



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4LEP / pdb_00004lep
Title : Structural insights into substrate recognition in proton dependent oligopeptide transporters
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Deposited on : 2013-06-26
Resolution : 3.20 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

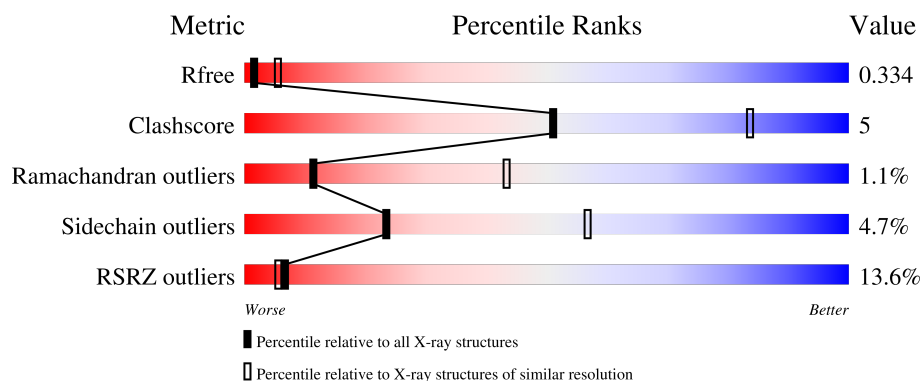
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	523	12% 71% 14% • 13%
1	B	523	11% 72% 13% • 14%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6858 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.

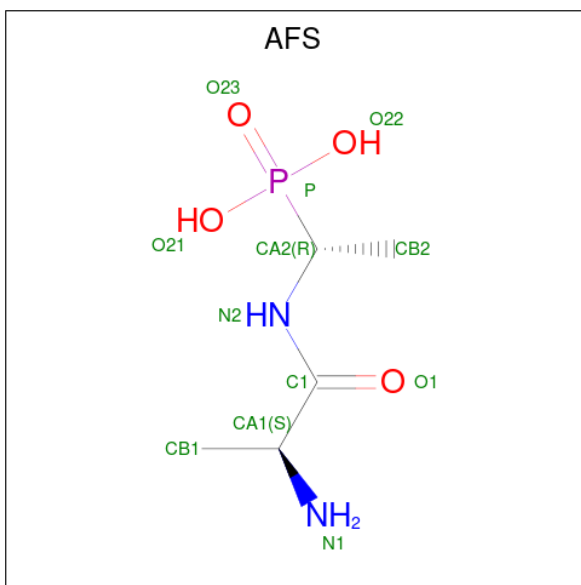
- Molecule 1 is a protein called Proton:oligopeptide symporter POT family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	0	0
			3434	2280	552	582	20			
1	B	452	Total	C	N	O	S	0	0	0
			3399	2255	543	581	20			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	517	ALA	-	expression tag	UNP Q8EHE6
A	518	GLU	-	expression tag	UNP Q8EHE6
A	519	ASN	-	expression tag	UNP Q8EHE6
A	520	LEU	-	expression tag	UNP Q8EHE6
A	521	TYR	-	expression tag	UNP Q8EHE6
A	522	PHE	-	expression tag	UNP Q8EHE6
A	523	GLN	-	expression tag	UNP Q8EHE6
B	517	ALA	-	expression tag	UNP Q8EHE6
B	518	GLU	-	expression tag	UNP Q8EHE6
B	519	ASN	-	expression tag	UNP Q8EHE6
B	520	LEU	-	expression tag	UNP Q8EHE6
B	521	TYR	-	expression tag	UNP Q8EHE6
B	522	PHE	-	expression tag	UNP Q8EHE6
B	523	GLN	-	expression tag	UNP Q8EHE6

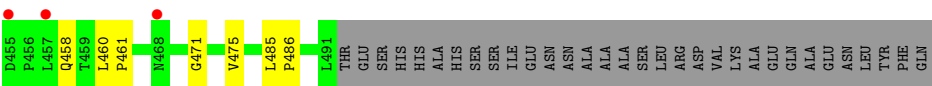
- Molecule 2 is N-[(1R)-1-phosphonoethyl]-L-alaninamide (CCD ID: AFS) (formula: C₅H₁₃N₂O₄P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			12	5	2	4	1		
2	B	1	Total	C	N	O	P	0	0
			12	5	2	4	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.24Å 107.73Å 205.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.86 – 3.20 29.86 – 3.20	Depositor EDS
% Data completeness (in resolution range)	77.6 (29.86-3.20) 77.4 (29.86-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 3.11Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.275 , 0.325 0.279 , 0.334	Depositor DCC
R_{free} test set	1258 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	81.4	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6858	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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: AFS, ZN

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	0.74	2/3517 (0.1%)	0.96	3/4787 (0.1%)
1	B	0.77	1/3480 (0.0%)	0.98	2/4738 (0.0%)
All	All	0.75	3/6997 (0.0%)	0.97	5/9525 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	319	THR	CA-CB	5.66	1.63	1.53
1	A	100	THR	CA-CB	5.26	1.59	1.53
1	B	13	SER	C-O	5.16	1.30	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	GLY	N-CA-C	-5.44	106.19	112.50
1	B	425	MET	N-CA-C	5.43	119.62	112.89
1	A	98	VAL	CA-C-N	5.28	126.44	119.84
1	A	98	VAL	C-N-CA	5.28	126.44	119.84
1	A	99	PRO	N-CA-C	5.05	122.88	112.47

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	3434	0	3450	37	0
1	B	3399	0	3420	35	0
2	A	12	0	11	0	0
2	B	12	0	11	0	0
3	A	1	0	0	0	0
All	All	6858	0	6892	72	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 5.

All (72) 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:101:GLU:OE1	1:A:105:PHE:CE2	1.86	1.28
1:A:101:GLU:OE1	1:A:105:PHE:CD2	2.28	0.85
1:A:389:TRP:HA	1:A:392:ILE:HD12	1.76	0.67
1:B:294:MET:HA	1:B:298:LEU:HB3	1.78	0.65
1:B:389:TRP:HA	1:B:392:ILE:HD12	1.77	0.65
1:B:55:SER:HB3	1:B:438:TYR:HA	1.79	0.65
1:B:425:MET:O	1:B:429:TYR:HB3	1.97	0.64
1:A:294:MET:HA	1:A:298:LEU:HB3	1.81	0.62
1:B:436:SER:HA	1:B:439:LEU:HD12	1.82	0.62
1:A:39:MET:HE1	1:A:53:VAL:HB	1.82	0.60
1:B:306:VAL:HG22	1:B:322:PRO:HG3	1.85	0.59
1:B:277:ALA:HB1	1:B:413:ILE:HD12	1.87	0.57
1:A:306:VAL:HG22	1:A:322:PRO:HG3	1.86	0.56
1:A:78:THR:HB	1:A:204:ALA:HB1	1.86	0.56
1:A:436:SER:HA	1:A:439:LEU:HD12	1.89	0.55
1:A:71:VAL:HA	1:A:75:ILE:HD12	1.89	0.55
1:A:55:SER:HB3	1:A:438:TYR:HA	1.91	0.53
1:A:77:GLY:HA2	1:A:206:TYR:HB2	1.92	0.52
1:B:385:LYS:HG3	1:B:460:LEU:HD13	1.91	0.52
1:A:329:ASN:HB3	1:A:330:PRO:HD3	1.92	0.51
1:A:132:TYR:N	1:A:133:GLU:HA	2.26	0.50
1:A:385:LYS:HG3	1:A:460:LEU:HD13	1.93	0.50
1:A:333:ILE:HG13	1:A:399:SER:HB2	1.93	0.50
1:A:439:LEU:O	1:A:442:VAL:HG22	2.12	0.50
1:B:39:MET:HE1	1:B:53:VAL:HG22	1.93	0.50
1:B:329:ASN:HB3	1:B:330:PRO:HD3	1.94	0.50
1:B:84:LEU:HD23	1:B:119:LEU:HD13	1.95	0.49
1:B:458:GLN:O	1:B:461:PRO:HD2	2.12	0.49
1:B:327:ALA:O	1:B:330:PRO:HD2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LEU:HD11	1:A:237:ILE:HD11	1.96	0.48
1:A:126:ASN:O	1:A:130:LYS:HG2	2.12	0.48
1:B:370:GLY:HA2	1:B:373:ILE:HD12	1.96	0.48
1:A:84:LEU:HD23	1:A:119:LEU:HD13	1.96	0.48
1:B:333:ILE:HG13	1:B:399:SER:HB2	1.96	0.47
1:A:107:PHE:CE2	1:A:239:GLU:HG2	2.48	0.47
1:B:306:VAL:HG13	1:B:322:PRO:HD3	1.96	0.47
1:B:195:ASN:O	1:B:199:MET:HG2	2.14	0.47
1:B:460:LEU:HB3	1:B:461:PRO:HD3	1.98	0.46
1:A:39:MET:HE2	1:A:50:ALA:HA	1.96	0.46
1:A:277:ALA:HB1	1:A:413:ILE:HD12	1.95	0.46
1:B:68:GLY:HA3	1:B:118:GLY:O	2.14	0.46
1:A:165:LYS:HD2	1:A:178:TRP:CZ3	2.50	0.46
1:A:327:ALA:O	1:A:330:PRO:HD2	2.16	0.46
1:B:485:LEU:N	1:B:486:PRO:HD2	2.31	0.46
1:A:97:THR:HA	1:A:179:HIS:HB3	1.98	0.45
1:A:127:LEU:O	1:A:131:ILE:HG13	2.16	0.45
1:A:458:GLN:O	1:A:461:PRO:HD2	2.17	0.45
1:B:168:VAL:HG12	1:B:177:GLY:H	1.81	0.45
1:B:206:TYR:HB3	1:B:207:GLY:H	1.62	0.45
1:B:219:SER:HA	1:B:222:ILE:HG22	1.99	0.45
1:A:425:MET:O	1:A:429:TYR:HB3	2.17	0.44
1:B:439:LEU:O	1:B:442:VAL:HG22	2.17	0.44
1:B:65:PRO:HB3	1:B:122:PRO:HD3	1.99	0.44
1:B:471:GLY:O	1:B:475:VAL:HG23	2.18	0.44
1:A:332:TRP:CE2	1:A:396:ALA:HB2	2.52	0.43
1:A:304:ARG:O	1:A:385:LYS:NZ	2.42	0.42
1:A:446:PHE:HB2	1:A:466:LEU:HD13	2.02	0.42
1:A:98:VAL:HG11	1:A:105:PHE:CE1	2.54	0.42
1:A:101:GLU:OE1	1:A:105:PHE:HE2	1.78	0.42
1:B:23:TRP:O	1:B:26:PHE:HB3	2.20	0.42
1:B:102:ASN:C	1:B:104:TRP:H	2.27	0.42
1:B:88:ILE:O	1:B:91:VAL:HG12	2.20	0.42
1:B:106:MET:HE2	1:B:106:MET:HB3	1.96	0.41
1:A:187:VAL:O	1:A:191:VAL:HG23	2.21	0.41
1:B:96:MET:HG2	1:B:109:ALA:HB1	2.03	0.41
1:A:485:LEU:N	1:A:486:PRO:HD2	2.36	0.41
1:B:71:VAL:HA	1:B:75:ILE:HD12	2.02	0.41
1:B:39:MET:HE2	1:B:50:ALA:HA	2.01	0.41
1:B:25:ARG:HH22	1:B:117:ASN:ND2	2.18	0.41
1:A:321:SER:HB3	1:A:324:GLN:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:ALA:HB2	1:A:466:LEU:HD22	2.03	0.40
1:B:332:TRP:HA	1:B:335:VAL:HG12	2.02	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	443/523 (85%)	412 (93%)	25 (6%)	6 (1%)	9	39
1	B	440/523 (84%)	406 (92%)	30 (7%)	4 (1%)	14	47
All	All	883/1046 (84%)	818 (93%)	55 (6%)	10 (1%)	11	43

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	PRO
1	B	99	PRO
1	B	177	GLY
1	A	177	GLY
1	A	409	GLY
1	B	409	GLY
1	A	319	THR
1	A	356	ILE
1	A	405	VAL
1	B	405	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	342/417 (82%)	325 (95%)	17 (5%)	22	55
1	B	340/417 (82%)	325 (96%)	15 (4%)	25	59
All	All	682/834 (82%)	650 (95%)	32 (5%)	23	57

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	ASP
1	A	82	MET
1	A	91	VAL
1	A	95	LEU
1	A	105	PHE
1	A	127	LEU
1	A	164	ILE
1	A	165	LYS
1	A	171	GLN
1	A	237	ILE
1	A	263	ILE
1	A	340	LEU
1	A	360	PHE
1	A	390	VAL
1	A	408	LEU
1	A	410	LEU
1	A	413	ILE
1	B	53	VAL
1	B	82	MET
1	B	91	VAL
1	B	102	ASN
1	B	127	LEU
1	B	164	ILE
1	B	165	LYS
1	B	171	GLN
1	B	237	ILE
1	B	340	LEU
1	B	360	PHE
1	B	390	VAL
1	B	408	LEU
1	B	410	LEU
1	B	413	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	261	HIS
1	A	437	GLN
1	B	41	GLN
1	B	261	HIS
1	B	292	GLN
1	B	329	ASN
1	B	458	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 3 ligands modelled in this entry, 1 is 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	$\# Z > 2$
2	AFS	B	601	-	9,11,11	3.29	5 (55%)	8,16,16	1.18	1 (12%)
2	AFS	A	601	-	9,11,11	2.51	3 (33%)	8,16,16	1.00	1 (12%)

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	AFS	B	601	-	-	2/14/14/14	-
2	AFS	A	601	-	-	0/14/14/14	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	AFS	P-O23	6.92	1.61	1.49
2	A	601	AFS	P-O21	5.10	1.62	1.54
2	A	601	AFS	P-O22	4.89	1.62	1.54
2	B	601	AFS	P-O22	4.59	1.62	1.54
2	B	601	AFS	P-CA2	-3.40	1.79	1.84
2	B	601	AFS	P-O21	-3.38	1.49	1.54
2	B	601	AFS	CB2-CA2	2.20	1.56	1.52
2	A	601	AFS	P-CA2	2.18	1.87	1.84

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	AFS	O22-P-O23	-2.61	106.91	113.45
2	A	601	AFS	O21-P-O23	-2.49	107.21	113.45

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	AFS	P-CA2-N2-C1
2	B	601	AFS	CB2-CA2-N2-C1

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

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	455/523 (86%)	0.69	65 (14%) 6 5	55, 98, 161, 186	0
1	B	452/523 (86%)	0.70	58 (12%) 7 6	54, 98, 161, 186	0
All	All	907/1046 (86%)	0.69	123 (13%) 7 5	54, 98, 161, 186	0

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	383	ASN	9.5
1	A	135	ASP	6.6
1	A	383	ASN	6.2
1	A	386	THR	6.1
1	B	320	TRP	5.8
1	A	203	LEU	5.5
1	A	207	GLY	5.4
1	B	402	GLU	5.4
1	A	406	SER	5.0
1	A	101	GLU	4.8
1	A	381	ALA	4.8
1	B	203	LEU	4.7
1	A	458	GLN	4.7
1	A	175	GLU	4.7
1	A	410	LEU	4.5
1	B	410	LEU	4.5
1	B	412	MET	4.4
1	A	204	ALA	4.4
1	B	134	GLY	4.3
1	A	451	GLN	4.3
1	A	133	GLU	4.1
1	A	375	GLY	4.1
1	B	102	ASN	4.0
1	A	376	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	380	PHE	4.0
1	A	382	VAL	3.9
1	B	451	GLN	3.9
1	A	171	GLN	3.9
1	A	194	GLY	3.9
1	B	202	SER	3.8
1	B	121	LYS	3.7
1	A	269	SER	3.7
1	B	415	ARG	3.7
1	B	305	ASN	3.6
1	B	411	ALA	3.6
1	B	310	PHE	3.5
1	B	380	PHE	3.4
1	B	206	TYR	3.4
1	B	205	ASN	3.4
1	B	426	MET	3.4
1	B	175	GLU	3.4
1	B	74	LYS	3.4
1	A	326	GLN	3.3
1	B	196	TYR	3.3
1	B	319	THR	3.2
1	B	117	ASN	3.2
1	A	131	ILE	3.1
1	B	386	THR	3.1
1	A	388	SER	3.1
1	A	297	SER	3.0
1	A	172	TYR	2.9
1	A	468	ASN	2.9
1	A	170	ALA	2.9
1	B	172	TYR	2.9
1	A	353	ASP	2.9
1	B	455	ASP	2.8
1	A	44	GLY	2.8
1	B	450	PRO	2.8
1	B	62	TYR	2.8
1	B	403	LEU	2.8
1	A	355	SER	2.8
1	A	357	ALA	2.7
1	B	306	VAL	2.7
1	A	305	ASN	2.7
1	A	319	THR	2.7
1	B	355	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	306	VAL	2.7
1	B	382	VAL	2.7
1	B	457	LEU	2.7
1	A	79	LYS	2.7
1	A	408	LEU	2.7
1	B	398	TYR	2.7
1	A	102	ASN	2.6
1	B	357	ALA	2.6
1	A	457	LEU	2.6
1	A	454	VAL	2.6
1	B	308	TRP	2.5
1	A	167	TYR	2.5
1	A	134	GLY	2.5
1	B	101	GLU	2.5
1	B	263	ILE	2.5
1	A	455	ASP	2.4
1	B	195	ASN	2.4
1	B	100	THR	2.4
1	B	200	HIS	2.4
1	B	468	ASN	2.4
1	A	426	MET	2.4
1	A	345	SER	2.4
1	A	205	ASN	2.4
1	B	111	GLY	2.3
1	A	411	ALA	2.3
1	A	308	TRP	2.3
1	A	402	GLU	2.3
1	A	379	GLN	2.3
1	A	58	ALA	2.2
1	B	12	HIS	2.2
1	A	449	VAL	2.2
1	A	163	TRP	2.2
1	A	398	TYR	2.2
1	A	453	LEU	2.2
1	A	268	PRO	2.2
1	B	409	GLY	2.2
1	B	64	SER	2.2
1	A	201	LYS	2.2
1	A	136	ASP	2.2
1	B	376	PHE	2.1
1	A	100	THR	2.1
1	B	44	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	133	GLU	2.1
1	B	399	SER	2.1
1	B	385	LYS	2.1
1	B	272	ALA	2.1
1	A	145	ILE	2.1
1	B	135	ASP	2.1
1	A	262	LEU	2.1
1	B	425	MET	2.1
1	A	416	TYR	2.1
1	B	132	TYR	2.1
1	B	295	SER	2.0
1	A	206	TYR	2.0
1	A	97	THR	2.0
1	A	492	THR	2.0
1	B	27	GLY	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
2	AFS	B	601	12/12	0.83	0.25	53,99,191,192	0
2	AFS	A	601	12/12	0.91	0.18	69,113,137,178	0
3	ZN	A	602	1/1	0.94	0.11	121,121,121,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.