



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2021 – 10:35 PM EDT

PDB ID : 1P8M
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, E.-D.; Christianson, D.W.
Deposited on : 2003-05-07
Resolution : 2.84 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

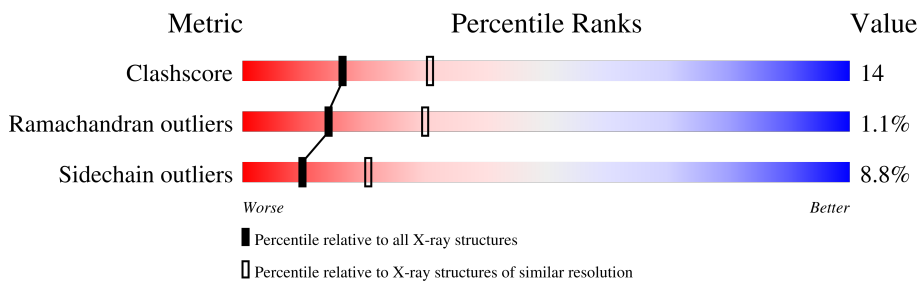
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	141614	1078 (2.86-2.82)
Ramachandran outliers	138981	1050 (2.86-2.82)
Sidechain outliers	138945	1051 (2.86-2.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	 71% 24% 5%
1	B	314	 70% 24% 5% •
1	C	314	 68% 28% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7239 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 Arginase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	6	0	0
			2397	1529	405	456	7			
1	B	314	Total	C	N	O	S	6	0	0
			2397	1529	405	456	7			
1	C	314	Total	C	N	O	S	6	0	0
			2397	1529	405	456	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	GLU	ASP	engineered mutation	UNP P07824
B	128	GLU	ASP	engineered mutation	UNP P07824
C	128	GLU	ASP	engineered mutation	UNP P07824

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	B	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	13	Total	O	0	0
			13	13		
3	B	17	Total	O	0	0
			17	17		

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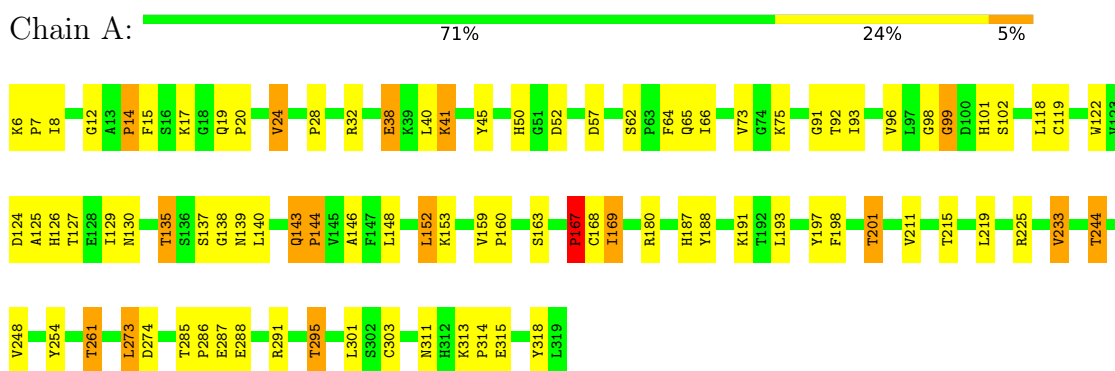
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	12	Total	O	0	0
			12	12		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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



V233	
D237	
P238	
V239	
F240	
T241	
T244	
V248	
G251	
Y254	
L258	
T261	
S271	
G272	
L273	
D274	
E277	
E287	
T295	
L299	
T300	
L301	
S302	
C303	
F304	
R308	
P314	
E315	
T316	
D317	
Y318	
L319	

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	88.49Å 88.49Å 113.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.97 – 2.84	Depositor
% Data completeness (in resolution range)	94.2 (28.97-2.84)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.286 , 0.298	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7239	wwPDB-VP
Average B, all atoms (Å ²)	74.0	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	0.76	17/2450 (0.7%)	0.79	0/3326
1	B	0.73	4/2450 (0.2%)	0.80	3/3326 (0.1%)
1	C	0.78	11/2450 (0.4%)	0.85	3/3326 (0.1%)
All	All	0.76	32/7350 (0.4%)	0.81	6/9978 (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	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	38	GLU	CB-CG	-9.93	1.33	1.52
1	B	99	GLY	N-CA	-8.51	1.33	1.46
1	B	33	LYS	CB-CG	-7.64	1.31	1.52
1	A	318	TYR	CE2-CZ	-7.60	1.28	1.38
1	A	197	TYR	CE2-CZ	-7.54	1.28	1.38
1	A	318	TYR	CE1-CZ	-7.43	1.28	1.38
1	A	45	TYR	CE2-CZ	-7.29	1.29	1.38
1	C	254	TYR	CE1-CZ	-7.25	1.29	1.38
1	A	197	TYR	CE1-CZ	-7.23	1.29	1.38
1	C	254	TYR	CG-CD2	-7.20	1.29	1.39
1	A	45	TYR	CE1-CZ	-7.13	1.29	1.38
1	C	254	TYR	CG-CD1	-7.02	1.30	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	156	PHE	CE2-CZ	-6.81	1.24	1.37
1	C	254	TYR	CE2-CZ	-6.68	1.29	1.38
1	C	213	GLU	CD-OE1	-6.68	1.18	1.25
1	A	318	TYR	CG-CD1	-6.62	1.30	1.39
1	A	45	TYR	CG-CD2	-6.54	1.30	1.39
1	B	38	GLU	CG-CD	-6.47	1.42	1.51
1	C	156	PHE	CG-CD1	-6.30	1.29	1.38
1	C	213	GLU	CD-OE2	-6.17	1.18	1.25
1	A	318	TYR	CG-CD2	-6.11	1.31	1.39
1	A	197	TYR	CG-CD1	-6.09	1.31	1.39
1	A	45	TYR	CG-CD1	-5.96	1.31	1.39
1	A	198	PHE	CE2-CZ	-5.95	1.26	1.37
1	A	197	TYR	CG-CD2	-5.92	1.31	1.39
1	C	156	PHE	CE1-CZ	-5.81	1.26	1.37
1	A	198	PHE	CE1-CZ	-5.72	1.26	1.37
1	A	198	PHE	CG-CD1	-5.54	1.30	1.38
1	C	156	PHE	CG-CD2	-5.46	1.30	1.38
1	A	198	PHE	CG-CD2	-5.33	1.30	1.38
1	C	38	GLU	CB-CG	-5.22	1.42	1.52
1	A	38	GLU	CB-CG	-5.17	1.42	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	99	GLY	N-CA-C	-11.22	85.04	113.10
1	C	213	GLU	OE1-CD-OE2	-8.39	113.24	123.30
1	B	6	LYS	C-N-CD	-7.00	105.20	120.60
1	C	237	ASP	CB-CG-OD2	6.53	124.17	118.30
1	B	89	LYS	CD-CE-NZ	-6.17	97.51	111.70
1	B	38	GLU	CB-CG-CD	-5.25	100.04	114.20

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 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	2397	0	2422	72	0
1	B	2397	0	2422	71	0
1	C	2397	0	2422	71	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	13	0	0	2	0
3	B	17	0	0	1	0
3	C	12	0	0	1	0
All	All	7239	0	7266	206	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 14.

All (206) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:THR:HG21	3:C:817:HOH:O	1.52	1.08
1:A:96:VAL:HG23	3:A:812:HOH:O	1.71	0.89
1:A:261:THR:HG22	1:A:303:CYS:SG	2.13	0.88
1:B:135:THR:HB	1:B:137:SER:O	1.74	0.88
1:B:140:LEU:O	1:B:144:PRO:HD3	1.74	0.87
1:C:135:THR:HB	1:C:137:SER:O	1.75	0.86
1:A:15:PHE:CZ	1:A:17:LYS:HB2	2.11	0.86
1:B:198:PHE:CE2	1:B:215:THR:HG22	2.11	0.86
1:A:140:LEU:O	1:A:144:PRO:HD3	1.74	0.85
1:B:261:THR:HG22	1:B:303:CYS:SG	2.16	0.85
1:A:135:THR:HB	1:A:137:SER:O	1.77	0.83
1:C:261:THR:CG2	1:C:303:CYS:SG	2.68	0.82
1:C:15:PHE:CZ	1:C:17:LYS:HB2	2.16	0.79
1:C:261:THR:HG23	1:C:303:CYS:SG	2.21	0.79
1:C:140:LEU:O	1:C:144:PRO:HD3	1.83	0.79
1:C:198:PHE:CE2	1:C:215:THR:HG22	2.18	0.78
1:B:198:PHE:HE2	1:B:215:THR:HG22	1.46	0.78
1:A:188:TYR:CD2	1:C:318:TYR:HB2	2.19	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:SER:HA	1:A:144:PRO:HG3	1.68	0.76
1:A:101:HIS:CE1	1:A:122:TRP:CZ2	2.75	0.74
1:B:102:SER:HA	1:B:144:PRO:HG3	1.68	0.74
1:B:15:PHE:CZ	1:B:17:LYS:HB2	2.22	0.73
1:A:40:LEU:HD23	1:A:301:LEU:HD23	1.71	0.73
1:A:7:PRO:HB2	1:A:92:THR:HA	1.70	0.73
1:C:211:VAL:O	1:C:215:THR:HG23	1.88	0.73
1:A:211:VAL:O	1:A:215:THR:HG23	1.91	0.70
1:C:153:LYS:HD3	1:C:167:PRO:HG2	1.72	0.70
1:C:101:HIS:CE1	1:C:122:TRP:CZ2	2.79	0.70
1:C:198:PHE:HE2	1:C:215:THR:HG22	1.57	0.70
1:A:180:ARG:HG3	1:A:248:VAL:HG11	1.72	0.69
1:A:153:LYS:HD3	1:A:167:PRO:HG2	1.73	0.69
1:B:180:ARG:HG3	1:B:248:VAL:HG11	1.74	0.69
1:B:211:VAL:O	1:B:215:THR:HG23	1.94	0.68
1:C:102:SER:HA	1:C:144:PRO:HG3	1.76	0.68
1:C:180:ARG:HG3	1:C:248:VAL:HG11	1.76	0.67
1:A:143:GLN:N	1:A:144:PRO:CD	2.58	0.67
1:A:32:ARG:HD3	3:A:820:HOH:O	1.95	0.66
1:B:143:GLN:N	1:B:144:PRO:CD	2.60	0.65
1:C:143:GLN:N	1:C:144:PRO:CD	2.59	0.65
1:B:101:HIS:CE1	1:B:122:TRP:CZ2	2.85	0.64
1:B:29:ALA:O	1:B:33:LYS:HG2	1.99	0.63
1:B:143:GLN:H	1:B:144:PRO:CD	2.12	0.63
1:B:30:ALA:HA	1:B:33:LYS:CG	2.28	0.62
1:A:15:PHE:CZ	1:A:73:VAL:HG22	2.34	0.62
1:A:143:GLN:H	1:A:144:PRO:CD	2.13	0.62
1:A:38:GLU:O	1:A:41:LYS:HB2	2.00	0.62
1:C:261:THR:HG21	1:C:299:LEU:O	1.99	0.61
1:A:66:ILE:HG22	1:A:138:GLY:CA	2.29	0.61
1:A:143:GLN:H	1:A:144:PRO:HD3	1.65	0.61
1:A:129:ILE:O	1:A:146:ALA:HB2	2.00	0.61
1:B:130:ASN:HB3	1:B:135:THR:HG23	1.83	0.60
1:C:143:GLN:H	1:C:144:PRO:HD3	1.65	0.60
1:C:187:HIS:O	1:C:191:LYS:HG2	2.02	0.59
1:B:66:ILE:HG22	1:B:138:GLY:CA	2.33	0.59
1:A:152:LEU:HD13	1:A:193:LEU:HD21	1.84	0.59
1:B:180:ARG:NH2	1:B:235:GLY:O	2.35	0.59
1:A:261:THR:CG2	1:A:303:CYS:SG	2.89	0.59
1:B:143:GLN:H	1:B:144:PRO:HD3	1.68	0.58
1:B:285:THR:OG1	1:B:288:GLU:HG3	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LEU:HD12	1:A:118:LEU:O	2.03	0.58
1:C:143:GLN:H	1:C:144:PRO:CD	2.15	0.58
1:C:261:THR:HG22	1:C:303:CYS:CA	2.34	0.57
1:B:75:LYS:HE3	3:B:822:HOH:O	2.05	0.57
1:A:130:ASN:HB3	1:A:135:THR:HG23	1.86	0.57
1:B:239:VAL:CG1	1:B:239:VAL:O	2.54	0.56
1:C:118:LEU:HD12	1:C:118:LEU:O	2.06	0.56
1:B:118:LEU:O	1:B:118:LEU:HD12	2.06	0.56
1:B:129:ILE:O	1:B:146:ALA:HB2	2.05	0.55
1:C:66:ILE:HG22	1:C:138:GLY:CA	2.36	0.55
1:A:12:GLY:HA3	1:A:52:ASP:OD1	2.07	0.55
1:C:261:THR:HG22	1:C:303:CYS:N	2.21	0.54
1:C:153:LYS:CD	1:C:167:PRO:HG2	2.37	0.54
1:B:284:LYS:O	1:B:285:THR:HG23	2.08	0.54
1:A:187:HIS:O	1:A:191:LYS:HG2	2.07	0.54
1:C:233:VAL:HG13	1:C:241:THR:HB	1.90	0.54
1:B:152:LEU:HD13	1:B:193:LEU:HD21	1.89	0.54
1:A:314:PRO:O	1:A:315:GLU:HB2	2.08	0.54
1:A:66:ILE:HG22	1:A:138:GLY:HA2	1.89	0.53
1:A:188:TYR:OH	1:C:317:ASP:OD1	2.24	0.53
1:C:125:ALA:HA	1:C:178:GLY:O	2.07	0.53
1:A:24:VAL:HB	1:A:99:GLY:HA2	1.90	0.52
1:B:64:PHE:O	1:B:65:GLN:HB2	2.08	0.52
1:C:64:PHE:O	1:C:65:GLN:HB2	2.09	0.52
1:B:233:VAL:HG12	1:B:244:THR:HG21	1.91	0.52
1:C:8:ILE:HD13	1:C:93:ILE:HB	1.91	0.52
1:B:12:GLY:HA3	1:B:52:ASP:OD1	2.10	0.52
1:B:261:THR:CG2	1:B:303:CYS:SG	2.96	0.52
1:C:152:LEU:HD13	1:C:193:LEU:HD21	1.91	0.52
1:C:261:THR:HG22	1:C:303:CYS:HA	1.93	0.51
1:C:130:ASN:HB3	1:C:135:THR:HG23	1.93	0.51
1:B:187:HIS:O	1:B:191:LYS:HG2	2.10	0.51
1:B:233:VAL:HG12	1:B:244:THR:CG2	2.41	0.51
1:B:118:LEU:HD12	1:B:118:LEU:C	2.32	0.50
1:C:125:ALA:HB1	1:C:180:ARG:HG2	1.93	0.50
1:B:30:ALA:HA	1:B:33:LYS:HG3	1.92	0.49
1:A:118:LEU:HD12	1:A:118:LEU:C	2.32	0.49
1:B:20:PRO:HD3	1:B:139:ASN:OD1	2.12	0.49
1:B:8:ILE:HD13	1:B:93:ILE:HB	1.95	0.49
1:A:14:PRO:HD3	1:A:28:PRO:HG2	1.93	0.48
1:B:29:ALA:O	1:B:33:LYS:HD3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ILE:HD13	1:A:93:ILE:HB	1.95	0.48
1:C:239:VAL:O	1:C:239:VAL:CG1	2.60	0.48
1:B:50:HIS:ND1	1:B:50:HIS:O	2.46	0.48
1:C:314:PRO:O	1:C:315:GLU:HB2	2.13	0.48
1:C:118:LEU:HD12	1:C:118:LEU:C	2.33	0.48
1:A:19:GLN:HB2	1:A:20:PRO:HD2	1.95	0.48
1:A:15:PHE:HZ	1:A:73:VAL:HG22	1.75	0.48
1:A:273:LEU:HD22	1:A:274:ASP:N	2.29	0.48
1:B:20:PRO:HD3	1:B:139:ASN:CG	2.34	0.48
1:B:259:TYR:CE1	1:C:205:LYS:HB2	2.49	0.47
1:A:188:TYR:HE2	1:C:319:LEU:HG	1.80	0.47
1:A:311:ASN:O	1:B:184:PRO:HA	2.14	0.47
1:A:233:VAL:HG12	1:A:244:THR:CG2	2.45	0.47
1:B:246:THR:O	1:B:246:THR:HG22	2.14	0.47
1:B:153:LYS:CD	1:B:167:PRO:HG2	2.45	0.46
1:B:314:PRO:O	1:B:315:GLU:HB2	2.15	0.46
1:C:302:SER:C	1:C:304:PHE:N	2.67	0.46
1:A:64:PHE:O	1:A:65:GLN:HB2	2.15	0.46
1:A:146:ALA:HA	1:A:152:LEU:HD23	1.98	0.46
1:C:146:ALA:HA	1:C:152:LEU:HD23	1.97	0.45
1:C:180:ARG:NH2	1:C:251:GLY:HA3	2.32	0.45
1:A:6:LYS:HD2	1:A:7:PRO:HD2	1.99	0.45
1:A:201:THR:CG2	1:C:308:ARG:HB2	2.47	0.44
1:A:285:THR:OG1	1:A:288:GLU:HG3	2.18	0.44
1:B:29:ALA:C	1:B:33:LYS:HD3	2.37	0.44
1:B:8:ILE:CD1	1:B:93:ILE:HB	2.48	0.44
1:B:273:LEU:HD22	1:B:274:ASP:N	2.33	0.44
1:A:6:LYS:HE2	1:A:91:GLY:O	2.18	0.44
1:A:313:LYS:HA	1:A:314:PRO:HD3	1.89	0.44
1:B:146:ALA:HA	1:B:152:LEU:HD23	1.99	0.44
1:C:226:PRO:HB3	1:C:271:SER:HB2	2.00	0.44
1:B:24:VAL:HB	1:B:99:GLY:HA2	1.99	0.44
1:A:40:LEU:CD2	1:A:301:LEU:HD23	2.43	0.44
1:B:19:GLN:HB2	1:B:20:PRO:HD2	2.00	0.44
1:B:308:ARG:O	1:C:180:ARG:O	2.35	0.44
1:B:30:ALA:C	1:B:33:LYS:HG2	2.39	0.43
1:C:148:LEU:O	1:C:169:ILE:HD13	2.17	0.43
1:A:153:LYS:CD	1:A:167:PRO:HG2	2.45	0.43
1:C:28:PRO:HD3	1:C:98:GLY:O	2.18	0.43
1:A:148:LEU:O	1:A:169:ILE:HD13	2.18	0.43
1:C:179:LEU:O	1:C:199:SER:HB2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:HIS:HA	1:C:190:ILE:CD1	2.48	0.43
1:A:143:GLN:N	1:A:144:PRO:HD2	2.32	0.43
1:B:125:ALA:HB1	1:B:180:ARG:HG2	1.99	0.43
1:B:153:LYS:HD2	1:B:167:PRO:HG2	2.01	0.43
1:C:254:TYR:CE1	1:C:258:LEU:HD11	2.54	0.43
1:C:244:THR:CG2	1:C:277:GLU:O	2.67	0.43
1:B:30:ALA:CA	1:B:33:LYS:HG2	2.49	0.43
1:A:124:ASP:HB3	1:A:126:HIS:O	2.19	0.43
1:B:143:GLN:N	1:B:144:PRO:HD2	2.33	0.43
1:C:230:SER:HA	1:C:274:ASP:HB2	2.01	0.43
1:C:244:THR:HG23	1:C:277:GLU:O	2.19	0.43
1:A:291:ARG:O	1:A:295:THR:HG22	2.19	0.42
1:C:12:GLY:HA3	1:C:52:ASP:OD1	2.18	0.42
1:C:187:HIS:HA	1:C:190:ILE:HD11	2.01	0.42
1:C:169:ILE:HD13	1:C:169:ILE:H	1.83	0.42
1:C:180:ARG:NH2	1:C:251:GLY:CA	2.82	0.42
1:A:129:ILE:O	1:A:130:ASN:C	2.57	0.42
1:C:20:PRO:HD3	1:C:139:ASN:CG	2.40	0.42
1:A:233:VAL:HG12	1:A:244:THR:HG21	2.01	0.42
1:C:19:GLN:HB2	1:C:20:PRO:HD2	2.02	0.42
1:A:6:LYS:HA	1:A:7:PRO:HD3	1.73	0.42
1:A:50:HIS:ND1	1:A:50:HIS:O	2.53	0.42
1:C:10:ILE:HD12	1:C:10:ILE:N	2.35	0.42
1:A:125:ALA:HB1	1:A:180:ARG:HG2	2.02	0.42
1:A:122:TRP:CD1	1:A:124:ASP:HB2	2.55	0.41
1:B:66:ILE:O	1:B:138:GLY:HA3	2.19	0.41
1:B:254:TYR:HE1	1:B:295:THR:HB	1.84	0.41
1:C:180:ARG:CZ	1:C:251:GLY:CA	2.98	0.41
1:B:6:LYS:HA	1:B:7:PRO:HD2	1.84	0.41
1:B:29:ALA:O	1:B:33:LYS:CD	2.68	0.41
1:C:40:LEU:HD23	1:C:301:LEU:HD23	2.01	0.41
1:C:159:VAL:HA	1:C:160:PRO:HD3	1.94	0.41
1:C:172:LYS:HA	1:C:172:LYS:HD3	1.82	0.41
1:C:149:LEU:HD23	1:C:149:LEU:HA	1.87	0.41
1:C:273:LEU:HD22	1:C:274:ASP:N	2.34	0.41
1:B:30:ALA:HA	1:B:33:LYS:HG2	2.02	0.41
1:A:7:PRO:O	1:A:8:ILE:HD13	2.20	0.41
1:B:40:LEU:HD23	1:B:301:LEU:HD23	2.02	0.41
1:B:237:ASP:HA	1:B:238:PRO:HD3	1.92	0.41
1:C:129:ILE:O	1:C:146:ALA:HB2	2.21	0.41
1:A:6:LYS:CD	1:A:7:PRO:HD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ALA:O	1:B:33:LYS:HG2	2.20	0.41
1:B:254:TYR:CE1	1:B:295:THR:HB	2.56	0.41
1:A:38:GLU:O	1:A:41:LYS:CB	2.66	0.41
1:A:201:THR:HG22	1:C:308:ARG:HB2	2.03	0.41
1:A:291:ARG:O	1:A:295:THR:CG2	2.69	0.41
1:B:129:ILE:O	1:B:130:ASN:C	2.59	0.41
1:B:226:PRO:HB3	1:B:271:SER:HB2	2.03	0.41
1:A:313:LYS:HE3	1:A:313:LYS:HB2	1.90	0.41
1:A:7:PRO:HB2	1:A:92:THR:CA	2.44	0.40
1:A:75:LYS:HA	1:A:75:LYS:HD2	1.91	0.40
1:A:159:VAL:HA	1:A:160:PRO:HD3	1.96	0.40
1:A:254:TYR:CE1	1:A:295:THR:HB	2.56	0.40
1:B:169:ILE:H	1:B:169:ILE:HD13	1.86	0.40
1:C:29:ALA:O	1:C:33:LYS:HB2	2.21	0.40
1:C:143:GLN:N	1:C:144:PRO:HD2	2.35	0.40
1:B:75:LYS:HA	1:B:75:LYS:HD2	1.93	0.40
1:B:159:VAL:HA	1:B:160:PRO:HD3	1.95	0.40
1:B:295:THR:O	1:B:299:LEU:HG	2.22	0.40
1:C:66:ILE:O	1:C:138:GLY:HA3	2.22	0.40
1:C:6:LYS:HD2	1:C:6:LYS:HA	1.93	0.40
1:B:21:ARG:HH11	1:B:282:LEU:HD13	1.87	0.40
1:C:180:ARG:CZ	1:C:251:GLY:HA3	2.51	0.40
1:A:20:PRO:HD3	1:A:139:ASN:CG	2.42	0.40
1:A:119:CYS:HB2	1:A:219:LEU:HD22	2.04	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	312/314 (99%)	296 (95%)	12 (4%)	4 (1%)	12 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	312/314 (99%)	296 (95%)	12 (4%)	4 (1%)	12	26
1	C	312/314 (99%)	295 (95%)	15 (5%)	2 (1%)	25	46
All	All	936/942 (99%)	887 (95%)	39 (4%)	10 (1%)	14	30

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	GLY
1	B	99	GLY
1	A	14	PRO
1	B	65	GLN
1	B	143	GLN
1	C	143	GLN
1	C	167	PRO
1	A	143	GLN
1	B	14	PRO
1	A	167	PRO

5.3.2 Protein sidechains [i](#)

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%)	243 (92%)	21 (8%)	12	25
1	B	264/264 (100%)	238 (90%)	26 (10%)	8	16
1	C	264/264 (100%)	241 (91%)	23 (9%)	10	21
All	All	792/792 (100%)	722 (91%)	70 (9%)	10	21

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	VAL
1	A	41	LYS
1	A	57	ASP
1	A	62	SER

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Mol	Chain	Res	Type
1	A	127	THR
1	A	135	THR
1	A	144	PRO
1	A	152	LEU
1	A	163	SER
1	A	167	PRO
1	A	168	CYS
1	A	169	ILE
1	A	201	THR
1	A	225	ARG
1	A	233	VAL
1	A	244	THR
1	A	261	THR
1	A	273	LEU
1	A	286	PRO
1	A	287	GLU
1	A	295	THR
1	B	14	PRO
1	B	24	VAL
1	B	33	LYS
1	B	56	VAL
1	B	57	ASP
1	B	62	SER
1	B	135	THR
1	B	144	PRO
1	B	152	LEU
1	B	163	SER
1	B	167	PRO
1	B	168	CYS
1	B	169	ILE
1	B	201	THR
1	B	222	ARG
1	B	223	LYS
1	B	224	LYS
1	B	225	ARG
1	B	226	PRO
1	B	233	VAL
1	B	244	THR
1	B	261	THR
1	B	273	LEU
1	B	286	PRO
1	B	287	GLU

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Mol	Chain	Res	Type
1	B	295	THR
1	C	14	PRO
1	C	24	VAL
1	C	56	VAL
1	C	57	ASP
1	C	62	SER
1	C	89	LYS
1	C	135	THR
1	C	144	PRO
1	C	152	LEU
1	C	163	SER
1	C	167	PRO
1	C	168	CYS
1	C	169	ILE
1	C	201	THR
1	C	222	ARG
1	C	225	ARG
1	C	226	PRO
1	C	231	PHE
1	C	244	THR
1	C	261	THR
1	C	273	LEU
1	C	287	GLU
1	C	295	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides 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.

No monomer is involved in short contacts.

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.