



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 10:11 AM EDT

PDB ID : 7YX7 / pdb_00007yx7
Title : Modified oligopeptidase B from *S. proteomaculans* in intermediate conformation with 1 spermine molecule at 1.72 Å resolution
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Deposited on : 2022-02-15
Resolution : 1.72 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

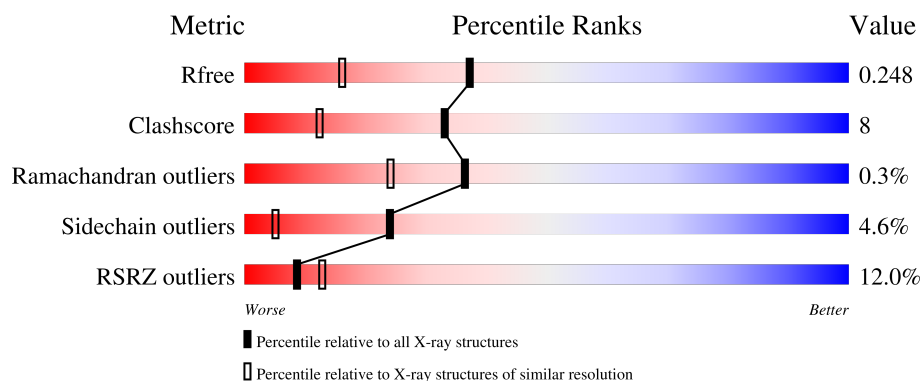
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	677	12% 86% 13%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6201 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.

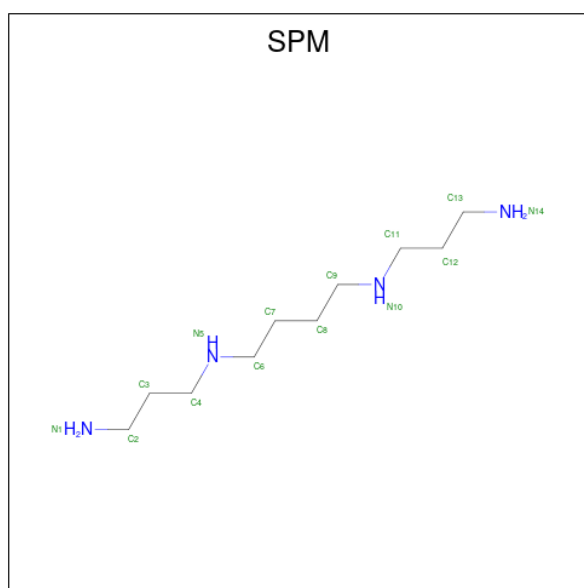
- Molecule 1 is a protein called Oligopeptidase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	677	5580	3542	939	1085	14	0	4	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	HIS	-	expression tag	UNP B3VI58
A	71	GLU	ILE	engineered mutation	UNP B3VI58
A	72	ASN	PRO	engineered mutation	UNP B3VI58
A	73	LEU	GLN	engineered mutation	UNP B3VI58
A	74	TYR	GLN	engineered mutation	UNP B3VI58
A	75	PHE	GLU	engineered mutation	UNP B3VI58
A	76	GLN	HIS	engineered mutation	UNP B3VI58

- Molecule 2 is SPERMINE (CCD ID: SPM) (formula: $C_{10}H_{26}N_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			14	10	4		

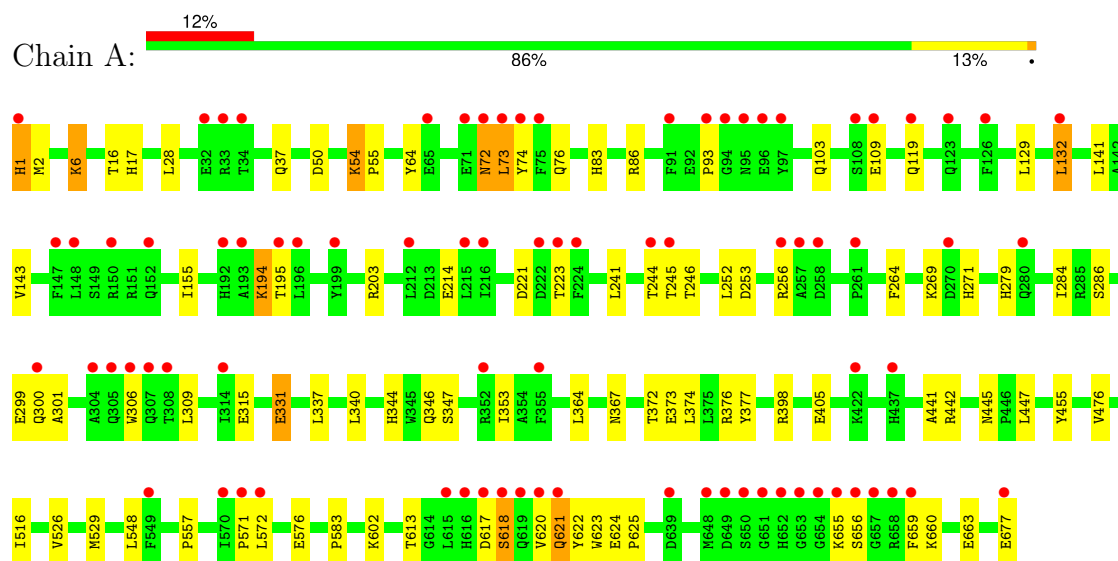
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	607	Total	O	0	0
			607	607		

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: Oligopeptidase B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.95Å 100.53Å 108.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.45 – 1.72 30.45 – 1.72	Depositor EDS
% Data completeness (in resolution range)	92.1 (30.45-1.72) 92.1 (30.45-1.72)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 1.72Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.199 , 0.239 0.210 , 0.248	Depositor DCC
R_{free} test set	3960 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	11.2	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6201	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.61% 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: SPM

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	1.05	2/5726 (0.0%)	1.20	0/7782

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	621	GLN	C-O	5.65	1.30	1.23
1	A	28	LEU	C-O	-5.18	1.17	1.24

There are no bond angle outliers.

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	5580	0	5288	85	0
2	A	14	0	26	7	0
3	A	607	0	0	9	1
All	All	6201	0	5314	85	1

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 8.

All (85) 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:132:LEU:CD2	1:A:141:LEU:HD11	1.60	1.28
1:A:221:ASP:OD1	1:A:223:THR:HG22	1.54	1.08
1:A:132:LEU:HD22	1:A:141:LEU:HD11	1.40	1.00
1:A:132:LEU:HD21	1:A:141:LEU:HD11	1.52	0.90
1:A:367:ASN:HD21	1:A:376:ARG:H	1.24	0.86
1:A:372[A]:THR:HG21	3:A:1361:HOH:O	1.80	0.81
1:A:398[B]:ARG:CG	1:A:398[B]:ARG:HH21	1.97	0.78
1:A:398[B]:ARG:HH21	1:A:398[B]:ARG:CB	1.96	0.78
1:A:441:ALA:H	1:A:445:ASN:HD21	1.35	0.75
1:A:64:TYR:HH	1:A:659:PHE:HD1	1.33	0.74
1:A:245:THR:HG23	1:A:246:THR:OG1	1.87	0.74
1:A:372[A]:THR:HG22	1:A:374:LEU:H	1.52	0.74
1:A:372[A]:THR:HG23	3:A:1126:HOH:O	1.87	0.73
1:A:398[B]:ARG:HH21	1:A:398[B]:ARG:HB2	1.53	0.72
1:A:223:THR:HG21	3:A:1364:HOH:O	1.89	0.71
1:A:244:THR:HB	3:A:1001:HOH:O	1.94	0.68
1:A:279:HIS:HE1	3:A:1167:HOH:O	1.76	0.67
1:A:86:ARG:HH11	1:A:103:GLN:HE21	1.42	0.66
1:A:271:HIS:HE1	1:A:286:SER:OG	1.82	0.62
1:A:331:GLU:HG2	1:A:364:LEU:CD1	2.30	0.61
1:A:398[B]:ARG:HH21	1:A:398[B]:ARG:HG3	1.66	0.60
1:A:64:TYR:OH	1:A:659:PHE:HD1	1.85	0.60
1:A:398[B]:ARG:HG3	1:A:398[B]:ARG:NH2	2.19	0.58
1:A:340:LEU:HB2	1:A:353:ILE:HG12	1.84	0.57
1:A:73:LEU:HD12	1:A:74:TYR:CE2	2.39	0.57
1:A:271:HIS:HD2	3:A:1228:HOH:O	1.86	0.57
1:A:455:TYR:OH	1:A:576:GLU:OE2	2.12	0.57
1:A:132:LEU:HD22	1:A:141:LEU:CD1	2.27	0.55
1:A:447:LEU:HD23	1:A:476:VAL:HB	1.86	0.55
1:A:344:HIS:HD2	1:A:346:GLN:H	1.56	0.53
1:A:617:ASP:OD1	1:A:618:SER:N	2.41	0.53
1:A:221:ASP:CG	1:A:223:THR:HG22	2.31	0.53
1:A:331:GLU:HG2	1:A:364:LEU:HD12	1.90	0.53
1:A:442:ARG:HH22	2:A:801:SPM:H112	1.73	0.53
1:A:340:LEU:HB2	1:A:353:ILE:CG1	2.40	0.52
1:A:398[B]:ARG:CG	1:A:398[B]:ARG:NH2	2.63	0.52
1:A:73:LEU:H	1:A:73:LEU:HD23	1.74	0.52
1:A:132:LEU:HD23	1:A:143:VAL:HG12	1.92	0.51
1:A:377:TYR:OH	1:A:398[A]:ARG:HD2	2.11	0.51
1:A:64:TYR:OH	1:A:659:PHE:CD1	2.60	0.51
1:A:132:LEU:CD2	1:A:141:LEU:CD1	2.57	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:HIS:CD2	1:A:346:GLN:H	2.28	0.50
1:A:17:HIS:HE1	1:A:583:PRO:O	1.95	0.49
1:A:516:ILE:HD11	1:A:526:VAL:HG21	1.94	0.49
1:A:613:THR:OG1	1:A:622:TYR:HA	2.12	0.49
1:A:442:ARG:NH2	2:A:801:SPM:H112	2.28	0.49
1:A:442:ARG:HH22	2:A:801:SPM:C11	2.25	0.48
1:A:245:THR:HG21	1:A:269:LYS:NZ	2.28	0.48
1:A:571:PRO:O	1:A:572:LEU:HB2	2.13	0.47
1:A:1:HIS:CD2	1:A:2:MET:HG3	2.50	0.47
1:A:442:ARG:HH12	2:A:801:SPM:H82	1.79	0.47
1:A:244:THR:CB	3:A:1001:HOH:O	2.60	0.46
1:A:344:HIS:CD2	1:A:347:SER:H	2.32	0.46
1:A:132:LEU:HD23	1:A:141:LEU:HD11	1.81	0.46
1:A:621:GLN:HB3	1:A:623:TRP:CZ3	2.51	0.46
1:A:132:LEU:CD2	1:A:143:VAL:HG12	2.46	0.46
1:A:73:LEU:HB2	1:A:93:PRO:HB3	1.97	0.46
1:A:86:ARG:HD2	1:A:103:GLN:HE21	1.79	0.45
1:A:557:PRO:O	1:A:620:VAL:HG11	2.15	0.45
1:A:143:VAL:O	1:A:155:ILE:HA	2.17	0.45
1:A:677:GLU:HA	1:A:677:GLU:OE1	2.17	0.45
1:A:83:HIS:HD2	3:A:1153:HOH:O	1.98	0.45
1:A:442:ARG:HH22	2:A:801:SPM:H81	1.83	0.44
1:A:624:GLU:HB2	1:A:625:PRO:HD3	2.00	0.44
1:A:54:LYS:N	1:A:55:PRO:CD	2.81	0.44
1:A:344:HIS:HE1	3:A:1078:HOH:O	2.01	0.43
1:A:622:TYR:C	1:A:622:TYR:CD1	2.97	0.43
1:A:659:PHE:CE2	1:A:663:GLU:OE1	2.72	0.43
1:A:617:ASP:CG	1:A:618:SER:N	2.77	0.42
1:A:1:HIS:HD2	1:A:2:MET:H	1.68	0.42
1:A:284:ILE:HG21	1:A:306:TRP:CH2	2.54	0.42
1:A:86:ARG:HH11	1:A:103:GLN:NE2	2.14	0.42
1:A:132:LEU:HD21	1:A:141:LEU:HD21	2.02	0.42
1:A:271:HIS:CE1	1:A:286:SER:OG	2.67	0.42
1:A:442:ARG:HH11	2:A:801:SPM:H22	1.84	0.42
1:A:72:ASN:ND2	1:A:405:GLU:CD	2.79	0.41
1:A:129:LEU:O	1:A:655:LYS:HE2	2.21	0.41
1:A:442:ARG:HH12	2:A:801:SPM:C8	2.33	0.41
1:A:344:HIS:HD2	1:A:347:SER:H	1.68	0.41
1:A:264:PHE:CE1	1:A:301:ALA:HA	2.56	0.40
1:A:1:HIS:CD2	1:A:2:MET:H	2.39	0.40
1:A:6:LYS:HB2	1:A:6:LYS:HE2	1.88	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ARG:HD2	1:A:214:GLU:OE1	2.21	0.40
1:A:529:MET:HE3	1:A:529:MET:HB2	1.91	0.40
1:A:253:ASP:OD2	1:A:256:ARG:NH1	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1413:HOH:O	3:A:1469:HOH:O[2_555]	2.19	0.01

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	679/677 (100%)	654 (96%)	23 (3%)	2 (0%)	36 24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	LYS
1	A	656	SER

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	595/591 (101%)	568 (96%)	27 (4%)	24 5

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	HIS
1	A	6	LYS
1	A	16	THR
1	A	37	GLN
1	A	50	ASP
1	A	54	LYS
1	A	72	ASN
1	A	73	LEU
1	A	76	GLN
1	A	109	GLU
1	A	119	GLN
1	A	132	LEU
1	A	194	LYS
1	A	195	THR
1	A	241	LEU
1	A	252	LEU
1	A	299	GLU
1	A	300	GLN
1	A	309	LEU
1	A	315	GLU
1	A	331	GLU
1	A	337	LEU
1	A	373	GLU
1	A	548	LEU
1	A	602	LYS
1	A	618	SER
1	A	660	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1	HIS
1	A	17	HIS
1	A	72	ASN
1	A	83	HIS
1	A	103	GLN
1	A	172	ASN

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Mol	Chain	Res	Type
1	A	181	ASN
1	A	192	HIS
1	A	220	GLN
1	A	262	GLN
1	A	271	HIS
1	A	279	HIS
1	A	300	GLN
1	A	339	GLN
1	A	344	HIS
1	A	367	ASN
1	A	445	ASN
1	A	502	GLN
1	A	518	GLN
1	A	544	GLN
1	A	585	GLN
1	A	616	HIS
1	A	652	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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
2	SPM	A	801	-	13,13,13	0.43	0	12,12,12	0.65	0

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
2	SPM	A	801	-	-	2/11/11/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	SPM	N10-C11-C12-C13
2	A	801	SPM	C8-C9-N10-C11

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	SPM	7	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	677/677 (100%)	0.76	81 (11%) 9 13	4, 13, 36, 79	4 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	654	GLY	8.9
1	A	257	ALA	6.5
1	A	657	GLY	6.2
1	A	620	VAL	5.8
1	A	650	SER	5.8
1	A	653	GLY	5.7
1	A	618	SER	5.6
1	A	74	TYR	5.5
1	A	93	PRO	5.1
1	A	652	HIS	5.1
1	A	1	HIS	5.0
1	A	651	GLY	4.7
1	A	193	ALA	4.6
1	A	617	ASP	4.3
1	A	306	TRP	4.3
1	A	73	LEU	4.3
1	A	655	LYS	4.2
1	A	616	HIS	4.2
1	A	659	PHE	4.2
1	A	656	SER	4.1
1	A	304	ALA	4.1
1	A	571	PRO	4.1
1	A	72	ASN	4.0
1	A	615	LEU	4.0
1	A	307	GLN	3.9
1	A	96	GLU	3.7
1	A	222	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	314	ILE	3.7
1	A	94	GLY	3.6
1	A	437	HIS	3.6
1	A	196	LEU	3.5
1	A	75	PHE	3.5
1	A	126	PHE	3.4
1	A	305	GLN	3.4
1	A	148	LEU	3.3
1	A	91	PHE	3.2
1	A	192	HIS	3.2
1	A	195	THR	3.2
1	A	572	LEU	3.1
1	A	215	LEU	3.0
1	A	549	PHE	3.0
1	A	199	TYR	2.9
1	A	570	ILE	2.9
1	A	108	SER	2.9
1	A	245	THR	2.8
1	A	422	LYS	2.8
1	A	355	PHE	2.8
1	A	639	ASP	2.8
1	A	658	ARG	2.8
1	A	619	GLN	2.8
1	A	132	LEU	2.8
1	A	150	ARG	2.7
1	A	256	ARG	2.7
1	A	95	ASN	2.7
1	A	109	GLU	2.7
1	A	97	TYR	2.6
1	A	223	THR	2.6
1	A	649	ASP	2.6
1	A	280	GLN	2.5
1	A	212	LEU	2.5
1	A	352	ARG	2.5
1	A	648	MET	2.5
1	A	32	GLU	2.4
1	A	244	THR	2.4
1	A	621	GLN	2.3
1	A	33	ARG	2.3
1	A	34	THR	2.3
1	A	224	PHE	2.3
1	A	261	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	258	ASP	2.2
1	A	270	ASP	2.2
1	A	65	GLU	2.1
1	A	152	GLN	2.1
1	A	308	THR	2.1
1	A	119	GLN	2.1
1	A	123	GLN	2.1
1	A	216	ILE	2.1
1	A	677	GLU	2.1
1	A	147	PHE	2.1
1	A	300	GLN	2.0
1	A	71[A]	GLU	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(Å ²)	Q<0.9
2	SPM	A	801	14/14	0.88	0.12	18,20,22,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.