



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:56 AM UTC

PDB ID : 6M7Z / pdb_00006m7z
Title : A divergent kinase lacking the glycine-rich loop regulates membrane ultra-structure of the Toxoplasma parasitophorous vacuole
Authors : Beraki, T.; Borek, D.M.; Reese, M.L.
Deposited on : 2018-08-21
Resolution : 2.50 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

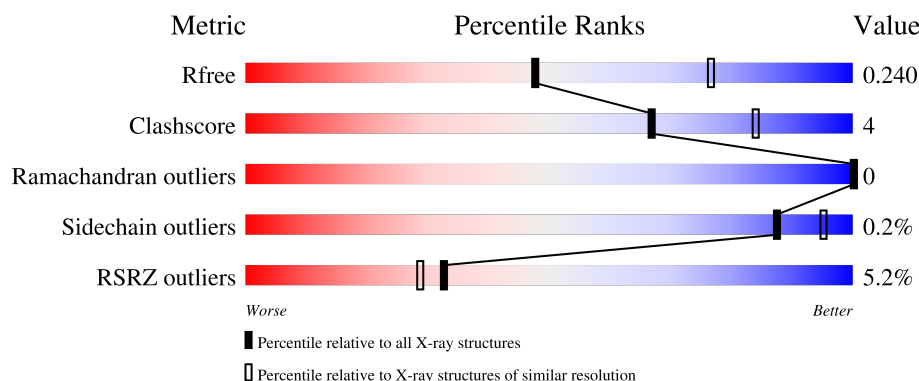
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	319	3% 76% 7% 18%
1	B	319	4% 74% 7% 19%
1	C	319	3% 77% 5% 18%
1	D	319	5% 73% 6% 20%
1	E	319	3% 76% 8% 16%

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Mol	Chain	Length	Quality of chain
1	F	319	7% 75% 6% 19%

2 Entry composition [i](#)

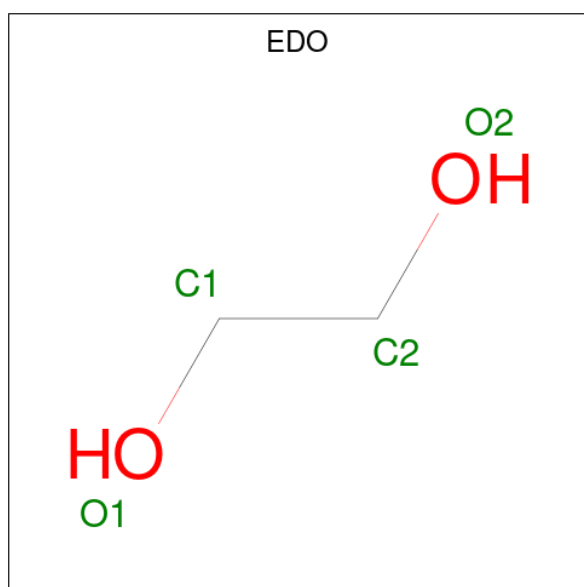
There are 4 unique types of molecules in this entry. The entry contains 12870 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bradyzoite pseudokinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	0	0
			2081	1333	360	382	6			
1	B	258	Total	C	N	O	S	0	0	0
			2043	1313	345	379	6			
1	C	262	Total	C	N	O	S	0	0	0
			2079	1334	356	383	6			
1	D	255	Total	C	N	O	S	0	1	0
			2023	1297	347	374	5			
1	E	269	Total	C	N	O	S	0	1	0
			2141	1370	373	392	6			
1	F	259	Total	C	N	O	S	0	0	0
			2055	1317	351	381	6			

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total Cl 1 1	0	0
3	E	2	Total Cl 2 2	0	0

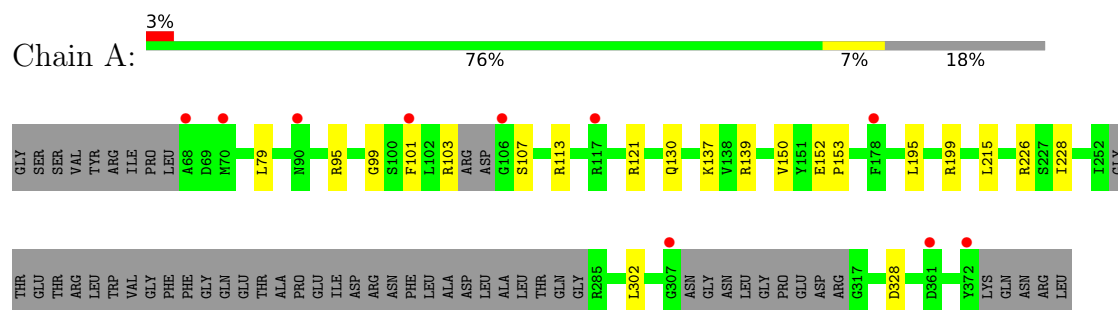
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total 46	O 46	0	0
4	B	51	Total 51	O 51	0	0
4	C	93	Total 93	O 93	0	0
4	D	67	Total 67	O 67	0	0
4	E	82	Total 82	O 82	0	0
4	F	42	Total 42	O 42	0	0

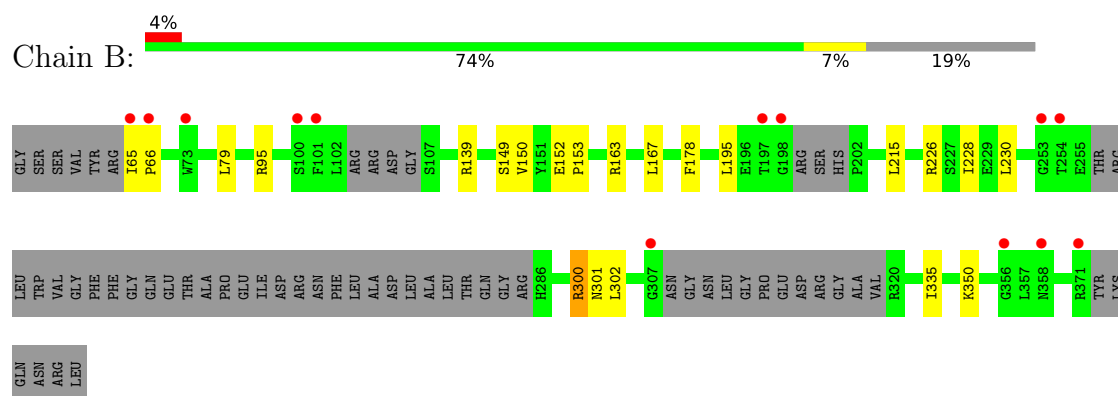
3 Residue-property plots [i](#)

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.

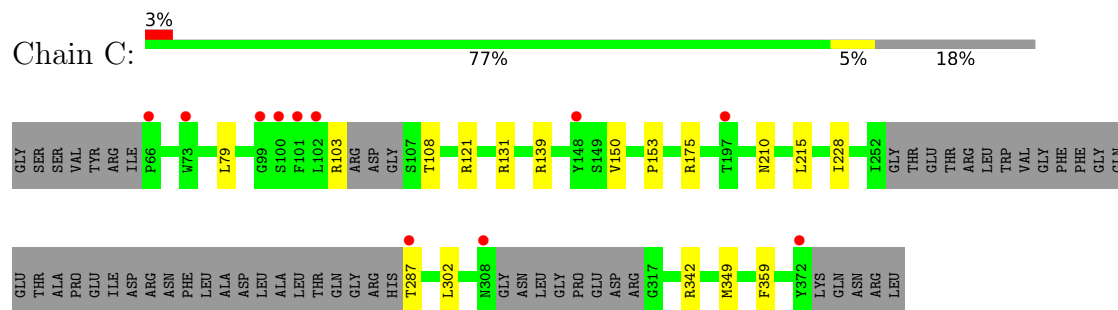
• Molecule 1: Bradyzoite pseudokinase 1



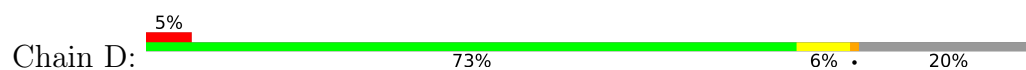
• Molecule 1: Bradyzoite pseudokinase 1

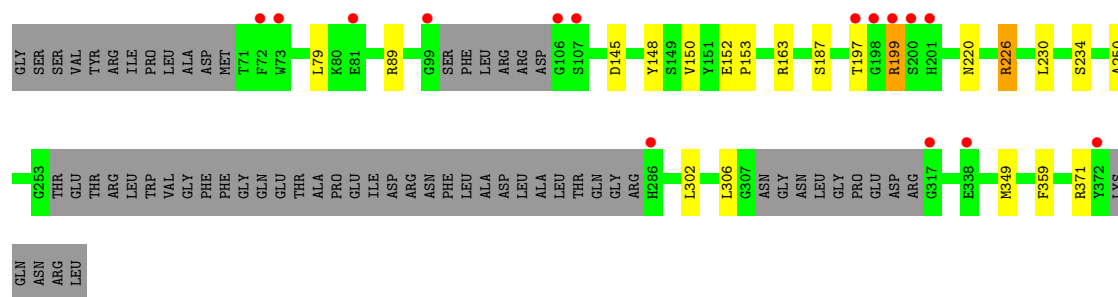


• Molecule 1: Bradyzoite pseudokinase 1

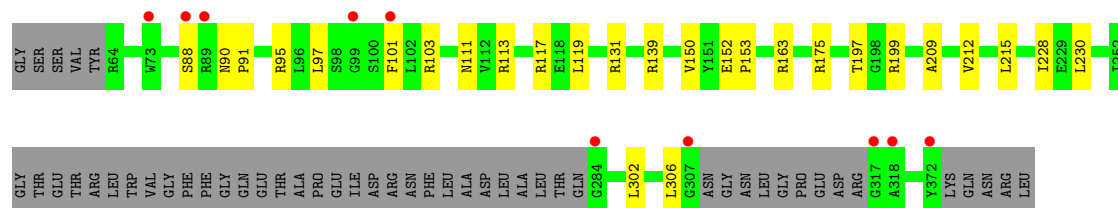
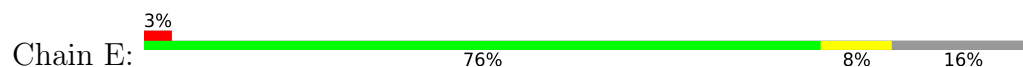


• Molecule 1: Bradyzoite pseudokinase 1

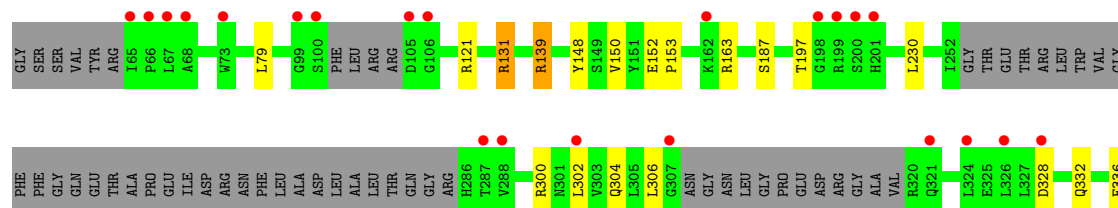
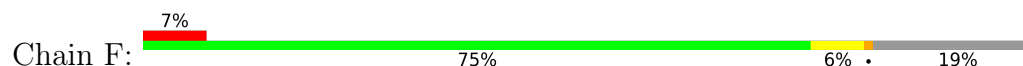




- Molecule 1: Bradyzoite pseudokinase 1



- Molecule 1: Bradyzoite pseudokinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.47Å 123.07Å 86.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.70 – 2.50 47.70 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.9 (47.70-2.50) 91.1 (47.70-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.63 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.197 , 0.239 0.200 , 0.240	Depositor DCC
R_{free} test set	2987 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.1	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	0.93	EDS
Total number of atoms	12870	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: CL, EDO

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	0.73	0/2122	0.95	0/2871
1	B	0.73	0/2082	0.94	0/2818
1	C	0.80	0/2120	0.94	0/2870
1	D	0.79	0/2066	0.96	0/2797
1	E	0.79	0/2187	0.97	0/2960
1	F	0.72	0/2096	0.94	0/2839
All	All	0.76	0/12673	0.95	0/17155

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	4
1	B	0	3
1	C	0	1
1	D	0	4
1	E	0	3
1	F	0	4
All	All	0	19

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	199	ARG	Sidechain
1	A	226	ARG	Sidechain
1	A	95	ARG	Sidechain
1	B	163	ARG	Sidechain
1	B	226	ARG	Sidechain
1	B	300	ARG	Sidechain
1	C	121	ARG	Sidechain
1	D	163	ARG	Sidechain
1	D	199	ARG	Sidechain
1	D	226	ARG	Sidechain
1	D	89	ARG	Sidechain
1	E	117	ARG	Sidechain
1	E	163	ARG	Sidechain
1	E	199	ARG	Sidechain
1	F	121	ARG	Sidechain
1	F	131	ARG	Sidechain
1	F	139	ARG	Sidechain
1	F	163	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	2081	0	2124	15	0
1	B	2043	0	2089	20	0
1	C	2079	0	2126	12	0
1	D	2023	0	2064	14	0
1	E	2141	0	2192	23	0
1	F	2055	0	2094	19	0
2	A	8	0	12	2	0
2	B	16	0	24	3	0
2	C	24	0	36	3	0
2	E	8	0	12	2	0
2	F	8	0	12	0	0
3	C	1	0	0	0	0
3	E	2	0	0	0	0
4	A	46	0	0	2	0
4	B	51	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	93	0	0	2	0
4	D	67	0	0	3	0
4	E	82	0	0	2	0
4	F	42	0	0	2	0
All	All	12870	0	12785	94	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 4.

All (94) 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:E:139:ARG:NH1	2:E:402:EDO:O1	1.97	0.97
1:D:199:ARG:HB3	1:E:197:THR:O	1.66	0.95
1:C:139:ARG:NH1	2:C:404:EDO:O1	2.04	0.91
1:B:350:LYS:NZ	4:B:501:HOH:O	2.02	0.90
1:B:178:PHE:HE2	1:E:97:LEU:HD12	1.35	0.90
1:E:113:ARG:O	1:F:304:GLN:NE2	2.15	0.77
1:A:139:ARG:NH1	2:A:402:EDO:O2	2.20	0.75
1:E:95:ARG:NH1	1:E:111:ASN:O	2.25	0.70
1:B:178:PHE:CE2	1:E:97:LEU:HD12	2.25	0.66
1:E:175:ARG:NH2	4:E:502:HOH:O	2.29	0.66
1:A:328:ASP:OD2	1:C:131:ARG:HD2	1.99	0.62
1:F:187:SER:HB2	4:F:501:HOH:O	1.99	0.62
1:A:195:LEU:CD1	1:A:302:LEU:CD1	2.79	0.60
1:F:300:ARG:NH2	1:F:332:GLN:HG2	2.18	0.58
1:B:139:ARG:HD2	2:B:404:EDO:H21	1.85	0.58
1:C:175:ARG:HD2	4:C:563:HOH:O	2.04	0.57
1:A:195:LEU:CD1	1:A:302:LEU:HD13	2.35	0.57
1:D:234:SER:OG	4:D:401:HOH:O	2.17	0.57
1:D:199:ARG:CB	1:E:197:THR:O	2.48	0.56
1:B:139:ARG:NH1	2:B:404:EDO:O2	2.38	0.56
1:B:230:LEU:HD12	1:B:301:ASN:OD1	2.06	0.55
1:E:88:SER:HB3	1:E:119:LEU:HD13	1.88	0.55
1:E:131:ARG:HD3	1:F:328:ASP:OD2	2.07	0.54
1:B:195:LEU:CD1	1:B:302:LEU:CD1	2.86	0.53
1:A:195:LEU:HD13	1:A:302:LEU:CD1	2.39	0.53
1:B:95:ARG:HH11	1:B:95:ARG:HG3	1.73	0.53
1:A:137:LYS:HD3	4:A:513:HOH:O	2.09	0.53
1:D:187:SER:HB2	4:D:401:HOH:O	2.08	0.52
1:B:167:LEU:HB2	2:B:401:EDO:H12	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:139:ARG:HH11	2:E:402:EDO:C1	2.23	0.51
1:F:336:GLU:HG2	1:F:341:ASN:CB	2.40	0.51
1:A:139:ARG:HH11	2:A:402:EDO:C2	2.23	0.50
1:D:220:ASN:ND2	4:D:402:HOH:O	2.29	0.50
1:B:195:LEU:CD1	1:B:302:LEU:HD13	2.40	0.50
1:B:65:ILE:N	1:B:66:PRO:CD	2.75	0.50
1:E:175:ARG:HD2	4:E:533:HOH:O	2.11	0.50
1:E:90:ASN:N	1:E:91:PRO:HD2	2.27	0.49
1:E:113:ARG:HB3	1:F:304:GLN:OE1	2.12	0.49
1:C:210:ASN:ND2	4:C:501:HOH:O	2.18	0.48
1:F:336:GLU:HG2	1:F:341:ASN:HB2	1.95	0.48
1:B:195:LEU:HD13	1:B:302:LEU:CD1	2.43	0.48
1:A:99:GLY:HA3	1:A:101:PHE:CE1	2.49	0.47
1:E:209:ALA:HA	1:E:212:VAL:HG22	1.97	0.47
1:E:302:LEU:C	1:E:302:LEU:HD23	2.40	0.47
1:F:139:ARG:NH1	1:F:148:TYR:CE2	2.83	0.47
1:F:300:ARG:HH21	1:F:332:GLN:HG2	1.80	0.47
1:D:302:LEU:C	1:D:302:LEU:HD23	2.40	0.46
1:F:187:SER:HB3	4:F:512:HOH:O	2.14	0.46
1:A:215:LEU:HD13	1:A:228:ILE:HD11	1.97	0.46
1:D:79:LEU:HD13	1:D:150:VAL:CG1	2.46	0.45
1:B:79:LEU:HD13	1:B:150:VAL:CG1	2.47	0.45
1:A:130:GLN:NE2	4:A:506:HOH:O	2.49	0.45
1:F:197:THR:HG22	1:F:197:THR:O	2.17	0.45
1:A:113:ARG:HD2	1:B:301:ASN:HD21	1.82	0.45
1:F:302:LEU:C	1:F:302:LEU:HD23	2.41	0.45
1:C:302:LEU:HD23	1:C:302:LEU:C	2.41	0.45
1:C:215:LEU:HD13	1:C:228:ILE:HD11	1.98	0.45
1:F:300:ARG:HH22	1:F:332:GLN:CD	2.24	0.44
1:F:302:LEU:CD2	1:F:306:LEU:HD12	2.46	0.44
1:C:103:ARG:HB2	1:C:108:THR:OG1	2.18	0.44
1:F:150:VAL:C	1:F:153:PRO:HD2	2.44	0.43
1:C:79:LEU:HD13	1:C:150:VAL:CG1	2.49	0.43
1:F:152:GLU:HB3	1:F:153:PRO:HD3	2.00	0.43
1:E:101:PHE:O	1:E:103:ARG:HG3	2.18	0.43
1:E:90:ASN:N	1:E:91:PRO:CD	2.82	0.43
1:C:153:PRO:HD3	2:C:404:EDO:H11	2.00	0.42
1:C:342:ARG:HH22	2:C:401:EDO:H12	1.84	0.42
1:D:349:MET:HE2	1:D:359:PHE:CD1	2.55	0.42
1:C:150:VAL:C	1:C:153:PRO:HD2	2.45	0.42
1:F:79:LEU:HD13	1:F:150:VAL:CG1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:LEU:HD13	1:A:150:VAL:CG1	2.50	0.42
1:D:152:GLU:HB3	1:D:153:PRO:HD3	2.02	0.42
1:B:152:GLU:HB3	1:B:153:PRO:HD3	2.02	0.42
1:C:349:MET:HE2	1:C:359:PHE:CD1	2.55	0.41
1:A:103:ARG:NH1	1:A:107:SER:O	2.53	0.41
1:D:226:ARG:HD2	1:D:250:ALA:HB3	2.02	0.41
1:B:95:ARG:HH11	1:B:95:ARG:CG	2.33	0.41
1:A:150:VAL:C	1:A:153:PRO:HD2	2.45	0.41
1:E:152:GLU:HB3	1:E:153:PRO:HD3	2.02	0.41
1:D:230:LEU:HD23	1:D:230:LEU:HA	1.96	0.41
1:B:150:VAL:C	1:B:153:PRO:HD2	2.46	0.41
1:D:197:THR:HG22	1:D:197:THR:O	2.21	0.41
1:F:230:LEU:HD23	1:F:230:LEU:HA	1.94	0.41
1:E:215:LEU:HD13	1:E:228:ILE:HD11	2.02	0.41
1:F:302:LEU:HD21	1:F:306:LEU:HD12	2.02	0.41
1:E:150:VAL:C	1:E:153:PRO:HD2	2.46	0.40
1:E:302:LEU:CD2	1:E:306:LEU:HD12	2.51	0.40
1:B:215:LEU:HD13	1:B:228:ILE:HD11	2.02	0.40
1:A:152:GLU:HB3	1:A:153:PRO:HD3	2.02	0.40
1:B:300:ARG:HD2	1:B:335:ILE:HD13	2.03	0.40
1:D:145:ASP:HB3	1:D:148:TYR:HD2	1.86	0.40
1:E:230:LEU:HD23	1:E:230:LEU:HA	1.97	0.40
1:B:149:SER:HB2	4:B:522:HOH:O	2.22	0.40
1:D:302:LEU:CD2	1:D:306:LEU:HD12	2.52	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	254/319 (80%)	251 (99%)	3 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	248/319 (78%)	245 (99%)	3 (1%)	0	100	100
1	C	254/319 (80%)	249 (98%)	5 (2%)	0	100	100
1	D	248/319 (78%)	245 (99%)	3 (1%)	0	100	100
1	E	264/319 (83%)	258 (98%)	6 (2%)	0	100	100
1	F	251/319 (79%)	248 (99%)	3 (1%)	0	100	100
All	All	1519/1914 (79%)	1496 (98%)	23 (2%)	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	231/278 (83%)	231 (100%)	0	100	100
1	B	229/278 (82%)	229 (100%)	0	100	100
1	C	232/278 (84%)	231 (100%)	1 (0%)	84	93
1	D	225/278 (81%)	224 (100%)	1 (0%)	84	93
1	E	238/278 (86%)	238 (100%)	0	100	100
1	F	230/278 (83%)	229 (100%)	1 (0%)	84	93
All	All	1385/1668 (83%)	1382 (100%)	3 (0%)	87	95

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	287	THR
1	D	371	ARG
1	F	131	ARG

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	286	HIS
1	A	301	ASN
1	A	323	HIS
1	A	341	ASN
1	B	323	HIS
1	C	323	HIS
1	D	341	ASN
1	E	201	HIS
1	F	341	ASN

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 19 ligands modelled in this entry, 3 are monoatomic - leaving 16 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	EDO	A	402	-	3,3,3	0.52	0	2,2,2	0.34	0
2	EDO	C	405	-	3,3,3	0.33	0	2,2,2	0.49	0
2	EDO	F	401	-	3,3,3	0.52	0	2,2,2	0.22	0
2	EDO	B	403	-	3,3,3	0.55	0	2,2,2	0.12	0
2	EDO	E	401	-	3,3,3	0.47	0	2,2,2	0.40	0
2	EDO	E	402	-	3,3,3	0.60	0	2,2,2	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	C	402	-	3,3,3	0.40	0	2,2,2	0.49	0
2	EDO	B	404	-	3,3,3	0.54	0	2,2,2	0.06	0
2	EDO	C	406	-	3,3,3	0.50	0	2,2,2	0.14	0
2	EDO	F	402	-	3,3,3	0.55	0	2,2,2	0.10	0
2	EDO	C	403	-	3,3,3	0.47	0	2,2,2	0.30	0
2	EDO	C	404	-	3,3,3	0.56	0	2,2,2	0.16	0
2	EDO	A	401	-	3,3,3	0.36	0	2,2,2	0.45	0
2	EDO	B	401	-	3,3,3	0.47	0	2,2,2	0.24	0
2	EDO	C	401	-	3,3,3	0.49	0	2,2,2	0.28	0
2	EDO	B	402	-	3,3,3	0.51	0	2,2,2	0.06	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	EDO	A	402	-	-	0/1/1/1	-
2	EDO	C	405	-	-	1/1/1/1	-
2	EDO	F	401	-	-	0/1/1/1	-
2	EDO	B	403	-	-	0/1/1/1	-
2	EDO	E	401	-	-	1/1/1/1	-
2	EDO	E	402	-	-	1/1/1/1	-
2	EDO	C	402	-	-	1/1/1/1	-
2	EDO	B	404	-	-	1/1/1/1	-
2	EDO	C	406	-	-	1/1/1/1	-
2	EDO	F	402	-	-	1/1/1/1	-
2	EDO	C	403	-	-	0/1/1/1	-
2	EDO	C	404	-	-	1/1/1/1	-
2	EDO	A	401	-	-	1/1/1/1	-
2	EDO	B	401	-	-	1/1/1/1	-
2	EDO	C	401	-	-	0/1/1/1	-
2	EDO	B	402	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	404	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	E	401	EDO	O1-C1-C2-O2
2	A	401	EDO	O1-C1-C2-O2
2	C	405	EDO	O1-C1-C2-O2
2	C	406	EDO	O1-C1-C2-O2
2	C	402	EDO	O1-C1-C2-O2
2	E	402	EDO	O1-C1-C2-O2
2	B	401	EDO	O1-C1-C2-O2
2	B	404	EDO	O1-C1-C2-O2
2	F	402	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	EDO	2	0
2	E	402	EDO	2	0
2	B	404	EDO	2	0
2	C	404	EDO	2	0
2	B	401	EDO	1	0
2	C	401	EDO	1	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	262/319 (82%)	0.19	10 (3%)	44 39	24, 46, 90, 110	0
1	B	258/319 (80%)	0.23	13 (5%)	34 30	23, 41, 86, 109	0
1	C	262/319 (82%)	-0.07	11 (4%)	40 36	21, 30, 73, 138	1 (0%)
1	D	255/319 (79%)	0.06	15 (5%)	28 24	20, 33, 81, 120	2 (0%)
1	E	269/319 (84%)	0.03	10 (3%)	45 40	18, 34, 80, 127	1 (0%)
1	F	259/319 (81%)	0.38	23 (8%)	15 13	25, 48, 95, 109	1 (0%)
All	All	1565/1914 (81%)	0.14	82 (5%)	33 29	18, 39, 86, 138	5 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	102	LEU	4.8
1	B	65	ILE	4.8
1	A	68	ALA	4.5
1	E	372	TYR	4.3
1	F	307	GLY	4.1
1	E	101	PHE	4.0
1	C	100	SER	3.8
1	C	101	PHE	3.8
1	E	73	TRP	3.7
1	B	198	GLY	3.5
1	D	73	TRP	3.5
1	D	72	PHE	3.5
1	D	106	GLY	3.4
1	C	99	GLY	3.4
1	B	356	GLY	3.4
1	C	308	ASN	3.3
1	D	317	GLY	3.3
1	F	200	SER	3.3
1	A	106	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	F	100	SER	3.2
1	F	201	HIS	3.2
1	B	371	ARG	3.1
1	A	178	PHE	3.1
1	C	73	TRP	3.1
1	D	199	ARG	3.1
1	F	65	ILE	3.1
1	F	68	ALA	3.1
1	D	201	HIS	3.1
1	B	358	ASN	3.1
1	C	372	TYR	3.0
1	C	66	PRO	3.0
1	E	317	GLY	2.9
1	B	254	THR	2.9
1	B	101	PHE	2.9
1	C	287	THR	2.9
1	D	372	TYR	2.8
1	D	81	GLU	2.8
1	A	361	ASP	2.8
1	D	107	SER	2.7
1	E	88	SER	2.7
1	F	73	TRP	2.7
1	F	99	GLY	2.6
1	F	67	LEU	2.6
1	E	307	GLY	2.6
1	A	101	PHE	2.6
1	A	372	TYR	2.6
1	B	307	GLY	2.5
1	C	148	TYR	2.5
1	C	197	THR	2.5
1	B	66	PRO	2.4
1	D	198	GLY	2.4
1	E	318	ALA	2.4
1	F	287	THR	2.4
1	D	286	HIS	2.4
1	A	117	ARG	2.3
1	E	89	ARG	2.3
1	F	288	VAL	2.3
1	F	326	LEU	2.3
1	D	197	THR	2.3
1	B	100	SER	2.3
1	D	200	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	73	TRP	2.3
1	F	324	LEU	2.3
1	A	70	MET	2.3
1	F	321	GLN	2.2
1	A	307	GLY	2.2
1	F	198	GLY	2.2
1	F	105	ASP	2.2
1	B	253	GLY	2.2
1	F	302	LEU	2.2
1	F	162	LYS	2.2
1	A	90	ASN	2.2
1	F	66	PRO	2.2
1	B	197	THR	2.1
1	E	284	GLY	2.1
1	F	106	GLY	2.1
1	F	340	GLY	2.1
1	D	338	GLU	2.1
1	F	328	ASP	2.1
1	F	199	ARG	2.0
1	D	99	GLY	2.0
1	E	99	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	EDO	C	403	4/4	0.81	0.23	54,59,68,76	0
2	EDO	C	401	4/4	0.82	0.15	56,57,61,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	B	403	4/4	0.85	0.16	41,53,64,65	0
2	EDO	F	402	4/4	0.87	0.26	53,58,59,85	0
3	CL	E	404	1/1	0.87	0.13	62,62,62,62	0
2	EDO	C	406	4/4	0.89	0.15	43,45,51,62	0
2	EDO	F	401	4/4	0.89	0.09	38,39,46,48	0
3	CL	E	403	1/1	0.91	0.10	49,49,49,49	0
2	EDO	B	404	4/4	0.91	0.17	42,49,54,55	0
2	EDO	C	405	4/4	0.93	0.10	32,32,36,36	0
2	EDO	C	402	4/4	0.93	0.09	31,38,45,48	0
2	EDO	B	402	4/4	0.94	0.08	36,39,40,41	0
2	EDO	A	401	4/4	0.94	0.10	32,43,49,51	0
2	EDO	B	401	4/4	0.94	0.10	34,41,50,52	0
2	EDO	C	404	4/4	0.95	0.13	24,39,51,55	0
2	EDO	A	402	4/4	0.96	0.09	29,42,43,47	0
3	CL	C	407	1/1	0.96	0.12	59,59,59,59	0
2	EDO	E	402	4/4	0.97	0.13	28,32,58,60	0
2	EDO	E	401	4/4	0.97	0.08	27,28,30,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.