



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:44 PM UTC

PDB ID : 2UXV / pdb_00002uxv
Title : SufI Protein from Escherichia Coli
Authors : Tarry, M.J.; Roversi, P.; Sargent, F.; Berks, B.C.; Lea, S.M.
Deposited on : 2007-03-30
Resolution : 2.61 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

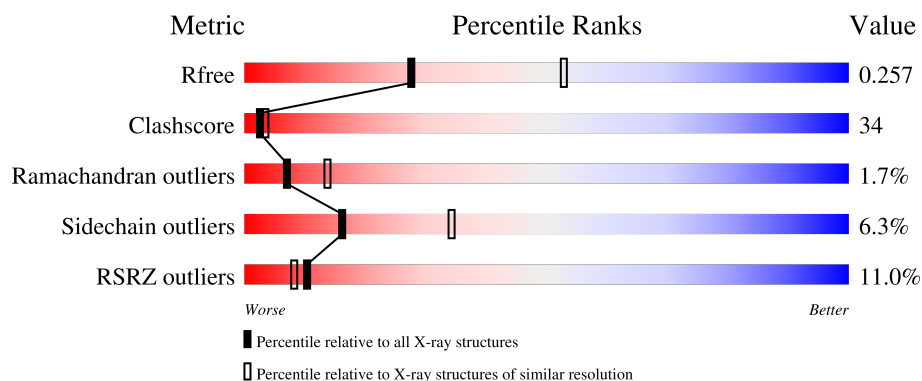
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	451	10% 49% 38% 5% • 6%
1	B	451	11% 47% 39% 6% • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6693 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 PROTEIN SUFI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	2	0
			3318	2108	587	609	14			
1	B	422	Total	C	N	O	S	0	1	0
			3313	2104	586	609	14			

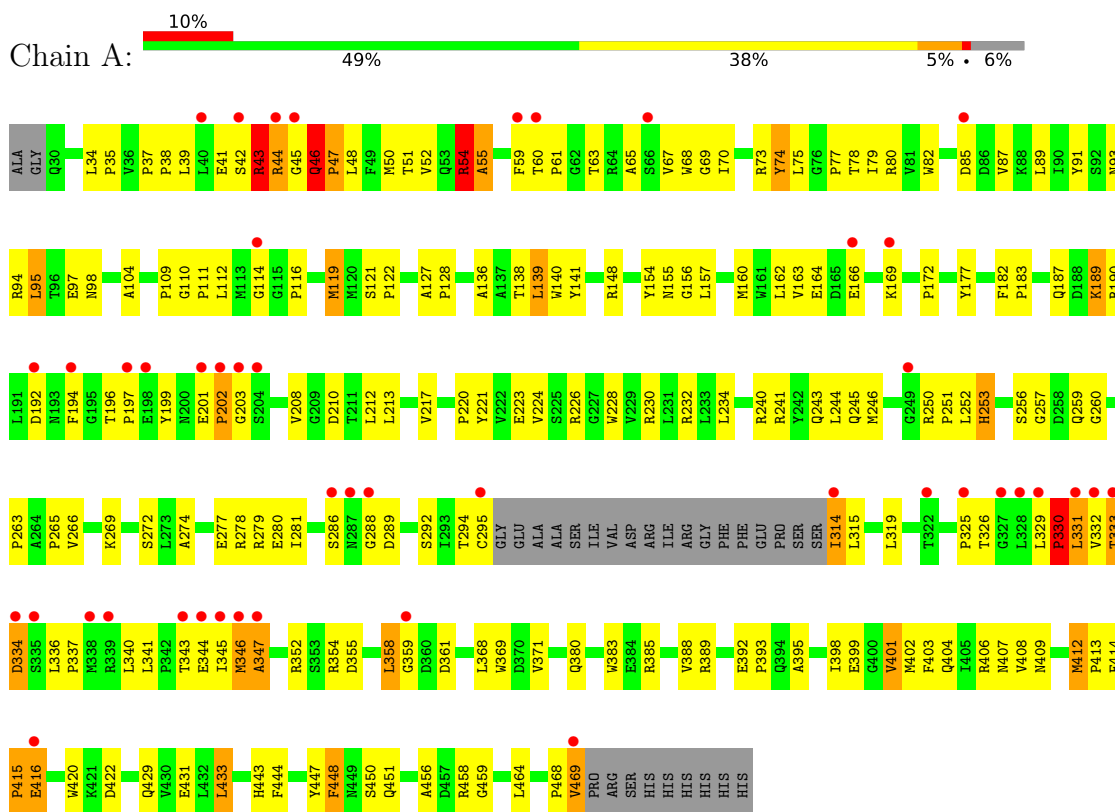
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	35	Total	O	0	0
			35	35		
2	B	27	Total	O	0	0
			27	27		

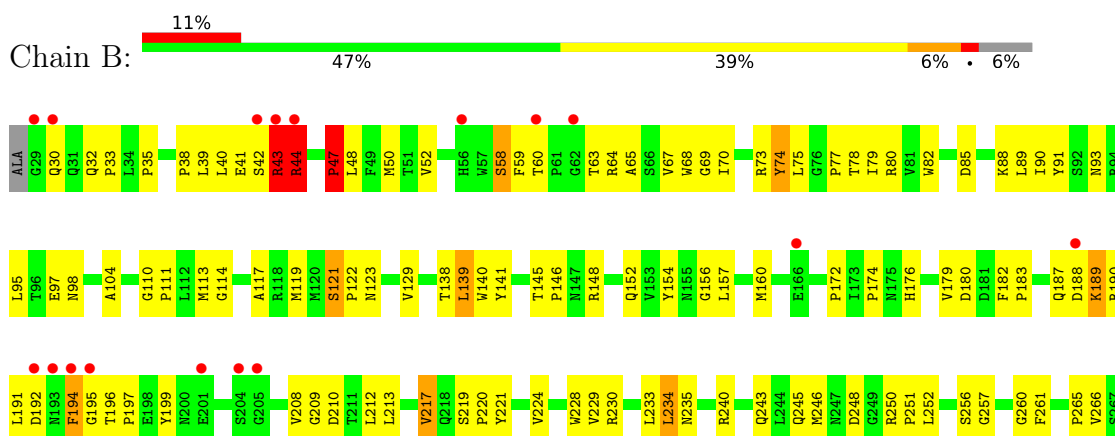
3 Residue-property plots

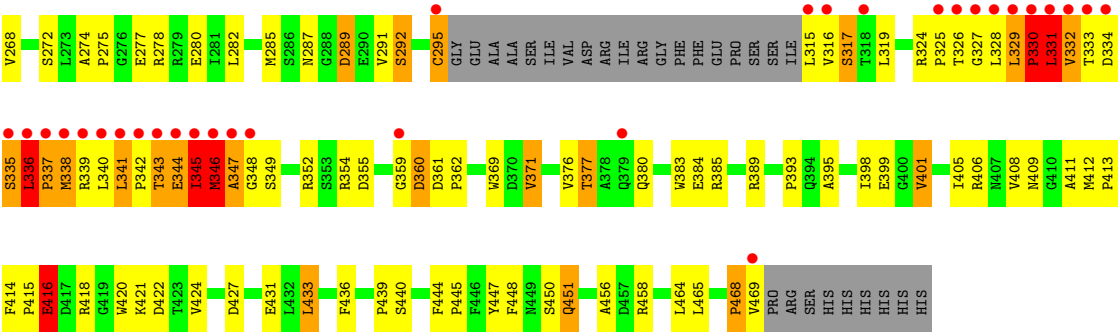
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: PROTEIN SUFI





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.12Å 48.88Å 131.96Å 90.00° 95.90° 90.00°	Depositor
Resolution (Å)	59.76 – 2.61 59.76 – 2.61	Depositor EDS
% Data completeness (in resolution range)	(Not available) (59.76-2.61) 98.4 (59.76-2.61)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.61Å)	Xtriage
Refinement program	TNT 5.13.1.0	Depositor
R, R_{free}	0.214 , (Not available) 0.227 , 0.257	Depositor DCC
R_{free} test set	1261 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	1.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6693	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.45 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1945e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.41	2/3411 (0.1%)	1.07	29/4660 (0.6%)
1	B	0.44	0/3406	1.08	32/4652 (0.7%)
All	All	0.43	2/6817 (0.0%)	1.08	61/9312 (0.7%)

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	A	0	3
1	B	0	2
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	47	PRO	N-CA	-5.22	1.40	1.47
1	A	55	ALA	CA-CB	-5.01	1.42	1.52

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	GLY	N-CA-C	-18.29	87.56	111.52
1	A	55	ALA	N-CA-C	12.92	127.34	108.34
1	B	338	MET	N-CA-C	-12.36	96.38	113.56
1	B	335	SER	N-CA-C	12.32	125.25	107.88
1	A	47	PRO	N-CA-C	-12.00	87.74	112.47
1	B	44	ARG	CB-CA-C	-10.69	94.23	110.26
1	A	54	ARG	CA-C-N	10.24	136.38	121.84
1	A	54	ARG	C-N-CA	10.24	136.38	121.84
1	A	334	ASP	N-CA-C	10.02	132.14	110.80
1	B	333	THR	N-CA-C	-9.71	101.45	113.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	GLN	N-CA-C	9.42	130.63	109.81
1	B	345	ILE	CA-C-N	8.39	136.80	121.70
1	B	345	ILE	C-N-CA	8.39	136.80	121.70
1	B	44	ARG	N-CA-C	8.29	121.77	109.59
1	B	30	GLN	N-CA-C	-8.23	98.30	110.48
1	A	333	THR	N-CA-C	8.13	120.81	110.24
1	B	325	PRO	N-CA-C	-8.10	97.11	112.01
1	A	415	PRO	CA-C-N	7.69	135.55	121.70
1	A	415	PRO	C-N-CA	7.69	135.55	121.70
1	B	44	ARG	CA-CB-CG	7.67	129.44	114.10
1	B	336	LEU	CA-C-N	7.12	128.74	119.84
1	B	336	LEU	C-N-CA	7.12	128.74	119.84
1	B	47	PRO	N-CA-C	-7.11	97.82	112.47
1	A	415	PRO	N-CA-C	7.09	127.08	112.47
1	B	345	ILE	N-CA-C	-7.03	94.72	109.34
1	A	347	ALA	N-CA-C	6.80	119.83	110.24
1	A	346	MET	N-CA-C	6.76	120.64	109.96
1	A	55	ALA	CB-CA-C	-6.58	106.94	116.54
1	A	45	GLY	N-CA-C	-6.57	97.61	113.18
1	A	87	VAL	N-CA-C	6.54	118.02	109.58
1	A	333	THR	CA-C-N	6.49	133.94	121.54
1	A	333	THR	C-N-CA	6.49	133.94	121.54
1	B	74	TYR	N-CA-C	-6.23	101.06	110.28
1	B	349	SER	N-CA-C	6.21	117.90	110.07
1	B	331	LEU	N-CA-C	6.02	123.62	110.80
1	B	416	GLU	N-CA-C	5.99	118.60	111.71
1	B	468	PRO	N-CA-C	-5.98	102.47	111.57
1	A	253	HIS	N-CA-C	5.87	117.86	108.41
1	A	114	GLY	CA-C-O	-5.81	116.38	122.01
1	B	344	GLU	N-CA-C	-5.81	102.86	110.53
1	A	74	TYR	N-CA-C	-5.81	101.03	110.20
1	B	336	LEU	N-CA-C	-5.59	101.87	110.32
1	A	330	PRO	CA-C-N	5.58	131.74	121.70
1	A	330	PRO	C-N-CA	5.58	131.74	121.70
1	B	189	LYS	N-CA-C	5.55	117.55	108.55
1	B	345	ILE	N-CA-CB	-5.40	102.31	111.23
1	B	187	GLN	N-CA-C	5.40	118.24	109.06
1	B	347	ALA	CA-C-N	5.39	131.41	121.70
1	B	347	ALA	C-N-CA	5.39	131.41	121.70
1	A	189	LYS	N-CA-C	5.35	117.81	109.52
1	B	43	ARG	N-CA-CB	5.32	119.47	110.49
1	B	325	PRO	CA-C-O	-5.32	115.49	122.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	345	ILE	CA-C-O	-5.29	114.16	120.78
1	A	187	GLN	N-CA-C	5.29	117.86	109.50
1	B	330	PRO	CA-C-N	5.14	131.36	121.54
1	B	330	PRO	C-N-CA	5.14	131.36	121.54
1	B	44	ARG	NE-CZ-NH1	5.12	126.62	121.50
1	A	448	PHE	N-CA-C	-5.10	100.86	109.07
1	A	450	SER	N-CA-C	-5.08	101.68	109.76
1	A	331	LEU	CA-C-N	5.04	130.77	121.70
1	A	331	LEU	C-N-CA	5.04	130.77	121.70

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	46	GLN	Peptide
1	A	54	ARG	Peptide,Mainchain
1	B	194	PHE	Sidechain
1	B	44	ARG	Mainchain

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	3318	0	3266	218	0
1	B	3313	0	3257	239	0
2	A	35	0	0	3	0
2	B	27	0	0	10	0
All	All	6693	0	6523	452	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 34.

All (452) 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:385:ARG:HH11	1:B:406:ARG:NE	1.46	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLU:HB3	1:A:326:THR:HG21	1.28	1.11
1:A:412:MET:HG3	1:A:413:PRO:HD2	1.27	1.10
1:B:228:TRP:HB2	1:B:332:VAL:HG13	1.30	1.10
1:B:265:PRO:HD3	1:B:340:LEU:HD13	1.32	1.10
1:B:345:ILE:HG12	1:B:346:MET:HG2	1.34	1.08
1:B:337:PRO:HB2	1:B:339:ARG:H	1.13	1.08
1:B:63:THR:HG21	1:B:195:GLY:HA3	1.33	1.08
1:B:228:TRP:CD1	1:B:332:VAL:HG22	1.95	1.01
1:B:385:ARG:NH1	1:B:406:ARG:HE	1.59	1.00
1:A:330:PRO:HG2	1:A:331:LEU:HA	1.40	1.00
1:B:337:PRO:HB3	1:B:339:ARG:HB3	1.43	1.00
1:A:395:ALA:HB3	1:A:451:GLN:HB3	1.41	0.99
1:A:406:ARG:HD2	1:A:433:LEU:HD22	1.43	0.98
1:A:55:ALA:HB1	1:A:67:VAL:H	1.28	0.97
1:A:385:ARG:HH12	1:A:406:ARG:NH2	1.62	0.97
1:B:345:ILE:HG12	1:B:346:MET:CG	1.94	0.97
1:B:174:PRO:HA	2:B:2011:HOH:O	1.63	0.96
1:A:347:ALA:HB2	1:A:404[B]:GLN:NE2	1.82	0.95
1:B:406:ARG:HD2	1:B:433:LEU:HD23	1.45	0.95
1:A:266:VAL:HG11	1:A:415:PRO:HB2	1.49	0.93
1:B:210:ASP:HA	1:B:315:LEU:HD23	1.50	0.92
1:B:377:THR:HG23	1:B:465:LEU:HD23	1.50	0.91
1:A:55:ALA:HB2	1:A:67:VAL:O	1.69	0.91
1:A:189:LYS:HG3	1:A:213:LEU:HD22	1.53	0.91
1:A:55:ALA:CB	1:A:67:VAL:H	1.85	0.89
1:A:329:LEU:HB3	1:A:330:PRO:HD2	1.55	0.89
1:B:369:TRP:CZ2	1:B:371:VAL:HG12	2.10	0.87
1:B:261:PHE:HB2	1:B:341:LEU:HD13	1.54	0.86
1:A:48:LEU:HB2	1:A:89:LEU:HD12	1.56	0.86
1:A:46:GLN:O	1:A:46:GLN:HG3	1.73	0.86
1:B:228:TRP:HD1	1:B:332:VAL:HG22	1.37	0.86
1:A:347:ALA:HB2	1:A:404[B]:GLN:HE21	1.40	0.86
1:A:352:ARG:HE	1:A:354:ARG:NH2	1.72	0.85
1:B:327:GLY:HA3	1:B:328:LEU:HG	1.58	0.85
1:A:52:VAL:HG12	1:A:95:LEU:HG	1.58	0.85
1:A:41:GLU:OE2	1:A:44:ARG:NH1	2.10	0.85
1:B:329:LEU:HB3	1:B:330:PRO:CD	2.08	0.83
1:A:246:MET:HE2	1:A:252:LEU:CD1	2.08	0.82
1:B:406:ARG:CD	1:B:433:LEU:HD23	2.09	0.82
1:B:329:LEU:HB3	1:B:330:PRO:HD2	1.63	0.81
1:A:352:ARG:HE	1:A:354:ARG:CZ	1.93	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:SER:HB2	1:B:280:GLU:HB2	1.61	0.81
1:B:385:ARG:HH11	1:B:406:ARG:HE	0.84	0.80
1:A:412:MET:HG3	1:A:413:PRO:CD	2.10	0.80
1:B:337:PRO:HB2	1:B:339:ARG:N	1.94	0.80
1:B:342:PRO:C	1:B:344:GLU:H	1.89	0.80
1:B:97[A]:GLU:OE2	1:B:148:ARG:HD3	1.81	0.79
1:B:176:HIS:CE1	1:B:179:VAL:HG21	2.17	0.79
1:A:369:TRP:CZ2	1:A:371:VAL:HG12	2.18	0.79
1:B:337:PRO:CB	1:B:339:ARG:H	1.93	0.79
1:A:226:ARG:NH2	1:A:329:LEU:HD21	1.97	0.78
1:A:330:PRO:HG2	1:A:331:LEU:HD23	1.66	0.78
1:A:385:ARG:NH1	1:A:406:ARG:NH2	2.32	0.77
1:A:223:GLU:HB3	1:A:326:THR:CG2	2.12	0.77
1:A:402:MET:HE2	1:A:420:TRP:CD2	2.20	0.77
1:A:288:GLY:O	1:A:325:PRO:HG2	1.84	0.76
1:B:393:PRO:HG3	1:B:427:ASP:OD1	1.85	0.76
1:A:330:PRO:CG	1:A:331:LEU:HA	2.14	0.76
1:B:342:PRO:O	1:B:344:GLU:N	2.17	0.76
1:A:414:PHE:CE2	1:B:251:PRO:HB3	2.20	0.76
1:A:41:GLU:HB3	1:A:80:ARG:HB3	1.66	0.76
1:B:48:LEU:HB2	1:B:89:LEU:HD12	1.68	0.76
1:A:55:ALA:CB	1:A:67:VAL:N	2.49	0.76
1:A:89:LEU:HD23	1:A:91:TYR:CE1	2.20	0.75
1:B:347:ALA:HB1	1:B:348:GLY:HA2	1.67	0.75
1:B:369:TRP:CH2	1:B:371:VAL:HG12	2.21	0.75
1:B:345:ILE:CG1	1:B:346:MET:HG2	2.13	0.75
1:A:210:ASP:HA	1:A:315:LEU:HD23	1.69	0.75
1:B:172:PRO:HG3	1:B:339:ARG:HH21	1.51	0.75
1:A:42:SER:OG	1:A:43:ARG:N	2.20	0.75
1:B:265:PRO:CD	1:B:340:LEU:HD13	2.14	0.75
1:A:54:ARG:HH21	1:A:54:ARG:HG3	1.53	0.74
1:B:70:ILE:HD12	1:B:141:TYR:HE1	1.52	0.74
1:B:35:PRO:HD3	1:B:217:VAL:HG11	1.69	0.74
1:A:50:MET:CE	1:A:79:ILE:HD11	2.18	0.74
1:B:32:GLN:N	2:B:2001:HOH:O	2.21	0.74
1:B:331:LEU:HD23	1:B:332:VAL:HA	1.69	0.72
1:A:43:ARG:NH1	1:A:46:GLN:HG2	2.05	0.72
1:A:50:MET:HE2	1:A:79:ILE:HD11	1.69	0.72
1:B:228:TRP:HD1	1:B:332:VAL:O	1.72	0.72
1:B:287:ASN:ND2	1:B:289:ASP:HB2	2.04	0.72
1:B:272:SER:HB3	2:B:2026:HOH:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:VAL:HG13	1:A:157:LEU:HD11	1.71	0.71
1:A:347:ALA:H	1:A:404[B]:GLN:HE22	1.35	0.71
1:A:55:ALA:HB2	1:A:67:VAL:C	2.15	0.71
1:B:40:LEU:HG	1:B:48:LEU:HD22	1.71	0.71
1:B:329:LEU:CB	1:B:330:PRO:HD2	2.20	0.71
1:A:43:ARG:HH11	1:A:46:GLN:HG2	1.57	0.70
1:A:228:TRP:NE1	1:A:334:ASP:O	2.25	0.70
1:A:468:PRO:HB2	1:A:469:VAL:HB	1.74	0.70
1:B:89:LEU:HD23	1:B:91:TYR:CE1	2.27	0.69
1:A:385:ARG:HH12	1:A:406:ARG:HH22	1.36	0.69
1:B:47:PRO:HB3	1:B:88:LYS:CG	2.22	0.69
1:A:380:GLN:HB3	1:A:469:VAL:HG23	1.74	0.69
1:B:50:MET:CE	1:B:79:ILE:HD11	2.22	0.69
1:B:219:SER:HB3	1:B:319:LEU:HD23	1.75	0.69
1:B:347:ALA:HB1	1:B:348:GLY:CA	2.22	0.69
1:A:59[A]:PHE:CD1	1:A:213:LEU:HD21	2.28	0.68
1:A:333:THR:HG22	1:A:334:ASP:H	1.57	0.68
1:A:329:LEU:HB3	1:A:330:PRO:CD	2.22	0.68
1:B:117:ALA:HB1	1:B:444:PHE:CE1	2.29	0.68
1:A:246:MET:HE2	1:A:252:LEU:HD13	1.73	0.68
1:A:402:MET:HG3	1:A:420:TRP:CE3	2.29	0.68
1:B:98:ASN:ND2	1:B:119:MET:HE3	2.09	0.68
1:A:59[A]:PHE:CE2	1:A:67:VAL:HG11	2.29	0.68
1:B:42:SER:C	1:B:43:ARG:HD2	2.19	0.68
1:B:228:TRP:HB2	1:B:332:VAL:CG1	2.17	0.68
1:A:352:ARG:NE	1:A:354:ARG:NH2	2.41	0.68
1:A:402:MET:HE2	1:A:420:TRP:CG	2.29	0.67
1:B:196:THR:CG2	1:B:197:PRO:HD2	2.25	0.67
1:B:395:ALA:HB3	1:B:451:GLN:HB2	1.75	0.67
1:B:328:LEU:O	1:B:329:LEU:O	2.12	0.67
1:B:337:PRO:CB	1:B:339:ARG:HB3	2.20	0.67
1:A:59[A]:PHE:HD1	1:A:213:LEU:HD21	1.58	0.67
1:A:406:ARG:HD2	1:A:433:LEU:CD2	2.21	0.67
1:A:59[A]:PHE:HE2	1:A:67:VAL:HG11	1.60	0.67
1:A:343:THR:CG2	1:A:345:ILE:HG23	2.25	0.67
1:B:58:SER:HB3	1:B:63:THR:O	1.94	0.67
1:B:70:ILE:CD1	1:B:141:TYR:HE1	2.08	0.67
1:B:192:ASP:OD1	1:B:194:PHE:HB2	1.94	0.67
1:A:343:THR:HG23	1:A:345:ILE:HG23	1.77	0.67
1:B:336:LEU:CD2	1:B:337:PRO:HD2	2.25	0.67
1:A:154:TYR:O	1:A:189:LYS:NZ	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ALA:N	1:A:404[B]:GLN:HE22	1.93	0.65
1:B:59:PHE:CD1	1:B:213:LEU:HD21	2.31	0.65
1:B:292:SER:OG	1:B:319:LEU:HD11	1.96	0.65
1:A:203:GLY:O	1:A:458:ARG:NH1	2.30	0.65
1:B:154:TYR:O	1:B:189:LYS:NZ	2.29	0.65
1:B:180:ASP:OD2	2:B:2011:HOH:O	2.15	0.65
1:B:192:ASP:OD2	1:B:196:THR:HB	1.96	0.65
1:B:274:ALA:O	1:B:277:GLU:HB2	1.96	0.64
1:B:355:ASP:OD2	1:B:389:ARG:NH1	2.26	0.64
1:A:395:ALA:CB	1:A:451:GLN:HB3	2.21	0.64
1:A:104:ALA:HB3	1:A:140:TRP:CD1	2.33	0.64
1:A:359:GLY:O	1:A:392:GLU:HB3	1.98	0.64
1:B:58:SER:HA	2:B:2002:HOH:O	1.97	0.64
1:A:82:TRP:O	1:A:85:ASP:HB2	1.99	0.63
1:B:385:ARG:NH1	1:B:406:ARG:NE	2.28	0.63
1:A:212:LEU:HD11	1:A:295:CYS:SG	2.39	0.63
1:B:190:ARG:NH1	1:B:210:ASP:OD2	2.30	0.63
1:B:282:LEU:HD12	1:B:336:LEU:HD21	1.81	0.63
1:A:232:ARG:HG2	1:A:280:GLU:HG2	1.79	0.63
1:A:241:ARG:NH2	1:A:272:SER:OG	2.32	0.63
1:A:263:PRO:HA	1:A:341:LEU:HB2	1.81	0.63
1:B:384:GLU:HG3	1:B:436:PHE:HE2	1.62	0.63
1:A:60:THR:HG23	1:A:61:PRO:HD2	1.81	0.63
1:B:172:PRO:HG3	1:B:339:ARG:NH2	2.14	0.63
1:B:47:PRO:HB3	1:B:88:LYS:HG3	1.80	0.62
1:A:172:PRO:HG2	1:A:337:PRO:HG3	1.81	0.62
1:B:113:MET:HB2	1:B:114:GLY:HA2	1.79	0.62
1:A:347:ALA:CB	1:A:433:LEU:HD23	2.30	0.62
1:B:261:PHE:HB2	1:B:341:LEU:CD1	2.29	0.61
1:B:82:TRP:O	1:B:85:ASP:HB2	2.00	0.61
1:A:192:ASP:OD2	1:A:196:THR:HB	2.01	0.61
1:B:327:GLY:HA2	1:B:328:LEU:HD23	1.82	0.61
1:B:468:PRO:HA	1:B:469:VAL:C	2.25	0.61
1:B:336:LEU:HD22	1:B:337:PRO:HD2	1.81	0.61
1:A:60:THR:CG2	1:A:61:PRO:HD2	2.31	0.60
1:A:468:PRO:CB	1:A:469:VAL:HB	2.31	0.60
1:A:35:PRO:HD3	1:A:217:VAL:HG11	1.82	0.60
1:A:208:VAL:HG13	1:A:240:ARG:NH2	2.17	0.60
1:B:341:LEU:HG	1:B:344:GLU:OE2	2.02	0.60
1:B:395:ALA:HB3	1:B:451:GLN:CB	2.32	0.60
1:B:399:GLU:O	1:B:401:VAL:HG13	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HD13	1:A:278:ARG:HG2	1.84	0.60
1:A:266:VAL:CG1	1:A:415:PRO:HB2	2.27	0.60
1:A:333:THR:CG2	1:A:334:ASP:H	2.15	0.60
1:A:347:ALA:HB1	1:A:433:LEU:CD2	2.32	0.60
1:B:63:THR:HG22	2:B:2002:HOH:O	2.02	0.59
1:A:67:VAL:O	1:A:68:TRP:HD1	1.87	0.58
1:B:326:THR:HB	1:B:327:GLY:HA3	1.85	0.58
1:A:226:ARG:HH21	1:A:329:LEU:HD21	1.66	0.58
1:B:234:LEU:HD12	1:B:235:ASN:N	2.18	0.58
1:B:138:THR:O	1:B:278:ARG:NE	2.37	0.58
1:A:333:THR:HG22	1:A:334:ASP:N	2.19	0.57
1:A:190:ARG:HG2	1:A:190:ARG:HH11	1.68	0.57
1:B:332:VAL:C	1:B:334:ASP:H	2.11	0.57
1:B:63:THR:HG23	1:B:194:PHE:O	2.04	0.57
1:B:230:ARG:NH1	1:B:280:GLU:OE1	2.32	0.57
1:A:39:LEU:HD21	1:A:80:ARG:CZ	2.35	0.57
1:B:332:VAL:HG23	1:B:334:ASP:CG	2.30	0.57
1:B:412:MET:HG2	1:B:413:PRO:HD2	1.87	0.57
1:A:274:ALA:O	1:A:277:GLU:HB2	2.05	0.57
1:B:377:THR:CG2	1:B:465:LEU:HD23	2.30	0.57
1:A:48:LEU:CB	1:A:89:LEU:HD12	2.32	0.56
1:A:221:TYR:HE2	1:A:223:GLU:CG	2.18	0.56
1:B:63:THR:HG21	1:B:195:GLY:CA	2.21	0.56
1:A:412:MET:CG	1:A:413:PRO:HD2	2.20	0.56
1:B:196:THR:HG22	1:B:197:PRO:HD2	1.86	0.56
1:B:395:ALA:CB	1:B:451:GLN:HB2	2.35	0.56
1:B:326:THR:HB	1:B:327:GLY:CA	2.36	0.56
1:B:361:ASP:HB3	1:B:362:PRO:HD2	1.88	0.56
1:A:256:SER:HB3	1:A:257:GLY:HA3	1.87	0.56
1:B:345:ILE:HG12	1:B:346:MET:HG3	1.83	0.56
1:B:327:GLY:HA3	1:B:328:LEU:CG	2.33	0.56
1:A:50:MET:HE2	1:A:79:ILE:CD1	2.37	0.55
1:B:385:ARG:HH12	1:B:406:ARG:HH21	1.54	0.55
1:B:104:ALA:HB3	1:B:140:TRP:CD1	2.42	0.55
1:B:334:ASP:O	1:B:335:SER:HB3	2.07	0.55
1:B:359:GLY:N	2:B:2021:HOH:O	2.39	0.55
1:B:190:ARG:HG2	1:B:209:GLY:HA2	1.89	0.55
1:A:398:ILE:HG23	1:A:448:PHE:HB3	1.89	0.55
1:A:82:TRP:HH2	1:A:169:LYS:HE2	1.72	0.54
1:B:414:PHE:HB3	1:B:415:PRO:HD2	1.89	0.54
1:A:399:GLU:O	1:A:401:VAL:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:VAL:HG12	1:B:268:VAL:HG13	1.90	0.54
1:B:41:GLU:HG3	1:B:80:ARG:HB3	1.90	0.54
1:A:208:VAL:HG13	1:A:240:ARG:HH21	1.73	0.54
1:A:330:PRO:CB	1:A:331:LEU:HA	2.37	0.54
1:A:98:ASN:ND2	1:A:121:SER:HA	2.23	0.54
1:A:192:ASP:OD1	1:A:194:PHE:N	2.41	0.54
1:B:60:THR:O	1:B:63:THR:HB	2.08	0.54
1:B:361:ASP:HB3	1:B:362:PRO:CD	2.37	0.54
1:A:468:PRO:HA	1:A:469:VAL:HG23	1.89	0.53
1:B:160:MET:SD	1:B:183:PRO:HB3	2.48	0.53
1:B:52:VAL:HG13	1:B:157:LEU:HD21	1.89	0.53
1:B:69:GLY:HA3	1:B:73:ARG:O	2.08	0.53
1:A:162:LEU:HD22	1:A:232:ARG:NH2	2.23	0.53
1:A:385:ARG:NH1	1:A:406:ARG:HH21	2.05	0.53
1:A:74:TYR:O	1:A:75:LEU:HB2	2.09	0.53
1:B:405:ILE:HD11	1:B:424:VAL:HG23	1.90	0.53
1:A:75:LEU:HA	1:A:156:GLY:O	2.09	0.53
1:A:228:TRP:CH2	1:A:336:LEU:HD21	2.45	0.53
1:B:47:PRO:HB3	1:B:88:LYS:HG2	1.90	0.53
1:A:330:PRO:CG	1:A:331:LEU:HD23	2.37	0.52
1:B:228:TRP:CD1	1:B:332:VAL:O	2.59	0.52
1:A:59[A]:PHE:HE2	1:A:67:VAL:CG1	2.21	0.52
1:B:337:PRO:C	1:B:339:ARG:N	2.63	0.52
1:A:347:ALA:CB	1:A:404[B]:GLN:NE2	2.64	0.52
1:B:341:LEU:HD23	1:B:344:GLU:HG3	1.91	0.52
1:B:360:ASP:OD1	1:B:360:ASP:N	2.37	0.52
1:A:402:MET:HE2	1:A:420:TRP:CE2	2.44	0.52
1:A:344:GLU:HG2	1:A:346:MET:HG2	1.91	0.52
1:B:59:PHE:HD1	1:B:213:LEU:HD21	1.70	0.51
1:B:154:TYR:CE2	1:B:199:TYR:HB2	2.44	0.51
1:A:263:PRO:HG3	1:A:344:GLU:HG3	1.92	0.51
1:B:113:MET:HB3	2:B:2005:HOH:O	2.10	0.51
1:B:174:PRO:O	1:B:230:ARG:HD2	2.11	0.51
1:B:384:GLU:HG3	1:B:436:PHE:CE2	2.44	0.51
1:A:136:ALA:HA	1:A:163:VAL:O	2.11	0.51
1:B:113:MET:CB	1:B:114:GLY:HA2	2.39	0.51
1:A:52:VAL:CG1	1:A:95:LEU:HG	2.37	0.51
1:A:70:ILE:HD12	1:A:141:TYR:HE1	1.75	0.51
1:B:383:TRP:HH2	1:B:420:TRP:HH2	1.57	0.51
1:A:55:ALA:HB1	1:A:67:VAL:N	2.10	0.51
1:A:456:ALA:C	1:A:458:ARG:H	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:ARG:NE	1:B:431:GLU:OE2	2.43	0.50
1:A:347:ALA:HB1	1:A:433:LEU:HD23	1.93	0.50
1:B:347:ALA:CB	1:B:348:GLY:CA	2.90	0.50
1:B:354:ARG:CZ	1:B:376:VAL:HG13	2.41	0.50
1:B:32:GLN:HB3	1:B:33:PRO:CD	2.41	0.50
1:B:383:TRP:CH2	1:B:420:TRP:HH2	2.29	0.50
1:A:35:PRO:HD3	1:A:217:VAL:CG1	2.42	0.50
1:B:52:VAL:HG12	1:B:95:LEU:CD1	2.41	0.50
1:A:407:ASN:HB2	1:A:431:GLU:HB2	1.93	0.50
1:B:176:HIS:ND1	1:B:179:VAL:HG21	2.27	0.50
1:A:228:TRP:CD1	1:A:334:ASP:O	2.65	0.50
1:B:39:LEU:HD21	1:B:80:ARG:CZ	2.41	0.50
1:B:245:GLN:HA	1:B:252:LEU:HD11	1.94	0.49
1:B:98:ASN:CG	1:B:119:MET:HE3	2.37	0.49
1:B:196:THR:HG23	1:B:197:PRO:HD2	1.94	0.49
1:B:74:TYR:O	1:B:75:LEU:HB2	2.12	0.49
1:B:385:ARG:NH1	1:B:406:ARG:HH21	2.11	0.49
1:A:278:ARG:NH2	1:A:422:ASP:OD2	2.46	0.49
1:B:189:LYS:HG2	1:B:191:LEU:HD21	1.93	0.49
1:A:74:TYR:CE1	1:A:156:GLY:HA3	2.48	0.49
1:B:210:ASP:HA	1:B:315:LEU:CD2	2.34	0.49
1:B:414:PHE:O	1:B:418:ARG:HG3	2.13	0.49
1:B:217:VAL:HG13	1:B:220:PRO:HG3	1.94	0.49
1:A:138:THR:O	1:A:278:ARG:NE	2.46	0.49
1:A:244:LEU:HA	1:A:294:THR:O	2.12	0.49
1:A:332:VAL:HG13	1:A:332:VAL:O	2.13	0.49
1:A:385:ARG:NH1	1:A:406:ARG:HH22	2.05	0.49
1:A:345:ILE:O	1:A:345:ILE:HG13	2.12	0.49
1:B:208:VAL:HG13	1:B:240:ARG:NH2	2.28	0.49
1:B:332:VAL:C	1:B:334:ASP:N	2.71	0.48
1:B:47:PRO:CG	1:B:90:ILE:HD12	2.42	0.48
1:A:344:GLU:HG2	1:A:344:GLU:O	2.13	0.48
1:A:358:LEU:HA	2:A:2023:HOH:O	2.12	0.48
1:B:329:LEU:CB	1:B:330:PRO:CD	2.80	0.48
1:B:337:PRO:C	1:B:339:ARG:H	2.16	0.48
1:B:415:PRO:O	1:B:418:ARG:HB2	2.13	0.48
1:A:203:GLY:HA2	1:A:458:ARG:HH22	1.79	0.48
1:A:414:PHE:HB3	1:A:415:PRO:HD2	1.94	0.48
1:A:109:PRO:HD2	1:A:112:LEU:HD12	1.95	0.48
1:A:65:ALA:HB1	1:A:196:THR:HG23	1.96	0.48
1:A:269:LYS:HG2	1:B:414:PHE:CZ	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:GLN:O	1:B:295:CYS:HA	2.14	0.48
1:B:384:GLU:CG	1:B:436:PHE:HE2	2.27	0.48
1:A:46:GLN:O	1:A:46:GLN:CG	2.53	0.47
1:A:82:TRP:HH2	1:A:169:LYS:CE	2.27	0.47
1:A:160:MET:SD	1:A:183:PRO:HB3	2.54	0.47
1:B:342:PRO:C	1:B:344:GLU:N	2.56	0.47
1:B:414:PHE:HB3	1:B:416:GLU:OE1	2.14	0.47
1:B:60:THR:HG23	1:B:63:THR:HB	1.97	0.47
1:A:420:TRP:N	2:A:2032:HOH:O	2.48	0.47
1:A:208:VAL:CG1	1:A:240:ARG:NH2	2.77	0.47
1:B:331:LEU:HD23	1:B:331:LEU:HA	1.33	0.47
1:B:176:HIS:N	2:B:2010:HOH:O	2.48	0.47
1:B:192:ASP:OD1	1:B:194:PHE:N	2.46	0.47
1:A:314:ILE:HG23	1:A:315:LEU:N	2.29	0.47
1:A:388:VAL:O	1:A:429:GLN:HA	2.15	0.47
1:B:257:GLY:HA3	1:B:260:GLY:O	2.14	0.47
1:A:51:THR:CG2	1:A:94:ARG:HG3	2.44	0.47
1:A:97:GLU:OE2	1:A:148:ARG:HD3	2.15	0.47
1:A:250:ARG:HG3	1:A:251:PRO:O	2.15	0.47
1:A:269:LYS:HG2	1:B:414:PHE:CE2	2.50	0.47
1:A:279:ARG:HG2	2:A:2019:HOH:O	2.14	0.47
1:A:154:TYR:CE2	1:A:199:TYR:HB2	2.50	0.46
1:A:355:ASP:OD2	1:A:389:ARG:NH2	2.35	0.46
1:B:469:VAL:HG12	1:B:469:VAL:O	2.13	0.46
1:A:253:HIS:CG	1:A:336:LEU:HD11	2.50	0.46
1:A:383:TRP:HH2	1:A:420:TRP:HH2	1.62	0.46
1:A:68:TRP:HB2	1:A:155:ASN:HB3	1.96	0.46
1:B:42:SER:O	1:B:43:ARG:HD2	2.15	0.46
1:B:336:LEU:HD23	1:B:337:PRO:HD2	1.97	0.46
1:B:337:PRO:HB3	1:B:339:ARG:CB	2.30	0.46
1:A:368:LEU:HD22	1:A:459:GLY:HA3	1.98	0.46
1:A:415:PRO:N	1:A:416:GLU:HB3	2.31	0.46
1:B:398:ILE:HG23	1:B:448:PHE:HB3	1.97	0.46
1:A:98:ASN:HB3	1:A:119:MET:HE3	1.97	0.46
1:A:414:PHE:CD2	1:B:251:PRO:HB3	2.50	0.46
1:B:208:VAL:CG1	1:B:240:ARG:NH2	2.79	0.46
1:B:110:GLY:N	1:B:111:PRO:CD	2.78	0.46
1:B:252:LEU:N	1:B:268:VAL:O	2.38	0.46
1:B:210:ASP:CA	1:B:315:LEU:HD23	2.34	0.45
1:A:78:THR:HG23	1:A:160:MET:HB3	1.98	0.45
1:A:392:GLU:HA	1:A:393:PRO:HD3	1.86	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:LEU:HA	1:B:156:GLY:O	2.15	0.45
1:B:316:VAL:HG23	1:B:317:SER:N	2.30	0.45
1:A:54:ARG:HH21	1:A:54:ARG:CG	2.27	0.45
1:A:138:THR:OG1	1:A:280:GLU:OE2	2.32	0.45
1:B:122:PRO:O	1:B:123:ASN:HB2	2.16	0.45
1:B:188:ASP:OD2	1:B:240:ARG:HB2	2.16	0.45
1:A:89:LEU:O	1:A:128:PRO:HD2	2.16	0.45
1:B:336:LEU:HA	1:B:337:PRO:HD2	1.27	0.45
1:B:352:ARG:HD3	1:B:354:ARG:NH2	2.31	0.45
1:B:332:VAL:HB	1:B:334:ASP:OD1	2.17	0.45
1:B:385:ARG:NH1	1:B:406:ARG:CZ	2.80	0.45
1:B:409:ASN:C	1:B:411:ALA:H	2.23	0.45
1:A:80:ARG:HD2	1:A:164:GLU:HG3	1.99	0.45
1:A:265:PRO:HD3	1:A:340:LEU:HG	1.99	0.45
1:A:63:THR:HB	1:A:194:PHE:O	2.16	0.45
1:A:182:PHE:HA	1:A:183:PRO:HD3	1.69	0.45
1:A:269:LYS:CG	1:B:414:PHE:CE2	2.99	0.45
1:B:212:LEU:HD11	1:B:295:CYS:SG	2.57	0.45
1:B:265:PRO:HG3	1:B:340:LEU:HD11	1.97	0.45
1:B:248:ASP:OD1	1:B:250:ARG:HG2	2.17	0.45
1:B:456:ALA:C	1:B:458:ARG:N	2.75	0.45
1:A:201:GLU:CD	1:A:201:GLU:H	2.25	0.45
1:A:402:MET:CE	1:A:420:TRP:CG	3.00	0.45
1:B:32:GLN:HB3	1:B:33:PRO:HD2	1.99	0.45
1:B:336:LEU:HD23	1:B:336:LEU:HA	1.71	0.44
1:B:414:PHE:HB3	1:B:415:PRO:CD	2.46	0.44
1:A:82:TRP:CZ3	1:A:164:GLU:HB3	2.52	0.44
1:B:93:ASN:HD22	1:B:121:SER:C	2.26	0.44
1:B:139:LEU:N	1:B:139:LEU:CD1	2.80	0.44
1:A:110:GLY:N	1:A:111:PRO:CD	2.79	0.44
1:B:440:SER:HB2	1:B:445:PRO:HA	1.99	0.44
1:A:228:TRP:CZ3	1:A:336:LEU:HD21	2.53	0.44
1:A:259:GLN:HB2	1:A:420:TRP:HB3	1.99	0.44
1:A:292:SER:HB3	1:A:319:LEU:CD1	2.48	0.44
1:A:292:SER:HB3	1:A:319:LEU:HD11	1.98	0.44
1:A:369:TRP:CH2	1:A:371:VAL:HG12	2.51	0.44
1:B:332:VAL:HG22	1:B:332:VAL:O	2.17	0.44
1:A:230:ARG:NH1	1:A:280:GLU:OE1	2.39	0.44
1:B:191:LEU:HD23	1:B:197:PRO:HA	1.98	0.44
1:A:202:PRO:HG2	1:A:203:GLY:H	1.83	0.44
1:B:221:TYR:CD2	1:B:221:TYR:C	2.96	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:LEU:O	1:B:341:LEU:C	2.61	0.44
1:B:456:ALA:C	1:B:458:ARG:H	2.25	0.44
1:B:38:PRO:HG2	1:B:77:PRO:HB3	2.00	0.43
1:A:415:PRO:HB2	1:A:416:GLU:HB2	2.00	0.43
1:B:341:LEU:HD23	1:B:344:GLU:CG	2.48	0.43
1:A:314:ILE:CG2	1:A:315:LEU:N	2.79	0.43
1:B:337:PRO:CB	1:B:339:ARG:CB	2.93	0.43
1:A:408:VAL:O	1:A:409:ASN:HB2	2.18	0.43
1:B:63:THR:CG2	1:B:64:ARG:N	2.81	0.43
1:A:456:ALA:C	1:A:458:ARG:N	2.77	0.43
1:B:88:LYS:HD3	1:B:129:VAL:HG22	2.00	0.43
1:A:34:LEU:HA	1:A:35:PRO:HD3	1.89	0.43
1:A:208:VAL:CG1	1:A:240:ARG:HH21	2.32	0.43
1:B:35:PRO:CD	1:B:217:VAL:CG1	2.96	0.43
1:B:60:THR:HG23	1:B:63:THR:CB	2.49	0.43
1:B:385:ARG:NH1	1:B:406:ARG:NH2	2.66	0.43
1:B:342:PRO:O	1:B:344:GLU:HG2	2.19	0.43
1:A:226:ARG:NH1	1:A:286:SER:O	2.47	0.43
1:B:331:LEU:CD2	1:B:332:VAL:HA	2.45	0.43
1:A:226:ARG:CZ	1:A:329:LEU:HD11	2.49	0.42
1:A:93:ASN:OD1	1:A:95:LEU:HB2	2.20	0.42
1:B:421:LYS:HD3	2:B:2027:HOH:O	2.19	0.42
1:A:69:GLY:HA3	1:A:73:ARG:O	2.20	0.42
1:A:98:ASN:HD22	1:A:122:PRO:HD3	1.84	0.42
1:B:408:VAL:O	1:B:409:ASN:HB2	2.19	0.42
1:A:217:VAL:HG13	1:A:220:PRO:HG3	2.01	0.42
1:A:221:TYR:HE2	1:A:223:GLU:HG3	1.84	0.42
1:A:82:TRP:CH2	1:A:169:LYS:CE	3.02	0.42
1:B:98:ASN:OD1	1:B:121:SER:HA	2.20	0.42
1:B:145:THR:HA	1:B:146:PRO:HD3	1.88	0.42
1:B:406:ARG:HD2	1:B:433:LEU:CD2	2.33	0.42
1:A:116:PRO:HG2	1:A:447:TYR:OH	2.20	0.42
1:A:402:MET:HE2	1:A:420:TRP:CD1	2.54	0.42
1:A:403:PHE:C	1:A:403:PHE:CD2	2.98	0.42
1:A:468:PRO:CA	1:A:469:VAL:HB	2.50	0.42
1:B:224:VAL:HG21	1:B:285:MET:HG3	2.00	0.42
1:A:414:PHE:HB3	1:A:416:GLU:OE1	2.20	0.42
1:A:37:PRO:O	1:A:177:TYR:OH	2.30	0.42
1:A:68:TRP:CH2	1:A:95:LEU:HD13	2.54	0.42
1:B:274:ALA:HB1	1:B:275:PRO:HD2	2.01	0.42
1:B:182:PHE:HA	1:B:183:PRO:HD3	1.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:O	1:B:68:TRP:HD1	2.03	0.42
1:A:139:LEU:N	1:A:139:LEU:CD1	2.83	0.41
1:A:333:THR:CG2	1:A:334:ASP:N	2.81	0.41
1:B:78:THR:HG23	1:B:160:MET:HB3	2.02	0.41
1:B:152:GLN:O	1:B:157:LEU:HB2	2.20	0.41
1:A:330:PRO:HD2	1:A:331:LEU:HD23	2.02	0.41
1:B:224:VAL:N	1:B:324:ARG:O	2.46	0.41
1:A:217:VAL:HG13	1:A:220:PRO:HB3	2.02	0.41
1:B:246:MET:HG3	1:B:250:ARG:HG3	2.02	0.41
1:A:217:VAL:CG1	1:A:220:PRO:HB3	2.50	0.41
1:B:334:ASP:OD2	1:B:335:SER:N	2.52	0.41
1:A:89:LEU:O	1:A:127:ALA:HA	2.20	0.41
1:B:224:VAL:HG11	1:B:229:VAL:HG11	2.03	0.41
1:B:265:PRO:HG3	1:B:340:LEU:CD1	2.50	0.41
1:A:369:TRP:CH2	1:A:371:VAL:CG1	3.04	0.41
1:B:148:ARG:O	1:B:152:GLN:HG3	2.19	0.41
1:A:196:THR:HA	1:A:197:PRO:HD3	1.90	0.41
1:A:224:VAL:O	1:A:326:THR:HG23	2.21	0.41
1:A:256:SER:CB	1:A:257:GLY:HA3	2.45	0.41
1:A:82:TRP:CE3	1:A:164:GLU:HB2	2.56	0.41
1:A:443:HIS:CD2	1:A:444:PHE:CD2	3.09	0.41
1:B:50:MET:HE2	1:B:79:ILE:HD11	2.00	0.41
1:B:331:LEU:HA	1:B:332:VAL:HA	1.60	0.41
1:B:65:ALA:HB1	1:B:196:THR:HG23	2.02	0.41
1:B:380:GLN:NE2	1:B:439:PRO:HD3	2.35	0.41
1:A:38:PRO:HG2	1:A:77:PRO:HB3	2.03	0.40
1:B:59:PHE:HZ	1:B:67:VAL:HG11	1.86	0.40
1:B:278:ARG:NH2	1:B:422:ASP:OD2	2.50	0.40
1:B:256:SER:HA	1:B:257:GLY:HA3	1.92	0.40
1:B:405:ILE:HD11	1:B:424:VAL:CG2	2.50	0.40
1:B:414:PHE:O	1:B:415:PRO:C	2.64	0.40
1:A:60:THR:CG2	1:A:61:PRO:CD	2.99	0.40
1:A:257:GLY:HA3	1:A:260:GLY:O	2.21	0.40
1:B:341:LEU:O	1:B:342:PRO:C	2.61	0.40
1:B:369:TRP:CH2	1:B:371:VAL:CG1	3.00	0.40
1:A:190:ARG:HG2	1:A:190:ARG:NH1	2.36	0.40
1:A:253:HIS:O	1:A:281:ILE:HA	2.22	0.40
1:A:383:TRP:CH2	1:A:420:TRP:HH2	2.39	0.40
1:B:265:PRO:HD2	1:B:338:MET:HA	2.03	0.40
1:B:47:PRO:HG3	1:B:90:ILE:CD1	2.51	0.40
1:B:399:GLU:HB2	1:B:447:TYR:H	1.87	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	419/451 (93%)	361 (86%)	51 (12%)	7 (2%)	7	14
1	B	419/451 (93%)	370 (88%)	42 (10%)	7 (2%)	7	14
All	All	838/902 (93%)	731 (87%)	93 (11%)	14 (2%)	7	14

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	330	PRO
1	A	416	GLU
1	B	329	LEU
1	B	330	PRO
1	B	331	LEU
1	B	343	THR
1	B	346	MET
1	A	43	ARG
1	A	44	ARG
1	A	47	PRO
1	B	337	PRO
1	A	202	PRO
1	B	47	PRO

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	364/386 (94%)	347 (95%)	17 (5%)	23	46
1	B	363/386 (94%)	334 (92%)	29 (8%)	11	23
All	All	727/772 (94%)	681 (94%)	46 (6%)	16	34

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ARG
1	A	54	ARG
1	A	95	LEU
1	A	119	MET
1	A	139	LEU
1	A	166	GLU
1	A	243	GLN
1	A	245	GLN
1	A	289	ASP
1	A	314	ILE
1	A	358	LEU
1	A	361	ASP
1	A	401	VAL
1	A	412	MET
1	A	433	LEU
1	A	464	LEU
1	A	469	VAL
1	B	43	ARG
1	B	44	ARG
1	B	58	SER
1	B	121	SER
1	B	139	LEU
1	B	217	VAL
1	B	233	LEU
1	B	234	LEU
1	B	289	ASP
1	B	291	VAL
1	B	292	SER
1	B	295	CYS
1	B	317	SER
1	B	330	PRO
1	B	332	VAL
1	B	336	LEU
1	B	341	LEU
1	B	343	THR

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Mol	Chain	Res	Type
1	B	345	ILE
1	B	346	MET
1	B	360	ASP
1	B	371	VAL
1	B	377	THR
1	B	401	VAL
1	B	416	GLU
1	B	433	LEU
1	B	450	SER
1	B	451	GLN
1	B	464	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	151	GLN
1	A	187	GLN
1	A	243	GLN
1	A	467	ASN
1	B	107	GLN
1	B	151	GLN
1	B	176	HIS
1	B	243	GLN
1	B	409	ASN
1	B	449	ASN
1	B	451	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	422/451 (93%)	0.62	45 (10%) 11 9	14, 30, 59, 79	3 (0%)
1	B	422/451 (93%)	0.62	48 (11%) 10 7	13, 29, 63, 79	1 (0%)
All	All	844/902 (93%)	0.62	93 (11%) 10 8	13, 30, 62, 79	4 (0%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	332	VAL	6.5
1	B	166	GLU	4.9
1	B	327	GLY	4.9
1	B	345	ILE	4.6
1	A	332	VAL	4.5
1	B	331	LEU	4.5
1	A	44	ARG	4.2
1	B	29	GLY	4.2
1	B	333	THR	3.9
1	B	194	PHE	3.9
1	B	43	ARG	3.8
1	B	342	PRO	3.7
1	B	469	VAL	3.7
1	B	336	LEU	3.7
1	B	341	LEU	3.7
1	B	330	PRO	3.7
1	B	326	THR	3.6
1	B	343	THR	3.4
1	A	192	ASP	3.4
1	A	59[A]	PHE	3.4
1	B	44	ARG	3.4
1	A	329	LEU	3.3
1	B	347	ALA	3.3
1	A	333	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	334	ASP	3.3
1	B	338	MET	3.3
1	B	325	PRO	3.2
1	B	344	GLU	3.2
1	A	338	MET	3.2
1	A	288	GLY	3.2
1	A	359	GLY	3.2
1	A	469	VAL	3.2
1	A	194	PHE	3.2
1	B	346	MET	3.1
1	A	345	ILE	3.1
1	B	328	LEU	3.1
1	A	344	GLU	3.1
1	A	114	GLY	3.1
1	A	328	LEU	3.0
1	B	339	ARG	3.0
1	A	45	GLY	2.9
1	B	337	PRO	2.9
1	A	331	LEU	2.8
1	A	346	MET	2.8
1	B	195	GLY	2.8
1	B	335	SER	2.8
1	B	340	LEU	2.8
1	B	359	GLY	2.8
1	B	192	ASP	2.8
1	B	201	GLU	2.7
1	B	329	LEU	2.7
1	A	66	SER	2.7
1	B	204	SER	2.7
1	A	166	GLU	2.7
1	B	42	SER	2.7
1	A	314	ILE	2.6
1	A	295	CYS	2.6
1	A	203	GLY	2.6
1	A	334	ASP	2.6
1	B	62	GLY	2.6
1	A	169	LYS	2.6
1	A	286	SER	2.5
1	A	325	PRO	2.5
1	A	287	ASN	2.5
1	B	315	LEU	2.5
1	B	56	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	197	PRO	2.5
1	B	30	GLN	2.4
1	B	205	GLY	2.4
1	A	202	PRO	2.3
1	B	379	GLN	2.3
1	B	316	VAL	2.3
1	A	416	GLU	2.3
1	A	327	GLY	2.3
1	A	322	THR	2.2
1	A	343	THR	2.2
1	A	198	GLU	2.2
1	A	85	ASP	2.2
1	A	335	SER	2.2
1	A	60	THR	2.2
1	B	60	THR	2.2
1	B	348	GLY	2.1
1	A	347	ALA	2.1
1	B	318	THR	2.1
1	A	40	LEU	2.1
1	B	295	CYS	2.1
1	A	339	ARG	2.1
1	B	193	ASN	2.1
1	A	204	SER	2.0
1	A	201	GLU	2.0
1	B	188	ASP	2.0
1	A	42	SER	2.0
1	A	249	GLY	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.