



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:43 PM UTC

PDB ID : 2IIU / pdb_00002iiu
Title : Crystal structure of a putative PhoU-like phosphate regulatory protein (NP_719307.1) from *Shewanella oneidensis* MR-1 at 2.28 Å resolution.
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2006-09-28
Resolution : 2.28 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

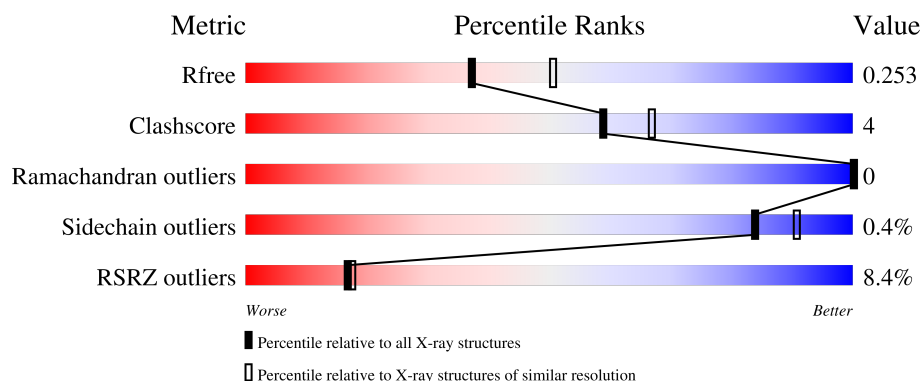
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

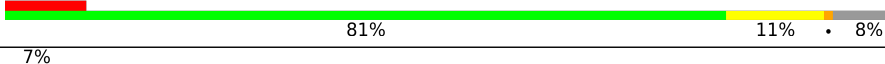
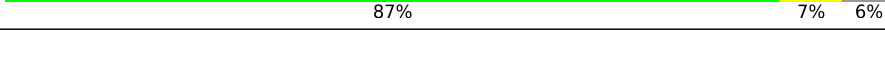
The reported resolution of this entry is 2.28 Å.

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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	227	
1	B	227	
1	C	227	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5045 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 Hypothetical protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	Se	0	3	0
			1618	1042	268	301	2	5			
1	B	209	Total	C	N	O	S	Se	0	0	0
			1585	1021	266	291	2	5			
1	C	214	Total	C	N	O	S	Se	0	0	0
			1623	1046	273	297	2	5			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q8EAX1
A	1	MSE	MET	modified residue	UNP Q8EAX1
A	22	MSE	MET	modified residue	UNP Q8EAX1
A	78	MSE	MET	modified residue	UNP Q8EAX1
A	161	MSE	MET	modified residue	UNP Q8EAX1
A	196	MSE	MET	modified residue	UNP Q8EAX1
A	222	MSE	MET	modified residue	UNP Q8EAX1
B	0	GLY	-	expression tag	UNP Q8EAX1
B	1	MSE	MET	modified residue	UNP Q8EAX1
B	22	MSE	MET	modified residue	UNP Q8EAX1
B	78	MSE	MET	modified residue	UNP Q8EAX1
B	161	MSE	MET	modified residue	UNP Q8EAX1
B	196	MSE	MET	modified residue	UNP Q8EAX1
B	222	MSE	MET	modified residue	UNP Q8EAX1
C	0	GLY	-	expression tag	UNP Q8EAX1
C	1	MSE	MET	modified residue	UNP Q8EAX1
C	22	MSE	MET	modified residue	UNP Q8EAX1
C	78	MSE	MET	modified residue	UNP Q8EAX1
C	161	MSE	MET	modified residue	UNP Q8EAX1
C	196	MSE	MET	modified residue	UNP Q8EAX1
C	222	MSE	MET	modified residue	UNP Q8EAX1

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



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

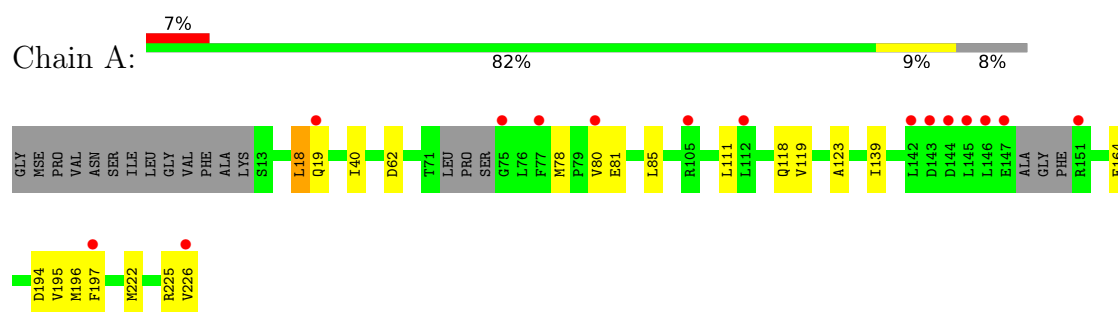
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	106	Total	O	0	0
			106	106		
3	B	63	Total	O	0	0
			63	63		
3	C	34	Total	O	0	0
			34	34		

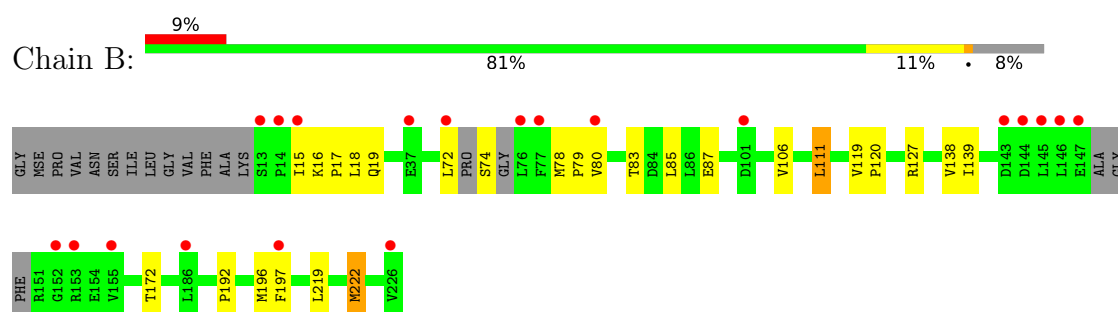
3 Residue-property plots [i](#)

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.

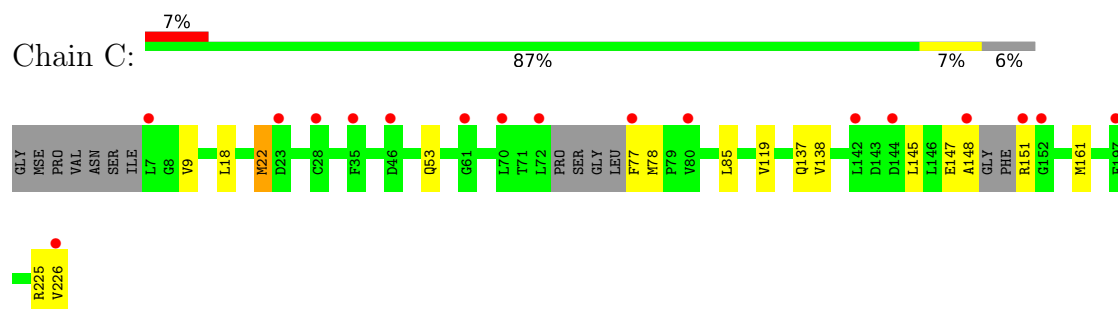
• Molecule 1: Hypothetical protein



• Molecule 1: Hypothetical protein



• Molecule 1: Hypothetical protein



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	99.88Å 99.88Å 397.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.46 – 2.28 47.46 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.46-2.28) 99.9 (47.46-2.28)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.65 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.200 , 0.247 0.208 , 0.253	Depositor DCC
R_{free} test set	1812 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5045	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% 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	1.06	3/1633 (0.2%)	0.99	0/2201
1	B	0.89	1/1598 (0.1%)	0.92	3/2152 (0.1%)
1	C	0.91	6/1638 (0.4%)	1.02	1/2209 (0.0%)
All	All	0.96	10/4869 (0.2%)	0.98	4/6562 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	137	GLN	CD-OE1	10.55	1.43	1.23
1	B	222	MSE	SE-CE	-8.01	1.71	1.95
1	C	53	GLN	CD-NE2	7.50	1.49	1.33
1	C	137	GLN	CD-NE2	7.00	1.48	1.33
1	C	53	GLN	CD-OE1	6.46	1.35	1.23
1	A	119	VAL	CA-CB	6.35	1.57	1.54
1	A	222	MSE	SE-CE	-6.27	1.76	1.95
1	C	119	VAL	CA-CB	5.31	1.57	1.54
1	C	22	MSE	SE-CE	-5.24	1.79	1.95
1	A	123	ALA	CA-CB	5.01	1.61	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	137	GLN	OE1-CD-NE2	7.63	130.23	122.60
1	B	78	MSE	CA-C-N	5.47	125.09	119.56
1	B	78	MSE	C-N-CA	5.47	125.09	119.56
1	B	111	LEU	CB-CA-C	-5.25	102.61	110.16

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	1618	0	1584	14	0
1	B	1585	0	1543	19	0
1	C	1623	0	1594	10	0
2	A	12	0	18	3	0
2	B	4	0	6	0	0
3	A	106	0	0	2	0
3	B	63	0	0	0	0
3	C	34	0	0	0	0
All	All	5045	0	4745	42	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 (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:C:18:LEU:CD1	1:C:85:LEU:HD22	2.13	0.78
1:A:18:LEU:HD13	1:A:85:LEU:HD22	1.71	0.72
1:B:72:LEU:O	1:B:74:SER:N	2.25	0.69
1:B:80:VAL:HG13	1:B:222:MSE:HE3	1.76	0.67
1:B:138:VAL:HG12	1:B:219:LEU:HD12	1.79	0.65
1:B:138:VAL:HG12	1:B:219:LEU:CD1	2.29	0.63
1:A:195:VAL:HG12	1:A:196:MSE:CE	2.29	0.62
1:A:111:LEU:HD22	1:A:197[B]:PHE:CE2	2.35	0.61
1:A:225:ARG:O	1:A:226:VAL:CB	2.49	0.60
1:C:225:ARG:O	1:C:226:VAL:CB	2.50	0.60
1:A:195:VAL:HG12	1:A:196:MSE:HE2	1.85	0.58
1:B:80:VAL:CG1	1:B:222:MSE:HE3	2.33	0.58
1:A:81:GLU:HA	2:A:229:EDO:H11	1.86	0.57
1:A:164:GLU:CD	1:C:9:VAL:HG13	2.30	0.56
1:A:19:GLN:HG2	1:A:139:ILE:HG22	1.86	0.56
1:B:80:VAL:HG11	1:B:85:LEU:HD21	1.88	0.55
2:A:227:EDO:H11	3:A:315:HOH:O	2.07	0.55
1:A:111:LEU:HD22	1:A:197[B]:PHE:HE2	1.73	0.54
1:B:83:THR:O	1:B:87:GLU:HG3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:MSE:HE3	1:A:80:VAL:O	2.08	0.52
1:C:22:MSE:O	1:C:22:MSE:HE3	2.09	0.52
1:C:18:LEU:HD13	1:C:85:LEU:HD22	1.91	0.50
1:A:78:MSE:HB2	2:A:229:EDO:H21	1.95	0.49
1:B:79:PRO:HB2	1:C:77:PHE:HB3	1.95	0.49
1:B:106:VAL:HG13	1:B:111:LEU:HB3	1.94	0.49
1:B:15:ILE:HG22	1:B:19:GLN:HG3	1.95	0.47
1:C:138:VAL:HG21	1:C:161:MSE:HG3	1.96	0.47
1:B:19:GLN:HG2	1:B:139:ILE:HG22	1.97	0.47
1:A:194:ASP:OD2	3:A:314:HOH:O	2.21	0.46
1:C:145:LEU:C	1:C:147:GLU:H	2.24	0.46
1:B:127:ARG:HB3	1:B:172:THR:HG22	1.99	0.45
1:C:148:ALA:C	1:C:151:ARG:N	2.75	0.45
1:B:138:VAL:CG1	1:B:219:LEU:HD12	2.45	0.44
1:A:196:MSE:HE2	1:A:196:MSE:HA	2.00	0.44
1:C:77:PHE:O	1:C:78:MSE:HG3	2.18	0.43
1:B:111:LEU:HD13	1:B:197:PHE:HD2	1.84	0.43
1:B:18:LEU:HD13	1:B:85:LEU:HD22	2.00	0.42
1:A:40:ILE:HG23	1:A:118:GLN:CD	2.45	0.42
1:B:16:LYS:HB3	1:B:17:PRO:HD3	2.02	0.42
1:B:119:VAL:HB	1:B:120:PRO:HD3	2.02	0.42
1:B:192:PRO:O	1:B:196:MSE:HG2	2.20	0.41
1:B:18:LEU:CD1	1:B:85:LEU:HD22	2.51	0.41

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	205/227 (90%)	204 (100%)	1 (0%)	0	100	100
1	B	202/227 (89%)	201 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	208/227 (92%)	204 (98%)	4 (2%)	0	100	100
All	All	615/681 (90%)	609 (99%)	6 (1%)	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	159/191 (83%)	157 (99%)	2 (1%)	61	76
1	B	152/191 (80%)	152 (100%)	0	100	100
1	C	156/191 (82%)	156 (100%)	0	100	100
All	All	467/573 (82%)	465 (100%)	2 (0%)	84	91

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	62	ASP

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	53	GLN
1	A	91	GLN
1	A	137	GLN
1	A	191	ASN
1	B	49	GLN
1	B	97	ASN
1	B	176	GLN
1	C	110	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](#)

4 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	$\# Z > 2$
2	EDO	A	228	-	3,3,3	0.47	0	2,2,2	0.14	0
2	EDO	A	227	-	3,3,3	0.34	0	2,2,2	0.47	0
2	EDO	A	229	-	3,3,3	0.52	0	2,2,2	0.22	0
2	EDO	B	227	-	3,3,3	0.98	0	2,2,2	0.14	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	228	-	-	0/1/1/1	-
2	EDO	A	227	-	-	1/1/1/1	-
2	EDO	A	229	-	-	1/1/1/1	-
2	EDO	B	227	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	227	EDO	O1-C1-C2-O2
2	A	229	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	227	EDO	1	0
2	A	229	EDO	2	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	203/227 (89%)	0.79	15 (7%)	20 21	19, 46, 70, 88	3 (1%)
1	B	204/227 (89%)	0.92	20 (9%)	13 14	36, 46, 67, 81	0
1	C	209/227 (92%)	1.02	17 (8%)	18 18	36, 46, 68, 79	0
All	All	616/681 (90%)	0.91	52 (8%)	17 17	19, 46, 71, 88	3 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	144	ASP	4.9
1	B	77	PHE	4.9
1	A	145	LEU	4.7
1	A	147	GLU	4.3
1	C	70	LEU	4.1
1	B	80	VAL	4.0
1	C	7	LEU	4.0
1	C	148	ALA	3.9
1	A	144	ASP	3.8
1	B	147	GLU	3.6
1	C	226	VAL	3.6
1	A	77	PHE	3.5
1	B	197	PHE	3.5
1	C	72	LEU	3.4
1	A	151	ARG	3.3
1	C	77	PHE	3.1
1	A	75	GLY	3.1
1	C	152	GLY	3.1
1	A	226	VAL	3.1
1	A	197[A]	PHE	3.0
1	A	80	VAL	2.9
1	B	76	LEU	2.9
1	B	153	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	151	ARG	2.9
1	B	146	LEU	2.7
1	B	15	ILE	2.7
1	A	146	LEU	2.6
1	C	28	CYS	2.6
1	B	37	GLU	2.6
1	C	35	PHE	2.6
1	B	143	ASP	2.5
1	C	23	ASP	2.4
1	B	101	ASP	2.3
1	B	72	LEU	2.2
1	B	226	VAL	2.3
1	C	80	VAL	2.3
1	B	152	GLY	2.2
1	C	61	GLY	2.2
1	B	155	VAL	2.2
1	B	14	PRO	2.2
1	C	46	ASP	2.2
1	B	13	SER	2.1
1	B	186	LEU	2.1
1	B	144	ASP	2.1
1	A	112	LEU	2.1
1	A	142	LEU	2.1
1	A	143	ASP	2.1
1	C	142	LEU	2.1
1	C	197	PHE	2.1
1	A	105	ARG	2.1
1	A	19	GLN	2.0
1	B	145	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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	227	4/4	0.76	0.22	38,52,57,61	0
2	EDO	A	229	4/4	0.77	0.17	62,69,72,72	0
2	EDO	A	227	4/4	0.90	0.24	38,48,49,50	0
2	EDO	A	228	4/4	0.93	0.28	46,52,53,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.