



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:24 AM UTC

PDB ID : 2ACI / pdb_00002aci
Title : Structure of D166A arginine deiminase
Authors : Galkin, A.; Lu, X.; Dunaway-Mariano, D.; Herzberg, O.
Deposited on : 2005-07-18
Resolution : 2.50 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

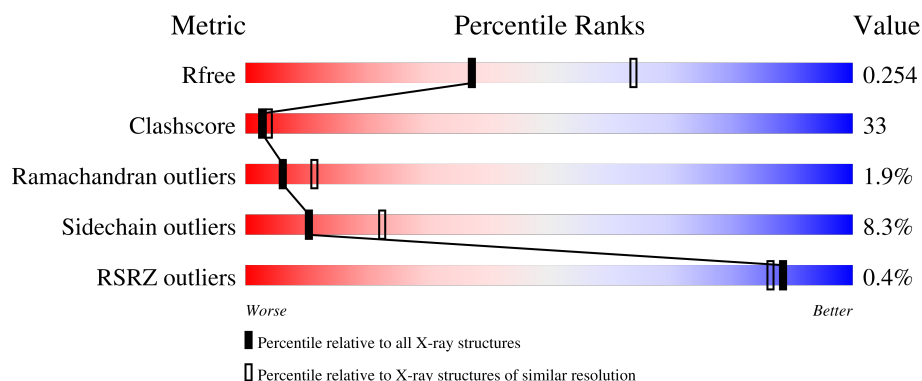
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

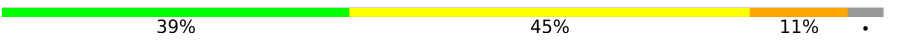
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	418	
1	B	418	
1	C	418	
1	D	418	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12953 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 Arginine deiminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	0	0
			3139	1988	546	588	17			
1	B	409	Total	C	N	O	S	0	0	0
			3193	2020	556	600	17			
1	C	403	Total	C	N	O	S	0	0	0
			3147	1992	547	591	17			
1	D	406	Total	C	N	O	S	0	0	0
			3174	2009	553	595	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	ALA	ASP	engineered mutation	UNP P13981
B	166	ALA	ASP	engineered mutation	UNP P13981
C	166	ALA	ASP	engineered mutation	UNP P13981
D	166	ALA	ASP	engineered mutation	UNP P13981

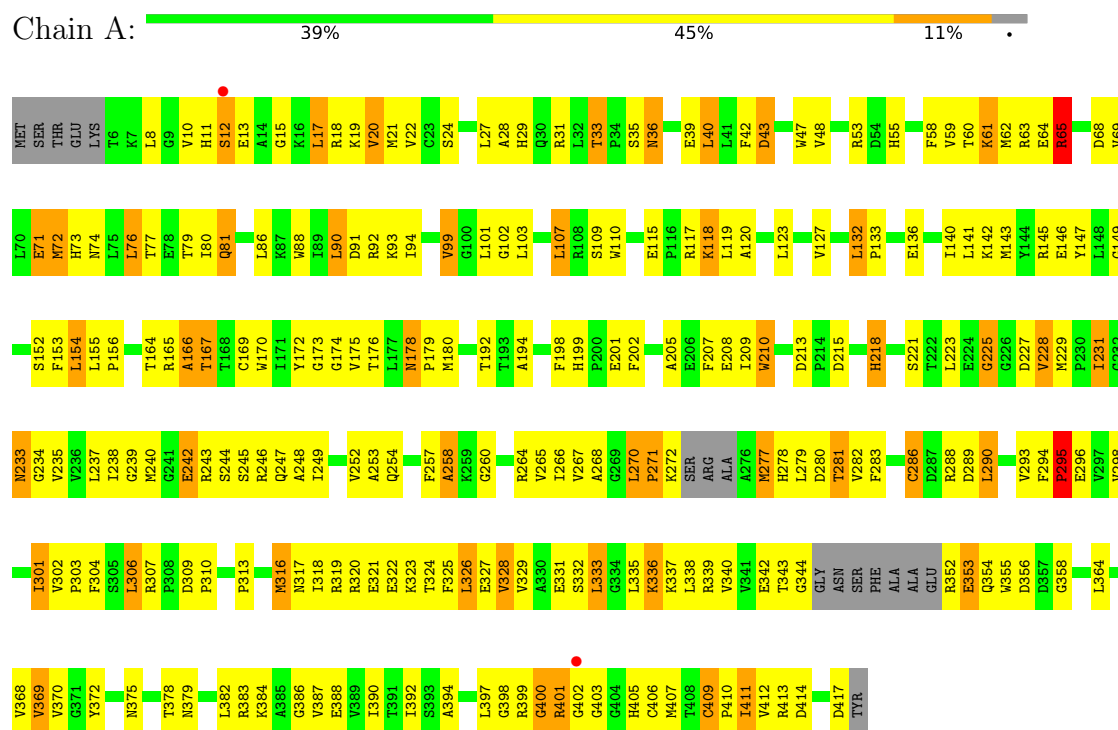
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	60	Total	O	0	0
			60	60		
2	B	90	Total	O	0	0
			90	90		
2	C	86	Total	O	0	0
			86	86		
2	D	64	Total	O	0	0
			64	64		

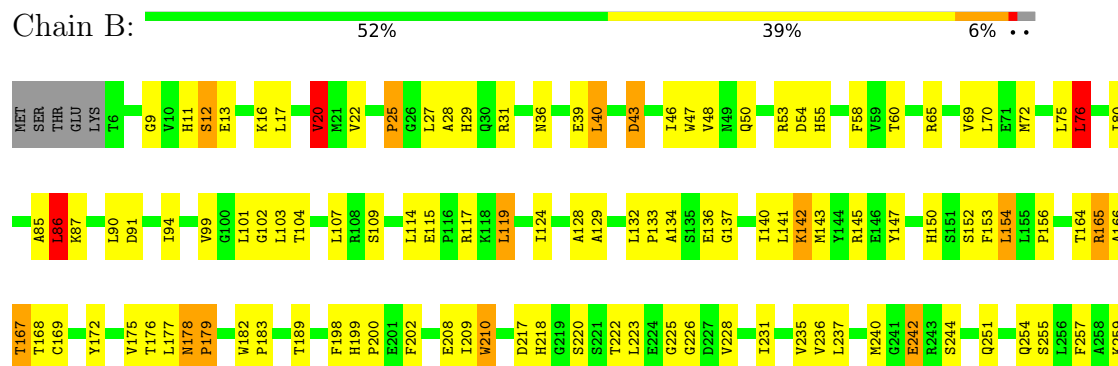
3 Residue-property plots

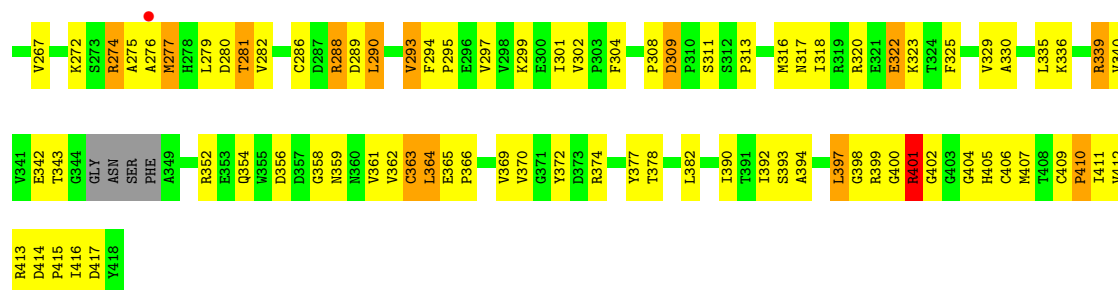
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



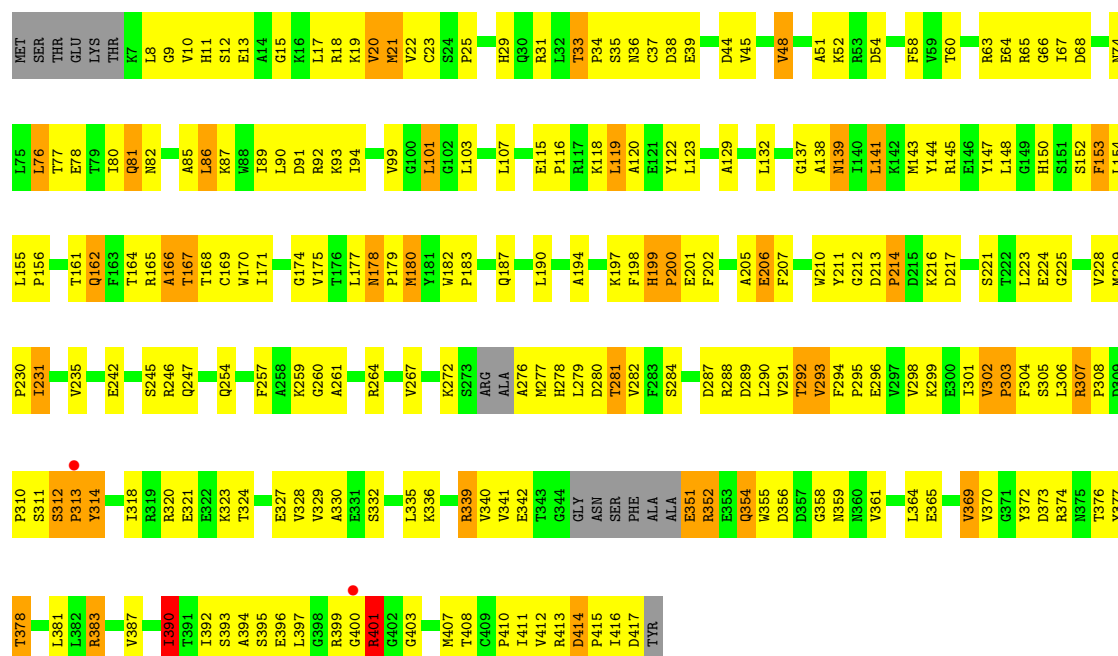
• Molecule 1: Arginine deiminase





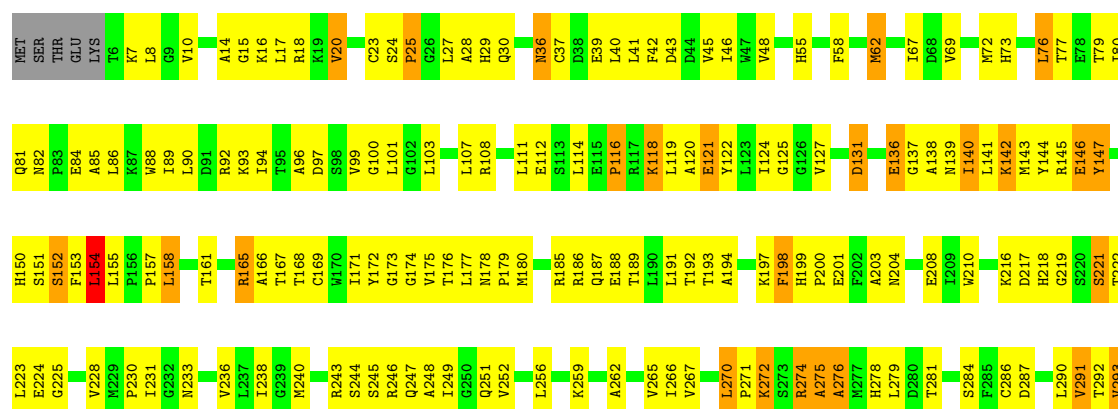
• Molecule 1: Arginine deiminase

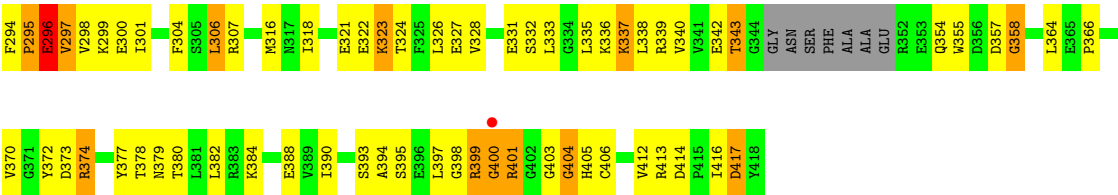
Chain C: 41% 45% 9% .



• Molecule 1: Arginine deiminase

Chain D: 41% 47% 9% .





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.20Å 123.90Å 150.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.50) 93.5 (20.00-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 2.88Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.198 , 0.272 0.182 , 0.254	Depositor DCC
R_{free} test set	1806 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12953	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.68% 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	0.88	3/3205 (0.1%)	1.30	35/4346 (0.8%)
1	B	0.92	0/3261	1.36	44/4422 (1.0%)
1	C	0.90	4/3213 (0.1%)	1.37	36/4356 (0.8%)
1	D	0.89	0/3242	1.35	39/4396 (0.9%)
All	All	0.90	7/12921 (0.1%)	1.34	154/17520 (0.9%)

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	C	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	303	PRO	N-CD	10.60	1.62	1.47
1	C	156	PRO	CA-C	-6.16	1.48	1.52
1	C	354	GLN	C-N	6.01	1.41	1.33
1	C	214	PRO	N-CD	-5.66	1.39	1.47
1	A	295	PRO	C-N	-5.46	1.26	1.33
1	A	411	ILE	CA-CB	-5.13	1.47	1.54
1	A	167	THR	CA-CB	-5.11	1.45	1.53

All (154) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	329	VAL	N-CA-C	-11.75	102.39	111.90
1	C	48	VAL	N-CA-C	11.19	121.16	110.42
1	A	178	ASN	N-CA-C	10.36	122.31	109.57
1	D	48	VAL	N-CA-C	10.19	120.83	110.23
1	C	354	GLN	OE1-CD-NE2	-10.13	112.47	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	302	VAL	CA-C-N	10.08	131.52	120.13
1	C	302	VAL	C-N-CA	10.08	131.52	120.13
1	D	178	ASN	N-CA-C	10.07	121.95	109.57
1	D	154	LEU	N-CA-C	-10.05	100.43	112.89
1	C	414	ASP	CA-C-N	9.81	131.10	120.11
1	C	414	ASP	C-N-CA	9.81	131.10	120.11
1	B	114	LEU	N-CA-C	9.65	122.28	110.41
1	C	166	ALA	N-CA-C	9.20	120.91	111.07
1	A	48	VAL	N-CA-C	8.89	119.69	110.36
1	D	136	GLU	N-CA-C	-8.78	101.17	112.23
1	C	178	ASN	N-CA-C	8.75	122.37	109.50
1	D	142	LYS	N-CA-C	-8.69	102.12	112.89
1	D	178	ASN	CA-C-N	8.66	130.67	119.84
1	D	178	ASN	C-N-CA	8.66	130.67	119.84
1	D	364	LEU	N-CA-C	-8.57	103.34	113.88
1	A	166	ALA	N-CA-C	8.46	120.58	111.36
1	A	65	ARG	N-CA-C	-8.44	102.54	112.92
1	B	166	ALA	N-CA-C	8.40	121.19	111.11
1	A	231	ILE	N-CA-C	8.13	119.93	112.29
1	D	417	ASP	N-CA-C	-7.91	103.58	113.23
1	A	277	MET	N-CA-C	7.90	120.91	109.14
1	B	142	LYS	N-CA-C	-7.86	102.72	111.28
1	A	99	VAL	N-CA-C	7.79	118.69	111.45
1	D	166	ALA	N-CA-C	7.74	120.46	111.02
1	B	294	PHE	CA-C-N	7.71	129.47	119.84
1	B	294	PHE	C-N-CA	7.71	129.47	119.84
1	C	167	THR	N-CA-C	7.54	120.97	111.69
1	B	228	VAL	N-CA-C	7.50	119.09	108.36
1	A	210	TRP	N-CA-C	7.50	120.40	111.33
1	B	154	LEU	N-CA-C	-7.45	102.31	111.40
1	A	228	VAL	N-CA-C	7.41	118.67	108.89
1	C	281	THR	N-CA-C	-7.38	104.43	113.50
1	B	210	TRP	N-CA-C	7.32	119.26	111.28
1	D	399	ARG	N-CA-C	-7.29	103.49	112.38
1	B	156	PRO	CA-C-N	7.17	127.09	119.85
1	B	156	PRO	C-N-CA	7.17	127.09	119.85
1	B	399	ARG	N-CA-C	-7.14	103.49	113.20
1	C	354	GLN	CG-CD-NE2	7.14	127.11	116.40
1	C	340	VAL	N-CA-C	7.12	118.08	108.11
1	B	199	HIS	CA-C-N	7.08	128.69	119.84
1	B	199	HIS	C-N-CA	7.08	128.69	119.84
1	B	48	VAL	N-CA-C	7.06	118.67	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	153	PHE	N-CA-C	7.03	120.18	108.73
1	D	374	ARG	N-CA-C	6.98	119.73	111.71
1	C	153	PHE	N-CA-C	6.93	119.72	108.63
1	B	167	THR	N-CA-C	6.83	118.52	111.14
1	A	72	MET	N-CA-C	6.83	118.72	111.28
1	A	153	PHE	N-CA-C	6.77	119.97	107.99
1	C	199	HIS	CA-C-N	6.74	128.26	119.84
1	C	199	HIS	C-N-CA	6.74	128.26	119.84
1	B	323	LYS	N-CA-C	-6.70	101.69	110.53
1	C	81	GLN	N-CA-C	-6.69	105.12	113.28
1	D	228	VAL	N-CA-C	6.63	118.95	108.87
1	D	167	THR	N-CA-C	6.59	120.92	113.01
1	D	158	LEU	CA-C-N	6.52	126.21	119.56
1	D	158	LEU	C-N-CA	6.52	126.21	119.56
1	D	275	ALA	N-CA-C	6.52	118.90	108.41
1	A	270	LEU	CA-C-N	6.51	127.98	119.84
1	A	270	LEU	C-N-CA	6.51	127.98	119.84
1	A	167	THR	N-CA-C	6.47	121.38	112.90
1	B	340	VAL	N-CA-C	6.43	117.39	108.12
1	C	182	TRP	CA-C-N	6.43	125.92	119.24
1	C	182	TRP	C-N-CA	6.43	125.92	119.24
1	D	25	PRO	N-CA-C	-6.41	101.18	111.38
1	D	198	PHE	N-CA-C	6.41	121.21	112.68
1	C	231	ILE	N-CA-C	6.39	117.17	110.72
1	C	310	PRO	N-CA-C	-6.36	106.40	114.35
1	A	286	CYS	N-CA-C	-6.29	107.45	114.62
1	D	120	ALA	N-CA-C	-6.26	104.37	111.07
1	D	267	VAL	CB-CA-C	-6.24	102.97	111.33
1	C	339	ARG	N-CA-C	-6.21	99.89	109.76
1	B	309	ASP	CA-C-N	-6.21	113.38	120.89
1	B	309	ASP	C-N-CA	-6.21	113.38	120.89
1	D	297	VAL	N-CA-C	6.19	117.30	111.48
1	B	115	GLU	CA-C-N	6.12	127.49	119.84
1	B	115	GLU	C-N-CA	6.12	127.49	119.84
1	C	390	ILE	CB-CA-C	-6.09	104.71	111.35
1	D	316	MET	N-CA-C	6.08	116.46	107.88
1	C	147	TYR	N-CA-C	5.99	120.66	113.23
1	B	20	VAL	CB-CA-C	-5.93	101.54	110.55
1	A	218	HIS	N-CA-C	5.92	119.56	112.93
1	B	339	ARG	N-CA-C	-5.91	100.37	109.76
1	B	313	PRO	N-CA-C	-5.87	105.88	113.57
1	A	36	ASN	N-CA-C	5.87	119.70	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	HIS	N-CA-C	5.87	118.11	110.43
1	D	279	LEU	N-CA-C	5.81	118.36	111.33
1	D	384	LYS	N-CA-C	-5.79	105.71	112.89
1	C	383	ARG	N-CA-C	-5.79	106.07	113.01
1	B	410	PRO	CA-C-N	-5.78	116.70	122.77
1	B	410	PRO	C-N-CA	-5.78	116.70	122.77
1	A	198	PHE	N-CA-C	5.76	120.43	113.41
1	C	359	ASN	N-CA-C	-5.75	103.05	110.19
1	D	274	ARG	N-CA-C	5.74	118.31	111.71
1	C	292	THR	N-CA-C	-5.72	102.02	110.48
1	C	393	SER	N-CA-C	5.68	117.94	109.25
1	D	89	ILE	N-CA-C	5.65	116.30	110.82
1	A	17	LEU	N-CA-C	-5.64	101.93	110.28
1	D	89	ILE	CB-CA-C	-5.61	104.70	112.04
1	C	161	THR	N-CA-C	-5.59	105.28	111.71
1	D	343	THR	N-CA-C	5.54	117.47	108.26
1	A	154	LEU	N-CA-C	-5.51	104.49	111.11
1	D	414	ASP	N-CA-C	5.51	118.66	110.39
1	A	295	PRO	CA-N-CD	-5.47	104.34	112.00
1	C	312	SER	CA-C-N	5.44	126.64	119.84
1	C	312	SER	C-N-CA	5.44	126.64	119.84
1	B	293	VAL	N-CA-C	5.43	116.48	108.45
1	C	20	VAL	CB-CA-C	-5.42	102.32	110.55
1	D	36	ASN	N-CA-C	5.41	121.09	113.02
1	B	359	ASN	N-CA-C	5.39	116.88	110.19
1	A	335	LEU	N-CA-C	-5.37	102.93	110.50
1	B	178	ASN	N-CA-C	5.35	121.63	109.81
1	D	118	LYS	N-CA-C	-5.34	105.90	112.90
1	D	358	GLY	N-CA-C	5.33	122.30	115.32
1	A	326	LEU	N-CA-C	-5.33	105.88	112.38
1	C	132	LEU	N-CA-C	5.31	116.52	109.93
1	B	417	ASP	N-CA-C	5.31	116.79	109.15
1	C	101	LEU	N-CA-C	5.30	117.75	111.33
1	D	204	ASN	N-CA-CB	-5.29	101.55	110.49
1	A	369	VAL	N-CA-C	5.28	115.88	108.17
1	B	393	SER	N-CA-C	5.27	118.14	109.76
1	A	399	ARG	N-CA-C	-5.26	102.55	110.70
1	B	364	LEU	N-CA-C	-5.25	107.42	113.88
1	A	136	GLU	N-CA-C	-5.25	105.97	112.38
1	C	378	THR	N-CA-C	-5.24	104.99	111.33
1	A	409	CYS	CA-C-N	-5.21	114.55	119.76
1	A	409	CYS	C-N-CA	-5.21	114.55	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	16	LYS	N-CA-C	5.21	118.11	110.30
1	B	237	LEU	CA-C-N	-5.20	115.77	122.99
1	B	237	LEU	C-N-CA	-5.20	115.77	122.99
1	B	322	GLU	N-CA-C	-5.19	107.00	113.38
1	B	86	LEU	N-CA-C	-5.18	105.53	111.07
1	B	76	LEU	N-CA-C	-5.17	105.54	111.07
1	A	254	GLN	N-CA-C	-5.16	105.55	111.07
1	A	11	HIS	CA-CB-CG	-5.15	108.65	113.80
1	A	39	GLU	N-CA-C	-5.15	106.97	113.20
1	B	335	LEU	N-CA-C	5.12	117.85	110.28
1	D	296	GLU	CA-C-N	-5.10	117.97	122.97
1	D	296	GLU	C-N-CA	-5.10	117.97	122.97
1	B	46	ILE	CB-CA-C	-5.10	102.97	110.82
1	A	316	MET	N-CA-C	5.09	116.93	109.24
1	C	11	HIS	N-CA-C	5.09	120.12	113.30
1	B	286	CYS	N-CA-C	-5.08	107.03	113.43
1	B	231	ILE	N-CA-C	5.07	119.31	112.98
1	D	69	VAL	CB-CA-C	-5.04	104.86	111.25
1	D	62	MET	N-CA-C	-5.03	106.32	112.90
1	A	81	GLN	N-CA-C	-5.02	107.17	113.15
1	A	90	LEU	N-CA-C	5.02	116.75	111.28
1	D	46	ILE	N-CA-C	5.00	116.50	108.89
1	A	91	ASP	N-CA-C	-5.00	105.95	111.71

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	144	TYR	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	3139	0	3128	270	0
1	B	3193	0	3177	166	0
1	C	3147	0	3132	221	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3174	0	3161	216	0
2	A	60	0	0	24	0
2	B	90	0	0	22	0
2	C	86	0	0	15	0
2	D	64	0	0	14	0
All	All	12953	0	12598	840	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 33.

All (840) 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:D:43:ASP:HB2	1:D:401:ARG:HH12	1.05	1.19
1:C:277:MET:HE2	1:C:281:THR:CG2	1.77	1.14
1:C:33:THR:HG22	1:C:35:SER:H	1.17	1.10
1:A:17:LEU:HD11	1:A:20:VAL:HG13	1.34	1.08
1:D:343:THR:HG21	1:D:358:GLY:N	1.69	1.07
1:D:343:THR:HG21	1:D:358:GLY:H	0.90	1.07
1:C:352:ARG:HB2	1:C:377:TYR:CE1	1.90	1.06
1:A:358:GLY:HA3	1:A:378:THR:HG21	1.34	1.04
1:C:180:MET:HA	1:C:180:MET:HE2	1.39	1.03
1:D:343:THR:CG2	1:D:358:GLY:H	1.71	1.03
1:C:150:HIS:HA	2:C:442:HOH:O	1.59	1.01
1:C:277:MET:HE2	1:C:281:THR:HG21	1.01	0.99
1:A:293:VAL:HG23	1:A:295:PRO:HD3	1.02	0.99
1:C:352:ARG:HB2	1:C:377:TYR:CZ	1.98	0.98
1:A:320:ARG:HD3	1:D:143:MET:CE	1.94	0.97
1:D:77:THR:O	1:D:80:ILE:HG22	1.65	0.97
1:B:145:ARG:HD3	1:B:152:SER:HB3	1.50	0.94
1:B:358:GLY:HA3	1:B:378:THR:HG21	1.51	0.93
1:A:293:VAL:HG23	1:A:295:PRO:CD	1.96	0.93
1:A:140:ILE:HA	1:A:143:MET:HE3	1.51	0.92
1:B:304:PHE:CE1	1:C:143:MET:HE3	2.05	0.92
1:C:33:THR:HB	1:C:36:ASN:ND2	1.85	0.92
1:D:43:ASP:HB2	1:D:401:ARG:NH1	1.84	0.91
1:D:7:LYS:H	1:D:7:LYS:HD2	1.36	0.90
1:D:144:TYR:HB3	1:D:150:HIS:HD2	1.34	0.90
1:B:176:THR:HG23	2:B:503:HOH:O	1.71	0.90
1:C:44:ASP:HB3	2:C:498:HOH:O	1.70	0.90
1:D:103:LEU:HD13	1:D:154:LEU:HD23	1.51	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:THR:HG22	2:B:498:HOH:O	1.73	0.87
1:A:227:ASP:HB3	1:A:239:GLY:H	1.37	0.87
1:A:290:LEU:HB3	2:A:478:HOH:O	1.75	0.87
1:A:229:MET:HE3	1:A:279:LEU:HD23	1.56	0.86
1:B:277:MET:HE1	1:B:297:VAL:HG21	1.58	0.85
1:B:304:PHE:HE1	1:C:143:MET:HE3	1.41	0.85
1:C:328:VAL:HG12	1:C:328:VAL:O	1.76	0.85
1:A:293:VAL:CG2	1:A:295:PRO:HD3	1.98	0.85
1:D:416:ILE:HG13	1:D:417:ASP:H	1.41	0.84
1:C:17:LEU:HD22	1:C:413:ARG:NH2	1.92	0.83
1:D:238:ILE:HD12	1:D:265:VAL:HG11	1.61	0.82
1:A:94:ILE:HG12	1:A:107:LEU:HD13	1.59	0.82
1:C:33:THR:HG22	1:C:35:SER:N	1.95	0.82
1:A:12:SER:O	1:A:413:ARG:HD2	1.80	0.82
1:B:390:ILE:HD12	2:B:479:HOH:O	1.81	0.81
1:C:277:MET:CE	1:C:281:THR:HG21	1.97	0.81
1:C:352:ARG:HB2	1:C:377:TYR:CD1	2.15	0.81
1:D:7:LYS:HD2	1:D:7:LYS:N	1.95	0.81
1:D:197:LYS:O	1:D:203:ALA:HB2	1.81	0.81
1:A:336:LYS:H	1:A:336:LYS:HD3	1.44	0.80
1:A:33:THR:HG22	1:A:36:ASN:H	1.45	0.80
1:A:202:PHE:HB2	2:A:463:HOH:O	1.81	0.80
1:A:320:ARG:HD3	1:D:143:MET:SD	2.20	0.80
1:B:404:GLY:HA3	2:B:483:HOH:O	1.81	0.80
1:C:314:TYR:H	1:C:314:TYR:HD2	1.30	0.79
1:D:25:PRO:HD3	1:D:55:HIS:CD2	2.17	0.79
1:B:397:LEU:HB3	1:B:407:MET:HE1	1.63	0.79
1:D:293:VAL:HG22	1:D:298:VAL:HG21	1.66	0.78
1:B:99:VAL:HG13	1:B:154:LEU:HD22	1.64	0.78
1:B:177:LEU:O	1:B:179:PRO:HD3	1.82	0.78
1:A:320:ARG:HD3	1:D:143:MET:HE2	1.64	0.78
1:A:320:ARG:CD	1:D:143:MET:HE2	2.14	0.78
1:A:90:LEU:HD22	1:A:94:ILE:CD1	2.14	0.77
1:A:324:THR:O	1:A:328:VAL:HG23	1.83	0.77
1:A:320:ARG:NE	1:D:143:MET:HE2	1.99	0.77
1:B:69:VAL:HG13	2:B:492:HOH:O	1.83	0.77
1:A:320:ARG:HE	1:D:143:MET:HE2	1.49	0.77
1:A:264:ARG:HH11	1:A:266:ILE:HD11	1.49	0.77
1:D:93:LYS:HE3	1:D:155:LEU:HD23	1.65	0.77
1:A:10:VAL:HB	1:A:170:TRP:O	1.86	0.76
1:A:165:ARG:O	1:A:225:GLY:HA3	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LEU:HD22	1:A:94:ILE:HD12	1.68	0.76
1:B:165:ARG:HD2	1:B:405:HIS:O	1.86	0.75
1:D:399:ARG:HG3	1:D:399:ARG:HH11	1.51	0.75
1:A:278:HIS:O	1:A:281:THR:HB	1.87	0.75
1:D:124:ILE:HG23	1:D:161:THR:HG21	1.69	0.74
1:C:33:THR:HG21	2:C:489:HOH:O	1.86	0.74
1:D:307:ARG:HG2	1:D:307:ARG:HH11	1.52	0.74
1:C:74:ASN:O	1:C:78:GLU:HG3	1.88	0.74
1:A:94:ILE:HG23	1:A:99:VAL:HG21	1.70	0.74
1:D:284:SER:HB2	1:D:292:THR:OG1	1.86	0.74
1:C:87:LYS:HD3	1:C:91:ASP:OD2	1.87	0.73
1:A:397:LEU:HD23	1:A:407:MET:HE1	1.69	0.73
1:B:352:ARG:HG3	1:B:377:TYR:CD2	2.23	0.73
1:C:21:MET:HB2	1:C:411:ILE:HD11	1.71	0.73
1:D:358:GLY:HA3	1:D:378:THR:HG21	1.69	0.73
1:A:264:ARG:NH1	1:A:332:SER:HB3	2.04	0.73
1:B:140:ILE:HA	1:B:143:MET:HE2	1.70	0.73
1:C:352:ARG:CB	1:C:377:TYR:CZ	2.72	0.72
1:D:41:LEU:HD13	1:D:185:ARG:HD2	1.71	0.72
1:D:145:ARG:HD3	1:D:152:SER:HB2	1.72	0.72
1:A:140:ILE:HD13	1:D:318:ILE:HG21	1.71	0.72
1:D:139:ASN:C	1:D:141:LEU:H	1.97	0.72
1:C:76:LEU:HD12	1:C:120:ALA:CB	2.20	0.72
1:B:242:GLU:HB2	1:B:275:ALA:O	1.90	0.72
1:D:153:PHE:HB3	1:D:155:LEU:O	1.89	0.72
1:D:80:ILE:HD11	1:D:119:LEU:HD23	1.70	0.72
1:D:307:ARG:HG2	1:D:307:ARG:NH1	2.04	0.72
1:C:301:ILE:O	1:C:303:PRO:HD3	1.90	0.72
1:B:374:ARG:HH11	1:B:374:ARG:HG3	1.55	0.71
1:D:25:PRO:HA	1:D:29:HIS:CE1	2.24	0.71
1:B:25:PRO:HA	1:B:29:HIS:CE1	2.25	0.71
1:A:402:GLY:HA2	2:A:470:HOH:O	1.90	0.71
1:B:142:LYS:HD2	1:B:145:ARG:NH2	2.05	0.71
1:D:24:SER:HA	1:D:55:HIS:CD2	2.25	0.70
1:B:290:LEU:HG	1:B:339:ARG:NH1	2.05	0.70
1:D:324:THR:O	1:D:328:VAL:HG23	1.91	0.70
1:C:76:LEU:HD12	1:C:120:ALA:HA	1.74	0.69
1:A:145:ARG:NH2	1:A:146:GLU:HG2	2.07	0.69
1:A:208:GLU:HG2	1:A:210:TRP:CZ3	2.27	0.69
1:B:274:ARG:HD3	1:B:297:VAL:HG22	1.74	0.69
1:C:277:MET:CE	1:C:281:THR:CG2	2.65	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:79:THR:OG1	1:D:199:HIS:HB2	1.93	0.69
1:C:314:TYR:N	1:C:314:TYR:CD2	2.58	0.69
1:C:33:THR:CG2	1:C:35:SER:H	1.99	0.69
1:B:54:ASP:HB2	1:B:397:LEU:HD21	1.74	0.69
1:B:277:MET:HE3	1:B:282:VAL:HG11	1.73	0.69
1:A:154:LEU:C	1:A:155:LEU:HD22	2.18	0.68
1:C:80:ILE:HD12	1:C:119:LEU:HD13	1.76	0.68
1:C:197:LYS:HD3	1:C:198:PHE:CE2	2.28	0.68
1:A:301:ILE:HD11	1:A:325:PHE:CD1	2.28	0.68
1:C:197:LYS:HD3	1:C:198:PHE:CZ	2.29	0.68
1:C:314:TYR:HD2	1:C:314:TYR:N	1.92	0.68
1:C:352:ARG:HB2	1:C:377:TYR:CE2	2.28	0.68
1:A:18:ARG:HB2	1:A:412:VAL:HG12	1.76	0.67
1:D:238:ILE:HD12	1:D:265:VAL:CG1	2.23	0.67
1:A:258:ALA:C	1:A:260:GLY:H	2.01	0.67
1:A:107:LEU:HD11	1:A:155:LEU:HD21	1.77	0.67
1:D:36:ASN:O	1:D:40:LEU:HD13	1.95	0.67
1:C:86:LEU:HD22	1:C:90:LEU:HG	1.76	0.66
1:A:117:ARG:HA	2:A:453:HOH:O	1.96	0.66
1:D:103:LEU:HD11	1:D:141:LEU:HD22	1.77	0.66
1:A:94:ILE:HG12	1:A:107:LEU:CD1	2.25	0.66
1:A:145:ARG:HG3	1:A:152:SER:HB3	1.78	0.66
1:A:298:VAL:HG11	1:A:326:LEU:HD21	1.76	0.66
1:A:145:ARG:HB3	1:A:145:ARG:NH1	2.10	0.66
1:A:229:MET:CE	1:A:279:LEU:HD23	2.26	0.66
1:C:264:ARG:HD2	1:C:305:SER:HB2	1.75	0.66
1:D:180:MET:O	1:D:186:ARG:NH1	2.29	0.66
1:D:122:TYR:CD1	1:D:127:VAL:HG22	2.29	0.66
1:C:324:THR:HG23	1:C:327:GLU:OE1	1.96	0.66
1:A:65:ARG:HH11	1:A:65:ARG:HB3	1.61	0.65
1:D:143:MET:O	1:D:146:GLU:HB3	1.96	0.65
1:D:290:LEU:HD12	2:D:447:HOH:O	1.95	0.65
1:A:302:VAL:HG21	1:D:147:TYR:HB2	1.76	0.65
1:A:145:ARG:CG	1:A:152:SER:HB3	2.26	0.65
1:C:145:ARG:HD3	1:C:152:SER:HB3	1.78	0.65
1:A:400:GLY:HA2	2:A:433:HOH:O	1.96	0.65
1:A:155:LEU:HD22	1:A:155:LEU:N	2.11	0.65
1:A:343:THR:HG21	1:A:378:THR:OG1	1.96	0.65
1:B:290:LEU:HG	1:B:339:ARG:HH12	1.62	0.65
1:C:260:GLY:HA2	2:C:476:HOH:O	1.97	0.65
1:D:62:MET:HB3	1:D:67:ILE:HD12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:326:LEU:HD11	1:D:340:VAL:HG11	1.79	0.64
1:A:306:LEU:HD11	1:A:318:ILE:HG12	1.80	0.64
1:A:323:LYS:HG2	2:A:473:HOH:O	1.98	0.64
1:B:99:VAL:HG12	1:B:103:LEU:HB2	1.78	0.64
1:C:303:PRO:HG2	1:C:321:GLU:HB2	1.79	0.64
1:B:397:LEU:HB3	1:B:407:MET:CE	2.27	0.64
1:D:138:ALA:O	1:D:142:LYS:HG3	1.97	0.64
1:C:174:GLY:HA3	1:C:210:TRP:CE2	2.33	0.64
1:A:99:VAL:HG12	1:A:103:LEU:HB2	1.80	0.64
1:D:296:GLU:HA	1:D:299:LYS:HE3	1.79	0.64
1:B:363:CYS:O	1:B:413:ARG:NH2	2.31	0.64
1:C:180:MET:HE1	1:C:224:GLU:HG3	1.80	0.64
1:C:399:ARG:HB3	2:C:498:HOH:O	1.97	0.64
1:B:58:PHE:CE1	1:B:370:VAL:HG11	2.32	0.63
1:B:58:PHE:HE1	1:B:370:VAL:HG11	1.62	0.63
1:D:151:SER:O	1:D:153:PHE:N	2.32	0.63
1:C:180:MET:HE2	1:C:180:MET:CA	2.22	0.63
1:C:320:ARG:HG2	1:C:320:ARG:HH11	1.62	0.63
1:A:47:TRP:HA	1:B:354:GLN:HE22	1.62	0.63
1:A:233:ASN:N	2:A:451:HOH:O	2.29	0.63
1:C:67:ILE:HG22	1:C:68:ASP:N	2.13	0.63
1:D:165:ARG:HD2	1:D:405:HIS:O	1.98	0.63
1:A:233:ASN:O	1:A:233:ASN:ND2	2.30	0.63
1:C:33:THR:HG23	1:C:34:PRO:HD2	1.81	0.63
1:C:372:TYR:OH	1:C:403:GLY:HA2	1.98	0.63
1:A:145:ARG:HB3	1:A:145:ARG:HH11	1.63	0.63
1:A:240:MET:HE2	1:A:246:ARG:HB3	1.80	0.63
1:B:86:LEU:HD22	1:B:90:LEU:HG	1.81	0.63
1:B:20:VAL:HB	2:B:492:HOH:O	1.99	0.63
1:A:320:ARG:CD	1:D:143:MET:CE	2.71	0.62
1:B:80:ILE:HG21	1:B:119:LEU:HD13	1.81	0.62
1:C:45:VAL:HG11	1:D:355:TRP:HE3	1.65	0.62
1:D:165:ARG:O	1:D:225:GLY:HA3	1.99	0.62
1:A:368:VAL:HG13	1:A:388:GLU:OE1	2.00	0.62
1:A:412:VAL:HG12	1:A:412:VAL:O	1.99	0.62
1:B:145:ARG:CD	1:B:152:SER:HB3	2.27	0.62
1:B:217:ASP:O	1:B:218:HIS:HB2	1.98	0.62
1:D:43:ASP:CB	1:D:401:ARG:HH12	1.97	0.62
1:C:33:THR:HB	1:C:36:ASN:HD21	1.61	0.62
1:D:96:ALA:O	1:D:100:GLY:HA2	1.99	0.62
1:C:13:GLU:CD	1:C:229:MET:HG2	2.25	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:TRP:NE1	1:A:92:ARG:NH2	2.48	0.61
1:C:352:ARG:CB	1:C:377:TYR:CE2	2.83	0.61
1:A:92:ARG:HG3	1:A:92:ARG:HH11	1.66	0.61
1:A:227:ASP:HB3	1:A:239:GLY:N	2.13	0.61
1:A:36:ASN:O	1:A:40:LEU:HB2	2.01	0.61
1:B:251:GLN:HG2	1:C:101:LEU:HD11	1.82	0.61
1:D:17:LEU:HD11	1:D:20:VAL:HG22	1.81	0.61
1:C:178:ASN:OD1	1:C:223:LEU:HD12	2.01	0.61
1:C:229:MET:HE3	1:C:279:LEU:HD23	1.82	0.61
1:C:9:GLY:O	1:C:412:VAL:HA	2.00	0.61
1:C:361:VAL:HB	1:C:369:VAL:HG12	1.83	0.61
1:D:125:GLY:O	1:D:157:PRO:HB3	2.01	0.61
1:A:77:THR:O	1:A:81:GLN:HG3	2.00	0.61
1:B:76:LEU:HD22	1:B:80:ILE:HG12	1.82	0.60
1:C:25:PRO:HA	1:C:29:HIS:CE1	2.37	0.60
1:C:328:VAL:O	1:C:328:VAL:CG1	2.47	0.60
1:B:325:PHE:O	1:B:329:VAL:HG23	2.01	0.60
1:D:41:LEU:CD1	1:D:185:ARG:HD2	2.31	0.60
1:A:272:LYS:HE2	1:A:277:MET:HE3	1.83	0.60
1:A:43:ASP:HA	1:A:401:ARG:NH2	2.16	0.60
1:B:94:ILE:HG22	2:B:498:HOH:O	2.02	0.60
1:D:321:GLU:HG3	1:D:328:VAL:HG11	1.82	0.60
1:C:352:ARG:HG3	1:C:377:TYR:CG	2.36	0.60
1:B:398:GLY:C	1:B:400:GLY:H	2.10	0.60
1:A:364:LEU:HD11	1:A:370:VAL:CG2	2.32	0.60
1:A:63:ARG:HG2	1:A:63:ARG:HH11	1.67	0.59
1:D:323:LYS:HB3	1:D:327:GLU:OE1	2.02	0.59
1:B:167:THR:HB	2:B:427:HOH:O	2.01	0.59
1:C:180:MET:HA	1:C:180:MET:CE	2.24	0.59
1:A:281:THR:HG22	1:A:282:VAL:HG13	1.84	0.59
1:B:400:GLY:C	1:B:402:GLY:H	2.11	0.59
1:A:319:ARG:HB3	2:A:421:HOH:O	2.02	0.59
1:B:145:ARG:HD3	1:B:152:SER:CB	2.30	0.59
1:B:169:CYS:HB2	1:B:176:THR:OG1	2.03	0.59
1:C:245:SER:HB2	1:C:247:GLN:OE1	2.03	0.59
1:A:313:PRO:HD2	2:A:471:HOH:O	2.02	0.59
1:A:320:ARG:HG2	1:A:320:ARG:HH11	1.67	0.59
1:A:286:CYS:N	2:A:478:HOH:O	2.35	0.59
1:A:319:ARG:HG3	1:A:320:ARG:N	2.18	0.59
1:C:187:GLN:O	1:C:190:LEU:HB3	2.02	0.59
1:C:165:ARG:O	1:C:225:GLY:HA3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:GLY:HA3	1:A:210:TRP:NE1	2.17	0.58
1:B:132:LEU:HD22	1:B:133:PRO:HD2	1.85	0.58
1:C:60:THR:O	1:C:64:GLU:HG2	2.04	0.58
1:D:157:PRO:O	1:D:158:LEU:HD23	2.02	0.58
1:A:303:PRO:HG2	1:A:321:GLU:HB2	1.86	0.58
1:C:76:LEU:HD12	1:C:120:ALA:CA	2.32	0.58
1:C:99:VAL:HG12	1:C:103:LEU:HB2	1.85	0.58
1:C:416:ILE:HG22	1:C:417:ASP:N	2.18	0.58
1:D:80:ILE:HD11	1:D:119:LEU:CD2	2.32	0.58
1:A:354:GLN:HG3	1:A:355:TRP:N	2.18	0.58
1:C:58:PHE:CD1	1:C:392:ILE:HD13	2.38	0.58
1:C:82:ASN:O	1:C:85:ALA:N	2.36	0.58
1:C:352:ARG:HG3	1:C:377:TYR:CD2	2.38	0.58
1:D:28:ALA:HB2	1:D:125:GLY:HA2	1.85	0.58
1:A:110:TRP:CZ3	1:A:127:VAL:HG11	2.39	0.58
1:B:11:HIS:HB2	1:B:415:PRO:HB3	1.86	0.58
1:B:362:VAL:HG12	2:B:483:HOH:O	2.03	0.58
1:C:10:VAL:HG21	1:C:410:PRO:HB2	1.86	0.58
1:C:45:VAL:HG11	1:D:355:TRP:CE3	2.39	0.58
1:C:180:MET:CE	1:C:224:GLU:HG3	2.33	0.58
1:C:167:THR:HG23	1:C:168:THR:HG23	1.86	0.58
1:A:295:PRO:HG2	1:A:342:GLU:OE2	2.04	0.58
1:C:400:GLY:C	1:C:401:ARG:HG3	2.29	0.58
1:B:295:PRO:O	1:B:299:LYS:HG2	2.04	0.57
1:C:17:LEU:HD22	1:C:413:ARG:HH22	1.65	0.57
1:C:205:ALA:HB3	1:C:207:PHE:HE2	1.69	0.57
1:C:293:VAL:CG1	1:C:294:PHE:N	2.66	0.57
1:D:372:TYR:CD2	1:D:394:ALA:HB2	2.39	0.57
1:B:255:SER:O	1:B:259:LYS:HD3	2.04	0.57
1:B:274:ARG:O	1:B:277:MET:HE2	2.04	0.57
1:C:294:PHE:HE1	1:C:296:GLU:HB3	1.69	0.57
1:A:247:GLN:HG2	1:D:97:ASP:O	2.04	0.57
1:A:288:ARG:HG3	1:A:289:ASP:N	2.19	0.57
1:A:379:ASN:O	1:A:383:ARG:HG3	2.04	0.57
1:D:168:THR:HG22	1:D:177:LEU:HA	1.86	0.57
1:A:242:GLU:HB3	1:A:243:ARG:NH1	2.19	0.57
1:A:264:ARG:NH1	1:A:266:ILE:HD11	2.19	0.57
1:B:320:ARG:HH12	1:B:322:GLU:HG2	1.70	0.57
1:A:295:PRO:O	1:A:296:GLU:C	2.48	0.57
1:C:369:VAL:HG23	1:C:387:VAL:CG1	2.35	0.57
1:B:277:MET:CE	1:B:297:VAL:HG21	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:ILE:HD13	1:C:107:LEU:HD12	1.86	0.57
1:D:188:GLU:HB3	2:D:470:HOH:O	2.04	0.57
1:A:43:ASP:HA	1:A:401:ARG:HH21	1.69	0.56
1:B:167:THR:HG22	1:B:168:THR:HG23	1.87	0.56
1:A:264:ARG:HH12	1:A:332:SER:HB3	1.70	0.56
1:D:223:LEU:HD23	1:D:224:GLU:N	2.19	0.56
1:A:240:MET:HE2	1:A:246:ARG:CB	2.35	0.56
1:A:320:ARG:HG2	1:A:320:ARG:NH1	2.19	0.56
1:A:342:GLU:O	1:A:343:THR:HG23	2.05	0.56
1:B:76:LEU:O	1:B:80:ILE:HG12	2.05	0.56
1:A:258:ALA:C	1:A:260:GLY:N	2.61	0.56
1:B:400:GLY:C	1:B:401:ARG:HG3	2.29	0.56
1:D:321:GLU:HG3	1:D:328:VAL:CG1	2.35	0.56
1:D:416:ILE:O	1:D:417:ASP:HB2	2.04	0.56
1:A:17:LEU:CD1	1:A:20:VAL:HG13	2.23	0.56
1:C:355:TRP:HE3	1:D:45:VAL:HG11	1.71	0.56
1:D:224:GLU:HG3	1:D:243:ARG:HB3	1.87	0.56
1:A:42:PHE:HZ	2:A:447:HOH:O	1.88	0.56
1:C:18:ARG:HD2	1:C:414:ASP:OD2	2.06	0.56
1:D:118:LYS:O	1:D:121:GLU:HB2	2.05	0.56
1:A:17:LEU:HD21	1:A:20:VAL:CG1	2.35	0.56
1:C:288:ARG:HG3	1:C:416:ILE:HG21	1.87	0.56
1:D:331:GLU:CD	1:D:331:GLU:C	2.74	0.56
1:C:17:LEU:HB2	1:C:413:ARG:NH1	2.21	0.56
1:C:33:THR:O	1:C:37:CYS:HB2	2.04	0.56
1:A:63:ARG:HG2	1:A:63:ARG:NH1	2.20	0.55
1:A:372:TYR:OH	1:A:403:GLY:HA2	2.06	0.55
1:D:165:ARG:NE	1:D:405:HIS:CE1	2.74	0.55
1:A:53:ARG:HD3	2:A:475:HOH:O	2.06	0.55
1:D:76:LEU:O	1:D:80:ILE:HB	2.07	0.55
1:D:165:ARG:CZ	1:D:405:HIS:CE1	2.88	0.55
1:C:293:VAL:HG13	1:C:298:VAL:HG21	1.88	0.55
1:C:416:ILE:CG2	1:C:417:ASP:H	2.19	0.55
1:A:145:ARG:HH22	1:A:146:GLU:HG2	1.71	0.55
1:B:142:LYS:HD2	1:B:145:ARG:HH22	1.70	0.55
1:C:354:GLN:HG3	1:C:355:TRP:O	2.06	0.55
1:B:43:ASP:HA	1:B:401:ARG:HH21	1.72	0.55
1:C:257:PHE:CG	1:C:308:PRO:HG3	2.41	0.55
1:A:101:LEU:HD12	1:A:102:GLY:N	2.22	0.55
1:D:233:ASN:ND2	1:D:332:SER:O	2.38	0.55
1:A:294:PHE:CD1	1:A:344:GLY:HA3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:GLY:C	1:A:355:TRP:HE1	2.15	0.55
1:C:178:ASN:HB3	1:C:223:LEU:O	2.07	0.55
1:C:320:ARG:HG2	1:C:320:ARG:NH1	2.21	0.55
1:A:164:THR:O	1:A:409:CYS:HB2	2.08	0.54
1:C:76:LEU:CD1	1:C:120:ALA:HA	2.37	0.54
1:A:343:THR:HG21	1:A:378:THR:CG2	2.38	0.54
1:A:167:THR:HG23	2:A:438:HOH:O	2.08	0.54
1:B:309:ASP:OD1	1:B:311:SER:HB3	2.06	0.54
1:C:139:ASN:HD22	1:C:139:ASN:N	2.05	0.54
1:D:145:ARG:HD3	1:D:152:SER:CB	2.36	0.54
1:C:383:ARG:HG3	1:C:383:ARG:NH1	2.23	0.54
1:A:237:LEU:O	1:A:238:ILE:HG13	2.08	0.54
1:D:169:CYS:O	1:D:175:VAL:HA	2.06	0.54
1:C:383:ARG:HG3	1:C:383:ARG:HH11	1.73	0.54
1:C:416:ILE:CG2	1:C:417:ASP:N	2.71	0.54
1:D:80:ILE:HD11	1:D:119:LEU:CG	2.38	0.54
1:A:33:THR:HG23	1:A:35:SER:H	1.73	0.54
1:C:352:ARG:CG	1:C:377:TYR:CD2	2.90	0.54
1:D:216:LYS:HD3	1:D:218:HIS:CE1	2.42	0.54
1:A:174:GLY:HA3	1:A:210:TRP:CE2	2.42	0.54
1:A:264:ARG:HH21	1:A:307:ARG:HH22	1.55	0.53
1:B:12:SER:OG	1:B:13:GLU:N	2.37	0.53
1:B:72:MET:HE1	2:B:433:HOH:O	2.07	0.53
1:B:86:LEU:HD22	1:B:90:LEU:CG	2.38	0.53
1:B:317:ASN:HD22	1:B:318:ILE:H	1.55	0.53
1:D:172:TYR:O	1:D:174:GLY:N	2.40	0.53
1:B:178:ASN:OD1	1:B:223:LEU:HD12	2.08	0.53
1:C:171:ILE:HG23	1:C:230:PRO:HB3	1.90	0.53
1:D:398:GLY:C	1:D:400:GLY:N	2.64	0.53
1:B:280:ASP:OD2	1:B:405:HIS:ND1	2.40	0.53
1:A:352:ARG:HD2	2:A:457:HOH:O	2.08	0.53
1:A:400:GLY:O	1:A:402:GLY:N	2.41	0.53
1:B:202:PHE:HZ	1:B:411:ILE:HD13	1.74	0.53
1:D:336:LYS:O	1:D:337:LYS:HB3	2.09	0.53
1:A:76:LEU:O	1:A:79:THR:HB	2.09	0.53
1:C:294:PHE:CE1	1:C:296:GLU:HB3	2.43	0.53
1:D:326:LEU:CD1	1:D:340:VAL:HG11	2.39	0.53
1:B:27:LEU:HD21	1:B:31:ARG:NH2	2.24	0.53
1:B:276:ALA:HB2	2:B:430:HOH:O	2.08	0.53
1:A:58:PHE:CE1	1:A:370:VAL:HG11	2.44	0.53
1:A:202:PHE:O	1:A:205:ALA:HB3	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:VAL:HG23	2:B:492:HOH:O	2.08	0.53
1:A:142:LYS:HA	1:A:145:ARG:HH11	1.73	0.53
1:A:169:CYS:SG	1:A:225:GLY:HA2	2.48	0.53
1:B:86:LEU:HD22	1:B:90:LEU:CD1	2.39	0.53
1:C:205:ALA:HB3	1:C:207:PHE:CE2	2.44	0.53
1:D:154:LEU:HD12	1:D:154:LEU:H	1.74	0.53
1:B:372:TYR:OH	1:B:407:MET:HE2	2.09	0.53
1:D:236:VAL:HG12	1:D:238:ILE:HG13	1.89	0.53
1:B:274:ARG:CD	1:B:297:VAL:HG22	2.39	0.53
1:B:277:MET:HB2	1:B:281:THR:HG21	1.90	0.53
1:D:8:LEU:HA	1:D:412:VAL:HG22	1.91	0.53
1:A:337:LYS:HD3	2:A:428:HOH:O	2.08	0.52
1:D:72:MET:CE	1:D:192:THR:OG1	2.57	0.52
1:A:199:HIS:CE1	1:A:201:GLU:HG3	2.45	0.52
1:D:171:ILE:HG23	1:D:230:PRO:HB3	1.91	0.52
1:A:80:ILE:HB	1:A:86:LEU:HD13	1.91	0.52
1:B:397:LEU:N	1:B:397:LEU:HD22	2.24	0.52
1:B:397:LEU:CB	1:B:407:MET:HE1	2.35	0.52
1:A:141:LEU:HD21	1:D:246:ARG:CZ	2.39	0.52
1:A:149:GLY:O	1:D:272:LYS:HG3	2.10	0.52
1:B:20:VAL:CG2	2:B:492:HOH:O	2.58	0.52
1:B:80:ILE:CB	2:B:475:HOH:O	2.58	0.52
1:B:374:ARG:HD2	1:B:394:ALA:HB3	1.91	0.52
1:A:10:VAL:CG2	1:A:170:TRP:HB2	2.40	0.52
1:A:209:ILE:O	1:A:209:ILE:HG22	2.10	0.52
1:A:33:THR:HB	1:A:36:ASN:ND2	2.25	0.52
1:C:400:GLY:O	1:C:401:ARG:HG3	2.09	0.52
1:A:309:ASP:HB2	1:A:317:ASN:HB2	1.92	0.52
1:B:167:THR:HG23	1:B:189:THR:HA	1.90	0.52
1:C:19:LYS:NZ	1:C:201:GLU:OE2	2.43	0.52
1:C:211:TYR:OH	1:C:221:SER:HB3	2.09	0.52
1:D:99:VAL:HG12	1:D:103:LEU:HB2	1.91	0.52
1:A:169:CYS:O	1:A:175:VAL:HA	2.09	0.51
1:A:199:HIS:CE1	1:A:201:GLU:CG	2.93	0.51
1:D:10:VAL:HG13	1:D:413:ARG:HB3	1.92	0.51
1:A:58:PHE:CZ	1:A:62:MET:HE2	2.45	0.51
1:B:58:PHE:CD1	1:B:392:ILE:HD13	2.45	0.51
1:B:235:VAL:HG12	1:B:236:VAL:N	2.24	0.51
1:B:330:ALA:HB2	2:B:449:HOH:O	2.10	0.51
1:D:217:ASP:O	1:D:218:HIS:HB2	2.09	0.51
1:A:18:ARG:HE	1:A:414:ASP:CG	2.19	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:165:ARG:HD3	1:D:405:HIS:ND1	2.25	0.51
1:A:17:LEU:HD21	1:A:20:VAL:HG11	1.91	0.51
1:A:218:HIS:O	1:A:221:SER:HB2	2.10	0.51
1:D:16:LYS:HD2	1:D:18:ARG:NH1	2.26	0.51
1:D:24:SER:HA	1:D:55:HIS:CG	2.46	0.51
1:A:60:THR:O	1:A:64:GLU:HG2	2.10	0.51
1:B:141:LEU:HD21	1:C:246:ARG:CZ	2.40	0.51
1:C:15:GLY:HA3	1:C:414:ASP:O	2.11	0.51
1:D:399:ARG:HG3	1:D:399:ARG:NH1	2.21	0.51
1:A:12:SER:O	1:A:413:ARG:CD	2.54	0.51
1:B:94:ILE:N	1:B:94:ILE:HD12	2.26	0.51
1:D:249:ILE:HD12	2:D:476:HOH:O	2.10	0.51
1:B:226:GLY:O	1:B:280:ASP:HB3	2.11	0.51
1:D:18:ARG:HG3	1:D:18:ARG:HH11	1.76	0.51
1:A:19:LYS:NZ	1:A:68:ASP:OD2	2.43	0.51
1:B:70:LEU:HB3	1:B:75:LEU:HD11	1.92	0.51
1:B:72:MET:HG2	1:B:124:ILE:HD11	1.92	0.51
1:A:253:ALA:CB	1:A:306:LEU:HD23	2.40	0.50
1:C:395:SER:OG	1:D:395:SER:HB3	2.11	0.50
1:A:248:ALA:O	1:A:252:VAL:HG23	2.11	0.50
1:A:338:LEU:O	1:A:340:VAL:HG23	2.11	0.50
1:C:54:ASP:HB2	1:C:397:LEU:CD1	2.42	0.50
1:A:142:LYS:HA	1:A:145:ARG:NH1	2.26	0.50
1:B:47:TRP:CD2	1:B:50:GLN:HB2	2.46	0.50
1:C:64:GLU:C	1:C:66:GLY:H	2.19	0.50
1:C:358:GLY:HA3	1:C:378:THR:HG21	1.92	0.50
1:A:372:TYR:CZ	1:A:403:GLY:HA2	2.47	0.50
1:B:17:LEU:HD21	1:B:20:VAL:CG1	2.41	0.50
1:C:355:TRP:CE3	1:D:45:VAL:HG11	2.46	0.50
1:A:179:PRO:HB3	1:A:218:HIS:CD2	2.46	0.50
1:B:317:ASN:ND2	1:B:318:ILE:N	2.59	0.50
1:A:10:VAL:HG23	1:A:170:TRP:CB	2.42	0.50
1:A:237:LEU:C	1:A:238:ILE:HG13	2.36	0.50
1:B:317:ASN:HD22	1:B:318:ILE:N	2.10	0.50
1:D:199:HIS:HE1	1:D:201:GLU:HG3	1.77	0.50
1:B:102:GLY:C	1:B:103:LEU:HD23	2.37	0.50
1:B:167:THR:CG2	1:B:189:THR:HA	2.42	0.50
1:D:27:LEU:HA	1:D:30:GLN:OE1	2.12	0.50
1:D:28:ALA:HB1	1:D:157:PRO:HB2	1.92	0.50
1:D:23:CYS:O	1:D:55:HIS:NE2	2.45	0.49
1:C:76:LEU:O	1:C:80:ILE:HG12	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:ILE:O	1:A:318:ILE:HG22	2.11	0.49
1:A:336:LYS:HG2	1:A:337:LYS:HG2	1.95	0.49
1:B:288:ARG:HG3	1:B:289:ASP:N	2.26	0.49
1:B:365:GLU:O	1:B:366:PRO:C	2.54	0.49
1:C:397:LEU:HB3	1:C:407:MET:HE1	1.94	0.49
1:A:22:VAL:O	1:A:71:GLU:HA	2.12	0.49
1:C:352:ARG:HB2	1:C:377:TYR:CG	2.47	0.49
1:D:94:ILE:N	1:D:94:ILE:HD12	2.27	0.49
1:D:295:PRO:HG2	1:D:342:GLU:OE2	2.12	0.49
1:A:10:VAL:HG23	1:A:170:TRP:HB3	1.94	0.49
1:A:72:MET:CE	1:A:192:THR:OG1	2.61	0.49
1:A:115:GLU:HG2	1:A:118:LYS:HB2	1.94	0.49
1:D:85:ALA:HA	1:D:198:PHE:CD2	2.48	0.49
1:C:122:TYR:O	1:C:123:LEU:C	2.54	0.49
1:D:42:PHE:O	1:D:43:ASP:C	2.55	0.49
1:D:131:ASP:OD2	1:D:131:ASP:N	2.44	0.49
1:C:307:ARG:NH1	1:C:307:ARG:HG2	2.27	0.49
1:C:153:PHE:HB2	2:C:436:HOH:O	2.11	0.49
1:A:18:ARG:HB2	1:A:412:VAL:O	2.13	0.49
1:A:90:LEU:HD22	1:A:94:ILE:HD11	1.95	0.49
1:C:373:ASP:OD2	1:C:374:ARG:N	2.45	0.49
1:D:72:MET:HE2	1:D:192:THR:OG1	2.12	0.49
1:D:169:CYS:SG	1:D:225:GLY:HA2	2.53	0.49
1:A:247:GLN:OE1	1:A:247:GLN:N	2.39	0.48
1:A:279:LEU:C	1:A:281:THR:H	2.21	0.48
1:B:80:ILE:HB	2:B:475:HOH:O	2.11	0.48
1:D:231:ILE:HD12	1:D:333:LEU:HD21	1.94	0.48
1:A:86:LEU:O	1:A:86:LEU:HG	2.13	0.48
1:C:313:PRO:HD2	1:C:314:TYR:CE2	2.49	0.48
1:D:328:VAL:HG12	1:D:328:VAL:O	2.13	0.48
1:A:58:PHE:CD1	1:A:392:ILE:HD13	2.48	0.48
1:A:199:HIS:ND1	1:A:201:GLU:HB2	2.27	0.48
1:A:258:ALA:O	1:A:260:GLY:N	2.45	0.48
1:D:175:VAL:HG22	1:D:176:THR:N	2.28	0.48
1:A:257:PHE:CD2	1:A:316:MET:HG2	2.48	0.48
1:C:10:VAL:HB	1:C:170:TRP:O	2.14	0.48
1:A:140:ILE:HD13	1:D:318:ILE:CG2	2.41	0.48
1:D:24:SER:HB2	2:D:425:HOH:O	2.12	0.48
1:A:19:LYS:HB3	1:A:412:VAL:HB	1.95	0.48
1:A:356:ASP:HB2	1:A:375:ASN:OD1	2.13	0.48
1:B:76:LEU:HD22	1:B:80:ILE:CG1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ARG:HG2	1:A:31:ARG:HH11	1.77	0.48
1:A:140:ILE:O	1:A:143:MET:HB2	2.13	0.48
1:C:169:CYS:SG	1:C:225:GLY:HA2	2.54	0.48
1:C:229:MET:HE3	1:C:279:LEU:CD2	2.43	0.48
1:D:39:GLU:C	1:D:40:LEU:HD12	2.38	0.48
1:D:88:TRP:NE1	1:D:92:ARG:CZ	2.77	0.48
1:A:272:LYS:CE	1:A:277:MET:HE3	2.44	0.47
1:A:323:LYS:HE2	1:A:327:GLU:O	2.14	0.47
1:B:374:ARG:HH11	1:B:374:ARG:CG	2.25	0.47
1:C:67:ILE:CG2	1:C:68:ASP:N	2.77	0.47
1:D:373:ASP:OD1	1:D:393:SER:HA	2.14	0.47
1:A:76:LEU:O	1:A:79:THR:N	2.47	0.47
1:C:153:PHE:HD2	2:C:436:HOH:O	1.96	0.47
1:C:183:PRO:HD2	2:C:441:HOH:O	2.13	0.47
1:D:37:CYS:HB2	1:D:42:PHE:O	2.14	0.47
1:A:28:ALA:CB	2:A:456:HOH:O	2.62	0.47
1:B:25:PRO:HA	1:B:29:HIS:HE1	1.77	0.47
1:C:242:GLU:CD	1:C:276:ALA:HA	2.40	0.47
1:A:242:GLU:OE1	1:A:243:ARG:NH1	2.47	0.47
1:A:343:THR:CG2	1:A:378:THR:OG1	2.61	0.47
1:A:384:LYS:C	1:A:386:GLY:H	2.22	0.47
1:A:10:VAL:HG21	1:A:170:TRP:HB2	1.95	0.47
1:C:264:ARG:HD3	1:C:307:ARG:CZ	2.44	0.47
1:A:17:LEU:HD11	1:A:20:VAL:CG1	2.24	0.47
1:A:71:GLU:OE2	1:A:73:HIS:ND1	2.48	0.47
1:A:213:ASP:OD1	1:A:215:ASP:HB2	2.15	0.47
1:A:233:ASN:C	1:A:235:VAL:H	2.22	0.47
1:A:293:VAL:HB	1:A:298:VAL:HG21	1.97	0.47
1:B:277:MET:HE3	1:B:282:VAL:CG1	2.42	0.47
1:D:25:PRO:HD3	1:D:55:HIS:CG	2.49	0.47
1:D:88:TRP:HE1	1:D:92:ARG:NH2	2.12	0.47
1:D:208:GLU:OE1	1:D:259:LYS:NZ	2.48	0.47
1:B:254:GLN:HB2	1:B:316:MET:HE2	1.95	0.47
1:A:155:LEU:N	1:A:155:LEU:CD2	2.76	0.47
1:A:165:ARG:NH2	1:A:405:HIS:CD2	2.82	0.47
1:A:394:ALA:HB1	1:A:398:GLY:HA3	1.96	0.47
1:A:132:LEU:HG	1:A:133:PRO:HD2	1.97	0.47
1:C:282:VAL:O	1:C:298:VAL:CG2	2.63	0.47
1:D:245:SER:C	2:D:476:HOH:O	2.57	0.47
1:D:416:ILE:HG23	1:D:417:ASP:N	2.29	0.47
1:A:33:THR:HG22	1:A:36:ASN:N	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:LEU:HD12	1:C:120:ALA:HB1	1.94	0.46
1:C:48:VAL:O	1:C:52:LYS:HG3	2.15	0.46
1:C:264:ARG:NH2	1:C:332:SER:HB2	2.30	0.46
1:C:381:LEU:HG	2:C:484:HOH:O	2.14	0.46
1:D:58:PHE:CE1	1:D:370:VAL:HG11	2.50	0.46
1:D:292:THR:O	1:D:293:VAL:HG23	2.16	0.46
1:D:343:THR:HG22	1:D:357:ASP:HA	1.97	0.46
1:C:370:VAL:HA	1:C:390:ILE:O	2.14	0.46
1:A:271:PRO:O	1:A:272:LYS:C	2.58	0.46
1:C:313:PRO:HD2	1:C:314:TYR:HE2	1.80	0.46
1:D:198:PHE:O	1:D:199:HIS:C	2.57	0.46
1:D:247:GLN:O	1:D:251:GLN:HG3	2.16	0.46
1:D:271:PRO:HG3	1:D:300:GLU:HB2	1.97	0.46
1:D:372:TYR:OH	1:D:403:GLY:HA2	2.15	0.46
1:A:88:TRP:CZ3	1:A:194:ALA:HB2	2.50	0.46
1:A:156:PRO:HD2	2:A:468:HOH:O	2.15	0.46
1:B:17:LEU:HD12	1:B:412:VAL:O	2.15	0.46
1:B:309:ASP:CG	1:B:311:SER:HB3	2.41	0.46
1:B:274:ARG:HB3	1:B:297:VAL:HG22	1.97	0.46
1:C:166:ALA:HA	1:C:225:GLY:HA3	1.98	0.46
1:C:397:LEU:HD13	2:C:419:HOH:O	2.15	0.46
1:D:15:GLY:O	1:D:413:ARG:NH1	2.43	0.46
1:B:208:GLU:HG2	1:B:210:TRP:CZ3	2.51	0.46
1:C:416:ILE:HG22	1:C:417:ASP:H	1.79	0.46
1:D:17:LEU:HD21	1:D:20:VAL:HG21	1.97	0.46
1:D:84:GLU:O	1:D:85:ALA:C	2.58	0.46
1:D:245:SER:O	1:D:249:ILE:HG13	2.15	0.46
1:B:302:VAL:HB	1:C:148:LEU:HD21	1.96	0.46
1:A:61:LYS:O	1:A:65:ARG:HG2	2.15	0.46
1:B:222:THR:O	1:B:244:SER:HA	2.16	0.46
1:B:411:ILE:HG22	1:B:412:VAL:HG23	1.98	0.46
1:C:45:VAL:O	1:C:45:VAL:HG12	2.16	0.46
1:C:352:ARG:HB2	1:C:377:TYR:CD2	2.51	0.46
1:C:372:TYR:CZ	1:C:403:GLY:HA2	2.51	0.46
1:D:293:VAL:O	1:D:295:PRO:HD3	2.16	0.46
1:C:65:ARG:O	1:C:65:ARG:HG3	2.15	0.46
1:C:137:GLY:O	1:C:138:ALA:C	2.59	0.46
1:D:90:LEU:HD13	1:D:108:ARG:HG3	1.98	0.46
1:D:140:ILE:O	1:D:140:ILE:HG13	2.14	0.46
1:D:278:HIS:O	1:D:281:THR:OG1	2.28	0.46
1:D:373:ASP:OD2	1:D:374:ARG:N	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:ILE:HD11	1:A:325:PHE:HB2	1.98	0.45
1:B:17:LEU:HD21	1:B:20:VAL:HG13	1.98	0.45
1:D:291:VAL:HG23	1:D:340:VAL:HG12	1.97	0.45
1:D:307:ARG:HH11	1:D:307:ARG:CG	2.19	0.45
1:A:301:ILE:HD12	1:A:301:ILE:C	2.41	0.45
1:A:59:VAL:HG13	1:A:69:VAL:HG11	1.98	0.45
1:B:36:ASN:O	1:B:40:LEU:HB2	2.16	0.45
1:B:176:THR:N	2:B:503:HOH:O	2.48	0.45
1:D:58:PHE:HE1	1:D:370:VAL:HG11	1.81	0.45
1:D:136:GLU:O	1:D:140:ILE:HG12	2.16	0.45
1:A:235:VAL:HG21	1:A:332:SER:HB2	1.97	0.45
1:B:372:TYR:OH	1:B:407:MET:CE	2.64	0.45
1:C:396:GLU:N	2:C:419:HOH:O	2.40	0.45
1:A:18:ARG:NH2	1:A:414:ASP:OD1	2.48	0.45
1:A:165:ARG:CZ	1:A:405:HIS:CE1	2.99	0.45
1:A:294:PHE:CE1	1:A:344:GLY:HA3	2.50	0.45
1:B:17:LEU:HD21	1:B:20:VAL:CG2	2.46	0.45
1:B:288:ARG:HG2	1:B:288:ARG:HH11	1.82	0.45
1:C:54:ASP:HB2	1:C:397:LEU:HD11	1.97	0.45
1:C:119:LEU:HD22	1:C:123:LEU:HG	1.99	0.45
1:C:267:VAL:HB	1:C:304:PHE:HB2	1.99	0.45
1:C:335:LEU:HD21	2:C:473:HOH:O	2.16	0.45
1:D:169:CYS:HA	2:D:448:HOH:O	2.16	0.45
1:D:290:LEU:HG	1:D:339:ARG:NH2	2.31	0.45
1:D:401:ARG:NH2	2:D:460:HOH:O	2.49	0.45
1:A:88:TRP:CE2	1:A:92:ARG:CZ	3.00	0.45
1:A:267:VAL:HB	1:A:304:PHE:HB2	1.99	0.45
1:C:307:ARG:HG2	1:C:307:ARG:HH11	1.81	0.45
1:A:264:ARG:HD3	1:A:266:ILE:CG1	2.47	0.45
1:A:304:PHE:HE1	1:D:143:MET:SD	2.40	0.45
1:B:103:LEU:HD13	1:B:154:LEU:HD21	1.99	0.45
1:B:132:LEU:CD2	1:B:133:PRO:HD2	2.47	0.45
1:D:179:PRO:HB3	2:D:469:HOH:O	2.17	0.45
1:A:233:ASN:O	1:A:235:VAL:N	2.50	0.45
1:A:286:CYS:HB2	2:A:478:HOH:O	2.17	0.45
1:A:290:LEU:C	2:A:478:HOH:O	2.59	0.45
1:B:164:THR:O	1:B:409:CYS:HB2	2.16	0.45
1:C:51:ALA:O	1:C:397:LEU:HD11	2.17	0.45
1:C:89:ILE:HD11	1:C:194:ALA:CB	2.47	0.45
1:C:93:LYS:HE3	1:C:155:LEU:HD12	1.99	0.45
1:C:306:LEU:HD12	1:C:306:LEU:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:355:TRP:CG	1:C:356:ASP:N	2.85	0.45
1:D:17:LEU:HB2	1:D:413:ARG:NH1	2.31	0.45
1:D:144:TYR:HB3	1:D:150:HIS:CD2	2.27	0.45
1:A:120:ALA:HB3	2:A:453:HOH:O	2.17	0.45
1:A:199:HIS:HE1	1:A:201:GLU:CG	2.30	0.45
1:B:169:CYS:SG	1:B:225:GLY:HA2	2.57	0.45
1:C:178:ASN:HA	1:C:179:PRO:HD3	1.83	0.45
1:D:158:LEU:O	1:D:161:THR:HG23	2.17	0.45
1:D:240:MET:HG3	2:D:476:HOH:O	2.17	0.45
1:A:257:PHE:CE2	1:A:316:MET:HG2	2.53	0.44
1:B:182:TRP:O	1:B:183:PRO:C	2.59	0.44
1:C:155:LEU:N	1:C:155:LEU:HD22	2.32	0.44
1:C:129:ALA:HA	1:C:154:LEU:HD11	1.98	0.44
1:D:256:LEU:HB3	1:D:262:ALA:HB3	1.98	0.44
1:A:166:ALA:HB1	1:A:180:MET:HE1	1.99	0.44
1:A:235:VAL:HB	2:A:451:HOH:O	2.18	0.44
1:C:94:ILE:HD13	1:C:107:LEU:CD1	2.48	0.44
1:C:199:HIS:CG	1:C:200:PRO:HD2	2.53	0.44
1:B:372:TYR:CD2	1:B:394:ALA:HB2	2.52	0.44
1:C:38:ASP:O	1:C:39:GLU:C	2.58	0.44
1:C:177:LEU:O	1:C:212:GLY:HA3	2.17	0.44
1:C:374:ARG:HH22	1:D:399:ARG:HH21	1.64	0.44
1:B:320:ARG:NH1	1:B:320:ARG:HG2	2.31	0.44
1:B:372:TYR:CG	1:B:394:ALA:HB2	2.51	0.44
1:C:180:MET:CA	1:C:180:MET:CE	2.92	0.44
1:A:33:THR:HB	1:A:36:ASN:HD21	1.81	0.44
1:A:307:ARG:NE	2:A:421:HOH:O	2.44	0.44
1:A:324:THR:O	1:A:327:GLU:HB2	2.17	0.44
1:A:331:GLU:C	1:A:333:LEU:N	2.75	0.44
1:A:384:LYS:C	1:A:386:GLY:N	2.74	0.44
1:B:86:LEU:HD22	1:B:90:LEU:HD11	2.00	0.44
1:D:82:ASN:HB3	1:D:85:ALA:CB	2.48	0.44
1:D:111:LEU:HA	1:D:114:LEU:CD1	2.47	0.44
1:D:291:VAL:HG21	1:D:338:LEU:HD13	2.00	0.44
1:D:400:GLY:C	1:D:401:ARG:HG3	2.41	0.44
1:A:13:GLU:HG3	1:A:229:MET:HB2	1.99	0.44
1:A:92:ARG:HH11	1:A:92:ARG:CG	2.30	0.44
1:B:20:VAL:CB	2:B:492:HOH:O	2.61	0.44
1:B:85:ALA:HB2	1:B:198:PHE:CG	2.53	0.44
1:B:209:ILE:HG22	1:B:209:ILE:O	2.16	0.44
1:B:364:LEU:HD12	2:B:479:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:374:ARG:NH2	1:D:399:ARG:HH21	2.16	0.44
1:A:29:HIS:ND1	1:A:29:HIS:N	2.66	0.44
1:B:80:ILE:HG21	1:B:119:LEU:CD1	2.47	0.44
1:C:210:TRP:CH2	1:C:261:ALA:HB2	2.52	0.44
1:C:295:PRO:HG3	1:C:342:GLU:HG2	1.99	0.44
1:A:27:LEU:HD21	1:A:31:ARG:HH21	1.83	0.44
1:A:264:ARG:HG2	1:A:265:VAL:N	2.33	0.44
1:A:331:GLU:C	1:A:333:LEU:H	2.26	0.44
1:B:254:GLN:HB2	1:B:316:MET:CE	2.48	0.44
1:D:14:ALA:O	1:D:366:PRO:HG3	2.18	0.44
1:D:41:LEU:CD2	2:D:460:HOH:O	2.65	0.44
1:D:107:LEU:O	1:D:111:LEU:HG	2.18	0.44
1:D:294:PHE:CD2	1:D:297:VAL:HG23	2.52	0.44
1:D:398:GLY:C	1:D:400:GLY:H	2.25	0.44
1:A:202:PHE:O	1:A:205:ALA:N	2.50	0.43
1:B:400:GLY:C	1:B:402:GLY:N	2.76	0.43
1:A:229:MET:SD	1:A:279:LEU:HD23	2.59	0.43
1:A:8:LEU:O	1:A:173:GLY:HA2	2.17	0.43
1:A:58:PHE:CE2	1:A:62:MET:HE2	2.53	0.43
1:A:233:ASN:O	1:A:235:VAL:HG23	2.18	0.43
1:A:290:LEU:HD23	1:A:339:ARG:O	2.19	0.43
1:C:33:THR:HG23	1:C:34:PRO:CD	2.47	0.43
1:C:397:LEU:N	1:C:397:LEU:HD12	2.33	0.43
1:A:239:GLY:O	1:A:244:SER:HB2	2.18	0.43
1:A:264:ARG:HD3	1:A:266:ILE:HG13	2.00	0.43
1:C:63:ARG:HD2	2:C:471:HOH:O	2.17	0.43
1:C:169:CYS:O	1:C:175:VAL:HA	2.18	0.43
1:C:216:LYS:HG2	1:C:217:ASP:N	2.32	0.43
1:A:280:ASP:HA	1:A:283:PHE:O	2.19	0.43
1:B:342:GLU:O	1:B:343:THR:HG22	2.17	0.43
1:A:20:VAL:HG12	1:A:410:PRO:HA	2.00	0.43
1:A:268:ALA:HB2	1:A:325:PHE:CE1	2.54	0.43
1:A:270:LEU:HA	1:A:271:PRO:HD3	1.73	0.43
1:A:329:VAL:CG1	1:A:338:LEU:HD11	2.49	0.43
1:A:387:VAL:HG12	1:A:388:GLU:N	2.33	0.43
1:B:31:ARG:HG2	1:B:31:ARG:HH11	1.84	0.43
1:B:342:GLU:C	1:B:343:THR:CG2	2.91	0.43
1:C:284:SER:O	1:C:291:VAL:HA	2.18	0.43
1:C:372:TYR:CG	1:C:394:ALA:HB2	2.53	0.43
1:A:356:ASP:C	1:A:356:ASP:OD1	2.61	0.43
1:A:364:LEU:HD11	1:A:370:VAL:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ILE:CG1	2:B:475:HOH:O	2.65	0.43
1:D:72:MET:O	1:D:73:HIS:C	2.61	0.43
1:D:174:GLY:HA3	1:D:210:TRP:CE2	2.53	0.43
1:D:275:ALA:O	1:D:276:ALA:C	2.62	0.43
1:A:15:GLY:HA3	1:A:414:ASP:O	2.19	0.43
1:C:33:THR:C	1:C:35:SER:N	2.76	0.43
1:C:267:VAL:N	1:C:304:PHE:O	2.49	0.43
1:D:284:SER:O	1:D:291:VAL:HA	2.19	0.43
1:A:74:ASN:O	1:A:77:THR:HB	2.18	0.43
1:C:80:ILE:HD12	1:C:119:LEU:CD1	2.47	0.43
1:C:141:LEU:HD22	1:C:141:LEU:HA	1.79	0.43
1:D:80:ILE:HD11	1:D:119:LEU:HG	2.00	0.43
1:D:219:GLY:C	1:D:221:SER:H	2.26	0.43
1:A:76:LEU:O	1:A:77:THR:C	2.62	0.43
1:A:303:PRO:CG	1:A:321:GLU:HB2	2.48	0.43
1:D:240:MET:CG	2:D:476:HOH:O	2.67	0.43
1:D:377:TYR:O	1:D:380:THR:HB	2.19	0.43
1:A:24:SER:HA	1:A:55:HIS:CE1	2.54	0.42
1:A:93:LYS:HE3	1:A:93:LYS:HB3	1.84	0.42
1:A:320:ARG:NH1	1:A:321:GLU:O	2.52	0.42
1:C:295:PRO:O	1:C:299:LYS:HG2	2.19	0.42
1:C:396:GLU:HA	1:C:399:ARG:HG3	2.00	0.42
1:D:80:ILE:HD13	1:D:80:ILE:HG21	1.72	0.42
1:D:274:ARG:HD2	2:D:479:HOH:O	2.18	0.42
1:A:143:MET:O	1:A:147:TYR:HB2	2.19	0.42
1:A:325:PHE:C	1:A:327:GLU:N	2.74	0.42
1:B:361:VAL:HG21	1:B:369:VAL:HG11	2.01	0.42
1:C:407:MET:HB3	2:C:428:HOH:O	2.18	0.42
1:D:88:TRP:NE1	1:D:92:ARG:NH2	2.67	0.42
1:B:251:GLN:CG	1:C:101:LEU:HD11	2.47	0.42
1:B:400:GLY:O	1:B:402:GLY:N	2.52	0.42
1:C:92:ARG:HG3	1:C:92:ARG:HH11	1.84	0.42
1:C:323:LYS:HB3	1:C:327:GLU:OE1	2.20	0.42
1:D:99:VAL:HG12	1:D:99:VAL:O	2.18	0.42
1:D:284:SER:HB2	1:D:292:THR:HG1	1.82	0.42
1:D:416:ILE:HD12	1:D:416:ILE:HA	1.84	0.42
1:A:65:ARG:HH11	1:A:65:ARG:CB	2.29	0.42
1:B:154:LEU:HD23	1:B:154:LEU:HA	1.90	0.42
1:C:414:ASP:HA	1:C:415:PRO:HD3	1.75	0.42
1:B:9:GLY:HA2	1:B:172:TYR:O	2.19	0.42
1:B:133:PRO:O	1:B:134:ALA:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:LYS:HE3	1:C:118:LYS:HB2	1.68	0.42
1:D:111:LEU:HA	1:D:114:LEU:HD12	2.01	0.42
1:D:416:ILE:HG13	1:D:417:ASP:N	2.20	0.42
1:A:21:MET:HE1	1:A:192:THR:HG23	2.00	0.42
1:B:136:GLU:O	1:B:137:GLY:C	2.61	0.42
1:B:147:TYR:HB3	1:C:302:VAL:HG21	2.00	0.42
1:B:267:VAL:HB	1:B:304:PHE:HB2	2.02	0.42
1:B:398:GLY:C	1:B:400:GLY:N	2.77	0.42
1:C:312:SER:HA	1:C:313:PRO:HD3	1.86	0.42
1:C:336:LYS:HE3	1:C:336:LYS:HB3	1.88	0.42
1:D:230:PRO:O	1:D:230:PRO:HG2	2.18	0.42
1:D:399:ARG:NH1	1:D:399:ARG:CG	2.83	0.42
1:B:80:ILE:HG23	1:B:86:LEU:HG	2.02	0.42
1:B:165:ARG:O	1:B:225:GLY:HA3	2.20	0.42
1:C:25:PRO:HG3	1:C:162:GLN:OE1	2.19	0.42
1:D:86:LEU:HD13	1:D:86:LEU:C	2.45	0.42
1:A:400:GLY:C	1:A:402:GLY:N	2.78	0.42
1:B:279:LEU:C	1:B:281:THR:H	2.27	0.42
1:C:31:ARG:HG2	1:C:31:ARG:HH11	1.85	0.42
1:C:80:ILE:CD1	1:C:119:LEU:HD13	2.46	0.42
1:C:282:VAL:O	1:C:298:VAL:HG21	2.19	0.42
1:D:186:ARG:HG2	1:D:186:ARG:HH11	1.84	0.42
1:A:170:TRP:CE2	1:A:411:ILE:HG23	2.55	0.42
1:A:286:CYS:CB	2:A:478:HOH:O	2.68	0.42
1:B:69:VAL:HA	2:B:492:HOH:O	2.20	0.42
1:D:266:ILE:HA	1:D:304:PHE:O	2.19	0.42
1:A:65:ARG:NH2	1:A:390:ILE:HD11	2.35	0.42
1:A:322:GLU:C	2:A:473:HOH:O	2.62	0.42
1:B:55:HIS:HE1	2:B:487:HOH:O	2.02	0.42
1:B:87:LYS:HG2	1:B:91:ASP:OD2	2.20	0.42
1:C:206:GLU:O	1:C:207:PHE:HB3	2.20	0.42
1:D:287:ASP:CB	1:D:290:LEU:HB2	2.50	0.42
1:A:245:SER:O	1:A:249:ILE:HG13	2.20	0.41
1:A:246:ARG:HG2	1:A:247:GLN:OE1	2.20	0.41
1:B:101:LEU:HD22	1:C:254:GLN:OE1	2.19	0.41
1:B:240:MET:HE3	1:C:150:HIS:NE2	2.35	0.41
1:B:317:ASN:ND2	1:B:318:ILE:H	2.17	0.41
1:C:23:CYS:O	1:C:162:GLN:HA	2.20	0.41
1:C:351:GLU:HB2	1:C:352:ARG:H	1.41	0.41
1:C:355:TRP:N	2:C:479:HOH:O	2.53	0.41
1:A:58:PHE:HE1	1:A:370:VAL:HG11	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:TRP:HE1	1:A:92:ARG:NH2	2.18	0.41
1:A:141:LEU:HD21	1:D:246:ARG:NE	2.35	0.41
1:B:50:GLN:OE1	1:B:53:ARG:NH2	2.40	0.41
1:D:248:ALA:O	1:D:252:VAL:HG23	2.20	0.41
1:D:306:LEU:CD1	1:D:306:LEU:N	2.84	0.41
1:A:228:VAL:HG22	1:A:238:ILE:HG12	2.01	0.41
1:A:231:ILE:HD13	1:A:237:LEU:HD11	2.02	0.41
1:B:43:ASP:CA	1:B:401:ARG:HH21	2.33	0.41
1:B:128:ALA:O	1:B:129:ALA:C	2.62	0.41
1:C:287:ASP:O	1:C:288:ARG:C	2.62	0.41
1:D:193:THR:O	1:D:194:ALA:C	2.62	0.41
1:D:222:THR:O	1:D:244:SER:HA	2.20	0.41
1:A:8:LEU:HD12	1:A:207:PHE:CD2	2.54	0.41
1:C:8:LEU:HA	1:C:412:VAL:HG22	2.02	0.41
1:C:231:ILE:HD11	1:C:235:VAL:HG11	2.02	0.41
1:C:278:HIS:O	1:C:279:LEU:C	2.63	0.41
1:C:267:VAL:HG23	1:C:306:LEU:CD1	2.51	0.41
1:C:292:THR:HA	1:C:341:VAL:O	2.21	0.41
1:D:379:ASN:HA	1:D:382:LEU:HD12	2.02	0.41
1:A:120:ALA:HA	1:A:123:LEU:HD12	2.02	0.41
1:C:12:SER:O	1:C:413:ARG:HD2	2.20	0.41
1:C:77:THR:O	1:C:81:GLN:HG3	2.20	0.41
1:D:188:GLU:O	1:D:191:LEU:N	2.54	0.41
1:A:63:ARG:HH11	1:A:63:ARG:CG	2.33	0.41
1:A:324:THR:O	1:A:327:GLU:C	2.64	0.41
1:A:353:GLU:OE2	1:A:353:GLU:HA	2.19	0.41
1:B:65:ARG:NH1	1:B:390:ILE:HD11	2.35	0.41
1:C:22:VAL:HG12	1:C:408:THR:HG22	2.02	0.41
1:C:169:CYS:SG	1:C:228:VAL:HB	2.61	0.41
1:D:370:VAL:HA	1:D:390:ILE:O	2.21	0.41
1:A:265:VAL:O	1:A:306:LEU:HB2	2.21	0.41
1:A:306:LEU:CD1	1:A:318:ILE:HG12	2.48	0.41
1:A:336:LYS:HD3	1:A:336:LYS:N	2.23	0.41
1:D:187:GLN:OE1	1:D:187:GLN:HA	2.20	0.41
1:D:270:LEU:HD12	1:D:301:ILE:HG12	2.02	0.41
1:D:354:GLN:HB2	2:D:465:HOH:O	2.21	0.41
1:D:403:GLY:O	1:D:404:GLY:C	2.64	0.41
1:A:142:LYS:O	1:A:146:GLU:HB2	2.21	0.41
1:B:99:VAL:HG13	1:B:154:LEU:CD2	2.44	0.41
1:B:141:LEU:HD21	1:C:246:ARG:NE	2.35	0.41
1:B:169:CYS:O	1:B:175:VAL:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:PHE:HB3	1:B:308:PRO:HB3	2.02	0.41
1:B:297:VAL:O	1:B:301:ILE:HG13	2.21	0.41
1:B:361:VAL:HB	1:B:369:VAL:HG13	2.02	0.41
1:B:374:ARG:CG	1:B:374:ARG:NH1	2.82	0.41
1:C:35:SER:C	1:C:37:CYS:H	2.28	0.41
1:C:199:HIS:HA	1:C:200:PRO:HD3	1.84	0.41
1:C:213:ASP:OD1	1:C:214:PRO:HD2	2.21	0.41
1:C:364:LEU:N	1:C:364:LEU:HD23	2.36	0.41
1:D:112:GLU:C	1:D:114:LEU:H	2.28	0.41
1:C:154:LEU:C	1:C:155:LEU:HD22	2.46	0.41
1:C:288:ARG:HE	1:C:288:ARG:HB2	1.68	0.41
1:D:138:ALA:O	1:D:141:LEU:HB3	2.20	0.41
1:D:324:THR:OG1	1:D:327:GLU:HG3	2.21	0.41
1:D:335:LEU:HD12	1:D:335:LEU:HA	1.93	0.41
1:A:110:TRP:HZ3	1:A:127:VAL:HG11	1.86	0.40
1:A:400:GLY:C	1:A:402:GLY:H	2.29	0.40
1:B:27:LEU:O	1:B:28:ALA:C	2.64	0.40
1:C:65:ARG:HE	1:C:65:ARG:HB2	1.51	0.40
1:C:293:VAL:HG13	1:C:298:VAL:CG2	2.50	0.40
1:A:24:SER:CA	1:A:55:HIS:CE1	3.05	0.40
1:C:199:HIS:O	1:C:202:PHE:N	2.48	0.40
1:D:322:GLU:O	1:D:323:LYS:O	2.39	0.40
1:A:103:LEU:HD13	1:A:141:LEU:HD11	2.03	0.40
1:A:240:MET:HE3	1:A:267:VAL:HG11	2.03	0.40
1:D:286:CYS:SG	1:D:292:THR:HG23	2.62	0.40
1:C:92:ARG:HG3	1:C:92:ARG:NH1	2.37	0.40
1:C:289:ASP:O	1:C:339:ARG:N	2.45	0.40
1:C:294:PHE:H	1:C:298:VAL:HG21	1.86	0.40
1:D:138:ALA:O	1:D:142:LYS:N	2.55	0.40
1:A:18:ARG:HH11	1:A:18:ARG:HG2	1.86	0.40
1:B:361:VAL:HB	1:B:369:VAL:CG1	2.52	0.40
1:D:249:ILE:CD1	2:D:476:HOH:O	2.68	0.40
1:D:388:GLU:OE2	1:D:390:ILE:HD11	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	396/418 (95%)	342 (86%)	45 (11%)	9 (2%)	5	8
1	B	405/418 (97%)	366 (90%)	36 (9%)	3 (1%)	18	34
1	C	397/418 (95%)	346 (87%)	45 (11%)	6 (2%)	8	16
1	D	402/418 (96%)	336 (84%)	53 (13%)	13 (3%)	3	4
All	All	1600/1672 (96%)	1390 (87%)	179 (11%)	31 (2%)	6	11

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	ASP
1	D	152	SER
1	D	276	ALA
1	D	323	LYS
1	A	172	TYR
1	A	225	GLY
1	B	272	LYS
1	C	164	THR
1	D	146	GLU
1	D	173	GLY
1	A	258	ALA
1	A	271	PRO
1	A	328	VAL
1	A	401	ARG
1	C	401	ARG
1	D	296	GLU
1	D	337	LYS
1	B	43	ASP
1	C	272	LYS
1	C	330	ALA
1	D	101	LEU
1	A	234	GLY

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Mol	Chain	Res	Type
1	C	313	PRO
1	D	116	PRO
1	B	401	ARG
1	D	400	GLY
1	C	365	GLU
1	D	404	GLY
1	D	140	ILE
1	A	400	GLY
1	D	137	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	342/354 (97%)	311 (91%)	31 (9%)	9	19
1	B	346/354 (98%)	312 (90%)	34 (10%)	7	16
1	C	343/354 (97%)	315 (92%)	28 (8%)	10	23
1	D	345/354 (98%)	324 (94%)	21 (6%)	17	35
All	All	1376/1416 (97%)	1262 (92%)	114 (8%)	10	22

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	SER
1	A	20	VAL
1	A	33	THR
1	A	40	LEU
1	A	61	LYS
1	A	65	ARG
1	A	71	GLU
1	A	76	LEU
1	A	107	LEU
1	A	109	SER
1	A	118	LYS
1	A	119	LEU

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Mol	Chain	Res	Type
1	A	132	LEU
1	A	176	THR
1	A	178	ASN
1	A	223	LEU
1	A	233	ASN
1	A	242	GLU
1	A	281	THR
1	A	290	LEU
1	A	295	PRO
1	A	301	ILE
1	A	306	LEU
1	A	310	PRO
1	A	333	LEU
1	A	336	LYS
1	A	353	GLU
1	A	369	VAL
1	A	382	LEU
1	A	406	CYS
1	A	417	ASP
1	B	12	SER
1	B	20	VAL
1	B	22	VAL
1	B	25	PRO
1	B	39	GLU
1	B	40	LEU
1	B	60	THR
1	B	76	LEU
1	B	86	LEU
1	B	107	LEU
1	B	109	SER
1	B	117	ARG
1	B	119	LEU
1	B	165	ARG
1	B	179	PRO
1	B	200	PRO
1	B	220	SER
1	B	242	GLU
1	B	274	ARG
1	B	277	MET
1	B	281	THR
1	B	288	ARG
1	B	290	LEU

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Mol	Chain	Res	Type
1	B	293	VAL
1	B	336	LYS
1	B	356	ASP
1	B	363	CYS
1	B	382	LEU
1	B	397	LEU
1	B	401	ARG
1	B	406	CYS
1	B	410	PRO
1	B	414	ASP
1	B	416	ILE
1	C	20	VAL
1	C	21	MET
1	C	33	THR
1	C	76	LEU
1	C	86	LEU
1	C	115	GLU
1	C	116	PRO
1	C	119	LEU
1	C	139	ASN
1	C	141	LEU
1	C	162	GLN
1	C	180	MET
1	C	200	PRO
1	C	206	GLU
1	C	259	LYS
1	C	280	ASP
1	C	290	LEU
1	C	293	VAL
1	C	307	ARG
1	C	311	SER
1	C	314	TYR
1	C	318	ILE
1	C	351	GLU
1	C	352	ARG
1	C	369	VAL
1	C	376	THR
1	C	390	ILE
1	C	401	ARG
1	D	20	VAL
1	D	76	LEU
1	D	81	GLN

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Mol	Chain	Res	Type
1	D	116	PRO
1	D	121	GLU
1	D	131	ASP
1	D	147	TYR
1	D	154	LEU
1	D	165	ARG
1	D	189	THR
1	D	200	PRO
1	D	221	SER
1	D	270	LEU
1	D	272	LYS
1	D	291	VAL
1	D	293	VAL
1	D	295	PRO
1	D	306	LEU
1	D	397	LEU
1	D	401	ARG
1	D	406	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	B	11	HIS
1	B	49	ASN
1	B	251	GLN
1	B	317	ASN
1	C	74	ASN
1	C	139	ASN
1	C	187	GLN
1	C	317	ASN
1	D	49	ASN
1	D	73	HIS
1	D	150	HIS
1	D	218	HIS
1	D	375	ASN

5.3.3 RNA ⓘ

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	402/418 (96%)	-0.41	2 (0%) 87 85	18, 41, 64, 78	0
1	B	409/418 (97%)	-0.56	1 (0%) 91 89	16, 30, 60, 78	0
1	C	403/418 (96%)	-0.50	2 (0%) 87 85	17, 36, 62, 78	0
1	D	406/418 (97%)	-0.43	1 (0%) 91 89	17, 39, 66, 79	0
All	All	1620/1672 (96%)	-0.47	6 (0%) 88 86	16, 37, 64, 79	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	SER	3.1
1	B	276	ALA	2.8
1	D	400	GLY	2.5
1	C	313	PRO	2.5
1	A	402	GLY	2.4
1	C	400	GLY	2.2

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 ⓘ

There are no such residues in this entry.