



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:37 PM GMT

PDB ID : 1PRT
Title : THE CRYSTAL STRUCTURE OF PERTUSSIS TOXIN
Authors : Stein, P.E.; Read, R.J.
Deposited on : 1993-11-22
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

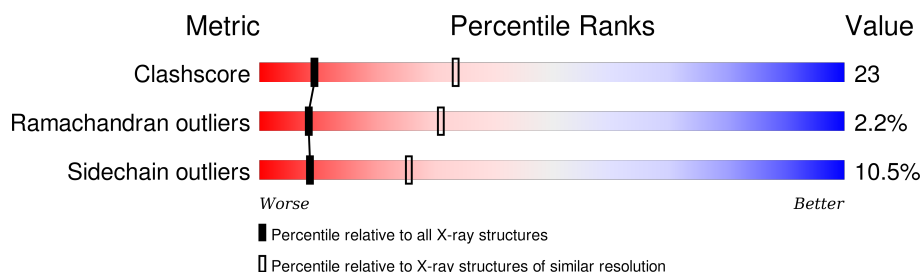
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)





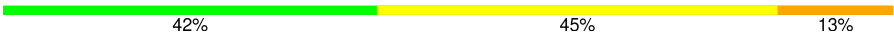
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	234	
1	G	234	
2	B	196	
2	H	196	
3	C	196	
3	I	196	
4	D	110	

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Mol	Chain	Length	Quality of chain
4	E	110	 54% 40% 5%
4	J	110	 65% 31% 5%
4	K	110	 49% 44% 7%
5	F	98	 54% 35% 10%
5	L	98	 42% 45% 13%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14504 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 PERTUSSIS TOXIN (SUBUNIT S1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	224	Total	C	N	O	S	0	0	0
			1769	1095	318	350	6			
1	G	224	Total	C	N	O	S	0	0	0
			1769	1095	318	350	6			

- Molecule 2 is a protein called PERTUSSIS TOXIN (SUBUNIT S2).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	196	Total	C	N	O	S	0	0	0
			1522	961	260	292	9			
2	H	196	Total	C	N	O	S	0	0	0
			1522	961	260	292	9			

- Molecule 3 is a protein called PERTUSSIS TOXIN (SUBUNIT S3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	196	Total	C	N	O	S	0	0	0
			1521	969	258	285	9			
3	I	196	Total	C	N	O	S	0	0	0
			1521	969	258	285	9			

- Molecule 4 is a protein called PERTUSSIS TOXIN (SUBUNIT S4).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	110	Total	C	N	O	S	0	0	0
			838	536	143	147	12			
4	E	110	Total	C	N	O	S	0	0	0
			838	536	143	147	12			
4	J	110	Total	C	N	O	S	0	0	0
			838	536	143	147	12			
4	K	110	Total	C	N	O	S	0	0	0
			838	536	143	147	12			

- Molecule 5 is a protein called PERTUSSIS TOXIN (SUBUNIT S5).

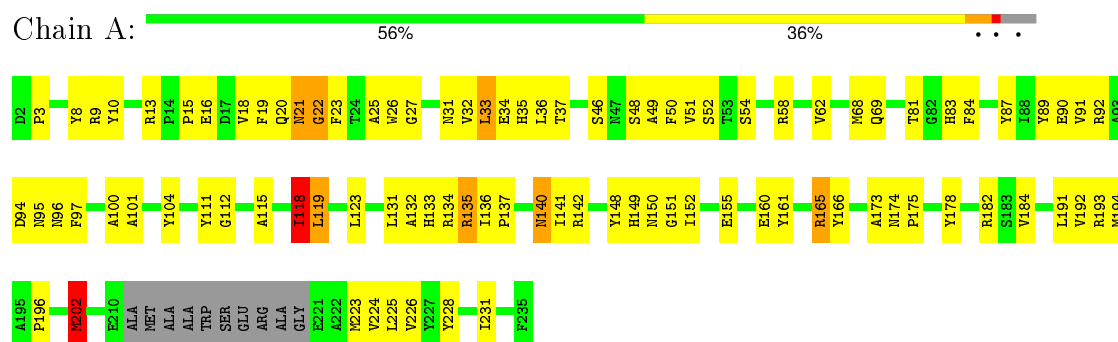
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	98	Total	C	N	O	S	0	0	0
			764	489	125	144	6			
5	L	98	Total	C	N	O	S	0	0	0
			764	489	125	144	6			

3 Residue-property plots

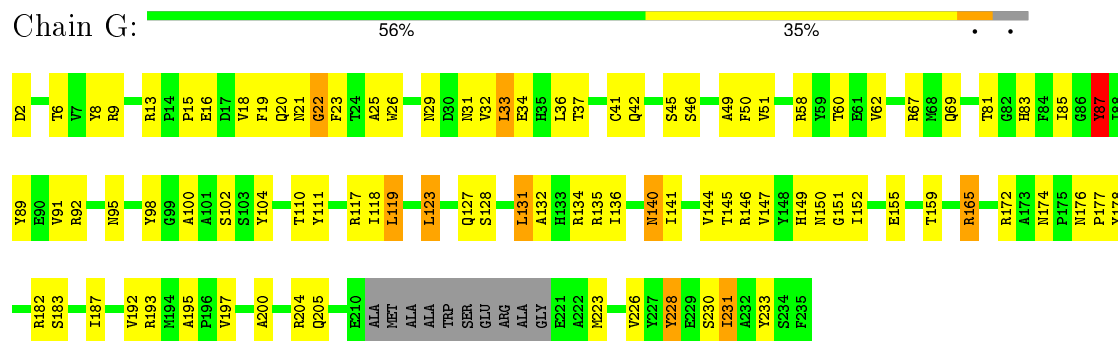
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

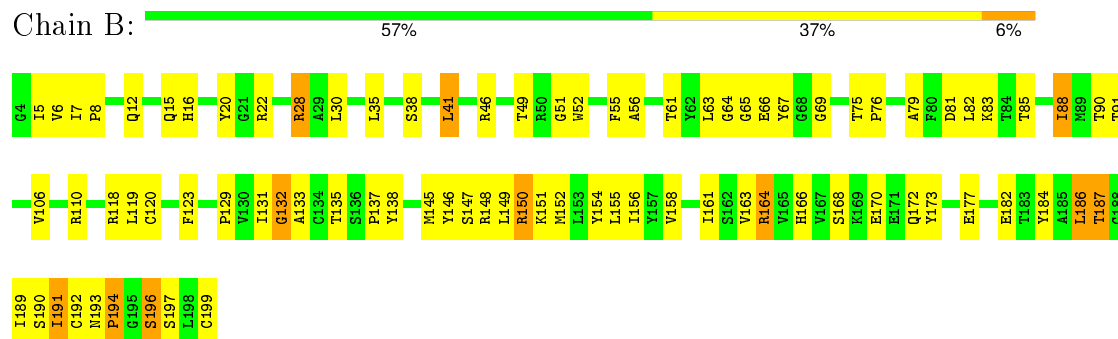
• Molecule 1: PERTUSSIS TOXIN (SUBUNIT S1)



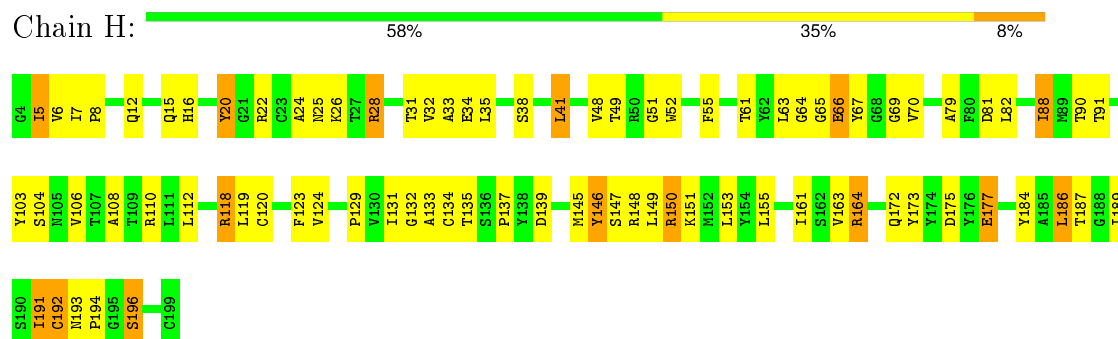
• Molecule 1: PERTUSSIS TOXIN (SUBUNIT S1)



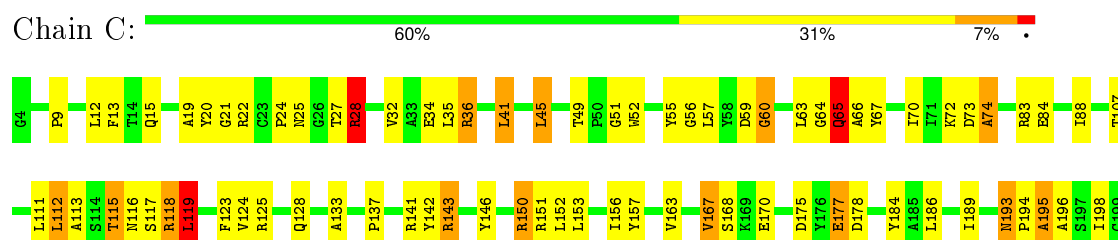
• Molecule 2: PERTUSSIS TOXIN (SUBUNIT S2)



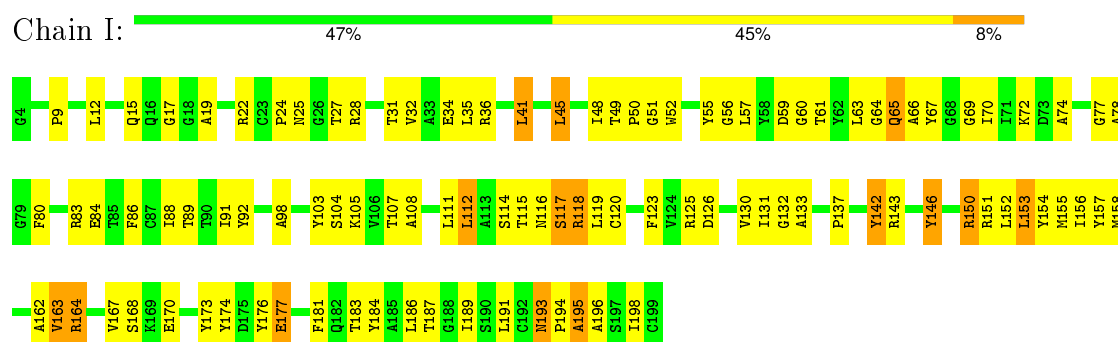
- Molecule 2: PERTUSSIS TOXIN (SUBUNIT S2)



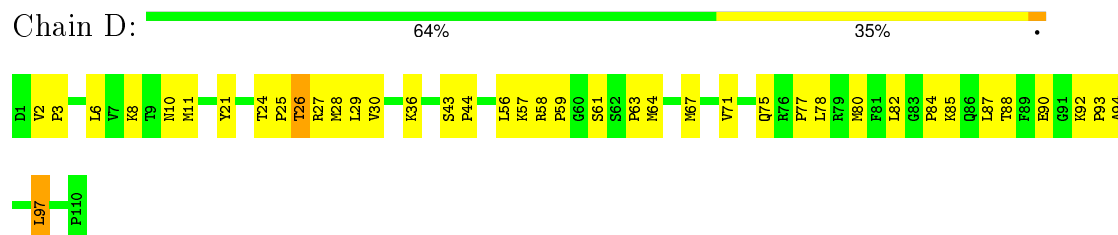
- Molecule 3: PERTUSSIS TOXIN (SUBUNIT S3)



- Molecule 3: PERTUSSIS TOXIN (SUBUNIT S3)

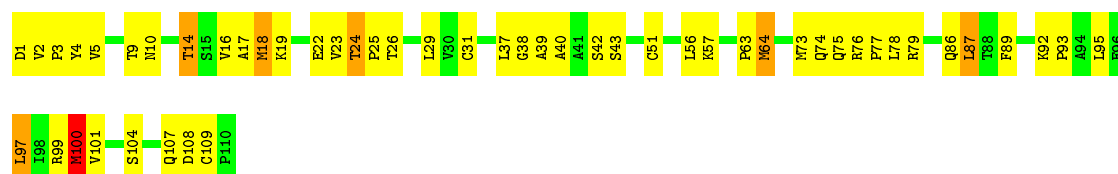


- Molecule 4: PERTUSSIS TOXIN (SUBUNIT S4)



- Molecule 4: PERTUSSIS TOXIN (SUBUNIT S4)

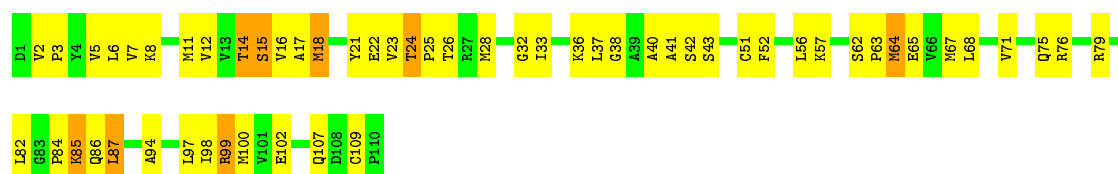




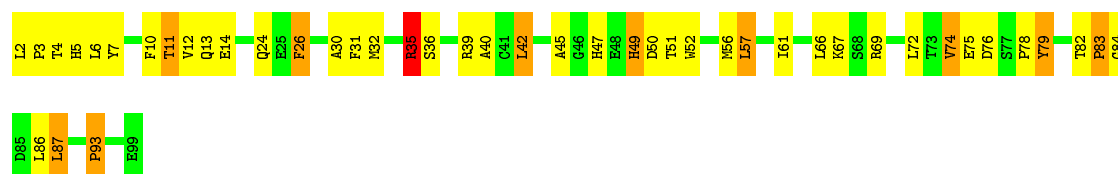
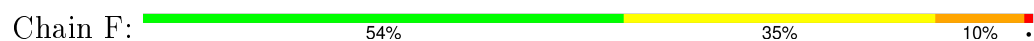
• Molecule 4: PERTUSSIS TOXIN (SUBUNIT S4)



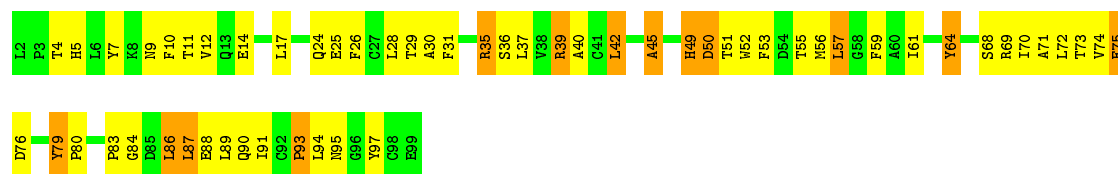
• Molecule 4: PERTUSSIS TOXIN (SUBUNIT S4)



• Molecule 5: PERTUSSIS TOXIN (SUBUNIT S5)



• Molecule 5: PERTUSSIS TOXIN (SUBUNIT S5)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	163.80 Å 98.20 Å 194.50 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.90	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.195 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14504	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.72	2/1809 (0.1%)	0.88	3/2457 (0.1%)
1	G	0.55	1/1809 (0.1%)	0.78	3/2457 (0.1%)
2	B	0.68	0/1558	0.90	2/2115 (0.1%)
2	H	0.65	0/1558	0.89	2/2115 (0.1%)
3	C	0.68	0/1557	0.87	1/2115 (0.0%)
3	I	0.64	0/1557	0.84	1/2115 (0.0%)
4	D	0.72	0/856	0.93	1/1155 (0.1%)
4	E	0.88	1/856 (0.1%)	0.98	3/1155 (0.3%)
4	J	0.66	0/856	0.89	2/1155 (0.2%)
4	K	0.62	0/856	0.87	1/1155 (0.1%)
5	F	0.69	0/782	0.91	1/1059 (0.1%)
5	L	0.61	0/782	0.88	0/1059
All	All	0.67	4/14836 (0.0%)	0.88	20/20112 (0.1%)

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	G	0	1
2	B	0	1
2	H	0	1
3	C	0	1
3	I	0	2
5	L	0	1
All	All	0	7

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	100	MET	C-N	15.84	1.70	1.34
1	A	202	MET	C-N	5.89	1.47	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	135	ARG	N-CA	5.03	1.56	1.46
1	G	134	ARG	C-O	5.00	1.32	1.23

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	100	MET	O-C-N	11.77	141.53	122.70
1	A	135	ARG	N-CA-CB	9.61	127.89	110.60
4	E	100	MET	CA-C-N	-9.22	96.92	117.20
1	G	135	ARG	NE-CZ-NH2	8.11	124.36	120.30
1	G	134	ARG	NE-CZ-NH2	8.03	124.31	120.30
2	H	191	ILE	N-CA-C	-6.21	94.22	111.00
4	J	18	MET	CG-SD-CE	6.15	110.05	100.20
4	E	100	MET	CG-SD-CE	6.13	110.01	100.20
1	A	202	MET	CG-SD-CE	6.11	109.98	100.20
2	B	191	ILE	N-CA-C	-6.05	94.65	111.00
1	A	194	MET	CG-SD-CE	5.98	109.77	100.20
4	K	18	MET	CG-SD-CE	5.98	109.77	100.20
4	D	30	VAL	N-CA-C	-5.90	95.07	111.00
2	B	132	GLY	N-CA-C	-5.66	98.96	113.10
3	C	119	LEU	N-CA-C	-5.51	96.11	111.00
2	H	192	CYS	N-CA-C	5.51	125.89	111.00
3	I	164	ARG	N-CA-C	-5.33	96.61	111.00
5	F	35	ARG	N-CA-C	5.16	124.93	111.00
4	J	78	LEU	CA-CB-CG	5.16	127.16	115.30
1	G	134	ARG	CA-C-N	5.13	128.48	117.20

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	138	TYR	Sidechain
3	C	28	ARG	Sidechain
1	G	87	TYR	Sidechain
2	H	146	TYR	Sidechain
3	I	103	TYR	Sidechain
3	I	146	TYR	Sidechain
5	L	64	TYR	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1769	0	1655	71	0
1	G	1769	0	1655	83	0
2	B	1522	0	1473	72	0
2	H	1522	0	1473	68	0
3	C	1521	0	1484	68	0
3	I	1521	0	1484	92	0
4	D	838	0	874	27	0
4	E	838	0	873	47	0
4	J	838	0	874	34	0
4	K	838	0	874	48	0
5	F	764	0	747	40	0
5	L	764	0	747	65	0
All	All	14504	0	14213	652	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 23.

All (652) 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)
4:E:100:MET:C	4:E:101:VAL:N	1.70	1.44
4:D:2:VAL:HG22	4:D:3:PRO:HD2	1.35	1.07
4:J:2:VAL:HG22	4:J:3:PRO:HD2	1.40	0.99
1:A:69:GLN:HG3	4:E:37:LEU:HD23	1.45	0.97
2:B:163:VAL:HG21	2:B:189:ILE:HG23	1.49	0.95
2:B:119:LEU:HB2	4:D:63:PRO:HB3	1.50	0.92
2:B:49:THR:HG21	2:B:64:GLY:HA3	1.55	0.89
1:G:31:ASN:HD22	1:G:34:GLU:HG3	1.37	0.88
4:K:18:MET:HE1	5:L:55:THR:CG2	2.04	0.88
1:G:31:ASN:ND2	1:G:34:GLU:HG3	1.90	0.86
1:A:54:SER:HB2	1:A:202:MET:HE3	1.58	0.85
1:A:152:ILE:H	1:A:152:ILE:HD12	1.42	0.85
3:I:119:LEU:HB2	4:K:63:PRO:HB3	1.58	0.85
1:G:91:VAL:HG11	1:G:136:ILE:HD11	1.58	0.85
4:K:18:MET:HE1	5:L:55:THR:HG22	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:PHE:CD1	1:A:131:LEU:HB3	2.13	0.83
4:E:100:MET:CA	4:E:101:VAL:N	2.41	0.83
3:I:49:THR:HG21	3:I:64:GLY:HA2	1.60	0.83
1:G:69:GLN:HG3	4:K:37:LEU:HD23	1.61	0.82
2:H:49:THR:HG21	2:H:64:GLY:HA3	1.62	0.82
1:A:54:SER:HB2	1:A:202:MET:CE	2.10	0.81
2:H:163:VAL:HG21	2:H:189:ILE:HG23	1.62	0.81
1:G:152:ILE:H	1:G:152:ILE:HD12	1.47	0.79
2:H:119:LEU:HB2	4:J:63:PRO:HB3	1.65	0.77
3:I:60:GLY:HA2	3:I:74:ALA:HB3	1.64	0.77
1:A:92:ARG:HB3	1:A:140:ASN:HD22	1.50	0.77
4:E:51:CYS:HB2	4:E:87:LEU:HD23	1.65	0.77
1:G:16:GLU:O	1:G:20:GLN:HG2	1.85	0.77
2:B:161:ILE:HG21	2:B:191:ILE:HD13	1.66	0.77
1:G:92:ARG:HB3	1:G:140:ASN:HD22	1.50	0.77
3:I:49:THR:HG23	3:I:55:TYR:HE2	1.51	0.76
2:B:145:MET:HA	5:F:24:GLN:HE22	1.50	0.76
1:A:111:TYR:CD2	1:A:182:ARG:HB2	2.20	0.76
4:K:18:MET:CE	5:L:55:THR:HG22	2.15	0.75
2:H:150:ARG:HG2	2:H:150:ARG:HH11	1.49	0.75
2:B:166:HIS:O	2:B:187:THR:HG22	1.85	0.75
1:G:111:TYR:CD2	1:G:182:ARG:HB2	2.22	0.75
2:H:137:PRO:HB3	2:H:146:TYR:CD2	2.22	0.74
1:A:91:VAL:HG11	1:A:136:ILE:HD11	1.69	0.74
1:G:31:ASN:HD21	1:G:183:SER:HB2	1.50	0.74
3:C:25:ASN:ND2	3:C:115:THR:HB	2.02	0.74
3:I:63:LEU:HD12	3:I:70:ILE:HB	1.71	0.73
1:A:31:ASN:HD22	1:A:34:GLU:HG3	1.54	0.73
2:H:148:ARG:HB3	5:L:17:LEU:HD21	1.70	0.73
1:A:31:ASN:ND2	1:A:34:GLU:HG3	2.01	0.73
2:B:63:LEU:HD12	2:B:63:LEU:H	1.53	0.72
2:B:52:TRP:CG	4:J:93:PRO:HB3	2.24	0.72
3:I:173:TYR:HB2	3:I:183:THR:HG22	1.69	0.72
2:B:193:ASN:O	2:B:196:SER:HB3	1.87	0.72
5:F:42:LEU:HD21	5:F:52:TRP:CZ3	2.24	0.72
3:C:115:THR:HG22	3:C:116:ASN:N	2.03	0.72
3:C:35:LEU:HD13	3:C:41:LEU:HD13	1.71	0.72
3:C:119:LEU:HB2	4:E:63:PRO:HB3	1.70	0.72
1:A:16:GLU:O	1:A:20:GLN:HG2	1.89	0.71
5:L:7:TYR:HB3	5:L:10:PHE:CD2	2.25	0.71
4:K:18:MET:CE	5:L:55:THR:CG2	2.69	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:63:LEU:H	2:H:63:LEU:HD12	1.57	0.70
1:G:51:VAL:HB	1:G:132:ALA:HB3	1.73	0.70
3:C:60:GLY:HA2	3:C:74:ALA:HB3	1.74	0.70
4:K:67:MET:O	4:K:71:VAL:HG22	1.91	0.70
1:A:32:VAL:O	1:A:36:LEU:HD13	1.92	0.70
4:K:62:SER:OG	4:K:65:GLU:HB2	1.92	0.69
2:B:28:ARG:HH21	2:B:177:GLU:HA	1.56	0.69
5:F:7:TYR:HB3	5:F:10:PHE:CD2	2.28	0.69
3:C:49:THR:HG21	3:C:64:GLY:HA2	1.74	0.69
3:I:25:ASN:ND2	3:I:115:THR:HB	2.08	0.69
1:A:152:ILE:N	1:A:152:ILE:HD12	2.09	0.68
1:A:95:ASN:O	1:A:174:ASN:HB2	1.94	0.68
1:G:22:GLY:HA2	1:G:141:ILE:HD11	1.76	0.68
1:A:83:HIS:CD2	1:A:150:ASN:HA	2.29	0.67
3:I:193:ASN:H	3:I:198:ILE:HB	1.58	0.67
4:D:2:VAL:CG2	4:D:3:PRO:HD2	2.18	0.67
3:C:51:GLY:HA2	3:C:65:GLN:OE1	1.95	0.67
5:F:7:TYR:HB2	5:F:72:LEU:HD23	1.77	0.66
2:B:154:TYR:O	2:B:158:VAL:HG22	1.95	0.66
3:I:56:GLY:O	3:I:57:LEU:HD12	1.95	0.66
2:B:145:MET:HB3	2:B:149:LEU:HD12	1.78	0.66
1:A:165:ARG:HH11	1:A:165:ARG:HB3	1.59	0.66
1:G:165:ARG:HH11	1:G:165:ARG:HB3	1.58	0.66
5:F:56:MET:HE2	5:F:86:LEU:HD21	1.77	0.66
1:G:195:ALA:HB1	4:K:75:GLN:HG2	1.77	0.66
2:H:155:LEU:HD22	5:L:61:ILE:HG23	1.76	0.65
5:F:74:VAL:HG23	5:F:75:GLU:O	1.95	0.65
2:H:12:GLN:O	2:H:88:ILE:HA	1.96	0.65
4:E:1:ASP:HA	4:E:86:GLN:HG3	1.79	0.65
3:I:194:PRO:HG2	3:I:195:ALA:H	1.62	0.65
2:H:35:LEU:HD12	2:H:41:LEU:HB3	1.79	0.65
1:G:83:HIS:CD2	1:G:150:ASN:HA	2.32	0.64
2:B:163:VAL:HG21	2:B:189:ILE:CG2	2.25	0.64
2:H:28:ARG:HE	2:H:177:GLU:HG3	1.62	0.64
3:C:193:ASN:H	3:C:198:ILE:HB	1.62	0.64
4:K:86:GLN:HA	4:K:94:ALA:O	1.96	0.64
5:F:6:LEU:HD23	5:F:7:TYR:N	2.12	0.64
1:A:83:HIS:HD2	1:A:150:ASN:HA	1.62	0.64
5:L:69:ARG:HG3	5:L:69:ARG:HH11	1.62	0.64
3:I:27:THR:HG22	3:I:89:THR:HA	1.80	0.64
1:G:81:THR:O	1:G:152:ILE:HD11	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:137:PRO:HB3	3:C:146:TYR:CD2	2.33	0.63
4:J:2:VAL:CG2	4:J:3:PRO:HD2	2.23	0.63
2:H:163:VAL:HG21	2:H:189:ILE:CG2	2.28	0.63
3:C:25:ASN:HD21	3:C:115:THR:HB	1.63	0.63
2:H:5:ILE:HG22	2:H:7:ILE:HD12	1.81	0.63
2:H:118:ARG:HG3	2:H:120:CYS:SG	2.39	0.63
1:A:36:LEU:CD2	1:A:104:TYR:HB2	2.29	0.63
3:C:52:TRP:HA	3:C:67:TYR:CE2	2.34	0.63
3:I:60:GLY:HA3	3:I:72:LYS:O	1.98	0.63
5:L:39:ARG:HB3	5:L:83:PRO:HG3	1.81	0.63
3:C:9:PRO:HD2	3:C:12:LEU:CD1	2.29	0.63
2:B:75:THR:HG23	2:B:76:PRO:HD2	1.80	0.63
2:H:148:ARG:H	2:H:148:ARG:HD2	1.64	0.62
2:H:28:ARG:HH21	2:H:177:GLU:HA	1.63	0.62
3:I:51:GLY:HA2	3:I:65:GLN:OE1	1.98	0.62
1:G:50:PHE:CD1	1:G:131:LEU:HB3	2.34	0.62
1:G:98:TYR:O	1:G:131:LEU:HD12	1.98	0.62
2:H:161:ILE:HG21	2:H:191:ILE:HD13	1.82	0.62
1:A:51:VAL:HB	1:A:132:ALA:HB3	1.81	0.62
5:F:10:PHE:CE1	5:F:32:MET:HB2	2.35	0.62
1:G:228:TYR:O	1:G:231:ILE:HD13	1.99	0.62
2:B:137:PRO:HB3	2:B:146:TYR:CD2	2.35	0.62
5:L:75:GLU:HG2	5:L:76:ASP:H	1.65	0.61
3:C:9:PRO:HD2	3:C:12:LEU:HD11	1.80	0.61
2:B:49:THR:HG22	2:B:65:GLY:N	2.15	0.61
5:F:7:TYR:HB3	5:F:10:PHE:CE2	2.35	0.61
4:E:79:ARG:HG3	4:E:109:CYS:SG	2.41	0.61
3:I:60:GLY:HA2	3:I:74:ALA:CB	2.30	0.61
2:H:148:ARG:HB3	5:L:17:LEU:CD2	2.31	0.61
3:I:150:ARG:HG3	3:I:151:ARG:N	2.15	0.61
3:I:52:TRP:HA	3:I:67:TYR:CE2	2.36	0.61
3:C:55:TYR:HD1	3:C:84:GLU:HB3	1.66	0.60
1:A:81:THR:O	1:A:152:ILE:HD11	2.01	0.60
3:I:49:THR:CG2	3:I:64:GLY:HA2	2.30	0.60
3:I:137:PRO:HB3	3:I:146:TYR:CD2	2.36	0.60
4:E:64:MET:CE	4:E:100:MET:CE	2.79	0.60
2:B:145:MET:HG2	5:F:24:GLN:OE1	2.01	0.60
5:L:9:ASN:OD1	5:L:97:TYR:HB3	2.02	0.60
1:A:15:PRO:HD3	1:A:87:TYR:CE2	2.37	0.60
1:G:58:ARG:O	1:G:62:VAL:HG23	2.01	0.60
2:B:150:ARG:HH11	2:B:150:ARG:HG2	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:112:LEU:HD12	3:C:113:ALA:N	2.17	0.60
3:C:52:TRP:CE2	3:C:66:ALA:HB1	2.37	0.60
1:G:36:LEU:CD2	1:G:104:TYR:HB2	2.31	0.59
3:C:63:LEU:HD12	3:C:70:ILE:HB	1.84	0.59
2:B:156:ILE:CD1	2:B:189:ILE:HD13	2.33	0.59
3:C:115:THR:HG22	3:C:116:ASN:H	1.67	0.59
4:D:80:MET:HB3	4:D:97:LEU:HD21	1.84	0.59
2:B:38:SER:OG	2:B:41:LEU:HB2	2.03	0.59
1:G:15:PRO:HD3	1:G:87:TYR:CE2	2.38	0.59
3:C:198:ILE:CD1	4:D:90:GLU:HG3	2.32	0.59
5:L:69:ARG:HG3	5:L:69:ARG:NH1	2.15	0.59
4:E:100:MET:C	4:E:101:VAL:CA	2.69	0.59
2:B:28:ARG:NH2	2:B:177:GLU:HA	2.18	0.59
4:D:26:THR:HG22	4:D:57:LYS:HG3	1.85	0.59
4:K:79:ARG:HG3	4:K:109:CYS:SG	2.43	0.58
1:G:18:VAL:HG12	1:G:141:ILE:HD13	1.85	0.58
2:H:150:ARG:HG2	2:H:150:ARG:NH1	2.18	0.58
1:G:50:PHE:HZ	1:G:178:TYR:CE2	2.22	0.58
2:B:49:THR:HG22	2:B:65:GLY:H	1.68	0.58
4:K:26:THR:HB	4:K:57:LYS:HZ2	1.69	0.58
4:E:64:MET:HE2	4:E:100:MET:HE2	1.85	0.58
4:K:26:THR:HB	4:K:57:LYS:NZ	2.19	0.58
2:H:137:PRO:HB3	2:H:146:TYR:HD2	1.68	0.58
5:F:86:LEU:HD23	5:F:86:LEU:O	2.04	0.57
1:G:26:TRP:HZ3	1:G:41:CYS:HA	1.69	0.57
2:B:30:LEU:HB2	2:B:88:ILE:HD13	1.86	0.57
2:B:51:GLY:O	2:B:65:GLY:HA3	2.05	0.57
3:C:119:LEU:CB	4:E:63:PRO:HB3	2.35	0.57
3:C:56:GLY:O	3:C:57:LEU:HD12	2.05	0.57
1:G:111:TYR:CD1	1:G:111:TYR:N	2.72	0.57
1:G:195:ALA:HB1	4:K:75:GLN:CG	2.34	0.57
3:C:168:SER:OG	3:C:170:GLU:HB2	2.04	0.57
3:I:49:THR:HG23	3:I:55:TYR:CE2	2.36	0.57
1:A:33:LEU:O	1:A:37:THR:HG22	2.03	0.57
3:I:9:PRO:HD2	3:I:12:LEU:CD1	2.35	0.56
3:C:143:ARG:HH11	3:C:143:ARG:HB2	1.70	0.56
2:H:20:TYR:HB3	4:J:4:TYR:HA	1.86	0.56
1:G:19:PHE:O	1:G:165:ARG:HD3	2.06	0.56
4:J:64:MET:CE	4:J:97:LEU:HD13	2.35	0.56
3:I:28:ARG:NH1	3:I:177:GLU:HA	2.20	0.56
3:C:142:TYR:CD1	3:C:142:TYR:N	2.73	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:163:VAL:HG22	3:I:191:LEU:HA	1.85	0.56
3:I:152:LEU:O	3:I:156:ILE:HG12	2.05	0.56
2:B:163:VAL:HG22	2:B:164:ARG:H	1.70	0.56
1:A:92:ARG:O	1:A:140:ASN:ND2	2.38	0.56
1:G:233:TYR:CZ	3:I:155:MET:HG3	2.41	0.56
1:G:233:TYR:CE2	3:I:155:MET:HG3	2.41	0.56
2:H:163:VAL:HG22	2:H:164:ARG:H	1.70	0.55
1:A:31:ASN:HD22	1:A:34:GLU:H	1.54	0.55
1:A:23:PHE:HB2	1:A:136:ILE:CG2	2.37	0.55
3:I:28:ARG:CZ	3:I:88:ILE:HD11	2.36	0.55
1:G:69:GLN:HG3	4:K:37:LEU:CD2	2.35	0.55
3:C:123:PHE:HE1	3:C:133:ALA:HB3	1.71	0.55
4:E:14:THR:HG23	5:F:93:PRO:HG3	1.88	0.55
4:K:52:PHE:CE1	4:K:68:LEU:HD22	2.42	0.55
1:A:18:VAL:HG11	1:A:89:TYR:CE2	2.41	0.55
5:L:57:LEU:HD22	5:L:61:ILE:HD11	1.89	0.55
3:I:104:SER:O	3:I:105:LYS:HB2	2.06	0.55
1:A:149:HIS:HA	1:A:155:GLU:O	2.06	0.55
4:J:37:LEU:HD12	4:J:37:LEU:N	2.21	0.55
3:C:41:LEU:HD22	3:C:45:LEU:HD22	1.88	0.55
4:K:2:VAL:HG13	4:K:3:PRO:HD2	1.88	0.54
2:H:145:MET:HA	5:L:24:GLN:HE22	1.72	0.54
2:H:106:VAL:HB	2:H:123:PHE:HB3	1.89	0.54
4:E:100:MET:HA	4:E:101:VAL:N	2.21	0.54
3:I:24:PRO:HG2	3:I:27:THR:OG1	2.06	0.54
2:H:90:THR:O	2:H:173:TYR:HA	2.08	0.54
4:J:2:VAL:HG12	4:J:84:PRO:HA	1.88	0.54
2:H:148:ARG:CB	5:L:17:LEU:HD21	2.36	0.54
1:G:26:TRP:CE3	1:G:46:SER:HB3	2.42	0.54
5:L:40:ALA:HB1	5:L:86:LEU:HB2	1.89	0.54
3:I:112:LEU:C	3:I:112:LEU:HD12	2.27	0.54
3:I:117:SER:O	4:K:63:PRO:HG2	2.08	0.54
3:I:119:LEU:CB	4:K:63:PRO:HB3	2.36	0.53
3:I:57:LEU:HD11	3:I:86:PHE:CD2	2.43	0.53
1:A:19:PHE:O	1:A:165:ARG:HD3	2.08	0.53
3:I:142:TYR:N	3:I:142:TYR:CD1	2.75	0.53
1:G:33:LEU:O	1:G:37:THR:HG22	2.09	0.53
1:A:152:ILE:H	1:A:152:ILE:CD1	2.16	0.53
2:B:149:LEU:HD22	2:B:186:LEU:HD13	1.89	0.53
5:F:75:GLU:HG2	5:F:76:ASP:H	1.73	0.53
2:B:12:GLN:O	2:B:88:ILE:HA	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:10:ASN:O	4:J:36:LYS:HA	2.09	0.53
3:I:50:PRO:HD2	3:I:55:TYR:OH	2.08	0.53
4:E:64:MET:CE	4:E:100:MET:HE3	2.39	0.53
1:G:111:TYR:HD2	1:G:182:ARG:HB2	1.72	0.53
5:L:7:TYR:HB3	5:L:10:PHE:CE2	2.43	0.53
3:I:198:ILE:CD1	4:J:90:GLU:HG3	2.39	0.53
4:D:67:MET:O	4:D:71:VAL:HG22	2.08	0.53
3:I:112:LEU:HD23	3:I:176:TYR:CE1	2.43	0.53
3:I:19:ALA:HA	3:I:83:ARG:O	2.08	0.53
4:E:51:CYS:SG	4:E:89:PHE:HB2	2.49	0.53
1:G:83:HIS:HD2	1:G:150:ASN:HA	1.74	0.53
4:E:64:MET:HE1	4:E:100:MET:HG2	1.90	0.52
1:G:111:TYR:CE2	1:G:182:ARG:HB2	2.44	0.52
2:H:15:GLN:HG3	2:H:16:HIS:H	1.74	0.52
3:I:133:ALA:HA	3:I:184:TYR:O	2.09	0.52
3:C:112:LEU:C	3:C:112:LEU:HD12	2.30	0.52
1:A:22:GLY:HA2	1:A:141:ILE:HD11	1.91	0.52
1:G:123:LEU:HD22	1:G:127:GLN:HG3	1.90	0.52
3:C:115:THR:CG2	3:C:116:ASN:N	2.72	0.52
3:I:193:ASN:O	3:I:198:ILE:HB	2.10	0.52
1:A:111:TYR:CE2	1:A:182:ARG:HB2	2.45	0.52
1:A:31:ASN:ND2	1:A:34:GLU:H	2.07	0.52
2:B:28:ARG:HE	2:B:177:GLU:HG3	1.74	0.52
2:H:147:SER:O	2:H:151:LYS:HG3	2.10	0.52
1:A:112:GLY:O	1:A:115:ALA:HB3	2.10	0.52
2:B:156:ILE:HD13	2:B:189:ILE:HD13	1.92	0.52
3:I:173:TYR:HB2	3:I:183:THR:CG2	2.39	0.52
2:H:35:LEU:CD1	2:H:41:LEU:HB3	2.39	0.52
5:L:25:GLU:HG2	5:L:26:PHE:N	2.25	0.52
2:B:163:VAL:HG22	2:B:164:ARG:N	2.25	0.52
2:B:163:VAL:HG23	2:B:190:SER:O	2.10	0.52
3:I:114:SER:HB2	3:I:117:SER:HB2	1.92	0.51
5:F:75:GLU:HB2	5:F:87:LEU:HD11	1.92	0.51
3:C:32:VAL:O	3:C:36:ARG:HB2	2.09	0.51
1:G:23:PHE:HB2	1:G:136:ILE:CG2	2.39	0.51
1:A:58:ARG:O	1:A:62:VAL:HG23	2.11	0.51
2:B:182:GLU:HG2	2:B:184:TYR:OH	2.11	0.51
2:H:52:TRP:O	2:H:67:TYR:HE2	1.93	0.51
5:L:7:TYR:O	5:L:71:ALA:HA	2.10	0.51
3:I:163:VAL:HG11	3:I:189:ILE:HG23	1.92	0.51
5:F:4:THR:HG22	5:F:5:HIS:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:55:PHE:CD1	2:H:55:PHE:N	2.78	0.51
5:F:10:PHE:CD1	5:F:32:MET:HB2	2.46	0.51
3:C:24:PRO:O	3:C:27:THR:OG1	2.27	0.51
1:G:149:HIS:HA	1:G:155:GLU:O	2.11	0.51
2:B:63:LEU:HD12	2:B:63:LEU:N	2.24	0.51
1:G:204:ARG:HA	1:G:223:MET:HE1	1.92	0.51
1:A:69:GLN:HG3	4:E:37:LEU:CD2	2.29	0.51
4:K:17:ALA:HB2	5:L:90:GLN:HB2	1.92	0.51
1:A:52:SER:OG	1:A:202:MET:HG3	2.10	0.51
3:C:32:VAL:HG12	3:C:57:LEU:HD22	1.93	0.51
5:L:68:SER:OG	5:L:91:ILE:HG23	2.11	0.51
1:A:84:PHE:O	1:A:148:TYR:HA	2.11	0.51
2:B:135:THR:CG2	2:B:146:TYR:HD1	2.24	0.51
4:K:52:PHE:CZ	4:K:68:LEU:HD22	2.46	0.51
5:L:64:TYR:CD1	5:L:64:TYR:C	2.84	0.51
2:B:135:THR:HG21	2:B:146:TYR:HA	1.93	0.50
4:J:37:LEU:HD12	4:J:37:LEU:H	1.76	0.50
5:L:72:LEU:H	5:L:72:LEU:HD23	1.75	0.50
2:H:28:ARG:NH2	2:H:177:GLU:HA	2.25	0.50
4:E:74:GLN:OE1	4:E:74:GLN:HA	2.11	0.50
2:H:193:ASN:HB2	5:L:14:GLU:OE1	2.10	0.50
1:G:50:PHE:HZ	1:G:178:TYR:CD2	2.29	0.50
2:B:148:ARG:HD2	2:B:148:ARG:H	1.75	0.50
5:L:9:ASN:HA	5:L:69:ARG:HB3	1.93	0.50
2:H:5:ILE:HG23	2:H:6:VAL:N	2.26	0.50
5:L:49:HIS:O	5:L:51:THR:N	2.45	0.50
2:H:119:LEU:HD21	2:H:153:LEU:HD23	1.94	0.50
3:I:52:TRP:H	3:I:66:ALA:CB	2.24	0.50
1:G:26:TRP:CZ3	1:G:41:CYS:HA	2.46	0.50
1:G:233:TYR:CE1	4:J:69:ARG:HD2	2.46	0.50
4:K:11:MET:SD	4:K:36:LYS:HB2	2.52	0.50
2:B:145:MET:HG2	5:F:24:GLN:CD	2.32	0.50
1:A:174:ASN:OD1	1:A:175:PRO:HD2	2.11	0.50
2:B:166:HIS:C	2:B:187:THR:HG22	2.32	0.50
5:L:75:GLU:HB2	5:L:87:LEU:HD11	1.94	0.50
2:B:55:PHE:CD1	2:B:55:PHE:N	2.79	0.50
4:K:18:MET:CE	5:L:55:THR:HG21	2.41	0.50
1:A:118:ILE:HG23	1:A:119:LEU:N	2.27	0.50
1:G:95:ASN:O	1:G:174:ASN:HB2	2.12	0.50
3:I:52:TRP:H	3:I:66:ALA:HB3	1.77	0.50
1:G:9:ARG:NH2	1:G:205:GLN:HB3	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:150:ARG:HG3	3:I:151:ARG:H	1.76	0.49
4:D:82:LEU:HD23	4:D:97:LEU:HD23	1.94	0.49
3:I:32:VAL:O	3:I:36:ARG:HB2	2.12	0.49
4:J:80:MET:HE2	4:J:82:LEU:HD21	1.94	0.49
5:F:39:ARG:HB3	5:F:83:PRO:HG3	1.94	0.49
3:I:193:ASN:N	3:I:198:ILE:HB	2.25	0.49
1:G:111:TYR:O	1:G:182:ARG:NH2	2.46	0.49
4:E:9:THR:CG2	4:E:10:ASN:N	2.75	0.49
1:A:90:GLU:HG2	1:A:142:ARG:NH1	2.28	0.49
4:D:93:PRO:HB3	2:H:52:TRP:CG	2.48	0.49
5:L:11:THR:O	5:L:30:ALA:HA	2.12	0.49
3:I:91:ILE:O	3:I:91:ILE:HG23	2.12	0.49
3:C:194:PRO:HG2	3:C:195:ALA:H	1.77	0.49
3:I:107:THR:O	3:I:123:PHE:HA	2.13	0.49
3:C:21:GLY:O	4:E:5:VAL:HG21	2.12	0.49
5:L:70:ILE:CD1	5:L:72:LEU:HD22	2.43	0.49
3:I:123:PHE:HE2	3:I:133:ALA:HB3	1.76	0.49
4:K:5:VAL:HG12	4:K:6:LEU:O	2.12	0.49
2:B:5:ILE:CG2	2:B:6:VAL:N	2.75	0.49
1:G:85:ILE:HA	1:G:147:VAL:O	2.13	0.49
4:E:64:MET:HE2	4:E:100:MET:CE	2.43	0.49
1:A:92:ARG:HB3	1:A:140:ASN:ND2	2.25	0.49
2:H:147:SER:O	2:H:151:LYS:HE2	2.12	0.49
4:K:14:THR:O	5:L:93:PRO:HG3	2.13	0.49
5:F:2:LEU:O	5:F:4:THR:N	2.46	0.49
4:D:28:MET:HG2	4:D:56:LEU:HD21	1.94	0.49
3:I:77:GLY:O	3:I:80:PHE:HB2	2.13	0.49
3:I:78:ALA:C	3:I:80:PHE:H	2.16	0.49
2:B:52:TRP:CD2	4:J:93:PRO:HB3	2.47	0.49
5:F:56:MET:CE	5:F:86:LEU:HD21	2.41	0.49
4:D:26:THR:O	4:D:27:ARG:HG2	2.13	0.49
4:K:107:GLN:H	4:K:107:GLN:CD	2.17	0.49
1:A:96:ASN:OD1	1:A:133:HIS:HB3	2.13	0.49
1:G:100:ALA:HA	1:G:131:LEU:HD11	1.95	0.48
1:A:26:TRP:CE3	1:A:46:SER:HB3	2.48	0.48
2:H:52:TRP:C	2:H:67:TYR:HE2	2.16	0.48
2:H:112:LEU:HD21	2:H:175:ASP:HB3	1.94	0.48
3:I:92:TYR:HB2	3:I:174:TYR:HE2	1.79	0.48
3:I:177:GLU:CD	4:K:79:ARG:HH22	2.17	0.48
3:I:25:ASN:HD21	3:I:115:THR:HB	1.78	0.48
4:D:64:MET:CE	4:D:97:LEU:HD13	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:150:ARG:HG3	3:C:151:ARG:N	2.28	0.48
4:K:32:GLY:C	4:K:33:ILE:HD12	2.34	0.48
4:J:86:GLN:HA	4:J:94:ALA:O	2.14	0.48
4:D:87:LEU:O	4:D:94:ALA:N	2.40	0.48
2:B:52:TRP:HE1	4:J:88:THR:HG23	1.78	0.48
3:C:49:THR:HG23	3:C:55:TYR:HE2	1.78	0.48
2:B:15:GLN:HG3	2:B:16:HIS:H	1.78	0.48
2:H:63:LEU:HD13	2:H:70:VAL:HB	1.96	0.48
2:B:118:ARG:HG3	2:B:120:CYS:SG	2.53	0.48
1:A:224:VAL:HG12	1:A:224:VAL:O	2.12	0.48
3:I:115:THR:HG22	3:I:116:ASN:N	2.28	0.48
3:I:9:PRO:HD2	3:I:12:LEU:HD11	1.95	0.48
3:I:104:SER:HA	3:I:164:ARG:HA	1.96	0.48
4:E:2:VAL:HG12	4:E:4:TYR:H	1.78	0.47
1:A:50:PHE:HD1	1:A:131:LEU:HB3	1.75	0.47
1:A:111:TYR:HD2	1:A:182:ARG:HB2	1.76	0.47
3:C:73:ASP:O	3:C:74:ALA:HB2	2.14	0.47
2:H:149:LEU:HD22	2:H:186:LEU:HD13	1.95	0.47
2:B:168:SER:OG	2:B:170:GLU:HG3	2.14	0.47
2:H:193:ASN:O	2:H:196:SER:HB3	2.13	0.47
2:H:25:ASN:O	2:H:26:LYS:HB2	2.14	0.47
4:K:98:ILE:O	4:K:99:ARG:HB2	2.14	0.47
4:E:26:THR:HB	4:E:57:LYS:NZ	2.30	0.47
4:K:51:CYS:HB2	4:K:87:LEU:HD23	1.97	0.47
5:L:73:THR:HB	5:L:88:GLU:HB3	1.96	0.47
1:A:8:TYR:N	1:A:8:TYR:CD1	2.83	0.47
5:F:32:MET:HE1	5:F:35:ARG:NH1	2.29	0.47
5:L:56:MET:HE2	5:L:86:LEU:HD21	1.96	0.47
1:G:149:HIS:CD2	1:G:151:GLY:HA2	2.49	0.47
4:E:107:GLN:H	4:E:107:GLN:CD	2.17	0.47
1:G:31:ASN:HD22	1:G:34:GLU:CG	2.16	0.47
4:K:12:VAL:HG13	4:K:37:LEU:HG	1.96	0.47
3:C:25:ASN:ND2	3:C:115:THR:CB	2.77	0.47
2:H:145:MET:HG2	5:L:24:GLN:CD	2.34	0.47
1:A:94:ASP:O	1:A:173:ALA:HA	2.13	0.47
2:B:164:ARG:CZ	2:B:199:CYS:HB3	2.44	0.47
2:H:148:ARG:N	2:H:148:ARG:HD2	2.27	0.47
1:G:87:TYR:N	1:G:87:TYR:CD1	2.81	0.47
3:I:133:ALA:HB1	3:I:186:LEU:HB2	1.96	0.47
4:K:57:LYS:HG3	5:L:49:HIS:NE2	2.30	0.47
5:L:75:GLU:HG2	5:L:76:ASP:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:108:ALA:HB2	2:H:123:PHE:CE2	2.49	0.47
4:J:16:VAL:O	4:J:17:ALA:HB2	2.13	0.47
3:I:111:LEU:HG	3:I:157:TYR:CE1	2.50	0.47
3:I:32:VAL:HG12	3:I:57:LEU:HD23	1.97	0.47
3:C:152:LEU:O	3:C:156:ILE:HG12	2.15	0.47
4:D:43:SER:HA	4:D:44:PRO:HD3	1.77	0.47
4:K:24:THR:HA	4:K:25:PRO:HD2	1.81	0.47
5:L:29:THR:HG23	5:L:37:LEU:HB3	1.97	0.47
1:G:92:ARG:CZ	1:G:140:ASN:ND2	2.78	0.46
3:C:107:THR:O	3:C:123:PHE:HA	2.15	0.46
5:L:57:LEU:HD22	5:L:61:ILE:CD1	2.44	0.46
2:B:41:LEU:HD12	2:B:41:LEU:HA	1.80	0.46
1:G:2:ASP:HB3	1:G:172:ARG:NH1	2.30	0.46
5:L:75:GLU:CG	5:L:76:ASP:H	2.27	0.46
1:A:50:PHE:HZ	1:A:178:TYR:CD2	2.33	0.46
1:G:32:VAL:O	1:G:36:LEU:HD13	2.15	0.46
2:H:51:GLY:O	2:H:65:GLY:HA3	2.14	0.46
5:F:35:ARG:HB2	5:F:36:SER:H	1.46	0.46
3:C:167:VAL:HG22	3:C:184:TYR:HB2	1.96	0.46
4:E:99:ARG:O	4:E:100:MET:HB3	2.16	0.46
5:F:7:TYR:HB2	5:F:72:LEU:CD2	2.46	0.46
5:L:7:TYR:HE1	5:L:74:VAL:HG22	1.81	0.46
3:C:60:GLY:HA3	3:C:72:LYS:O	2.16	0.46
5:F:78:PRO:HG2	5:F:79:TYR:CE2	2.51	0.46
1:G:8:TYR:CD1	1:G:8:TYR:N	2.84	0.46
3:I:154:TYR:CE1	3:I:158:MET:HG3	2.51	0.46
5:F:57:LEU:O	5:F:61:ILE:HG13	2.16	0.46
4:K:18:MET:HE1	5:L:55:THR:HG21	1.91	0.46
3:C:13:PHE:CD2	3:C:41:LEU:HD21	2.52	0.46
1:G:19:PHE:HZ	1:G:144:VAL:HG12	1.81	0.46
5:L:14:GLU:O	5:L:28:LEU:HA	2.16	0.45
3:I:63:LEU:O	3:I:69:GLY:HA2	2.16	0.45
1:G:200:ALA:O	1:G:204:ARG:HB2	2.16	0.45
5:F:69:ARG:HH11	5:F:69:ARG:HG3	1.81	0.45
3:I:45:LEU:HA	3:I:48:ILE:HG22	1.98	0.45
3:I:12:LEU:HD22	3:I:89:THR:O	2.17	0.45
2:B:145:MET:HB3	2:B:149:LEU:CD1	2.43	0.45
2:B:52:TRP:O	2:B:67:TYR:HE2	2.00	0.45
4:J:87:LEU:O	4:J:94:ALA:N	2.50	0.45
4:E:97:LEU:HD11	4:E:100:MET:SD	2.55	0.45
5:L:53:PHE:O	5:L:57:LEU:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:28:ARG:NH1	3:C:177:GLU:HA	2.31	0.45
4:J:64:MET:HE1	4:J:97:LEU:HD13	1.98	0.45
5:L:26:PHE:CE1	5:L:42:LEU:HB3	2.52	0.45
4:E:17:ALA:N	4:E:31:CYS:O	2.49	0.45
4:E:40:ALA:C	4:E:42:SER:N	2.68	0.45
1:G:6:THR:HA	1:G:91:VAL:O	2.17	0.45
3:C:52:TRP:H	3:C:66:ALA:HB3	1.82	0.45
2:B:90:THR:O	2:B:173:TYR:HA	2.16	0.45
1:A:25:ALA:HB1	1:A:48:SER:O	2.16	0.45
2:B:131:ILE:HG13	2:B:132:GLY:H	1.82	0.45
1:G:110:THR:HB	1:G:111:TYR:CD1	2.52	0.45
3:I:28:ARG:NH2	3:I:34:GLU:OE1	2.50	0.45
2:B:155:LEU:HD22	5:F:61:ILE:HG23	1.99	0.45
2:B:156:ILE:HD11	2:B:189:ILE:HD13	1.98	0.45
5:F:49:HIS:O	5:F:51:THR:N	2.50	0.45
1:A:191:LEU:HA	1:A:191:LEU:HD23	1.71	0.45
2:H:193:ASN:HB2	5:L:14:GLU:CD	2.37	0.45
1:A:27:GLY:O	1:A:49:ALA:HA	2.17	0.45
3:C:141:ARG:HB3	3:C:142:TYR:CD1	2.52	0.44
2:H:145:MET:HB2	2:H:149:LEU:HD12	1.98	0.44
1:A:68:MET:HB2	1:A:84:PHE:CE2	2.52	0.44
3:I:118:ARG:HG3	3:I:120:CYS:SG	2.57	0.44
4:E:64:MET:CE	4:E:100:MET:HE2	2.45	0.44
5:F:75:GLU:N	5:F:87:LEU:HD11	2.32	0.44
3:I:153:LEU:HA	3:I:153:LEU:HD23	1.76	0.44
2:B:131:ILE:HD12	2:B:131:ILE:HA	1.85	0.44
4:J:28:MET:SD	4:J:56:LEU:HD21	2.57	0.44
3:C:28:ARG:CZ	3:C:88:ILE:HD11	2.47	0.44
4:J:11:MET:SD	4:J:36:LYS:HB2	2.58	0.44
2:H:52:TRP:CH2	2:H:66:GLU:HB3	2.52	0.44
4:K:64:MET:HE1	4:K:100:MET:HG2	1.99	0.44
2:H:79:ALA:C	2:H:81:ASP:H	2.19	0.44
1:A:149:HIS:CD2	1:A:151:GLY:HA2	2.53	0.44
1:A:111:TYR:CD1	1:A:111:TYR:N	2.86	0.44
2:H:7:ILE:HA	2:H:8:PRO:HD3	1.85	0.44
3:I:31:THR:HG21	3:I:177:GLU:HB2	2.00	0.44
5:L:88:GLU:HG2	5:L:89:LEU:N	2.32	0.44
4:E:40:ALA:O	4:E:42:SER:N	2.50	0.44
1:G:102:SER:HB3	1:G:177:PRO:HG3	2.00	0.44
3:I:35:LEU:HD13	3:I:41:LEU:HD13	1.99	0.44
5:L:74:VAL:O	5:L:74:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:167:VAL:CG2	3:C:184:TYR:HB2	2.47	0.44
3:C:34:GLU:HG3	3:C:178:ASP:OD1	2.18	0.44
1:G:144:VAL:HG22	1:G:145:THR:N	2.33	0.44
4:J:37:LEU:CD1	4:J:37:LEU:H	2.31	0.44
4:D:92:LYS:HB3	4:D:93:PRO:HD2	1.98	0.44
1:G:25:ALA:HB1	1:G:49:ALA:HA	1.98	0.44
2:B:56:ALA:O	2:B:85:THR:HA	2.18	0.44
3:C:118:ARG:H	3:C:118:ARG:HG2	1.57	0.44
4:J:81:PHE:CD1	4:J:81:PHE:N	2.85	0.44
4:E:24:THR:HA	4:E:25:PRO:HD2	1.82	0.44
4:E:40:ALA:C	4:E:42:SER:H	2.21	0.44
4:E:76:ARG:HA	4:E:77:PRO:HD3	1.71	0.44
2:H:133:ALA:HA	2:H:184:TYR:O	2.17	0.44
5:F:40:ALA:HA	5:F:84:GLY:O	2.18	0.44
4:K:18:MET:HE2	5:L:55:THR:HG22	1.99	0.43
2:B:35:LEU:HD12	2:B:41:LEU:HB3	2.00	0.43
1:A:142:ARG:HB2	1:A:166:TYR:CE1	2.53	0.43
4:K:15:SER:OG	5:L:93:PRO:HD3	2.17	0.43
4:K:8:LYS:HG2	4:K:82:LEU:HD12	1.99	0.43
1:A:13:ARG:HH11	1:A:13:ARG:HG3	1.83	0.43
4:D:75:GLN:NE2	4:D:75:GLN:HA	2.33	0.43
5:L:94:LEU:O	5:L:95:ASN:HB2	2.19	0.43
4:J:10:ASN:ND2	4:J:109:CYS:HB3	2.33	0.43
2:B:133:ALA:HA	2:B:184:TYR:O	2.19	0.43
1:A:34:GLU:HA	1:A:37:THR:HG22	2.00	0.43
5:L:10:PHE:O	5:L:70:ILE:HG12	2.19	0.43
3:I:78:ALA:O	3:I:80:PHE:N	2.51	0.43
4:D:11:MET:O	4:D:78:LEU:N	2.48	0.43
4:E:18:MET:HE2	4:E:18:MET:HB3	1.84	0.43
5:F:4:THR:HG22	5:F:5:HIS:H	1.84	0.43
4:D:24:THR:HA	4:D:25:PRO:HD3	1.70	0.43
1:A:20:GLN:HB2	1:A:21:ASN:OD1	2.19	0.43
3:C:177:GLU:CD	4:E:79:ARG:HH22	2.21	0.43
4:E:92:LYS:HB3	4:E:93:PRO:HD2	2.00	0.43
1:G:92:ARG:CZ	1:G:140:ASN:HD21	2.31	0.43
2:H:38:SER:OG	2:H:41:LEU:HB2	2.17	0.43
2:H:119:LEU:O	2:H:134:CYS:HA	2.18	0.43
2:B:193:ASN:HB2	5:F:14:GLU:OE1	2.19	0.43
4:D:8:LYS:HG3	4:D:82:LEU:HD12	2.00	0.43
3:I:107:THR:HG22	3:I:108:ALA:N	2.34	0.43
4:D:85:LYS:HB3	4:D:85:LYS:HE3	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:ALA:C	2:B:81:ASP:H	2.21	0.43
2:H:5:ILE:HG22	2:H:7:ILE:CD1	2.48	0.43
4:K:21:TYR:HB3	4:K:24:THR:O	2.19	0.43
4:D:11:MET:SD	4:D:36:LYS:HB2	2.59	0.43
3:C:111:LEU:HG	3:C:157:TYR:CE2	2.54	0.43
3:C:141:ARG:HG2	3:C:142:TYR:CE1	2.54	0.43
5:L:40:ALA:HA	5:L:84:GLY:O	2.19	0.43
2:H:32:VAL:HG12	2:H:33:ALA:N	2.33	0.43
1:G:13:ARG:O	1:G:89:TYR:OH	2.30	0.42
3:I:32:VAL:HG12	3:I:57:LEU:CD2	2.49	0.42
3:C:198:ILE:HD12	4:D:90:GLU:HG3	1.99	0.42
3:I:167:VAL:CG2	3:I:184:TYR:HB2	2.49	0.42
4:D:28:MET:SD	4:D:56:LEU:HD21	2.59	0.42
4:E:26:THR:O	4:E:56:LEU:HB2	2.19	0.42
4:K:21:TYR:N	4:K:28:MET:HA	2.33	0.42
3:I:181:PHE:CD1	3:I:181:PHE:C	2.92	0.42
4:J:104:SER:HB3	4:J:108:ASP:OD1	2.19	0.42
3:I:12:LEU:HD11	3:I:174:TYR:CE2	2.54	0.42
3:I:45:LEU:HD12	3:I:45:LEU:HA	1.66	0.42
2:H:131:ILE:HG13	2:H:132:GLY:H	1.84	0.42
1:G:195:ALA:HB2	5:L:94:LEU:HD12	2.01	0.42
4:E:107:GLN:CD	4:E:107:GLN:N	2.72	0.42
1:G:176:ASN:HA	1:G:177:PRO:HD3	1.83	0.42
5:L:35:ARG:HB2	5:L:36:SER:H	1.54	0.42
2:B:75:THR:CG2	2:B:76:PRO:HD2	2.48	0.42
3:I:142:TYR:CE2	3:I:187:THR:HG22	2.54	0.42
1:A:3:PRO:HB3	1:A:173:ALA:HB3	2.01	0.42
4:D:10:ASN:O	4:D:36:LYS:HA	2.19	0.42
4:K:76:ARG:HD2	4:K:102:GLU:OE2	2.19	0.42
2:H:31:THR:OG1	2:H:34:GLU:HB2	2.20	0.42
3:C:186:LEU:HD21	3:C:189:ILE:HD11	2.01	0.42
1:G:152:ILE:HA	4:K:41:ALA:HB3	2.02	0.42
3:C:60:GLY:HA2	3:C:74:ALA:CB	2.47	0.42
1:G:18:VAL:HG11	1:G:89:TYR:CE2	2.54	0.42
2:H:5:ILE:CG2	2:H:6:VAL:N	2.81	0.42
3:I:162:ALA:O	3:I:163:VAL:HG23	2.19	0.42
4:J:10:ASN:HB3	4:J:37:LEU:HD13	2.01	0.42
2:B:15:GLN:HG3	2:B:16:HIS:N	2.34	0.42
4:K:28:MET:SD	4:K:56:LEU:HD21	2.60	0.42
4:D:11:MET:O	4:D:77:PRO:HA	2.20	0.42
3:C:19:ALA:HA	3:C:83:ARG:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:193:ARG:O	1:G:197:VAL:HA	2.20	0.42
3:I:194:PRO:C	3:I:196:ALA:H	2.22	0.42
2:B:147:SER:O	2:B:151:LYS:HG3	2.20	0.42
1:A:95:ASN:C	1:A:97:PHE:H	2.23	0.42
3:I:193:ASN:HA	3:I:194:PRO:HD2	1.90	0.42
3:I:142:TYR:HD1	3:I:142:TYR:N	2.15	0.42
4:E:74:GLN:O	4:E:75:GLN:HB2	2.20	0.42
4:J:50:PHE:N	4:J:50:PHE:CD1	2.86	0.42
4:J:82:LEU:HB3	4:J:95:LEU:HB3	2.01	0.42
5:F:11:THR:O	5:F:30:ALA:HA	2.20	0.42
1:G:117:ARG:HD3	1:G:117:ARG:HA	1.89	0.42
1:A:54:SER:CB	1:A:202:MET:CE	2.92	0.42
3:C:25:ASN:HD21	3:C:115:THR:CB	2.29	0.42
1:G:87:TYR:HB2	1:G:89:TYR:CE1	2.55	0.42
3:I:131:ILE:HG13	3:I:132:GLY:H	1.85	0.42
5:L:4:THR:HG22	5:L:5:HIS:N	2.35	0.42
2:B:182:GLU:HB3	2:B:184:TYR:CE2	2.55	0.42
4:E:2:VAL:HG13	4:E:3:PRO:HD2	2.02	0.42
2:B:46:ARG:HH11	2:B:46:ARG:HG3	1.85	0.42
2:B:106:VAL:HB	2:B:123:PHE:HB3	2.00	0.42
3:C:45:LEU:HA	3:C:45:LEU:HD12	1.72	0.41
5:F:10:PHE:CE1	5:F:32:MET:CB	3.01	0.41
1:G:13:ARG:HH11	1:G:13:ARG:HG3	1.85	0.41
2:H:41:LEU:HD12	2:H:41:LEU:HA	1.94	0.41
4:J:51:CYS:SG	4:J:89:PHE:HB2	2.59	0.41
4:J:40:ALA:O	4:J:46:ALA:HB2	2.19	0.41
2:B:83:LYS:HE3	4:J:2:VAL:O	2.20	0.41
1:A:32:VAL:HG12	1:A:32:VAL:O	2.21	0.41
2:H:106:VAL:HA	2:H:124:VAL:O	2.20	0.41
2:B:118:ARG:NH1	2:B:173:TYR:CE1	2.89	0.41
5:L:50:ASP:C	5:L:52:TRP:N	2.72	0.41
2:H:103:TYR:N	2:H:103:TYR:CD1	2.88	0.41
1:A:225:LEU:HD23	1:A:225:LEU:HA	1.86	0.41
1:A:69:GLN:HE21	4:E:38:GLY:H	1.67	0.41
3:C:25:ASN:ND2	3:C:115:THR:CG2	2.84	0.41
2:B:38:SER:CB	2:B:41:LEU:HB2	2.51	0.41
3:C:194:PRO:C	3:C:196:ALA:H	2.24	0.41
3:I:55:TYR:HD1	3:I:84:GLU:HB3	1.84	0.41
5:L:74:VAL:O	5:L:74:VAL:CG2	2.69	0.41
2:H:15:GLN:HG3	2:H:16:HIS:N	2.36	0.41
5:L:59:PHE:CE2	5:L:91:ILE:HG13	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:20:TYR:CD1	3:C:83:ARG:HA	2.55	0.41
4:K:85:LYS:H	4:K:85:LYS:HD3	1.84	0.41
1:A:100:ALA:HA	1:A:131:LEU:HD11	2.02	0.41
2:B:52:TRP:CD1	4:J:93:PRO:HB3	2.54	0.41
4:K:87:LEU:O	4:K:94:ALA:N	2.53	0.41
5:L:39:ARG:HB3	5:L:83:PRO:CG	2.50	0.41
3:C:21:GLY:C	4:E:5:VAL:HG21	2.41	0.41
3:I:168:SER:C	3:I:170:GLU:H	2.24	0.41
3:C:56:GLY:C	3:C:57:LEU:HD12	2.41	0.41
2:H:175:ASP:O	4:J:99:ARG:NH2	2.54	0.41
1:A:35:HIS:ND1	1:A:50:PHE:HB2	2.35	0.41
1:A:50:PHE:HZ	1:A:178:TYR:CE2	2.39	0.41
1:G:152:ILE:HD12	1:G:152:ILE:N	2.26	0.41
1:G:110:THR:CG2	1:G:111:TYR:CE1	3.04	0.41
5:L:97:TYR:CD1	5:L:97:TYR:N	2.88	0.41
3:C:28:ARG:HD2	3:C:175:ASP:O	2.21	0.41
1:G:37:THR:HG21	1:G:187:ILE:HD13	2.02	0.41
2:B:152:MET:O	2:B:156:ILE:HG12	2.21	0.41
5:F:26:PHE:HE1	5:F:42:LEU:HD12	1.85	0.41
1:G:98:TYR:CE2	1:G:174:ASN:ND2	2.89	0.41
3:C:63:LEU:CD1	3:C:70:ILE:HB	2.48	0.41
4:K:107:GLN:N	4:K:107:GLN:CD	2.74	0.41
4:E:2:VAL:HG21	4:E:95:LEU:HD21	2.03	0.41
5:F:79:TYR:O	5:F:82:THR:HB	2.21	0.41
2:H:131:ILE:HD11	2:H:184:TYR:CD2	2.56	0.41
3:I:125:ARG:HB2	3:I:130:VAL:HG21	2.03	0.41
1:G:119:LEU:HD12	1:G:119:LEU:HA	1.81	0.41
1:G:146:ARG:HB3	1:G:159:THR:OG1	2.21	0.41
3:C:125:ARG:O	3:C:128:GLN:HG2	2.20	0.41
1:A:160:GLU:C	1:A:161:TYR:CD1	2.94	0.41
1:A:193:ARG:HH11	1:A:223:MET:HG3	1.86	0.41
4:K:40:ALA:O	4:K:42:SER:N	2.54	0.41
1:G:127:GLN:O	1:G:128:SER:C	2.60	0.41
4:E:19:LYS:CG	4:E:29:LEU:HB2	2.51	0.41
4:E:64:MET:HB3	4:E:64:MET:HE2	1.76	0.40
5:L:51:THR:HG22	5:L:55:THR:OG1	2.21	0.40
1:G:149:HIS:HE1	4:K:38:GLY:O	2.04	0.40
3:C:124:VAL:HG13	3:C:128:GLN:O	2.21	0.40
4:E:104:SER:OG	4:E:108:ASP:OD1	2.39	0.40
1:G:152:ILE:CD1	1:G:152:ILE:H	2.23	0.40
2:H:177:GLU:CD	4:J:79:ARG:NH2	2.74	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:168:SER:C	3:C:170:GLU:H	2.24	0.40
3:I:112:LEU:HD23	3:I:176:TYR:CD1	2.56	0.40
3:I:167:VAL:HG22	3:I:184:TYR:HB2	2.03	0.40
3:I:131:ILE:HG13	3:I:132:GLY:N	2.37	0.40
1:G:42:GLN:HB3	1:G:45:SER:HB3	2.04	0.40
5:L:79:TYR:HA	5:L:80:PRO:HD2	1.89	0.40
4:D:2:VAL:HG12	4:D:84:PRO:HA	2.02	0.40
2:H:104:SER:O	2:H:106:VAL:HG13	2.20	0.40
3:I:108:ALA:HB2	3:I:123:PHE:CE1	2.56	0.40
5:L:45:ALA:HB3	5:L:50:ASP:OD2	2.21	0.40
2:B:7:ILE:HA	2:B:8:PRO:HD3	1.83	0.40
4:E:73:MET:CE	5:F:66:LEU:HD11	2.51	0.40
5:F:66:LEU:O	5:F:67:LYS:HB2	2.22	0.40
1:G:67:ARG:HD2	1:G:67:ARG:HA	1.91	0.40
2:B:193:ASN:HB2	5:F:14:GLU:CD	2.42	0.40
3:I:153:LEU:HD23	3:I:156:ILE:HD11	2.04	0.40
5:L:26:PHE:HE1	5:L:42:LEU:HD13	1.86	0.40
1:G:204:ARG:HA	1:G:223:MET:CE	2.51	0.40
5:F:78:PRO:HG2	5:F:79:TYR:CZ	2.56	0.40
1:A:101:ALA:O	1:A:104:TYR:N	2.54	0.40
3:I:48:ILE:HG13	3:I:48:ILE:O	2.21	0.40
4:D:58:ARG:HA	4:D:59:PRO:HD3	1.83	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	220/234 (94%)	197 (90%)	20 (9%)	3 (1%)	14	44
1	G	220/234 (94%)	194 (88%)	24 (11%)	2 (1%)	21	57
2	B	194/196 (99%)	176 (91%)	15 (8%)	3 (2%)	13	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	194/196 (99%)	173 (89%)	17 (9%)	4 (2%)	9	32
3	C	194/196 (99%)	168 (87%)	21 (11%)	5 (3%)	7	26
3	I	194/196 (99%)	171 (88%)	18 (9%)	5 (3%)	7	26
4	D	108/110 (98%)	99 (92%)	9 (8%)	0	100	100
4	E	108/110 (98%)	95 (88%)	9 (8%)	4 (4%)	4	17
4	J	108/110 (98%)	98 (91%)	10 (9%)	0	100	100
4	K	108/110 (98%)	97 (90%)	5 (5%)	6 (6%)	2	7
5	F	96/98 (98%)	83 (86%)	8 (8%)	5 (5%)	2	8
5	L	96/98 (98%)	81 (84%)	11 (12%)	4 (4%)	3	13
All	All	1840/1888 (98%)	1632 (89%)	167 (9%)	41 (2%)	8	31

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	F	45	ALA
5	F	49	HIS
5	L	49	HIS
5	L	50	ASP
2	B	110	ARG
3	C	65	GLN
4	E	22	GLU
5	F	47	HIS
5	F	50	ASP
1	G	22	GLY
1	G	29	ASN
2	H	48	VAL
2	H	69	GLY
3	I	126	ASP
4	K	22	GLU
4	K	85	LYS
2	B	69	GLY
3	C	59	ASP
4	E	100	MET
2	H	24	ALA
3	I	98	ALA
1	A	135	ARG
3	C	195	ALA
2	H	110	ARG
3	I	59	ASP

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Mol	Chain	Res	Type
5	L	45	ALA
5	L	75	GLU
2	B	194	PRO
3	C	60	GLY
4	E	39	ALA
3	I	195	ALA
3	I	17	GLY
4	K	99	ARG
4	E	23	VAL
1	A	22	GLY
1	A	118	ILE
5	F	3	PRO
4	K	7	VAL
4	K	23	VAL
3	C	74	ALA
4	K	84	PRO

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	185/190 (97%)	167 (90%)	18 (10%)	10	30
1	G	185/190 (97%)	170 (92%)	15 (8%)	15	39
2	B	163/163 (100%)	144 (88%)	19 (12%)	7	19
2	H	163/163 (100%)	140 (86%)	23 (14%)	4	12
3	C	155/155 (100%)	136 (88%)	19 (12%)	6	17
3	I	155/155 (100%)	139 (90%)	16 (10%)	9	26
4	D	94/94 (100%)	87 (93%)	7 (7%)	17	44
4	E	94/94 (100%)	85 (90%)	9 (10%)	10	31
4	J	94/94 (100%)	88 (94%)	6 (6%)	22	53
4	K	94/94 (100%)	86 (92%)	8 (8%)	13	37
5	F	83/83 (100%)	70 (84%)	13 (16%)	3	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	L	83/83 (100%)	73 (88%)	10 (12%)	6	18
All	All	1548/1558 (99%)	1385 (90%)	163 (10%)	8	25

All (163) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	10	TYR
1	A	21	ASN
1	A	33	LEU
1	A	118	ILE
1	A	119	LEU
1	A	123	LEU
1	A	134	ARG
1	A	137	PRO
1	A	140	ASN
1	A	165	ARG
1	A	184	VAL
1	A	192	VAL
1	A	196	PRO
1	A	202	MET
1	A	226	VAL
1	A	228	TYR
1	A	231	ILE
2	B	20	TYR
2	B	22	ARG
2	B	28	ARG
2	B	41	LEU
2	B	61	THR
2	B	66	GLU
2	B	82	LEU
2	B	88	ILE
2	B	91	THR
2	B	129	PRO
2	B	150	ARG
2	B	164	ARG
2	B	172	GLN
2	B	186	LEU
2	B	187	THR
2	B	192	CYS
2	B	194	PRO
2	B	196	SER

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Mol	Chain	Res	Type
2	B	197	SER
3	C	15	GLN
3	C	22	ARG
3	C	28	ARG
3	C	36	ARG
3	C	41	LEU
3	C	45	LEU
3	C	65	GLN
3	C	112	LEU
3	C	115	THR
3	C	117	SER
3	C	118	ARG
3	C	119	LEU
3	C	143	ARG
3	C	150	ARG
3	C	153	LEU
3	C	163	VAL
3	C	167	VAL
3	C	177	GLU
3	C	193	ASN
4	D	6	LEU
4	D	21	TYR
4	D	26	THR
4	D	29	LEU
4	D	61	SER
4	D	88	THR
4	D	97	LEU
4	E	14	THR
4	E	16	VAL
4	E	18	MET
4	E	24	THR
4	E	43	SER
4	E	64	MET
4	E	78	LEU
4	E	87	LEU
4	E	97	LEU
5	F	11	THR
5	F	12	VAL
5	F	13	GLN
5	F	26	PHE
5	F	31	PHE
5	F	35	ARG

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Mol	Chain	Res	Type
5	F	42	LEU
5	F	57	LEU
5	F	74	VAL
5	F	79	TYR
5	F	83	PRO
5	F	87	LEU
5	F	93	PRO
1	G	21	ASN
1	G	33	LEU
1	G	60	THR
1	G	87	TYR
1	G	118	ILE
1	G	119	LEU
1	G	123	LEU
1	G	131	LEU
1	G	140	ASN
1	G	165	ARG
1	G	192	VAL
1	G	226	VAL
1	G	228	TYR
1	G	230	SER
1	G	231	ILE
2	H	5	ILE
2	H	20	TYR
2	H	22	ARG
2	H	28	ARG
2	H	41	LEU
2	H	61	THR
2	H	66	GLU
2	H	82	LEU
2	H	88	ILE
2	H	91	THR
2	H	118	ARG
2	H	129	PRO
2	H	135	THR
2	H	139	ASP
2	H	150	ARG
2	H	164	ARG
2	H	172	GLN
2	H	177	GLU
2	H	186	LEU
2	H	187	THR

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Mol	Chain	Res	Type
2	H	192	CYS
2	H	194	PRO
2	H	196	SER
3	I	15	GLN
3	I	22	ARG
3	I	41	LEU
3	I	45	LEU
3	I	61	THR
3	I	65	GLN
3	I	112	LEU
3	I	117	SER
3	I	118	ARG
3	I	142	TYR
3	I	143	ARG
3	I	150	ARG
3	I	153	LEU
3	I	163	VAL
3	I	177	GLU
3	I	193	ASN
4	J	2	VAL
4	J	10	ASN
4	J	26	THR
4	J	79	ARG
4	J	88	THR
4	J	97	LEU
4	K	14	THR
4	K	15	SER
4	K	16	VAL
4	K	24	THR
4	K	43	SER
4	K	64	MET
4	K	87	LEU
4	K	97	LEU
5	L	12	VAL
5	L	31	PHE
5	L	35	ARG
5	L	39	ARG
5	L	42	LEU
5	L	57	LEU
5	L	79	TYR
5	L	86	LEU
5	L	87	LEU

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Mol	Chain	Res	Type
5	L	93	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	42	GLN
1	A	66	HIS
1	A	95	ASN
1	A	140	ASN
2	B	12	GLN
2	B	15	GLN
2	B	105	ASN
2	B	172	GLN
3	C	25	ASN
4	D	10	ASN
4	D	75	GLN
4	E	47	HIS
5	F	24	GLN
1	G	31	ASN
1	G	42	GLN
1	G	66	HIS
1	G	83	HIS
1	G	95	ASN
1	G	140	ASN
1	G	149	HIS
2	H	12	GLN
2	H	15	GLN
2	H	93	ASN
2	H	172	GLN
3	I	25	ASN
3	I	172	GLN
4	J	10	ASN
4	K	47	HIS
5	L	24	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.