Full wwPDB X-ray Structure Validation Report ⓘ

Sep 12, 2023 – 04:02 PM EDT

PDB ID	:	4O4K
Title	:	DNA Double-Strand Break Repair Pathway Choice Is Directed by Distinct MRE11 Nuclease Activities
Authors	:	Shibata, A.; Moiani, D.; Arvai, A.S.; Perry, J.; Harding, S.M.; Genois, M.; Maity, R.; Rossum-Fikkert, S.; Kertokalio, A.; Romoli, F.; Ismail, A.; Ismalaj, E.; Petricci, E.; Neale, M.J.; Bristow, R.G.; Masson, J.; Wyman, C.; Jeggo, P.A.; Tainer, J.A.
Deposited on	:	2013-12-18
Resolution	:	2.10 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

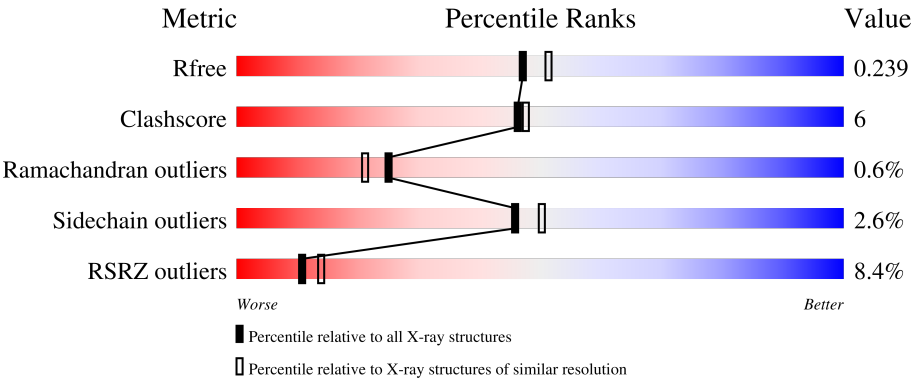
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	336	6% 84% 14%
1	B	336	10% 80% 15%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	2PK	A	403	-	-	-	X
3	2PK	B	403	-	-	X	X

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5318 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 Exonuclease, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	1	0
			2629	1692	449	482	6			
1	B	321	Total	C	N	O	S	0	0	0
			2524	1629	423	466	6			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9X1X0
A	-10	GLY	-	expression tag	UNP Q9X1X0
A	-9	SER	-	expression tag	UNP Q9X1X0
A	-8	ASP	-	expression tag	UNP Q9X1X0
A	-7	LYS	-	expression tag	UNP Q9X1X0
A	-6	ILE	-	expression tag	UNP Q9X1X0
A	-5	HIS	-	expression tag	UNP Q9X1X0
A	-4	HIS	-	expression tag	UNP Q9X1X0
A	-3	HIS	-	expression tag	UNP Q9X1X0
A	-2	HIS	-	expression tag	UNP Q9X1X0
A	-1	HIS	-	expression tag	UNP Q9X1X0
A	0	HIS	-	expression tag	UNP Q9X1X0
A	1	VAL	-	expression tag	UNP Q9X1X0
B	-11	MET	-	expression tag	UNP Q9X1X0
B	-10	GLY	-	expression tag	UNP Q9X1X0
B	-9	SER	-	expression tag	UNP Q9X1X0
B	-8	ASP	-	expression tag	UNP Q9X1X0
B	-7	LYS	-	expression tag	UNP Q9X1X0
B	-6	ILE	-	expression tag	UNP Q9X1X0
B	-5	HIS	-	expression tag	UNP Q9X1X0
B	-4	HIS	-	expression tag	UNP Q9X1X0
B	-3	HIS	-	expression tag	UNP Q9X1X0
B	-2	HIS	-	expression tag	UNP Q9X1X0
B	-1	HIS	-	expression tag	UNP Q9X1X0
B	0	HIS	-	expression tag	UNP Q9X1X0

Continued on next page...

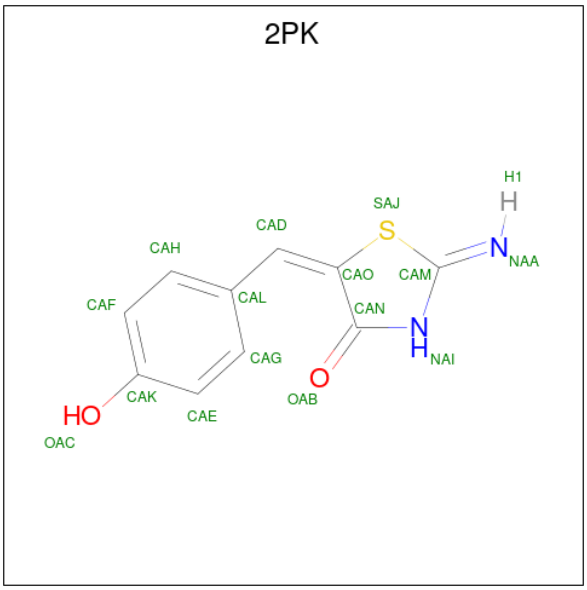
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	VAL	-	expression tag	UNP Q9X1X0

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

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

- Molecule 3 is (5 {E})-2-azanylidene-5-[(4-hydroxyphenyl)methylidene]-1,3-thiazolidin-4-one (three-letter code: 2PK) (formula: $C_{10}H_8N_2O_2S$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O S 15 10 2 2 1	0	0
3	B	1	Total C N O S 15 10 2 2 1	0	0

- Molecule 4 is water.

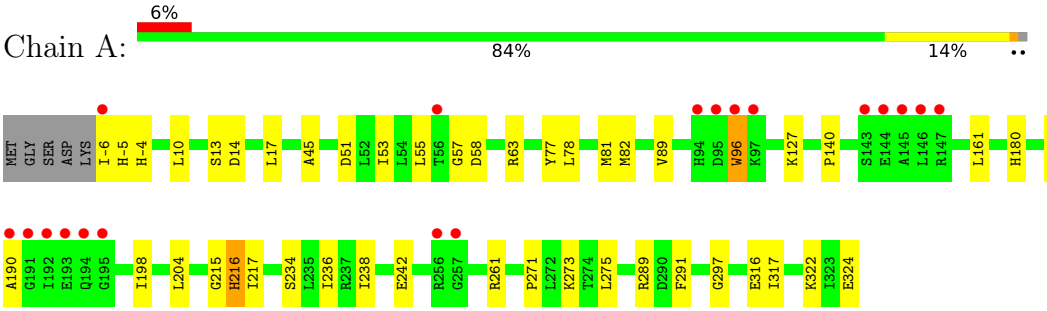
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	76	Total O 76 76	0	0
4	B	55	Total O 55 55	0	0

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 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.

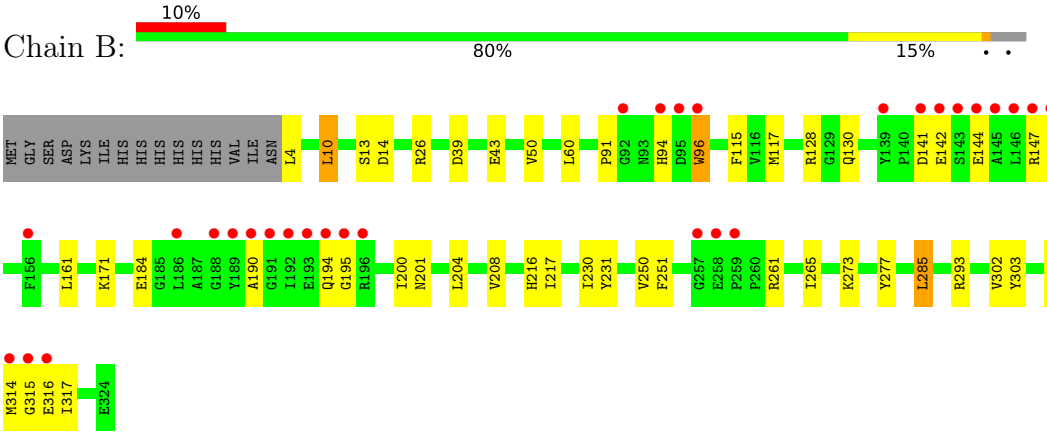
• Molecule 1: Exonuclease, putative

Chain A:



• Molecule 1: Exonuclease, putative

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	47.77Å 113.15Å 80.89Å 90.00° 101.47° 90.00°	Depositor
Resolution (Å)	37.41 – 2.10 37.41 – 2.10	Depositor EDS
% Data completeness (in resolution range)	88.7 (37.41-2.10) 88.6 (37.41-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.45 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.193 , 0.240 0.195 , 0.239	Depositor DCC
R_{free} test set	2000 reflections (4.32%)	wwPDB-VP
Wilson B-factor (Å^2)	40.8	Xtriage
Anisotropy	0.594	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 49.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5318	wwPDB-VP
Average B, all atoms (Å^2)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.94% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, 2PK

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	0.41	0/2694	0.57	0/3650
1	B	0.37	0/2580	0.58	0/3495
All	All	0.39	0/5274	0.57	0/7145

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2629	0	2593	31	0
1	B	2524	0	2493	32	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	15	0	9	3	0
3	B	15	0	9	7	0
4	A	76	0	0	5	0
4	B	55	0	0	2	0
All	All	5318	0	5104	65	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 6.

All (65) 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)
3:A:403:2PK:H9	3:A:403:2PK:OAB	1.78	0.84
3:B:403:2PK:H9	3:B:403:2PK:OAB	1.79	0.83
1:B:26:ARG:NH2	1:B:314:MET:O	2.15	0.79
1:A:242:GLU:OE1	4:A:526:HOH:O	2.05	0.73
1:B:285:LEU:HD12	1:B:313:LEU:HD21	1.70	0.71
1:A:236:ILE:O	4:A:526:HOH:O	2.09	0.68
1:B:316:GLU:HG3	1:B:317:ILE:HG13	1.76	0.66
1:B:10:LEU:HB2	1:B:50:VAL:HG11	1.77	0.65
1:A:234:SER:OG	4:A:526:HOH:O	2.13	0.64
1:B:194:GLN:N	1:B:195:GLY:HA2	2.15	0.61
1:B:96:TRP:H	1:B:96:TRP:HD1	1.48	0.61
1:A:57:GLY:HA2	1:A:180:HIS:CD2	2.35	0.61
1:B:147:ARG:N	1:B:148:LYS:HA	2.16	0.60
1:B:60:LEU:HB2	3:B:403:2PK:H3	1.82	0.60
1:B:60:LEU:H	3:B:403:2PK:H3	1.66	0.60
1:A:140:PRO:HD2	1:A:198:ILE:HD13	1.84	0.59
1:A:77:TYR:O	1:A:81:MET:HG3	2.03	0.58
1:A:127:LYS:NZ	4:A:511:HOH:O	2.35	0.58
1:A:316:GLU:HG3	1:A:317:ILE:HG13	1.86	0.57
1:B:91:PRO:HA	3:B:403:2PK:H8	1.85	0.57
1:B:13:SER:OG	1:B:14:ASP:OD2	2.22	0.57
1:B:141:ASP:HA	1:B:142:GLU:HB2	1.85	0.57
1:B:293:ARG:NE	4:B:532:HOH:O	2.39	0.55
1:B:115:PHE:HB3	1:B:117:MET:SD	2.47	0.54
1:B:190:ALA:HB1	4:B:524:HOH:O	2.09	0.52
1:B:94:HIS:CE1	3:B:403:2PK:H4	2.45	0.51
1:A:-5:HIS:O	1:A:-4:HIS:ND1	2.44	0.51
1:B:201:ASN:HB3	1:B:204:LEU:HD23	1.93	0.50
1:B:141:ASP:HB3	1:B:142:GLU:C	2.32	0.49
1:B:60:LEU:HB2	3:B:403:2PK:CAD	2.42	0.48
1:B:161:LEU:HD13	1:B:204:LEU:HB3	1.95	0.48
1:A:140:PRO:HD2	1:A:198:ILE:CD1	2.44	0.47
1:B:251:PHE:HB2	1:B:265:ILE:HD11	1.97	0.47
1:A:13:SER:OG	1:A:14:ASP:OD2	2.32	0.46
1:A:51:ASP:HA	1:A:127:LYS:HB3	1.97	0.46
1:A:186:LEU:HB2	1:A:190:ALA:HB2	1.97	0.46
1:B:311:PRO:O	1:B:313:LEU:N	2.48	0.46
1:A:273:LYS:HG2	1:A:291:PHE:CZ	2.51	0.46
1:A:53:ILE:HD13	1:A:81:MET:HB3	1.97	0.46
1:B:184:GLU:N	1:B:200:ILE:O	2.39	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ARG:HA	3:A:403:2PK:SAJ	2.57	0.45
1:A:55:LEU:HD12	1:A:89:VAL:HG22	1.98	0.45
1:A:58:ASP:HA	3:A:403:2PK:OAB	2.17	0.45
1:A:10:LEU:HD22	1:A:45:ALA:HB2	1.99	0.45
1:B:96:TRP:CD1	1:B:96:TRP:N	2.83	0.45
1:B:277:TYR:O	1:B:303:TYR:HA	2.17	0.44
1:A:-6:ILE:HB	1:A:-5:HIS:H	1.56	0.44
1:B:142:GLU:H	1:B:144:GLU:N	2.16	0.43
1:B:312:ASP:CG	1:B:315:GLY:H	2.22	0.43
1:A:289:ARG:HE	1:A:316:GLU:CD	2.22	0.43
1:A:78:LEU:O	1:A:82:MET:HG3	2.18	0.43
1:B:94:HIS:HE1	3:B:403:2PK:H4	1.81	0.42
1:B:230:ILE:HG21	1:B:250:VAL:HG21	2.02	0.42
1:B:128:ARG:HB2	1:B:130:GLN:HE22	1.85	0.42
1:A:187:ALA:HA	1:A:188:GLY:HA2	1.60	0.41
1:A:236:ILE:HG13	1:A:238:ILE:HD13	2.02	0.41
1:A:96:TRP:CD1	1:A:96:TRP:N	2.87	0.41
1:A:238:ILE:HG12	4:A:526:HOH:O	2.19	0.41
1:A:215:GLY:O	1:A:216:HIS:HB3	2.20	0.41
1:B:39:ASP:O	1:B:43:GLU:HG3	2.20	0.41
1:A:57:GLY:HA2	1:A:180:HIS:HD2	1.81	0.41
1:A:161:LEU:HD13	1:A:204:LEU:HB3	2.03	0.41
1:B:217:ILE:HG12	1:B:231:TYR:CE1	2.55	0.41
1:A:271:PRO:O	1:A:297:GLY:HA3	2.21	0.41
1:A:322:LYS:HG2	1:A:324:GLU:H	1.86	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	330/336 (98%)	313 (95%)	16 (5%)	1 (0%)	41 41

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	319/336 (95%)	299 (94%)	17 (5%)	3 (1%)	17 12
All	All	649/672 (97%)	612 (94%)	33 (5%)	4 (1%)	25 21

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	216	HIS
1	B	216	HIS
1	B	312	ASP
1	B	208	VAL

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	279/298 (94%)	274 (98%)	5 (2%)	59 65
1	B	265/298 (89%)	256 (97%)	9 (3%)	37 39
All	All	544/596 (91%)	530 (97%)	14 (3%)	46 50

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LEU
1	A	96	TRP
1	A	217	ILE
1	A	261	ARG
1	A	275	LEU
1	B	4	LEU
1	B	10	LEU
1	B	96	TRP
1	B	171	LYS
1	B	261	ARG
1	B	273	LYS
1	B	285	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	302	VAL
1	B	310	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	HIS
1	B	180	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	2PK	A	403	-	16,16,16	3.08	4 (25%)	18,22,22	5.29	6 (33%)
3	2PK	B	403	-	16,16,16	3.52	4 (25%)	18,22,22	4.94	6 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	2PK	A	403	-	-	2/4/16/16	0/2/2/2
3	2PK	B	403	-	-	2/4/16/16	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	2PK	CAM-SAJ	-10.47	1.61	1.74
3	B	403	2PK	CAN-CAO	-7.98	1.35	1.48
3	A	403	2PK	CAN-CAO	-7.80	1.35	1.48
3	A	403	2PK	CAM-SAJ	-7.56	1.64	1.74
3	A	403	2PK	CAO-SAJ	-4.17	1.65	1.73
3	B	403	2PK	CAO-SAJ	-3.11	1.67	1.73
3	B	403	2PK	CAL-CAD	-2.89	1.41	1.46
3	A	403	2PK	CAL-CAD	-2.83	1.41	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403	2PK	CAD-CAO-CAN	18.00	134.72	120.47
3	B	403	2PK	CAD-CAO-CAN	17.04	133.95	120.47
3	A	403	2PK	CAD-CAO-SAJ	-7.78	119.26	129.22
3	B	403	2PK	CAD-CAO-SAJ	-7.23	119.97	129.22
3	A	403	2PK	CAO-CAN-NAI	7.01	116.06	110.22
3	B	403	2PK	CAO-CAN-NAI	5.70	114.97	110.22
3	A	403	2PK	CAN-CAO-SAJ	-5.22	105.94	109.84
3	B	403	2PK	CAN-CAO-SAJ	-5.08	106.05	109.84
3	A	403	2PK	CAM-SAJ-CAO	4.84	96.28	89.50
3	B	403	2PK	CAM-SAJ-CAO	4.71	96.11	89.50
3	A	403	2PK	OAB-CAN-NAI	-3.01	118.08	125.08
3	B	403	2PK	OAB-CAN-NAI	-2.44	119.40	125.08

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	2PK	CAL-CAD-CAO-CAN
3	B	403	2PK	CAL-CAD-CAO-CAN
3	A	403	2PK	CAL-CAD-CAO-SAJ

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	B	403	2PK	CAL-CAD-CAO-SAJ

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	2PK	3	0
3	B	403	2PK	7	0

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

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, 95^{th} 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	331/336 (98%)	0.35	21 (6%) 20 24	30, 48, 134, 201	0
1	B	321/336 (95%)	0.65	34 (10%) 6 7	31, 58, 152, 215	0
All	All	652/672 (97%)	0.50	55 (8%) 11 14	30, 52, 141, 215	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	190	ALA	18.6
1	B	189	TYR	16.6
1	A	192	ILE	11.8
1	B	191	GLY	11.2
1	A	143	SER	10.8
1	B	192	ILE	10.4
1	A	190	ALA	9.8
1	B	146	LEU	9.8
1	B	145	ALA	9.8
1	A	191	GLY	9.4
1	A	195	GLY	8.8
1	B	257	GLY	6.9
1	A	189	TYR	6.3
1	A	-6	ILE	6.3
1	A	145	ALA	6.2
1	B	148	LYS	6.0
1	B	196	ARG	5.4
1	A	144	GLU	5.4
1	B	194	GLN	5.3
1	B	95	ASP	5.1
1	B	312	ASP	4.9
1	B	195	GLY	4.5
1	B	188	GLY	4.5
1	B	193	GLU	4.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	313	LEU	4.3
1	B	139	TYR	4.3
1	B	316	GLU	4.2
1	B	143	SER	4.2
1	B	314	MET	4.0
1	B	141	ASP	3.9
1	A	146	LEU	3.7
1	B	315	GLY	3.7
1	A	96	TRP	3.6
1	A	194	GLN	3.5
1	B	153	PHE	3.5
1	B	149	ASN	3.3
1	B	96	TRP	3.3
1	A	193	GLU	3.3
1	B	144	GLU	3.2
1	A	94	HIS	2.9
1	A	257	GLY	2.9
1	A	188	GLY	2.8
1	B	147	ARG	2.8
1	A	147	ARG	2.8
1	B	259	PRO	2.8
1	B	156	PHE	2.7
1	A	95	ASP	2.6
1	A	256	ARG	2.6
1	B	142	GLU	2.6
1	A	97	LYS	2.4
1	B	186	LEU	2.3
1	B	258	GLU	2.2
1	B	92	GLY	2.2
1	B	94	HIS	2.0
1	A	56	THR	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	B-factors(Å^2)	Q<0.9
3	2PK	A	403	15/15	0.56	0.69	36,58,83,196	15
3	2PK	B	403	15/15	0.74	0.55	35,52,74,226	15
2	MN	B	402	1/1	0.84	0.26	64,64,64,64	1
2	MN	A	402	1/1	0.89	0.13	52,52,52,52	1
2	MN	B	401	1/1	0.90	0.08	69,69,69,69	1
2	MN	A	401	1/1	0.95	0.04	63,63,63,63	1

6.5 Other polymers (i)

There are no such residues in this entry.