



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 05:13 AM GMT

PDB ID : 1QX5
Title : Crystal structure of apoCalmodulin
Authors : Schumacher, M.A.; Crum, M.; Miller, M.C.
Deposited on : 2003-09-04
Resolution : 2.54 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

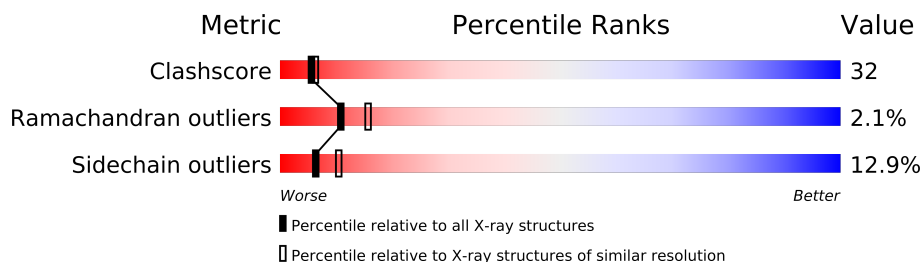
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.54 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	79885	4284 (2.58-2.50)
Ramachandran outliers	78287	4193 (2.58-2.50)
Sidechain outliers	78261	4195 (2.58-2.50)

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	B	148	
1	D	148	
1	I	148	
1	J	148	
1	K	148	
1	R	148	
1	T	148	
1	Y	148	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9333 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Calmodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	145	Total	C	N	O	S	0	0	0
			1146	702	184	251	9			
1	I	145	Total	C	N	O	S	0	0	0
			1146	702	184	251	9			
1	B	145	Total	C	N	O	S	0	0	0
			1146	702	184	251	9			
1	J	144	Total	C	N	O	S	0	0	0
			1138	698	183	248	9			
1	K	143	Total	C	N	O	S	0	0	0
			1129	693	181	246	9			
1	T	145	Total	C	N	O	S	0	0	0
			1146	702	184	251	9			
1	R	147	Total	C	N	O	S	0	0	0
			1161	711	187	254	9			
1	Y	145	Total	C	N	O	S	0	0	0
			1146	702	184	251	9			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	26	Total	O	0	0
			26	26		
2	D	24	Total	O	0	0
			24	24		
2	I	30	Total	O	0	0
			30	30		
2	J	15	Total	O	0	0
			15	15		
2	K	20	Total	O	0	0
			20	20		
2	R	30	Total	O	0	0
			30	30		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	T	7	Total 7	O 7	0	0
2	Y	23	Total 23	O 23	0	0

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.

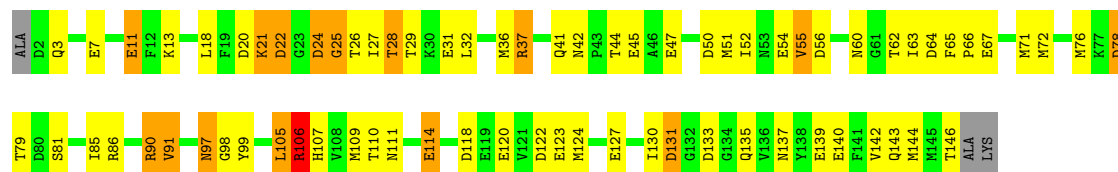
• Molecule 1: Calmodulin

Chain D: 



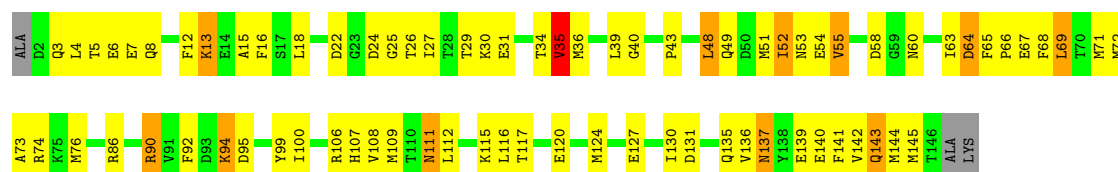
• Molecule 1: Calmodulin

Chain I: 



• Molecule 1: Calmodulin

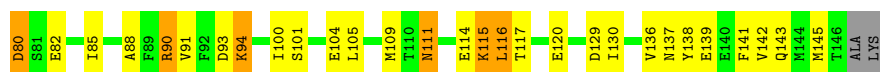
Chain B: 



• Molecule 1: Calmodulin

Chain J: 





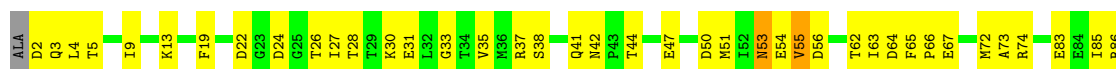
• Molecule 1: Calmodulin

Chain K:



• Molecule 1: Calmodulin

Chain T:



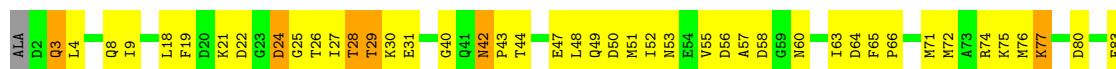
• Molecule 1: Calmodulin

Chain R:



• Molecule 1: Calmodulin

Chain Y:



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	146.00Å 146.00Å 78.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	66.39 – 2.54	Depositor
% Data completeness (in resolution range)	98.5 (66.39-2.54)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.227 , 0.277	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9333	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

5 Model quality i

5.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	B	0.83	2/1158 (0.2%)	0.83	0/1555
1	D	0.87	1/1158 (0.1%)	0.87	2/1555 (0.1%)
1	I	1.01	0/1158	1.06	8/1555 (0.5%)
1	J	0.68	0/1150	0.73	1/1544 (0.1%)
1	K	0.82	0/1141	0.82	1/1532 (0.1%)
1	R	0.93	1/1173 (0.1%)	0.92	1/1573 (0.1%)
1	T	0.59	0/1158	0.68	1/1555 (0.1%)
1	Y	0.77	0/1158	0.87	4/1555 (0.3%)
All	All	0.82	4/9254 (0.0%)	0.85	18/12424 (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	D	0	1
1	R	0	1
1	Y	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	83	GLU	CG-CD	6.67	1.61	1.51
1	R	83	GLU	CG-CD	5.80	1.60	1.51
1	B	35	VAL	CA-CB	5.67	1.66	1.54
1	B	15	ALA	CA-CB	5.19	1.63	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	106	ARG	NE-CZ-NH2	-7.73	116.44	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	37	ARG	NE-CZ-NH2	-7.70	116.45	120.30
1	Y	131	ASP	CB-CG-OD2	-7.38	111.66	118.30
1	I	131	ASP	CB-CG-OD1	7.36	124.93	118.30
1	R	129	ASP	CB-CG-OD1	-6.13	112.79	118.30
1	D	90	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	Y	129	ASP	CB-CG-OD2	-6.05	112.85	118.30
1	I	90	ARG	NE-CZ-NH1	-6.02	117.29	120.30
1	I	91	VAL	CB-CA-C	-5.97	100.05	111.40
1	D	90	ARG	NE-CZ-NH2	-5.93	117.34	120.30
1	Y	106	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	T	93	ASP	CB-CG-OD1	-5.85	113.03	118.30
1	Y	106	ARG	NE-CZ-NH1	5.83	123.22	120.30
1	J	93	ASP	CB-CG-OD2	5.72	123.45	118.30
1	I	37	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	I	131	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	K	37	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	I	130	ILE	CB-CA-C	-5.19	101.22	111.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	92	PHE	Sidechain
1	R	99	TYR	Sidechain
1	Y	138	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1146	0	1070	89	0
1	D	1146	0	1070	83	0
1	I	1146	0	1070	63	0
1	J	1138	0	1066	67	0
1	K	1129	0	1058	93	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	R	1161	0	1088	77	0
1	T	1146	0	1070	74	0
1	Y	1146	0	1070	66	0
2	B	26	0	0	1	0
2	D	24	0	0	2	0
2	I	30	0	0	2	0
2	J	15	0	0	2	0
2	K	20	0	0	2	0
2	R	30	0	0	1	0
2	T	7	0	0	2	0
2	Y	23	0	0	2	0
All	All	9333	0	8562	566	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 32.

All (566) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:94:LYS:HG3	1:K:94:LYS:HG2	1.39	1.04
1:K:4:LEU:HD21	1:K:73:ALA:HA	1.39	1.04
1:D:44:THR:HG21	1:I:131:ASP:OD2	1.59	1.03
1:R:124:MET:HA	1:R:144:MET:HE1	1.42	1.01
1:J:90:ARG:HG2	1:J:90:ARG:HH11	1.24	1.01
1:Y:55:VAL:HG11	1:Y:63:ILE:HG23	1.42	1.00
1:D:55:VAL:HG13	1:D:56:ASP:H	1.30	0.93
1:Y:97:ASN:ND2	1:Y:99:TYR:H	1.67	0.91
1:R:55:VAL:HG11	1:R:63:ILE:HG23	1.51	0.90
1:B:29:THR:HG23	1:B:52:ILE:HG13	1.55	0.88
1:I:28:THR:HG23	1:I:31:GLU:OE2	1.71	0.88
1:B:5:THR:HG23	1:B:8:GLN:H	1.37	0.88
1:I:20:ASP:HA	1:I:21:LYS:HZ1	1.38	0.86
1:B:137:ASN:HD21	1:B:139:GLU:HB2	1.38	0.85
1:K:97:ASN:ND2	1:K:99:TYR:H	1.75	0.84
1:R:5:THR:H	1:R:8:GLN:HG3	1.42	0.84
1:K:28:THR:OG1	1:K:31:GLU:HG3	1.78	0.84
1:K:97:ASN:ND2	2:K:164:HOH:O	2.11	0.84
1:K:54:GLU:HG2	1:K:74:ARG:HH22	1.42	0.84
1:B:94:LYS:HG3	1:K:94:LYS:CG	2.08	0.83
1:B:31:GLU:O	1:B:35:VAL:HG12	1.78	0.83
1:B:4:LEU:HD21	1:B:73:ALA:HA	1.58	0.83
1:R:26:THR:HA	1:R:63:ILE:O	1.80	0.82
1:T:55:VAL:HG11	1:T:63:ILE:HG13	1.59	0.82

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:124:MET:HA	1:K:144:MET:HE1	1.61	0.81
1:J:27:ILE:HB	1:J:31:GLU:HG3	1.62	0.81
1:D:55:VAL:HG21	1:D:67:GLU:CD	2.01	0.81
1:T:27:ILE:HG22	1:T:63:ILE:HG22	1.63	0.79
1:B:94:LYS:CG	1:K:94:LYS:HG2	2.13	0.79
1:I:97:ASN:ND2	2:I:171:HOH:O	2.17	0.78
1:I:20:ASP:HA	1:I:21:LYS:NZ	1.99	0.78
1:D:44:THR:CG2	1:I:131:ASP:OD2	2.32	0.77
1:I:55:VAL:HG12	1:I:56:ASP:N	1.99	0.77
1:J:44:THR:HB	1:J:47:GLU:HG3	1.67	0.77
1:I:91:VAL:O	1:R:94:LYS:HE2	1.85	0.76
1:B:117:THR:HG23	1:B:120:GLU:OE1	1.87	0.75
1:T:22:ASP:HB2	1:T:24:ASP:OD1	1.86	0.75
1:R:44:THR:HG22	1:R:47:GLU:HG3	1.66	0.75
1:K:55:VAL:HG22	1:K:67:GLU:HB3	1.66	0.75
1:B:109:MET:HE1	1:K:108:VAL:HG12	1.69	0.75
1:B:94:LYS:HE2	1:K:94:LYS:HG2	1.70	0.74
1:Y:119:GLU:O	1:Y:123:GLU:HG3	1.86	0.74
1:B:116:LEU:HD21	1:B:145:MET:SD	2.27	0.74
1:J:29:THR:HB	1:J:52:ILE:HD13	1.68	0.73
1:K:140:GLU:O	1:K:144:MET:HG3	1.88	0.73
1:Y:140:GLU:O	1:Y:144:MET:HG2	1.89	0.73
1:R:5:THR:OG1	1:R:7:GLU:HB3	1.89	0.72
1:K:29:THR:HG23	1:K:52:ILE:HD12	1.70	0.72
1:J:7:GLU:HG2	1:J:8:GLN:N	2.03	0.72
1:J:90:ARG:NE	1:T:138:TYR:OH	2.23	0.72
1:I:97:ASN:ND2	1:I:99:TYR:H	1.85	0.72
1:K:124:MET:CA	1:K:144:MET:HE1	2.18	0.72
1:R:31:GLU:O	1:R:35:VAL:HG23	1.87	0.72
1:K:25:GLY:O	1:K:64:ASP:HA	1.89	0.71
1:J:42:ASN:H	1:J:42:ASN:HD22	1.38	0.71
1:I:22:ASP:HB2	1:I:24:ASP:OD1	1.90	0.71
1:D:48:LEU:HD23	1:D:51:MET:HE3	1.72	0.71
1:T:44:THR:HG23	1:T:47:GLU:OE1	1.91	0.71
1:D:30:LYS:HE2	1:D:133:ASP:OD1	1.91	0.70
1:B:131:ASP:OD2	1:J:44:THR:HG21	1.91	0.70
1:J:54:GLU:OE2	1:J:71:MET:HE2	1.90	0.70
1:B:137:ASN:ND2	1:B:139:GLU:HB2	2.07	0.70
1:D:44:THR:HG22	1:D:47:GLU:H	1.55	0.70
1:K:42:ASN:HD22	1:K:42:ASN:H	1.40	0.70
1:R:44:THR:HG21	1:Y:131:ASP:OD1	1.92	0.70
1:D:24:ASP:O	1:D:26:THR:N	2.21	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:117:THR:OG1	1:J:120:GLU:HG3	1.91	0.70
1:I:55:VAL:HG11	1:I:63:ILE:HG12	1.73	0.70
1:B:130:ILE:HD12	1:B:136:VAL:CG2	2.22	0.69
1:I:24:ASP:CG	1:I:25:GLY:H	1.95	0.69
1:B:130:ILE:HD12	1:B:136:VAL:HG21	1.74	0.69
1:K:64:ASP:HB2	1:K:67:GLU:HG3	1.73	0.69
1:B:107:HIS:O	1:B:111:ASN:HB3	1.93	0.69
1:K:4:LEU:HA	1:K:8:GLN:OE1	1.93	0.69
1:R:29:THR:OG1	1:R:52:ILE:HG13	1.93	0.68
1:B:86:ARG:O	1:B:90:ARG:HG2	1.94	0.68
1:R:119:GLU:O	1:R:123:GLU:HG3	1.93	0.68
1:D:55:VAL:HG13	1:D:56:ASP:N	2.07	0.68
1:D:7:GLU:O	1:D:11:GLU:HG3	1.93	0.68
1:D:70:THR:O	1:D:74:ARG:HD3	1.94	0.67
1:D:63:ILE:HG13	1:D:67:GLU:HG3	1.75	0.67
1:R:44:THR:HG21	1:Y:131:ASP:CG	2.15	0.67
1:D:92:PHE:O	1:D:94:LYS:HD2	1.95	0.67
1:B:108:VAL:HG12	1:B:109:MET:CE	2.24	0.66
1:J:130:ILE:HG22	2:J:160:HOH:O	1.93	0.66
1:D:34:THR:HB	2:D:155:HOH:O	1.94	0.66
1:K:124:MET:CB	1:K:144:MET:HE1	2.25	0.66
2:J:156:HOH:O	1:T:37:ARG:HD3	1.94	0.66
1:T:97:ASN:H	1:T:97:ASN:HD22	1.44	0.66
1:D:127:GLU:HB3	1:D:144:MET:HE1	1.78	0.66
1:R:55:VAL:CG1	1:R:63:ILE:HG23	2.25	0.66
1:J:90:ARG:CG	1:J:90:ARG:HH11	2.01	0.65
1:T:26:THR:HA	1:T:63:ILE:O	1.97	0.65
1:R:77:LYS:NZ	1:R:77:LYS:HA	2.11	0.65
1:R:55:VAL:HG12	1:R:56:ASP:N	2.11	0.65
1:T:33:GLY:O	1:T:37:ARG:HG2	1.97	0.65
1:B:68:PHE:HA	1:B:71:MET:HE3	1.79	0.65
1:B:74:ARG:NH1	1:B:74:ARG:HB3	2.13	0.64
1:K:4:LEU:HD11	1:K:73:ALA:HB2	1.80	0.64
1:Y:131:ASP:HB2	1:Y:135:GLN:O	1.96	0.64
1:T:94:LYS:HD2	1:T:94:LYS:H	1.63	0.64
1:D:48:LEU:HD23	1:D:51:MET:CE	2.28	0.64
1:I:140:GLU:O	1:I:144:MET:HG3	1.98	0.63
1:D:44:THR:HG22	1:D:46:ALA:H	1.63	0.63
1:T:95:ASP:O	1:T:97:ASN:N	2.31	0.63
1:T:130:ILE:HG12	1:T:136:VAL:HG22	1.81	0.63
1:Y:120:GLU:O	1:Y:124:MET:HG3	1.97	0.63
1:J:55:VAL:HG22	1:J:67:GLU:HB3	1.80	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:91:VAL:O	1:I:91:VAL:HG12	1.97	0.63
1:R:130:ILE:N	1:R:130:ILE:HD12	2.14	0.63
1:I:86:ARG:O	1:I:90:ARG:HG3	1.98	0.63
1:D:56:ASP:HB2	1:D:62:THR:O	1.99	0.63
1:K:55:VAL:HG13	1:K:67:GLU:CD	2.19	0.62
1:B:5:THR:HG22	1:B:8:GLN:CD	2.18	0.62
1:R:18:LEU:HD11	1:R:125:ILE:HG21	1.81	0.62
1:I:51:MET:O	1:I:54:GLU:HB3	1.99	0.62
1:J:64:ASP:HB3	1:J:66:PRO:HD2	1.79	0.62
1:K:65:PHE:HB3	1:K:66:PRO:HD3	1.82	0.62
1:R:44:THR:HG23	1:R:46:ALA:H	1.65	0.62
1:D:76:MET:HE2	1:D:79:THR:HB	1.82	0.62
1:B:139:GLU:OE1	1:B:142:VAL:HG11	2.00	0.62
1:B:92:PHE:HB2	1:B:100:ILE:HG21	1.82	0.62
1:R:27:ILE:O	1:R:63:ILE:HD12	2.00	0.61
1:I:110:THR:O	1:I:114:GLU:HB3	2.00	0.61
1:B:22:ASP:HB2	1:B:24:ASP:OD1	2.00	0.61
1:Y:72:MET:O	1:Y:76:MET:HB2	2.00	0.61
1:D:21:LYS:HD3	1:D:22:ASP:H	1.64	0.61
1:Y:48:LEU:HD23	1:Y:51:MET:HE3	1.81	0.61
1:T:86:ARG:HD2	1:T:90:ARG:NH2	2.15	0.61
1:T:104:GLU:O	1:T:108:VAL:HG23	2.01	0.61
1:K:54:GLU:HG2	1:K:74:ARG:NH2	2.14	0.60
1:K:107:HIS:O	1:K:111:ASN:HB3	2.01	0.60
1:R:97:ASN:ND2	1:R:99:TYR:H	1.98	0.60
1:R:130:ILE:HG22	1:R:131:ASP:N	2.15	0.60
1:Y:64:ASP:HB3	1:Y:66:PRO:HD2	1.83	0.60
1:R:28:THR:HG23	1:R:31:GLU:OE1	2.02	0.60
1:D:120:GLU:O	1:D:124:MET:HG3	2.00	0.60
1:B:139:GLU:HA	1:B:142:VAL:HG12	1.83	0.60
1:T:28:THR:HG22	1:T:30:LYS:H	1.66	0.60
1:T:93:ASP:OD1	1:T:93:ASP:C	2.39	0.60
1:I:41:GLN:OE1	1:I:79:THR:HG21	2.02	0.60
1:I:97:ASN:HD22	1:I:98:GLY:N	1.99	0.60
1:R:5:THR:N	1:R:8:GLN:HG3	2.14	0.59
1:J:55:VAL:HG23	1:J:71:MET:HE3	1.83	0.59
1:Y:48:LEU:O	1:Y:52:ILE:HG13	2.02	0.59
1:J:94:LYS:HG3	1:T:94:LYS:HG3	1.85	0.59
1:Y:56:ASP:OD1	1:Y:60:ASN:HB2	2.01	0.59
1:R:77:LYS:HA	1:R:77:LYS:HZ3	1.68	0.59
1:B:109:MET:HE2	1:B:109:MET:HA	1.84	0.59
1:Y:87:GLU:OE2	1:Y:90:ARG:NH1	2.35	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:64:ASP:HB3	1:I:66:PRO:HD2	1.84	0.59
1:D:55:VAL:HG21	1:D:67:GLU:OE1	2.02	0.59
1:K:29:THR:HG23	1:K:52:ILE:CD1	2.32	0.59
1:K:74:ARG:HB2	1:K:75:LYS:HZ2	1.68	0.59
1:T:27:ILE:HG22	1:T:63:ILE:CG2	2.32	0.59
1:B:65:PHE:HB3	1:B:66:PRO:HD3	1.85	0.59
1:K:90:ARG:HH11	1:K:90:ARG:HG2	1.67	0.59
1:K:12:PHE:CZ	1:K:76:MET:HG3	2.38	0.58
1:R:77:LYS:HD2	1:R:77:LYS:N	2.19	0.58
1:T:64:ASP:HB3	1:T:66:PRO:HD2	1.86	0.58
1:T:28:THR:HB	1:T:31:GLU:HG3	1.86	0.58
1:T:107:HIS:O	1:T:111:ASN:HB3	2.04	0.58
1:Y:97:ASN:C	1:Y:97:ASN:HD22	2.07	0.58
1:I:55:VAL:HG12	1:I:56:ASP:H	1.65	0.58
1:B:115:LYS:HD2	1:K:85:ILE:HD11	1.86	0.58
1:B:5:THR:HG22	1:B:8:GLN:NE2	2.18	0.58
1:T:27:ILE:HG23	1:T:28:THR:N	2.19	0.58
1:D:21:LYS:HD2	1:D:31:GLU:OE1	2.03	0.58
1:K:117:THR:OG1	1:K:120:GLU:HG3	2.04	0.57
1:Y:97:ASN:HD22	1:Y:98:GLY:N	2.01	0.57
1:B:115:LYS:CD	1:K:85:ILE:HD11	2.35	0.57
1:B:43:PRO:HB2	1:B:48:LEU:HD13	1.84	0.57
1:R:124:MET:HA	1:R:144:MET:CE	2.28	0.57
1:Y:26:THR:HG22	1:Y:64:ASP:OD1	2.04	0.57
1:T:64:ASP:HB2	1:T:67:GLU:HG3	1.85	0.57
1:D:63:ILE:HD11	1:D:71:MET:CE	2.35	0.57
1:J:137:ASN:ND2	1:J:139:GLU:HB2	2.19	0.57
1:D:130:ILE:HB	1:D:136:VAL:HG22	1.87	0.57
1:I:146:THR:HG21	1:R:82:GLU:HG3	1.87	0.57
1:B:24:ASP:OD2	1:B:26:THR:HB	2.04	0.57
1:B:40:GLY:HA2	1:K:107:HIS:CD2	2.39	0.57
1:Y:28:THR:HG23	1:Y:31:GLU:OE1	2.03	0.57
1:D:48:LEU:HA	1:D:51:MET:HE2	1.86	0.57
1:I:32:LEU:HD22	1:I:52:ILE:CD1	2.35	0.56
1:D:44:THR:CG2	1:D:46:ALA:H	2.19	0.56
1:R:29:THR:HA	1:R:52:ILE:HD11	1.86	0.56
1:T:28:THR:HB	1:T:31:GLU:CG	2.35	0.56
1:R:44:THR:HG21	1:Y:131:ASP:OD2	2.04	0.56
1:I:85:ILE:HD11	1:R:115:LYS:HD2	1.88	0.56
1:K:124:MET:HG2	1:K:144:MET:CE	2.36	0.56
1:J:27:ILE:HG13	1:J:63:ILE:HG23	1.87	0.56
1:I:131:ASP:OD1	1:I:133:ASP:N	2.38	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:55:VAL:HG23	1:J:71:MET:CE	2.35	0.56
1:I:97:ASN:C	1:I:97:ASN:HD22	2.08	0.56
1:Y:43:PRO:HB2	1:Y:48:LEU:HG	1.88	0.55
1:J:111:ASN:ND2	1:J:115:LYS:HE3	2.21	0.55
1:D:87:GLU:OE1	2:D:164:HOH:O	2.18	0.55
1:K:4:LEU:HD23	1:K:4:LEU:N	2.21	0.55
1:Y:44:THR:OG1	1:Y:47:GLU:HG3	2.06	0.55
1:T:117:THR:O	1:T:120:GLU:N	2.39	0.55
1:J:28:THR:O	1:J:31:GLU:HG2	2.07	0.55
1:J:29:THR:O	1:J:52:ILE:HD11	2.07	0.55
1:Y:26:THR:HA	1:Y:63:ILE:O	2.07	0.55
1:D:56:ASP:OD1	1:D:60:ASN:HB3	2.06	0.55
1:K:55:VAL:HG21	1:K:63:ILE:HD12	1.88	0.55
1:B:13:LYS:HA	1:B:65:PHE:CE1	2.42	0.55
1:K:19:PHE:HD2	1:K:35:VAL:HG22	1.72	0.55
1:B:13:LYS:HA	1:B:65:PHE:HE1	1.72	0.54
1:J:137:ASN:HD21	1:J:139:GLU:HB2	1.71	0.54
1:T:119:GLU:O	1:T:123:GLU:HB2	2.07	0.54
1:Y:42:ASN:H	1:Y:42:ASN:HD22	1.55	0.54
1:D:32:LEU:HG	1:D:36:MET:CE	2.38	0.54
1:T:116:LEU:HD11	1:T:145:MET:SD	2.47	0.54
1:T:97:ASN:ND2	1:T:99:TYR:H	2.04	0.54
1:D:76:MET:HA	1:D:76:MET:CE	2.37	0.54
1:R:97:ASN:ND2	2:R:158:HOH:O	2.39	0.54
1:B:108:VAL:HG12	1:B:109:MET:HE3	1.89	0.54
1:Y:97:ASN:HD22	1:Y:99:TYR:H	1.52	0.54
1:I:97:ASN:C	1:I:97:ASN:ND2	2.60	0.54
1:J:64:ASP:CB	1:J:66:PRO:HD2	2.37	0.54
1:D:41:GLN:O	1:D:43:PRO:HD3	2.08	0.54
1:K:55:VAL:HG11	1:K:63:ILE:CB	2.38	0.54
1:D:63:ILE:HD11	1:D:71:MET:HE3	1.90	0.54
1:T:97:ASN:ND2	2:T:154:HOH:O	2.32	0.54
1:D:36:MET:HE3	1:D:48:LEU:HD21	1.90	0.53
1:Y:97:ASN:ND2	2:Y:163:HOH:O	2.40	0.53
1:J:55:VAL:HG12	1:J:56:ASP:N	2.22	0.53
1:T:115:LYS:O	1:T:116:LEU:HD12	2.08	0.53
1:I:106:ARG:C	1:I:106:ARG:HD2	2.28	0.53
1:Y:117:THR:OG1	1:Y:120:GLU:HG3	2.07	0.53
1:K:91:VAL:HG12	1:K:91:VAL:O	2.08	0.53
1:R:109:MET:CE	1:R:109:MET:HA	2.39	0.53
1:D:44:THR:HG22	1:D:46:ALA:N	2.23	0.53
1:I:105:LEU:HD22	1:I:109:MET:HG2	1.91	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:7:GLU:O	1:K:11:GLU:HG3	2.08	0.53
1:D:124:MET:HA	1:D:144:MET:CE	2.38	0.53
1:Y:56:ASP:C	1:Y:58:ASP:H	2.12	0.53
1:I:120:GLU:O	1:I:124:MET:HG3	2.09	0.53
1:J:55:VAL:HG13	1:J:67:GLU:CD	2.29	0.53
1:B:140:GLU:O	1:B:144:MET:HG3	2.08	0.53
1:J:138:TYR:CE2	1:T:98:GLY:HA2	2.44	0.53
1:T:97:ASN:H	1:T:97:ASN:ND2	2.07	0.53
1:R:109:MET:HA	1:R:109:MET:HE3	1.91	0.53
1:D:41:GLN:HG3	1:D:41:GLN:O	2.08	0.52
1:Y:77:LYS:HA	1:Y:77:LYS:NZ	2.23	0.52
1:I:26:THR:HA	1:I:63:ILE:O	2.09	0.52
1:B:117:THR:OG1	1:B:120:GLU:HB2	2.09	0.52
1:K:36:MET:O	1:K:39:LEU:HB2	2.09	0.52
1:D:122:ASP:O	1:D:126:ARG:HG3	2.10	0.52
1:D:48:LEU:HA	1:D:51:MET:CE	2.39	0.52
1:B:76:MET:HE2	1:B:76:MET:N	2.24	0.52
1:T:93:ASP:OD1	1:T:93:ASP:O	2.28	0.52
1:B:94:LYS:HE2	1:K:94:LYS:CG	2.39	0.52
1:D:44:THR:HG23	2:I:156:HOH:O	2.08	0.52
1:R:44:THR:HG23	1:R:46:ALA:N	2.24	0.52
1:Y:49:GLN:HA	1:Y:52:ILE:HD12	1.92	0.52
1:J:136:VAL:HG21	1:J:141:PHE:HE1	1.75	0.52
1:D:47:GLU:O	1:D:51:MET:HG3	2.09	0.52
1:R:44:THR:HG22	1:R:47:GLU:H	1.75	0.52
1:Y:47:GLU:O	1:Y:51:MET:HG3	2.10	0.52
1:Y:18:LEU:HD12	1:Y:19:PHE:CE2	2.44	0.52
1:Y:71:MET:O	1:Y:74:ARG:HG2	2.10	0.52
1:D:82:GLU:HB2	1:Y:146:THR:HG21	1.92	0.51
1:J:117:THR:HG23	1:J:120:GLU:OE1	2.10	0.51
1:Y:53:ASN:HA	1:Y:57:ALA:HB2	1.93	0.51
1:D:116:LEU:O	1:D:116:LEU:HD23	2.10	0.51
1:J:90:ARG:CG	1:J:90:ARG:NH1	2.67	0.51
1:K:55:VAL:C	1:K:57:ALA:H	2.13	0.51
1:J:65:PHE:N	1:J:66:PRO:CD	2.74	0.51
1:Y:94:LYS:HD2	1:Y:94:LYS:N	2.26	0.51
1:J:100:ILE:O	1:T:135:GLN:HA	2.11	0.51
1:K:74:ARG:HB2	1:K:75:LYS:NZ	2.25	0.51
1:R:48:LEU:HD23	1:R:51:MET:HE3	1.92	0.51
1:J:80:ASP:OD2	1:J:82:GLU:HB3	2.10	0.51
1:J:90:ARG:HG2	1:J:90:ARG:NH1	2.04	0.51
1:R:29:THR:HG23	1:R:52:ILE:HG13	1.91	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:T:19:PHE:HD2	1:T:35:VAL:HG22	1.76	0.51
1:K:55:VAL:HG11	1:K:63:ILE:CG1	2.40	0.51
1:K:16:PHE:HD1	1:K:68:PHE:CD1	2.29	0.51
1:I:24:ASP:CG	1:I:25:GLY:N	2.64	0.51
1:Y:18:LEU:HD11	1:Y:125:ILE:HG21	1.93	0.51
1:D:64:ASP:OD1	1:D:66:PRO:HD2	2.11	0.50
1:B:95:ASP:N	1:B:95:ASP:OD2	2.44	0.50
1:K:49:GLN:C	1:K:51:MET:H	2.12	0.50
1:K:28:THR:HG23	1:K:31:GLU:OE1	2.10	0.50
1:Y:44:THR:HG23	1:Y:47:GLU:OE2	2.11	0.50
1:B:5:THR:HG22	1:B:8:GLN:CG	2.42	0.50
1:I:97:ASN:HD21	1:I:99:TYR:HB2	1.76	0.50
1:K:56:ASP:C	1:K:58:ASP:H	2.14	0.50
1:T:51:MET:O	1:T:55:VAL:HG23	2.11	0.50
1:Y:42:ASN:ND2	1:Y:42:ASN:O	2.44	0.50
1:K:146:THR:CG2	1:K:146:THR:O	2.59	0.50
1:K:117:THR:O	1:K:118:ASP:C	2.50	0.50
1:D:56:ASP:CG	1:D:61:GLY:H	2.15	0.50
1:I:55:VAL:HG13	1:I:67:GLU:CD	2.32	0.50
1:T:97:ASN:N	1:T:97:ASN:HD22	2.06	0.50
1:T:121:VAL:O	1:T:125:ILE:HG13	2.12	0.50
1:R:58:ASP:OD1	1:R:60:ASN:HB2	2.10	0.50
1:D:24:ASP:OD2	1:D:26:THR:HG23	2.10	0.50
1:I:105:LEU:CD2	1:I:109:MET:HG2	2.42	0.50
1:D:112:LEU:O	1:D:116:LEU:HD22	2.11	0.50
1:R:97:ASN:HD22	1:R:97:ASN:H	1.60	0.50
1:Y:97:ASN:C	1:Y:97:ASN:ND2	2.65	0.50
1:D:32:LEU:HG	1:D:36:MET:HE1	1.94	0.49
1:R:29:THR:CG2	1:R:52:ILE:HG13	2.42	0.49
1:B:36:MET:HE1	1:B:48:LEU:HD12	1.93	0.49
1:B:135:GLN:HG2	2:K:159:HOH:O	2.12	0.49
1:B:108:VAL:HG12	1:B:109:MET:HE2	1.92	0.49
1:R:25:GLY:O	1:R:64:ASP:HA	2.12	0.49
1:Y:116:LEU:HD21	1:Y:145:MET:SD	2.52	0.49
1:T:5:THR:O	1:T:9:ILE:HG13	2.12	0.49
1:J:116:LEU:HD13	1:J:120:GLU:HB2	1.94	0.49
1:J:130:ILE:O	1:J:130:ILE:HG23	2.12	0.49
1:R:130:ILE:HD12	1:R:130:ILE:H	1.75	0.49
1:Y:18:LEU:HD12	1:Y:19:PHE:CZ	2.47	0.49
1:B:135:GLN:HG3	1:K:99:TYR:HB3	1.94	0.49
1:K:55:VAL:HG13	1:K:67:GLU:OE1	2.13	0.49
1:R:44:THR:HG22	1:R:47:GLU:CG	2.38	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:T:125:ILE:HG23	1:T:130:ILE:HD11	1.93	0.49
1:J:142:VAL:HG13	1:J:143:GLN:N	2.27	0.49
1:B:5:THR:CG2	1:B:8:GLN:H	2.19	0.49
1:K:12:PHE:HZ	1:K:76:MET:HG3	1.78	0.49
1:R:48:LEU:HA	1:R:51:MET:HE2	1.94	0.49
1:R:107:HIS:ND1	1:R:111:ASN:ND2	2.41	0.49
1:K:55:VAL:HG11	1:K:63:ILE:HB	1.94	0.49
1:B:141:PHE:O	1:B:145:MET:HG3	2.13	0.49
1:R:130:ILE:HG13	1:R:136:VAL:CG2	2.43	0.49
1:Y:107:HIS:O	1:Y:111:ASN:HB2	2.12	0.49
1:B:65:PHE:CZ	1:B:69:LEU:HG	2.48	0.49
1:T:117:THR:H	1:T:120:GLU:HB2	1.78	0.49
1:J:139:GLU:O	1:J:142:VAL:HG12	2.13	0.48
1:K:91:VAL:O	1:K:91:VAL:CG1	2.61	0.48
1:B:99:TYR:HB3	1:K:135:GLN:HG3	1.95	0.48
1:B:63:ILE:HD12	1:B:71:MET:HE1	1.94	0.48
1:Y:125:ILE:HG23	1:Y:130:ILE:HD11	1.94	0.48
1:K:48:LEU:O	1:K:52:ILE:HG12	2.13	0.48
1:I:72:MET:HA	1:I:72:MET:HE2	1.95	0.48
1:B:3:GLN:HE21	1:B:4:LEU:H	1.62	0.48
1:T:115:LYS:C	1:T:116:LEU:HD12	2.34	0.48
1:Y:53:ASN:CA	1:Y:57:ALA:HB2	2.43	0.48
1:R:64:ASP:HB3	1:R:66:PRO:HD2	1.94	0.48
1:K:28:THR:HG1	1:K:31:GLU:HG3	1.77	0.48
1:J:114:GLU:C	1:J:116:LEU:H	2.16	0.48
1:I:123:GLU:O	1:I:127:GLU:HG2	2.14	0.48
1:B:145:MET:CE	1:K:85:ILE:HD12	2.44	0.47
1:B:40:GLY:HA2	1:K:107:HIS:NE2	2.29	0.47
1:Y:28:THR:HG23	1:Y:31:GLU:CD	2.34	0.47
1:R:65:PHE:HB3	1:R:66:PRO:HD3	1.96	0.47
1:D:64:ASP:O	1:D:67:GLU:HG2	2.15	0.47
1:D:22:ASP:HB3	1:D:24:ASP:OD1	2.14	0.47
1:R:130:ILE:CG2	1:R:134:GLY:HA2	2.45	0.47
1:D:86:ARG:HD3	1:D:90:ARG:NH1	2.29	0.47
1:R:74:ARG:HG3	1:R:75:LYS:N	2.29	0.47
1:R:3:GLN:HE21	1:R:3:GLN:HB2	1.51	0.47
1:R:55:VAL:HG23	1:R:71:MET:HE1	1.97	0.47
1:T:31:GLU:O	1:T:35:VAL:HG23	2.15	0.47
1:B:111:ASN:O	1:B:115:LYS:HG3	2.15	0.47
1:T:116:LEU:HD21	1:T:124:MET:CE	2.45	0.47
1:R:126:ARG:HG2	1:R:126:ARG:HH11	1.80	0.47
1:T:72:MET:HA	1:T:72:MET:HE2	1.95	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:137:ASN:ND2	1:B:139:GLU:H	2.12	0.47
1:I:91:VAL:O	1:I:91:VAL:CG1	2.63	0.47
1:D:130:ILE:HD11	1:D:134:GLY:HA2	1.96	0.47
1:I:131:ASP:OD2	1:I:135:GLN:HB2	2.15	0.46
1:J:116:LEU:CD1	1:J:120:GLU:HB2	2.45	0.46
1:I:44:THR:OG1	1:I:47:GLU:HG3	2.14	0.46
1:I:142:VAL:O	1:I:146:THR:HG23	2.15	0.46
1:D:137:ASN:OD1	1:D:139:GLU:HB2	2.14	0.46
1:T:130:ILE:HG22	1:T:131:ASP:N	2.31	0.46
1:K:117:THR:HG23	1:K:120:GLU:OE2	2.16	0.46
1:K:52:ILE:CG2	1:K:56:ASP:HB3	2.45	0.46
1:D:102:ALA:HB2	1:Y:136:VAL:CG2	2.45	0.46
1:Y:22:ASP:HB2	1:Y:24:ASP:OD2	2.15	0.46
1:B:74:ARG:CB	1:B:74:ARG:HH11	2.28	0.46
1:Y:117:THR:HG23	1:Y:120:GLU:OE2	2.15	0.46
1:R:18:LEU:HD11	1:R:125:ILE:CG2	2.44	0.46
1:K:97:ASN:OD1	1:K:99:TYR:HB2	2.16	0.46
1:B:115:LYS:HD2	1:K:85:ILE:CD1	2.46	0.46
1:D:8:GLN:NE2	1:D:76:MET:HG2	2.30	0.46
1:J:88:ALA:HB1	1:T:112:LEU:HD22	1.96	0.46
1:J:22:ASP:OD1	1:J:22:ASP:N	2.49	0.46
1:I:107:HIS:O	1:I:111:ASN:HB3	2.14	0.46
1:J:29:THR:HG23	1:J:61:GLY:C	2.35	0.46
1:Y:8:GLN:O	1:Y:9:ILE:C	2.52	0.46
1:Y:27:ILE:HG13	1:Y:27:ILE:O	2.16	0.46
1:J:42:ASN:O	1:J:42:ASN:ND2	2.49	0.46
1:R:14:GLU:O	1:R:18:LEU:HB2	2.16	0.46
1:R:86:ARG:HD2	1:R:90:ARG:NH2	2.30	0.46
1:Y:129:ASP:N	1:Y:129:ASP:OD2	2.46	0.46
1:D:117:THR:OG1	1:D:120:GLU:HG3	2.16	0.46
1:R:130:ILE:CG2	1:R:131:ASP:N	2.79	0.46
1:D:130:ILE:HD13	1:D:130:ILE:O	2.16	0.46
1:J:91:VAL:O	1:J:91:VAL:HG12	2.16	0.46
1:D:13:LYS:HA	1:D:65:PHE:CE1	2.51	0.46
1:K:52:ILE:HG22	1:K:52:ILE:O	2.16	0.46
1:D:116:LEU:CD2	1:D:116:LEU:N	2.78	0.46
1:R:121:VAL:O	1:R:125:ILE:HG13	2.16	0.45
1:R:4:LEU:HD11	1:R:73:ALA:HA	1.99	0.45
1:R:137:ASN:OD1	1:R:139:GLU:HB2	2.16	0.45
1:D:64:ASP:OD1	1:D:67:GLU:HG2	2.17	0.45
1:K:21:LYS:HG3	1:K:22:ASP:H	1.81	0.45
1:K:124:MET:CG	1:K:144:MET:CE	2.93	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:T:54:GLU:HG2	1:T:74:ARG:HH22	1.82	0.45
1:J:101:SER:HB3	1:J:104:GLU:HB2	1.99	0.45
1:Y:92:PHE:O	1:Y:94:LYS:HD2	2.16	0.45
1:B:109:MET:HE1	1:K:108:VAL:CG1	2.42	0.45
1:B:12:PHE:CE2	1:B:72:MET:HB3	2.52	0.45
1:R:27:ILE:HG13	1:R:63:ILE:HD13	1.99	0.45
1:I:27:ILE:HB	1:I:31:GLU:HB2	1.99	0.45
1:K:64:ASP:HB3	1:K:66:PRO:HD2	1.98	0.45
1:J:105:LEU:HD12	1:J:109:MET:HG2	1.99	0.45
1:I:142:VAL:HG11	1:R:86:ARG:NH1	2.32	0.45
1:Y:42:ASN:N	1:Y:42:ASN:HD22	2.11	0.45
1:K:88:ALA:O	1:K:91:VAL:HB	2.17	0.45
1:B:64:ASP:OD1	1:B:67:GLU:HG3	2.17	0.45
1:D:2:ASP:HB3	1:D:77:LYS:NZ	2.32	0.45
1:I:55:VAL:HG21	1:I:63:ILE:CD1	2.46	0.45
1:D:124:MET:HA	1:D:144:MET:HE1	1.97	0.45
1:I:37:ARG:HA	1:I:41:GLN:O	2.17	0.45
1:J:137:ASN:HA	1:T:98:GLY:O	2.17	0.45
1:R:81:SER:O	1:R:85:ILE:HG13	2.18	0.44
1:T:28:THR:H	1:T:31:GLU:HB2	1.82	0.44
1:J:54:GLU:CD	1:J:71:MET:HE2	2.38	0.44
1:I:29:THR:OG1	1:I:52:ILE:HG12	2.16	0.44
1:I:11:GLU:HA	1:I:11:GLU:OE2	2.17	0.44
1:K:55:VAL:HG11	1:K:63:ILE:HG13	1.98	0.44
1:T:93:ASP:OD2	1:T:97:ASN:ND2	2.50	0.44
1:B:58:ASP:C	1:B:60:ASN:H	2.20	0.44
1:J:31:GLU:H	1:J:31:GLU:HG2	1.50	0.44
1:D:13:LYS:HD2	1:D:65:PHE:CE1	2.53	0.44
1:B:30:LYS:HD3	1:B:30:LYS:HA	1.84	0.44
1:J:85:ILE:HG22	1:T:142:VAL:HG22	2.00	0.44
1:K:42:ASN:HD22	1:K:42:ASN:N	2.10	0.44
1:K:120:GLU:O	1:K:124:MET:HG3	2.18	0.44
1:D:18:LEU:HA	1:D:18:LEU:HD23	1.83	0.44
1:R:56:ASP:OD1	1:R:60:ASN:HB3	2.17	0.44
1:J:90:ARG:CD	1:T:138:TYR:OH	2.66	0.44
1:T:19:PHE:HE2	1:T:38:SER:CB	2.30	0.44
1:D:24:ASP:HB2	1:D:26:THR:HG22	1.99	0.44
1:I:118:ASP:C	1:I:118:ASP:OD1	2.56	0.44
1:K:5:THR:O	1:K:8:GLN:N	2.51	0.44
1:R:72:MET:HE1	1:R:75:LYS:HD2	2.00	0.44
1:J:16:PHE:C	1:J:16:PHE:CD2	2.92	0.44
1:D:52:ILE:N	1:D:52:ILE:HD12	2.32	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:107:HIS:HE1	1:K:84:GLU:O	2.00	0.43
1:B:63:ILE:CD1	1:B:71:MET:HE1	2.48	0.43
1:B:25:GLY:O	1:B:64:ASP:HA	2.17	0.43
1:I:137:ASN:OD1	1:I:139:GLU:HB2	2.18	0.43
1:J:47:GLU:O	1:J:51:MET:HG3	2.17	0.43
1:J:141:PHE:O	1:J:145:MET:HG2	2.18	0.43
1:D:33:GLY:O	1:D:37:ARG:HG3	2.17	0.43
1:B:63:ILE:HD12	1:B:71:MET:CE	2.48	0.43
1:T:117:THR:OG1	1:T:120:GLU:HG3	2.18	0.43
1:B:58:ASP:C	1:B:60:ASN:N	2.71	0.43
1:T:47:GLU:O	1:T:51:MET:HG3	2.18	0.43
1:Y:4:LEU:HA	1:Y:8:GLN:OE1	2.18	0.43
1:K:41:GLN:O	1:K:43:PRO:HD3	2.18	0.43
1:D:25:GLY:O	1:D:64:ASP:HA	2.19	0.43
1:J:58:ASP:OD2	1:J:60:ASN:HB2	2.18	0.43
1:I:55:VAL:HG23	1:I:71:MET:CE	2.48	0.43
1:K:27:ILE:HG13	1:K:63:ILE:HG23	1.99	0.43
1:D:107:HIS:CE1	1:Y:40:GLY:CA	3.02	0.43
1:K:5:THR:O	1:K:6:GLU:C	2.57	0.43
1:K:26:THR:HA	1:K:63:ILE:O	2.18	0.43
1:I:135:GLN:HE21	1:R:99:TYR:HB3	1.84	0.43
1:D:64:ASP:CG	1:D:66:PRO:HD2	2.39	0.43
1:T:87:GLU:OE2	1:T:90:ARG:NH1	2.52	0.43
1:J:139:GLU:O	1:J:143:GLN:HG2	2.19	0.43
1:R:48:LEU:O	1:R:51:MET:HB2	2.18	0.43
1:D:55:VAL:CG1	1:D:56:ASP:H	2.14	0.43
1:K:55:VAL:HG21	1:K:63:ILE:CD1	2.48	0.43
1:T:111:ASN:C	1:T:111:ASN:HD22	2.22	0.43
1:Y:53:ASN:O	1:Y:57:ALA:HB2	2.19	0.43
1:D:116:LEU:C	1:D:116:LEU:HD23	2.37	0.43
1:I:76:MET:C	1:I:78:ASP:H	2.22	0.43
1:R:26:THR:HG22	1:R:63:ILE:O	2.19	0.42
1:I:97:ASN:HD22	1:I:99:TYR:H	1.63	0.42
1:B:69:LEU:HA	1:B:69:LEU:HD23	1.83	0.42
1:R:50:ASP:O	1:R:54:GLU:HB2	2.18	0.42
1:K:69:LEU:HA	1:K:69:LEU:HD12	1.83	0.42
1:Y:3:GLN:CD	1:Y:3:GLN:N	2.72	0.42
1:J:19:PHE:HD1	1:J:35:VAL:HG22	1.84	0.42
1:D:55:VAL:O	1:D:57:ALA:N	2.50	0.42
1:B:111:ASN:HD22	1:B:111:ASN:C	2.22	0.42
1:D:30:LYS:HE2	1:D:133:ASP:CG	2.39	0.42
1:K:49:GLN:C	1:K:51:MET:N	2.72	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:T:4:LEU:HD21	1:T:73:ALA:HA	2.02	0.42
1:D:121:VAL:O	1:D:125:ILE:HG13	2.19	0.42
1:J:27:ILE:CG1	1:J:63:ILE:HG23	2.48	0.42
1:I:55:VAL:CG2	1:I:71:MET:HE3	2.49	0.42
1:R:44:THR:CG2	1:R:47:GLU:H	2.32	0.42
1:B:55:VAL:HG12	1:B:67:GLU:HG2	2.01	0.42
1:B:115:LYS:HD3	1:K:85:ILE:HD11	2.02	0.42
1:D:117:THR:N	1:D:120:GLU:OE2	2.53	0.42
1:T:116:LEU:HD23	1:T:121:VAL:HG22	2.00	0.42
1:R:130:ILE:N	1:R:130:ILE:CD1	2.82	0.42
1:T:65:PHE:N	1:T:66:PRO:CD	2.83	0.42
1:R:51:MET:O	1:R:54:GLU:HB3	2.20	0.42
1:R:55:VAL:HG13	1:R:67:GLU:CD	2.40	0.42
1:B:131:ASP:OD2	1:J:44:THR:CG2	2.66	0.42
1:B:74:ARG:NH1	1:B:74:ARG:CB	2.82	0.42
1:R:141:PHE:O	1:R:145:MET:HG3	2.19	0.42
1:D:36:MET:HE3	1:D:48:LEU:CD2	2.49	0.42
1:D:63:ILE:HD11	1:D:71:MET:HE1	2.02	0.42
1:I:56:ASP:HB2	1:I:62:THR:O	2.19	0.42
1:R:29:THR:OG1	1:R:52:ILE:CG1	2.65	0.42
1:B:92:PHE:HB2	1:B:100:ILE:CG2	2.49	0.42
1:B:51:MET:HA	1:B:54:GLU:HB3	2.01	0.42
1:K:124:MET:HB3	1:K:144:MET:HE1	2.01	0.42
1:K:55:VAL:HG12	1:K:56:ASP:N	2.35	0.42
1:Y:80:ASP:OD1	1:Y:83:GLU:HG2	2.19	0.42
1:T:37:ARG:HA	1:T:41:GLN:O	2.20	0.42
1:Y:29:THR:OG1	1:Y:52:ILE:HD13	2.20	0.42
1:T:56:ASP:OD1	1:T:62:THR:O	2.37	0.42
1:K:92:PHE:CE2	1:K:105:LEU:HA	2.54	0.42
1:D:68:PHE:O	1:D:72:MET:HG2	2.19	0.42
1:B:90:ARG:HB2	2:B:156:HOH:O	2.20	0.41
1:R:21:LYS:HG3	1:R:22:ASP:N	2.35	0.41
1:T:86:ARG:HD2	1:T:90:ARG:HH22	1.84	0.41
1:J:45:GLU:HA	1:J:45:GLU:OE1	2.20	0.41
1:I:131:ASP:OD1	1:I:131:ASP:C	2.59	0.41
1:Y:65:PHE:N	1:Y:66:PRO:CD	2.83	0.41
1:I:55:VAL:CG1	1:I:63:ILE:HG12	2.48	0.41
1:T:94:LYS:HD3	2:T:155:HOH:O	2.19	0.41
1:J:109:MET:HE1	1:T:112:LEU:HB3	2.02	0.41
1:J:111:ASN:O	1:J:114:GLU:HB3	2.20	0.41
1:B:130:ILE:HG23	1:B:136:VAL:CG2	2.50	0.41
1:D:130:ILE:HD13	1:D:131:ASP:O	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:36:MET:SD	1:K:72:MET:HE3	2.61	0.41
1:B:34:THR:HG23	1:K:103:ALA:HB2	2.02	0.41
1:D:108:VAL:O	1:D:108:VAL:HG23	2.20	0.41
1:B:142:VAL:HG13	1:B:143:GLN:N	2.35	0.41
1:K:90:ARG:HG2	1:K:90:ARG:NH1	2.34	0.41
1:B:12:PHE:HZ	1:B:76:MET:SD	2.43	0.41
1:Y:48:LEU:HA	1:Y:51:MET:HE2	2.02	0.41
1:T:116:LEU:HD21	1:T:124:MET:HE1	2.02	0.41
1:B:130:ILE:CD1	1:B:136:VAL:CG2	2.97	0.41
1:J:142:VAL:CG1	1:J:143:GLN:N	2.84	0.41
1:I:32:LEU:HD22	1:I:52:ILE:HD11	2.02	0.41
1:J:90:ARG:HD2	1:T:138:TYR:CZ	2.56	0.41
1:B:107:HIS:CE1	1:K:84:GLU:O	2.74	0.41
1:I:65:PHE:O	1:I:66:PRO:C	2.58	0.41
1:K:12:PHE:CE2	1:K:76:MET:HG3	2.56	0.41
1:Y:118:ASP:HB2	2:Y:153:HOH:O	2.21	0.41
1:B:116:LEU:HD11	1:B:124:MET:SD	2.61	0.41
1:B:74:ARG:HB3	1:B:74:ARG:HH11	1.82	0.41
1:T:94:LYS:HD2	1:T:94:LYS:N	2.35	0.41
1:Y:24:ASP:CG	1:Y:25:GLY:H	2.25	0.41
1:B:16:PHE:CE1	1:B:27:ILE:HD13	2.56	0.41
1:J:38:SER:HB2	1:T:106:ARG:HD3	2.02	0.41
1:D:55:VAL:HG22	1:D:56:ASP:N	2.35	0.40
1:K:32:LEU:HD12	1:K:52:ILE:HD11	2.04	0.40
1:D:24:ASP:OD1	1:D:24:ASP:N	2.53	0.40
1:Y:56:ASP:C	1:Y:58:ASP:N	2.73	0.40
1:B:49:GLN:HA	1:B:52:ILE:HB	2.02	0.40
1:R:5:THR:HG1	1:R:7:GLU:HB3	1.83	0.40
1:T:19:PHE:CE2	1:T:38:SER:HB3	2.56	0.40
1:I:97:ASN:HD21	1:I:99:TYR:H	1.66	0.40
1:T:108:VAL:O	1:T:112:LEU:HB3	2.20	0.40
1:Y:77:LYS:HA	1:Y:77:LYS:HZ3	1.86	0.40
1:B:130:ILE:HD13	1:B:130:ILE:N	2.36	0.40
1:T:53:ASN:HA	1:T:53:ASN:HD22	1.60	0.40
1:Y:131:ASP:HB3	1:Y:133:ASP:H	1.85	0.40
1:J:115:LYS:HD2	1:T:85:ILE:HD11	2.04	0.40
1:I:47:GLU:O	1:I:51:MET:HG3	2.22	0.40
1:I:32:LEU:O	1:I:36:MET:HG2	2.22	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	B	143/148 (97%)	131 (92%)	11 (8%)	1 (1%)	30	49
1	D	143/148 (97%)	131 (92%)	9 (6%)	3 (2%)	11	16
1	I	143/148 (97%)	130 (91%)	10 (7%)	3 (2%)	11	16
1	J	142/148 (96%)	125 (88%)	14 (10%)	3 (2%)	11	16
1	K	141/148 (95%)	123 (87%)	12 (8%)	6 (4%)	4	4
1	R	145/148 (98%)	135 (93%)	5 (3%)	5 (3%)	6	7
1	T	143/148 (97%)	128 (90%)	13 (9%)	2 (1%)	16	27
1	Y	143/148 (97%)	129 (90%)	13 (9%)	1 (1%)	30	49
All	All	1143/1184 (96%)	1032 (90%)	87 (8%)	24 (2%)	11	16

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	22	ASP
1	I	55	VAL
1	T	96	GLY
1	R	55	VAL
1	R	147	ALA
1	D	25	GLY
1	K	75	LYS
1	Y	24	ASP
1	D	56	ASP
1	I	25	GLY
1	J	115	LYS
1	K	21	LYS
1	K	22	ASP
1	K	59	GLY
1	K	79	THR
1	R	24	ASP
1	J	56	ASP
1	K	56	ASP
1	I	24	ASP

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Mol	Chain	Res	Type
1	J	55	VAL
1	T	55	VAL
1	R	22	ASP
1	R	25	GLY
1	B	55	VAL

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	B	125/126 (99%)	106 (85%)	19 (15%)	4	7
1	D	125/126 (99%)	105 (84%)	20 (16%)	3	6
1	I	125/126 (99%)	105 (84%)	20 (16%)	3	6
1	J	124/126 (98%)	106 (86%)	18 (14%)	5	8
1	K	123/126 (98%)	113 (92%)	10 (8%)	17	29
1	R	126/126 (100%)	112 (89%)	14 (11%)	9	15
1	T	125/126 (99%)	115 (92%)	10 (8%)	17	30
1	Y	125/126 (99%)	107 (86%)	18 (14%)	5	8
All	All	998/1008 (99%)	869 (87%)	129 (13%)	6	10

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	6	GLU
1	D	18	LEU
1	D	21	LYS
1	D	24	ASP
1	D	26	THR
1	D	34	THR
1	D	49	GLN
1	D	58	ASP
1	D	60	ASN
1	D	69	LEU
1	D	75	LYS
1	D	76	MET

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Mol	Chain	Res	Type
1	D	80	ASP
1	D	90	ARG
1	D	94	LYS
1	D	100	ILE
1	D	114	GLU
1	D	116	LEU
1	D	130	ILE
1	D	143	GLN
1	I	3	GLN
1	I	7	GLU
1	I	11	GLU
1	I	13	LYS
1	I	18	LEU
1	I	21	LYS
1	I	22	ASP
1	I	28	THR
1	I	42	ASN
1	I	45	GLU
1	I	50	ASP
1	I	60	ASN
1	I	78	ASP
1	I	81	SER
1	I	97	ASN
1	I	105	LEU
1	I	106	ARG
1	I	114	GLU
1	I	122	ASP
1	I	143	GLN
1	B	6	GLU
1	B	7	GLU
1	B	13	LYS
1	B	18	LEU
1	B	35	VAL
1	B	39	LEU
1	B	48	LEU
1	B	52	ILE
1	B	53	ASN
1	B	64	ASP
1	B	69	LEU
1	B	90	ARG
1	B	94	LYS
1	B	106	ARG

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Mol	Chain	Res	Type
1	B	111	ASN
1	B	112	LEU
1	B	127	GLU
1	B	137	ASN
1	B	143	GLN
1	J	4	LEU
1	J	7	GLU
1	J	13	LYS
1	J	18	LEU
1	J	31	GLU
1	J	41	GLN
1	J	42	ASN
1	J	50	ASP
1	J	60	ASN
1	J	69	LEU
1	J	70	THR
1	J	77	LYS
1	J	80	ASP
1	J	90	ARG
1	J	94	LYS
1	J	111	ASN
1	J	116	LEU
1	J	129	ASP
1	K	4	LEU
1	K	22	ASP
1	K	24	ASP
1	K	32	LEU
1	K	41	GLN
1	K	42	ASN
1	K	60	ASN
1	K	97	ASN
1	K	106	ARG
1	K	135	GLN
1	T	2	ASP
1	T	3	GLN
1	T	13	LYS
1	T	42	ASN
1	T	50	ASP
1	T	53	ASN
1	T	83	GLU
1	T	97	ASN
1	T	111	ASN

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Mol	Chain	Res	Type
1	T	122	ASP
1	R	3	GLN
1	R	7	GLU
1	R	8	GLN
1	R	18	LEU
1	R	22	ASP
1	R	26	THR
1	R	28	THR
1	R	60	ASN
1	R	69	LEU
1	R	77	LYS
1	R	97	ASN
1	R	109	MET
1	R	114	GLU
1	R	116	LEU
1	Y	3	GLN
1	Y	21	LYS
1	Y	28	THR
1	Y	29	THR
1	Y	30	LYS
1	Y	42	ASN
1	Y	50	ASP
1	Y	75	LYS
1	Y	77	LYS
1	Y	84	GLU
1	Y	94	LYS
1	Y	97	ASN
1	Y	101	SER
1	Y	105	LEU
1	Y	114	GLU
1	Y	118	ASP
1	Y	135	GLN
1	Y	144	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	3	GLN
1	D	60	ASN
1	D	111	ASN
1	I	3	GLN
1	I	8	GLN

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Mol	Chain	Res	Type
1	I	53	ASN
1	I	60	ASN
1	I	97	ASN
1	I	135	GLN
1	I	143	GLN
1	B	3	GLN
1	B	41	GLN
1	B	53	ASN
1	B	107	HIS
1	B	111	ASN
1	B	137	ASN
1	J	41	GLN
1	J	42	ASN
1	J	60	ASN
1	J	107	HIS
1	J	111	ASN
1	J	137	ASN
1	K	42	ASN
1	K	49	GLN
1	K	53	ASN
1	K	60	ASN
1	K	97	ASN
1	K	135	GLN
1	T	3	GLN
1	T	42	ASN
1	T	53	ASN
1	T	60	ASN
1	T	97	ASN
1	T	111	ASN
1	T	135	GLN
1	R	3	GLN
1	R	49	GLN
1	R	53	ASN
1	R	60	ASN
1	R	97	ASN
1	R	111	ASN
1	R	143	GLN
1	Y	3	GLN
1	Y	42	ASN
1	Y	53	ASN
1	Y	97	ASN
1	Y	111	ASN

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Mol	Chain	Res	Type
1	Y	135	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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.