



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:11 PM GMT

PDB ID : 1P8S
Title : Structural and Functional Importance of First-Shell Metal Ligands in the Binuclear Manganese Cluster of Arginase I.
Authors : Cama, E.; Emig, F.A.; Ash, D.E.; Christianson, D.W.
Deposited on : 2003-05-07
Resolution : 3.20 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>

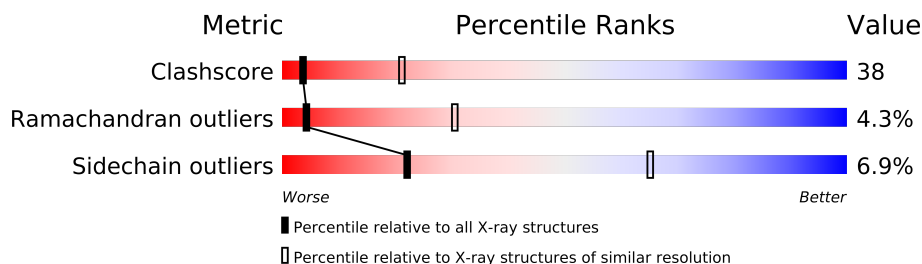
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.20 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	1078 (3.26-3.14)
Ramachandran outliers	78287	1059 (3.26-3.14)
Sidechain outliers	78261	1058 (3.26-3.14)

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	A	314	
1	B	314	
1	C	314	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7221 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 Arginase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2394	1527	405	454	8			
1	B	314	Total	C	N	O	S	0	0	0
			2394	1527	405	454	8			
1	C	314	Total	C	N	O	S	0	0	0
			2394	1527	405	454	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	232	CYS	ASP	ENGINEERED	UNP P07824
B	232	CYS	ASP	ENGINEERED	UNP P07824
C	232	CYS	ASP	ENGINEERED	UNP P07824

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mn	0	0
			2	2		
2	A	2	Total	Mn	0	0
			2	2		
2	C	2	Total	Mn	0	0
			2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	12	Total	O	0	0
			12	12		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	10	Total	O	0	0
			10	10		

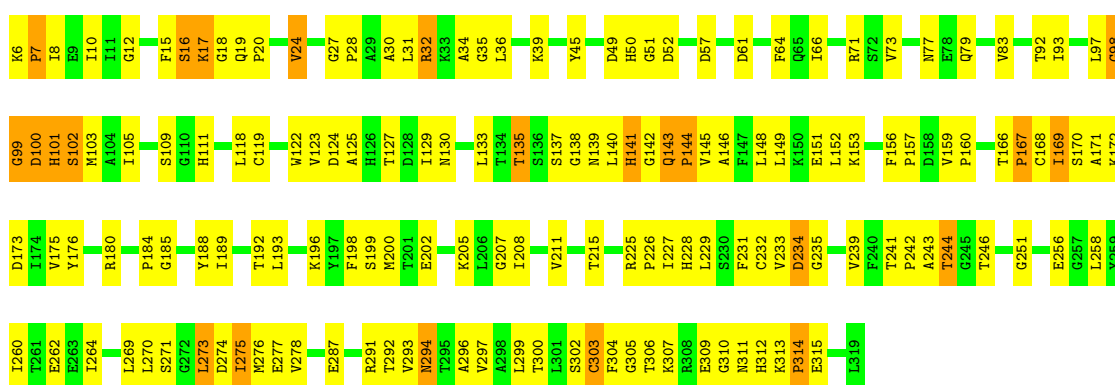
3 Residue-property plots

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.

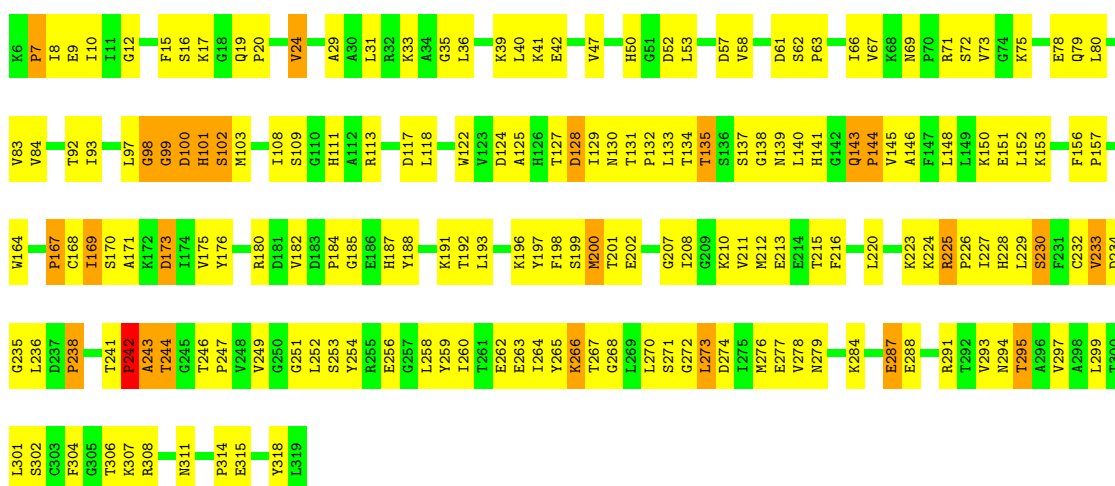
• Molecule 1: Arginase 1

Chain A:



• Molecule 1: Arginase 1

Chain B:



• Molecule 1: Arginase 1

Chain C:



P314	T244	P167	G6	K6
L319	G245	C168	L97	P7
	G246	I169	G98	I8
			G99	E9
	L252	I174	D100	I10
	S253	V175	H101	I11
	S254		S102	G12
	R255	G178	M103	
	G256	L179	A104	F15
	G257	R180	I105	S16
	L258			K17
	G259	D183	I108	G18
	L260	P184		I19
	E261	G185	H111	P20
	E262	E186		
	G263	H187	L118	G23
	T264	I188	C119	V24
	V265	I189	V120	E25
	G266	I190		K26
	T267	K191	A125	G27
	G268	T192	H126	P28
	L269	L193	T127	
	L270		D128	L31
	S271	S199	I129	R32
	G272	M200	N130	K33
	L273	T201	T131	A34
	D274	E202	P132	G35
	L275	V203	L133	L36
	N276	D204	T134	V37
	E277	K205	T135	E38
	N278		S136	K39
	N279	V211	S137	L40
		M212	G138	
			N139	Y45
	L282	T215	L140	
	G283	S217	G142	D52
	K284	L218	P143	
	E287	L219	P144	F55
	E288	L220	V145	S56
	N289	G221	A146	D57
	T290		F147	
	K291	P226	L148	F64
	T292	L227	L149	G65
	V293	H228	K150	I66
	N294	L229	E151	V67
	T295	S230	L152	
	A296	F231	K153	V73
	V297	C232		
	A298	V233	F156	N77
	L299	D234	P157	E78
		G235	D158	Q79
	G303	L236	V159	
	F304	D237	P160	V83
	G305	P238	G161	E86
	K307	V239	F162	
		F240	S163	N90
	N311	T241	W164	G91
	K312	P242	V165	T92
	V313	A243	T166	I02

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	90.94Å 90.94Å 105.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.77 – 3.20	Depositor
% Data completeness (in resolution range)	92.9 (29.77-3.20)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.286 , 0.315	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7221	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN

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	0.49	0/2447	0.77	1/3322 (0.0%)
1	B	0.51	0/2447	0.80	2/3322 (0.1%)
1	C	0.53	0/2447	0.81	1/3322 (0.0%)
All	All	0.51	0/7341	0.79	4/9966 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	99	GLY	N-CA-C	10.12	138.39	113.10
1	B	99	GLY	N-CA-C	8.88	135.29	113.10
1	C	99	GLY	N-CA-C	8.43	134.16	113.10
1	B	242	PRO	CA-N-CD	-5.05	104.43	111.50

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	98	GLY	Peptide
1	B	98	GLY	Peptide
1	C	98	GLY	Peptide

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	A	2394	0	2421	186	0
1	B	2394	0	2421	194	0
1	C	2394	0	2421	192	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	11	0	0	0	0
3	B	12	0	0	0	0
3	C	10	0	0	5	0
All	All	7221	0	7263	556	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 38.

All (556) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:227:ILE:HG13	1:A:270:LEU:HA	1.33	1.08
1:A:311:ASN:O	1:B:184:PRO:HA	1.53	1.05
1:B:238:PRO:O	1:B:242:PRO:HD3	1.56	1.05
1:C:211:VAL:O	1:C:215:THR:HG23	1.63	0.99
1:C:143:GLN:H	1:C:144:PRO:HD3	1.29	0.97
1:A:143:GLN:H	1:A:144:PRO:CD	1.77	0.96
1:A:12:GLY:HA3	1:A:52:ASP:OD1	1.67	0.94
1:B:16:SER:HB3	1:B:24:VAL:HG23	1.52	0.92
1:B:258:LEU:O	1:B:262:GLU:HG3	1.71	0.90
1:A:307:LYS:H	1:A:311:ASN:HD21	1.21	0.87
1:B:243:ALA:HB1	1:B:279:ASN:O	1.74	0.87
1:A:227:ILE:CG1	1:A:270:LEU:HA	2.03	0.86

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:12:GLY:HA3	1:B:52:ASP:OD1	1.74	0.85
1:B:102:SER:HA	1:B:144:PRO:HG3	1.58	0.85
1:A:153:LYS:HD2	1:A:167:PRO:HG2	1.57	0.85
1:B:103:MET:HB2	1:B:276:MET:CE	2.07	0.85
1:C:135:THR:HB	1:C:137:SER:O	1.77	0.84
1:C:32:ARG:HH11	1:C:32:ARG:HG3	1.43	0.84
1:B:143:GLN:H	1:B:144:PRO:HD3	1.40	0.83
1:C:256:GLU:O	1:C:260:ILE:HG13	1.76	0.83
1:A:100:ASP:HA	1:A:277:GLU:HG2	1.60	0.82
1:A:258:LEU:O	1:A:262:GLU:HG3	1.78	0.82
1:B:256:GLU:O	1:B:260:ILE:HG13	1.80	0.82
1:A:16:SER:HB3	1:A:24:VAL:HG23	1.60	0.81
1:A:233:VAL:HG13	1:A:241:THR:HB	1.59	0.81
1:A:143:GLN:H	1:A:144:PRO:HD3	1.44	0.80
1:C:103:MET:HB2	1:C:276:MET:CE	2.12	0.80
1:A:300:THR:HA	1:A:303:CYS:SG	2.22	0.80
1:B:100:ASP:HA	1:B:277:GLU:CD	2.02	0.80
1:B:173:ASP:HB3	1:B:225:ARG:HH22	1.45	0.80
1:A:102:SER:HA	1:A:144:PRO:HG3	1.64	0.79
1:B:93:ILE:HG23	1:B:271:SER:O	1.83	0.78
1:C:143:GLN:H	1:C:144:PRO:CD	1.97	0.78
1:C:100:ASP:HA	1:C:277:GLU:CD	2.04	0.78
1:B:71:ARG:O	1:B:75:LYS:HG3	1.84	0.78
1:C:148:LEU:HD22	1:C:168:CYS:SG	2.23	0.77
1:A:36:LEU:HD13	1:A:297:VAL:HG22	1.64	0.77
1:A:50:HIS:ND1	1:A:50:HIS:O	2.17	0.77
1:C:143:GLN:N	1:C:144:PRO:CD	2.46	0.77
1:A:103:MET:HB2	1:A:276:MET:CE	2.15	0.77
1:C:67:VAL:CG1	1:C:140:LEU:HD21	2.15	0.77
1:A:227:ILE:HG12	1:A:269:LEU:C	2.05	0.77
1:C:129:ILE:HG13	1:C:145:VAL:HB	1.67	0.77
1:A:229:LEU:HD22	1:A:264:ILE:CD1	2.16	0.76
1:B:118:LEU:HD12	1:B:118:LEU:C	2.06	0.76
1:A:8:ILE:HD13	1:A:93:ILE:HB	1.67	0.76
1:B:175:VAL:HG22	1:B:196:LYS:HB3	1.68	0.75
1:C:143:GLN:N	1:C:144:PRO:HD3	2.01	0.75
1:A:135:THR:HB	1:A:137:SER:O	1.86	0.75
1:C:291:ARG:O	1:C:295:THR:HG23	1.84	0.75
1:A:133:LEU:HD11	1:A:156:PHE:HA	1.68	0.75
1:A:119:CYS:SG	1:A:227:ILE:HG22	2.27	0.75
1:A:175:VAL:HG13	1:A:196:LYS:O	1.85	0.75
1:C:119:CYS:SG	1:C:219:LEU:HD13	2.27	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:207:GLY:O	1:B:211:VAL:HG23	1.87	0.74
1:C:32:ARG:NH1	1:C:32:ARG:HG3	1.99	0.74
1:C:16:SER:HB3	1:C:24:VAL:HG23	1.70	0.74
1:B:117:ASP:O	1:B:225:ARG:NH1	2.22	0.73
1:B:211:VAL:O	1:B:215:THR:HG23	1.89	0.73
1:C:203:VAL:O	1:C:203:VAL:HG12	1.87	0.73
1:B:227:ILE:HB	1:B:270:LEU:HA	1.71	0.73
1:A:275:ILE:CG2	1:A:275:ILE:O	2.36	0.73
1:C:288:GLU:HG2	1:C:291:ARG:NH2	2.04	0.73
1:B:314:PRO:O	1:B:315:GLU:HB2	1.88	0.73
1:C:19:GLN:HB2	1:C:20:PRO:HD2	1.72	0.72
1:C:307:LYS:H	1:C:311:ASN:HD21	1.38	0.72
1:C:130:ASN:O	1:C:146:ALA:HB2	1.90	0.72
1:B:124:ASP:OD2	1:B:125:ALA:N	2.21	0.72
1:B:232:CYS:O	1:B:235:GLY:N	2.20	0.72
1:B:66:ILE:HG22	1:B:138:GLY:HA2	1.72	0.72
1:A:227:ILE:HD12	1:A:264:ILE:CG2	2.20	0.72
1:B:103:MET:HB2	1:B:276:MET:HE3	1.71	0.72
1:A:103:MET:HB2	1:A:276:MET:HE3	1.71	0.71
1:B:143:GLN:H	1:B:144:PRO:CD	2.03	0.71
1:B:101:HIS:HD2	1:B:276:MET:HG3	1.55	0.71
1:A:19:GLN:HB2	1:A:20:PRO:HD2	1.71	0.71
1:B:173:ASP:CB	1:B:225:ARG:HH22	2.04	0.71
1:A:275:ILE:HD11	1:A:296:ALA:O	1.90	0.71
1:B:233:VAL:HG21	1:B:278:VAL:HG22	1.73	0.71
1:C:79:GLN:O	1:C:83:VAL:HG23	1.88	0.71
1:A:198:PHE:CD2	1:A:211:VAL:HG13	2.26	0.71
1:B:103:MET:HB2	1:B:276:MET:HE1	1.72	0.70
1:A:135:THR:HG21	1:A:143:GLN:HE21	1.56	0.70
1:C:15:PHE:CZ	1:C:17:LYS:HB2	2.26	0.70
1:C:129:ILE:CG1	1:C:145:VAL:HB	2.22	0.70
1:A:198:PHE:CE1	1:A:215:THR:HG22	2.27	0.70
1:A:135:THR:CG2	1:A:143:GLN:HE21	2.05	0.70
1:B:308:ARG:NH1	1:C:200:MET:HB3	2.06	0.70
1:B:236:LEU:HB3	1:B:252:LEU:O	1.92	0.69
1:B:232:CYS:SG	1:B:234:ASP:CG	2.70	0.69
1:A:229:LEU:HD22	1:A:264:ILE:HD12	1.74	0.69
1:B:232:CYS:SG	1:B:234:ASP:OD2	2.51	0.69
1:A:143:GLN:N	1:A:144:PRO:CD	2.51	0.69
1:C:12:GLY:HA3	1:C:52:ASP:OD1	1.92	0.68
1:B:295:THR:O	1:B:299:LEU:HG	1.92	0.68
1:A:227:ILE:HD12	1:A:264:ILE:HG23	1.73	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:233:VAL:HG13	1:B:241:THR:HB	1.74	0.68
1:C:103:MET:HB2	1:C:276:MET:HE3	1.74	0.68
1:C:8:ILE:HD13	1:C:93:ILE:HB	1.74	0.68
1:C:129:ILE:HB	1:C:145:VAL:HB	1.75	0.68
1:C:100:ASP:HA	1:C:277:GLU:OE1	1.93	0.68
1:B:175:VAL:HA	1:B:196:LYS:O	1.93	0.68
1:C:287:GLU:HG3	1:C:288:GLU:N	2.08	0.68
1:A:180:ARG:NH2	1:A:235:GLY:O	2.27	0.68
1:A:256:GLU:O	1:A:260:ILE:HG13	1.94	0.68
1:C:153:LYS:HD2	1:C:167:PRO:HG2	1.76	0.68
1:A:242:PRO:HD2	1:A:292:THR:OG1	1.94	0.67
1:B:19:GLN:HB2	1:B:20:PRO:HD2	1.77	0.67
1:A:127:THR:OG1	1:A:129:ILE:HG22	1.94	0.67
1:A:260:ILE:O	1:A:264:ILE:HG13	1.94	0.67
1:A:227:ILE:HG12	1:A:269:LEU:O	1.94	0.67
1:A:118:LEU:C	1:A:118:LEU:HD12	2.15	0.67
1:A:105:ILE:HD11	1:A:144:PRO:HA	1.76	0.67
1:C:200:MET:HE1	1:C:252:LEU:HD21	1.76	0.67
1:B:16:SER:CB	1:B:24:VAL:HG23	2.24	0.66
1:C:67:VAL:HG13	1:C:140:LEU:HD21	1.77	0.66
1:C:230:SER:HA	1:C:274:ASP:HB2	1.78	0.66
1:A:111:HIS:CE1	1:A:271:SER:HB3	2.31	0.66
1:B:108:ILE:HG12	1:B:228:HIS:NE2	2.11	0.66
1:B:318:TYR:HB2	1:C:188:TYR:CD2	2.31	0.66
1:C:103:MET:HB2	1:C:276:MET:HE1	1.76	0.66
1:B:227:ILE:O	1:B:270:LEU:HD12	1.96	0.66
1:B:180:ARG:NH2	1:B:235:GLY:O	2.29	0.66
1:B:288:GLU:HG2	1:B:291:ARG:NH2	2.09	0.66
1:C:129:ILE:HG12	1:C:129:ILE:O	1.95	0.65
1:A:233:VAL:HG13	1:A:241:THR:CB	2.27	0.65
1:A:275:ILE:CD1	1:A:296:ALA:O	2.45	0.64
1:B:127:THR:OG1	1:B:129:ILE:HG22	1.97	0.64
1:A:20:PRO:HD3	1:A:139:ASN:OD1	1.97	0.64
1:C:200:MET:CE	1:C:252:LEU:HD21	2.27	0.64
1:B:111:HIS:CE1	1:B:271:SER:HB3	2.33	0.64
1:A:139:ASN:HB3	1:A:141:HIS:CD2	2.32	0.64
1:A:275:ILE:HG22	1:A:275:ILE:O	1.98	0.64
1:C:67:VAL:HG11	1:C:140:LEU:HD21	1.79	0.64
1:B:238:PRO:O	1:B:242:PRO:CD	2.41	0.64
1:C:200:MET:HE2	1:C:252:LEU:HD23	1.79	0.64
1:A:180:ARG:CZ	1:A:251:GLY:HA2	2.27	0.64
1:B:101:HIS:CD2	1:B:276:MET:HG3	2.32	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:230:SER:HA	1:B:274:ASP:HB2	1.78	0.63
1:C:233:VAL:HG21	1:C:278:VAL:HG22	1.80	0.63
1:C:129:ILE:CB	1:C:145:VAL:HB	2.29	0.63
1:C:100:ASP:OD1	1:C:102:SER:N	2.31	0.63
1:B:143:GLN:N	1:B:144:PRO:CD	2.60	0.63
1:B:187:HIS:O	1:B:191:LYS:HG2	1.99	0.63
1:C:135:THR:HG21	1:C:143:GLN:NE2	2.13	0.63
1:A:233:VAL:CG1	1:A:241:THR:HB	2.28	0.63
1:C:7:PRO:HB2	1:C:92:THR:HG22	1.80	0.63
1:A:211:VAL:O	1:A:215:THR:HG23	1.99	0.63
1:B:233:VAL:HG21	1:B:278:VAL:CG2	2.29	0.63
1:B:223:LYS:CE	1:B:225:ARG:HD2	2.29	0.63
1:C:200:MET:HE2	1:C:252:LEU:CD2	2.28	0.62
1:C:135:THR:HG21	1:C:143:GLN:HE21	1.63	0.62
1:B:66:ILE:HG22	1:B:138:GLY:CA	2.28	0.62
1:C:129:ILE:HD12	1:C:145:VAL:HG11	1.81	0.62
1:A:139:ASN:HB3	1:A:141:HIS:HD2	1.64	0.62
1:C:102:SER:HA	1:C:144:PRO:HG3	1.80	0.62
1:A:8:ILE:HD11	1:A:93:ILE:HD12	1.81	0.62
1:C:19:GLN:HB3	1:C:141:HIS:CD2	2.34	0.62
1:C:233:VAL:CG1	1:C:241:THR:HB	2.30	0.62
1:A:123:VAL:HG11	1:A:231:PHE:CE2	2.35	0.62
1:C:83:VAL:O	1:C:86:GLU:HB3	1.98	0.61
1:B:254:TYR:CE1	1:B:295:THR:HB	2.35	0.61
1:C:265:TYR:HD1	3:C:802:HOH:O	1.84	0.61
1:B:150:LYS:HG3	1:B:169:ILE:O	2.00	0.61
1:B:111:HIS:HE1	1:B:271:SER:HB3	1.66	0.61
1:A:305:GLY:O	1:A:306:THR:C	2.38	0.61
1:C:180:ARG:NH2	1:C:235:GLY:O	2.33	0.61
1:B:79:GLN:O	1:B:83:VAL:HG23	2.01	0.61
1:C:131:THR:C	1:C:133:LEU:H	2.03	0.61
1:B:273:LEU:HD22	1:B:274:ASP:N	2.17	0.60
1:C:234:ASP:HB3	1:C:244:THR:HG21	1.82	0.60
1:C:148:LEU:O	1:C:167:PRO:HA	2.01	0.60
1:C:187:HIS:O	1:C:191:LYS:HG2	2.00	0.60
1:A:275:ILE:HD11	1:A:300:THR:OG1	2.01	0.60
1:B:130:ASN:HB3	1:B:135:THR:HG23	1.82	0.60
1:A:31:LEU:HD13	1:A:97:LEU:HD22	1.84	0.60
1:A:103:MET:HB2	1:A:276:MET:HE1	1.84	0.60
1:C:233:VAL:HG13	1:C:241:THR:HB	1.83	0.60
1:B:199:SER:O	1:B:202:GLU:N	2.32	0.60
1:A:100:ASP:HA	1:A:277:GLU:CG	2.30	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:232:CYS:SG	1:C:234:ASP:CG	2.81	0.60
1:A:228:HIS:HE1	1:A:274:ASP:OD2	1.85	0.59
1:C:288:GLU:HG2	1:C:291:ARG:HH21	1.66	0.59
1:A:227:ILE:CG1	1:A:270:LEU:CA	2.79	0.59
1:B:101:HIS:O	1:B:103:MET:N	2.35	0.59
1:C:118:LEU:C	1:C:118:LEU:HD12	2.23	0.59
1:C:127:THR:CB	1:C:129:ILE:HG22	2.33	0.59
1:A:135:THR:OG1	1:A:143:GLN:NE2	2.33	0.59
1:B:233:VAL:C	1:B:235:GLY:H	2.04	0.59
1:B:307:LYS:H	1:B:311:ASN:HD21	1.49	0.59
1:C:217:SER:O	1:C:221:GLY:HA3	2.03	0.58
1:C:297:VAL:CG1	3:C:804:HOH:O	2.51	0.58
1:C:7:PRO:CB	1:C:92:THR:HG22	2.34	0.58
1:C:226:PRO:HB3	1:C:271:SER:HB2	1.85	0.58
1:A:198:PHE:HA	1:A:202:GLU:HG2	1.85	0.58
1:A:149:LEU:HD12	1:A:152:LEU:HD13	1.86	0.58
1:A:185:GLY:O	1:A:188:TYR:HB3	2.04	0.58
1:C:130:ASN:O	1:C:146:ALA:CB	2.51	0.58
1:B:20:PRO:HD3	1:B:139:ASN:OD1	2.04	0.58
1:B:173:ASP:HB3	1:B:225:ARG:NH2	2.18	0.57
1:B:146:ALA:HA	1:B:152:LEU:CD2	2.34	0.57
1:C:139:ASN:HB3	1:C:141:HIS:CD2	2.39	0.57
1:C:212:MET:HE2	1:C:260:ILE:HA	1.86	0.57
1:C:261:THR:HG22	1:C:303:CYS:SG	2.44	0.57
1:A:229:LEU:HD22	1:A:264:ILE:HD13	1.86	0.57
1:C:139:ASN:HB3	1:C:141:HIS:HD2	1.68	0.57
1:A:16:SER:CB	1:A:24:VAL:HG23	2.34	0.57
1:B:263:GLU:OE2	1:B:266:LYS:HD3	2.04	0.57
1:A:135:THR:CG2	1:A:143:GLN:NE2	2.68	0.57
1:B:223:LYS:HE3	1:B:225:ARG:HD2	1.86	0.57
1:B:308:ARG:HH11	1:C:200:MET:HB3	1.67	0.56
1:B:262:GLU:OE1	1:C:201:THR:HG22	2.05	0.56
1:B:129:ILE:HG23	1:B:129:ILE:O	2.05	0.56
1:A:7:PRO:HB2	1:A:92:THR:HA	1.86	0.56
1:A:312:HIS:HB3	1:B:184:PRO:O	2.05	0.56
1:C:125:ALA:HA	1:C:178:GLY:O	2.06	0.56
1:B:131:THR:H	1:B:134:THR:HG1	1.52	0.56
1:B:100:ASP:HA	1:B:277:GLU:OE1	2.06	0.56
1:A:129:ILE:HG13	1:A:145:VAL:HB	1.87	0.56
1:B:156:PHE:HB2	1:B:157:PRO:HD2	1.86	0.56
1:A:35:GLY:O	1:A:39:LYS:HG3	2.06	0.56
1:A:140:LEU:O	1:A:144:PRO:HD3	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:310:GLY:HA3	1:B:182:VAL:HB	1.88	0.55
1:A:184:PRO:HA	1:C:311:ASN:O	2.06	0.55
1:A:229:LEU:CD2	1:A:264:ILE:HD12	2.34	0.55
1:A:77:ASN:OD1	1:A:103:MET:HA	2.05	0.55
1:A:143:GLN:H	1:A:144:PRO:HD2	1.65	0.55
1:B:131:THR:O	1:B:133:LEU:N	2.39	0.55
1:B:293:VAL:HG12	1:B:294:ASN:N	2.22	0.55
1:B:185:GLY:O	1:B:188:TYR:HB3	2.06	0.55
1:B:284:LYS:HB2	1:B:288:GLU:OE1	2.07	0.55
1:B:131:THR:C	1:B:133:LEU:H	2.10	0.55
1:A:10:ILE:N	1:A:10:ILE:HD12	2.21	0.55
1:A:36:LEU:HD21	1:A:97:LEU:HD21	1.89	0.55
1:C:200:MET:CE	1:C:252:LEU:CD2	2.85	0.55
1:A:8:ILE:CD1	1:A:93:ILE:HD12	2.36	0.55
1:B:192:THR:HG22	1:B:193:LEU:HD12	1.88	0.55
1:A:199:SER:O	1:A:202:GLU:HB2	2.07	0.54
1:B:151:GLU:OE1	1:B:151:GLU:N	2.40	0.54
1:A:123:VAL:HG11	1:A:231:PHE:CZ	2.43	0.54
1:C:32:ARG:HH11	1:C:32:ARG:CG	2.15	0.54
1:C:200:MET:HA	1:C:200:MET:CE	2.38	0.54
1:B:254:TYR:HE1	1:B:295:THR:HB	1.73	0.54
1:A:314:PRO:O	1:A:315:GLU:HB2	2.07	0.54
1:C:260:ILE:O	1:C:264:ILE:HG13	2.08	0.54
1:B:39:LYS:HD3	1:B:301:LEU:HD11	1.90	0.54
1:C:297:VAL:HG12	3:C:804:HOH:O	2.07	0.54
1:A:299:LEU:O	1:A:302:SER:HB3	2.08	0.53
1:A:124:ASP:OD1	1:A:232:CYS:HB2	2.07	0.53
1:A:15:PHE:CZ	1:A:73:VAL:HA	2.42	0.53
1:B:39:LYS:O	1:B:42:GLU:HB2	2.09	0.53
1:C:127:THR:OG1	1:C:129:ILE:HG22	2.08	0.53
1:B:137:SER:HB2	1:B:139:ASN:ND2	2.23	0.53
1:C:242:PRO:CG	1:C:288:GLU:HB3	2.38	0.53
1:B:15:PHE:CZ	1:B:17:LYS:HB2	2.44	0.53
1:B:8:ILE:HD13	1:B:93:ILE:HB	1.91	0.53
1:B:131:THR:C	1:B:133:LEU:N	2.62	0.53
1:A:64:PHE:CD2	1:A:159:VAL:HG22	2.43	0.53
1:A:15:PHE:CZ	1:A:17:LYS:HB2	2.44	0.53
1:A:244:THR:HG23	1:A:277:GLU:O	2.08	0.53
1:B:124:ASP:CG	1:B:232:CYS:HB3	2.29	0.53
1:A:192:THR:HG22	1:A:193:LEU:HD12	1.90	0.53
1:A:105:ILE:O	1:A:109:SER:HB2	2.09	0.53
1:C:233:VAL:HG21	1:C:278:VAL:CG2	2.38	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:174:ILE:HG12	1:C:175:VAL:N	2.24	0.52
1:B:135:THR:HB	1:B:137:SER:O	2.09	0.52
1:A:79:GLN:O	1:A:83:VAL:HG23	2.09	0.52
1:C:120:VAL:HG22	1:C:228:HIS:HB3	1.92	0.52
1:A:31:LEU:CD1	1:A:97:LEU:HD22	2.39	0.52
1:B:223:LYS:HE2	1:B:225:ARG:HD2	1.91	0.52
1:A:6:LYS:O	1:A:45:TYR:HD2	1.93	0.52
1:A:258:LEU:HD23	1:A:299:LEU:HD23	1.91	0.52
1:C:105:ILE:HD11	1:C:144:PRO:HA	1.89	0.52
1:B:137:SER:C	1:B:139:ASN:H	2.13	0.52
1:C:307:LYS:H	1:C:311:ASN:ND2	2.05	0.52
1:A:98:GLY:HA3	1:A:276:MET:SD	2.49	0.52
1:B:12:GLY:CA	1:B:52:ASP:OD1	2.54	0.52
1:A:133:LEU:CD1	1:A:156:PHE:HA	2.39	0.52
1:B:229:LEU:HD22	1:B:264:ILE:HD12	1.92	0.52
1:C:133:LEU:HD11	1:C:156:PHE:HA	1.91	0.52
1:B:265:TYR:O	1:B:268:GLY:N	2.40	0.51
1:B:233:VAL:C	1:B:235:GLY:N	2.63	0.51
1:A:101:HIS:O	1:A:103:MET:N	2.43	0.51
1:C:199:SER:O	1:C:200:MET:C	2.45	0.51
1:C:258:LEU:HD23	1:C:299:LEU:HD23	1.93	0.51
1:B:129:ILE:HG13	1:B:145:VAL:HB	1.93	0.51
1:C:232:CYS:HG	1:C:234:ASP:CG	2.14	0.51
1:C:131:THR:O	1:C:133:LEU:N	2.44	0.51
1:A:130:ASN:O	1:A:146:ALA:HB2	2.11	0.51
1:A:234:ASP:HB3	1:A:244:THR:HG21	1.93	0.51
1:C:240:PHE:CD1	1:C:254:TYR:HB2	2.46	0.51
1:A:311:ASN:O	1:B:184:PRO:CA	2.44	0.51
1:A:100:ASP:OD1	1:A:102:SER:N	2.43	0.51
1:C:233:VAL:HG13	1:C:241:THR:CB	2.41	0.50
1:A:16:SER:O	1:A:18:GLY:N	2.44	0.50
1:C:77:ASN:OD1	1:C:103:MET:HA	2.11	0.50
1:B:198:PHE:CD2	1:B:211:VAL:HG13	2.45	0.50
1:C:233:VAL:HG13	1:C:241:THR:CG2	2.42	0.50
1:B:199:SER:O	1:B:200:MET:C	2.48	0.50
1:B:140:LEU:HD23	1:B:143:GLN:NE2	2.26	0.50
1:A:124:ASP:OD1	1:A:232:CYS:CB	2.60	0.50
1:B:212:MET:O	1:B:216:PHE:CD2	2.64	0.50
1:B:210:LYS:O	1:B:213:GLU:HB3	2.12	0.50
1:A:156:PHE:HB2	1:A:157:PRO:HD2	1.94	0.50
1:C:185:GLY:O	1:C:188:TYR:HB3	2.11	0.50
1:B:15:PHE:CZ	1:B:73:VAL:HA	2.47	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:212:MET:O	1:C:216:PHE:CD2	2.65	0.50
1:B:148:LEU:HB3	1:B:169:ILE:CD1	2.42	0.50
1:C:239:VAL:O	1:C:239:VAL:HG12	2.12	0.50
1:C:203:VAL:CG1	1:C:203:VAL:O	2.59	0.49
1:A:307:LYS:N	1:A:311:ASN:HD21	1.99	0.49
1:A:8:ILE:CD1	1:A:93:ILE:HB	2.39	0.49
1:C:258:LEU:HA	1:C:261:THR:OG1	2.12	0.49
1:B:131:THR:N	1:B:134:THR:OG1	2.35	0.49
1:C:31:LEU:HD13	1:C:97:LEU:HD22	1.94	0.49
1:C:237:ASP:CG	1:C:238:PRO:HD2	2.33	0.49
1:A:153:LYS:HD2	1:A:167:PRO:CG	2.37	0.49
1:B:226:PRO:HB3	1:B:271:SER:HB2	1.94	0.49
1:C:32:ARG:HD3	1:C:37:VAL:HG21	1.95	0.49
1:C:131:THR:C	1:C:133:LEU:N	2.66	0.49
1:B:67:VAL:HG13	1:B:140:LEU:HD21	1.95	0.49
1:B:66:ILE:O	1:B:138:GLY:HA3	2.12	0.49
1:A:19:GLN:HB3	1:A:141:HIS:CD2	2.48	0.49
1:B:19:GLN:CB	1:B:20:PRO:HD2	2.41	0.49
1:A:207:GLY:O	1:A:211:VAL:HG23	2.13	0.49
1:B:7:PRO:CB	1:B:92:THR:HG22	2.42	0.49
1:A:124:ASP:OD2	1:A:125:ALA:N	2.45	0.49
1:A:229:LEU:O	1:A:273:LEU:HD23	2.13	0.48
1:A:135:THR:HG21	1:A:143:GLN:NE2	2.24	0.48
1:C:200:MET:O	1:C:201:THR:C	2.50	0.48
1:C:229:LEU:HD22	1:C:264:ILE:HD12	1.94	0.48
1:C:200:MET:HA	1:C:200:MET:HE3	1.94	0.48
1:A:293:VAL:HG12	1:A:294:ASN:N	2.29	0.48
1:A:18:GLY:HA3	1:A:100:ASP:CG	2.34	0.48
1:B:10:ILE:N	1:B:10:ILE:HD12	2.28	0.48
1:C:73:VAL:HG12	1:C:162:PHE:HZ	1.78	0.48
1:C:16:SER:O	1:C:18:GLY:N	2.47	0.48
1:B:146:ALA:HA	1:B:152:LEU:HD23	1.94	0.48
1:B:197:TYR:C	1:B:198:PHE:CD1	2.87	0.48
1:B:244:THR:O	1:B:247:PRO:HD3	2.14	0.48
1:B:168:CYS:SG	1:B:169:ILE:HG23	2.54	0.48
1:B:199:SER:O	1:B:201:THR:N	2.46	0.48
1:B:180:ARG:NH2	1:B:251:GLY:CA	2.77	0.48
1:C:232:CYS:SG	1:C:234:ASP:OD2	2.72	0.48
1:C:149:LEU:HD12	1:C:152:LEU:HD13	1.95	0.47
1:C:16:SER:C	1:C:18:GLY:H	2.15	0.47
1:C:192:THR:HG22	1:C:193:LEU:HD12	1.95	0.47
1:B:143:GLN:O	1:B:144:PRO:C	2.53	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:275:ILE:CD1	1:A:300:THR:OG1	2.62	0.47
1:A:16:SER:C	1:A:18:GLY:H	2.16	0.47
1:B:229:LEU:HD22	1:B:264:ILE:CD1	2.44	0.47
1:A:36:LEU:CD2	1:A:97:LEU:HD21	2.45	0.47
1:B:180:ARG:NH2	1:B:251:GLY:HA2	2.30	0.47
1:B:198:PHE:CE1	1:B:215:THR:HG22	2.50	0.47
1:C:86:GLU:O	1:C:86:GLU:HG3	2.12	0.47
1:C:188:TYR:O	1:C:189:ILE:C	2.53	0.47
1:A:173:ASP:CB	1:A:225:ARG:HH22	2.26	0.47
1:A:64:PHE:CE2	1:A:159:VAL:HG22	2.49	0.47
1:B:7:PRO:HB2	1:B:92:THR:HG22	1.95	0.47
1:A:27:GLY:O	1:A:30:ALA:HB3	2.14	0.47
1:B:232:CYS:SG	1:B:234:ASP:OD1	2.72	0.47
1:A:234:ASP:OD1	1:A:234:ASP:C	2.52	0.47
1:B:308:ARG:NH2	1:C:204:ASP:OD2	2.44	0.47
1:B:143:GLN:O	1:B:145:VAL:N	2.48	0.47
1:C:148:LEU:HA	1:C:148:LEU:HD23	1.59	0.47
1:B:53:LEU:HD22	1:B:83:VAL:HG21	1.97	0.47
1:C:237:ASP:CG	1:C:238:PRO:CD	2.83	0.47
1:A:205:LYS:O	1:A:205:LYS:HG2	2.15	0.47
1:A:180:ARG:HB3	1:A:200:MET:HG2	1.97	0.46
1:A:307:LYS:H	1:A:311:ASN:ND2	2.00	0.46
1:B:287:GLU:HG3	1:B:288:GLU:N	2.31	0.46
1:C:7:PRO:HB2	1:C:92:THR:CB	2.45	0.46
1:C:111:HIS:CE1	1:C:271:SER:HB3	2.51	0.46
1:A:32:ARG:NE	1:A:49:ASP:OD2	2.34	0.46
1:A:300:THR:C	1:A:302:SER:H	2.17	0.46
1:C:201:THR:OG1	1:C:202:GLU:N	2.49	0.46
1:C:212:MET:CE	1:C:260:ILE:HG23	2.45	0.46
1:C:258:LEU:O	1:C:262:GLU:N	2.47	0.46
1:B:122:TRP:HB3	1:B:176:TYR:CD1	2.50	0.46
1:A:273:LEU:HD22	1:A:274:ASP:N	2.29	0.46
1:A:15:PHE:CD2	1:A:103:MET:SD	3.08	0.46
1:C:8:ILE:CD1	1:C:93:ILE:HB	2.42	0.46
1:C:7:PRO:HB2	1:C:92:THR:CG2	2.45	0.46
1:A:61:ASP:OD1	1:A:71:ARG:HB2	2.15	0.46
1:B:124:ASP:OD1	1:B:232:CYS:HB2	2.16	0.46
1:C:243:ALA:HB1	1:C:279:ASN:O	2.16	0.46
1:B:98:GLY:HA3	1:B:276:MET:SD	2.56	0.46
1:C:108:ILE:HG22	1:C:108:ILE:O	2.16	0.46
1:C:233:VAL:HG13	1:C:241:THR:HG21	1.98	0.46
1:B:124:ASP:CG	1:B:232:CYS:CB	2.84	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:58:VAL:HG21	1:B:75:LYS:HE3	1.98	0.45
1:C:287:GLU:O	1:C:291:ARG:N	2.40	0.45
1:C:108:ILE:HG12	1:C:228:HIS:CD2	2.51	0.45
1:A:101:HIS:C	1:A:103:MET:N	2.69	0.45
1:C:103:MET:CB	1:C:276:MET:HE1	2.45	0.45
1:A:16:SER:C	1:A:18:GLY:N	2.69	0.45
1:B:101:HIS:C	1:B:103:MET:N	2.68	0.45
1:C:67:VAL:HG13	1:C:140:LEU:CD2	2.45	0.45
1:C:16:SER:C	1:C:18:GLY:N	2.70	0.45
1:B:108:ILE:HG12	1:B:228:HIS:CD2	2.51	0.45
1:C:265:TYR:CD1	3:C:802:HOH:O	2.56	0.45
1:B:109:SER:O	1:B:113:ARG:HG3	2.16	0.45
1:C:129:ILE:HD12	1:C:145:VAL:CG1	2.45	0.45
1:B:41:LYS:HG2	1:B:47:VAL:HB	1.97	0.45
1:B:200:MET:CE	1:B:252:LEU:HD23	2.46	0.45
1:B:139:ASN:O	1:B:143:GLN:NE2	2.42	0.45
1:B:291:ARG:O	1:B:295:THR:HG23	2.17	0.45
1:C:133:LEU:HD21	1:C:157:PRO:CD	2.47	0.45
1:C:313:LYS:HE3	1:C:313:LYS:HB2	1.78	0.45
1:A:100:ASP:OD1	1:A:102:SER:CB	2.65	0.45
1:A:111:HIS:HE1	1:A:271:SER:HB3	1.77	0.45
1:B:216:PHE:O	1:B:220:LEU:N	2.46	0.45
1:B:122:TRP:HB3	1:B:176:TYR:HD1	1.81	0.45
1:A:300:THR:C	1:A:302:SER:N	2.70	0.45
1:C:137:SER:HB2	1:C:139:ASN:ND2	2.31	0.45
1:A:142:GLY:O	1:A:143:GLN:HG2	2.17	0.45
1:B:152:LEU:CD1	1:B:193:LEU:HD21	2.46	0.45
1:B:35:GLY:O	1:B:39:LYS:HG3	2.17	0.45
1:C:168:CYS:SG	1:C:169:ILE:HG23	2.57	0.45
1:A:32:ARG:HG3	1:A:32:ARG:NH1	2.32	0.45
1:B:78:GLU:HB2	1:B:164:TRP:NE1	2.31	0.45
1:A:270:LEU:HD23	1:A:304:PHE:CE1	2.52	0.45
1:A:100:ASP:HB2	1:A:277:GLU:OE2	2.17	0.45
1:C:275:ILE:O	1:C:275:ILE:HG22	2.17	0.44
1:C:34:ALA:HB2	1:C:293:VAL:HG11	1.99	0.44
1:B:9:GLU:OE2	1:B:50:HIS:HD2	1.99	0.44
1:B:130:ASN:CB	1:B:135:THR:HG23	2.47	0.44
1:B:101:HIS:C	1:B:103:MET:H	2.20	0.44
1:A:159:VAL:HA	1:A:160:PRO:HD3	1.85	0.44
1:A:100:ASP:CA	1:A:277:GLU:HG2	2.41	0.44
1:B:129:ILE:HD12	1:B:145:VAL:HG11	1.99	0.44
1:B:263:GLU:O	1:B:267:THR:HG23	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:273:LEU:HD11	1:A:303:CYS:SG	2.58	0.44
1:C:127:THR:HB	1:C:129:ILE:CG2	2.47	0.44
1:A:28:PRO:C	1:A:30:ALA:H	2.21	0.44
1:C:10:ILE:HD12	1:C:10:ILE:N	2.33	0.44
1:C:64:PHE:CZ	1:C:158:ASP:O	2.71	0.44
1:A:97:LEU:HA	1:A:275:ILE:HG22	2.00	0.44
1:B:234:ASP:OD1	1:B:234:ASP:C	2.56	0.44
1:A:101:HIS:C	1:A:103:MET:H	2.22	0.44
1:A:105:ILE:O	1:A:109:SER:CB	2.65	0.44
1:B:233:VAL:HG12	1:B:244:THR:CG2	2.48	0.44
1:B:53:LEU:CD2	1:B:83:VAL:HG21	2.48	0.44
1:A:7:PRO:CB	1:A:92:THR:HG22	2.48	0.44
1:A:275:ILE:HG23	1:A:275:ILE:O	2.15	0.43
1:A:27:GLY:O	1:A:30:ALA:N	2.45	0.43
1:B:102:SER:HB3	1:B:141:HIS:HA	1.99	0.43
1:C:55:PHE:CZ	1:C:103:MET:CE	3.02	0.43
1:C:73:VAL:HG12	1:C:162:PHE:CZ	2.52	0.43
1:A:122:TRP:HB3	1:A:176:TYR:CD1	2.54	0.43
1:A:111:HIS:ND1	1:A:118:LEU:HD22	2.33	0.43
1:C:287:GLU:O	1:C:290:THR:N	2.52	0.43
1:C:111:HIS:ND1	1:C:228:HIS:ND1	2.62	0.43
1:C:23:GLY:O	1:C:26:LYS:HB2	2.17	0.43
1:A:239:VAL:HG12	1:A:239:VAL:O	2.18	0.43
1:A:229:LEU:CD2	1:A:264:ILE:CD1	2.92	0.43
1:B:262:GLU:O	1:B:265:TYR:HB3	2.18	0.43
1:C:284:LYS:HB2	1:C:288:GLU:OE1	2.17	0.43
1:B:131:THR:O	1:B:134:THR:N	2.43	0.43
1:C:267:THR:OG1	1:C:269:LEU:HB2	2.19	0.43
1:A:152:LEU:CD1	1:A:193:LEU:HD21	2.49	0.43
1:C:24:VAL:HB	1:C:99:GLY:HA2	2.00	0.43
1:C:151:GLU:OE1	1:C:151:GLU:N	2.45	0.43
1:A:143:GLN:O	1:A:146:ALA:N	2.48	0.43
1:A:118:LEU:C	1:A:118:LEU:CD1	2.87	0.43
1:A:34:ALA:HB2	1:A:293:VAL:HG11	2.01	0.43
1:A:262:GLU:OE1	1:B:201:THR:CG2	2.67	0.43
1:B:148:LEU:O	1:B:167:PRO:HA	2.19	0.43
1:B:302:SER:C	1:B:304:PHE:N	2.72	0.43
1:B:150:LYS:HD2	1:B:170:SER:HB3	2.01	0.42
1:B:124:ASP:OD1	1:B:232:CYS:CB	2.67	0.42
1:B:241:THR:N	1:B:242:PRO:CD	2.81	0.42
1:A:143:GLN:N	1:A:144:PRO:HD2	2.29	0.42
1:B:127:THR:O	1:B:128:ASP:C	2.58	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:32:ARG:HD3	1:C:37:VAL:CG2	2.49	0.42
1:B:58:VAL:HG23	1:B:72:SER:HA	1.99	0.42
1:A:151:GLU:N	1:A:151:GLU:OE1	2.47	0.42
1:B:61:ASP:OD2	1:B:69:ASN:HA	2.19	0.42
1:A:312:HIS:HA	1:B:184:PRO:HB3	2.01	0.42
1:A:66:ILE:HG22	1:A:138:GLY:HA2	2.01	0.42
1:B:31:LEU:HD13	1:B:97:LEU:HD22	2.00	0.42
1:B:233:VAL:HG12	1:B:244:THR:HG21	2.01	0.42
1:B:233:VAL:O	1:B:235:GLY:N	2.53	0.42
1:B:153:LYS:HD2	1:B:167:PRO:HG2	2.00	0.42
1:C:6:LYS:O	1:C:45:TYR:HD2	2.03	0.42
1:A:313:LYS:HE3	1:A:313:LYS:HB2	1.84	0.42
1:B:259:TYR:CE1	1:C:205:LYS:HB2	2.54	0.42
1:A:148:LEU:O	1:A:167:PRO:HA	2.20	0.42
1:B:40:LEU:CD2	1:B:301:LEU:HD23	2.50	0.42
1:C:31:LEU:CD1	1:C:97:LEU:HD22	2.50	0.42
1:C:149:LEU:HA	1:C:149:LEU:HD23	1.80	0.42
1:C:27:GLY:O	1:C:28:PRO:C	2.57	0.42
1:B:262:GLU:OE1	1:C:201:THR:CG2	2.67	0.42
1:B:306:THR:HA	1:B:311:ASN:HD21	1.84	0.42
1:C:19:GLN:CB	1:C:20:PRO:HD2	2.45	0.42
1:A:50:HIS:CG	1:A:50:HIS:O	2.72	0.42
1:B:273:LEU:CD2	1:B:274:ASP:N	2.83	0.42
1:B:311:ASN:O	1:C:184:PRO:HA	2.20	0.42
1:C:128:ASP:O	1:C:144:PRO:HD2	2.20	0.42
1:A:208:ILE:HA	1:A:208:ILE:HD12	1.88	0.41
1:B:307:LYS:N	1:B:311:ASN:HD21	2.18	0.41
1:C:232:CYS:O	1:C:235:GLY:N	2.45	0.41
1:A:309:GLU:O	1:B:182:VAL:HG21	2.20	0.41
1:B:139:ASN:HB3	1:B:141:HIS:CD2	2.55	0.41
1:A:19:GLN:HB2	1:A:20:PRO:CD	2.44	0.41
1:B:29:ALA:O	1:B:33:LYS:HB2	2.19	0.41
1:A:143:GLN:O	1:A:144:PRO:C	2.59	0.41
1:B:258:LEU:HD23	1:B:258:LEU:HA	1.85	0.41
1:B:175:VAL:HG13	1:B:196:LYS:O	2.20	0.41
1:C:279:ASN:ND2	1:C:282:LEU:HD12	2.35	0.41
1:B:62:SER:HA	1:B:63:PRO:HD3	1.87	0.41
1:C:164:TRP:CZ2	1:C:165:VAL:HG22	2.55	0.41
1:A:227:ILE:HG12	1:A:270:LEU:N	2.35	0.41
1:A:100:ASP:HA	1:A:277:GLU:CD	2.41	0.41
1:C:150:LYS:HA	1:C:167:PRO:HB3	2.03	0.41
1:C:203:VAL:HG22	1:C:211:VAL:HG21	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:100:ASP:OD1	1:C:102:SER:CB	2.67	0.41
1:C:19:GLN:HB2	1:C:20:PRO:CD	2.47	0.41
1:B:208:ILE:HA	1:B:208:ILE:HD12	1.85	0.41
1:A:275:ILE:HD13	1:A:296:ALA:O	2.21	0.41
1:C:129:ILE:O	1:C:129:ILE:HG23	2.21	0.41
1:C:305:GLY:O	1:C:307:LYS:HG3	2.21	0.41
1:C:6:LYS:HA	1:C:7:PRO:HD3	1.79	0.41
1:C:183:ASP:O	1:C:184:PRO:C	2.56	0.41
1:A:73:VAL:HG21	1:A:140:LEU:HD12	2.03	0.41
1:A:166:THR:O	1:A:167:PRO:C	2.59	0.41
1:A:122:TRP:N	1:A:175:VAL:O	2.44	0.41
1:C:306:THR:HA	1:C:311:ASN:ND2	2.36	0.41
1:A:10:ILE:H	1:A:10:ILE:HD12	1.85	0.41
1:A:168:CYS:SG	1:A:169:ILE:HG23	2.61	0.41
1:A:208:ILE:HB	1:A:256:GLU:HG2	2.02	0.41
1:B:229:LEU:O	1:B:273:LEU:HD23	2.21	0.41
1:B:36:LEU:HD13	1:B:297:VAL:HG22	2.02	0.41
1:C:66:ILE:O	1:C:138:GLY:HA3	2.21	0.41
1:A:170:SER:O	1:A:172:LYS:N	2.54	0.41
1:A:232:CYS:SG	1:A:234:ASP:CG	2.99	0.40
1:B:80:LEU:HD23	1:B:103:MET:O	2.21	0.40
1:C:40:LEU:HD23	1:C:40:LEU:HA	1.86	0.40
1:C:231:PHE:HZ	1:C:236:LEU:HD21	1.86	0.40
1:B:80:LEU:O	1:B:84:VAL:HG23	2.21	0.40
1:B:253:SER:OG	1:B:256:GLU:HG3	2.21	0.40
1:C:6:LYS:HG2	3:C:806:HOH:O	2.22	0.40
1:C:118:LEU:HD12	1:C:118:LEU:O	2.21	0.40
1:C:66:ILE:HG22	1:C:138:GLY:HA2	2.03	0.40
1:A:198:PHE:CZ	1:A:215:THR:CG2	3.04	0.40
1:C:152:LEU:CD1	1:C:193:LEU:HD21	2.51	0.40
1:C:129:ILE:HG13	1:C:145:VAL:CB	2.44	0.40
1:C:86:GLU:O	1:C:90:ASN:ND2	2.54	0.40
1:B:229:LEU:N	1:B:272:GLY:O	2.41	0.40
1:C:35:GLY:O	1:C:39:LYS:HG3	2.22	0.40
1:B:236:LEU:HD23	1:B:252:LEU:HB2	2.04	0.40
1:C:127:THR:O	1:C:128:ASP:C	2.60	0.40
1:A:242:PRO:O	1:A:244:THR:N	2.53	0.40
1:A:233:VAL:HG21	1:A:278:VAL:HG22	2.02	0.40
1:C:73:VAL:HG21	1:C:140:LEU:HD12	2.02	0.40
1:C:175:VAL:HG23	1:C:219:LEU:HD21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	312/314 (99%)	261 (84%)	38 (12%)	13 (4%)	4	31
1	B	312/314 (99%)	270 (86%)	29 (9%)	13 (4%)	4	31
1	C	312/314 (99%)	266 (85%)	32 (10%)	14 (4%)	4	29
All	All	936/942 (99%)	797 (85%)	99 (11%)	40 (4%)	4	30

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	GLY
1	A	143	GLN
1	B	99	GLY
1	C	99	GLY
1	A	171	ALA
1	A	189	ILE
1	B	102	SER
1	B	143	GLN
1	B	266	LYS
1	C	149	LEU
1	A	102	SER
1	A	243	ALA
1	A	291	ARG
1	B	243	ALA
1	C	17	LYS
1	C	199	SER
1	A	17	LYS
1	B	200	MET
1	B	249	VAL
1	C	100	ASP
1	C	143	GLN
1	A	294	ASN
1	B	128	ASP
1	B	144	PRO
1	B	171	ALA

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Mol	Chain	Res	Type
1	C	157	PRO
1	C	188	TYR
1	C	189	ILE
1	C	203	VAL
1	A	16	SER
1	A	51	GLY
1	C	132	PRO
1	C	180	ARG
1	B	132	PRO
1	A	167	PRO
1	A	144	PRO
1	B	167	PRO
1	C	129	ILE
1	C	160	PRO
1	B	242	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	264/264 (100%)	246 (93%)	18 (7%)	22	65
1	B	264/264 (100%)	245 (93%)	19 (7%)	21	63
1	C	264/264 (100%)	246 (93%)	18 (7%)	22	65
All	All	792/792 (100%)	737 (93%)	55 (7%)	22	65

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	PRO
1	A	24	VAL
1	A	32	ARG
1	A	57	ASP
1	A	100	ASP
1	A	101	HIS
1	A	135	THR
1	A	141	HIS

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Mol	Chain	Res	Type
1	A	169	ILE
1	A	226	PRO
1	A	234	ASP
1	A	244	THR
1	A	246	THR
1	A	273	LEU
1	A	275	ILE
1	A	287	GLU
1	A	303	CYS
1	A	314	PRO
1	B	7	PRO
1	B	24	VAL
1	B	57	ASP
1	B	100	ASP
1	B	101	HIS
1	B	135	THR
1	B	169	ILE
1	B	173	ASP
1	B	224	LYS
1	B	225	ARG
1	B	230	SER
1	B	233	VAL
1	B	238	PRO
1	B	242	PRO
1	B	244	THR
1	B	246	THR
1	B	273	LEU
1	B	287	GLU
1	B	295	THR
1	C	9	GLU
1	C	32	ARG
1	C	57	ASP
1	C	100	ASP
1	C	101	HIS
1	C	135	THR
1	C	160	PRO
1	C	169	ILE
1	C	174	ILE
1	C	184	PRO
1	C	200	MET
1	C	246	THR
1	C	253	SER

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Mol	Chain	Res	Type
1	C	273	LEU
1	C	287	GLU
1	C	295	THR
1	C	313	LYS
1	C	314	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	115	HIS
1	A	141	HIS
1	A	143	GLN
1	A	228	HIS
1	A	311	ASN
1	B	50	HIS
1	B	79	GLN
1	B	90	ASN
1	B	141	HIS
1	B	311	ASN
1	C	79	GLN
1	C	90	ASN
1	C	141	HIS
1	C	143	GLN
1	C	311	ASN

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 ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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.