



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:16 PM UTC

PDB ID : 5JXT / pdb_00005jxt
Title : Crystal structure of MtISWI bound with histone H4 tail
Authors : Chen, Z.; Yan, L.
Deposited on : 2016-05-13
Resolution : 3.01 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

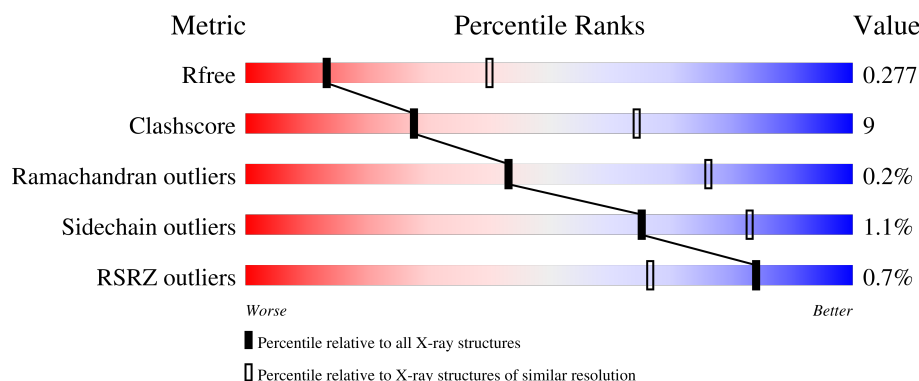
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	349	70% 23% 6%
1	B	349	73% 21% 6%
1	C	349	71% 22% 7%
1	D	349	62% 22% 16%
1	E	349	73% 17% 10%

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Mol	Chain	Length	Quality of chain
1	F	349	 2% 70% 22% 7%
1	G	349	 70% 18% 12%
1	H	349	 70% 21% 8%
1	I	349	 69% 17% 14%
1	J	349	 71% 19% 9%
1	K	349	 68% 19% 12%
1	L	349	 67% 21% 11%
1	M	349	 74% 19% 7%
1	N	349	 75% 16% 9%
1	O	349	 69% 22% 8%
1	P	349	 2% 73% 17% 9%
2	Q	21	 19% 5% 76%
2	R	21	 19% 5% 76%
2	S	21	 19% 5% 76%
2	T	21	 24% 76%
2	U	21	 19% 5% 76%
2	V	21	 19% 5% 76%
2	W	21	 14% 5% 81%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 40399 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 Chromatin-remodeling complex ATPase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	328	Total	C	N	O	S	0	0	0
			2599	1636	454	497	12			
1	E	313	Total	C	N	O	S	0	0	0
			2484	1566	437	469	12			
1	F	324	Total	C	N	O	S	0	0	0
			2578	1623	450	493	12			
1	G	308	Total	C	N	O	S	0	0	0
			2439	1536	430	461	12			
1	J	317	Total	C	N	O	S	0	0	0
			2515	1583	439	481	12			
1	N	319	Total	C	N	O	S	0	0	0
			2536	1596	442	486	12			
1	O	322	Total	C	N	O	S	0	0	0
			2563	1613	447	491	12			
1	P	316	Total	C	N	O	S	0	0	0
			2517	1585	439	481	12			
1	C	324	Total	C	N	O	S	0	0	0
			2574	1623	446	493	12			
1	H	321	Total	C	N	O	S	0	0	0
			2549	1602	446	489	12			
1	K	306	Total	C	N	O	S	0	0	0
			2427	1529	426	460	12			
1	L	311	Total	C	N	O	S	0	0	0
			2464	1551	434	467	12			
1	A	327	Total	C	N	O	S	0	0	0
			2597	1636	453	496	12			
1	I	301	Total	C	N	O	S	0	0	0
			2388	1506	420	450	12			
1	D	292	Total	C	N	O	S	0	0	0
			2326	1467	411	437	11			
1	M	324	Total	C	N	O	S	0	0	0
			2574	1622	448	492	12			

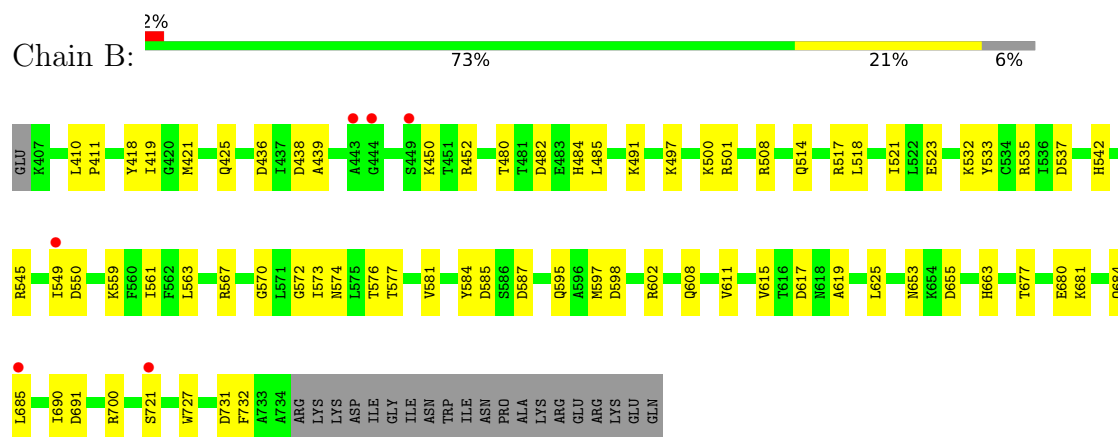
- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	R	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	S	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	U	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	Q	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	T	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	V	5	Total	C	N	O	0	0	0
			39	23	11	5			
2	W	4	Total	C	N	O	0	0	0
			35	21	10	4			

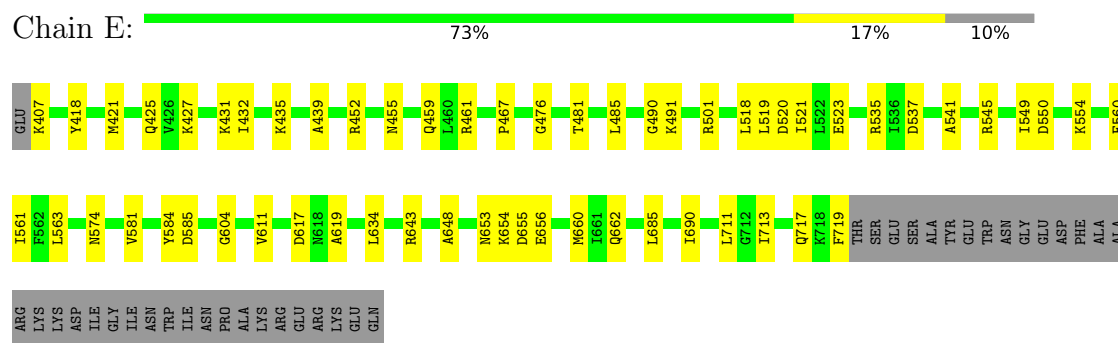
3 Residue-property plots

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.

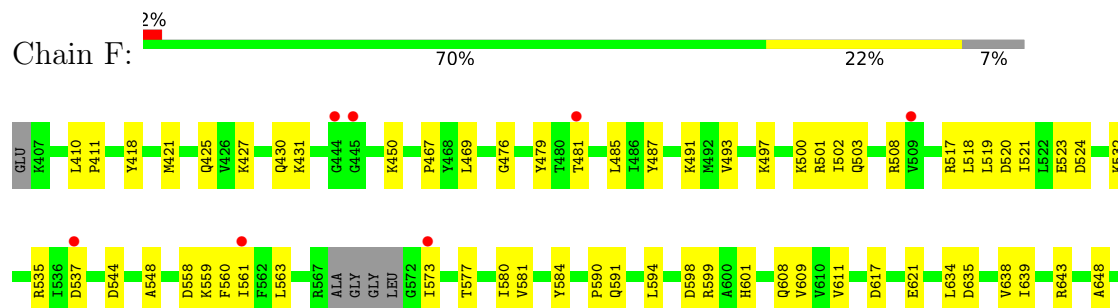
- Molecule 1: Chromatin-remodeling complex ATPase-like protein



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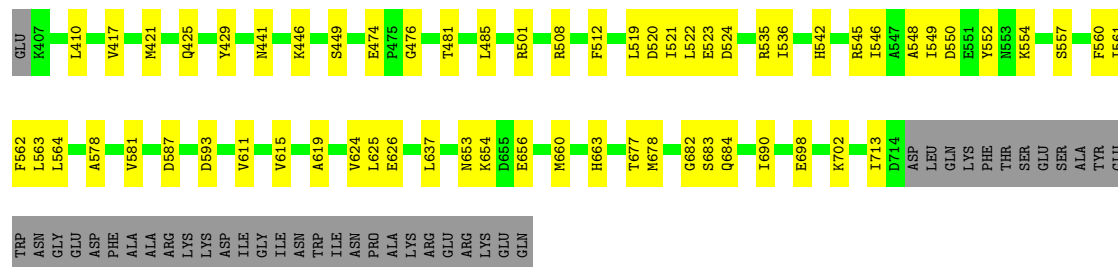


- Molecule 1: Chromatin-remodeling complex ATPase-like protein

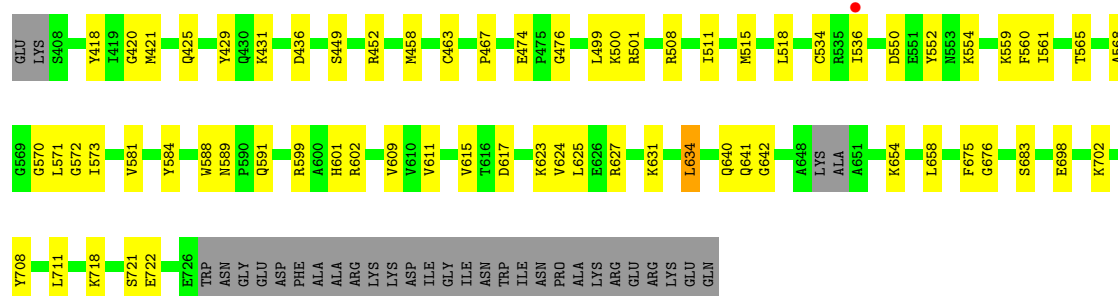




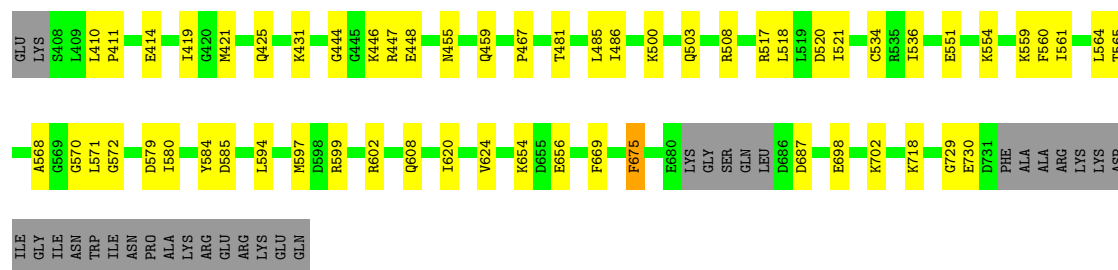
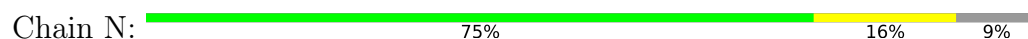
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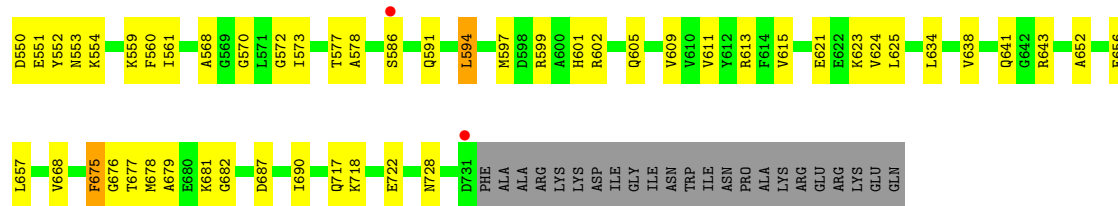


- Molecule 1: Chromatin-remodeling complex ATPase-like protein

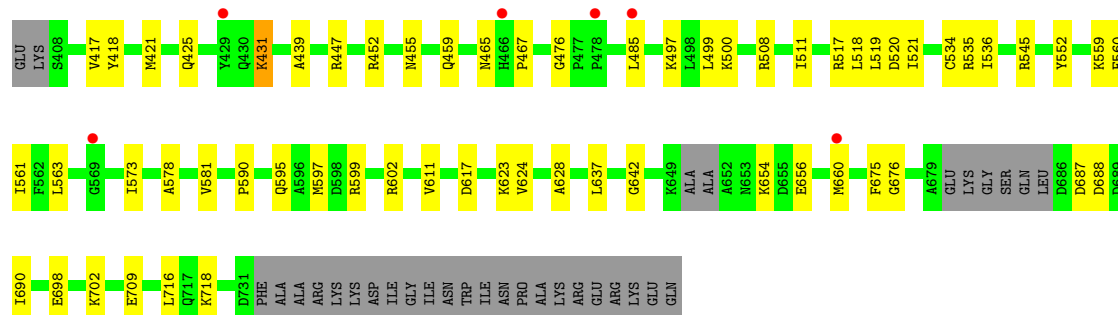
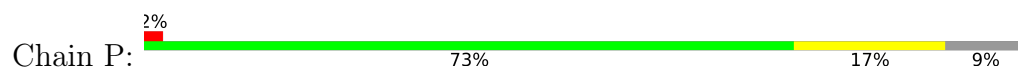


- Molecule 1: Chromatin-remodeling complex ATPase-like protein

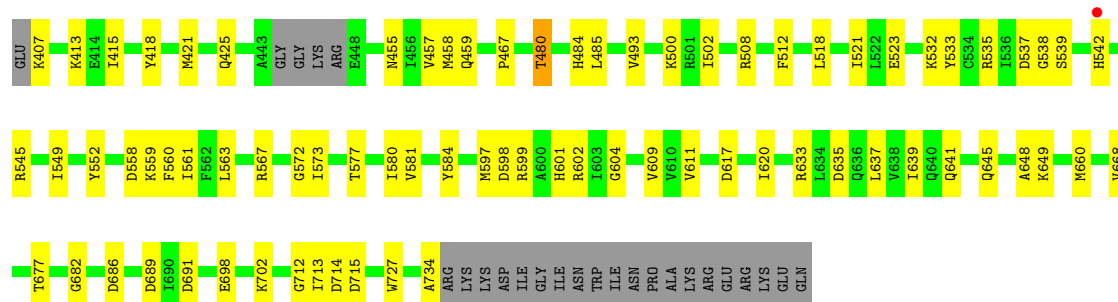




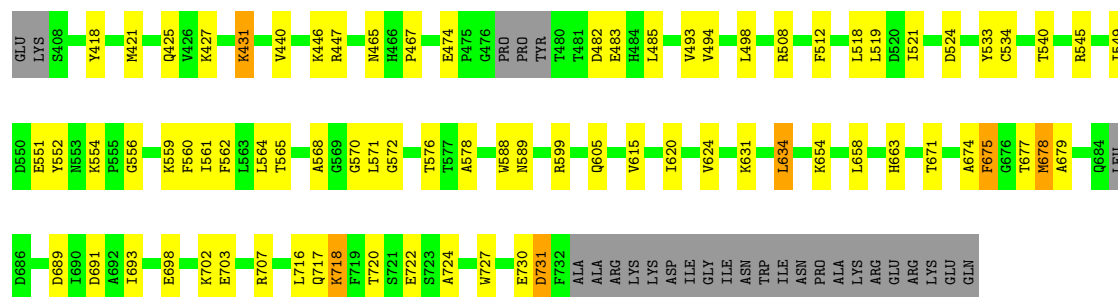
• Molecule 1: Chromatin-remodeling complex ATPase-like protein



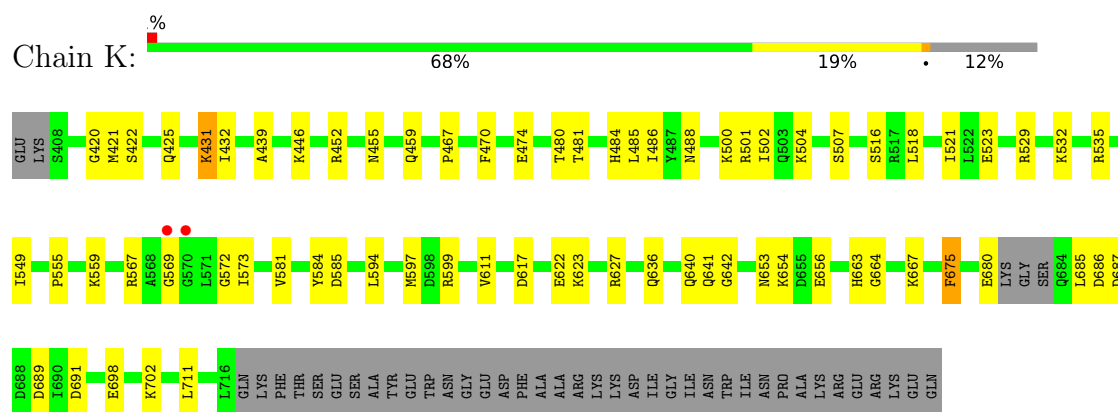
• Molecule 1: Chromatin-remodeling complex ATPase-like protein



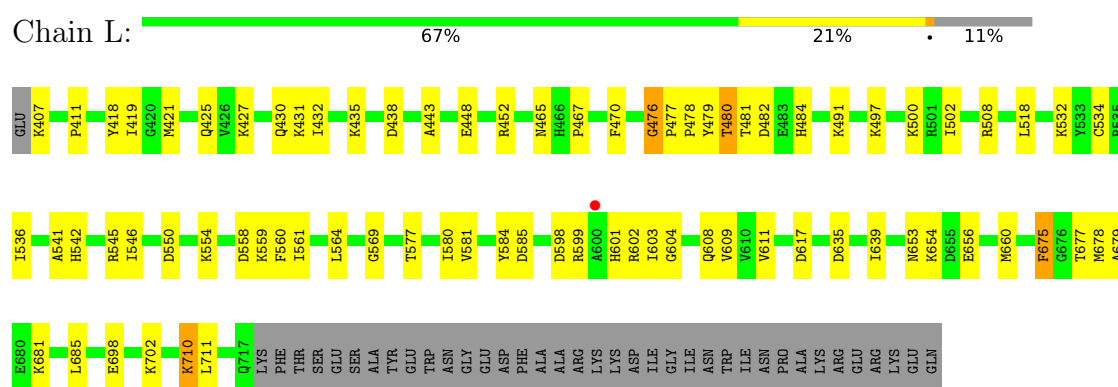
• Molecule 1: Chromatin-remodeling complex ATPase-like protein



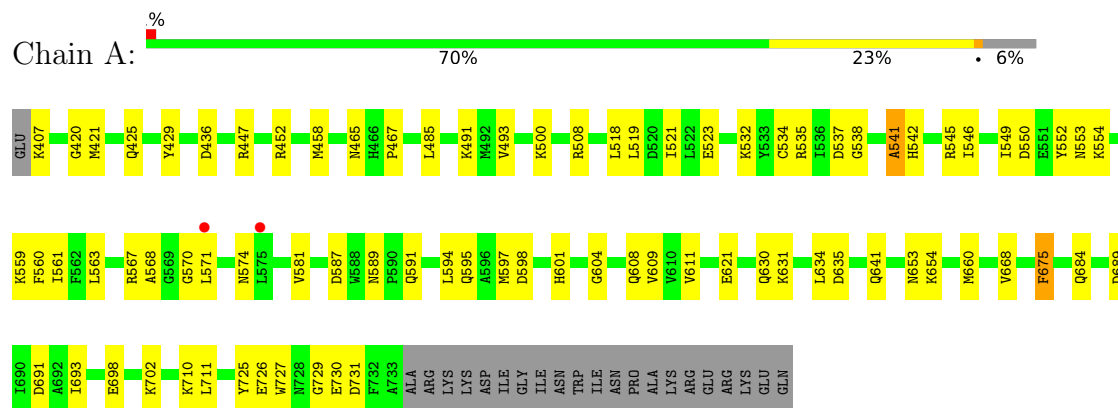
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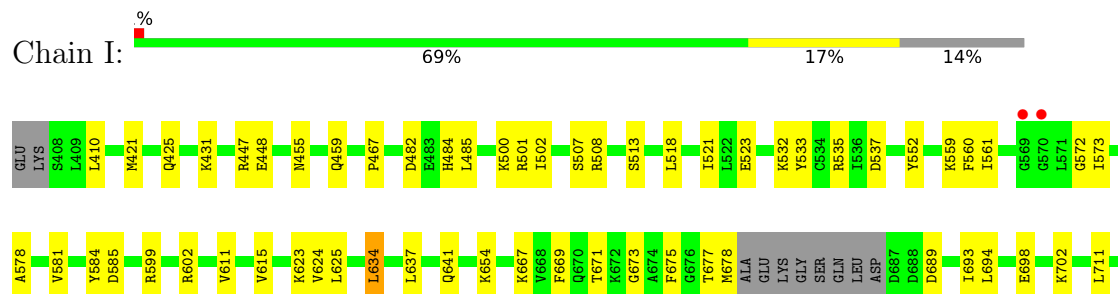
• Molecule 1: Chromatin-remodeling complex ATPase-like protein

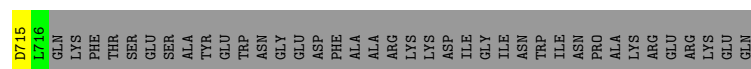


• Molecule 1: Chromatin-remodeling complex ATPase-like protein

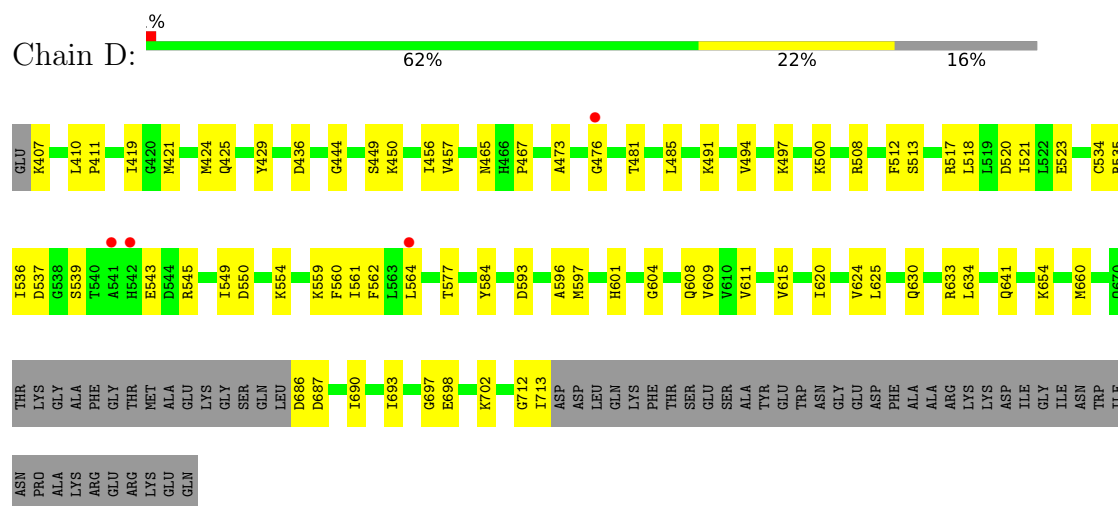


• Molecule 1: Chromatin-remodeling complex ATPase-like protein

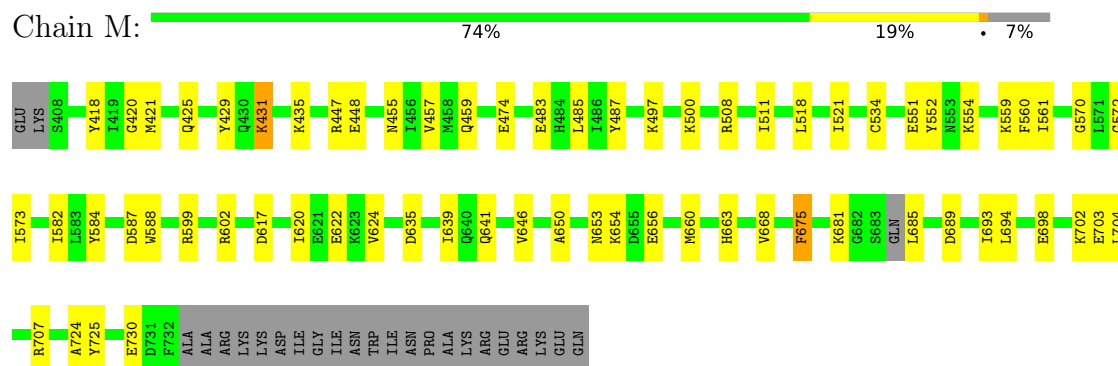




- Molecule 1: Chromatin-remodeling complex ATPase-like protein



- Molecule 1: Chromatin-remodeling complex ATPase-like protein



- Molecule 2: Histone H4

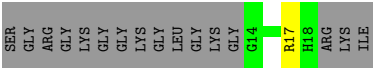


- Molecule 2: Histone H4

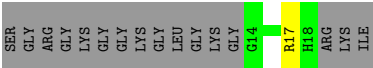


- Molecule 2: Histone H4





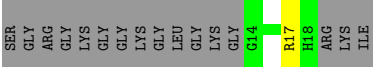
● Molecule 2: Histone H4



● Molecule 2: Histone H4



● Molecule 2: Histone H4



● Molecule 2: Histone H4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	102.24Å 119.32Å 132.50Å 89.59° 105.96° 93.08°	Depositor
Resolution (Å)	42.00 – 3.01 42.00 – 3.01	Depositor EDS
% Data completeness (in resolution range)	94.2 (42.00-3.01) 94.2 (42.00-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.218 , 0.276 0.219 , 0.277	Depositor DCC
R_{free} test set	5661 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	71.6	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	40399	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% 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	0.16	0/2637	0.46	0/3549
1	B	0.17	0/2639	0.44	0/3552
1	C	0.16	0/2613	0.43	1/3518 (0.0%)
1	D	0.19	0/2359	0.46	2/3174 (0.1%)
1	E	0.15	0/2520	0.42	2/3389 (0.1%)
1	F	0.13	0/2617	0.46	2/3521 (0.1%)
1	G	0.15	0/2474	0.43	2/3328 (0.1%)
1	H	0.15	0/2584	0.36	0/3472
1	I	0.11	0/2422	0.31	0/3259
1	J	0.12	0/2551	0.39	2/3432 (0.1%)
1	K	0.12	0/2461	0.35	0/3312
1	L	0.14	0/2499	0.45	2/3362 (0.1%)
1	M	0.15	0/2613	0.34	0/3516
1	N	0.11	0/2574	0.32	0/3465
1	O	0.11	0/2601	0.34	2/3500 (0.1%)
1	P	0.10	0/2554	0.33	2/3436 (0.1%)
2	Q	0.10	0/39	0.58	0/49
2	R	0.09	0/39	0.35	0/49
2	S	0.12	0/39	0.48	0/49
2	T	0.14	0/39	0.49	0/49
2	U	0.12	0/39	0.41	0/49
2	V	0.13	0/39	0.51	0/49
2	W	0.11	0/35	0.33	0/44
All	All	0.14	0/40987	0.40	17/55123 (0.0%)

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	1
1	H	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	O	0	1
All	All	0	3

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	476	GLY	CA-C-N	9.02	129.67	120.38
1	F	476	GLY	C-N-CA	9.02	129.67	120.38
1	J	476	GLY	CA-C-N	8.56	125.91	119.66
1	J	476	GLY	C-N-CA	8.56	125.91	119.66
1	L	476	GLY	CA-C-N	6.35	126.92	120.38
1	L	476	GLY	C-N-CA	6.35	126.92	120.38
1	D	476	GLY	CA-C-N	6.01	126.57	120.38
1	D	476	GLY	C-N-CA	6.01	126.57	120.38
1	G	476	GLY	CA-C-N	5.89	126.45	120.38
1	G	476	GLY	C-N-CA	5.89	126.45	120.38
1	C	682	GLY	N-CA-C	5.72	119.89	110.97
1	E	476	GLY	CA-C-N	5.61	126.16	120.38
1	E	476	GLY	C-N-CA	5.61	126.16	120.38
1	P	476	GLY	CA-C-N	5.33	125.87	120.38
1	P	476	GLY	C-N-CA	5.33	125.87	120.38
1	O	476	GLY	CA-C-N	5.04	125.57	120.38
1	O	476	GLY	C-N-CA	5.04	125.57	120.38

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	570	GLY	Peptide
1	H	675	PHE	Peptide
1	O	681	LYS	Peptide

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	2597	0	2607	72	0
1	B	2599	0	2603	57	0
1	C	2574	0	2579	56	0
1	D	2326	0	2364	57	0
1	E	2484	0	2521	50	0
1	F	2578	0	2580	63	0
1	G	2439	0	2476	56	0
1	H	2549	0	2553	64	0
1	I	2388	0	2422	43	0
1	J	2515	0	2532	46	0
1	K	2427	0	2456	55	0
1	L	2464	0	2499	58	0
1	M	2574	0	2580	47	0
1	N	2536	0	2539	46	0
1	O	2563	0	2571	62	0
1	P	2517	0	2522	46	0
2	Q	39	0	40	4	0
2	R	39	0	40	2	0
2	S	39	0	40	1	0
2	T	39	0	40	0	0
2	U	39	0	40	1	0
2	V	39	0	40	1	0
2	W	35	0	37	4	0
All	All	40399	0	40681	732	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 9.

All (732) 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:F:532:LYS:NZ	1:F:558:ASP:OD2	1.84	1.10
1:D:481:THR:HG22	1:M:447:ARG:HH12	1.19	1.04
1:M:681:LYS:NZ	1:M:689:ASP:OD2	1.91	1.03
1:N:481:THR:HG22	2:R:17:ARG:HH12	1.21	1.02
1:E:523:GLU:OE1	1:E:535:ARG:NH1	1.92	1.02
1:A:508:ARG:NH1	1:A:559:LYS:O	1.93	1.00
1:E:654:LYS:NZ	1:H:446:LYS:O	1.98	0.97
1:B:567:ARG:NH1	1:B:587:ASP:OD2	2.00	0.95
1:E:461:ARG:HH12	1:E:585:ASP:HB2	1.31	0.93
1:A:538:GLY:HA2	1:A:545:ARG:HH11	1.29	0.93
1:B:501:ARG:NH1	1:K:691:ASP:OD1	2.02	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:481:THR:HG22	1:N:447:ARG:HH12	1.33	0.91
1:J:570:GLY:O	1:J:599:ARG:NH1	2.06	0.88
1:C:535:ARG:NH2	1:C:537:ASP:OD2	2.06	0.87
1:B:598:ASP:OD2	1:B:727:TRP:HB2	1.75	0.87
1:F:481:THR:HG22	1:O:447:ARG:HH12	1.38	0.86
1:O:551:GLU:OE2	1:O:554:LYS:NZ	2.11	0.84
1:M:681:LYS:HD3	1:M:685:LEU:HD11	1.59	0.83
1:B:691:ASP:OD1	1:K:501:ARG:NH1	2.11	0.82
1:G:481:THR:HG22	1:P:447:ARG:HH12	1.44	0.81
1:D:497:LYS:NZ	1:D:500:LYS:HD2	1.96	0.81
1:E:432:ILE:HA	1:E:435:LYS:HE2	1.63	0.80
1:H:718:LYS:NZ	1:H:722:GLU:OE1	2.12	0.80
1:D:523:GLU:OE1	1:D:535:ARG:NH1	2.15	0.79
1:A:421:MET:HE2	1:A:425:GLN:HB3	1.65	0.79
1:N:572:GLY:HA2	1:N:602:ARG:HH12	1.47	0.78
1:G:523:GLU:OE1	1:G:535:ARG:NH1	2.15	0.78
1:E:421:MET:HE2	1:E:425:GLN:HB3	1.66	0.77
1:H:474:GLU:OE2	2:W:17:ARG:NH1	2.15	0.77
1:I:671:THR:HG22	1:I:673:GLY:H	1.50	0.77
1:A:538:GLY:HA2	1:A:545:ARG:NH1	2.00	0.76
1:B:535:ARG:NE	1:B:537:ASP:OD1	2.18	0.75
1:H:551:GLU:OE2	1:H:554:LYS:NZ	2.14	0.74
1:G:481:THR:HG22	1:P:447:ARG:NH1	2.01	0.74
1:A:542:HIS:CE1	1:A:546:ILE:HD11	2.23	0.74
1:N:421:MET:HE2	1:N:425:GLN:HB3	1.70	0.74
1:L:681:LYS:HD2	1:L:685:LEU:HD23	1.69	0.73
1:B:523:GLU:OE1	1:B:535:ARG:NH1	2.19	0.73
1:O:538:GLY:HA3	1:O:568:ALA:HB1	1.70	0.73
1:H:570:GLY:O	1:H:599:ARG:NH1	2.20	0.73
1:J:425:GLN:HG2	1:J:467:PRO:HG3	1.70	0.73
1:O:675:PHE:O	1:O:679:ALA:N	2.19	0.73
1:L:480:THR:O	1:L:484:HIS:NE2	2.18	0.73
1:A:598:ASP:OD2	1:A:727:TRP:HB2	1.90	0.72
1:D:421:MET:HE2	1:D:425:GLN:HB3	1.70	0.71
1:A:571:LEU:HD12	1:A:595:GLN:HE21	1.55	0.71
1:D:497:LYS:HZ2	1:D:500:LYS:HD2	1.55	0.71
1:N:446:LYS:O	1:L:654:LYS:NZ	2.23	0.70
1:G:562:PHE:CE2	1:G:564:LEU:HD21	2.26	0.70
1:E:427:LYS:HE2	1:E:431:LYS:NZ	2.07	0.70
1:A:535:ARG:NE	1:A:537:ASP:OD1	2.24	0.70
1:M:421:MET:HE2	1:M:425:GLN:HB3	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:712:GLY:HA2	1:I:637:LEU:HD21	1.74	0.69
1:O:508:ARG:NH2	1:O:552:TYR:O	2.21	0.69
1:F:591:GLN:NE2	1:F:726:GLU:O	2.26	0.69
1:L:421:MET:HE2	1:L:425:GLN:HB3	1.75	0.69
1:C:421:MET:HE2	1:C:425:GLN:HB3	1.74	0.69
1:F:598:ASP:OD2	1:F:727:TRP:HB2	1.93	0.68
1:P:421:MET:HE2	1:P:425:GLN:HB3	1.76	0.68
1:C:598:ASP:OD2	1:C:727:TRP:HB2	1.93	0.68
1:F:481:THR:HG22	1:O:447:ARG:NH1	2.09	0.68
1:K:481:THR:CG2	2:Q:17:ARG:HH12	2.07	0.67
1:P:675:PHE:HB3	1:D:421:MET:H	1.59	0.67
1:N:624:VAL:HG22	1:L:660:MET:HE3	1.77	0.67
1:H:465:ASN:HB3	1:H:518:LEU:HD13	1.75	0.67
1:K:572:GLY:H	1:K:599:ARG:HH12	1.42	0.67
1:G:660:MET:HE3	1:M:624:VAL:HG22	1.75	0.66
1:D:485:LEU:HD11	1:D:521:ILE:HG23	1.76	0.66
1:B:485:LEU:HD11	1:B:521:ILE:HG23	1.77	0.66
1:F:523:GLU:OE1	1:F:535:ARG:NH1	2.27	0.66
1:B:418:TYR:HB3	1:B:617:ASP:HB2	1.78	0.66
1:M:508:ARG:NH2	1:M:552:TYR:O	2.20	0.66
1:B:439:ALA:HB2	1:B:452:ARG:NH1	2.11	0.65
1:I:421:MET:HE2	1:I:425:GLN:HB3	1.78	0.65
1:K:421:MET:HE2	1:K:425:GLN:HB3	1.77	0.65
1:E:501:ARG:NH1	1:H:691:ASP:OD1	2.30	0.65
1:F:427:LYS:HE2	1:F:431:LYS:HZ3	1.62	0.65
1:O:465:ASN:HB3	1:O:518:LEU:HD13	1.79	0.64
1:P:698:GLU:HG2	1:P:702:LYS:HE3	1.78	0.64
1:H:698:GLU:HG2	1:H:702:LYS:HE3	1.78	0.64
1:K:686:ASP:N	1:K:689:ASP:OD2	2.30	0.64
1:L:438:ASP:OD2	1:L:452:ARG:NH2	2.30	0.64
1:B:615:VAL:HG21	1:B:625:LEU:HD22	1.80	0.64
1:E:660:MET:HE3	1:H:624:VAL:HG22	1.79	0.64
1:O:421:MET:HE2	1:O:425:GLN:HB3	1.80	0.64
1:H:524:ASP:OD2	2:W:17:ARG:HG3	1.97	0.64
1:K:523:GLU:OE1	1:K:535:ARG:NH1	2.31	0.64
1:F:524:ASP:OD2	1:O:447:ARG:HG3	1.98	0.63
1:I:572:GLY:H	1:I:599:ARG:HH12	1.46	0.63
1:F:508:ARG:NH2	1:F:577:THR:OG1	2.32	0.63
1:N:481:THR:HG22	2:R:17:ARG:NH1	2.04	0.63
1:I:523:GLU:OE2	1:I:533:TYR:OH	2.12	0.63
1:B:574:ASN:OD1	1:B:576:THR:HG23	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:ASP:HB3	1:A:554:LYS:HE3	1.81	0.63
1:N:698:GLU:HG2	1:N:702:LYS:HE3	1.80	0.63
1:B:421:MET:HE2	1:B:425:GLN:HB3	1.81	0.62
1:A:508:ARG:NH2	1:A:552:TYR:O	2.30	0.62
1:B:572:GLY:HA2	1:B:602:ARG:HH12	1.64	0.62
1:G:474:GLU:OE1	1:P:447:ARG:NH2	2.33	0.62
1:G:536:ILE:HD11	1:G:548:ALA:HB3	1.81	0.62
1:A:727:TRP:CD1	1:A:729:GLY:H	2.18	0.62
1:G:653:ASN:OD1	1:G:654:LYS:N	2.33	0.62
1:C:485:LEU:HD11	1:C:521:ILE:HG23	1.82	0.62
1:L:432:ILE:HA	1:L:435:LYS:HE3	1.82	0.62
1:B:542:HIS:HA	1:B:545:ARG:HB2	1.81	0.61
1:J:722:GLU:OE1	1:H:589:ASN:ND2	2.32	0.61
1:F:532:LYS:HB3	1:F:559:LYS:HD3	1.82	0.61
1:H:421:MET:HE2	1:H:425:GLN:HB3	1.82	0.61
1:C:572:GLY:HA2	1:C:602:ARG:HH12	1.63	0.61
1:H:508:ARG:NH2	1:H:552:TYR:O	2.26	0.61
1:K:572:GLY:H	1:K:599:ARG:NH1	1.98	0.61
1:N:551:GLU:OE2	1:N:554:LYS:NZ	2.32	0.61
1:B:581:VAL:HB	1:B:611:VAL:HG22	1.83	0.61
1:G:421:MET:HE2	1:G:425:GLN:HB3	1.83	0.61
1:L:536:ILE:HB	1:L:564:LEU:HD23	1.82	0.61
1:A:591:GLN:NE2	1:A:726:GLU:O	2.26	0.61
1:B:411:PRO:O	1:B:608:GLN:NE2	2.26	0.60
1:N:508:ARG:NH1	1:N:559:LYS:O	2.33	0.60
1:F:501:ARG:HH12	1:I:694:LEU:HD12	1.65	0.60
1:J:624:VAL:HG22	1:C:660:MET:HE3	1.83	0.60
1:B:532:LYS:HB3	1:B:559:LYS:HD3	1.84	0.60
1:I:535:ARG:NE	1:I:537:ASP:OD1	2.31	0.60
1:G:550:ASP:O	1:G:554:LYS:HG3	2.02	0.60
1:O:487:TYR:HE1	1:A:684:GLN:HB2	1.66	0.60
1:A:485:LEU:HD11	1:A:521:ILE:HG23	1.83	0.60
1:G:690:ILE:HG13	1:M:497:LYS:HG2	1.82	0.60
1:N:421:MET:H	1:L:675:PHE:HB3	1.67	0.60
1:P:573:ILE:O	1:P:602:ARG:NH1	2.35	0.59
1:A:538:GLY:HA3	1:A:568:ALA:HB1	1.84	0.59
1:J:421:MET:HE2	1:J:425:GLN:HB3	1.84	0.59
1:A:571:LEU:HD13	1:A:595:GLN:HG2	1.83	0.59
1:K:481:THR:HG22	2:Q:17:ARG:HH12	1.66	0.59
1:D:444:GLY:C	1:D:449:SER:HB2	2.26	0.59
1:F:450:LYS:HB3	1:I:654:LYS:HG2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:ASP:OD2	1:A:452:ARG:NE	2.36	0.59
1:N:675:PHE:HB2	1:L:421:MET:H	1.67	0.59
1:O:535:ARG:NE	1:O:537:ASP:OD1	2.31	0.59
1:I:425:GLN:HG2	1:I:467:PRO:HG3	1.85	0.59
1:E:490:GLY:HA2	1:H:678:MET:SD	2.42	0.59
1:L:675:PHE:O	1:L:679:ALA:N	2.32	0.59
1:E:535:ARG:NE	1:E:537:ASP:OD1	2.36	0.59
1:G:449:SER:HB2	1:M:654:LYS:NZ	2.18	0.58
1:P:656:GLU:OE2	1:P:718:LYS:HE3	2.03	0.58
1:K:581:VAL:HB	1:K:611:VAL:HG22	1.85	0.58
1:K:653:ASN:OD1	1:K:654:LYS:N	2.37	0.58
1:E:685:LEU:HD11	1:H:494:VAL:HA	1.84	0.58
1:N:729:GLY:H	1:K:567:ARG:HH21	1.50	0.58
1:O:578:ALA:O	1:O:605:GLN:NE2	2.35	0.58
1:E:481:THR:HG22	1:N:447:ARG:NH1	2.12	0.58
1:J:698:GLU:HG2	1:J:702:LYS:HE3	1.85	0.58
1:J:425:GLN:NE2	1:J:463:CYS:O	2.37	0.58
1:C:413:LYS:HE2	1:C:415:ILE:HD11	1.85	0.58
1:C:698:GLU:HG2	1:C:702:LYS:HE3	1.85	0.58
1:H:485:LEU:HD11	1:H:521:ILE:HG23	1.84	0.58
1:K:480:THR:O	1:K:484:HIS:NE2	2.30	0.58
1:E:427:LYS:HE2	1:E:431:LYS:HZ3	1.67	0.58
1:F:719:PHE:HD1	1:A:541:ALA:HA	1.69	0.58
1:K:572:GLY:N	1:K:599:ARG:HH12	2.02	0.58
1:K:597:MET:HE2	1:K:611:VAL:HG11	1.86	0.58
1:L:482:ASP:HB2	1:L:484:HIS:HD2	1.69	0.58
1:N:425:GLN:HG2	1:N:467:PRO:HG3	1.86	0.58
1:K:685:LEU:HB3	1:K:689:ASP:OD2	2.04	0.58
1:M:572:GLY:C	1:M:602:ARG:HH12	2.12	0.58
1:B:497:LYS:HE2	1:K:687:ASP:HB3	1.86	0.57
1:F:421:MET:HE2	1:F:425:GLN:HB3	1.86	0.57
1:P:465:ASN:HB3	1:P:518:LEU:HD13	1.85	0.57
1:P:687:ASP:O	1:P:690:ILE:HG22	2.04	0.57
1:H:474:GLU:CD	2:W:17:ARG:HH12	2.11	0.57
1:L:518:LEU:HD11	1:L:584:TYR:CE1	2.39	0.57
1:D:518:LEU:HD21	1:D:584:TYR:CD1	2.38	0.57
1:O:687:ASP:HA	1:O:690:ILE:HG22	1.85	0.57
1:F:591:GLN:HG2	1:F:728:ASN:HB2	1.85	0.57
1:N:570:GLY:O	1:N:599:ARG:NH2	2.37	0.57
1:K:698:GLU:HG2	1:K:702:LYS:HE3	1.87	0.57
1:J:436:ASP:OD1	1:J:452:ARG:NH2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:534:CYS:HB2	1:H:562:PHE:CD1	2.39	0.57
1:D:534:CYS:HB2	1:D:562:PHE:CD1	2.40	0.56
1:L:508:ARG:NH2	1:L:577:THR:OG1	2.37	0.56
1:G:536:ILE:HB	1:G:564:LEU:HD22	1.86	0.56
1:B:438:ASP:HB2	1:B:452:ARG:HH22	1.70	0.56
1:F:485:LEU:HD11	1:F:521:ILE:HG23	1.87	0.56
1:F:501:ARG:NH1	1:I:694:LEU:HD12	2.20	0.56
1:I:581:VAL:HB	1:I:611:VAL:HG22	1.88	0.56
1:P:623:LYS:HG2	1:D:660:MET:HA	1.87	0.56
1:K:664:GLY:HA2	1:K:667:LYS:HE2	1.87	0.56
1:F:427:LYS:HE2	1:F:431:LYS:NZ	2.21	0.56
1:N:565:THR:OG1	1:N:568:ALA:HB2	2.05	0.55
1:B:721:SER:HB3	1:C:542:HIS:CE1	2.41	0.55
1:P:439:ALA:HB2	1:P:452:ARG:HE	1.71	0.55
1:C:633:ARG:HH12	1:C:714:ASP:CG	2.15	0.55
1:L:550:ASP:O	1:L:554:LYS:HG3	2.07	0.55
1:A:581:VAL:HB	1:A:611:VAL:HG22	1.88	0.55
1:G:615:VAL:HG21	1:G:625:LEU:HD22	1.88	0.55
1:G:663:HIS:HE1	1:M:622:GLU:OE2	1.89	0.55
1:H:534:CYS:SG	1:H:559:LYS:HG3	2.47	0.55
1:O:678:MET:HE1	1:A:493:VAL:HG21	1.89	0.55
1:C:538:GLY:HA2	1:C:545:ARG:NH1	2.21	0.55
1:H:576:THR:O	1:H:605:GLN:NE2	2.39	0.55
1:A:436:ASP:OD2	1:A:452:ARG:NH2	2.38	0.55
1:D:411:PRO:O	1:D:608:GLN:NE2	2.35	0.55
1:G:542:HIS:CE1	1:G:546:ILE:HD11	2.41	0.55
1:N:517:ARG:HA	1:N:520:ASP:OD2	2.07	0.55
1:F:418:TYR:HB3	1:F:617:ASP:HB2	1.89	0.55
1:G:519:LEU:HD22	1:G:563:LEU:HB3	1.89	0.55
1:C:645:GLN:HG3	1:H:720:THR:HG21	1.88	0.55
1:C:523:GLU:OE1	1:C:535:ARG:NH1	2.36	0.55
1:A:594:LEU:HD21	1:A:725:TYR:C	2.32	0.55
1:F:535:ARG:NE	1:F:537:ASP:OD1	2.27	0.54
1:G:677:THR:HB	1:M:420:GLY:H	1.72	0.54
1:E:550:ASP:O	1:E:554:LYS:HG3	2.08	0.54
1:O:501:ARG:NH1	1:A:691:ASP:OD1	2.41	0.54
1:P:687:ASP:HB3	1:D:497:LYS:HE2	1.90	0.54
1:H:565:THR:OG1	1:H:568:ALA:HB2	2.08	0.54
1:E:523:GLU:CD	1:E:535:ARG:HH11	2.10	0.54
1:G:536:ILE:HD12	1:G:545:ARG:HA	1.89	0.54
1:O:421:MET:H	1:A:675:PHE:HB3	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:545:ARG:O	1:C:549:ILE:HG12	2.07	0.54
1:A:552:TYR:HD1	1:A:559:LYS:HB2	1.72	0.54
1:E:461:ARG:NH1	1:E:585:ASP:HB2	2.12	0.54
1:G:698:GLU:HG2	1:G:702:LYS:HE3	1.90	0.54
1:C:538:GLY:HA2	1:C:545:ARG:HH11	1.73	0.54
1:L:698:GLU:HG2	1:L:702:LYS:HE3	1.89	0.54
1:N:421:MET:HB3	1:N:425:GLN:HB2	1.89	0.54
1:O:424:MET:HE2	1:O:467:PRO:HB2	1.91	0.53
1:C:508:ARG:NH2	1:C:552:TYR:O	2.30	0.53
1:C:518:LEU:HD11	1:C:584:TYR:CE1	2.43	0.53
1:A:553:ASN:ND2	1:A:574:ASN:O	2.40	0.53
1:E:653:ASN:OD1	1:E:656:GLU:HB2	2.08	0.53
1:O:623:LYS:HG2	1:A:660:MET:HA	1.91	0.53
1:O:485:LEU:HD11	1:O:521:ILE:HG23	1.91	0.53
1:L:542:HIS:CE1	1:L:546:ILE:HD11	2.43	0.53
1:P:709:GLU:HA	1:D:633:ARG:HD3	1.89	0.53
1:O:417:VAL:HG23	1:A:693:ILE:HG22	1.91	0.53
1:A:567:ARG:NE	1:A:587:ASP:OD2	2.42	0.53
1:H:425:GLN:HG2	1:H:467:PRO:HG3	1.90	0.53
1:M:483:GLU:HB3	1:M:487:TYR:CE2	2.44	0.53
1:O:487:TYR:CE1	1:A:684:GLN:HB2	2.44	0.53
1:B:450:LYS:HD2	1:K:654:LYS:HE2	1.91	0.53
1:J:508:ARG:NH2	1:J:552:TYR:O	2.34	0.53
1:D:550:ASP:O	1:D:554:LYS:HG3	2.08	0.53
1:M:431:LYS:HG2	1:M:435:LYS:HE3	1.90	0.53
1:G:521:ILE:HA	1:P:447:ARG:HG2	1.91	0.53
1:J:615:VAL:HG21	1:J:625:LEU:HD22	1.91	0.53
1:N:654:LYS:NZ	1:L:443:ALA:HB1	2.24	0.53
1:K:572:GLY:CA	1:K:599:ARG:HH12	2.22	0.53
1:B:700:ARG:HH12	1:K:617:ASP:CG	2.16	0.53
1:N:485:LEU:HD11	1:N:521:ILE:HG23	1.90	0.53
1:O:570:GLY:O	1:O:599:ARG:NH2	2.42	0.53
1:O:597:MET:HE2	1:O:611:VAL:HG11	1.90	0.53
1:B:700:ARG:NH1	1:K:617:ASP:OD2	2.32	0.52
1:P:418:TYR:HB3	1:P:617:ASP:HB2	1.92	0.52
1:I:615:VAL:HG21	1:I:625:LEU:HD22	1.90	0.52
1:G:563:LEU:O	1:G:564:LEU:HD23	2.09	0.52
1:N:572:GLY:HA3	1:K:711:LEU:HD21	1.91	0.52
1:M:518:LEU:HD21	1:M:584:TYR:CD1	2.44	0.52
1:E:634:LEU:HD11	1:H:716:LEU:HB3	1.90	0.52
1:J:458:MET:HE3	1:H:731:ASP:HB3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:481:THR:HG22	2:S:17:ARG:HH12	1.74	0.52
1:P:595:GLN:O	1:P:599:ARG:NE	2.43	0.52
1:H:689:ASP:O	1:H:693:ILE:HG13	2.09	0.52
1:A:597:MET:HE2	1:A:611:VAL:HG11	1.91	0.52
1:D:481:THR:HG22	1:M:447:ARG:NH1	2.04	0.52
1:F:491:LYS:HZ1	1:F:621:GLU:CD	2.17	0.52
1:O:419:ILE:HD12	1:O:491:LYS:HG2	1.92	0.52
1:O:643:ARG:HH21	1:O:717:GLN:HG2	1.74	0.52
1:D:686:ASP:O	1:D:690:ILE:N	2.39	0.52
1:B:535:ARG:HA	1:B:563:LEU:O	2.09	0.52
1:C:645:GLN:O	1:C:649:LYS:HG2	2.10	0.52
1:H:521:ILE:HA	2:W:17:ARG:HG2	1.92	0.52
1:D:698:GLU:HG2	1:D:702:LYS:HE3	1.92	0.52
1:B:482:ASP:H	1:B:484:HIS:CE1	2.28	0.52
1:G:508:ARG:NH1	1:G:557:SER:HB3	2.24	0.52
1:J:420:GLY:H	1:C:677:THR:HB	1.74	0.52
1:O:656:GLU:OE2	1:O:718:LYS:HD2	2.09	0.52
1:P:508:ARG:HB2	1:P:578:ALA:HA	1.92	0.52
1:C:455:ASN:O	1:C:459:GLN:HG2	2.10	0.52
1:P:508:ARG:NH2	1:P:552:TYR:O	2.25	0.52
1:I:502:ILE:HG23	1:I:507:SER:HB2	1.91	0.52
1:E:425:GLN:OE1	1:E:490:GLY:N	2.41	0.51
1:E:439:ALA:HB2	1:E:452:ARG:NH1	2.26	0.51
1:C:567:ARG:HH11	1:C:734:ALA:HB2	1.73	0.51
1:E:485:LEU:HD11	1:E:521:ILE:HG23	1.92	0.51
1:O:601:HIS:HB2	1:O:609:VAL:HG11	1.91	0.51
1:N:654:LYS:HZ1	1:L:443:ALA:HB1	1.74	0.51
1:F:581:VAL:HB	1:F:611:VAL:HG22	1.91	0.51
1:O:586:SER:OG	1:O:613:ARG:NH1	2.36	0.51
1:I:508:ARG:NH2	1:I:552:TYR:O	2.23	0.51
1:I:572:GLY:H	1:I:599:ARG:NH1	2.09	0.51
1:E:545:ARG:O	1:E:549:ILE:HG12	2.10	0.51
1:F:532:LYS:HB2	1:F:559:LYS:HA	1.91	0.51
1:J:515:MET:HA	1:H:730:GLU:HG2	1.93	0.51
1:C:532:LYS:NZ	1:C:558:ASP:OD2	2.43	0.51
1:O:508:ARG:NH1	1:O:559:LYS:O	2.42	0.51
1:P:485:LEU:HD11	1:P:521:ILE:HG23	1.93	0.51
1:D:508:ARG:NH2	1:D:577:THR:OG1	2.42	0.51
1:D:597:MET:HE2	1:D:611:VAL:HG11	1.92	0.51
1:C:532:LYS:HB3	1:C:559:LYS:HD3	1.92	0.51
1:A:601:HIS:HB2	1:A:609:VAL:HG11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:648:ALA:HB2	1:I:634:LEU:HB3	1.93	0.51
1:O:641:GLN:HB2	1:A:641:GLN:HG3	1.92	0.51
1:I:523:GLU:OE1	1:I:535:ARG:NH1	2.38	0.51
1:D:536:ILE:HG12	1:D:562:PHE:CZ	2.46	0.51
1:E:461:ARG:NH2	1:E:585:ASP:OD2	2.40	0.51
1:F:713:ILE:HG13	1:I:641:GLN:HG3	1.93	0.51
1:M:418:TYR:HB3	1:M:617:ASP:HB2	1.93	0.51
1:E:574:ASN:HD22	1:L:603:ILE:HD12	1.76	0.50
1:B:677:THR:HB	1:K:420:GLY:H	1.76	0.50
1:P:455:ASN:O	1:P:459:GLN:HG2	2.11	0.50
1:P:597:MET:HE2	1:P:611:VAL:HG11	1.93	0.50
1:H:512:PHE:CE2	1:H:564:LEU:HD13	2.45	0.50
1:L:411:PRO:O	1:L:608:GLN:NE2	2.38	0.50
1:L:710:LYS:HA	1:L:710:LYS:NZ	2.26	0.50
1:F:591:GLN:HE21	1:F:728:ASN:N	2.08	0.50
1:J:627:ARG:O	1:J:631:LYS:HG2	2.11	0.50
1:C:601:HIS:HB2	1:C:609:VAL:HG11	1.92	0.50
1:A:534:CYS:SG	1:A:559:LYS:HG3	2.51	0.50
1:F:573:ILE:HG13	1:F:599:ARG:HD3	1.93	0.50
1:J:711:LEU:HD21	1:H:572:GLY:HA3	1.92	0.50
1:O:573:ILE:CD1	1:O:599:ARG:HH11	2.25	0.50
1:C:637:LEU:HD22	1:C:713:ILE:HD12	1.93	0.50
1:H:671:THR:HG21	1:H:674:ALA:O	2.11	0.50
1:A:730:GLU:HG2	1:A:731:ASP:H	1.77	0.50
1:E:711:LEU:HD13	1:E:719:PHE:CD2	2.47	0.50
1:F:497:LYS:NZ	1:F:500:LYS:HD2	2.26	0.50
1:P:417:VAL:HG13	1:D:693:ILE:HG21	1.94	0.50
1:H:674:ALA:HB3	1:H:679:ALA:HB2	1.93	0.50
1:I:455:ASN:O	1:I:459:GLN:HG2	2.11	0.50
1:H:703:GLU:OE1	1:H:707:ARG:NH2	2.45	0.50
1:F:601:HIS:HB2	1:F:609:VAL:HG11	1.94	0.50
1:K:425:GLN:HG2	1:K:467:PRO:HG3	1.92	0.50
1:F:493:VAL:HG21	1:I:678:MET:HE1	1.93	0.50
1:H:534:CYS:HB2	1:H:562:PHE:CE1	2.47	0.50
1:L:542:HIS:O	1:L:546:ILE:HG13	2.11	0.50
1:I:532:LYS:HB3	1:I:559:LYS:HD3	1.93	0.50
1:D:497:LYS:HZ3	1:D:500:LYS:HD2	1.72	0.50
1:E:541:ALA:O	1:E:545:ARG:HG3	2.11	0.49
1:I:698:GLU:HG2	1:I:702:LYS:HE3	1.94	0.49
1:N:411:PRO:O	1:N:608:GLN:NE2	2.37	0.49
1:O:572:GLY:HA2	1:O:602:ARG:HH12	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:572:GLY:N	1:I:599:ARG:HH12	2.09	0.49
1:P:497:LYS:NZ	1:P:500:LYS:HD2	2.28	0.49
1:L:534:CYS:SG	1:L:559:LYS:HG3	2.52	0.49
1:I:518:LEU:HD21	1:I:584:TYR:CD1	2.47	0.49
1:G:542:HIS:HA	1:G:545:ARG:HE	1.78	0.49
1:K:555:PRO:HG2	1:A:608:GLN:HB2	1.94	0.49
1:E:662:GLN:NE2	1:H:440:VAL:O	2.36	0.49
1:O:652:ALA:HB1	1:O:657:LEU:HG	1.93	0.49
1:K:549:ILE:HD11	1:K:573:ILE:HD12	1.93	0.49
1:P:425:GLN:HG2	1:P:467:PRO:HG3	1.93	0.49
1:C:458:MET:HE1	1:L:448:GLU:HA	1.94	0.49
1:L:407:LYS:HE3	1:L:603:ILE:HG23	1.95	0.49
1:F:430:GLN:HG2	1:I:669:PHE:HD1	1.78	0.49
1:F:719:PHE:CD1	1:A:541:ALA:HA	2.46	0.49
1:C:581:VAL:HB	1:C:611:VAL:HG22	1.95	0.49
1:G:446:LYS:NZ	1:M:724:ALA:HA	2.27	0.49
1:J:474:GLU:OE1	2:U:17:ARG:NH2	2.41	0.49
1:N:455:ASN:O	1:N:459:GLN:HG2	2.13	0.49
1:I:410:LEU:HD12	1:I:410:LEU:H	1.77	0.49
1:M:485:LEU:HD11	1:M:521:ILE:HG23	1.93	0.49
1:M:570:GLY:O	1:M:599:ARG:NH2	2.24	0.49
1:F:487:TYR:HD1	1:F:493:VAL:HG22	1.78	0.48
1:O:534:CYS:SG	1:O:559:LYS:HG3	2.52	0.48
1:G:512:PHE:CE2	1:G:564:LEU:HD12	2.48	0.48
1:G:581:VAL:HB	1:G:611:VAL:HG22	1.95	0.48
1:O:643:ARG:HE	1:O:717:GLN:HG2	1.78	0.48
1:A:630:GLN:O	1:A:634:LEU:HG	2.13	0.48
1:D:630:GLN:O	1:D:634:LEU:HG	2.13	0.48
1:D:693:ILE:O	1:D:697:GLY:N	2.30	0.48
1:F:479:TYR:CD2	1:O:448:GLU:HB3	2.48	0.48
1:G:508:ARG:HH12	1:G:557:SER:HB3	1.76	0.48
1:O:678:MET:O	1:O:682:GLY:HA3	2.13	0.48
1:M:560:PHE:CD2	1:M:561:ILE:HG13	2.48	0.48
1:G:519:LEU:HD23	1:G:522:LEU:HD12	1.94	0.48
1:G:637:LEU:HD22	1:G:713:ILE:HD12	1.94	0.48
1:E:581:VAL:HB	1:E:611:VAL:HG22	1.95	0.48
1:K:636:GLN:O	1:K:640:GLN:HB2	2.12	0.48
1:O:483:GLU:HB3	1:O:487:TYR:HE2	1.79	0.48
1:C:480:THR:O	1:C:484:HIS:NE2	2.43	0.48
1:H:427:LYS:HE2	1:H:431:LYS:HZ3	1.78	0.48
1:M:573:ILE:HG12	1:M:599:ARG:HD2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:698:GLU:HG2	1:F:702:LYS:HE3	1.96	0.48
1:G:653:ASN:HB3	1:G:656:GLU:HG3	1.94	0.48
1:P:517:ARG:HA	1:P:520:ASP:OD2	2.14	0.48
1:B:436:ASP:OD2	1:B:452:ARG:NH2	2.46	0.48
1:H:674:ALA:C	1:H:679:ALA:HB2	2.39	0.48
1:N:421:MET:H	1:L:675:PHE:CB	2.27	0.48
1:A:425:GLN:HG2	1:A:467:PRO:HG3	1.95	0.48
1:M:653:ASN:HB2	1:M:656:GLU:HG3	1.96	0.48
1:G:545:ARG:O	1:G:549:ILE:HG12	2.14	0.47
1:J:571:LEU:O	1:H:707:ARG:HD3	2.14	0.47
1:O:668:VAL:HG11	1:A:429:TYR:CD1	2.49	0.47
1:P:581:VAL:HB	1:P:611:VAL:HG22	1.96	0.47
1:K:474:GLU:OE1	2:Q:17:ARG:NH2	2.39	0.47
1:B:514:GLN:HB2	1:B:585:ASP:OD2	2.15	0.47
1:B:533:TYR:HA	1:B:561:ILE:O	2.14	0.47
1:E:518:LEU:HD21	1:E:584:TYR:CD1	2.49	0.47
1:G:501:ARG:HH12	1:M:694:LEU:HD12	1.78	0.47
1:J:641:GLN:HG3	1:C:713:ILE:HG21	1.96	0.47
1:C:573:ILE:HB	1:C:599:ARG:HH11	1.79	0.47
1:F:654:LYS:NZ	1:I:448:GLU:HB3	2.29	0.47
1:N:730:GLU:OE1	1:K:516:SER:N	2.44	0.47
1:D:534:CYS:HG	1:D:559:LYS:HE2	1.79	0.47
1:E:685:LEU:HG	1:H:493:VAL:HG12	1.96	0.47
1:F:721:SER:HB3	1:A:542:HIS:CE1	2.49	0.47
1:E:660:MET:HE1	1:H:588:TRP:CZ3	2.49	0.47
1:O:615:VAL:HG21	1:O:625:LEU:HD22	1.95	0.47
1:L:598:ASP:O	1:L:602:ARG:HD2	2.14	0.47
1:G:654:LYS:HE2	1:M:448:GLU:O	2.15	0.47
1:J:565:THR:OG1	1:J:568:ALA:HB2	2.13	0.47
1:J:634:LEU:HB3	1:C:648:ALA:HB2	1.97	0.47
1:J:718:LYS:NZ	1:J:722:GLU:HG2	2.30	0.47
1:A:523:GLU:OE1	1:A:535:ARG:NH1	2.33	0.47
1:G:626:GLU:HG2	1:M:704:LEU:HD11	1.97	0.47
1:J:518:LEU:HD21	1:J:584:TYR:CD1	2.50	0.47
1:N:656:GLU:OE2	1:N:718:LYS:HE3	2.15	0.47
1:H:482:ASP:CG	1:H:483:GLU:H	2.21	0.47
1:A:595:GLN:NE2	1:A:730:GLU:O	2.48	0.47
1:O:483:GLU:OE1	1:O:529:ARG:NH2	2.48	0.47
1:G:421:MET:H	1:M:675:PHE:HB3	1.79	0.47
1:E:634:LEU:HD12	1:H:716:LEU:HD13	1.96	0.46
1:H:447:ARG:HH12	1:L:481:THR:CG2	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:491:LYS:HD2	1:E:584:TYR:CE1	2.50	0.46
1:L:425:GLN:HG2	1:L:467:PRO:HG3	1.97	0.46
1:E:418:TYR:HB3	1:E:617:ASP:HB2	1.96	0.46
1:G:485:LEU:HD11	1:G:521:ILE:HG23	1.97	0.46
1:G:536:ILE:HG23	1:G:545:ARG:HG2	1.98	0.46
1:O:448:GLU:H	1:O:448:GLU:HG2	1.53	0.46
1:O:573:ILE:HG12	1:O:599:ARG:HH11	1.80	0.46
1:I:560:PHE:CD2	1:I:561:ILE:HG13	2.50	0.46
1:O:572:GLY:C	1:O:602:ARG:HH12	2.23	0.46
1:G:449:SER:HB2	1:M:654:LYS:HZ3	1.79	0.46
1:O:483:GLU:HB3	1:O:487:TYR:CE2	2.51	0.46
1:K:518:LEU:HD11	1:K:584:TYR:CE1	2.51	0.46
1:B:410:LEU:HD12	1:B:410:LEU:H	1.81	0.46
1:H:545:ARG:O	1:H:549:ILE:HG12	2.15	0.46
1:K:486:ILE:HD12	1:K:529:ARG:HH21	1.81	0.46
1:L:560:PHE:CD2	1:L:561:ILE:HG13	2.51	0.46
1:A:436:ASP:CG	1:A:452:ARG:HE	2.24	0.46
1:A:491:LYS:NZ	1:A:621:GLU:OE2	2.45	0.46
1:A:595:GLN:CD	1:A:727:TRP:HE1	2.23	0.46
1:E:425:GLN:HG2	1:E:467:PRO:HG3	1.97	0.46
1:F:485:LEU:HD22	1:O:447:ARG:NH1	2.31	0.46
1:J:573:ILE:HG12	1:J:599:ARG:HH11	1.80	0.46
1:C:686:ASP:OD2	1:C:689:ASP:N	2.47	0.46
1:L:419:ILE:HD13	1:L:491:LYS:HA	1.97	0.46
1:A:500:LYS:HE3	1:A:500:LYS:HB2	1.71	0.46
1:D:593:ASP:O	1:D:597:MET:HG3	2.15	0.46
1:G:542:HIS:HA	1:G:545:ARG:HH11	1.81	0.46
1:O:429:TYR:CD1	1:A:668:VAL:HG11	2.50	0.46
1:G:619:ALA:O	1:M:663:HIS:HD2	1.99	0.46
1:J:581:VAL:HB	1:J:611:VAL:HG22	1.98	0.46
1:L:678:MET:HB3	1:L:678:MET:HE3	1.81	0.46
1:L:581:VAL:HB	1:L:611:VAL:HG22	1.97	0.46
1:A:549:ILE:HG22	1:A:553:ASN:ND2	2.32	0.46
1:D:410:LEU:HD12	1:D:410:LEU:H	1.81	0.46
1:D:535:ARG:NE	1:D:537:ASP:OD1	2.41	0.46
1:N:444:GLY:HA3	1:N:448:GLU:OE1	2.17	0.45
1:N:414:GLU:OE2	1:N:597:MET:HE1	2.16	0.45
1:P:590:PRO:HG3	1:P:628:ALA:HB1	1.97	0.45
1:P:536:ILE:HD13	1:P:545:ARG:HB2	1.98	0.45
1:P:660:MET:HE3	1:D:624:VAL:HG22	1.99	0.45
1:M:551:GLU:OE2	1:M:554:LYS:NZ	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:654:LYS:HZ2	1:J:658:LEU:HD22	1.82	0.45
1:G:536:ILE:CD1	1:G:545:ARG:HA	2.47	0.45
1:J:500:LYS:HB2	1:J:500:LYS:HE3	1.72	0.45
1:J:721:SER:HB2	1:H:631:LYS:HD3	1.98	0.45
1:N:500:LYS:HE3	1:N:500:LYS:HB2	1.75	0.45
1:L:432:ILE:HD11	1:L:470:PHE:CZ	2.51	0.45
1:A:407:LYS:HZ1	1:A:604:GLY:N	2.14	0.45
1:D:513:SER:OG	1:D:518:LEU:HD23	2.17	0.45
1:M:587:ASP:OD1	1:M:588:TRP:N	2.50	0.45
1:N:410:LEU:HD12	1:N:410:LEU:H	1.82	0.45
1:P:688:ASP:HB3	1:H:556:GLY:HA2	1.99	0.45
1:H:678:MET:HE3	1:H:678:MET:HB3	1.60	0.45
1:L:710:LYS:HA	1:L:710:LYS:HZ2	1.81	0.45
1:F:532:LYS:HG3	1:F:558:ASP:CG	2.42	0.45
1:F:691:ASP:OD1	1:I:501:ARG:NH1	2.50	0.45
1:J:534:CYS:SG	1:J:559:LYS:HG3	2.57	0.45
1:O:553:ASN:HA	1:O:577:THR:HG21	1.98	0.45
1:C:533:TYR:HA	1:C:561:ILE:O	2.17	0.45
1:H:560:PHE:CD2	1:H:561:ILE:HG13	2.52	0.45
1:L:532:LYS:NZ	1:L:558:ASP:OD2	2.40	0.45
1:A:521:ILE:HG12	1:I:447:ARG:HB2	1.99	0.45
1:D:465:ASN:HB3	1:D:518:LEU:HD13	1.99	0.45
1:B:549:ILE:HG13	1:B:550:ASP:H	1.81	0.45
1:G:410:LEU:H	1:G:410:LEU:HD12	1.82	0.45
1:P:642:GLY:HA2	1:D:641:GLN:NE2	2.32	0.45
1:C:660:MET:HE1	1:H:727:TRP:CH2	2.52	0.45
1:M:646:VAL:O	1:M:650:ALA:N	2.50	0.45
1:E:520:ASP:OD1	1:E:535:ARG:NH2	2.47	0.45
1:G:520:ASP:OD1	1:G:535:ARG:NH1	2.46	0.45
1:N:572:GLY:CA	1:N:602:ARG:HH12	2.24	0.45
1:M:698:GLU:HG2	1:M:702:LYS:HE3	1.99	0.45
1:D:407:LYS:NZ	1:D:604:GLY:H	2.14	0.44
1:C:418:TYR:HB3	1:C:617:ASP:HB2	1.99	0.44
1:B:518:LEU:HD21	1:B:584:TYR:CD1	2.52	0.44
1:B:597:MET:HE2	1:B:611:VAL:HG11	1.99	0.44
1:G:524:ASP:OD2	1:P:447:ARG:HG3	2.17	0.44
1:J:588:TRP:CZ3	1:J:589:ASN:HB3	2.52	0.44
1:H:508:ARG:HB2	1:H:578:ALA:HA	1.98	0.44
1:D:687:ASP:HA	1:D:690:ILE:HG22	1.98	0.44
1:B:436:ASP:OD2	1:B:452:ARG:CZ	2.65	0.44
1:F:502:ILE:HG23	1:F:580:ILE:HD12	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:534:CYS:SG	1:N:559:LYS:HG3	2.58	0.44
1:L:653:ASN:HB2	1:L:656:GLU:H	1.82	0.44
1:A:698:GLU:HG2	1:A:702:LYS:HE3	2.00	0.44
1:F:635:ASP:O	1:F:639:ILE:HG13	2.17	0.44
1:D:534:CYS:HB2	1:D:562:PHE:HD1	1.82	0.44
1:D:560:PHE:CD2	1:D:561:ILE:HG13	2.52	0.44
1:D:562:PHE:CD2	1:D:564:LEU:HD11	2.53	0.44
1:K:585:ASP:OD1	1:K:585:ASP:N	2.51	0.44
1:A:532:LYS:HB3	1:A:559:LYS:HD3	1.99	0.44
1:B:653:ASN:ND2	1:B:655:ASP:HB2	2.32	0.44
1:F:654:LYS:HZ1	1:I:448:GLU:HB3	1.81	0.44
1:G:417:VAL:HG23	1:M:693:ILE:HG22	2.00	0.44
1:G:560:PHE:CD2	1:G:561:ILE:HG13	2.53	0.44
1:O:425:GLN:HG2	1:O:467:PRO:HG3	2.00	0.44
1:J:591:GLN:NE2	1:H:717:GLN:O	2.41	0.44
1:O:550:ASP:O	1:O:554:LYS:HG3	2.18	0.44
1:M:500:LYS:HB2	1:M:500:LYS:HE3	1.83	0.44
1:B:595:GLN:HE22	1:B:731:ASP:HA	1.83	0.44
1:F:518:LEU:HD21	1:F:584:TYR:CD1	2.53	0.44
1:N:536:ILE:HB	1:N:564:LEU:CD2	2.47	0.44
1:C:533:TYR:OH	1:K:446:LYS:NZ	2.50	0.44
1:D:419:ILE:HD13	1:D:491:LYS:HA	2.00	0.44
1:D:539:SER:O	1:D:543:GLU:HB3	2.18	0.44
1:B:419:ILE:HD13	1:B:491:LYS:HA	2.00	0.43
1:B:497:LYS:CE	1:K:687:ASP:HB3	2.46	0.43
1:B:508:ARG:NH2	1:B:577:THR:OG1	2.51	0.43
1:E:421:MET:O	1:H:674:ALA:HB1	2.18	0.43
1:F:411:PRO:O	1:F:608:GLN:NE2	2.37	0.43
1:F:660:MET:HE3	1:I:624:VAL:HG22	1.99	0.43
1:G:683:SER:OG	1:G:684:GLN:N	2.50	0.43
1:O:677:THR:HB	1:A:420:GLY:H	1.83	0.43
1:A:567:ARG:NH2	1:A:589:ASN:HB3	2.34	0.43
1:D:421:MET:HE1	1:D:429:TYR:CD2	2.53	0.43
1:O:624:VAL:HG22	1:A:660:MET:HE3	2.01	0.43
1:P:690:ILE:CD1	1:D:494:VAL:HG13	2.47	0.43
1:K:502:ILE:HG23	1:K:507:SER:HB2	2.01	0.43
1:A:458:MET:HE2	1:A:458:MET:HB3	1.53	0.43
1:I:585:ASP:N	1:I:585:ASP:OD1	2.51	0.43
1:E:560:PHE:CD2	1:E:561:ILE:HG13	2.53	0.43
1:F:634:LEU:O	1:F:638:VAL:HG23	2.18	0.43
1:O:486:ILE:HD12	1:O:529:ARG:HH21	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:624:VAL:HG22	1:D:660:MET:HE3	2.00	0.43
1:P:637:LEU:HD21	1:D:712:GLY:HA2	2.00	0.43
1:L:477:PRO:HA	1:L:478:PRO:C	2.44	0.43
1:A:710:LYS:HD3	1:A:710:LYS:HA	1.83	0.43
1:I:689:ASP:O	1:I:693:ILE:HG13	2.18	0.43
1:B:595:GLN:NE2	1:B:731:ASP:HA	2.34	0.43
1:B:663:HIS:HE1	1:K:622:GLU:OE2	2.02	0.43
1:O:523:GLU:HG2	1:O:533:TYR:CE1	2.54	0.43
1:L:418:TYR:HB3	1:L:617:ASP:HB2	2.01	0.43
1:E:519:LEU:HD22	1:E:563:LEU:HB3	2.00	0.43
1:F:520:ASP:OD1	1:F:535:ARG:NH1	2.37	0.43
1:G:446:LYS:HE2	1:M:725:TYR:HB3	2.01	0.43
1:H:654:LYS:O	1:H:658:LEU:HB2	2.17	0.43
1:K:675:PHE:HD1	1:K:675:PHE:HA	1.71	0.43
1:I:711:LEU:HD22	1:I:715:ASP:OD2	2.18	0.43
1:J:560:PHE:CD2	1:J:561:ILE:HG13	2.54	0.43
1:E:643:ARG:HD3	1:E:717:GLN:HE22	1.84	0.43
1:J:429:TYR:CD1	1:C:668:VAL:HG11	2.53	0.43
1:J:501:ARG:NH1	1:C:691:ASP:OD1	2.46	0.43
1:C:537:ASP:C	1:C:539:SER:H	2.26	0.43
1:B:681:LYS:HG3	1:B:685:LEU:HD22	2.01	0.43
1:E:407:LYS:NZ	1:E:604:GLY:H	2.17	0.43
1:L:479:TYR:O	1:L:480:THR:HG23	2.18	0.43
1:A:545:ARG:O	1:A:549:ILE:HG13	2.18	0.43
1:F:410:LEU:HD22	1:F:608:GLN:HB2	2.01	0.43
1:F:469:LEU:HA	1:F:517:ARG:HH21	1.84	0.43
1:O:718:LYS:HD3	1:O:722:GLU:OE1	2.19	0.43
1:P:534:CYS:SG	1:P:559:LYS:HG3	2.58	0.43
1:A:465:ASN:HB3	1:A:518:LEU:HD13	1.99	0.43
1:M:511:ILE:HG12	1:M:582:ILE:HB	2.00	0.43
1:N:419:ILE:HA	1:L:677:THR:HG21	2.01	0.42
1:N:669:PHE:HD2	1:L:430:GLN:HG2	1.84	0.42
1:K:455:ASN:O	1:K:459:GLN:HG2	2.19	0.42
1:F:560:PHE:CD2	1:F:561:ILE:HG13	2.54	0.42
1:J:418:TYR:HB3	1:J:617:ASP:HB2	1.99	0.42
1:J:499:LEU:HD21	1:J:511:ILE:HD11	2.01	0.42
1:B:549:ILE:HD13	1:B:573:ILE:HD13	2.00	0.42
1:K:481:THR:HG22	2:Q:17:ARG:NH1	2.33	0.42
1:K:485:LEU:HD11	1:K:521:ILE:HG23	2.01	0.42
1:B:500:LYS:HE3	1:B:500:LYS:HB2	1.82	0.42
1:B:532:LYS:HB2	1:B:559:LYS:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:624:VAL:HG22	1:M:660:MET:HE3	2.00	0.42
1:J:572:GLY:HA2	1:J:602:ARG:HH12	1.85	0.42
1:J:642:GLY:HA2	1:C:641:GLN:HE22	1.85	0.42
1:L:585:ASP:N	1:L:585:ASP:OD1	2.52	0.42
1:A:571:LEU:HD23	1:A:571:LEU:HA	1.63	0.42
1:F:723:SER:HB2	1:F:726:GLU:HG3	2.02	0.42
1:P:654:LYS:HE2	1:D:450:LYS:HE3	2.00	0.42
1:H:620:ILE:O	1:H:624:VAL:HG23	2.19	0.42
1:K:623:LYS:O	1:K:627:ARG:HG2	2.18	0.42
1:I:500:LYS:HB2	1:I:500:LYS:HE3	1.78	0.42
1:D:512:PHE:CE2	1:D:596:ALA:HB1	2.54	0.42
1:F:660:MET:HA	1:I:623:LYS:HG2	2.00	0.42
1:J:601:HIS:HB2	1:J:609:VAL:HG11	2.01	0.42
1:J:623:LYS:HG2	1:C:660:MET:HA	2.01	0.42
1:N:585:ASP:N	1:N:585:ASP:OD1	2.51	0.42
1:N:687:ASP:HB3	1:L:497:LYS:HE2	2.01	0.42
1:C:637:LEU:HD22	1:C:713:ILE:CD1	2.50	0.42
1:C:649:LYS:HA	1:H:724:ALA:HB2	2.01	0.42
1:H:552:TYR:CE1	1:H:562:PHE:HB2	2.55	0.42
1:L:541:ALA:O	1:L:545:ARG:N	2.50	0.42
1:D:517:ARG:HA	1:D:520:ASP:OD2	2.19	0.42
1:E:535:ARG:HA	1:E:563:LEU:O	2.19	0.42
1:E:648:ALA:HB2	1:H:634:LEU:HB3	2.01	0.42
1:G:587:ASP:N	1:G:593:ASP:OD2	2.41	0.42
1:C:635:ASP:O	1:C:639:ILE:HG12	2.19	0.42
1:O:638:VAL:HG13	1:A:641:GLN:HE21	1.85	0.42
1:L:601:HIS:HB2	1:L:609:VAL:HG11	2.02	0.42
1:A:571:LEU:CD1	1:A:595:GLN:HG2	2.48	0.42
1:A:653:ASN:OD1	1:A:654:LYS:N	2.51	0.42
1:I:485:LEU:HD11	1:I:521:ILE:HG23	2.01	0.42
1:M:534:CYS:SG	1:M:559:LYS:HG3	2.59	0.42
1:B:450:LYS:HB3	1:K:654:LYS:HE2	2.01	0.42
1:B:480:THR:O	1:B:484:HIS:CE1	2.73	0.42
1:B:684:GLN:NE2	1:K:486:ILE:HG21	2.35	0.42
1:E:455:ASN:O	1:E:459:GLN:HG2	2.20	0.42
1:F:535:ARG:HA	1:F:563:LEU:O	2.19	0.42
1:D:545:ARG:O	1:D:549:ILE:HG12	2.20	0.42
1:D:601:HIS:HB2	1:D:609:VAL:HG11	2.02	0.42
1:E:660:MET:HE1	1:H:588:TRP:CH2	2.55	0.42
1:G:690:ILE:HG13	1:M:497:LYS:CG	2.48	0.42
1:N:503:GLN:HB2	1:N:560:PHE:CZ	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:716:LEU:HB3	1:D:634:LEU:HD23	2.02	0.42
1:K:432:ILE:HD11	1:K:470:PHE:CZ	2.55	0.42
1:D:424:MET:HE2	1:D:473:ALA:HB1	2.02	0.42
1:M:703:GLU:CD	1:M:707:ARG:HH21	2.28	0.42
1:E:427:LYS:HE2	1:E:431:LYS:HZ1	1.84	0.41
1:P:520:ASP:OD1	1:P:535:ARG:NH2	2.47	0.41
1:I:482:ASP:HB2	1:I:484:HIS:HD2	1.85	0.41
1:M:483:GLU:HB3	1:M:487:TYR:HE2	1.82	0.41
1:F:719:PHE:O	1:A:545:ARG:HD2	2.20	0.41
1:C:425:GLN:HG2	1:C:467:PRO:HG3	2.01	0.41
1:H:533:TYR:HA	1:H:561:ILE:O	2.20	0.41
1:L:407:LYS:NZ	1:L:604:GLY:H	2.18	0.41
1:L:427:LYS:HE2	1:L:431:LYS:HZ3	1.84	0.41
1:I:667:LYS:O	1:I:671:THR:OG1	2.31	0.41
1:B:570:GLY:HA3	1:B:573:ILE:HD12	2.01	0.41
1:E:690:ILE:HD11	1:H:498:LEU:HD13	2.02	0.41
1:G:678:MET:O	1:G:682:GLY:HA3	2.19	0.41
1:N:620:ILE:O	1:N:624:VAL:HG23	2.21	0.41
1:K:431:LYS:HB2	1:K:431:LYS:NZ	2.35	0.41
1:K:569:GLY:O	1:K:599:ARG:HD2	2.20	0.41
1:D:436:ASP:CG	1:D:456:ILE:HD11	2.45	0.41
1:B:680:GLU:O	1:B:681:LYS:HD2	2.20	0.41
1:N:518:LEU:HD21	1:N:584:TYR:CD1	2.55	0.41
1:C:535:ARG:HA	1:C:563:LEU:O	2.20	0.41
1:H:418:TYR:HA	1:H:615:VAL:O	2.20	0.41
1:L:502:ILE:HG23	1:L:580:ILE:HD12	2.01	0.41
1:M:474:GLU:OE1	2:V:17:ARG:NH2	2.34	0.41
1:F:503:GLN:HB2	1:F:560:PHE:CZ	2.56	0.41
1:F:643:ARG:NH2	1:F:720:THR:O	2.45	0.41
1:G:429:TYR:CD1	1:M:668:VAL:HG11	2.56	0.41
1:J:536:ILE:HG22	1:J:568:ALA:HB1	2.02	0.41
1:O:491:LYS:HZ1	1:O:621:GLU:CD	2.28	0.41
1:O:560:PHE:CD2	1:O:561:ILE:HG13	2.55	0.41
1:P:431:LYS:HB2	1:P:431:LYS:NZ	2.35	0.41
1:M:421:MET:HE1	1:M:429:TYR:HD2	1.85	0.41
1:M:455:ASN:O	1:M:459:GLN:HG2	2.21	0.41
1:N:560:PHE:CD2	1:N:561:ILE:HG13	2.56	0.41
1:A:519:LEU:HD22	1:A:563:LEU:HB3	2.02	0.41
1:I:552:TYR:HE2	1:I:578:ALA:HB2	1.86	0.41
1:E:653:ASN:ND2	1:E:655:ASP:HB2	2.35	0.41
1:C:500:LYS:HB2	1:C:500:LYS:HE3	1.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:597:MET:HE2	1:C:611:VAL:HG11	2.03	0.41
1:L:635:ASP:O	1:L:639:ILE:HG12	2.21	0.41
1:D:457:VAL:HG13	1:D:620:ILE:HD11	2.01	0.41
1:A:447:ARG:H	1:A:447:ARG:HG3	1.65	0.41
1:A:560:PHE:CD2	1:A:561:ILE:HG13	2.55	0.41
1:B:450:LYS:HB3	1:K:654:LYS:HG2	2.02	0.41
1:E:619:ALA:O	1:H:663:HIS:HD2	2.04	0.41
1:F:544:ASP:O	1:F:548:ALA:N	2.52	0.41
1:F:590:PRO:O	1:F:594:LEU:HD13	2.21	0.41
1:G:542:HIS:CA	1:G:545:ARG:HH11	2.34	0.41
1:J:573:ILE:HG12	1:J:599:ARG:NH1	2.36	0.41
1:J:708:TYR:CE1	1:H:571:LEU:HB2	2.56	0.41
1:P:499:LEU:HD21	1:P:511:ILE:HD11	2.02	0.41
1:C:508:ARG:NH2	1:C:577:THR:OG1	2.49	0.41
1:C:560:PHE:CD2	1:C:561:ILE:HG13	2.55	0.41
1:K:653:ASN:HB3	1:K:656:GLU:HG3	2.03	0.41
1:L:500:LYS:HE3	1:L:500:LYS:HB2	1.88	0.41
1:L:569:GLY:O	1:L:599:ARG:HD2	2.21	0.41
1:A:631:LYS:O	1:A:635:ASP:HB2	2.21	0.41
1:A:689:ASP:O	1:A:693:ILE:HG13	2.21	0.41
1:D:534:CYS:SG	1:D:559:LYS:HE2	2.61	0.41
1:J:550:ASP:O	1:J:554:LYS:HG3	2.21	0.41
1:P:519:LEU:HD22	1:P:563:LEU:HB3	2.03	0.41
1:C:512:PHE:HZ	1:C:599:ARG:HD2	1.86	0.41
1:K:422:SER:OG	1:K:488:ASN:O	2.36	0.41
1:A:535:ARG:HA	1:A:563:LEU:O	2.21	0.41
1:I:573:ILE:O	1:I:602:ARG:NH1	2.54	0.41
1:D:425:GLN:HG2	1:D:467:PRO:HG3	2.02	0.41
1:D:654:LYS:H	1:D:654:LYS:HG2	1.56	0.41
1:F:521:ILE:HA	1:O:447:ARG:HG2	2.03	0.40
1:N:551:GLU:HA	1:N:554:LYS:HD2	2.03	0.40
1:B:595:GLN:HE22	1:B:732:PHE:H	1.69	0.40
1:G:552:TYR:HE2	1:G:578:ALA:HB2	1.87	0.40
1:J:623:LYS:HD2	1:J:623:LYS:HA	1.97	0.40
1:J:683:SER:HA	1:C:493:VAL:HG11	2.02	0.40
1:O:591:GLN:HA	1:O:594:LEU:HB2	2.02	0.40
1:K:532:LYS:HB2	1:K:559:LYS:HA	2.03	0.40
1:L:685:LEU:HD12	1:L:685:LEU:HA	1.87	0.40
1:B:595:GLN:HE22	1:B:732:PHE:N	2.19	0.40
1:F:519:LEU:O	1:F:523:GLU:N	2.52	0.40
1:N:579:ASP:OD2	1:N:580:ILE:HG13	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:654:LYS:HE3	1:P:654:LYS:HB2	1.89	0.40
1:C:457:VAL:HG13	1:C:620:ILE:HD11	2.04	0.40
1:C:712:GLY:O	1:C:715:ASP:OD2	2.39	0.40
1:H:519:LEU:HD11	1:H:564:LEU:C	2.47	0.40
1:L:432:ILE:O	1:L:435:LYS:HG2	2.21	0.40
1:M:457:VAL:HG13	1:M:620:ILE:HD11	2.02	0.40
1:B:497:LYS:HA	1:B:497:LYS:HD2	1.85	0.40
1:D:615:VAL:HG21	1:D:625:LEU:HD22	2.03	0.40
1:B:517:ARG:NH1	1:J:449:SER:HA	2.37	0.40
1:B:567:ARG:NE	1:B:732:PHE:HB3	2.37	0.40
1:B:619:ALA:O	1:K:663:HIS:HD2	2.05	0.40
1:F:425:GLN:HG2	1:F:467:PRO:HG3	2.03	0.40
1:N:669:PHE:CD2	1:L:430:GLN:HG2	2.56	0.40
1:O:668:VAL:HG11	1:A:429:TYR:HD1	1.86	0.40
1:P:535:ARG:HA	1:P:563:LEU:O	2.20	0.40
1:P:560:PHE:CD2	1:P:561:ILE:HG13	2.57	0.40
1:C:407:LYS:NZ	1:C:604:GLY:H	2.20	0.40
1:C:502:ILE:HG23	1:C:580:ILE:HD12	2.04	0.40
1:K:439:ALA:HB2	1:K:452:ARG:CZ	2.52	0.40
1:K:500:LYS:HB2	1:K:500:LYS:HE3	1.81	0.40
1:L:465:ASN:HB3	1:L:518:LEU:HD12	2.04	0.40
1:L:476:GLY:HA2	1:L:477:PRO:HD3	1.88	0.40
1:I:513:SER:OG	1:I:518:LEU:HD23	2.22	0.40
1:M:635:ASP:O	1:M:639:ILE:HG12	2.21	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	325/349 (93%)	309 (95%)	14 (4%)	2 (1%)	21 56

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	326/349 (93%)	309 (95%)	17 (5%)	0	100	100
1	C	320/349 (92%)	307 (96%)	13 (4%)	0	100	100
1	D	288/349 (82%)	279 (97%)	9 (3%)	0	100	100
1	E	311/349 (89%)	294 (94%)	17 (6%)	0	100	100
1	F	320/349 (92%)	310 (97%)	10 (3%)	0	100	100
1	G	306/349 (88%)	297 (97%)	9 (3%)	0	100	100
1	H	315/349 (90%)	307 (98%)	6 (2%)	2 (1%)	21	56
1	I	297/349 (85%)	287 (97%)	10 (3%)	0	100	100
1	J	313/349 (90%)	302 (96%)	10 (3%)	1 (0%)	36	70
1	K	302/349 (86%)	293 (97%)	8 (3%)	1 (0%)	36	70
1	L	309/349 (88%)	297 (96%)	12 (4%)	0	100	100
1	M	320/349 (92%)	311 (97%)	9 (3%)	0	100	100
1	N	315/349 (90%)	307 (98%)	8 (2%)	0	100	100
1	O	318/349 (91%)	307 (96%)	10 (3%)	1 (0%)	36	70
1	P	310/349 (89%)	298 (96%)	11 (4%)	1 (0%)	36	70
2	Q	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	R	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	S	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	T	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	U	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	V	3/21 (14%)	2 (67%)	1 (33%)	0	100	100
2	W	2/21 (10%)	2 (100%)	0	0	100	100
All	All	5015/5731 (88%)	4828 (96%)	179 (4%)	8 (0%)	43	76

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	731	ASP
1	O	676	GLY
1	J	676	GLY
1	A	541	ALA
1	H	540	THR
1	K	642	GLY
1	A	711	LEU

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Mol	Chain	Res	Type
1	P	676	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	275/294 (94%)	274 (100%)	1 (0%)	84	90
1	B	274/294 (93%)	273 (100%)	1 (0%)	84	90
1	C	273/294 (93%)	272 (100%)	1 (0%)	84	90
1	D	248/294 (84%)	247 (100%)	1 (0%)	84	90
1	E	264/294 (90%)	263 (100%)	1 (0%)	84	90
1	F	273/294 (93%)	272 (100%)	1 (0%)	84	90
1	G	259/294 (88%)	258 (100%)	1 (0%)	84	90
1	H	270/294 (92%)	264 (98%)	6 (2%)	45	74
1	I	254/294 (86%)	250 (98%)	4 (2%)	55	79
1	J	268/294 (91%)	264 (98%)	4 (2%)	57	80
1	K	258/294 (88%)	252 (98%)	6 (2%)	44	74
1	L	262/294 (89%)	258 (98%)	4 (2%)	57	80
1	M	273/294 (93%)	269 (98%)	4 (2%)	57	80
1	N	269/294 (92%)	264 (98%)	5 (2%)	50	76
1	O	273/294 (93%)	267 (98%)	6 (2%)	45	74
1	P	268/294 (91%)	267 (100%)	1 (0%)	84	90
2	Q	3/12 (25%)	3 (100%)	0	100	100
2	R	3/12 (25%)	3 (100%)	0	100	100
2	S	3/12 (25%)	3 (100%)	0	100	100
2	T	3/12 (25%)	3 (100%)	0	100	100
2	U	3/12 (25%)	3 (100%)	0	100	100
2	V	3/12 (25%)	3 (100%)	0	100	100
2	W	3/12 (25%)	3 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	4282/4788 (89%)	4235 (99%)	47 (1%)	65 83

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	690	ILE
1	E	713	ILE
1	F	685	LEU
1	G	441	ASN
1	J	431	LYS
1	J	634	LEU
1	J	640	GLN
1	J	675	PHE
1	N	431	LYS
1	N	486	ILE
1	N	571	LEU
1	N	594	LEU
1	N	675	PHE
1	O	431	LYS
1	O	448	GLU
1	O	594	LEU
1	O	634	LEU
1	O	675	PHE
1	O	728	ASN
1	P	431	LYS
1	C	480	THR
1	H	431	LYS
1	H	634	LEU
1	H	675	PHE
1	H	677	THR
1	H	678	MET
1	H	718	LYS
1	K	431	LYS
1	K	504	LYS
1	K	594	LEU
1	K	641	GLN
1	K	675	PHE
1	K	680	GLU
1	L	480	THR
1	L	675	PHE
1	L	710	LYS
1	L	711	LEU

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Mol	Chain	Res	Type
1	A	675	PHE
1	I	431	LYS
1	I	634	LEU
1	I	675	PHE
1	I	677	THR
1	D	713	ILE
1	M	431	LYS
1	M	641	GLN
1	M	675	PHE
1	M	730	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	416	ASN
1	B	484	HIS
1	B	514	GLN
1	B	595	GLN
1	B	640	GLN
1	E	455	ASN
1	E	618	ASN
1	E	636	GLN
1	E	645	GLN
1	F	505	GLN
1	F	591	GLN
1	F	641	GLN
1	F	717	GLN
1	G	488	ASN
1	G	514	GLN
1	G	542	HIS
1	G	589	ASN
1	J	640	GLN
1	N	455	ASN
1	N	459	GLN
1	O	589	ASN
1	P	589	ASN
2	U	18	HIS
1	C	416	ASN
1	C	505	GLN
1	C	542	HIS
1	C	636	GLN
1	H	484	HIS

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Mol	Chain	Res	Type
1	K	488	ASN
1	L	505	GLN
1	L	589	ASN
1	L	641	GLN
1	A	416	ASN
1	A	542	HIS
1	A	574	ASN
1	A	595	GLN
1	A	641	GLN
1	A	717	GLN
1	I	488	ASN
1	I	601	HIS
1	I	630	GLN
1	I	636	GLN
1	D	488	ASN
1	M	591	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	327/349 (93%)	0.01	2 (0%) 85 69	40, 74, 108, 133	0
1	B	328/349 (93%)	-0.04	6 (1%) 67 44	38, 68, 111, 145	0
1	C	324/349 (92%)	-0.20	1 (0%) 90 79	38, 61, 90, 139	0
1	D	292/349 (83%)	0.00	4 (1%) 73 51	38, 76, 125, 169	0
1	E	313/349 (89%)	-0.17	0 100 100	40, 71, 104, 132	0
1	F	324/349 (92%)	0.12	8 (2%) 58 35	42, 80, 128, 177	0
1	G	308/349 (88%)	-0.08	0 100 100	45, 75, 107, 142	0
1	H	321/349 (91%)	-0.21	0 100 100	40, 64, 97, 121	0
1	I	301/349 (86%)	-0.24	2 (0%) 84 66	35, 58, 109, 145	0
1	J	317/349 (90%)	-0.32	1 (0%) 90 79	36, 55, 91, 124	0
1	K	306/349 (87%)	-0.23	2 (0%) 84 66	42, 63, 99, 137	0
1	L	311/349 (89%)	-0.07	1 (0%) 90 79	46, 69, 105, 139	0
1	M	324/349 (92%)	-0.19	0 100 100	38, 65, 98, 132	0
1	N	319/349 (91%)	-0.32	0 100 100	39, 63, 92, 132	0
1	O	322/349 (92%)	-0.08	2 (0%) 85 69	45, 75, 101, 128	0
1	P	316/349 (90%)	0.17	6 (1%) 66 43	46, 92, 125, 151	0
2	Q	5/21 (23%)	-0.22	0 100 100	58, 66, 78, 80	0
2	R	5/21 (23%)	-0.28	0 100 100	63, 71, 74, 88	0
2	S	5/21 (23%)	0.09	0 100 100	76, 77, 81, 113	0
2	T	5/21 (23%)	-0.39	0 100 100	53, 56, 85, 87	0
2	U	5/21 (23%)	-0.34	0 100 100	58, 65, 72, 91	0
2	V	5/21 (23%)	-0.32	0 100 100	64, 66, 79, 79	0
2	W	4/21 (19%)	0.01	0 100 100	79, 81, 81, 85	0
All	All	5087/5731 (88%)	-0.12	35 (0%) 84 66	35, 69, 109, 177	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	444	GLY	5.6
1	P	485	LEU	4.5
1	I	570	GLY	4.3
1	K	570	GLY	4.1
1	F	445	GLY	3.7
1	B	444	GLY	3.4
1	B	549	ILE	3.3
1	F	481	THR	3.3
1	B	685	LEU	3.3
1	L	600	ALA	3.3
1	F	561	ILE	3.2
1	K	569	GLY	3.1
1	D	564	LEU	2.9
1	P	478	PRO	2.7
1	F	509	VAL	2.5
1	D	542	HIS	2.4
1	A	575	LEU	2.4
1	P	429	TYR	2.4
1	I	569	GLY	2.3
1	B	443	ALA	2.3
1	C	542	HIS	2.3
1	P	660	MET	2.3
1	O	731	ASP	2.2
1	D	541	ALA	2.2
1	F	734	ALA	2.2
1	J	536	ILE	2.2
1	P	466	HIS	2.1
1	F	537	ASP	2.1
1	F	573	ILE	2.1
1	B	449	SER	2.1
1	B	721	SER	2.1
1	O	586	SER	2.0
1	P	569	GLY	2.0
1	D	476	GLY	2.0
1	A	571	LEU	2.0

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

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.