



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:31 PM GMT

PDB ID : 4FBQ
Title : Crystal structure of a covalently fused Nbs1-Mre11 complex with two manganese ions per active site
Authors : Schiller, C.B.; Lammens, K.; Hopfner, K.P.
Deposited on : 2012-05-23
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

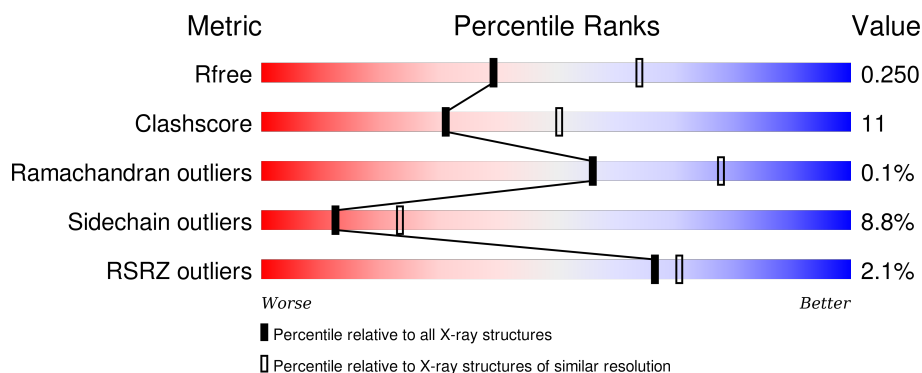
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	472	<div> <div>3%</div> <div>63%</div> <div>23%</div> <div>•</div> <div>11%</div> </div>
1	B	472	<div> <div>%</div> <div>64%</div> <div>19%</div> <div>•</div> <div>15%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6776 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 DNA repair and telomere maintenance protein nbs1, DNA repair protein rad32 CHIMERIC PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	418	Total	C	N	O	S	87	0	0
			3387	2157	591	631	8			
1	B	402	Total	C	N	O	S	70	0	0
			3240	2065	561	606	8			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP O43070
A	-3	PRO	-	EXPRESSION TAG	UNP O43070
A	-2	LEU	-	EXPRESSION TAG	UNP O43070
A	-1	GLY	-	EXPRESSION TAG	UNP O43070
A	0	SER	-	EXPRESSION TAG	UNP O43070
A	532	VAL	-	LINKER	UNP Q09683
A	533	ASP	-	LINKER	UNP Q09683
A	534	GLY	-	LINKER	UNP Q09683
A	535	SER	-	LINKER	UNP Q09683
A	536	ALA	-	LINKER	UNP Q09683
A	537	GLY	-	LINKER	UNP Q09683
A	538	SER	-	LINKER	UNP Q09683
A	539	ALA	-	LINKER	UNP Q09683
A	540	GLY	-	LINKER	UNP Q09683
A	541	SER	-	LINKER	UNP Q09683
B	-4	GLY	-	EXPRESSION TAG	UNP O43070
B	-3	PRO	-	EXPRESSION TAG	UNP O43070
B	-2	LEU	-	EXPRESSION TAG	UNP O43070
B	-1	GLY	-	EXPRESSION TAG	UNP O43070
B	0	SER	-	EXPRESSION TAG	UNP O43070
B	532	VAL	-	LINKER	UNP Q09683
B	533	ASP	-	LINKER	UNP Q09683
B	534	GLY	-	LINKER	UNP Q09683
B	535	SER	-	LINKER	UNP Q09683

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Chain	Residue	Modelled	Actual	Comment	Reference
B	536	ALA	-	LINKER	UNP Q09683
B	537	GLY	-	LINKER	UNP Q09683
B	538	SER	-	LINKER	UNP Q09683
B	539	ALA	-	LINKER	UNP Q09683
B	540	GLY	-	LINKER	UNP Q09683
B	541	SER	-	LINKER	UNP Q09683

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

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	65	Total O 65 65	0	0
3	B	80	Total O 80 80	0	0

Q1408
F1409
Y1410
L1411
K1412
LYS

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.26 Å 79.03 Å 222.97 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.41 – 2.50 47.41 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.3 (47.41-2.50) 98.3 (47.41-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.96 (at 2.51 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.213 , 0.257 0.200 , 0.250	Depositor DCC
R_{free} test set	1992 reflections (5.45%)	DCC
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 36564 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6776	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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	0.47	0/3460	0.60	1/4685 (0.0%)
1	B	0.46	0/3313	0.56	0/4497
All	All	0.46	0/6773	0.58	1/9182 (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	478	ASP	CB-CG-OD2	5.16	122.95	118.30

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	3387	0	3358	74	0
1	B	3240	0	3200	66	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	65	0	0	1	0
3	B	80	0	0	1	0
All	All	6776	0	6558	138	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 11.

All (138) 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:1209:ASP:O	1:A:1212:ARG:HG3	1.39	1.19
1:B:1292:LEU:HD21	1:B:1294:ILE:HD11	1.16	1.12
1:B:1292:LEU:CD2	1:B:1294:ILE:HD11	1.88	1.03
1:A:1323:SER:HA	1:A:1325:ILE:H	1.26	0.99
1:B:1164:ASN:HB3	1:B:1193:ARG:HG2	1.52	0.89
1:B:1164:ASN:HB3	1:B:1193:ARG:CG	2.09	0.82
1:B:1016:ASN:O	1:B:1179:THR:HG21	1.79	0.82
1:B:1328:MET:HA	1:B:1383:GLY:HA3	1.63	0.78
1:B:1209:ASP:O	1:B:1212:ARG:HG3	1.83	0.77
1:A:1209:ASP:O	1:A:1212:ARG:CG	2.30	0.73
1:A:1323:SER:HA	1:A:1325:ILE:N	2.02	0.72
1:A:1163:GLU:O	1:A:1164:ASN:ND2	2.22	0.72
1:B:1016:ASN:O	1:B:1179:THR:CG2	2.38	0.71
1:B:1015:GLU:HB3	3:B:1642:HOH:O	1.91	0.69
1:B:1354:TRP:O	1:B:1358:GLN:HG2	1.93	0.69
1:A:1323:SER:OG	1:A:1324:SER:HA	1.93	0.68
1:A:1148:ILE:O	1:A:1152:THR:HG23	1.95	0.67
1:B:495:TYR:CE1	1:B:1212:ARG:HD3	2.31	0.65
1:A:1392:PHE:O	1:A:1395:ARG:HG3	1.97	0.65
1:A:1026:PRO:HG2	1:A:1066:ILE:HA	1.78	0.65
1:A:1164:ASN:O	1:A:1166:ASN:N	2.32	0.63
1:A:1209:ASP:OD1	1:A:1210:LEU:N	2.30	0.63
1:B:487:LEU:O	1:B:1237:SER:HB2	1.99	0.62
1:A:1344:GLU:O	1:A:1348:THR:HG23	2.00	0.62
1:A:1099:LEU:HD21	1:A:1175:GLN:HB2	1.81	0.62
1:B:1180:LYS:HB3	1:B:1215:TRP:CE3	2.35	0.62
1:B:1292:LEU:HD21	1:B:1294:ILE:CD1	2.10	0.62
1:A:1029:GLY:O	1:A:1032:GLU:HG3	1.99	0.62
1:A:1099:LEU:CD2	1:A:1175:GLN:HB2	2.31	0.60
1:A:1213:ASP:O	1:A:1297:LYS:NZ	2.30	0.60
1:B:495:TYR:CD1	1:B:1212:ARG:NE	2.70	0.59
1:B:1216:PHE:CE2	1:B:1294:ILE:HD12	2.38	0.59
1:B:1171:PRO:HB3	1:B:1184:TYR:CZ	2.37	0.59
1:B:489:ASN:ND2	1:B:1241:ASP:OD2	2.35	0.59
1:B:1164:ASN:CB	1:B:1193:ARG:CG	2.81	0.58
1:B:1210:LEU:HB3	1:B:1211:TYR:CD2	2.38	0.58
1:B:1171:PRO:HB3	1:B:1184:TYR:CE2	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:496:ILE:HD13	1:A:1206:LEU:HD13	1.85	0.57
1:A:1314:MET:CE	1:A:1375:ARG:HD3	2.34	0.57
1:B:1222:HIS:CD2	1:B:1250:HIS:HB2	2.39	0.57
1:A:521:PHE:CE2	1:A:1097:GLU:HB2	2.40	0.57
1:B:1368:LYS:HB2	1:B:1369:PRO:HD2	1.86	0.56
1:B:1131:HIS:HD2	1:B:1157:TYR:OH	1.88	0.56
1:A:1163:GLU:HG3	1:A:1166:ASN:N	2.20	0.56
1:A:1063:GLY:O	1:A:1222:HIS:HD2	1.89	0.55
1:A:1364:VAL:HG12	1:A:1365:GLU:OE2	2.07	0.55
1:A:1354:TRP:O	1:A:1358:GLN:HG2	2.07	0.54
1:A:1169:VAL:HG12	1:A:1184:TYR:HD1	1.72	0.54
1:A:1316:ASP:OD1	1:A:1377:ARG:HB2	2.07	0.54
1:A:1148:ILE:HD13	1:B:1077:LEU:HD23	1.89	0.53
1:B:1131:HIS:CD2	1:B:1187:SER:HB3	2.43	0.53
1:A:1171:PRO:HB3	1:A:1184:TYR:CZ	2.44	0.53
1:B:1210:LEU:HB3	1:B:1211:TYR:HD2	1.73	0.53
1:A:1387:GLU:OE2	1:A:1395:ARG:NH1	2.41	0.52
1:A:1375:ARG:HG3	1:A:1375:ARG:HH11	1.74	0.52
1:A:1229:THR:HG22	1:A:1230:PRO:HD2	1.91	0.52
1:A:1097:GLU:HG2	1:A:1099:LEU:HD13	1.93	0.51
1:A:1379:ASP:OD1	1:A:1381:THR:OG1	2.24	0.51
1:A:1326:PRO:O	1:A:1329:VAL:HG12	2.11	0.51
1:A:1131:HIS:CD2	1:A:1187:SER:HB3	2.45	0.51
1:B:1164:ASN:CB	1:B:1193:ARG:HG2	2.33	0.50
1:B:1031:GLY:O	1:B:1038:GLY:HA2	2.10	0.50
1:A:1209:ASP:C	1:A:1212:ARG:HG3	2.26	0.50
1:B:1180:LYS:HB3	1:B:1215:TRP:CZ3	2.47	0.50
1:A:1375:ARG:NH2	1:A:1408:GLN:HB2	2.27	0.50
1:A:520:ASN:OD1	1:A:522:LYS:HG2	2.12	0.50
1:A:1031:GLY:O	1:A:1038:GLY:HA2	2.11	0.50
1:A:1180:LYS:HB3	1:A:1215:TRP:CE3	2.47	0.49
1:A:1133:ASN:ND2	1:A:1134:HIS:HD2	2.10	0.49
1:A:1336:LEU:HD11	1:A:1387:GLU:HB2	1.94	0.49
1:B:1032:GLU:OE2	1:B:1073:SER:OG	2.23	0.49
1:A:1235:PRO:HG2	1:A:1238:PHE:CZ	2.48	0.49
1:A:1135:ASP:O	1:A:1136:ASP:C	2.51	0.49
1:B:1338:TYR:OH	1:B:1342:LYS:HE2	2.12	0.48
1:A:1375:ARG:NH1	1:A:1405:ASP:O	2.46	0.48
1:B:1375:ARG:NH2	1:B:1408:GLN:OE1	2.46	0.48
1:A:1135:ASP:O	1:A:1136:ASP:O	2.30	0.48
1:A:524:PHE:CD2	1:A:1122:ASN:ND2	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1094:CYS:SG	1:B:1096:LEU:HB2	2.54	0.47
1:A:1052:ALA:CB	1:A:1060:ILE:HD11	2.45	0.47
1:B:1196:HIS:O	1:B:1200:GLU:HG2	2.15	0.46
1:B:1332:LYS:O	1:B:1336:LEU:HD13	2.15	0.46
1:A:1284:THR:HG21	1:A:1314:MET:HB3	1.98	0.46
1:A:1303:LYS:HE3	3:A:1618:HOH:O	2.14	0.46
1:A:1261:ASN:O	1:A:1265:LYS:N	2.45	0.46
1:A:1312:PHE:CD1	1:A:1313:ILE:N	2.84	0.46
1:A:1092:LYS:HA	1:A:1093:PRO:HD2	1.80	0.46
1:A:1034:ASP:HA	1:A:1035:PRO:HD3	1.74	0.46
1:B:1026:PRO:HA	1:B:1274:THR:OG1	2.15	0.46
1:B:1238:PHE:CD1	1:B:1238:PHE:N	2.84	0.46
1:B:1019:ARG:NH1	1:B:1055:ARG:O	2.39	0.45
1:A:1180:LYS:HB3	1:A:1215:TRP:CZ3	2.52	0.45
1:B:1243:TYR:O	1:B:1266:PHE:HB3	2.17	0.45
1:B:1146:LEU:HB3	1:B:1157:TYR:CD2	2.52	0.45
1:A:1375:ARG:HG3	1:A:1375:ARG:NH1	2.31	0.45
1:A:1047:GLU:OE1	1:A:1308:THR:HG23	2.16	0.45
1:B:1267:THR:HG1	1:B:1299:PHE:HZ	1.64	0.45
1:B:1131:HIS:CD2	1:B:1157:TYR:OH	2.69	0.44
1:A:1234:LEU:HA	1:A:1235:PRO:HD3	1.81	0.44
1:A:1146:LEU:HD22	1:A:1146:LEU:HA	1.77	0.44
1:A:1208:PRO:HG2	1:A:1215:TRP:CD1	2.52	0.44
1:B:1034:ASP:HA	1:B:1035:PRO:HD3	1.78	0.44
1:A:1312:PHE:C	1:A:1312:PHE:CD1	2.91	0.43
1:A:1069:ASP:O	1:A:1072:PRO:HD3	2.18	0.43
1:B:1211:TYR:HB3	1:B:1214:GLU:HB2	2.00	0.43
1:B:1264:GLN:O	1:B:1265:LYS:HB2	2.18	0.43
1:A:1016:ASN:OD1	1:A:1177:GLY:HA3	2.18	0.43
1:B:1161:VAL:HG11	1:B:1189:VAL:CG2	2.48	0.43
1:A:525:GLN:HB2	1:A:525:GLN:HE21	1.64	0.43
1:B:1331:ASN:ND2	1:B:1383:GLY:O	2.49	0.43
1:A:1056:ASP:OD1	1:A:1092:LYS:HE2	2.18	0.42
1:B:1055:ARG:NH2	1:B:1302:GLU:OE1	2.53	0.42
1:A:521:PHE:CZ	1:A:1097:GLU:HB2	2.54	0.42
1:A:1131:HIS:HD2	1:A:1157:TYR:OH	2.01	0.42
1:A:1131:HIS:ND1	1:A:1131:HIS:C	2.73	0.42
1:B:1395:ARG:O	1:B:1399:ARG:NH2	2.53	0.42
1:B:1380:TYR:CD1	1:B:1409:PHE:HB3	2.55	0.42
1:B:1342:LYS:HD2	1:B:1342:LYS:HA	1.60	0.42
1:A:1287:LYS:HB2	1:A:1309:VAL:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1065:ASP:OD1	1:B:1134:HIS:HB2	2.20	0.41
1:B:1331:ASN:HD21	1:B:1384:TYR:HA	1.85	0.41
1:A:1071:LYS:HA	1:A:1072:PRO:HD2	1.87	0.41
1:B:1048:ILE:HA	1:B:1048:ILE:HD13	1.80	0.41
1:B:1328:MET:CA	1:B:1383:GLY:HA3	2.40	0.41
1:B:1017:THR:HA	1:B:1294:ILE:O	2.20	0.41
1:B:1026:PRO:HG2	1:B:1066:ILE:HA	2.03	0.41
1:A:1278:SER:O	1:A:1283:GLU:HG3	2.21	0.41
1:B:1328:MET:HG3	1:B:1328:MET:O	2.21	0.41
1:A:1026:PRO:HA	1:A:1274:THR:OG1	2.21	0.41
1:B:1299:PHE:CD2	1:B:1299:PHE:O	2.74	0.41
1:B:1217:ASN:ND2	1:B:1242:PHE:HD2	2.19	0.41
1:B:1178:PHE:CD2	1:B:1179:THR:HG22	2.55	0.40
1:B:1331:ASN:O	1:B:1334:GLU:N	2.51	0.40
1:A:1048:ILE:HD11	1:A:1306:LEU:HD21	2.03	0.40
1:B:1411:LEU:HA	1:B:1411:LEU:HD23	1.86	0.40
1:A:1151:VAL:HB	1:B:1078:TYR:CE1	2.57	0.40
1:A:1332:LYS:HE2	1:A:1386:THR:O	2.21	0.40
1:B:1238:PHE:HD1	1:B:1238:PHE:N	2.20	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	406/472 (86%)	386 (95%)	19 (5%)	1 (0%)	52	75
1	B	394/472 (84%)	376 (95%)	18 (5%)	0	100	100
All	All	800/944 (85%)	762 (95%)	37 (5%)	1 (0%)	56	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1162	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	379/420 (90%)	345 (91%)	34 (9%)	12	22
1	B	362/420 (86%)	331 (91%)	31 (9%)	13	24
All	All	741/840 (88%)	676 (91%)	65 (9%)	12	23

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	490	LEU
1	A	493	VAL
1	A	496	ILE
1	A	518	ARG
1	A	525	GLN
1	A	1041	SER
1	A	1099	LEU
1	A	1100	SER
1	A	1109	THR
1	A	1113	ASN
1	A	1122	ASN
1	A	1131	HIS
1	A	1142	ARG
1	A	1146	LEU
1	A	1160	ARG
1	A	1163	GLU
1	A	1209	ASP
1	A	1210	LEU
1	A	1213	ASP
1	A	1229	THR
1	A	1259	SER
1	A	1266	PHE
1	A	1267	THR

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Mol	Chain	Res	Type
1	A	1277	THR
1	A	1315	LYS
1	A	1319	LEU
1	A	1325	ILE
1	A	1348	THR
1	A	1349	GLU
1	A	1361	VAL
1	A	1365	GLU
1	A	1371	LEU
1	A	1381	THR
1	A	1397	VAL
1	B	482	GLU
1	B	496	ILE
1	B	1017	THR
1	B	1032	GLU
1	B	1099	LEU
1	B	1112	CYS
1	B	1122	ASN
1	B	1142	ARG
1	B	1146	LEU
1	B	1151	VAL
1	B	1157	TYR
1	B	1179	THR
1	B	1193	ARG
1	B	1209	ASP
1	B	1214	GLU
1	B	1266	PHE
1	B	1294	ILE
1	B	1295	THR
1	B	1297	LYS
1	B	1298	ASP
1	B	1315	LYS
1	B	1319	LEU
1	B	1335	VAL
1	B	1361	VAL
1	B	1365	GLU
1	B	1367	GLU
1	B	1368	LYS
1	B	1381	THR
1	B	1393	SER
1	B	1397	VAL
1	B	1404	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	525	GLN
1	A	1046	ASN
1	A	1120	ASN
1	A	1122	ASN
1	A	1164	ASN
1	A	1222	HIS
1	B	1120	ASN
1	B	1131	HIS
1	B	1300	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 [i](#)

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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	418/472 (88%)	0.06	13 (3%) 52 57	21, 43, 80, 109	26 (6%)
1	B	402/472 (85%)	-0.07	4 (0%) 84 86	22, 43, 80, 101	18 (4%)
All	All	820/944 (86%)	-0.01	17 (2%) 67 71	21, 43, 81, 109	44 (5%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	525	GLN	4.9
1	A	520	ASN	4.3
1	A	521	PHE	4.2
1	A	481	PHE	4.1
1	B	1141	GLY	3.9
1	A	1109	THR	3.8
1	A	518	ARG	3.7
1	A	1112	CYS	3.6
1	B	1164	ASN	3.4
1	A	523	LYS	3.1
1	A	519	LYS	3.0
1	A	477	ASP	2.9
1	A	480	ALA	2.9
1	A	486	ARG	2.8
1	B	481	PHE	2.4
1	B	495	TYR	2.0
1	A	1164	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	B	1501	1/1	0.98	0.15	1.43	36,36,36,36	0
2	MN	B	1502	1/1	0.98	0.15	1.21	42,42,42,42	0
2	MN	A	1501	1/1	0.98	0.12	-0.33	49,49,49,49	0
2	MN	A	1502	1/1	0.98	0.12	-0.58	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.