



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:19 AM UTC

PDB ID : 5CJQ / pdb_00005cjq
Title : Crystal structure of a trimeric influenza hemagglutinin stem in complex with an broadly neutralizing antibody CR9114
Authors : Zhu, X.; Wilson, I.A.
Deposited on : 2015-07-14
Resolution : 3.60 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

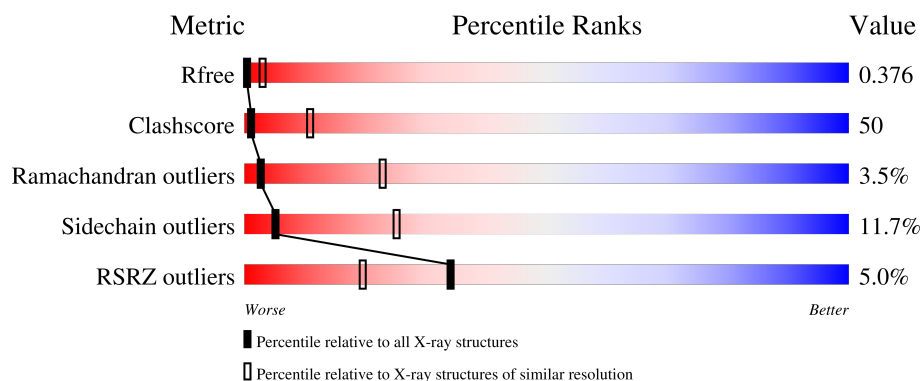
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	L	215	 9% 29% 43% 21% 5% •
2	H	230	 3% 47% 37% 10% • 6%
3	A	66	 3% 27% 44% 8% 21%
4	B	193	 2% 29% 41% 8% • 20%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4812 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 CR9114 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			

- Molecule 2 is a protein called CR9114 heavy chain.

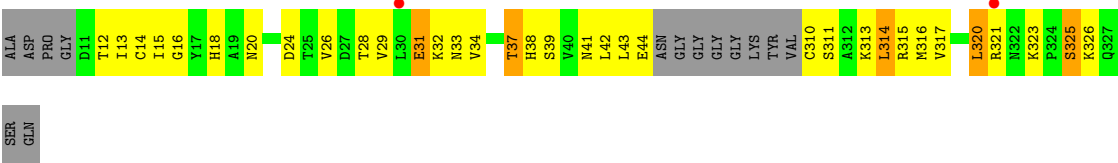
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1613	1017	269	320	7			

- Molecule 3 is a protein called Designed influenza hemagglutinin stem #4900, HA1.

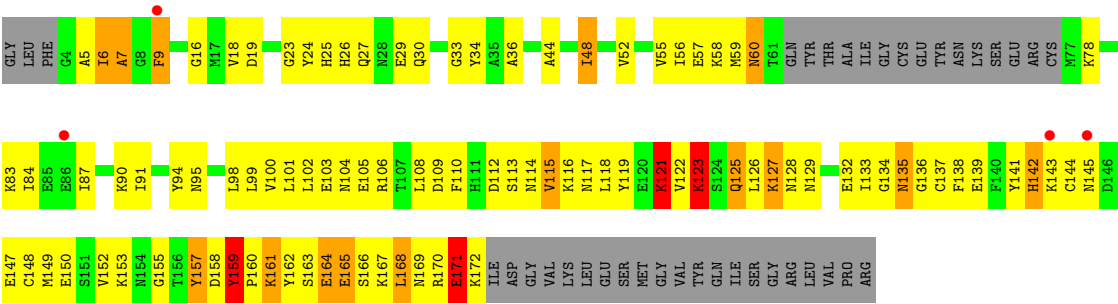
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	52	Total	C	N	O	S	0	0	0
			397	242	72	80	3			

- Molecule 4 is a protein called Designed influenza hemagglutinin stem #4900, HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	154	Total	C	N	O	S	0	0	0
			1234	767	208	251	8			



● Molecule 4: Designed influenza hemagglutinin stem #4900, HA2



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	157.70Å 157.70Å 202.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.85 – 3.60 38.85 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (38.85-3.60) 99.6 (38.85-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 3.55Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.346 , 0.369 0.351 , 0.376	Depositor DCC
R_{free} test set	550 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	138.2	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 155.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.030 for -2/3*h-1/3*k+2/3*l,-1/3*h-2/3*k-2/3*l,2/3*h-2/3*k+1/3*l 0.003 for -h,1/3*h-1/3*k+2/3*l,2/3*h+4/3*k+1/3*l 0.000 for -1/3*h+1/3*k-2/3*l,-k,-4/3*h-2/3*k+1/3*l	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	4812	wwPDB-VP
Average B, all atoms (Å ²)	158.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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	L	0.93	0/1606	2.03	70/2193 (3.2%)
2	H	0.88	9/1652 (0.5%)	1.49	34/2251 (1.5%)
3	A	0.79	0/399	1.47	7/539 (1.3%)
4	B	0.48	0/1256	1.08	5/1687 (0.3%)
All	All	0.81	9/4913 (0.2%)	1.61	116/6670 (1.7%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	L	0	2
2	H	0	2
All	All	0	4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	206	LYS	CA-CB	11.23	1.71	1.53
2	H	206	LYS	CA-C	7.08	1.60	1.52
2	H	207	VAL	N-CA	6.66	1.54	1.46
2	H	188	SER	CA-C	6.14	1.61	1.52
2	H	207	VAL	CA-CB	-5.62	1.47	1.54
2	H	206	LYS	CB-CG	5.32	1.68	1.52
2	H	206	LYS	CD-CE	5.29	1.68	1.52
2	H	171	GLN	N-CA	5.16	1.52	1.45
2	H	201	LYS	CA-CB	5.12	1.61	1.53

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	206	LYS	CB-CG-CD	14.55	144.77	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	142	GLY	N-CA-C	14.20	146.82	113.18
1	L	54	ARG	CA-C-N	13.63	133.82	119.78
1	L	54	ARG	C-N-CA	13.63	133.82	119.78
2	H	206	LYS	N-CA-C	-12.03	88.61	108.34
1	L	58	VAL	N-CA-C	11.19	118.26	108.63
1	L	64	GLY	N-CA-C	10.25	126.96	111.18
1	L	151	ASP	CB-CA-C	-9.81	103.89	117.23
1	L	118	PHE	CA-C-N	9.43	130.10	120.38
1	L	118	PHE	C-N-CA	9.43	130.10	120.38
1	L	136	ILE	N-CA-C	-9.32	95.12	108.17
2	H	196	CYS	N-CA-C	9.10	123.72	109.07
2	H	184	VAL	CB-CA-C	8.85	119.44	109.89
1	L	37	GLN	CB-CA-C	-8.79	94.52	109.72
1	L	170	ASN	N-CA-CB	-8.65	100.08	114.27
1	L	188	HIS	N-CA-C	8.65	123.86	110.20
1	L	139	PHE	CB-CA-C	-8.62	92.90	109.37
2	H	185	PRO	CA-C-O	8.32	131.82	121.67
2	H	206	LYS	CA-C-O	-8.16	111.59	120.24
1	L	67	SER	N-CA-C	8.07	122.98	108.69
1	L	170	ASN	N-CA-C	8.00	129.28	113.29
1	L	168	SER	N-CA-C	7.98	120.06	111.36
3	A	31	GLU	N-CA-C	7.95	121.83	109.52
2	H	189	LEU	CA-C-N	7.94	128.74	120.00
2	H	189	LEU	C-N-CA	7.94	128.74	120.00
1	L	73	LEU	N-CA-C	7.80	121.61	108.90
2	H	195	ILE	N-CA-CB	-7.78	100.16	111.52
1	L	140	TYR	CA-C-N	7.76	145.62	127.00
1	L	140	TYR	C-N-CA	7.76	145.62	127.00
1	L	39	PHE	N-CA-C	-7.70	100.39	110.39
1	L	140	TYR	C-N-CD	-7.68	103.70	120.60
1	L	66	LYS	CG-CD-CE	-7.40	94.27	111.30
1	L	120	PRO	N-CA-C	-7.38	99.62	111.14
2	H	199	ASN	N-CA-C	7.37	121.21	108.76
2	H	113	SER	N-CA-C	-7.31	99.21	110.10
3	A	314	LEU	N-CA-C	7.29	119.19	110.44
4	B	123	LYS	N-CA-C	-7.27	103.39	111.82
2	H	185	PRO	CA-C-N	7.21	131.27	120.31
2	H	185	PRO	C-N-CA	7.21	131.27	120.31
1	L	121	SER	N-CA-C	-7.15	99.90	110.48
1	L	3	ALA	N-CA-C	7.11	121.22	112.54
1	L	39	PHE	CB-CA-C	7.08	119.44	108.61
1	L	58	VAL	CA-C-N	7.06	127.05	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	58	VAL	C-N-CA	7.06	127.05	119.78
1	L	7	PRO	CA-C-N	6.98	126.80	119.05
1	L	7	PRO	C-N-CA	6.98	126.80	119.05
2	H	186	SER	N-CA-C	-6.96	103.74	111.82
1	L	139	PHE	N-CA-C	6.91	120.67	109.40
1	L	86	TYR	CA-CB-CG	6.88	126.29	113.90
2	H	210	ARG	CG-CD-NE	6.83	127.03	112.00
1	L	201	THR	N-CA-C	6.81	120.35	109.24
4	B	142	HIS	N-CA-C	-6.81	98.66	108.60
1	L	184	GLN	N-CA-C	-6.80	103.87	111.28
2	H	208	ASP	N-CA-C	6.80	120.01	109.07
1	L	163	THR	CB-CA-C	6.75	119.24	109.11
1	L	14	THR	CA-C-N	-6.69	112.87	119.76
1	L	14	THR	C-N-CA	-6.69	112.87	119.76
1	L	128	ASN	CA-C-N	-6.64	112.38	122.81
1	L	128	ASN	C-N-CA	-6.64	112.38	122.81
1	L	132	LEU	CA-C-N	-6.61	114.31	123.10
1	L	132	LEU	C-N-CA	-6.61	114.31	123.10
1	L	172	TYR	N-CA-C	6.56	120.59	110.42
1	L	163	THR	N-CA-C	-6.53	100.71	110.24
1	L	163	THR	CA-C-N	6.49	126.47	119.78
1	L	163	THR	C-N-CA	6.49	126.47	119.78
2	H	190	GLY	N-CA-C	6.48	121.06	112.77
4	B	159	TYR	CA-C-N	-6.46	111.76	119.84
4	B	159	TYR	C-N-CA	-6.46	111.76	119.84
1	L	70	SER	N-CA-C	6.45	119.69	109.50
2	H	11	VAL	N-CA-C	6.42	117.36	108.12
3	A	323	LYS	CA-C-N	6.39	126.30	119.85
3	A	323	LYS	C-N-CA	6.39	126.30	119.85
2	H	171	GLN	CG-CD-NE2	-6.29	106.96	116.40
2	H	185	PRO	N-CA-C	6.24	121.30	111.38
1	L	5	THR	N-CA-C	6.20	118.63	108.52
1	L	185	TRP	N-CA-C	6.20	118.11	111.36
1	L	95(A)	LYS	N-CA-C	-6.09	102.99	111.39
2	H	184	VAL	N-CA-C	-6.07	102.16	109.01
1	L	35	TRP	CA-CB-CG	6.05	125.09	113.60
2	H	206	LYS	O-C-N	-5.95	116.31	123.27
2	H	210	ARG	CB-CG-CD	5.94	124.97	111.30
1	L	127	ALA	N-CA-C	5.92	120.12	113.01
1	L	139	PHE	CA-CB-CG	5.83	119.63	113.80
1	L	185	TRP	CA-CB-CG	-5.80	102.58	113.60
2	H	101	ASP	N-CA-C	5.77	121.22	113.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	150	ALA	N-CA-C	-5.68	99.99	109.24
1	L	186	LYS	CB-CG-CD	5.63	124.24	111.30
2	H	160	THR	N-CA-C	5.61	120.99	114.04
2	H	206	LYS	CB-CA-C	5.58	119.07	110.14
2	H	197	ASN	CA-C-N	-5.56	115.70	123.10
2	H	197	ASN	C-N-CA	-5.56	115.70	123.10
1	L	169	ASN	N-CA-C	-5.55	106.54	113.15
4	B	155	GLY	N-CA-C	-5.51	107.71	115.32
1	L	145	THR	N-CA-C	5.48	118.33	109.40
1	L	37	GLN	CA-C-N	-5.44	115.10	122.77
1	L	37	GLN	C-N-CA	-5.44	115.10	122.77
1	L	134	CYS	N-CA-C	-5.44	99.87	108.73
3	A	33	ASN	N-CA-CB	5.43	119.73	111.70
2	H	207	VAL	N-CA-CB	5.41	119.41	111.52
1	L	73	LEU	CA-CB-CG	-5.38	97.47	116.30
3	A	38	HIS	N-CA-C	5.38	117.94	108.75
1	L	136	ILE	CB-CA-C	5.37	118.84	110.83
2	H	196	CYS	CB-CA-C	-5.35	100.46	109.72
1	L	114	SER	N-CA-C	-5.35	100.52	109.24
1	L	2	SER	CA-C-N	5.30	129.57	120.72
1	L	2	SER	C-N-CA	5.30	129.57	120.72
2	H	150	VAL	N-CA-C	5.22	116.00	108.48
3	A	37	THR	N-CA-C	5.19	117.02	111.36
2	H	157	GLY	CA-C-N	5.18	128.18	120.31
2	H	157	GLY	C-N-CA	5.18	128.18	120.31
1	L	193	CYS	CA-C-N	-5.12	115.88	123.05
1	L	193	CYS	C-N-CA	-5.12	115.88	123.05
1	L	189	ARG	N-CA-C	5.10	121.66	110.80
1	L	33	VAL	CA-C-N	-5.08	114.84	122.81
1	L	33	VAL	C-N-CA	-5.08	114.84	122.81
2	H	166	PHE	N-CA-C	5.06	117.49	110.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	203	SER	Peptide
2	H	206	LYS	Mainchain
1	L	129	LYS	Peptide
1	L	188	HIS	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	L	1568	0	1517	219	0
2	H	1613	0	1568	124	0
3	A	397	0	401	40	0
4	B	1234	0	1153	119	0
All	All	4812	0	4639	469	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 50.

All (469) 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)
2:H:114:ALA:HB1	2:H:146:PHE:CE2	1.56	1.40
2:H:114:ALA:CB	2:H:146:PHE:HE2	1.43	1.32
2:H:114:ALA:CB	2:H:146:PHE:CE2	2.21	1.19
1:L:140:TYR:CE1	1:L:171:LYS:HD2	1.82	1.14
2:H:146:PHE:HB2	2:H:175:LEU:HD11	1.28	1.10
1:L:124:GLU:HG2	1:L:129:LYS:HB2	1.45	0.98
3:A:316:MET:HE1	4:B:55:VAL:HG11	1.51	0.93
4:B:121:LYS:HZ3	4:B:122:VAL:HG22	1.33	0.92
2:H:114:ALA:HB1	2:H:146:PHE:HE2	0.92	0.91
1:L:136:ILE:HG22	1:L:139:PHE:CZ	2.08	0.89
2:H:120:SER:HB2	2:H:143:LYS:HD3	1.55	0.88
1:L:51:ASN:HA	1:L:66:LYS:HD2	1.56	0.88
1:L:21:ILE:HD12	1:L:86:TYR:CZ	2.10	0.86
1:L:142:GLY:HA2	1:L:172:TYR:CZ	2.11	0.85
2:H:195:ILE:HG23	2:H:210:ARG:HE	1.39	0.85
1:L:144:VAL:HG21	1:L:197:HIS:CD2	2.10	0.85
2:H:168:ALA:HA	2:H:178:LEU:HG	1.59	0.84
4:B:121:LYS:HG3	4:B:122:VAL:HG13	1.60	0.83
1:L:135:LEU:HD22	2:H:166:PHE:CE2	2.15	0.81
2:H:165:THR:HG22	2:H:180:SER:HB3	1.61	0.81
2:H:163:VAL:HG13	2:H:182:VAL:HB	1.61	0.81
1:L:149:LYS:HD3	1:L:152:SER:HA	1.61	0.81
2:H:13:LYS:HG3	2:H:14:PRO:HD2	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:150:ALA:HB1	1:L:188:HIS:CE1	2.16	0.80
1:L:150:ALA:HB1	1:L:188:HIS:HE1	1.46	0.80
4:B:164:GLU:OE1	4:B:164:GLU:N	2.14	0.80
1:L:146:VAL:HG22	1:L:148:TRP:HE1	1.47	0.79
4:B:162:TYR:CE2	4:B:165:GLU:HG2	2.17	0.79
1:L:144:VAL:HG21	1:L:197:HIS:HD2	1.45	0.79
1:L:184:GLN:HE22	1:L:188:HIS:CE1	1.99	0.79
1:L:130:ALA:HB3	1:L:180:LEU:O	1.83	0.79
4:B:122:VAL:O	4:B:126:LEU:N	2.17	0.78
3:A:39:SER:OG	3:A:315:ARG:CZ	2.32	0.77
1:L:196:THR:HA	1:L:201:THR:HA	1.66	0.77
2:H:209:LYS:C	2:H:210:ARG:HH11	1.92	0.77
1:L:75:ILE:HD11	1:L:86:TYR:HE2	1.50	0.76
1:L:37:GLN:HE22	1:L:47:LEU:HB2	1.50	0.76
1:L:45:LYS:HZ3	1:L:47:LEU:HD13	1.51	0.76
2:H:195:ILE:HG23	2:H:210:ARG:NE	2.00	0.76
2:H:201:LYS:HD2	2:H:201:LYS:H	1.50	0.76
3:A:325:SER:O	3:A:326:LYS:HG2	1.85	0.76
1:L:110:LYS:HD2	1:L:111:ALA:N	2.00	0.75
1:L:181:THR:HG22	1:L:182:PRO:HD2	1.67	0.75
1:L:188:HIS:CG	1:L:189:ARG:HD2	2.23	0.74
3:A:321:ARG:HD3	4:B:108:LEU:HG	1.69	0.74
4:B:158:ASP:OD1	4:B:159:TYR:N	2.20	0.74
1:L:122:SER:O	1:L:126:GLN:NE2	2.20	0.74
1:L:11:VAL:HG11	1:L:21:ILE:HD11	1.68	0.73
4:B:9:PHE:HZ	4:B:115:VAL:HG11	1.54	0.73
1:L:124:GLU:OE2	1:L:130:ALA:HA	1.88	0.72
1:L:195:VAL:O	1:L:202:VAL:N	2.19	0.72
3:A:26:VAL:HG13	3:A:315:ARG:HH11	1.54	0.71
4:B:162:TYR:CD2	4:B:165:GLU:HG2	2.25	0.71
1:L:28:ILE:HD11	1:L:69:THR:O	1.90	0.71
1:L:118:PHE:HB3	2:H:124:LEU:HD22	1.71	0.71
2:H:114:ALA:HB3	2:H:146:PHE:HE2	1.50	0.70
1:L:37:GLN:HE22	1:L:47:LEU:CB	2.04	0.70
2:H:182:VAL:HG22	2:H:184:VAL:HG23	1.74	0.70
2:H:146:PHE:HB2	2:H:175:LEU:CD1	2.17	0.70
1:L:49:TYR:O	1:L:53:GLN:HB2	1.90	0.69
1:L:188:HIS:HA	1:L:189:ARG:HH11	1.58	0.69
1:L:144:VAL:HG11	1:L:196:THR:O	1.93	0.69
1:L:144:VAL:HG12	1:L:145:THR:H	1.58	0.69
1:L:35:TRP:CZ2	1:L:73:LEU:HD13	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:28:THR:HB	4:B:105:GLU:HG2	1.74	0.68
4:B:129:ASN:O	4:B:141:TYR:N	2.16	0.68
4:B:122:VAL:HA	4:B:125:GLN:HB2	1.75	0.68
1:L:75:ILE:HD11	1:L:86:TYR:CE2	2.29	0.68
2:H:209:LYS:N	2:H:210:ARG:NH1	2.41	0.67
4:B:23:GLY:HA3	4:B:36:ALA:HA	1.76	0.67
2:H:146:PHE:CB	2:H:175:LEU:HD11	2.18	0.67
1:L:61:ARG:HH22	1:L:82:ASP:CG	2.02	0.67
2:H:13:LYS:CG	2:H:14:PRO:HD2	2.24	0.66
2:H:195:ILE:HG22	2:H:210:ARG:HH21	1.60	0.66
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.77	0.66
3:A:26:VAL:HG13	3:A:315:ARG:NH1	2.11	0.66
1:L:151:ASP:N	1:L:189:ARG:HD3	2.11	0.66
2:H:11:VAL:HG13	2:H:110:THR:O	1.94	0.66
2:H:149:PRO:O	2:H:200:HIS:HB2	1.96	0.66
2:H:11:VAL:HG22	2:H:110:THR:HB	1.77	0.65
2:H:123:PRO:HD3	2:H:209:LYS:HG2	1.78	0.65
1:L:118:PHE:CD1	2:H:124:LEU:HB3	2.31	0.65
1:L:166:LYS:CE	1:L:170:ASN:OD1	2.44	0.65
2:H:148:GLU:HG3	2:H:149:PRO:HA	1.78	0.64
1:L:35:TRP:HZ2	1:L:73:LEU:HD22	1.62	0.64
1:L:146:VAL:CG2	1:L:148:TRP:HE1	2.10	0.64
1:L:6:GLN:NE2	1:L:88:CYS:H	1.94	0.64
1:L:184:GLN:HE22	1:L:188:HIS:CD2	2.15	0.64
1:L:5:THR:O	1:L:23:CYS:HA	1.98	0.64
1:L:47:LEU:HD12	1:L:58:VAL:HG22	1.79	0.64
1:L:166:LYS:HE3	1:L:170:ASN:OD1	1.98	0.64
2:H:148:GLU:HG3	2:H:149:PRO:CA	2.28	0.64
1:L:47:LEU:HD12	1:L:58:VAL:CG2	2.27	0.64
2:H:2:VAL:HG11	2:H:94:ARG:NH1	2.12	0.64
4:B:103:GLU:OE1	4:B:103:GLU:N	2.31	0.64
2:H:209:LYS:N	2:H:210:ARG:HH11	1.96	0.63
1:L:188:HIS:CD2	1:L:189:ARG:NH1	2.66	0.63
1:L:35:TRP:CZ2	1:L:73:LEU:HD22	2.33	0.63
1:L:33:VAL:HG11	1:L:66:LYS:HZ3	1.64	0.63
3:A:39:SER:HB2	3:A:316:MET:O	1.99	0.63
4:B:103:GLU:HA	4:B:106:ARG:CG	2.29	0.63
1:L:136:ILE:CG2	1:L:139:PHE:CZ	2.82	0.62
1:L:156:LYS:O	1:L:159:VAL:HG12	1.99	0.62
1:L:37:GLN:NE2	1:L:47:LEU:HB2	2.15	0.62
1:L:14:THR:HG22	1:L:106(A):LEU:HD12	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:170:ASN:C	1:L:171:LYS:HD3	2.24	0.62
1:L:144:VAL:HG11	1:L:197:HIS:HA	1.81	0.62
1:L:141:PRO:C	1:L:172:TYR:CE2	2.78	0.61
1:L:171:LYS:HD3	1:L:171:LYS:N	2.14	0.61
2:H:3:GLN:OE1	2:H:4:LEU:N	2.32	0.61
2:H:204:ASN:OD1	2:H:204:ASN:N	2.33	0.61
1:L:33:VAL:HG11	1:L:66:LYS:NZ	2.15	0.61
1:L:33:VAL:HG22	1:L:89:ALA:O	2.00	0.61
1:L:110:LYS:HZ1	1:L:112:ALA:HA	1.65	0.61
2:H:200:HIS:N	2:H:205:THR:O	2.33	0.61
2:H:150:VAL:HG23	2:H:200:HIS:HB3	1.82	0.61
2:H:201:LYS:H	2:H:201:LYS:CD	2.10	0.61
4:B:9:PHE:HZ	4:B:115:VAL:CG1	2.14	0.61
1:L:52:ASP:OD1	1:L:53:GLN:NE2	2.33	0.60
1:L:85:GLU:HG2	1:L:102:THR:O	2.01	0.60
4:B:162:TYR:HE2	4:B:165:GLU:HG2	1.65	0.60
1:L:28:ILE:HG22	1:L:33:VAL:HG23	1.83	0.60
1:L:124:GLU:HG2	1:L:129:LYS:CB	2.28	0.60
2:H:186:SER:O	2:H:189:LEU:HG	2.02	0.60
1:L:23:CYS:N	1:L:71:ALA:O	2.34	0.60
4:B:121:LYS:CG	4:B:122:VAL:HG13	2.31	0.60
1:L:33:VAL:HG23	1:L:90:ALA:HB2	1.83	0.60
1:L:27(A):SER:HA	1:L:30:ARG:HH11	1.65	0.60
2:H:184:VAL:HG13	2:H:185:PRO:HD2	1.83	0.60
1:L:54:ARG:NH1	1:L:58:VAL:O	2.35	0.59
1:L:184:GLN:O	1:L:184:GLN:NE2	2.35	0.59
2:H:201:LYS:HD2	2:H:201:LYS:N	2.16	0.59
2:H:114:ALA:CB	2:H:146:PHE:CZ	2.83	0.59
1:L:144:VAL:HG12	1:L:145:THR:N	2.15	0.59
2:H:170:LEU:HD13	2:H:176:TYR:CE1	2.38	0.59
4:B:102:LEU:HB3	4:B:106:ARG:NH1	2.17	0.59
4:B:94:TYR:O	4:B:98:LEU:HG	2.03	0.59
1:L:186:LYS:HG2	1:L:187:SER:N	2.18	0.58
2:H:195:ILE:CG2	2:H:210:ARG:HH21	2.16	0.58
1:L:6:GLN:HE22	1:L:88:CYS:H	1.52	0.58
1:L:83:GLU:OE2	1:L:106:VAL:CG2	2.51	0.58
4:B:103:GLU:HA	4:B:106:ARG:HG2	1.85	0.58
4:B:119:TYR:O	4:B:123:LYS:HE2	2.03	0.58
1:L:83:GLU:HG3	1:L:106:VAL:HG22	1.85	0.58
1:L:66:LYS:HB3	1:L:71:ALA:HB2	1.85	0.58
1:L:175:SER:OG	2:H:166:PHE:CD2	2.54	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:13:ILE:HA	4:B:26:HIS:HA	1.86	0.58
1:L:118:PHE:HE2	1:L:135:LEU:HD12	1.69	0.58
1:L:160:GLU:HB3	2:H:169:VAL:HG21	1.86	0.58
4:B:19:ASP:HB2	4:B:36:ALA:HB3	1.85	0.58
1:L:115:VAL:HG22	1:L:136:ILE:HG23	1.86	0.57
1:L:188:HIS:ND1	1:L:189:ARG:HD2	2.19	0.57
1:L:142:GLY:HA2	1:L:172:TYR:CE2	2.39	0.57
4:B:112:ASP:O	4:B:116:LYS:HG3	2.05	0.57
2:H:155:ASN:HD21	2:H:195:ILE:H	1.51	0.57
1:L:37:GLN:HG2	1:L:45:LYS:O	2.04	0.56
2:H:13:LYS:HG3	2:H:113:SER:HA	1.86	0.56
1:L:83:GLU:OE2	1:L:106:VAL:HG23	2.05	0.56
1:L:135:LEU:HD22	2:H:166:PHE:CD2	2.39	0.56
1:L:35:TRP:HZ2	1:L:73:LEU:HD13	1.70	0.56
1:L:197:HIS:CE1	1:L:198:GLU:HG2	2.41	0.56
1:L:147:ALA:C	1:L:148:TRP:HD1	2.14	0.56
2:H:113:SER:O	2:H:114:ALA:HB2	2.05	0.56
3:A:43:LEU:HD21	3:A:314:LEU:HD12	1.88	0.56
1:L:65:SER:C	1:L:66:LYS:HG2	2.31	0.56
2:H:148:GLU:HB2	2:H:176:TYR:CE2	2.41	0.56
1:L:37:GLN:CG	1:L:45:LYS:HG2	2.36	0.56
1:L:39:PHE:CE1	1:L:42:THR:HG21	2.41	0.56
2:H:94:ARG:O	2:H:100(D):MET:HA	2.06	0.56
3:A:16:GLY:HA2	4:B:9:PHE:CE1	2.41	0.56
1:L:27(A):SER:HA	1:L:30:ARG:NH1	2.21	0.56
1:L:166:LYS:HE3	1:L:170:ASN:O	2.06	0.56
1:L:51:ASN:CA	1:L:66:LYS:HD2	2.34	0.55
1:L:35:TRP:NE1	1:L:73:LEU:HD13	2.21	0.55
3:A:321:ARG:NH1	4:B:108:LEU:HD23	2.20	0.55
4:B:121:LYS:HG2	4:B:122:VAL:N	2.21	0.55
1:L:47:LEU:O	1:L:55:PRO:HD2	2.05	0.55
1:L:23:CYS:O	1:L:71:ALA:N	2.33	0.55
1:L:35:TRP:HE1	1:L:73:LEU:HD13	1.71	0.55
1:L:179:SER:O	1:L:180:LEU:HG	2.07	0.55
4:B:59:MET:SD	4:B:59:MET:N	2.78	0.55
4:B:133:ILE:HG12	4:B:137:CYS:C	2.30	0.55
4:B:165:GLU:O	4:B:169:ASN:N	2.33	0.55
1:L:38:GLN:NE2	1:L:44:PRO:HG3	2.22	0.55
3:A:24:ASP:HB3	3:A:315:ARG:NH2	2.22	0.55
2:H:36:TRP:CE2	2:H:80:MET:HB2	2.42	0.55
1:L:183:GLU:O	1:L:187:SER:OG	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:183:GLU:HG3	1:L:184:GLN:N	2.22	0.55
2:H:129:LYS:HG3	2:H:135:THR:O	2.07	0.55
3:A:39:SER:OG	3:A:315:ARG:NE	2.39	0.55
4:B:78:LYS:H	4:B:78:LYS:HD2	1.72	0.55
3:A:310:CYS:SG	3:A:311:SER:N	2.80	0.54
1:L:34:ASN:O	1:L:89:ALA:N	2.39	0.54
4:B:118:LEU:O	4:B:121:LYS:HD3	2.07	0.54
1:L:110:LYS:HD3	1:L:197:HIS:CE1	2.42	0.54
2:H:40:ALA:HB3	2:H:43:GLN:CD	2.32	0.54
1:L:48:ILE:HD11	1:L:51:ASN:O	2.07	0.54
1:L:110:LYS:HD2	1:L:111:ALA:H	1.72	0.54
1:L:148:TRP:CZ2	1:L:176:SER:HB3	2.42	0.54
2:H:11:VAL:HG11	2:H:147:PRO:HG3	1.90	0.54
1:L:36:TYR:O	1:L:87:TYR:N	2.37	0.54
1:L:35:TRP:CE2	1:L:73:LEU:HD13	2.43	0.54
2:H:192:GLN:NE2	2:H:193:THR:O	2.41	0.54
4:B:30:GLN:HE22	4:B:145:ASN:HB3	1.72	0.54
1:L:61:ARG:NH2	1:L:82:ASP:OD2	2.41	0.54
1:L:127:ALA:O	1:L:129:LYS:HE3	2.08	0.53
3:A:16:GLY:HA2	4:B:9:PHE:HE1	1.73	0.53
2:H:95:HIS:HB3	2:H:100(B):SER:OG	2.08	0.53
3:A:311:SER:HB3	3:A:313:LYS:NZ	2.24	0.53
1:L:96:ALA:HB3	2:H:47:TRP:CG	2.43	0.53
1:L:186:LYS:CG	1:L:187:SER:N	2.72	0.53
1:L:7:PRO:O	1:L:102:THR:OG1	2.25	0.53
2:H:93:ALA:HB1	2:H:100(D):MET:HB3	1.90	0.53
2:H:159:LEU:HD21	2:H:194:TYR:HD1	1.73	0.53
1:L:168:SER:C	1:L:170:ASN:H	2.16	0.53
2:H:121:VAL:HG23	2:H:209:LYS:HB2	1.90	0.53
4:B:90:LYS:HZ2	4:B:94:TYR:HE2	1.57	0.53
1:L:50:SER:HB2	1:L:53:GLN:HG2	1.90	0.53
1:L:151:ASP:CG	1:L:189:ARG:HE	2.17	0.53
2:H:198:VAL:O	2:H:207:VAL:N	2.32	0.53
1:L:35:TRP:CH2	1:L:47:LEU:HD23	2.44	0.53
2:H:124:LEU:N	2:H:139:GLY:O	2.26	0.52
4:B:164:GLU:H	4:B:164:GLU:CD	2.12	0.52
1:L:21:ILE:HD12	1:L:86:TYR:OH	2.07	0.52
1:L:37:GLN:NE2	1:L:47:LEU:HD13	2.24	0.52
2:H:4:LEU:HG	2:H:102:VAL:CG1	2.38	0.52
2:H:126:PRO:HA	2:H:137:ALA:O	2.09	0.52
4:B:159:TYR:C	4:B:161:LYS:N	2.67	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:2:VAL:HG11	2:H:94:ARG:HH12	1.74	0.52
1:L:132:LEU:HD11	1:L:185:TRP:CZ2	2.44	0.52
2:H:119:PRO:HD3	2:H:200:HIS:HE1	1.72	0.52
2:H:166:PHE:C	2:H:178:LEU:HD21	2.34	0.52
4:B:134:GLY:O	4:B:135:ASN:HB3	2.09	0.52
1:L:136:ILE:CG2	1:L:139:PHE:HZ	2.23	0.52
3:A:15:ILE:HG13	4:B:24:TYR:HD1	1.75	0.52
2:H:4:LEU:HB3	2:H:22:CYS:SG	2.51	0.51
1:L:146:VAL:HG22	1:L:148:TRP:NE1	2.20	0.51
1:L:113:PRO:HA	1:L:138:ASP:HB3	1.93	0.51
4:B:159:TYR:CZ	4:B:163:SER:HA	2.46	0.51
2:H:119:PRO:HD3	2:H:200:HIS:CE1	2.45	0.51
2:H:123:PRO:HB2	2:H:211:VAL:HG23	1.92	0.51
1:L:37:GLN:NE2	1:L:45:LYS:HG2	2.25	0.51
3:A:16:GLY:CA	4:B:9:PHE:HE1	2.24	0.51
4:B:25:HIS:HD2	4:B:33:GLY:C	2.19	0.51
1:L:35:TRP:CZ3	1:L:86:TYR:CD1	2.99	0.51
1:L:147:ALA:C	1:L:148:TRP:CD1	2.89	0.51
2:H:184:VAL:HG11	2:H:194:TYR:OH	2.11	0.51
2:H:210:ARG:N	2:H:210:ARG:HD2	2.25	0.51
1:L:66:LYS:NZ	1:L:71:ALA:HB1	2.26	0.50
2:H:155:ASN:ND2	2:H:195:ILE:H	2.09	0.50
3:A:34:VAL:HG21	3:A:321:ARG:HH12	1.76	0.50
2:H:33:ALA:HB2	2:H:98:TYR:O	2.11	0.50
2:H:162:GLY:C	2:H:182:VAL:HG23	2.36	0.50
4:B:163:SER:O	4:B:167:LYS:HG2	2.10	0.50
1:L:33:VAL:CG2	1:L:90:ALA:HB2	2.41	0.50
1:L:35:TRP:HE1	1:L:73:LEU:CD1	2.24	0.50
1:L:73:LEU:HD23	1:L:74:ALA:N	2.27	0.50
4:B:6:ILE:HG23	4:B:7:ALA:H	1.76	0.50
4:B:165:GLU:HA	4:B:168:LEU:HG	1.92	0.50
1:L:144:VAL:CG1	1:L:196:THR:O	2.59	0.50
4:B:25:HIS:HB2	4:B:34:TYR:HD1	1.76	0.50
1:L:204:LYS:HD2	1:L:205:THR:H	1.77	0.50
4:B:102:LEU:HB3	4:B:106:ARG:HH12	1.76	0.50
2:H:11:VAL:HG21	2:H:147:PRO:HG2	1.93	0.50
1:L:140:TYR:HA	1:L:172:TYR:HE2	1.76	0.49
2:H:145:TYR:HB2	2:H:200:HIS:CE1	2.47	0.49
1:L:132:LEU:HD11	1:L:185:TRP:HZ2	1.77	0.49
1:L:151:ASP:H	1:L:189:ARG:HD3	1.76	0.49
4:B:25:HIS:HB2	4:B:34:TYR:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:133:ILE:HG12	4:B:137:CYS:O	2.12	0.49
1:L:37:GLN:HE21	1:L:45:LYS:HG2	1.77	0.49
4:B:166:SER:O	4:B:170:ARG:HB2	2.12	0.49
1:L:85:GLU:HG2	1:L:102:THR:C	2.38	0.49
1:L:114:SER:H	1:L:138:ASP:HB3	1.77	0.49
2:H:199:ASN:HA	2:H:205:THR:O	2.12	0.49
4:B:142:HIS:HD2	4:B:143:LYS:H	1.60	0.49
1:L:117:LEU:HD11	1:L:132:LEU:HB3	1.95	0.49
1:L:132:LEU:HD13	1:L:178:LEU:HB3	1.94	0.49
1:L:204:LYS:CD	1:L:205:THR:H	2.26	0.49
2:H:136:ALA:O	2:H:183:THR:HG23	2.13	0.49
4:B:9:PHE:CE2	4:B:115:VAL:HB	2.48	0.49
2:H:196:CYS:C	2:H:210:ARG:HH22	2.21	0.49
4:B:150:GLU:OE2	4:B:153:LYS:NZ	2.41	0.49
1:L:95:LEU:HA	2:H:61:GLN:OE1	2.13	0.48
1:L:37:GLN:HG3	1:L:45:LYS:HG2	1.94	0.48
1:L:121:SER:O	1:L:124:GLU:HB3	2.12	0.48
1:L:65:SER:O	1:L:66:LYS:HG2	2.13	0.48
1:L:128:ASN:HB2	1:L:129:LYS:HD3	1.96	0.48
2:H:209:LYS:H	2:H:210:ARG:NH1	2.10	0.48
1:L:139:PHE:HB3	1:L:172:TYR:CZ	2.48	0.48
4:B:87:ILE:O	4:B:91:ILE:HG13	2.14	0.48
4:B:121:LYS:HZ3	4:B:122:VAL:CG2	2.15	0.48
2:H:125:ALA:O	2:H:127:SER:N	2.41	0.48
4:B:24:TYR:CE1	4:B:121:LYS:HE3	2.48	0.48
2:H:184:VAL:CG1	2:H:185:PRO:HD2	2.43	0.48
2:H:201:LYS:HA	2:H:204:ASN:HA	1.95	0.48
4:B:90:LYS:NZ	4:B:94:TYR:HE2	2.12	0.48
4:B:114:ASN:O	4:B:117:ASN:HB2	2.13	0.48
2:H:73:ILE:HG13	2:H:74:PHE:N	2.28	0.47
2:H:195:ILE:HG22	2:H:210:ARG:NH2	2.27	0.47
4:B:159:TYR:CE1	4:B:161:LYS:HE2	2.49	0.47
2:H:165:THR:HG22	2:H:180:SER:CB	2.38	0.47
1:L:44:PRO:HD2	2:H:103:TRP:CE3	2.49	0.47
4:B:109:ASP:O	4:B:113:SER:HB3	2.14	0.47
2:H:148:GLU:HG3	2:H:149:PRO:CB	2.44	0.47
4:B:127:LYS:HB3	4:B:127:LYS:HE3	1.35	0.47
4:B:133:ILE:HD11	4:B:139:GLU:N	2.29	0.47
1:L:85:GLU:C	1:L:85:GLU:OE1	2.58	0.47
2:H:74:PHE:CZ	3:A:42:LEU:HA	2.49	0.47
2:H:159:LEU:HD21	2:H:194:TYR:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:154:TRP:CZ3	2:H:196:CYS:SG	3.08	0.47
4:B:52:VAL:O	4:B:56:ILE:HG13	2.14	0.47
1:L:182:PRO:O	1:L:186:LYS:N	2.39	0.47
4:B:128:ASN:ND2	4:B:170:ARG:HH22	2.12	0.47
3:A:13:ILE:HG13	4:B:26:HIS:HB3	1.97	0.47
1:L:114:SER:N	1:L:138:ASP:HB3	2.30	0.46
4:B:59:MET:O	4:B:60:ASN:HB3	2.16	0.46
4:B:84:ILE:O	4:B:87:ILE:HB	2.15	0.46
1:L:14:THR:HG22	1:L:106(A):LEU:CD1	2.43	0.46
3:A:15:ILE:HD11	4:B:121:LYS:NZ	2.30	0.46
3:A:313:LYS:HE2	3:A:313:LYS:HB2	1.80	0.46
4:B:127:LYS:HE2	4:B:159:TYR:CD2	2.50	0.46
1:L:140:TYR:HE1	1:L:171:LYS:HD2	1.68	0.46
2:H:195:ILE:CG2	2:H:210:ARG:NH2	2.77	0.46
4:B:127:LYS:HB3	4:B:128:ASN:H	1.58	0.46
4:B:30:GLN:NE2	4:B:145:ASN:HB3	2.31	0.46
1:L:148:TRP:CE3	1:L:178:LEU:HB2	2.50	0.46
2:H:4:LEU:HG	2:H:102:VAL:HG12	1.98	0.46
3:A:12:THR:HA	4:B:138:PHE:O	2.16	0.46
4:B:9:PHE:HE2	4:B:115:VAL:HB	1.80	0.46
4:B:159:TYR:CE2	4:B:163:SER:HB3	2.50	0.46
1:L:160:GLU:CB	2:H:169:VAL:HG21	2.45	0.46
3:A:18:HIS:O	3:A:320:LEU:HD11	2.16	0.46
4:B:16:GLY:HA3	4:B:34:TYR:CE2	2.51	0.46
1:L:35:TRP:HB3	1:L:88:CYS:HA	1.98	0.45
1:L:118:PHE:CE2	1:L:135:LEU:HD12	2.51	0.45
1:L:35:TRP:HA	1:L:87:TYR:O	2.16	0.45
2:H:195:ILE:HG23	2:H:210:ARG:CZ	2.46	0.45
4:B:115:VAL:HA	4:B:118:LEU:HB2	1.99	0.45
1:L:122:SER:O	1:L:126:GLN:HG2	2.16	0.45
1:L:133:VAL:HG11	2:H:179:SER:OG	2.16	0.45
1:L:142:GLY:HA2	1:L:172:TYR:CE1	2.51	0.45
2:H:155:ASN:HD22	2:H:159:LEU:HD22	1.81	0.45
4:B:25:HIS:CD2	4:B:33:GLY:H	2.35	0.45
4:B:58:LYS:HD2	4:B:58:LYS:N	2.32	0.45
1:L:134:CYS:SG	1:L:148:TRP:CZ2	3.10	0.45
2:H:209:LYS:C	2:H:210:ARG:HD2	2.41	0.45
1:L:58:VAL:HA	1:L:59:PRO:HD3	1.80	0.45
4:B:157:TYR:OH	4:B:161:LYS:NZ	2.35	0.45
3:A:28:THR:CB	4:B:105:GLU:HG2	2.44	0.45
4:B:159:TYR:CE2	4:B:163:SER:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:47:LEU:HD12	1:L:58:VAL:HG21	1.99	0.45
4:B:122:VAL:CG2	4:B:126:LEU:HG	2.47	0.45
4:B:163:SER:OG	4:B:164:GLU:OE1	2.33	0.45
1:L:117:LEU:HD13	1:L:134:CYS:SG	2.57	0.45
1:L:139:PHE:CB	1:L:172:TYR:CZ	3.00	0.45
1:L:184:GLN:NE2	1:L:188:HIS:CD2	2.84	0.44
2:H:152:VAL:HG22	2:H:198:VAL:HG13	1.99	0.44
4:B:56:ILE:O	4:B:59:MET:HG2	2.17	0.44
2:H:32:TYR:CZ	2:H:94:ARG:NH2	2.84	0.44
3:A:24:ASP:HB3	3:A:315:ARG:CZ	2.46	0.44
3:A:42:LEU:HD11	3:A:316:MET:HE3	1.98	0.44
1:L:34:ASN:N	1:L:89:ALA:O	2.49	0.44
1:L:132:LEU:CG	1:L:185:TRP:HZ2	2.31	0.44
1:L:183:GLU:OE1	1:L:186:LYS:HE2	2.17	0.44
1:L:188:HIS:HB3	1:L:189:ARG:HB2	1.98	0.44
4:B:135:ASN:OD1	4:B:136:GLY:N	2.51	0.44
1:L:37:GLN:OE1	1:L:47:LEU:HD22	2.18	0.44
1:L:139:PHE:HB2	1:L:172:TYR:CE2	2.53	0.44
1:L:17:GLN:HG2	1:L:18:ARG:H	1.81	0.44
1:L:47:LEU:HD21	1:L:62:PHE:CD1	2.53	0.44
1:L:85:GLU:OE1	1:L:87:TYR:CE2	2.70	0.44
2:H:150:VAL:HA	2:H:200:HIS:CB	2.48	0.44
1:L:17:GLN:HG2	1:L:18:ARG:N	2.32	0.44
3:A:29:VAL:N	4:B:101:LEU:HB3	2.33	0.44
4:B:144:CYS:O	4:B:149:MET:HE2	2.18	0.44
4:B:118:LEU:C	4:B:121:LYS:HB3	2.43	0.43
4:B:129:ASN:ND2	4:B:159:TYR:CZ	2.86	0.43
4:B:168:LEU:O	4:B:171:GLU:N	2.35	0.43
1:L:132:LEU:CD1	1:L:185:TRP:CZ2	3.01	0.43
3:A:314:LEU:HD22	4:B:100:VAL:HG21	1.99	0.43
1:L:113:PRO:HA	1:L:138:ASP:O	2.19	0.43
1:L:129:LYS:HZ2	1:L:129:LYS:HG3	1.73	0.43
1:L:181:THR:O	1:L:184:GLN:HB3	2.18	0.43
4:B:44:ALA:HA	4:B:110:PHE:HZ	1.82	0.43
4:B:127:LYS:NZ	4:B:159:TYR:CE2	2.80	0.43
1:L:54:ARG:HD2	1:L:58:VAL:O	2.18	0.43
2:H:114:ALA:HB2	2:H:146:PHE:CZ	2.54	0.43
4:B:48:ILE:O	4:B:52:VAL:HG23	2.18	0.43
2:H:80:MET:HE3	2:H:80:MET:HB3	1.85	0.43
1:L:92:ASP:CG	1:L:95:LEU:HD13	2.44	0.43
4:B:171:GLU:HG2	4:B:172:LYS:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:66:LYS:HB3	1:L:71:ALA:CB	2.47	0.43
1:L:140:TYR:CD1	1:L:171:LYS:HB3	2.54	0.43
4:B:148:CYS:O	4:B:152:VAL:HG23	2.18	0.43
4:B:166:SER:HA	4:B:169:ASN:HB2	2.01	0.43
3:A:15:ILE:HD11	4:B:121:LYS:HZ1	1.83	0.43
3:A:15:ILE:N	3:A:15:ILE:HD12	2.34	0.43
1:L:163:THR:HA	1:L:164:PRO:HD3	1.85	0.43
1:L:110:LYS:NZ	1:L:112:ALA:HA	2.33	0.42
1:L:132:LEU:CD1	1:L:185:TRP:HZ2	2.32	0.42
1:L:147:ALA:O	1:L:148:TRP:HD1	2.01	0.42
4:B:6:ILE:HG12	4:B:7:ALA:N	2.34	0.42
2:H:71:ALA:HA	2:H:78:ALA:HA	2.01	0.42
1:L:6:GLN:HE22	1:L:88:CYS:N	2.17	0.42
1:L:54:ARG:HA	1:L:55:PRO:HD3	1.84	0.42
1:L:140:TYR:CD1	1:L:171:LYS:HD2	2.47	0.42
2:H:74:PHE:HZ	3:A:41:ASN:O	2.03	0.42
2:H:156:SER:O	2:H:156:SER:OG	2.18	0.42
4:B:129:ASN:CG	4:B:159:TYR:CE1	2.98	0.42
4:B:165:GLU:HA	4:B:165:GLU:OE1	2.19	0.42
4:B:125:GLN:CD	4:B:157:TYR:HB3	2.45	0.42
1:L:168:SER:C	1:L:170:ASN:N	2.77	0.42
1:L:60:ASP:OD1	1:L:60:ASP:N	2.53	0.42
1:L:127:ALA:O	1:L:128:ASN:CG	2.63	0.42
1:L:27:ASP:O	1:L:30:ARG:HG2	2.20	0.42
1:L:133:VAL:HG12	1:L:135:LEU:CD2	2.49	0.42
4:B:121:LYS:CG	4:B:122:VAL:N	2.82	0.42
1:L:110:LYS:HZ3	1:L:197:HIS:CE1	2.37	0.42
4:B:127:LYS:CE	4:B:129:ASN:HD21	2.33	0.42
1:L:20:THR:HG23	1:L:72:SER:OG	2.19	0.41
4:B:95:ASN:O	4:B:99:LEU:HG	2.20	0.41
3:A:31:GLU:OE1	3:A:34:VAL:HG22	2.20	0.41
4:B:164:GLU:N	4:B:164:GLU:CD	2.73	0.41
1:L:148:TRP:HH2	1:L:176:SER:O	2.03	0.41
3:A:14:CYS:O	4:B:25:HIS:N	2.41	0.41
1:L:108:GLN:HG2	1:L:140:TYR:CE2	2.55	0.41
1:L:110:LYS:CE	1:L:197:HIS:ND1	2.83	0.41
2:H:175:LEU:HD12	2:H:175:LEU:HA	1.21	0.41
1:L:148:TRP:O	1:L:154:PRO:HA	2.21	0.41
1:L:48:ILE:HD11	1:L:52:ASP:HA	2.01	0.41
1:L:63:SER:C	1:L:73:LEU:HD11	2.45	0.41
1:L:66:LYS:CE	1:L:71:ALA:HB1	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:166:LYS:HE2	1:L:170:ASN:OD1	2.19	0.41
2:H:178:LEU:HD23	2:H:179:SER:H	1.86	0.41
2:H:209:LYS:O	2:H:210:ARG:NH1	2.54	0.41
1:L:135:LEU:CD2	2:H:166:PHE:CE2	2.96	0.41
2:H:144:ASP:CG	2:H:144:ASP:O	2.64	0.41
3:A:31:GLU:OE1	3:A:321:ARG:NH2	2.53	0.41
2:H:114:ALA:HB3	2:H:146:PHE:CE2	2.36	0.41
4:B:142:HIS:CD2	4:B:143:LYS:H	2.36	0.41
1:L:18:ARG:HG2	1:L:76:SER:HA	2.02	0.41
1:L:139:PHE:HD1	1:L:172:TYR:O	2.04	0.41
1:L:140:TYR:HA	1:L:172:TYR:CE2	2.53	0.41
1:L:151:ASP:OD1	1:L:189:ARG:NE	2.54	0.41
1:L:194:GLN:NE2	1:L:203:GLU:OE1	2.50	0.41
1:L:204:LYS:HD2	1:L:204:LYS:HA	1.43	0.41
2:H:12:LYS:NZ	2:H:17:SER:O	2.34	0.41
2:H:96:GLY:N	2:H:100(C):GLY:O	2.54	0.41
2:H:148:GLU:HG3	2:H:149:PRO:HB3	2.01	0.41
2:H:155:ASN:ND2	2:H:159:LEU:HD22	2.36	0.41
2:H:196:CYS:H	2:H:210:ARG:NH1	2.18	0.41
3:A:321:ARG:HD3	4:B:108:LEU:CG	2.44	0.41
4:B:165:GLU:OE1	4:B:168:LEU:HD12	2.20	0.41
1:L:110:LYS:HZ3	1:L:197:HIS:CG	2.39	0.40
2:H:11:VAL:HG11	2:H:147:PRO:CG	2.51	0.40
2:H:143:LYS:O	2:H:144:ASP:OD1	2.39	0.40
4:B:27:GLN:OE1	4:B:27:GLN:HA	2.21	0.40
3:A:317:VAL:HG12	4:B:104:ASN:OD1	2.21	0.40
4:B:143:LYS:HA	4:B:143:LYS:HD3	1.62	0.40
1:L:110:LYS:HD2	1:L:110:LYS:C	2.42	0.40
1:L:181:THR:HG22	1:L:182:PRO:CD	2.45	0.40
4:B:160:PRO:O	4:B:161:LYS:C	2.64	0.40
1:L:132:LEU:HD12	1:L:132:LEU:N	2.36	0.40
2:H:151:THR:HG22	2:H:199:ASN:O	2.21	0.40
3:A:20:ASN:ND2	3:A:37:THR:HB	2.36	0.40
4:B:127:LYS:NZ	4:B:128:ASN:OD1	2.37	0.40
4:B:30:GLN:HE22	4:B:145:ASN:CB	2.33	0.40
4:B:159:TYR:HA	4:B:161:LYS:HB2	2.02	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	L	209/215 (97%)	200 (96%)	4 (2%)	5 (2%)	4	29
2	H	213/230 (93%)	204 (96%)	6 (3%)	3 (1%)	9	38
3	A	48/66 (73%)	44 (92%)	3 (6%)	1 (2%)	5	31
4	B	150/193 (78%)	112 (75%)	25 (17%)	13 (9%)	0	7
All	All	620/704 (88%)	560 (90%)	38 (6%)	22 (4%)	3	23

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	107	GLY
1	L	128	ASN
2	H	189	LEU
2	H	204	ASN
3	A	325	SER
4	B	5	ALA
4	B	147	GLU
4	B	161	LYS
1	L	106(A)	LEU
1	L	180	LEU
1	L	189	ARG
4	B	9	PHE
4	B	121	LYS
4	B	135	ASN
4	B	157	TYR
4	B	159	TYR
4	B	171	GLU
4	B	6	ILE
4	B	7	ALA
4	B	132	GLU
2	H	149	PRO
4	B	60	ASN

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	L	175/179 (98%)	145 (83%)	30 (17%)	2	13
2	H	181/193 (94%)	165 (91%)	16 (9%)	9	33
3	A	48/56 (86%)	45 (94%)	3 (6%)	16	44
4	B	133/167 (80%)	119 (90%)	14 (10%)	6	29
All	All	537/595 (90%)	474 (88%)	63 (12%)	5	25

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	20	THR
1	L	26	SER
1	L	28	ILE
1	L	30	ARG
1	L	32	SER
1	L	33	VAL
1	L	37	GLN
1	L	45	LYS
1	L	48	ILE
1	L	53	GLN
1	L	61	ARG
1	L	63	SER
1	L	66	LYS
1	L	73	LEU
1	L	106	VAL
1	L	106(A)	LEU
1	L	108	GLN
1	L	124	GLU
1	L	136	ILE
1	L	139	PHE
1	L	144	VAL
1	L	156	LYS
1	L	161	THR
1	L	167	GLN

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Mol	Chain	Res	Type
1	L	181	THR
1	L	183	GLU
1	L	184	GLN
1	L	200	SER
1	L	204	LYS
1	L	205	THR
2	H	19	LYS
2	H	87	THR
2	H	102	VAL
2	H	116	THR
2	H	135	THR
2	H	138	LEU
2	H	143	LYS
2	H	156	SER
2	H	160	THR
2	H	163	VAL
2	H	164	HIS
2	H	173	SER
2	H	178	LEU
2	H	200	HIS
2	H	201	LYS
2	H	204	ASN
3	A	32	LYS
3	A	44	GLU
3	A	320	LEU
4	B	18	VAL
4	B	29	GLU
4	B	48	ILE
4	B	57	GLU
4	B	83	LYS
4	B	115	VAL
4	B	121	LYS
4	B	123	LYS
4	B	125	GLN
4	B	127	LYS
4	B	164	GLU
4	B	165	GLU
4	B	168	LEU
4	B	171	GLU

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

Mol	Chain	Res	Type
1	L	6	GLN
1	L	34	ASN
1	L	37	GLN
1	L	38	GLN
1	L	103	GLN
1	L	184	GLN
1	L	188	HIS
2	H	155	ASN
2	H	200	HIS
4	B	50	ASN
4	B	129	ASN
4	B	145	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	L	211/215 (98%)	0.72	19 (9%)	15 11	104, 181, 245, 333	0
2	H	217/230 (94%)	0.43	7 (3%)	50 29	74, 135, 238, 337	0
3	A	52/66 (78%)	0.39	2 (3%)	44 25	91, 123, 170, 200	0
4	B	154/193 (79%)	0.29	4 (2%)	57 33	90, 141, 191, 220	0
All	All	634/704 (90%)	0.49	32 (5%)	34 19	74, 160, 234, 337	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	165	THR	3.9
1	L	35	TRP	3.4
4	B	143	LYS	3.1
1	L	86	TYR	3.0
4	B	145	ASN	2.9
1	L	171	LYS	2.8
2	H	61	GLN	2.8
2	H	203	SER	2.7
1	L	75	ILE	2.6
1	L	197	HIS	2.6
4	B	86	GLU	2.6
2	H	137	ALA	2.6
4	B	9	PHE	2.5
1	L	144	VAL	2.4
2	H	207	VAL	2.4
1	L	202	VAL	2.4
1	L	193	CYS	2.3
1	L	137	SER	2.3
3	A	321	ARG	2.3
1	L	94	SER	2.3
1	L	191	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	L	78	LEU	2.2
2	H	163	VAL	2.2
1	L	13	GLY	2.2
1	L	112	ALA	2.1
1	L	185	TRP	2.1
3	A	30	LEU	2.1
2	H	128	SER	2.1
1	L	62	PHE	2.1
1	L	19	VAL	2.1
1	L	64	GLY	2.0
1	L	46	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.