



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:50 AM UTC

PDB ID : 6BX7 / pdb_00006bx7
Title : Crystal Structure of Human Protocadherin-1 EC1-4
Authors : Modak, D.; Sotomayor, M.
Deposited on : 2017-12-17
Resolution : 2.85 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

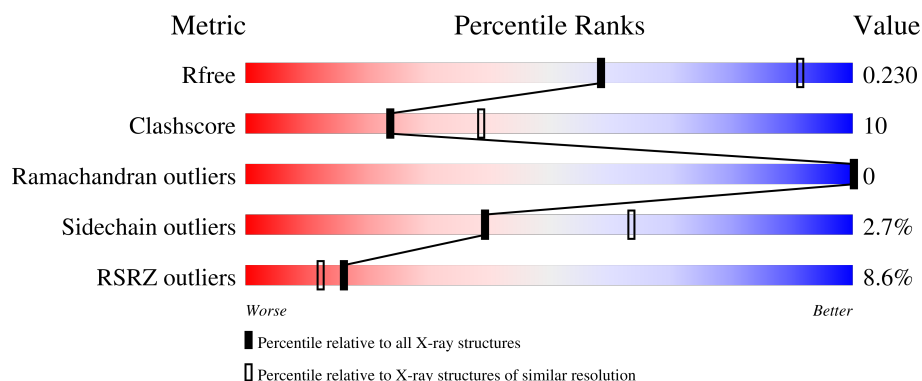
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	455	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3339 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.

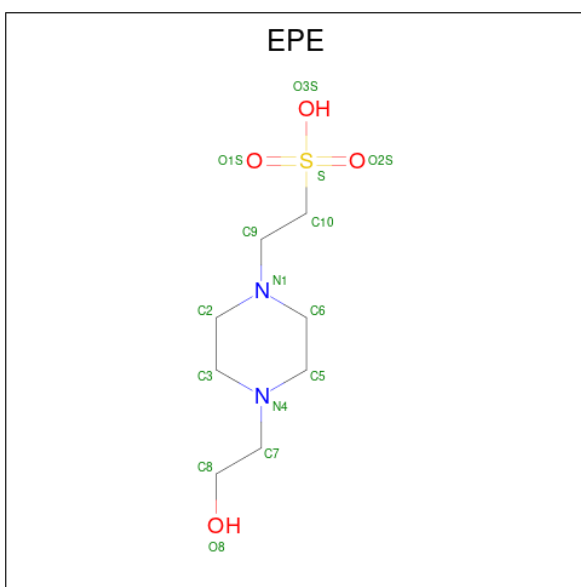
- Molecule 1 is a protein called Protocadherin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	0	0
			3288	2039	563	680	6			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q08174
A	447	LEU	-	expression tag	UNP Q08174
A	448	GLU	-	expression tag	UNP Q08174
A	449	HIS	-	expression tag	UNP Q08174
A	450	HIS	-	expression tag	UNP Q08174
A	451	HIS	-	expression tag	UNP Q08174
A	452	HIS	-	expression tag	UNP Q08174
A	453	HIS	-	expression tag	UNP Q08174
A	454	HIS	-	expression tag	UNP Q08174

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	Ca	0	0
			11	11		

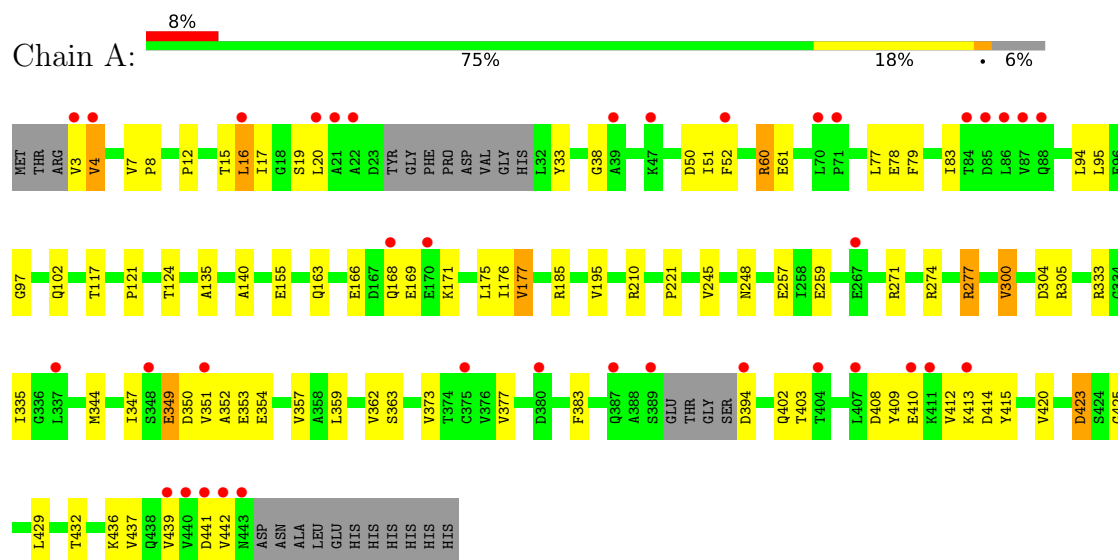
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		

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: Protocadherin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	146.98Å 146.98Å 147.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.26 – 2.85 48.26 – 2.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.26-2.85) 99.9 (48.26-2.85)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.221 , 0.243 (Not available) , 0.230	Depositor DCC
R_{free} test set	1120 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3339	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, CA

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	0.87	1/3338 (0.0%)	1.10	8/4548 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	300	VAL	N-CA	-6.24	1.39	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	423	ASP	CB-CA-C	-7.78	94.94	110.42
1	A	277	ARG	CB-CA-C	-6.24	99.08	110.63
1	A	394	ASP	CA-CB-CG	5.68	118.28	112.60
1	A	414	ASP	CA-CB-CG	5.42	118.02	112.60
1	A	350	ASP	CA-CB-CG	5.28	117.88	112.60
1	A	117	THR	CB-CA-C	5.14	118.53	111.23
1	A	121	PRO	CA-C-N	5.09	128.90	121.31
1	A	121	PRO	C-N-CA	5.09	128.90	121.31

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	185	ARG	Sidechain
1	A	210	ARG	Sidechain
1	A	277	ARG	Sidechain
1	A	38	GLY	Peptide
1	A	60	ARG	Sidechain

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	3288	0	3224	66	0
2	A	15	0	18	1	0
3	A	11	0	0	0	0
4	A	25	0	0	0	0
All	All	3339	0	3242	66	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 10.

All (66) 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:349:GLU:OE1	1:A:409:TYR:HB2	1.53	1.08
1:A:16:LEU:HA	1:A:52:PHE:CD1	1.92	1.03
1:A:349:GLU:OE1	1:A:441:ASP:HA	1.75	0.84
1:A:409:TYR:HA	1:A:412:VAL:CG2	2.07	0.84
1:A:124:THR:HG23	1:A:177:VAL:HG21	1.64	0.77
1:A:347:ILE:O	1:A:439:VAL:HA	1.90	0.71
1:A:423:ASP:HB3	1:A:425:GLY:H	1.56	0.71
1:A:349:GLU:OE1	1:A:409:TYR:CB	2.37	0.71
1:A:354:GLU:HB2	1:A:402:GLN:HE21	1.57	0.69
1:A:333:ARG:NH2	1:A:359:LEU:HD23	2.10	0.67
1:A:16:LEU:HA	1:A:52:PHE:CE1	2.31	0.66
1:A:94:LEU:C	1:A:94:LEU:HD12	2.24	0.63
1:A:60:ARG:NH1	1:A:61:GLU:OE2	2.33	0.61
1:A:16:LEU:HB2	1:A:52:PHE:CE1	2.36	0.60
1:A:415:TYR:HB2	1:A:437:VAL:HG22	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ILE:HD11	1:A:351:VAL:HG11	1.84	0.60
1:A:409:TYR:HA	1:A:412:VAL:HG21	1.84	0.59
1:A:166:GLU:OE1	1:A:171:LYS:HG2	2.03	0.58
1:A:168:GLN:O	1:A:169:GLU:HG2	2.02	0.58
1:A:19:SER:HB3	1:A:50:ASP:HB3	1.87	0.57
1:A:77:LEU:C	1:A:77:LEU:HD23	2.30	0.56
1:A:349:GLU:CD	1:A:442:VAL:H	2.13	0.56
1:A:83:ILE:HD13	1:A:95:LEU:HD23	1.87	0.56
1:A:409:TYR:CG	1:A:441:ASP:HB2	2.42	0.54
1:A:353:GLU:O	1:A:354:GLU:HG2	2.08	0.54
1:A:51:ILE:C	1:A:52:PHE:CD1	2.88	0.52
1:A:155:GLU:HG3	2:A:501:EPE:H62	1.92	0.52
1:A:51:ILE:O	1:A:52:PHE:CD1	2.63	0.52
1:A:163:GLN:HB2	1:A:176:ILE:HD11	1.93	0.51
1:A:354:GLU:HB2	1:A:402:GLN:NE2	2.26	0.51
1:A:77:LEU:HD23	1:A:78:GLU:N	2.25	0.51
1:A:274:ARG:HG2	1:A:274:ARG:HH21	1.74	0.51
1:A:423:ASP:OD2	1:A:429:LEU:HB2	2.11	0.51
1:A:335:ILE:CG2	1:A:359:LEU:HD22	2.41	0.51
1:A:51:ILE:O	1:A:52:PHE:HD1	1.94	0.50
1:A:79:PHE:CZ	1:A:97:GLY:HA3	2.47	0.50
1:A:20:LEU:HD22	1:A:33:TYR:CE1	2.48	0.49
1:A:412:VAL:HG23	1:A:439:VAL:HG21	1.94	0.49
1:A:245:VAL:HG21	1:A:300:VAL:HG11	1.94	0.48
1:A:409:TYR:HA	1:A:412:VAL:HG22	1.95	0.48
1:A:19:SER:HB3	1:A:50:ASP:CB	2.43	0.48
1:A:135:ALA:HB3	1:A:195:VAL:HG21	1.95	0.48
1:A:16:LEU:HA	1:A:52:PHE:HD1	1.69	0.47
1:A:383:PHE:CE1	1:A:403:THR:HG22	2.49	0.47
1:A:259:GLU:OE2	1:A:305:ARG:NH2	2.47	0.47
1:A:60:ARG:NH2	1:A:140:ALA:HB2	2.30	0.47
1:A:409:TYR:CD2	1:A:441:ASP:HB2	2.51	0.46
1:A:166:GLU:OE2	1:A:171:LYS:HE3	2.16	0.46
1:A:408:ASP:C	1:A:410:GLU:H	2.24	0.46
1:A:33:TYR:CE1	1:A:83:ILE:HD11	2.51	0.45
1:A:3:VAL:HG22	1:A:4:VAL:H	1.81	0.45
1:A:412:VAL:CG2	1:A:439:VAL:HG21	2.47	0.45
1:A:335:ILE:HG23	1:A:359:LEU:HB2	2.00	0.44
1:A:257:GLU:O	1:A:304:ASP:HA	2.18	0.43
1:A:8:PRO:HA	1:A:102:GLN:HB2	2.01	0.43
1:A:17:ILE:HG12	1:A:51:ILE:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:LEU:CB	1:A:52:PHE:CE1	3.01	0.42
1:A:19:SER:HB3	1:A:50:ASP:HA	2.01	0.42
1:A:221:PRO:HA	1:A:248:ASN:O	2.20	0.42
1:A:413:LYS:N	1:A:439:VAL:HG22	2.35	0.41
1:A:79:PHE:CE1	1:A:97:GLY:HA3	2.55	0.41
1:A:12:PRO:O	1:A:15:THR:OG1	2.37	0.41
1:A:362:VAL:CG1	1:A:373:VAL:HG11	2.51	0.41
1:A:351:VAL:O	1:A:352:ALA:C	2.64	0.41
1:A:344:MET:HE2	1:A:436:LYS:HE3	2.02	0.41
1:A:420:VAL:HG22	1:A:432:THR:HG22	2.03	0.41

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	423/455 (93%)	390 (92%)	33 (8%)	0	100 100

There are no Ramachandran outliers to report.

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	370/392 (94%)	360 (97%)	10 (3%)	39 63

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	7	VAL
1	A	16	LEU
1	A	175	LEU
1	A	177	VAL
1	A	271	ARG
1	A	349	GLU
1	A	357	VAL
1	A	363	SER
1	A	377	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	196	GLN
1	A	315	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 12 ligands modelled in this entry, 11 are monoatomic - leaving 1 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	EPE	A	501	-	15,15,15	1.99	1 (6%)	19,20,20	1.45	2 (10%)

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	EPE	A	501	-	-	3/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	EPE	C10-S	-7.15	1.67	1.77

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	EPE	O2S-S-C10	3.78	112.44	106.73
2	A	501	EPE	O3S-S-O2S	-2.14	106.04	111.40

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	EPE	C8-C7-N4-C3
2	A	501	EPE	C9-C10-S-O2S
2	A	501	EPE	C9-C10-S-O3S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	EPE	1	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, 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	429/455 (94%)	0.27	37 (8%) 16 12	39, 67, 136, 163	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	LYS	5.3
1	A	86	LEU	5.0
1	A	3	VAL	4.5
1	A	87	VAL	4.3
1	A	70	LEU	4.2
1	A	440	VAL	3.8
1	A	443	ASN	3.6
1	A	411	LYS	3.5
1	A	394	ASP	3.4
1	A	442	VAL	3.3
1	A	71	PRO	3.3
1	A	84	THR	3.2
1	A	21	ALA	3.1
1	A	337	LEU	3.1
1	A	348	SER	3.0
1	A	439	VAL	3.0
1	A	22	ALA	2.9
1	A	410	GLU	2.8
1	A	39	ALA	2.8
1	A	20	LEU	2.7
1	A	413	LYS	2.7
1	A	389	SER	2.7
1	A	85	ASP	2.6
1	A	16	LEU	2.5
1	A	375	CYS	2.5
1	A	170	GLU	2.4
1	A	387	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	380	ASP	2.3
1	A	4	VAL	2.3
1	A	351	VAL	2.2
1	A	407	LEU	2.2
1	A	168	GLN	2.2
1	A	267	GLU	2.2
1	A	52	PHE	2.1
1	A	441	ASP	2.1
1	A	88	GLN	2.0
1	A	404	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 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	EPE	A	501	15/15	0.87	0.19	97,106,112,112	0
3	CA	A	511	1/1	0.90	0.08	110,110,110,110	0
3	CA	A	512	1/1	0.94	0.05	62,62,62,62	0
3	CA	A	502	1/1	0.98	0.03	75,75,75,75	0
3	CA	A	508	1/1	0.98	0.03	66,66,66,66	0
3	CA	A	506	1/1	1.00	0.01	45,45,45,45	0
3	CA	A	507	1/1	1.00	0.03	40,40,40,40	0
3	CA	A	503	1/1	1.00	0.02	60,60,60,60	0
3	CA	A	509	1/1	1.00	0.03	49,49,49,49	0
3	CA	A	510	1/1	1.00	0.02	53,53,53,53	0
3	CA	A	504	1/1	1.00	0.01	50,50,50,50	0
3	CA	A	505	1/1	1.00	0.03	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.