR:CG2	1:A:30:PRO:HD3	2.11	0.79
1:B:44:ALA:CB	2:B:295:HOH:O	2.28	0.79
1:A:2:TYR:CE1	1:A:4:GLY:O	2.35	0.79
1:B:77:LEU:HA	2:B:204:HOH:O	1.81	0.79
1:A:37:THR:HG23	2:A:313:HOH:O	1.80	0.79
1:A:53:ASN:HD22	1:A:62:TRP:HD1	1.31	0.78
1:B:65:THR:HG22	1:B:67:ASP:N	1.97	0.78
1:A:13:SER:O	1:A:51:CYS:SG	2.41	0.77
1:B:28:MET:CE	1:B:32:ARG:HG2	2.10	0.77
1:B:1:CYS:O	1:B:2:TYR:HB2	1.85	0.76
2:A:229:HOH:O	1:B:48:MET:HG3	1.84	0.76
1:B:23:GLN:O	2:B:290:HOH:O	2.03	0.76
1:B:71:ARG:HB2	2:B:239:HOH:O	1.86	0.76
1:A:53:ASN:ND2	1:A:57:ASP:CB	2.45	0.75
1:A:38:PRO:O	1:A:42:PRO:HA	1.86	0.75
1:A:53:ASN:HD22	1:A:62:TRP:CD1	2.03	0.75
1:B:25:TRP:O	1:B:33:HIS:NE2	2.18	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:CYS:O	1:B:52:ARG:HD2	1.87	0.74
1:B:2:TYR:O	1:B:80:CYS:N	2.20	0.74
1:B:16:THR:HA	1:B:75:CYS:HA	1.68	0.74
1:A:5:ASP:OD1	1:A:7:GLN:HG3	1.88	0.74
1:A:25:TRP:CZ2	1:A:52:ARG:HB2	2.23	0.74
1:A:5:ASP:HA	1:A:58:LYS:O	1.88	0.73
1:B:37:THR:HB	1:B:38:PRO:HD2	1.70	0.73
1:B:22:CYS:HB3	2:B:290:HOH:O	1.86	0.73
1:B:45:GLY:HA3	1:B:52:ARG:HE	1.51	0.73
1:A:25:TRP:HZ2	1:A:52:ARG:HB2	1.53	0.73
1:A:26:SER:HB3	2:A:287:HOH:O	1.88	0.73
1:B:3:HIS:HA	1:B:80:CYS:O	1.91	0.70
1:B:28:MET:HE3	1:B:32:ARG:CG	2.14	0.70
1:A:26:SER:HA	2:A:304:HOH:O	1.90	0.70
1:A:31:HIS:ND1	1:A:71:ARG:HA	2.06	0.69
1:A:35:LYS:NZ	1:A:64:PHE:HZ	1.90	0.69
1:A:7:GLN:OE1	1:A:55:ASP:O	2.10	0.69
1:A:33:HIS:CD2	1:A:35:LYS:H	2.09	0.69
1:A:55:ASP:C	1:A:57:ASP:H	2.01	0.69
1:B:23:GLN:NE2	1:B:27:SER:O	2.25	0.69
1:A:44:ALA:CB	1:A:46:LEU:HD13	2.21	0.69
1:A:47:THR:HB	2:A:236:HOH:O	1.92	0.69
1:A:65:THR:N	1:A:71:ARG:O	2.24	0.69
1:A:23:GLN:HE21	1:A:27:SER:C	2.01	0.68
1:A:55:ASP:OD1	1:A:62:TRP:NE1	2.24	0.68
1:B:45:GLY:O	1:B:46:LEU:HD12	1.93	0.67
1:B:34:GLN:CG	2:B:301:HOH:O	2.41	0.67
1:A:62:TRP:HB2	1:A:72:TRP:HZ3	1.60	0.67
1:A:62:TRP:HB2	1:A:72:TRP:CE3	2.30	0.67
1:B:35:LYS:HA	1:B:41:TYR:CE2	2.31	0.66
1:B:40:ASN:O	1:B:41:TYR:CG	2.49	0.66
1:B:2:TYR:CE2	1:B:61:PRO:HD2	2.30	0.66
1:A:2:TYR:CB	1:A:78:LYS:O	2.43	0.65
1:A:51:CYS:O	1:A:52:ARG:CD	2.34	0.65
1:A:47:THR:O	1:A:48:MET:C	2.40	0.65
1:B:41:TYR:HB2	1:B:46:LEU:CD2	2.26	0.65
1:A:20:LYS:HB3	1:A:65:THR:HG23	1.78	0.65
1:A:21:LYS:O	1:A:66:THR:CG2	2.42	0.64
1:B:22:CYS:HB2	1:B:49:ASN:HB2	1.79	0.64
1:A:2:TYR:CD1	1:A:79:LYS:HG3	2.33	0.64
1:B:24:SER:CB	2:B:247:HOH:O	2.46	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:ASN:HA	1:B:62:TRP:NE1	2.12	0.64
1:A:33:HIS:HD2	1:A:35:LYS:H	1.46	0.63
1:A:20:LYS:HB3	1:A:65:THR:CG2	2.28	0.63
1:A:20:LYS:HD2	1:A:65:THR:HG21	1.80	0.63
1:B:2:TYR:CE1	1:B:4:GLY:O	2.52	0.63
1:B:38:PRO:HG3	1:B:46:LEU:O	1.99	0.62
1:B:30:PRO:CB	2:B:260:HOH:O	2.27	0.62
1:A:9:TYR:CE1	1:A:11:GLY:HA3	2.35	0.61
1:B:25:TRP:HA	2:B:195:HOH:O	2.00	0.61
1:B:48:MET:HB2	1:B:50:TYR:CZ	2.34	0.61
1:B:2:TYR:HD2	1:B:77:LEU:HD12	1.65	0.61
1:A:11:GLY:O	1:A:52:ARG:NH1	2.27	0.61
1:A:2:TYR:HE1	1:A:4:GLY:O	1.82	0.61
1:B:63:CYS:O	1:B:72:TRP:CB	2.46	0.61
1:B:67:ASP:OD1	1:B:68:PRO:HD2	2.01	0.61
1:A:5:ASP:HB2	2:A:275:HOH:O	2.01	0.61
1:A:64:PHE:CD1	1:A:72:TRP:HB3	2.36	0.61
1:A:1:CYS:CB	2:A:203:HOH:O	2.49	0.61
1:B:2:TYR:CD2	1:B:77:LEU:HD12	2.35	0.61
1:A:9:TYR:CZ	1:A:11:GLY:HA3	2.36	0.60
1:B:2:TYR:HB3	1:B:79:LYS:HA	1.82	0.60
1:A:1:CYS:HB3	2:A:203:HOH:O	2.01	0.60
1:A:44:ALA:HB3	1:A:46:LEU:CD1	2.27	0.60
1:B:57:ASP:OD1	1:B:58:LYS:N	2.34	0.60
1:B:24:SER:HB2	2:B:247:HOH:O	2.02	0.59
1:B:41:TYR:N	1:B:42:PRO:HD3	2.17	0.59
1:A:27:SER:OG	1:A:28:MET:N	2.36	0.58
1:B:62:TRP:CD1	1:B:62:TRP:N	2.68	0.58
1:A:4:GLY:O	1:A:5:ASP:C	2.46	0.58
1:A:64:PHE:HA	1:A:72:TRP:CB	2.33	0.58
1:B:3:HIS:CA	1:B:80:CYS:O	2.52	0.58
1:A:2:TYR:CE1	1:A:79:LYS:HG3	2.39	0.58
1:A:78:LYS:CD	1:A:79:LYS:O	2.46	0.58
1:B:4:GLY:O	1:B:79:LYS:HE3	2.04	0.58
1:B:34:GLN:H	1:B:35:LYS:HE3	1.69	0.58
1:A:58:LYS:C	2:A:279:HOH:O	2.47	0.57
1:B:46:LEU:HD11	1:B:54:PRO:HG3	1.86	0.57
1:A:20:LYS:HD3	1:A:65:THR:HG21	1.87	0.57
1:A:24:SER:C	1:A:26:SER:N	2.61	0.57
1:A:30:PRO:HB2	1:A:31:HIS:CD2	2.39	0.57
1:B:25:TRP:O	1:B:37:THR:HG22	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ALA:O	1:A:46:LEU:N	2.38	0.57
1:A:57:ASP:O	2:A:275:HOH:O	2.18	0.56
1:B:64:PHE:HD1	1:B:71:ARG:O	1.88	0.56
1:B:46:LEU:HD13	2:B:295:HOH:O	2.04	0.56
1:A:22:CYS:O	1:A:49:ASN:HB2	2.06	0.56
1:A:16:THR:HB	1:A:73:GLU:HG2	1.88	0.56
1:A:37:THR:HG1	1:A:40:ASN:HD22	1.46	0.56
1:A:7:GLN:OE1	1:A:56:ALA:CA	2.46	0.55
1:A:15:THR:HG22	1:A:20:LYS:O	2.06	0.55
1:A:9:TYR:OH	1:A:13:SER:HB2	2.07	0.55
1:A:62:TRP:CB	1:A:72:TRP:CZ3	2.85	0.55
1:A:23:GLN:HE22	1:A:28:MET:HA	1.71	0.55
1:A:26:SER:N	2:A:304:HOH:O	2.38	0.55
1:B:15:THR:HA	1:B:20:LYS:O	2.06	0.55
1:B:23:GLN:HG2	1:B:24:SER:N	2.22	0.55
1:A:65:THR:HB	1:A:70:VAL:HG12	1.88	0.55
1:A:52:ARG:O	1:A:61:PRO:HA	2.07	0.54
1:B:38:PRO:HB3	1:B:46:LEU:O	2.06	0.54
1:B:41:TYR:CB	1:B:46:LEU:HD22	2.35	0.54
1:B:41:TYR:N	1:B:42:PRO:CD	2.70	0.54
1:B:65:THR:HG21	1:B:70:VAL:HB	1.88	0.54
1:B:2:TYR:CE2	1:B:61:PRO:CD	2.90	0.54
1:A:65:THR:HB	1:A:70:VAL:O	2.08	0.54
1:B:66:THR:O	1:B:68:PRO:HD3	2.08	0.54
1:B:13:SER:O	1:B:50:TYR:HA	2.08	0.54
1:A:65:THR:CB	1:A:70:VAL:HG12	2.37	0.53
1:A:5:ASP:HB2	1:A:57:ASP:O	2.09	0.53
1:A:24:SER:C	1:A:26:SER:H	2.15	0.53
1:B:9:TYR:CZ	1:B:11:GLY:HA3	2.43	0.53
1:B:23:GLN:CD	1:B:30:PRO:HG2	2.34	0.53
1:A:26:SER:CA	2:A:304:HOH:O	2.50	0.53
1:A:70:VAL:HG13	2:A:282:HOH:O	2.09	0.53
1:B:14:SER:HB2	1:B:49:ASN:ND2	2.24	0.53
1:A:22:CYS:HB2	1:A:49:ASN:O	2.09	0.53
1:A:31:HIS:CE1	1:A:71:ARG:CA	2.82	0.53
1:A:7:GLN:OE1	1:A:55:ASP:C	2.52	0.52
1:A:1:CYS:SG	1:A:2:TYR:N	2.82	0.52
1:A:25:TRP:HE1	1:A:52:ARG:CG	2.23	0.52
1:A:7:GLN:HG2	1:A:53:ASN:O	2.09	0.52
1:A:9:TYR:O	1:A:10:ARG:HD3	2.10	0.51
1:A:37:THR:HG1	1:A:40:ASN:ND2	2.05	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:PRO:HG2	2:A:311:HOH:O	2.08	0.51
1:B:32:ARG:NH1	2:B:225:HOH:O	2.44	0.51
1:A:37:THR:CG2	2:A:313:HOH:O	2.48	0.51
1:A:62:TRP:CZ3	1:A:64:PHE:CE2	2.97	0.51
1:A:67:ASP:OD2	1:A:70:VAL:HG23	2.11	0.51
1:A:73:GLU:CB	2:A:282:HOH:O	2.37	0.51
1:B:48:MET:CB	1:B:50:TYR:HE1	2.20	0.51
1:A:44:ALA:CB	1:A:46:LEU:CD1	2.89	0.51
1:A:65:THR:N	1:A:72:TRP:HA	2.26	0.51
1:A:62:TRP:HZ3	1:A:64:PHE:CE2	2.30	0.50
1:B:2:TYR:HE1	1:B:4:GLY:O	1.93	0.50
1:B:2:TYR:HE2	1:B:61:PRO:HD2	1.75	0.50
1:B:20:LYS:HD3	1:B:67:ASP:HB2	1.94	0.50
1:A:53:ASN:ND2	1:A:62:TRP:CD1	2.79	0.49
1:B:44:ALA:HB3	2:B:295:HOH:O	2.03	0.49
1:B:52:ARG:HA	1:B:61:PRO:HB3	1.94	0.49
1:B:19:GLY:HA3	2:B:231:HOH:O	2.13	0.49
1:B:31:HIS:CE1	1:B:71:ARG:HA	2.47	0.49
1:B:38:PRO:CG	1:B:46:LEU:O	2.61	0.49
1:A:76:ASN:N	2:A:307:HOH:O	2.45	0.49
1:B:53:ASN:H	1:B:53:ASN:HD22	1.60	0.49
1:B:15:THR:HG22	1:B:20:LYS:O	2.13	0.49
1:B:78:LYS:HG3	1:B:79:LYS:N	2.28	0.48
1:A:62:TRP:CB	1:A:72:TRP:HZ3	2.24	0.48
1:B:1:CYS:O	1:B:2:TYR:CB	2.56	0.48
1:B:9:TYR:CE1	1:B:77:LEU:HD13	2.48	0.48
1:B:55:ASP:OD1	1:B:55:ASP:N	2.43	0.48
1:B:78:LYS:HG3	1:B:79:LYS:H	1.79	0.48
1:A:25:TRP:HE1	1:A:52:ARG:HG2	1.78	0.48
1:B:2:TYR:CZ	1:B:61:PRO:HD3	2.48	0.48
1:B:7:GLN:O	1:B:10:ARG:NH1	2.41	0.48
1:B:23:GLN:HG3	1:B:30:PRO:HG2	1.95	0.48
1:B:40:ASN:C	1:B:41:TYR:CG	2.92	0.48
1:B:75:CYS:SG	1:B:77:LEU:HD21	2.54	0.48
1:A:2:TYR:O	1:A:80:CYS:N	2.46	0.47
1:A:55:ASP:C	1:A:57:ASP:N	2.70	0.47
1:B:2:TYR:OH	1:B:60:GLY:HA3	2.13	0.47
1:B:22:CYS:HB2	1:B:49:ASN:O	2.13	0.47
1:A:65:THR:H	1:A:72:TRP:HA	1.80	0.47
1:A:33:HIS:HD2	1:A:35:LYS:N	2.12	0.47
1:A:38:PRO:O	1:A:42:PRO:CA	2.60	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HA	2:A:275:HOH:O	2.13	0.47
1:B:35:LYS:HA	1:B:41:TYR:CD2	2.48	0.47
1:B:23:GLN:CG	1:B:30:PRO:HG2	2.45	0.47
1:B:53:ASN:HB3	1:B:62:TRP:CD1	2.50	0.47
1:A:21:LYS:HG2	1:A:22:CYS:N	2.30	0.47
1:B:13:SER:O	1:B:50:TYR:HB3	2.13	0.47
1:B:24:SER:OG	1:B:27:SER:CB	2.63	0.47
1:B:44:ALA:O	1:B:46:LEU:N	2.48	0.47
1:B:64:PHE:HA	1:B:71:ARG:O	2.15	0.47
1:A:41:TYR:HB2	1:A:46:LEU:HD22	1.96	0.47
1:B:34:GLN:HG3	2:B:301:HOH:O	2.08	0.47
1:A:48:MET:CE	1:B:48:MET:HE1	2.42	0.47
1:A:25:TRP:O	2:A:304:HOH:O	2.21	0.46
1:A:37:THR:O	1:A:41:TYR:N	2.34	0.46
1:A:13:SER:HB3	1:A:51:CYS:SG	2.55	0.46
1:B:16:THR:HG21	1:B:73:GLU:HG3	1.98	0.46
1:B:34:GLN:HB3	2:B:301:HOH:O	2.14	0.46
1:A:16:THR:HG21	1:A:20:LYS:CB	2.38	0.46
1:B:46:LEU:CD1	2:B:295:HOH:O	2.61	0.46
1:A:23:GLN:NE2	1:A:27:SER:C	2.71	0.46
1:B:23:GLN:HE21	1:B:27:SER:C	2.24	0.46
1:A:9:TYR:HB3	2:A:261:HOH:O	2.16	0.46
1:A:5:ASP:OD1	1:A:6:GLY:N	2.49	0.45
1:A:16:THR:OG1	1:A:20:LYS:HB2	2.16	0.45
1:A:35:LYS:HE3	1:A:64:PHE:CE2	2.51	0.45
1:B:67:ASP:O	1:B:70:VAL:O	2.34	0.45
2:A:315:HOH:O	1:B:12:THR:HG23	2.17	0.45
1:B:54:PRO:HD2	1:B:62:TRP:CZ2	2.51	0.45
1:B:65:THR:CG2	1:B:67:ASP:HB3	2.46	0.45
1:B:71:ARG:O	1:B:72:TRP:HB3	2.16	0.45
1:A:68:PRO:HD3	2:A:222:HOH:O	2.15	0.45
1:B:2:TYR:HB3	1:B:78:LYS:O	2.16	0.45
1:B:34:GLN:CB	2:B:301:HOH:O	2.64	0.45
1:A:1:CYS:HB2	2:A:203:HOH:O	2.14	0.45
1:A:20:LYS:HA	2:A:189:HOH:O	2.16	0.45
1:A:47:THR:O	1:A:48:MET:O	2.35	0.45
1:B:21:LYS:HB2	1:B:21:LYS:HE3	1.79	0.45
1:B:38:PRO:CB	1:B:46:LEU:O	2.65	0.45
1:A:26:SER:CB	2:A:287:HOH:O	2.55	0.44
1:B:45:GLY:O	1:B:52:ARG:HD3	2.17	0.44
1:A:13:SER:O	1:A:51:CYS:N	2.44	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:MET:HB3	1:A:49:ASN:H	1.31	0.44
1:B:53:ASN:HA	1:B:62:TRP:HE1	1.79	0.44
1:B:19:GLY:CA	2:B:231:HOH:O	2.66	0.44
1:B:4:GLY:HA2	1:B:79:LYS:HE2	1.99	0.44
1:B:57:ASP:OD1	1:B:60:GLY:N	2.48	0.44
1:B:45:GLY:HA3	1:B:52:ARG:NE	2.29	0.44
1:B:57:ASP:OD1	1:B:58:LYS:HG2	2.18	0.44
1:A:64:PHE:CE1	1:A:71:ARG:HG2	2.53	0.43
1:B:50:TYR:CD1	1:B:50:TYR:N	2.86	0.43
1:B:78:LYS:CG	1:B:79:LYS:N	2.82	0.43
1:A:63:CYS:O	1:A:72:TRP:CB	2.58	0.43
1:B:54:PRO:HB3	2:B:295:HOH:O	2.18	0.43
1:A:23:GLN:CG	1:A:24:SER:N	2.82	0.43
1:B:31:HIS:NE2	1:B:70:VAL:O	2.52	0.43
1:B:53:ASN:C	1:B:55:ASP:N	2.75	0.43
1:A:40:ASN:HB3	1:A:41:TYR:CD1	2.53	0.43
1:B:4:GLY:HA2	1:B:79:LYS:CE	2.49	0.43
1:A:2:TYR:O	1:A:80:CYS:HB2	2.18	0.42
1:A:67:ASP:O	1:A:70:VAL:O	2.37	0.42
1:B:29:THR:N	1:B:30:PRO:CD	2.80	0.42
1:B:47:THR:O	1:B:48:MET:O	2.37	0.42
1:A:5:ASP:CG	1:A:7:GLN:HG3	2.44	0.42
1:A:10:ARG:C	2:A:315:HOH:O	2.62	0.42
1:A:41:TYR:HD2	2:A:298:HOH:O	2.02	0.42
1:B:16:THR:C	1:B:18:THR:H	2.26	0.42
1:B:24:SER:OG	1:B:27:SER:N	2.53	0.42
1:A:2:TYR:CD1	1:A:2:TYR:C	2.98	0.42
1:A:47:THR:HG23	2:B:271:HOH:O	2.20	0.42
1:A:23:GLN:HE22	1:A:28:MET:CA	2.32	0.42
1:B:53:ASN:C	1:B:55:ASP:H	2.27	0.41
1:A:17:THR:O	1:A:18:THR:CG2	2.68	0.41
1:B:18:THR:HG23	2:B:314:HOH:O	2.20	0.41
1:B:39:GLU:O	1:B:42:PRO:HD3	2.20	0.41
1:B:47:THR:HB	1:B:48:MET:H	1.46	0.41
1:A:16:THR:HG23	1:A:20:LYS:HB2	1.94	0.41
1:A:68:PRO:CD	2:A:222:HOH:O	2.69	0.41
1:B:30:PRO:CG	2:B:260:HOH:O	2.63	0.41
1:A:37:THR:HB	1:A:38:PRO:HD2	2.02	0.41
1:B:23:GLN:HB3	1:B:64:PHE:O	2.20	0.41
1:B:52:ARG:C	1:B:61:PRO:HA	2.46	0.41
1:A:17:THR:O	1:A:18:THR:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LYS:HB2	2:A:267:HOH:O	2.20	0.41
1:B:10:ARG:HH11	1:B:10:ARG:HD2	1.55	0.41
1:B:60:GLY:O	1:B:62:TRP:CD1	2.55	0.41
1:B:2:TYR:CZ	1:B:61:PRO:CD	3.04	0.41
1:A:35:LYS:NZ	1:A:64:PHE:CZ	2.77	0.40
1:A:53:ASN:CG	1:A:55:ASP:O	2.59	0.40
1:A:15:THR:HG22	1:A:21:LYS:HA	2.03	0.40
1:A:35:LYS:HE3	1:A:64:PHE:HE2	1.87	0.40
1:B:34:GLN:HG2	2:B:301:HOH:O	2.12	0.40
1:B:4:GLY:C	1:B:6:GLY:N	2.75	0.40
1:B:14:SER:O	1:B:21:LYS:HA	2.21	0.40
1:B:24:SER:OG	1:B:27:SER:HB2	2.21	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	76/88 (86%)	54 (71%)	17 (22%)	5 (7%)	1	0
1	B	76/88 (86%)	57 (75%)	14 (18%)	5 (7%)	1	0
All	All	152/176 (86%)	111 (73%)	31 (20%)	10 (7%)	1	0

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	48	MET
1	B	48	MET
1	A	17	THR
1	B	30	PRO
1	B	72	TRP
1	A	71	ARG

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Mol	Chain	Res	Type
1	B	11	GLY
1	B	45	GLY
1	A	2	TYR
1	A	45	GLY

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.