



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 05:13 PM UTC

PDB ID : 3CWC / pdb_00003cwc
Title : Crystal structure of putative glycerate kinase 2 from Salmonella typhimurium LT2
Authors : Osipiuk, J.; Zhou, M.; Holzle, D.; Anderson, W.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2008-04-21
Resolution : 2.23 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

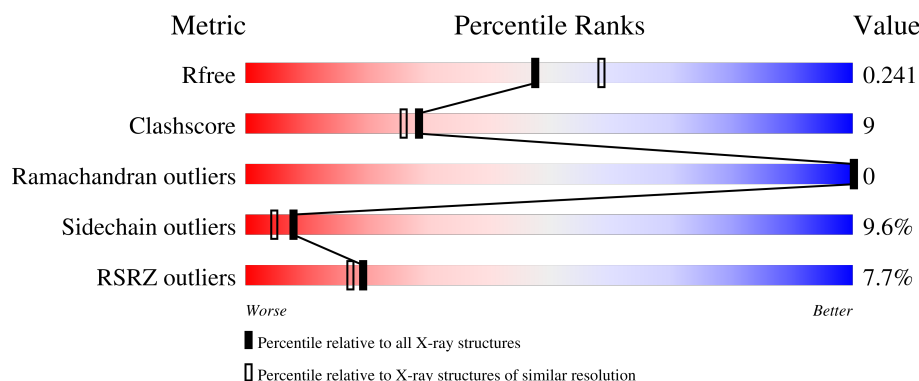
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	383	
1	B	383	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5567 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 Putative glycerate kinase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	Se	0	0	0
			2685	1673	477	523	5	7			
1	B	374	Total	C	N	O	S	Se	0	0	0
			2720	1695	483	530	5	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q8ZME0
A	-1	ASN	-	expression tag	UNP Q8ZME0
A	0	ALA	-	expression tag	UNP Q8ZME0
B	-2	SER	-	expression tag	UNP Q8ZME0
B	-1	ASN	-	expression tag	UNP Q8ZME0
B	0	ALA	-	expression tag	UNP Q8ZME0

- 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	B	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	B	1	Total Cl 1 1	0	0

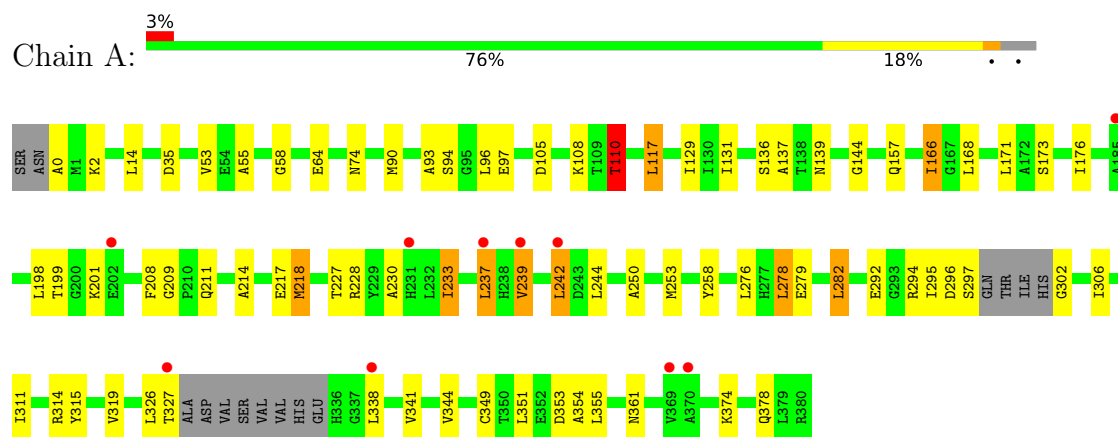
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	70	Total O 70 70	0	0
4	B	83	Total O 83 83	0	0

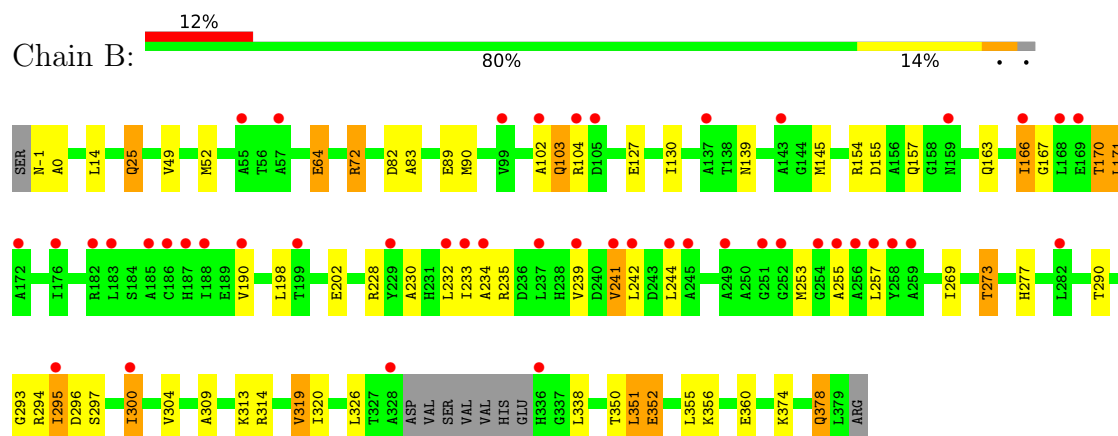
3 Residue-property plots

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: Putative glycerate kinase 2



• Molecule 1: Putative glycerate kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.73Å 59.55Å 103.61Å 90.00° 106.01° 90.00°	Depositor
Resolution (Å)	49.81 – 2.23 49.81 – 2.23	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.81-2.23) 99.8 (49.81-2.23)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.22Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.188 , 0.237 0.192 , 0.241	Depositor DCC
R_{free} test set	1834 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	46.4	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5567	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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	0.95	1/2713 (0.0%)	1.02	2/3663 (0.1%)
1	B	1.03	2/2750 (0.1%)	1.04	0/3718
All	All	0.99	3/5463 (0.1%)	1.03	2/7381 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	290	THR	CA-CB	6.67	1.62	1.53
1	B	300	ILE	CA-CB	5.28	1.61	1.54
1	A	341	VAL	N-CA	-5.02	1.40	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110	THR	CB-CA-C	-6.84	96.86	109.37
1	A	374	LYS	N-CA-C	5.77	117.25	111.07

There are no chirality outliers.

There are no planarity outliers.

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	2685	0	2699	47	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2720	0	2731	49	0
2	A	4	0	6	0	0
2	B	4	0	6	0	0
3	B	1	0	0	0	0
4	A	70	0	0	3	0
4	B	83	0	0	3	0
All	All	5567	0	5442	96	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 9.

All (96) 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:295:ILE:HG22	1:A:326:LEU:HD21	1.31	1.10
1:A:295:ILE:HA	1:A:326:LEU:HD23	1.58	0.83
1:B:309:ALA:HB2	1:B:338:LEU:HD23	1.65	0.77
1:B:155:ASP:HB2	1:B:170:THR:HG22	1.71	0.72
1:A:242:LEU:H	1:A:242:LEU:CD1	2.03	0.71
1:A:242:LEU:HD12	1:A:242:LEU:N	2.08	0.69
1:A:295:ILE:HG22	1:A:326:LEU:CD2	2.16	0.68
1:A:90:MSE:HE2	1:A:94:SER:HB2	1.75	0.68
1:A:242:LEU:H	1:A:242:LEU:HD12	1.58	0.68
1:B:269:ILE:HD13	1:B:304:VAL:HG22	1.76	0.67
1:B:350:THR:HG22	1:B:350:THR:O	1.95	0.66
1:B:25:GLN:HG2	4:B:460:HOH:O	1.95	0.66
1:B:52:MSE:CE	1:B:130:ILE:HD13	2.25	0.66
1:B:64:GLU:OE1	1:B:72:ARG:HB3	1.96	0.66
1:B:309:ALA:HB2	1:B:338:LEU:CD2	2.25	0.66
1:B:232:LEU:HD22	1:B:235:ARG:HH21	1.61	0.65
1:A:294:ARG:HD3	4:A:448:HOH:O	1.98	0.62
1:A:244:LEU:HD11	1:A:258:TYR:CG	2.35	0.62
1:A:278:LEU:HD13	1:A:311:ILE:HD12	1.80	0.62
1:A:208:PHE:O	1:A:211:GLN:HG2	1.99	0.61
1:A:319:VAL:HG22	1:A:338:LEU:CD2	2.32	0.59
1:A:230:ALA:HB1	1:A:242:LEU:HG	1.84	0.59
1:B:297:SER:O	1:B:300:ILE:HG22	2.02	0.58
1:B:49:VAL:HG23	1:B:89:GLU:HG3	1.84	0.57
1:A:279:GLU:OE2	1:A:315:TYR:OH	2.20	0.57
1:B:171:LEU:HD13	1:B:232:LEU:HD12	1.86	0.57
1:B:232:LEU:HD22	1:B:235:ARG:NH2	2.21	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ASP:OD1	1:A:218:MSE:HE2	2.06	0.56
1:A:242:LEU:CD1	1:A:242:LEU:N	2.67	0.56
1:B:14:LEU:HD23	1:B:355:LEU:HD22	1.89	0.55
1:A:53:VAL:HG13	1:A:58:GLY:C	2.33	0.54
1:A:233:ILE:HG23	1:A:239:VAL:HG13	1.90	0.54
1:B:351:LEU:O	1:B:352:GLU:HB3	2.07	0.53
1:A:176:ILE:HD12	1:A:176:ILE:O	2.09	0.53
1:A:250:ALA:HB3	1:A:253:MSE:HB3	1.90	0.52
1:B:350:THR:O	1:B:350:THR:CG2	2.58	0.52
1:A:108:LYS:CE	1:A:166:ILE:HG21	2.40	0.52
1:B:374:LYS:O	1:B:378:GLN:HG2	2.09	0.52
1:A:344:VAL:HG13	1:A:361:ASN:HB3	1.91	0.52
1:B:52:MSE:CE	1:B:130:ILE:CD1	2.87	0.52
1:A:349:CYS:HA	1:A:354:ALA:HB3	1.92	0.51
1:A:353:ASP:HB2	4:A:419:HOH:O	2.09	0.51
1:B:356:LYS:NZ	1:B:360:GLU:OE1	2.42	0.51
1:A:110:THR:HG23	4:A:424:HOH:O	2.11	0.51
1:B:155:ASP:CB	1:B:170:THR:HG22	2.40	0.51
1:A:129:ILE:HG22	1:A:131:ILE:HG13	1.93	0.51
1:B:166:ILE:HG22	1:B:167:GLY:N	2.25	0.51
1:B:294:ARG:HB2	1:B:300:ILE:HD13	1.93	0.50
1:B:52:MSE:HE3	1:B:130:ILE:CD1	2.41	0.50
1:A:0:ALA:HB3	1:A:35:ASP:OD2	2.11	0.50
1:A:319:VAL:HG22	1:A:338:LEU:HD22	1.94	0.49
1:A:295:ILE:HA	1:A:326:LEU:CD2	2.37	0.49
1:B:313:LYS:HG2	1:B:314:ARG:N	2.27	0.49
1:B:230:ALA:HB1	1:B:242:LEU:HG	1.94	0.49
1:A:311:ILE:O	1:A:314:ARG:HB3	2.13	0.48
1:B:277:HIS:HD2	4:B:511:HOH:O	1.96	0.48
1:A:278:LEU:HD22	1:A:282:LEU:HD22	1.96	0.48
1:B:230:ALA:HB3	1:B:242:LEU:HD21	1.95	0.48
1:B:102:ALA:C	1:B:103:GLN:HG2	2.39	0.48
1:B:269:ILE:O	1:B:273:THR:HB	2.14	0.48
1:A:237:LEU:HB3	1:A:239:VAL:HG12	1.96	0.48
1:B:241:VAL:HG22	1:B:255:ALA:HB1	1.94	0.48
1:A:93:ALA:HB2	1:A:117:LEU:HD22	1.96	0.47
1:B:295:ILE:CG2	1:B:296:ASP:N	2.77	0.47
1:A:96:LEU:HD13	1:A:137:ALA:HA	1.96	0.47
1:A:176:ILE:HG12	1:A:237:LEU:HD11	1.96	0.47
1:B:293:GLY:O	1:B:294:ARG:HD3	2.14	0.47
1:B:230:ALA:CB	1:B:242:LEU:CD2	2.94	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:ILE:HG22	1:B:296:ASP:N	2.31	0.46
1:A:105:ASP:OD1	1:A:218:MSE:CE	2.63	0.46
1:B:241:VAL:CG2	1:B:255:ALA:HB1	2.46	0.46
1:A:237:LEU:HD13	1:A:237:LEU:HA	1.73	0.46
1:B:90:MSE:HE2	1:B:253:MSE:HE1	1.97	0.46
1:B:228:ARG:CZ	1:B:228:ARG:HB2	2.47	0.45
1:B:14:LEU:HD23	1:B:355:LEU:CD2	2.47	0.45
1:B:104:ARG:HB2	4:B:468:HOH:O	2.16	0.44
1:B:145:MSE:HE1	1:B:257:LEU:HG	2.00	0.44
1:A:326:LEU:HD22	1:A:326:LEU:N	2.33	0.43
1:A:292:GLU:OE1	1:A:302:GLY:HA2	2.18	0.43
1:B:319:VAL:HG22	1:B:338:LEU:HD22	2.00	0.43
1:A:209:GLY:C	1:A:214:ALA:HB3	2.44	0.42
1:A:90:MSE:HE1	1:A:136:SER:HB2	2.02	0.42
1:B:234:ALA:HA	1:B:239:VAL:O	2.20	0.42
1:A:14:LEU:HD23	1:A:355:LEU:HD22	2.02	0.42
1:B:82:ASP:O	1:B:83:ALA:HB3	2.20	0.41
1:A:55:ALA:HB1	1:A:276:LEU:HD22	2.02	0.41
1:B:228:ARG:HB2	1:B:228:ARG:NH2	2.36	0.41
1:B:309:ALA:CB	1:B:338:LEU:HD23	2.43	0.41
1:B:154:ARG:O	1:B:171:LEU:HD23	2.21	0.41
1:A:110:THR:HG22	1:A:144:GLY:HA3	2.01	0.41
1:B:253:MSE:HG2	1:B:257:LEU:HD12	2.02	0.41
1:A:296:ASP:O	1:A:297:SER:C	2.64	0.41
1:A:344:VAL:CG1	1:A:361:ASN:HB3	2.51	0.41
1:B:295:ILE:HD12	1:B:295:ILE:HA	1.97	0.41
1:A:199:THR:HG21	1:A:227:THR:HA	2.03	0.40
1:B:-1:ASN:O	1:B:0:ALA:C	2.64	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	363/383 (95%)	353 (97%)	10 (3%)	0	100	100
1	B	370/383 (97%)	357 (96%)	13 (4%)	0	100	100
All	All	733/766 (96%)	710 (97%)	23 (3%)	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	269/275 (98%)	242 (90%)	27 (10%)	7	4
1	B	273/275 (99%)	248 (91%)	25 (9%)	8	5
All	All	542/550 (98%)	490 (90%)	52 (10%)	8	5

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	64	GLU
1	A	74	ASN
1	A	97	GLU
1	A	110	THR
1	A	117	LEU
1	A	139	ASN
1	A	157	GLN
1	A	166	ILE
1	A	168	LEU
1	A	171	LEU
1	A	173	SER
1	A	198	LEU
1	A	201	LYS
1	A	217	GLU
1	A	218	MSE
1	A	228	ARG
1	A	233	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	237	LEU
1	A	239	VAL
1	A	242	LEU
1	A	278	LEU
1	A	282	LEU
1	A	306	ILE
1	A	327	THR
1	A	351	LEU
1	A	378	GLN
1	B	25	GLN
1	B	64	GLU
1	B	72	ARG
1	B	103	GLN
1	B	127	GLU
1	B	139	ASN
1	B	157	GLN
1	B	163	GLN
1	B	166	ILE
1	B	170	THR
1	B	171	LEU
1	B	190	VAL
1	B	198	LEU
1	B	202	GLU
1	B	233	ILE
1	B	241	VAL
1	B	244	LEU
1	B	273	THR
1	B	295	ILE
1	B	319	VAL
1	B	320	ILE
1	B	326	LEU
1	B	351	LEU
1	B	352	GLU
1	B	378	GLN

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	62	HIS
1	A	74	ASN
1	A	103	GLN
1	A	163	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	25	GLN
1	B	74	ASN
1	B	120	HIS
1	B	147	GLN
1	B	159	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 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	EDO	A	402	-	3,3,3	0.55	0	2,2,2	0.67	0
2	EDO	B	401	-	3,3,3	0.51	0	2,2,2	0.36	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	-	-	1/1/1/1	-
2	EDO	B	401	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	362/383 (94%)	0.42	10 (2%) 55 55	28, 47, 62, 67	0
1	B	367/383 (95%)	0.79	46 (12%) 8 7	29, 45, 62, 69	0
All	All	729/766 (95%)	0.61	56 (7%) 19 17	28, 46, 62, 69	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	257	LEU	3.8
1	B	300	ILE	3.7
1	B	185	ALA	3.6
1	B	239	VAL	3.6
1	B	241	VAL	3.2
1	B	168	LEU	3.1
1	B	172	ALA	2.9
1	B	252	GLY	2.9
1	B	237	LEU	2.9
1	B	244	LEU	2.8
1	B	242	LEU	2.8
1	B	102	ALA	2.7
1	B	256	ALA	2.7
1	A	369	VAL	2.7
1	B	176	ILE	2.7
1	B	255	ALA	2.6
1	B	159	ASN	2.6
1	B	169	GLU	2.6
1	B	190	VAL	2.6
1	B	183	LEU	2.5
1	B	188	ILE	2.5
1	B	259	ALA	2.5
1	B	282	LEU	2.5
1	A	185	ALA	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	232	LEU	2.5
1	A	202	GLU	2.4
1	A	237	LEU	2.4
1	A	242	LEU	2.4
1	A	239	VAL	2.3
1	B	258	TYR	2.3
1	B	104	ARG	2.3
1	B	229	TYR	2.3
1	B	328	ALA	2.3
1	B	55	ALA	2.3
1	B	245	ALA	2.3
1	B	187	HIS	2.3
1	B	336	HIS	2.3
1	B	99	VAL	2.3
1	B	186	CYS	2.2
1	A	231	HIS	2.2
1	B	137	ALA	2.2
1	B	199	THR	2.2
1	B	182	ARG	2.2
1	B	254	GLY	2.2
1	A	338	LEU	2.2
1	B	233	ILE	2.2
1	B	143	ALA	2.2
1	B	234	ALA	2.2
1	B	166	ILE	2.1
1	B	57	ALA	2.1
1	B	249	ALA	2.1
1	B	105	ASP	2.1
1	A	370	ALA	2.1
1	B	251	GLY	2.0
1	A	327	THR	2.0
1	B	295	ILE	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	EDO	B	401	4/4	0.75	0.17	63,65,65,65	0
2	EDO	A	402	4/4	0.80	0.18	65,66,66,66	0
3	CL	B	403	1/1	0.99	0.15	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.