



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 08:08 pm BST

PDB ID : 1ONO
Title : IspC Mn²⁺ complex
Authors : Steinbacher, S.; Kaiser, J.; Eisenreich, W.; Huber, R.; Bacher, A.; Rohdich, F.
Deposited on : 2003-02-28
Resolution : 2.50 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) : 2.11

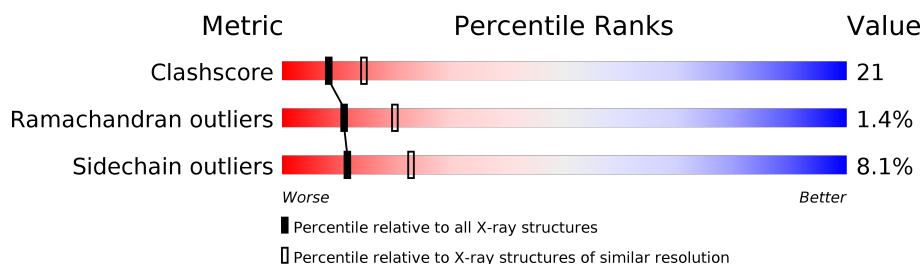
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 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	398	
1	B	398	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6115 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 1-deoxy-D-xylulose 5-phosphate reductoisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			2955	1847	518	564	26			
1	B	389	Total	C	N	O	S	0	0	0
			2955	1847	518	564	26			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	96	Total	O	0	0
			96	96		
3	B	107	Total	O	0	0
			107	107		

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.39 Å 52.25 Å 108.33 Å 90.00° 92.52° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	98.7 (20.00-2.50)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.216 , 0.287	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6115	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.48	0/2998	0.69	1/4062 (0.0%)
1	B	0.46	0/2998	0.70	0/4062
All	All	0.47	0/5996	0.70	1/8124 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	LEU	CA-CB-CG	5.17	127.19	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2955	0	2989	133	0
1	B	2955	0	2989	128	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	96	0	0	7	0
3	B	107	0	0	7	0
All	All	6115	0	5978	251	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 21.

All (251) 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:386:ARG:HB3	1:A:390:ARG:HH21	1.43	0.83
1:B:386:ARG:HB3	1:B:390:ARG:HH21	1.45	0.81
1:B:199:ALA:HA	1:B:354:THR:HG22	1.68	0.75
1:B:115:ARG:HG3	1:B:115:ARG:HH11	1.52	0.75
1:A:185:GLY:HA3	1:A:228:LYS:HE3	1.68	0.75
1:A:251:HIS:HE1	1:A:255:VAL:H	1.35	0.74
1:B:172:ASP:OD2	1:B:175:GLN:HB2	1.88	0.73
1:B:185:GLY:HA3	1:B:228:LYS:HE3	1.69	0.73
1:B:251:HIS:HE1	1:B:255:VAL:H	1.38	0.72
1:B:288:ASN:HB3	3:B:1070:HOH:O	1.91	0.71
1:B:170:TYR:CD1	3:B:1029:HOH:O	2.43	0.70
1:A:84:ALA:O	1:A:88:MET:HG2	1.91	0.70
1:A:361:LEU:O	1:A:365:GLU:HG3	1.93	0.69
1:B:136:MET:CE	1:B:169:GLY:HA3	2.23	0.69
1:A:387:GLU:O	1:A:391:LYS:HG2	1.93	0.69
1:B:236:ARG:NH2	1:B:326:GLU:OE2	2.26	0.68
1:B:170:TYR:N	1:B:170:TYR:CD2	2.59	0.68
1:A:217:LYS:HD2	1:A:343:VAL:CG1	2.24	0.68
1:B:170:TYR:N	1:B:170:TYR:HD2	1.92	0.68
1:B:226:MET:HE2	1:B:230:LEU:HG	1.76	0.68
1:A:257:HIS:HE1	3:A:1027:HOH:O	1.77	0.67
1:B:47:LEU:HD11	1:B:72:GLN:HG3	1.77	0.67
1:A:170:TYR:N	1:A:170:TYR:CD2	2.64	0.66
1:A:304:ALA:HB1	1:B:304:ALA:HB1	1.77	0.66
1:A:377:ASP:O	1:A:381:VAL:HG23	1.96	0.66
1:A:375:VAL:O	1:A:379:LEU:HG	1.96	0.65
1:B:384:ASN:O	1:B:388:VAL:HG23	1.97	0.65
1:A:190:PHE:CZ	1:A:206:ALA:HB2	2.31	0.65
1:A:236:ARG:NH2	1:A:326:GLU:OE2	2.30	0.65
1:B:216:ARG:N	1:B:219:SER:HG	1.95	0.65
1:B:10:THR:HG22	1:B:41:ARG:HG2	1.78	0.65
1:B:387:GLU:O	1:B:391:LYS:HG2	1.96	0.65
1:A:10:THR:HG22	1:A:41:ARG:HG2	1.78	0.64
1:B:199:ALA:HA	1:B:354:THR:CG2	2.28	0.64
1:A:251:HIS:CE1	1:A:255:VAL:H	2.15	0.64
1:A:384:ASN:O	1:A:388:VAL:HG23	1.97	0.64
1:B:377:ASP:O	1:B:381:VAL:HG23	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:O	1:A:141:GLN:HG2	1.98	0.64
1:B:115:ARG:NH1	1:B:115:ARG:HG3	2.11	0.63
1:B:177:GLY:HA2	1:B:263:GLN:HE21	1.64	0.63
1:B:251:HIS:CE1	1:B:255:VAL:H	2.17	0.63
1:A:172:ASP:OD2	1:A:175:GLN:HB2	1.99	0.62
1:A:236:ARG:NH2	1:A:243:ALA:HB2	2.15	0.62
1:B:389:ALA:O	1:B:393:VAL:HG23	2.00	0.62
1:A:295:LYS:HE3	3:A:1080:HOH:O	2.00	0.61
1:B:130:THR:HB	1:B:378:VAL:HG11	1.83	0.61
1:A:292:SER:OG	1:A:294:VAL:HG23	2.00	0.61
1:B:298:ASP:O	1:B:302:LEU:HD13	2.01	0.61
1:B:226:MET:O	1:B:230:LEU:HG	2.01	0.61
1:A:190:PHE:HZ	1:A:206:ALA:HB2	1.64	0.60
1:A:217:LYS:HD2	1:A:343:VAL:HG11	1.84	0.60
1:B:375:VAL:O	1:B:379:LEU:HG	2.01	0.60
1:B:285:ALA:HB2	1:B:290:VAL:HG13	1.84	0.59
1:B:217:LYS:NZ	1:B:217:LYS:HB3	2.16	0.59
1:B:66:LYS:O	1:B:70:GLN:HB2	2.03	0.59
1:A:386:ARG:HB3	1:A:390:ARG:NH2	2.15	0.59
1:B:386:ARG:HB3	1:B:390:ARG:NH2	2.14	0.59
1:A:156:ILE:O	1:A:160:LEU:HG	2.03	0.58
1:A:47:LEU:HD11	1:A:72:GLN:HG3	1.85	0.58
1:A:348:ALA:HB3	1:A:350:GLN:HE21	1.68	0.58
1:B:164:ILE:HD13	1:B:173:LEU:HD23	1.85	0.58
1:A:136:MET:CE	1:A:169:GLY:HA3	2.33	0.58
1:A:266:SER:CB	1:B:270:GLN:HE21	2.17	0.57
1:A:199:ALA:HA	1:A:354:THR:HG22	1.86	0.57
1:B:271:LEU:N	1:B:271:LEU:HD23	2.19	0.57
1:A:298:ASP:O	1:A:302:LEU:HD13	2.03	0.56
1:B:94:VAL:O	1:B:118:LYS:HE3	2.04	0.56
1:B:216:ARG:HG2	3:B:1065:HOH:O	2.04	0.56
1:A:196:ARG:HA	1:A:196:ARG:CZ	2.36	0.56
1:B:361:LEU:O	1:B:365:GLU:HG3	2.06	0.56
1:B:130:THR:HB	1:B:378:VAL:CG1	2.34	0.56
1:A:17:THR:O	1:A:21:VAL:HG23	2.05	0.55
1:A:66:LYS:O	1:A:70:GLN:HB2	2.07	0.54
1:B:137:ASP:O	1:B:141:GLN:HG2	2.08	0.54
1:B:163:PRO:HG2	1:B:176:ASN:ND2	2.22	0.54
1:B:190:PHE:CZ	1:B:206:ALA:HB2	2.43	0.54
1:B:217:LYS:HD2	1:B:343:VAL:CG1	2.38	0.54
1:B:37:LYS:HE3	1:B:58:ASP:OD2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLY:H	1:A:329:GLN:HE21	1.55	0.54
1:A:7:LEU:HB3	1:A:101:ILE:CD1	2.38	0.54
1:A:130:THR:HB	1:A:378:VAL:CG1	2.38	0.54
1:A:64:LEU:O	1:A:64:LEU:HD23	2.08	0.54
1:B:135:PHE:O	1:B:139:VAL:HG23	2.08	0.54
1:B:368:ASP:HB3	3:B:1054:HOH:O	2.06	0.54
1:B:89:ALA:HB1	1:B:113:ALA:HB2	1.89	0.54
1:B:146:LEU:O	1:B:168:LEU:HD12	2.08	0.54
1:A:270:GLN:HE21	1:B:266:SER:CB	2.21	0.54
1:A:125:LYS:HG2	1:A:234:GLU:OE2	2.09	0.53
1:A:217:LYS:NZ	1:A:217:LYS:HB3	2.22	0.53
1:B:136:MET:HE2	1:B:169:GLY:HA3	1.90	0.53
1:B:53:TYR:HD2	1:B:79:LEU:HD11	1.73	0.53
1:A:216:ARG:N	1:A:219:SER:HG	2.07	0.53
1:A:226:MET:CE	1:A:335:LEU:HD13	2.39	0.53
1:B:131:CYS:HB3	1:B:134:LEU:HD12	1.91	0.53
1:A:170:TYR:HD2	1:A:170:TYR:N	2.06	0.53
1:A:305:LEU:HB2	1:B:305:LEU:HB2	1.91	0.53
1:B:62:ALA:O	1:B:66:LYS:HG3	2.09	0.53
1:A:359:LEU:O	1:A:363:VAL:HG23	2.08	0.52
1:A:270:GLN:C	1:A:271:LEU:HD23	2.29	0.52
1:B:348:ALA:O	1:B:349:GLN:HB2	2.11	0.51
1:A:394:MET:C	1:A:396:LEU:H	2.13	0.51
1:A:39:VAL:O	1:A:43:VAL:HG23	2.10	0.51
1:A:338:ALA:O	1:A:342:THR:HG23	2.11	0.51
1:A:389:ALA:O	1:A:393:VAL:HG23	2.10	0.51
1:A:199:ALA:HA	1:A:354:THR:CG2	2.40	0.51
1:B:156:ILE:HD13	1:B:181:ILE:CG2	2.41	0.51
1:B:395:ARG:O	1:B:395:ARG:HD2	2.09	0.51
1:A:130:THR:HB	1:A:378:VAL:HG11	1.92	0.51
1:A:257:HIS:HD2	1:A:270:GLN:OE1	1.94	0.50
1:A:372:PRO:HA	1:A:377:ASP:HB3	1.93	0.50
1:B:382:ASP:O	1:B:386:ARG:HG3	2.10	0.50
1:A:72:GLN:HE21	1:A:72:GLN:HA	1.76	0.50
1:B:359:LEU:O	1:B:363:VAL:HG23	2.12	0.50
1:A:152:GLU:O	1:A:156:ILE:HG13	2.12	0.50
1:A:132:GLY:H	1:A:329:GLN:NE2	2.09	0.50
1:B:66:LYS:HG2	1:B:78:VAL:HB	1.93	0.50
1:A:18:LEU:HD22	1:A:49:PHE:CD1	2.47	0.50
1:A:93:ASP:N	1:A:93:ASP:OD1	2.42	0.50
1:B:270:GLN:C	1:B:271:LEU:HD23	2.32	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ARG:O	1:A:45:GLN:HG3	2.12	0.49
1:A:89:ALA:HB1	1:A:113:ALA:HB2	1.93	0.49
1:A:168:LEU:O	1:A:170:TYR:CE2	2.65	0.49
1:A:196:ARG:HA	1:A:196:ARG:NE	2.28	0.49
1:A:35:ALA:HB3	1:A:42:MET:HE2	1.93	0.49
1:A:329:GLN:N	1:A:371:GLU:OE1	2.44	0.49
1:B:236:ARG:NH2	1:B:243:ALA:HB2	2.27	0.49
1:B:250:ILE:O	1:B:307:PHE:HA	2.12	0.49
1:A:53:TYR:HD2	1:A:79:LEU:HD11	1.78	0.49
1:B:196:ARG:CZ	1:B:196:ARG:HA	2.42	0.49
3:A:1030:HOH:O	1:B:296:PRO:HG3	2.13	0.49
1:B:171:ALA:HB3	3:B:1088:HOH:O	2.13	0.48
1:A:36:GLY:HA2	1:A:57:ASP:OD1	2.13	0.48
1:B:348:ALA:HB3	1:B:350:GLN:HE21	1.77	0.48
1:B:194:PRO:O	1:B:197:ASP:HB2	2.13	0.48
1:A:250:ILE:O	1:A:307:PHE:HA	2.13	0.48
1:B:120:ILE:HD12	1:B:120:ILE:N	2.28	0.48
1:B:390:ARG:HG3	3:B:1102:HOH:O	2.12	0.48
1:B:64:LEU:HD23	1:B:64:LEU:O	2.13	0.48
1:A:18:LEU:HD22	1:A:49:PHE:CE1	2.49	0.48
1:B:386:ARG:O	1:B:390:ARG:HB2	2.14	0.48
1:B:75:ARG:O	1:B:75:ARG:HG3	2.13	0.48
1:A:115:ARG:NH1	1:A:115:ARG:HG3	2.28	0.48
1:A:120:ILE:HB	1:A:146:LEU:HD23	1.96	0.48
1:A:108:LEU:HD23	1:A:375:VAL:HG21	1.96	0.48
1:A:346:PHE:CG	1:A:356:ILE:HD11	2.49	0.47
1:A:299:PHE:CB	1:B:249:LEU:HD11	2.44	0.47
1:A:88:MET:HE2	1:A:88:MET:HA	1.94	0.47
1:B:30:VAL:HG11	1:B:33:LEU:HD21	1.96	0.47
1:A:226:MET:HG2	1:A:339:ASN:ND2	2.30	0.47
1:A:115:ARG:HD3	3:A:1064:HOH:O	2.14	0.47
1:A:72:GLN:NE2	1:A:72:GLN:HA	2.29	0.47
1:A:249:LEU:HD11	1:B:299:PHE:CB	2.45	0.47
1:A:88:MET:HE2	1:A:91:LEU:HD13	1.95	0.47
1:B:333:THR:HG21	1:B:378:VAL:HG13	1.97	0.47
1:A:395:ARG:O	1:A:395:ARG:HD2	2.15	0.47
1:A:143:LYS:CE	1:A:170:TYR:HE1	2.28	0.46
1:A:156:ILE:HD13	1:A:181:ILE:CG2	2.45	0.46
1:B:173:LEU:HD11	1:B:241:ALA:HB2	1.97	0.46
1:B:201:MET:HA	1:B:205:GLN:OE1	2.14	0.46
1:B:359:LEU:HD22	1:B:392:GLU:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLN:NE2	1:A:141:GLN:HA	2.30	0.46
1:A:217:LYS:HD2	1:A:343:VAL:HG12	1.94	0.46
1:A:66:LYS:HG2	1:A:78:VAL:HB	1.98	0.46
1:B:17:THR:O	1:B:21:VAL:HG23	2.15	0.46
1:B:182:LEU:HD23	1:B:259:MET:CE	2.46	0.46
1:A:228:LYS:HD3	1:A:228:LYS:HA	1.83	0.46
1:A:136:MET:HE2	1:A:169:GLY:HA3	1.97	0.46
1:A:318:LEU:O	1:A:322:MET:HG3	2.16	0.46
1:A:382:ASP:O	1:A:386:ARG:HG3	2.16	0.46
1:A:90:ALA:O	1:A:91:LEU:C	2.54	0.46
1:B:278:THR:HB	1:B:279:PRO:CD	2.46	0.46
1:A:46:CYS:O	1:A:50:SER:HA	2.17	0.45
1:B:348:ALA:O	1:B:349:GLN:CB	2.65	0.45
1:B:35:ALA:HB3	1:B:42:MET:HE2	1.97	0.45
1:A:276:MET:HE3	3:A:1075:HOH:O	2.16	0.45
1:A:370:ARG:NH2	3:A:1044:HOH:O	2.49	0.45
1:A:39:VAL:HG11	1:A:64:LEU:HD22	1.99	0.45
1:A:168:LEU:O	1:A:170:TYR:CD2	2.70	0.45
1:B:132:GLY:H	1:B:329:GLN:HE21	1.65	0.45
1:B:361:LEU:HD22	1:B:365:GLU:HG3	1.97	0.45
1:B:84:ALA:O	1:B:88:MET:HG2	2.17	0.45
1:A:226:MET:HE3	1:A:230:LEU:HG	1.98	0.45
1:B:120:ILE:HB	1:B:146:LEU:HD23	1.99	0.45
1:A:348:ALA:O	1:A:349:GLN:HB2	2.17	0.45
1:A:62:ALA:O	1:A:66:LYS:HG3	2.16	0.45
1:A:66:LYS:NZ	3:A:1035:HOH:O	2.50	0.45
1:B:217:LYS:HD2	1:B:343:VAL:HG11	1.99	0.45
1:B:41:ARG:O	1:B:45:GLN:HG3	2.17	0.44
1:B:224:THR:HG22	1:B:353:PHE:CZ	2.52	0.44
1:B:7:LEU:HA	1:B:34:VAL:HB	1.98	0.44
1:B:46:CYS:O	1:B:50:SER:HA	2.18	0.44
1:A:299:PHE:HB2	1:B:249:LEU:HD11	1.99	0.44
1:B:39:VAL:HG11	1:B:64:LEU:HD22	2.00	0.44
1:A:7:LEU:HB3	1:A:101:ILE:HG13	2.00	0.44
1:B:39:VAL:HG21	1:B:64:LEU:HD13	1.99	0.44
1:A:218:ILE:HG13	1:A:219:SER:N	2.33	0.44
1:A:226:MET:HE1	1:A:335:LEU:HD13	2.00	0.44
1:A:271:LEU:HD23	1:A:271:LEU:N	2.33	0.44
1:A:338:ALA:HB2	1:A:363:VAL:HG11	1.99	0.43
1:B:363:VAL:HG21	1:B:389:ALA:HB2	1.99	0.43
1:A:115:ARG:HH11	1:A:115:ARG:HG3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:CYS:HB2	1:A:135:PHE:CD1	2.54	0.43
1:A:386:ARG:HG2	1:A:386:ARG:HH11	1.84	0.43
1:A:124:ASN:ND2	1:A:127:SER:H	2.17	0.43
1:A:153:HIS:CE1	1:A:234:GLU:HB2	2.54	0.43
1:A:249:LEU:HD11	1:B:299:PHE:HB2	2.01	0.43
1:A:361:LEU:HD22	1:A:365:GLU:CD	2.38	0.43
1:A:266:SER:HB3	1:B:270:GLN:HE21	1.84	0.43
1:A:182:LEU:HD23	1:A:259:MET:CE	2.50	0.42
1:B:394:MET:C	1:B:396:LEU:H	2.22	0.42
1:B:342:THR:OG1	1:B:356:ILE:HG23	2.19	0.42
1:B:259:MET:HG2	1:B:269:ALA:HB2	2.01	0.42
1:B:37:LYS:NZ	1:B:57:ASP:OD2	2.53	0.42
1:A:333:THR:HG21	1:A:378:VAL:HG13	1.99	0.42
1:B:143:LYS:HZ2	1:B:170:TYR:HE1	1.68	0.42
1:B:16:SER:O	1:B:20:VAL:HG23	2.19	0.42
1:A:164:ILE:HD11	1:A:176:ASN:HB2	2.01	0.42
1:A:394:MET:C	1:A:396:LEU:N	2.73	0.42
1:B:190:PHE:HZ	1:B:206:ALA:HB2	1.84	0.42
1:B:217:LYS:HZ2	1:B:217:LYS:HB3	1.85	0.42
1:B:182:LEU:HD23	1:B:259:MET:HE3	2.02	0.42
1:B:329:GLN:N	1:B:371:GLU:OE1	2.53	0.42
1:B:53:TYR:CD2	1:B:79:LEU:HD11	2.52	0.42
1:A:7:LEU:C	1:A:101:ILE:HG13	2.40	0.42
1:A:174:GLU:HG3	1:A:175:GLN:N	2.35	0.42
1:B:251:HIS:CE1	1:B:256:ILE:H	2.38	0.41
1:B:220:VAL:O	1:B:224:THR:HG23	2.20	0.41
1:B:362:SER:O	1:B:363:VAL:C	2.56	0.41
1:A:195:LEU:O	1:A:197:ASP:N	2.54	0.41
1:A:386:ARG:O	1:A:390:ARG:HB2	2.20	0.41
1:B:338:ALA:HB2	1:B:363:VAL:HG11	2.02	0.41
1:B:65:LEU:HD12	1:B:69:LEU:HG	2.03	0.41
1:A:278:THR:HB	1:A:279:PRO:CD	2.50	0.41
1:B:226:MET:O	1:B:226:MET:HE2	2.20	0.41
1:A:120:ILE:HD12	1:A:120:ILE:N	2.36	0.41
1:B:7:LEU:C	1:B:101:ILE:HG13	2.41	0.41
1:B:194:PRO:HG2	1:B:197:ASP:OD1	2.20	0.41
1:B:226:MET:HE1	1:B:335:LEU:HD13	2.03	0.41
1:B:115:ARG:CG	1:B:115:ARG:HH11	2.25	0.41
1:A:7:LEU:O	1:A:100:ALA:HB3	2.21	0.41
1:B:195:LEU:O	1:B:197:ASP:N	2.54	0.41
1:A:277:ARG:HB3	1:A:292:SER:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:PHE:HZ	1:B:267:VAL:HG11	1.86	0.41
1:A:226:MET:HG2	1:A:339:ASN:CG	2.41	0.41
1:A:118:LYS:O	1:A:120:ILE:HD12	2.22	0.40
1:A:108:LEU:N	1:A:109:PRO:CD	2.84	0.40
1:B:118:LYS:O	1:B:120:ILE:HD12	2.22	0.40
1:B:90:ALA:O	1:B:91:LEU:C	2.60	0.40
1:B:65:LEU:CD1	1:B:69:LEU:HG	2.51	0.40
1:A:346:PHE:CD2	1:A:356:ILE:HD11	2.57	0.40
1:B:329:GLN:HG3	3:B:1011:HOH:O	2.21	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	385/398 (97%)	353 (92%)	26 (7%)	6 (2%)	9	17
1	B	385/398 (97%)	355 (92%)	25 (6%)	5 (1%)	12	21
All	All	770/796 (97%)	708 (92%)	51 (7%)	11 (1%)	11	20

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	197	ASP
1	B	197	ASP
1	A	258	SER
1	B	258	SER
1	B	349	GLN
1	A	349	GLN
1	A	373	GLN
1	B	373	GLN
1	A	367	MET

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Mol	Chain	Res	Type
1	B	367	MET
1	A	372	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	320/328 (98%)	293 (92%)	27 (8%)	11	21
1	B	320/328 (98%)	295 (92%)	25 (8%)	12	24
All	All	640/656 (98%)	588 (92%)	52 (8%)	11	23

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	40	THR
1	A	58	ASP
1	A	64	LEU
1	A	70	GLN
1	A	91	LEU
1	A	92	GLU
1	A	162	GLN
1	A	168	LEU
1	A	170	TYR
1	A	174	GLU
1	A	182	LEU
1	A	197	ASP
1	A	216	ARG
1	A	217	LYS
1	A	219	SER
1	A	249	LEU
1	A	258	SER
1	A	271	LEU
1	A	297	LEU
1	A	318	LEU

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Mol	Chain	Res	Type
1	A	335	LEU
1	A	346	PHE
1	A	354	THR
1	A	361	LEU
1	A	369	MET
1	A	370	ARG
1	B	4	LEU
1	B	64	LEU
1	B	91	LEU
1	B	92	GLU
1	B	162	GLN
1	B	168	LEU
1	B	170	TYR
1	B	174	GLU
1	B	197	ASP
1	B	216	ARG
1	B	217	LYS
1	B	219	SER
1	B	238	LEU
1	B	249	LEU
1	B	258	SER
1	B	271	LEU
1	B	297	LEU
1	B	314	ARG
1	B	318	LEU
1	B	320	LEU
1	B	335	LEU
1	B	354	THR
1	B	361	LEU
1	B	369	MET
1	B	370	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	70	GLN
1	A	71	GLN
1	A	72	GLN
1	A	124	ASN
1	A	141	GLN
1	A	162	GLN

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Mol	Chain	Res	Type
1	A	165	GLN
1	A	175	GLN
1	A	176	ASN
1	A	251	HIS
1	A	257	HIS
1	A	263	GLN
1	A	270	GLN
1	A	329	GLN
1	A	350	GLN
1	B	3	GLN
1	B	70	GLN
1	B	71	GLN
1	B	72	GLN
1	B	124	ASN
1	B	141	GLN
1	B	145	GLN
1	B	162	GLN
1	B	165	GLN
1	B	175	GLN
1	B	176	ASN
1	B	251	HIS
1	B	257	HIS
1	B	263	GLN
1	B	270	GLN
1	B	329	GLN
1	B	350	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 2 ligands modelled in this entry, 2 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.