



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:50 PM UTC

PDB ID : 4OGE / pdb_00004oge
Title : Crystal structure of the Type II-C Cas9 enzyme from *Actinomyces naeslundii*
Authors : Jiang, F.; Ma, E.; Lin, S.; Doudna, J.A.
Deposited on : 2014-01-15
Resolution : 2.20 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

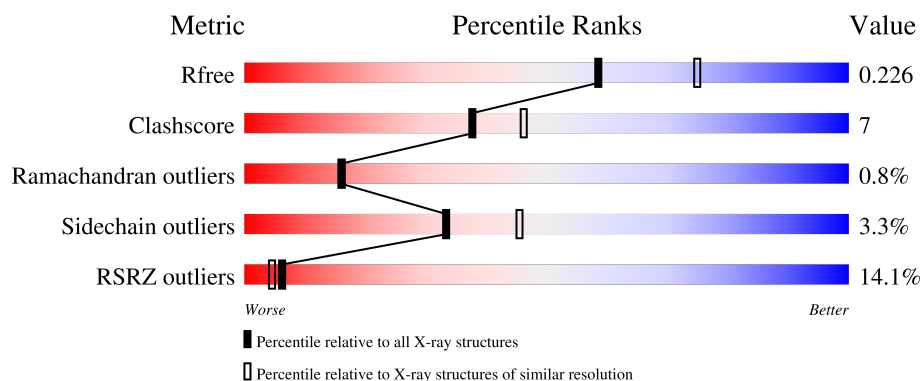
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	1101	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SPD	A	1205	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7799 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 HNH endonuclease domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	977	Total	C	N	O	S	0	0	0
			7432	4620	1392	1391	29			

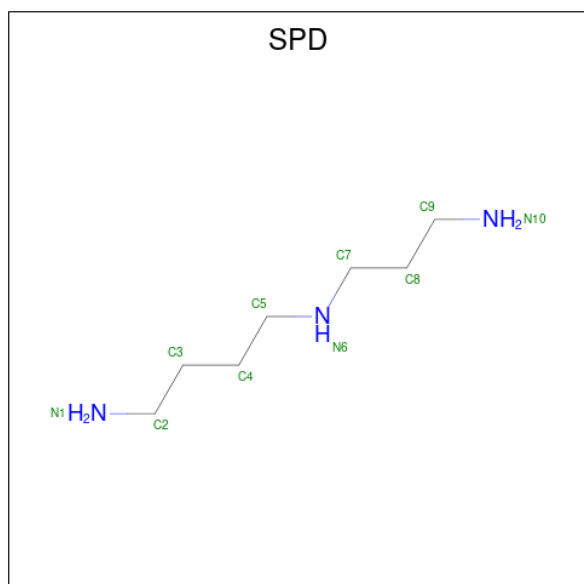
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Mg	0	0
			3	3		

- Molecule 4 is SPERMIDINE (CCD ID: SPD) (formula: C₇H₁₉N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			10	7	3		
4	A	1	Total	C	N	0	0
			10	7	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	343	Total	O	0	0
			343	343		

- Molecule 1: HNH endonuclease domain protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.42Å 133.03Å 80.69Å 90.00° 96.22° 90.00°	Depositor
Resolution (Å)	68.69 – 2.20 68.69 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.9 (68.69-2.20) 97.9 (68.69-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.186 , 0.226 0.187 , 0.226	Depositor DCC
R_{free} test set	3927 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7799	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: MG, ZN, SPD

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	0.50	0/7575	0.83	3/10287 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	507	VAL	N-CA-C	7.65	117.70	110.74
1	A	237	VAL	N-CA-C	-6.24	104.59	113.07
1	A	716	ASP	N-CA-C	-5.81	101.69	110.28

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	7432	0	7179	103	0
2	A	1	0	0	0	0
3	A	3	0	0	0	0
4	A	20	0	38	13	0
5	A	343	0	0	6	0
All	All	7799	0	7217	103	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 7.

All (103) 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:1031:THR:HG23	1:A:1033:SER:H	1.47	0.80
1:A:757:ARG:HB2	4:A:1205:SPD:H72	1.63	0.80
1:A:632:ILE:HD12	1:A:658:ILE:HG23	1.70	0.72
1:A:757:ARG:HH21	4:A:1205:SPD:C4	2.04	0.70
1:A:710:ARG:NH1	5:A:1584:HOH:O	2.23	0.70
1:A:1055:ASN:ND2	1:A:1059:GLU:OE2	2.24	0.70
1:A:857:CYS:O	1:A:861:ARG:HG2	1.92	0.70
1:A:539:GLN:HG3	1:A:548:ILE:HD11	1.72	0.70
1:A:757:ARG:HH21	4:A:1205:SPD:H41	1.57	0.70
1:A:319:ASP:OD2	1:A:652:ARG:NH2	2.26	0.68
1:A:1066:ALA:HB3	1:A:1069:VAL:HG13	1.76	0.67
1:A:378:TRP:O	1:A:386:ARG:NH1	2.29	0.66
1:A:71:ARG:HB3	1:A:397:PRO:HG3	1.76	0.66
1:A:1025:ARG:NH2	5:A:1357:HOH:O	2.29	0.64
1:A:660:ARG:NH2	1:A:667:ASP:OD2	2.29	0.64
1:A:382:ASP:OD1	1:A:382:ASP:N	2.31	0.64
1:A:640:LYS:HE3	1:A:647:SER:HB3	1.83	0.61
1:A:761:ARG:HD3	1:A:773:TRP:CZ2	2.36	0.61
1:A:500:GLU:HG3	1:A:501:VAL:HG13	1.83	0.60
1:A:925:SER:HA	4:A:1206:SPD:H72	1.83	0.60
1:A:952:LEU:O	1:A:955:ARG:HG3	2.01	0.60
1:A:95:PHE:HB3	1:A:143:ILE:HD11	1.84	0.59
1:A:757:ARG:CB	4:A:1205:SPD:H72	2.32	0.59
1:A:477:SER:O	1:A:480:ARG:HG2	2.01	0.59
1:A:301:ASN:ND2	5:A:1417:HOH:O	2.35	0.59
1:A:628:VAL:O	1:A:632:ILE:HG12	2.02	0.58
1:A:251:PRO:HG2	1:A:256:VAL:HG11	1.84	0.58
1:A:1025:ARG:NH1	5:A:1534:HOH:O	2.36	0.58
1:A:410:PRO:HA	1:A:413:ASP:HB2	1.85	0.58
1:A:754:LEU:O	4:A:1205:SPD:H81	2.04	0.57
1:A:242:LEU:O	1:A:244:ALA:N	2.38	0.56
1:A:352:ARG:NH2	1:A:683:ASN:OD1	2.36	0.56
1:A:507:VAL:CG1	1:A:511:PHE:HB2	2.36	0.56
1:A:97:ILE:HG13	1:A:98:PRO:HD3	1.88	0.55
1:A:708:ALA:HB1	4:A:1205:SPD:H82	1.87	0.55
1:A:242:LEU:O	1:A:245:VAL:N	2.36	0.54
1:A:638:TRP:HB2	1:A:654:LYS:HD3	1.89	0.54
1:A:421:HIS:NE2	1:A:512:THR:O	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:925:SER:HA	4:A:1206:SPD:C7	2.37	0.54
1:A:258:ARG:HE	1:A:260:ALA:HA	1.74	0.53
1:A:446:ASP:OD2	1:A:454:ARG:HD2	2.08	0.53
1:A:279:GLU:OE1	1:A:435:THR:HG22	2.09	0.53
1:A:381:ALA:HB3	1:A:386:ARG:HD2	1.89	0.52
1:A:699:MET:HG2	1:A:701:TYR:CZ	2.44	0.52
1:A:383:SER:HA	1:A:386:ARG:HD3	1.92	0.52
1:A:725:LYS:N	1:A:726:GLY:HA3	2.25	0.51
1:A:401:GLU:O	1:A:405:ILE:HG12	2.10	0.51
1:A:672:GLU:H	1:A:672:GLU:CD	2.18	0.51
1:A:757:ARG:HH21	4:A:1205:SPD:H42	1.75	0.50
1:A:849:ARG:NH2	1:A:904:ASP:OD2	2.40	0.50
1:A:674:SER:O	1:A:702:ARG:NH1	2.45	0.50
1:A:91:ARG:HA	1:A:95:PHE:O	2.13	0.49
1:A:666:GLU:OE2	5:A:1387:HOH:O	2.20	0.48
1:A:845:GLN:HB3	1:A:907:GLU:O	2.14	0.47
1:A:403:ALA:HB1	1:A:406:ILE:HB	1.95	0.47
1:A:604:ARG:HD2	1:A:604:ARG:C	2.39	0.47
1:A:257:SER:HA	1:A:258:ARG:HA	1.45	0.47
1:A:65:LEU:O	1:A:69:ALA:HB3	2.14	0.47
1:A:237:VAL:O	1:A:241:LEU:HG	2.15	0.47
1:A:507:VAL:HG12	1:A:511:PHE:HB2	1.95	0.47
1:A:509:ASP:N	1:A:510:GLY:HA3	2.30	0.47
1:A:389:MET:HE1	1:A:402:CYS:HB3	1.97	0.46
1:A:258:ARG:HB2	1:A:258:ARG:HH11	1.81	0.46
1:A:12:LEU:HB2	1:A:499:PRO:HA	1.97	0.46
1:A:794:MET:O	1:A:798:THR:HG23	2.16	0.46
1:A:1031:THR:HG22	1:A:1034:LYS:HB2	1.98	0.46
1:A:378:TRP:CZ2	1:A:386:ARG:HG2	2.51	0.45
1:A:1013:ARG:O	1:A:1016:GLU:HB2	2.16	0.45
1:A:371:ILE:HD11	1:A:416:LYS:O	2.16	0.45
1:A:293:ARG:HB2	1:A:341:ARG:HB2	1.98	0.45
1:A:715:ILE:HB	1:A:798:THR:HG21	1.98	0.45
1:A:258:ARG:HH21	1:A:261:PRO:HD3	1.81	0.45
1:A:880:ARG:NH1	1:A:885:GLU:HB2	2.31	0.45
1:A:89:VAL:O	1:A:93:LEU:HB2	2.16	0.45
1:A:926:ILE:O	4:A:1206:SPD:H52	2.17	0.45
1:A:156:VAL:HA	1:A:157:GLU:C	2.43	0.44
1:A:229:ALA:HA	1:A:230:ARG:HA	1.54	0.44
1:A:431:ARG:O	1:A:435:THR:HG23	2.18	0.43
1:A:758:SER:HB2	4:A:1205:SPD:H92	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:PHE:CD1	1:A:332:LYS:HE3	2.54	0.43
1:A:1057:ILE:O	1:A:1062:GLY:HA2	2.19	0.43
1:A:711:LYS:C	4:A:1205:SPD:H31	2.44	0.43
1:A:712:ALA:HB2	4:A:1205:SPD:H42	2.01	0.43
1:A:656:GLU:O	1:A:660:ARG:HG3	2.19	0.43
1:A:232:GLY:HA3	1:A:233:VAL:HA	1.59	0.43
1:A:416:LYS:O	1:A:419:SER:HB3	2.19	0.42
1:A:156:VAL:HG13	1:A:158:SER:OG	2.20	0.42
1:A:298:LYS:HA	1:A:299:GLY:HA2	1.64	0.42
1:A:392:TYR:OH	1:A:403:ALA:HB2	2.19	0.42
1:A:401:GLU:O	1:A:405:ILE:N	2.52	0.42
1:A:1079:ARG:HG3	1:A:1096:THR:O	2.19	0.42
1:A:483:LYS:O	1:A:487:ARG:HG2	2.19	0.42
1:A:419:SER:O	1:A:421:HIS:ND1	2.53	0.41
1:A:596:GLY:HA3	1:A:656:GLU:HG2	2.03	0.41
1:A:166:SER:HA	1:A:167:PRO:HD3	1.87	0.41
1:A:471:ALA:HA	1:A:472:PRO:HD3	1.87	0.41
1:A:820:LEU:HD22	1:A:948:THR:HG23	2.03	0.41
1:A:848:ASP:HB3	1:A:909:PRO:HG3	2.02	0.41
1:A:935:GLU:OE2	1:A:988:THR:OG1	2.28	0.41
1:A:271:ARG:HG2	5:A:1552:HOH:O	2.21	0.40
1:A:417:LEU:HD23	1:A:417:LEU:HA	1.74	0.40
1:A:389:MET:O	1:A:393:LEU:HG	2.21	0.40
1:A:994:VAL:O	1:A:997:ASP:HB2	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	965/1101 (88%)	914 (95%)	43 (4%)	8 (1%)	16 16

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	ARG
1	A	508	ARG
1	A	407	ALA
1	A	414	GLN
1	A	923	GLY
1	A	162	PRO
1	A	231	GLN
1	A	245	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	759/931 (82%)	734 (97%)	25 (3%)	33	45

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ARG
1	A	86	LEU
1	A	90	LEU
1	A	93	LEU
1	A	143	ILE
1	A	245	VAL
1	A	246	PHE
1	A	258	ARG
1	A	359	ILE
1	A	382	ASP
1	A	386	ARG
1	A	401	GLU
1	A	402	CYS
1	A	408	GLU
1	A	475	ASN
1	A	512	THR
1	A	601	VAL

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Mol	Chain	Res	Type
1	A	640	LYS
1	A	652	ARG
1	A	672	GLU
1	A	675	MET
1	A	791	ARG
1	A	867	GLU
1	A	1042	LEU
1	A	1069	VAL

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

Mol	Chain	Res	Type
1	A	267	GLN
1	A	321	GLN
1	A	336	HIS
1	A	368	GLN
1	A	503	HIS
1	A	531	ASN
1	A	586	GLN
1	A	687	HIS
1	A	814	GLN

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 oligosaccharides 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$
4	SPD	A	1206	-	9,9,9	0.44	0	8,8,8	0.72	0
4	SPD	A	1205	-	9,9,9	0.46	0	8,8,8	1.41	2 (25%)

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
4	SPD	A	1206	-	-	3/7/7/7	-
4	SPD	A	1205	-	-	5/7/7/7	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1205	SPD	C4-C5-N6	-2.40	105.62	112.07
4	A	1205	SPD	C7-C8-C9	-2.39	105.64	114.17

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1206	SPD	N6-C7-C8-C9
4	A	1205	SPD	C4-C5-N6-C7
4	A	1205	SPD	N6-C7-C8-C9
4	A	1206	SPD	C4-C5-N6-C7
4	A	1205	SPD	C2-C3-C4-C5
4	A	1206	SPD	C8-C7-N6-C5
4	A	1205	SPD	N1-C2-C3-C4
4	A	1205	SPD	C8-C7-N6-C5

There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1206	SPD	3	0
4	A	1205	SPD	10	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	977/1101 (88%)	0.34	138 (14%) 6 4	20, 42, 147, 199	0

All (138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	511	PHE	6.7
1	A	227	ILE	6.7
1	A	65	LEU	5.9
1	A	509	ASP	5.7
1	A	507	VAL	5.4
1	A	98	PRO	5.2
1	A	233	VAL	5.1
1	A	149	TRP	5.0
1	A	168	PHE	4.9
1	A	10	HIS	4.7
1	A	825	ALA	4.6
1	A	156	VAL	4.6
1	A	225	ARG	4.6
1	A	89	VAL	4.4
1	A	510	GLY	4.4
1	A	94	GLY	4.4
1	A	97	ILE	4.3
1	A	140	VAL	4.2
1	A	162	PRO	4.1
1	A	153	TYR	4.1
1	A	143	ILE	4.1
1	A	347	THR	4.1
1	A	237	VAL	4.0
1	A	409	LEU	4.0
1	A	160	LEU	4.0
1	A	138	MET	3.9
1	A	76	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	230	ARG	3.8
1	A	93	LEU	3.8
1	A	410	PRO	3.8
1	A	69	ALA	3.7
1	A	642	THR	3.7
1	A	241	LEU	3.7
1	A	144	ALA	3.7
1	A	159	LEU	3.6
1	A	139	ALA	3.5
1	A	90	LEU	3.5
1	A	95	PHE	3.5
1	A	518	ASP	3.5
1	A	163	ALA	3.4
1	A	822	ASP	3.4
1	A	152	PRO	3.4
1	A	512	THR	3.4
1	A	297	THR	3.3
1	A	827	THR	3.3
1	A	382	ASP	3.3
1	A	229	ALA	3.3
1	A	169	MET	3.3
1	A	157	GLU	3.3
1	A	246	PHE	3.3
1	A	402	CYS	3.3
1	A	81	THR	3.2
1	A	165	GLU	3.2
1	A	142	HIS	3.2
1	A	514	GLU	3.1
1	A	407	ALA	3.1
1	A	547	TYR	3.1
1	A	68	ILE	3.1
1	A	245	VAL	3.1
1	A	141	ARG	3.1
1	A	67	GLY	3.1
1	A	251	PRO	3.1
1	A	352	ARG	3.1
1	A	508	ARG	3.1
1	A	161	SER	3.0
1	A	255	ALA	3.0
1	A	419	SER	3.0
1	A	150	ARG	3.0
1	A	151	ASN	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	253	GLY	3.0
1	A	824	ASN	3.0
1	A	516	MET	3.0
1	A	299	GLY	3.0
1	A	155	LYS	3.0
1	A	239	LYS	2.9
1	A	167	PRO	2.9
1	A	145	ARG	2.9
1	A	148	GLY	2.9
1	A	823	GLY	2.9
1	A	640	LYS	2.9
1	A	74	ARG	2.9
1	A	244	ALA	2.9
1	A	298	LYS	2.8
1	A	228	CYS	2.8
1	A	826	HIS	2.8
1	A	137	SER	2.8
1	A	254	SER	2.8
1	A	147	ARG	2.8
1	A	243	ARG	2.8
1	A	158	SER	2.7
1	A	513	SER	2.6
1	A	238	CYS	2.6
1	A	48	ASP	2.6
1	A	146	HIS	2.6
1	A	75	LEU	2.6
1	A	523	ALA	2.6
1	A	1049	ASN	2.6
1	A	66	SER	2.6
1	A	829	ASN	2.5
1	A	415	ALA	2.5
1	A	416	LYS	2.5
1	A	72	ALA	2.5
1	A	96	PRO	2.5
1	A	73	ARG	2.4
1	A	517	ALA	2.4
1	A	727	ARG	2.4
1	A	257	SER	2.4
1	A	353	SER	2.4
1	A	418	ASP	2.4
1	A	234	SER	2.4
1	A	404	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	249	ASP	2.3
1	A	235	PRO	2.3
1	A	346	HIS	2.3
1	A	520	ARG	2.3
1	A	231	GLN	2.3
1	A	248	ALA	2.3
1	A	84	GLN	2.3
1	A	728	LYS	2.3
1	A	250	SER	2.3
1	A	400	SER	2.3
1	A	170	LYS	2.3
1	A	86	LEU	2.2
1	A	397	PRO	2.2
1	A	258	ARG	2.2
1	A	384	GLU	2.2
1	A	398	THR	2.2
1	A	828	VAL	2.2
1	A	768	GLY	2.2
1	A	271	ARG	2.1
1	A	939	PRO	2.1
1	A	525	ARG	2.1
1	A	226	LYS	2.1
1	A	78	HIS	2.1
1	A	515	ARG	2.1
1	A	354	ALA	2.0
1	A	403	ALA	2.0
1	A	240	GLN	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 oligosaccharides 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, 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	B-factors($\text{\AA}^2$)	Q<0.9
4	SPD	A	1206	10/10	0.80	0.21	41,70,89,90	0
4	SPD	A	1205	10/10	0.86	0.16	26,45,49,53	0
3	MG	A	1203	1/1	0.86	0.14	53,53,53,53	0
3	MG	A	1204	1/1	0.98	0.04	14,14,14,14	0
2	ZN	A	1201	1/1	0.99	0.02	28,28,28,28	0
3	MG	A	1202	1/1	0.99	0.02	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.