



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:10 AM UTC

PDB ID : 8GV2 / pdb_00008gv2
Title : Crystal structure of anti-FX IgG fab without FAST-Ig mutations
Authors : Koga, H.; Yamano, T.; Fukami, T.A.; Sampei, Z.; Shiraiwa, H.; Torizawa, T.
Deposited on : 2022-09-14
Resolution : 1.27 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

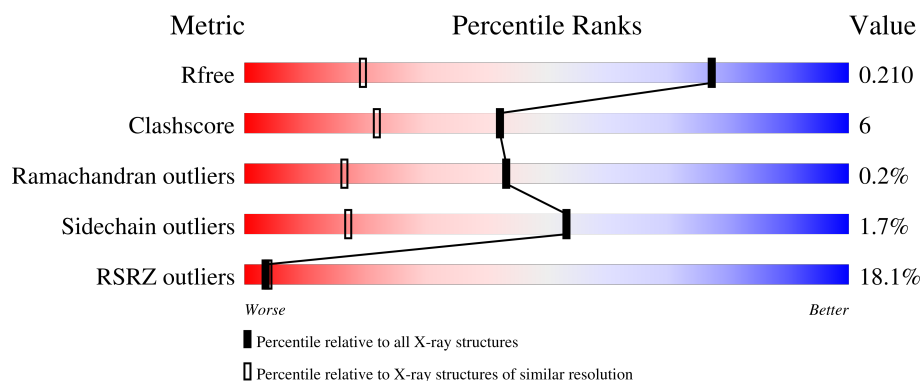
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.27 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	2836 (1.30-1.26)
Clashscore	190562	2911 (1.30-1.26)
Ramachandran outliers	187476	2841 (1.30-1.26)
Sidechain outliers	187428	2840 (1.30-1.26)
RSRZ outliers	180081	2832 (1.30-1.26)

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	224	
2	B	213	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	B	304	-	-	X	-
3	EDO	B	306	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3830 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 Anti-factor X IgG fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	10	0
			1678	1054	276	337	11			

- Molecule 2 is a protein called Anti-factor X IgG fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	9	0
			1612	1008	266	332	6			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			4	2	2		

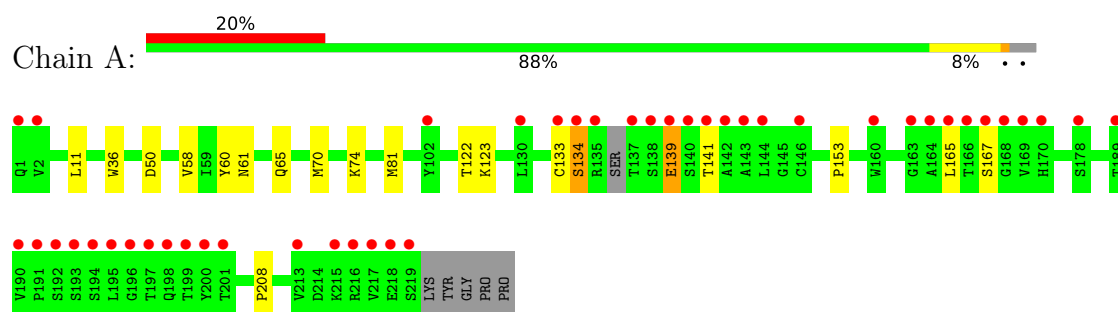
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	224	Total	O	0	0
			224	224		
4	B	220	Total	O	0	0
			220	220		

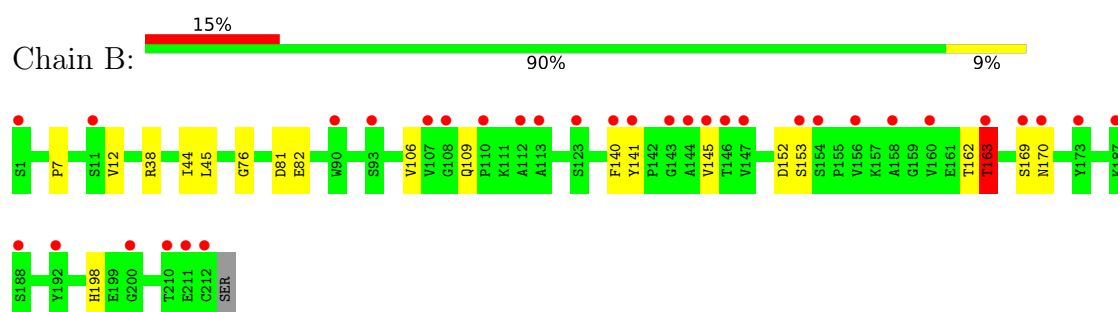
3 Residue-property plots [i](#)

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: Anti-factor X IgG fab heavy chain



- Molecule 2: Anti-factor X IgG fab light chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.51Å 62.29Å 81.59Å 90.00° 106.15° 90.00°	Depositor
Resolution (Å)	49.63 – 1.27 49.63 – 1.27	Depositor EDS
% Data completeness (in resolution range)	87.6 (49.63-1.27) 87.6 (49.63-1.27)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 1.28Å)	Xtriage
Refinement program	BUSTER 2.11.8 (8-JUN-2022)	Depositor
R, R_{free}	0.195 , 0.215 0.190 , 0.210	Depositor DCC
R_{free} test set	4672 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3830	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% 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: 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	0/1742	1.00	2/2369 (0.1%)
2	B	0.88	0/1677	1.04	4/2296 (0.2%)
All	All	0.92	0/3419	1.02	6/4665 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	ASN	N-CA-C	-6.51	99.29	109.25
2	B	140	PHE	CA-CB-CG	5.91	119.71	113.80
1	A	139	GLU	CB-CG-CD	5.22	121.48	112.60
2	B	169	SER	CA-C-N	5.22	131.52	121.18
2	B	169	SER	C-N-CA	5.22	131.52	121.18
2	B	163	THR	N-CA-CB	-5.05	102.61	110.29

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	1678	0	1635	22	0
2	B	1612	0	1552	20	0
3	A	60	0	90	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	36	0	54	14	0
4	A	224	0	0	3	2
4	B	220	0	0	2	0
All	All	3830	0	3331	42	2

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 6.

All (42) 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:81[B]:MET:CE	1:A:81[B]:MET:SD	2.18	1.32
2:B:44:ILE:HD11	3:B:304:EDO:C2	1.82	1.09
2:B:44:ILE:HD11	3:B:304:EDO:H21	1.41	1.01
1:A:81[B]:MET:HE3	1:A:81[B]:MET:HB2	1.54	0.90
2:B:44:ILE:HD11	3:B:304:EDO:H22	1.59	0.84
1:A:81[B]:MET:CE	1:A:81[B]:MET:HB2	2.12	0.80
1:A:133:CYS:HB2	3:B:303:EDO:H22	1.77	0.67
2:B:109:GLN:HB2	2:B:141:TYR:CE1	2.30	0.66
2:B:38:ARG:NH2	3:B:306:EDO:H11	2.10	0.66
2:B:44:ILE:CD1	3:B:304:EDO:H21	2.25	0.63
2:B:198:HIS:HD2	4:B:553:HOH:O	1.83	0.62
2:B:81:ASP:HA	3:B:306:EDO:H12	1.83	0.61
1:A:60[B]:TYR:OH	4:A:401:HOH:O	2.07	0.60
1:A:60[B]:TYR:CE2	1:A:65:GLN:HG3	2.37	0.59
1:A:36:TRP:HD1	1:A:70[A]:MET:HE3	1.68	0.59
2:B:76:GLY:HA3	3:B:305:EDO:H22	1.83	0.59
1:A:81[B]:MET:CE	1:A:81[B]:MET:CG	2.81	0.58
1:A:81[B]:MET:CE	1:A:81[B]:MET:CB	2.82	0.56
1:A:139:GLU:HG3	1:A:141:THR:H	1.72	0.55
1:A:50:ASP:O	1:A:58[A]:VAL:HG13	2.08	0.54
1:A:11:LEU:HD23	1:A:122[B]:THR:CG2	2.38	0.54
1:A:11:LEU:HD23	1:A:122[B]:THR:HG22	1.90	0.53
1:A:123:LYS:HA	3:A:308:EDO:H21	1.94	0.50
2:B:162:THR:HG21	3:B:309:EDO:H11	1.92	0.50
2:B:152:ASP:O	2:B:153:SER:HB3	2.12	0.49
3:A:315:EDO:H22	2:B:45:LEU:HD23	1.96	0.47
1:A:65:GLN:OE1	4:A:402:HOH:O	2.21	0.46
2:B:163:THR:HG21	4:B:420:HOH:O	2.16	0.45
1:A:65:GLN:NE2	4:A:412:HOH:O	2.48	0.45
1:A:74:LYS:HE2	3:A:307:EDO:H11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:82[A]:GLU:HG2	2:B:106:VAL:HG23	1.99	0.44
1:A:36:TRP:CD1	1:A:70[A]:MET:HE3	2.49	0.44
2:B:7:PRO:HB2	3:B:307:EDO:H12	1.99	0.44
2:B:81:ASP:OD1	3:B:306:EDO:C2	2.66	0.43
2:B:7:PRO:CB	3:B:307:EDO:H12	2.48	0.43
2:B:81:ASP:OD1	3:B:306:EDO:H21	2.18	0.43
1:A:36:TRP:CZ3	1:A:81[B]:MET:HG2	2.55	0.41
1:A:153:PRO:HD2	1:A:208:PRO:HB2	2.01	0.41
1:A:60[B]:TYR:CD1	1:A:65:GLN:HA	2.56	0.41
2:B:38:ARG:NH2	3:B:306:EDO:C1	2.82	0.41
1:A:60[B]:TYR:CD2	1:A:65:GLN:HG3	2.56	0.40
2:B:12:VAL:O	2:B:106:VAL:HA	2.22	0.40

All (2) 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)
4:A:597:HOH:O	4:A:597:HOH:O[2_656]	2.02	0.18
4:A:454:HOH:O	4:A:454:HOH:O[2_656]	2.16	0.04

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	224/224 (100%)	220 (98%)	3 (1%)	1 (0%)	30	8
2	B	219/213 (103%)	214 (98%)	5 (2%)	0	100	100
All	All	443/437 (101%)	434 (98%)	8 (2%)	1 (0%)	43	16

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	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	192/194 (99%)	189 (98%)	3 (2%)	55	19
2	B	182/178 (102%)	179 (98%)	3 (2%)	55	19
All	All	374/372 (100%)	368 (98%)	6 (2%)	53	19

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	SER
1	A	165	LEU
1	A	167	SER
2	B	145	VAL
2	B	163	THR
2	B	170	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	170	HIS
1	A	198	GLN
2	B	37	GLN
2	B	171	ASN
2	B	195	GLN
2	B	198	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](#)

24 ligands are 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	$\# Z > 2$
3	EDO	A	315	-	3,3,3	0.33	0	2,2,2	0.93	0
3	EDO	A	302	-	3,3,3	0.08	0	2,2,2	0.49	0
3	EDO	A	312	-	3,3,3	0.15	0	2,2,2	0.76	0
3	EDO	A	311	-	3,3,3	0.41	0	2,2,2	0.19	0
3	EDO	A	306	-	3,3,3	0.17	0	2,2,2	0.58	0
3	EDO	A	310	-	3,3,3	0.36	0	2,2,2	0.09	0
3	EDO	A	314	-	3,3,3	0.42	0	2,2,2	0.27	0
3	EDO	B	302	-	3,3,3	0.17	0	2,2,2	0.52	0
3	EDO	A	301	-	3,3,3	0.29	0	2,2,2	0.27	0
3	EDO	B	307	-	3,3,3	0.39	0	2,2,2	0.41	0
3	EDO	A	303	-	3,3,3	0.53	0	2,2,2	0.23	0
3	EDO	B	306	-	3,3,3	0.54	0	2,2,2	0.49	0
3	EDO	B	305	-	3,3,3	0.31	0	2,2,2	0.30	0
3	EDO	A	308	-	3,3,3	0.27	0	2,2,2	0.10	0
3	EDO	A	307	-	3,3,3	0.24	0	2,2,2	0.51	0
3	EDO	A	309	-	3,3,3	0.16	0	2,2,2	0.42	0
3	EDO	B	303	-	3,3,3	0.23	0	2,2,2	0.41	0
3	EDO	A	313	-	3,3,3	0.26	0	2,2,2	0.32	0
3	EDO	A	304	-	3,3,3	0.26	0	2,2,2	0.20	0
3	EDO	A	305	-	3,3,3	0.24	0	2,2,2	0.36	0
3	EDO	B	308	-	3,3,3	0.24	0	2,2,2	0.38	0
3	EDO	B	301	-	3,3,3	0.18	0	2,2,2	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	EDO	B	309	-	3,3,3	0.29	0	2,2,2	0.16	0
3	EDO	B	304	-	3,3,3	0.52	0	2,2,2	0.70	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
3	EDO	A	315	-	-	1/1/1/1	-
3	EDO	A	302	-	-	0/1/1/1	-
3	EDO	A	312	-	-	1/1/1/1	-
3	EDO	A	311	-	-	1/1/1/1	-
3	EDO	A	306	-	-	1/1/1/1	-
3	EDO	A	310	-	-	0/1/1/1	-
3	EDO	A	314	-	-	0/1/1/1	-
3	EDO	B	302	-	-	1/1/1/1	-
3	EDO	A	301	-	-	0/1/1/1	-
3	EDO	B	307	-	-	0/1/1/1	-
3	EDO	A	303	-	-	1/1/1/1	-
3	EDO	B	306	-	-	1/1/1/1	-
3	EDO	B	305	-	-	1/1/1/1	-
3	EDO	A	308	-	-	0/1/1/1	-
3	EDO	A	307	-	-	0/1/1/1	-
3	EDO	A	309	-	-	1/1/1/1	-
3	EDO	B	303	-	-	0/1/1/1	-
3	EDO	A	313	-	-	0/1/1/1	-
3	EDO	A	304	-	-	1/1/1/1	-
3	EDO	A	305	-	-	1/1/1/1	-
3	EDO	B	308	-	-	1/1/1/1	-
3	EDO	B	301	-	-	0/1/1/1	-
3	EDO	B	309	-	-	1/1/1/1	-
3	EDO	B	304	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	304	EDO	O1-C1-C2-O2
3	A	305	EDO	O1-C1-C2-O2
3	A	306	EDO	O1-C1-C2-O2
3	B	302	EDO	O1-C1-C2-O2
3	B	304	EDO	O1-C1-C2-O2
3	B	305	EDO	O1-C1-C2-O2
3	B	306	EDO	O1-C1-C2-O2
3	A	309	EDO	O1-C1-C2-O2
3	A	311	EDO	O1-C1-C2-O2
3	A	312	EDO	O1-C1-C2-O2
3	B	309	EDO	O1-C1-C2-O2
3	A	315	EDO	O1-C1-C2-O2
3	B	308	EDO	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	315	EDO	1	0
3	B	307	EDO	2	0
3	B	306	EDO	5	0
3	B	305	EDO	1	0
3	A	308	EDO	1	0
3	A	307	EDO	1	0
3	B	303	EDO	1	0
3	B	309	EDO	1	0
3	B	304	EDO	4	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	218/224 (97%)	1.01	45 (20%) 2 3	8, 23, 61, 71	10 (4%)
2	B	212/213 (99%)	0.91	33 (15%) 5 6	11, 23, 41, 60	9 (4%)
All	All	430/437 (98%)	0.96	78 (18%) 3 4	8, 23, 52, 71	19 (4%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	195	LEU	6.4
2	B	144	ALA	5.8
1	A	191	PRO	5.3
1	A	133	CYS	5.0
1	A	165	LEU	4.7
1	A	194	SER	4.7
1	A	219	SER	4.6
1	A	140	SER	4.6
1	A	200	TYR	4.4
1	A	137	THR	4.3
1	A	141	THR	4.2
2	B	173	TYR	4.2
2	B	145	VAL	4.2
1	A	190	VAL	4.1
2	B	170	ASN	3.9
1	A	135	ARG	3.8
1	A	166	THR	3.7
2	B	141	TYR	3.7
1	A	163	GLY	3.5
1	A	164	ALA	3.5
2	B	143	GLY	3.5
1	A	168	GLY	3.4
2	B	212	CYS	3.4
1	A	197	THR	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	90	TRP	3.4
1	A	134	SER	3.3
1	A	196	GLY	3.3
2	B	163	THR	3.3
1	A	142	ALA	3.2
1	A	146	CYS	3.2
1	A	217	VAL	3.2
2	B	110	PRO	3.2
1	A	138	SER	3.1
1	A	143	ALA	3.1
1	A	144	LEU	3.0
2	B	153	SER	3.0
1	A	169	VAL	3.0
1	A	139	GLU	2.9
1	A	193	SER	2.9
2	B	108	GLY	2.8
2	B	200	GLY	2.8
1	A	198	GLN	2.8
2	B	156	VAL	2.8
2	B	210	THR	2.8
1	A	160	TRP	2.8
2	B	187	LYS	2.7
2	B	1	SER	2.7
1	A	218	GLU	2.7
1	A	213	VAL	2.6
2	B	112	ALA	2.6
1	A	192	SER	2.6
1	A	1	GLN	2.6
2	B	93	SER	2.6
2	B	188	SER	2.5
2	B	211	GLU	2.5
1	A	102	TYR	2.5
2	B	160	VAL	2.5
2	B	113	ALA	2.5
2	B	169	SER	2.4
1	A	199	THR	2.4
1	A	201	THR	2.4
2	B	146	THR	2.4
2	B	192	TYR	2.4
1	A	189	THR	2.4
2	B	140	PHE	2.4
2	B	107	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	170	HIS	2.3
1	A	178	SER	2.3
1	A	215	LYS	2.2
2	B	158	ALA	2.2
1	A	216	ARG	2.2
1	A	130	LEU	2.1
2	B	123	SER	2.1
2	B	154	SER	2.1
1	A	2	VAL	2.1
1	A	167	SER	2.1
2	B	147	VAL	2.1
2	B	11	SER	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
3	EDO	B	303	4/4	0.61	0.32	72,72,72,72	0
3	EDO	B	306	4/4	0.66	0.21	41,42,42,42	0
3	EDO	A	315	4/4	0.67	0.17	38,38,39,40	0
3	EDO	A	313	4/4	0.71	0.18	49,49,49,50	0
3	EDO	A	312	4/4	0.74	0.18	30,32,34,36	0
3	EDO	B	308	4/4	0.74	0.18	60,60,60,60	0
3	EDO	A	303	4/4	0.75	0.16	30,33,36,38	0
3	EDO	A	310	4/4	0.75	0.17	40,42,43,44	0
3	EDO	A	304	4/4	0.76	0.16	61,61,61,61	0
3	EDO	A	308	4/4	0.78	0.18	55,55,55,55	0
3	EDO	A	306	4/4	0.80	0.18	47,48,48,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	B	305	4/4	0.81	0.14	41,43,44,45	0
3	EDO	B	309	4/4	0.82	0.15	54,54,54,54	0
3	EDO	A	302	4/4	0.83	0.13	29,32,35,38	0
3	EDO	A	311	4/4	0.83	0.15	35,37,38,40	0
3	EDO	A	307	4/4	0.84	0.13	33,35,36,37	0
3	EDO	B	304	4/4	0.85	0.14	26,29,32,34	0
3	EDO	B	307	4/4	0.87	0.13	34,36,38,38	0
3	EDO	B	301	4/4	0.88	0.12	35,35,36,36	0
3	EDO	B	302	4/4	0.88	0.12	30,33,34,36	0
3	EDO	A	309	4/4	0.88	0.13	41,41,42,42	0
3	EDO	A	305	4/4	0.91	0.11	39,40,41,42	0
3	EDO	A	301	4/4	0.92	0.10	30,32,34,36	0
3	EDO	A	314	4/4	0.93	0.10	26,29,29,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.