



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:18 PM UTC

PDB ID : 6EMG / pdb_00006emg
Title : Crystal structure of Rrp1 from Chaetomium thermophilum in space group P6322
Authors : Ahmed, Y.L.; Sinning, I.
Deposited on : 2017-10-02
Resolution : 2.24 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

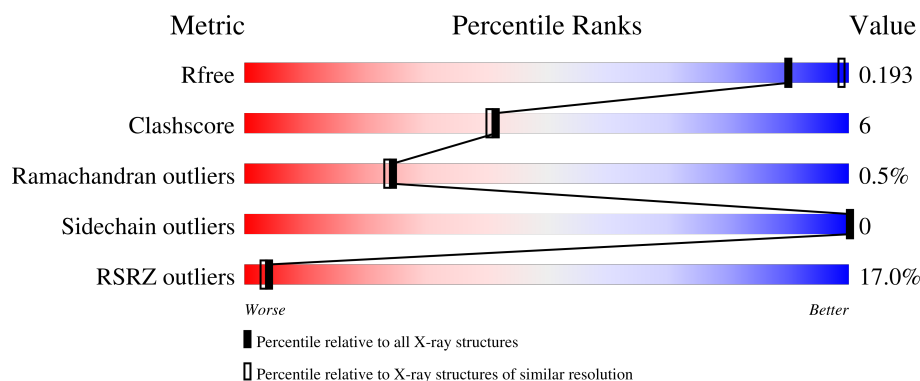
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	290	8% 73% 7% 20%
1	B	290	19% 69% 8% 22%

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	A	305	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3966 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 G0S4M2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	1	0
			1875	1195	332	337	11			
1	B	226	Total	C	N	O	S	0	1	0
			1839	1176	323	330	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	283	GLY	-	expression tag	UNP G0S4M2
A	284	SER	-	expression tag	UNP G0S4M2
A	285	HIS	-	expression tag	UNP G0S4M2
A	286	HIS	-	expression tag	UNP G0S4M2
A	287	HIS	-	expression tag	UNP G0S4M2
A	288	HIS	-	expression tag	UNP G0S4M2
A	289	HIS	-	expression tag	UNP G0S4M2
A	290	HIS	-	expression tag	UNP G0S4M2
B	283	GLY	-	expression tag	UNP G0S4M2
B	284	SER	-	expression tag	UNP G0S4M2
B	285	HIS	-	expression tag	UNP G0S4M2
B	286	HIS	-	expression tag	UNP G0S4M2
B	287	HIS	-	expression tag	UNP G0S4M2
B	288	HIS	-	expression tag	UNP G0S4M2
B	289	HIS	-	expression tag	UNP G0S4M2
B	290	HIS	-	expression tag	UNP G0S4M2

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



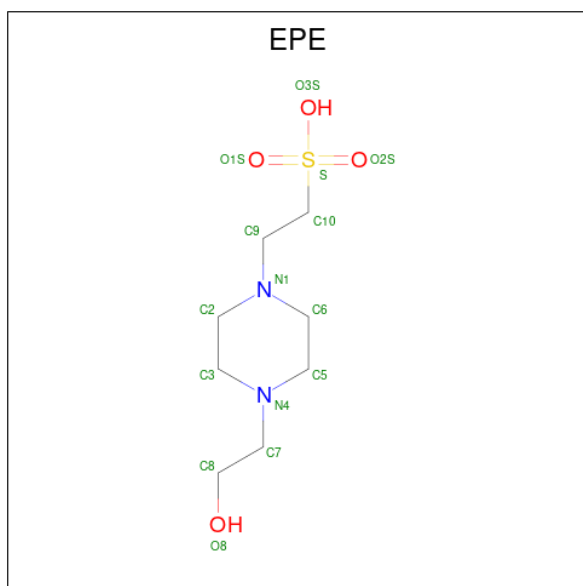
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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



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

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	141	Total O 141 141	0	0

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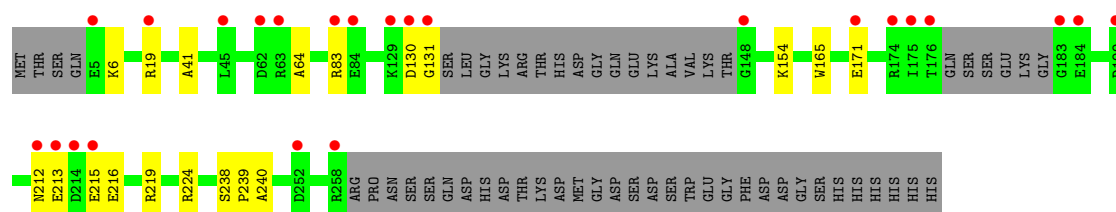
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	48	Total	O	0	0
			48	48		

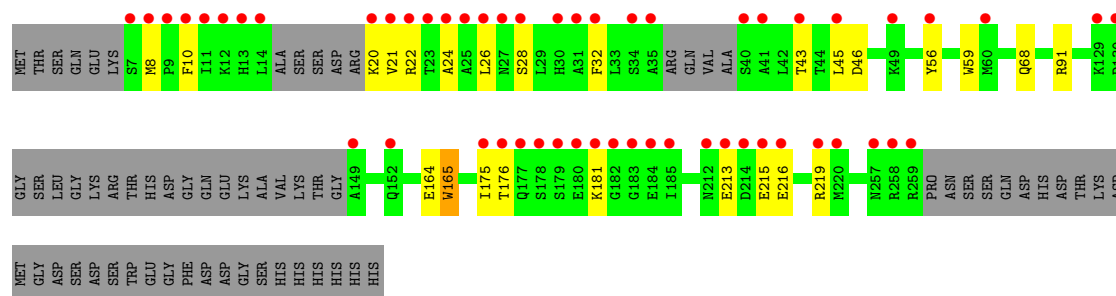
3 Residue-property plots [i](#)

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: G0S4M2



• Molecule 1: G0S4M2



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	214.26Å 214.26Å 124.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.22 – 2.24 49.22 – 2.24	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.22-2.24) 100.0 (49.22-2.24)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.187 , 0.207 (Not available) , 0.193	Depositor DCC
R_{free} test set	3988 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3966	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.54% 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: PO4, EDO, EPE

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	0.16	0/1910	0.40	1/2581 (0.0%)
1	B	0.16	0/1874	0.41	0/2532
All	All	0.16	0/3784	0.40	1/5113 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	ASP	N-CA-C	-7.29	102.47	112.03

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	1875	0	1905	17	0
1	B	1839	0	1865	33	0
2	A	10	0	0	0	0
2	B	10	0	0	1	0
3	A	28	0	42	0	0
4	A	15	0	17	3	0
5	A	141	0	0	4	0
5	B	48	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3966	0	3829	49	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 6.

All (49) 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:B:21:VAL:O	1:B:22:ARG:NH1	1.73	1.19
1:B:21:VAL:C	1:B:22:ARG:NH1	2.33	0.86
1:A:83:ARG:HH22	1:A:131:GLY:HA3	1.41	0.83
1:B:22:ARG:HH11	1:B:22:ARG:HA	1.49	0.78
1:B:21:VAL:HG13	1:B:22:ARG:HH12	1.49	0.78
1:B:22:ARG:HH11	1:B:22:ARG:CA	1.97	0.77
1:A:171:GLU:OE1	5:A:401:HOH:O	2.15	0.64
1:B:21:VAL:C	1:B:22:ARG:HH12	2.00	0.61
1:A:216:GLU:OE1	1:A:219:ARG:NH2	2.32	0.60
1:A:19:ARG:NH1	5:A:404:HOH:O	2.34	0.60
1:B:56[B]:TYR:OH	2:B:301:PO4:O3	2.21	0.59
1:B:43:THR:HG23	1:B:46:ASP:H	1.69	0.58
1:A:215:GLU:OE1	5:A:402:HOH:O	2.17	0.57
1:B:22:ARG:NH1	1:B:22:ARG:HB2	2.21	0.55
1:A:6:LYS:NZ	1:A:41:ALA:O	2.41	0.54
1:A:83:ARG:HH22	1:A:131:GLY:CA	2.16	0.54
1:A:83:ARG:CZ	1:A:154:LYS:HE3	2.39	0.52
1:B:216:GLU:HG2	1:B:219:ARG:NH2	2.25	0.52
1:A:83:ARG:NH2	1:A:131:GLY:HA3	2.21	0.51
1:A:212:ASN:HD22	1:A:215:GLU:HG2	1.75	0.51
1:B:165:TRP:CH2	1:B:175:ILE:HD11	2.46	0.51
1:B:165:TRP:HH2	1:B:175:ILE:HD11	1.76	0.50
1:A:64:ALA:N	5:A:406:HOH:O	2.45	0.49
1:B:8:MET:HG3	1:B:32:PHE:HE2	1.77	0.49
1:A:212:ASN:ND2	1:A:215:GLU:HG2	2.28	0.49
1:B:22:ARG:NH1	1:B:22:ARG:CB	2.76	0.49
1:B:20:LYS:O	1:B:21:VAL:HG12	2.14	0.48
1:A:239:PRO:HD2	4:A:310:EPE:H31	1.97	0.47
1:B:22:ARG:NH1	1:B:22:ARG:CA	2.72	0.47
1:B:56[B]:TYR:HA	1:B:59:TRP:HB3	1.96	0.47
1:B:213:GLU:HB2	1:B:219:ARG:HD3	1.96	0.47
1:B:32:PHE:CD1	1:B:32:PHE:C	2.92	0.47
1:B:32:PHE:C	1:B:32:PHE:HD1	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:LEU:CD2	1:B:181:LYS:HB2	2.46	0.46
1:B:43:THR:HG22	1:B:46:ASP:CG	2.40	0.46
1:B:22:ARG:HB2	1:B:22:ARG:CZ	2.46	0.46
1:A:240:ALA:H	4:A:310:EPE:H32	1.80	0.46
1:A:238:SER:CB	4:A:310:EPE:H22	2.46	0.46
1:A:213:GLU:O	1:A:219:ARG:HD3	2.17	0.44
1:B:22:ARG:HG3	1:B:26:LEU:HD12	1.99	0.44
1:B:175:ILE:HG22	1:B:176:THR:HG23	2.00	0.44
1:B:215:GLU:H	1:B:215:GLU:HG2	1.52	0.43
1:B:24:ALA:O	1:B:28:SER:N	2.39	0.43
1:B:21:VAL:C	1:B:22:ARG:HH11	2.11	0.43
1:B:10:PHE:CZ	1:B:32:PHE:HD2	2.37	0.42
1:B:10:PHE:CE2	1:B:32:PHE:HD2	2.37	0.42
1:B:91:ARG:NH1	1:B:164:GLU:OE2	2.53	0.41
1:B:22:ARG:HG3	1:B:26:LEU:CD1	2.51	0.41
1:A:224:ARG:HD3	1:B:68:GLN:HB2	2.02	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	227/290 (78%)	225 (99%)	1 (0%)	1 (0%)	30	30
1	B	219/290 (76%)	213 (97%)	5 (2%)	1 (0%)	24	23
All	All	446/580 (77%)	438 (98%)	6 (1%)	2 (0%)	24	30

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	TRP
1	B	165	TRP

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	201/251 (80%)	201 (100%)	0	100	100
1	B	198/251 (79%)	198 (100%)	0	100	100
All	All	399/502 (80%)	399 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	13	HIS
1	A	187	GLN
1	A	212	ASN
1	B	30	HIS
1	B	187	GLN
1	B	223	GLN
1	B	234	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

12 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	303	-	3,3,3	0.42	0	2,2,2	0.40	0
2	PO4	A	301	-	4,4,4	0.96	0	6,6,6	0.44	0
2	PO4	B	301	-	4,4,4	0.97	0	6,6,6	0.46	0
4	EPE	A	310	-	15,15,15	0.86	1 (6%)	19,20,20	1.64	4 (21%)
3	EDO	A	308	-	3,3,3	0.43	0	2,2,2	0.34	0
3	EDO	A	307	-	3,3,3	0.44	0	2,2,2	0.34	0
3	EDO	A	309	-	3,3,3	0.43	0	2,2,2	0.34	0
3	EDO	A	304	-	3,3,3	0.43	0	2,2,2	0.33	0
3	EDO	A	306	-	3,3,3	0.41	0	2,2,2	0.42	0
2	PO4	A	302	-	4,4,4	0.98	0	6,6,6	0.51	0
2	PO4	B	302	-	4,4,4	1.05	0	6,6,6	0.41	0
3	EDO	A	305	-	3,3,3	0.43	0	2,2,2	0.38	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	303	-	-	0/1/1/1	-
4	EPE	A	310	-	-	1/9/19/19	0/1/1/1
3	EDO	A	308	-	-	1/1/1/1	-
3	EDO	A	307	-	-	1/1/1/1	-
3	EDO	A	309	-	-	0/1/1/1	-
3	EDO	A	304	-	-	0/1/1/1	-
3	EDO	A	306	-	-	0/1/1/1	-
3	EDO	A	305	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	310	EPE	C10-S	2.82	1.81	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	310	EPE	C5-N4-C3	3.51	116.41	108.84
4	A	310	EPE	C7-N4-C3	3.24	119.87	111.24
4	A	310	EPE	C7-N4-C5	2.54	118.01	111.24
4	A	310	EPE	O3S-S-C10	2.14	110.19	106.00

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	310	EPE	C8-C7-N4-C5
3	A	307	EDO	O1-C1-C2-O2
3	A	308	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	PO4	1	0
4	A	310	EPE	3	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	232/290 (80%)	0.29	24 (10%)	12 11	32, 49, 91, 146	1 (0%)
1	B	226/290 (77%)	1.29	54 (23%)	2 1	33, 68, 167, 201	1 (0%)
All	All	458/580 (78%)	0.79	78 (17%)	4 3	32, 58, 148, 201	2 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	VAL	10.7
1	A	176	THR	9.7
1	B	180	GLU	8.6
1	B	259	ARG	8.5
1	B	41	ALA	8.5
1	B	40	SER	8.4
1	B	23	THR	8.3
1	B	20	LYS	8.2
1	B	35	ALA	8.2
1	A	183	GLY	7.3
1	A	131	GLY	7.2
1	B	130	ASP	7.1
1	A	148	GLY	7.1
1	B	25	ALA	6.8
1	B	9	PRO	6.5
1	B	13	HIS	5.9
1	A	130	ASP	5.8
1	B	182	GLY	5.7
1	B	56[A]	TYR	5.6
1	B	179	SER	5.5
1	B	177	GLN	5.5
1	B	213	GLU	5.5
1	B	12	LYS	5.4
1	B	24	ALA	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	149	ALA	5.4
1	B	181	LYS	5.4
1	B	7	SER	5.1
1	B	45	LEU	5.0
1	B	14	LEU	5.0
1	B	183	GLY	4.9
1	A	258	ARG	4.8
1	B	32	PHE	4.8
1	B	129	LYS	4.7
1	B	8	MET	4.7
1	B	26	LEU	4.6
1	B	212	ASN	4.4
1	B	215	GLU	4.3
1	A	212	ASN	3.9
1	B	214	ASP	3.9
1	B	10	PHE	3.9
1	A	199	ASP	3.8
1	B	220	MET	3.8
1	B	22	ARG	3.8
1	B	34	SER	3.7
1	A	184	GLU	3.6
1	A	213	GLU	3.6
1	B	176	THR	3.4
1	B	60	MET	3.3
1	B	49	LYS	3.3
1	B	152	GLN	3.3
1	A	63	ARG	3.2
1	A	215	GLU	3.2
1	B	30	HIS	3.1
1	A	45	LEU	3.1
1	A	175	ILE	2.9
1	B	43	THR	2.9
1	B	184	GLU	2.9
1	B	27	ASN	2.8
1	B	178	SER	2.8
1	A	83	ARG	2.8
1	B	258	ARG	2.8
1	A	62	ASP	2.8
1	A	252	ASP	2.7
1	B	175	ILE	2.7
1	B	11	ILE	2.6
1	A	129	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	171	GLU	2.5
1	B	185	ILE	2.5
1	A	174	ARG	2.4
1	B	31	ALA	2.4
1	A	214	ASP	2.4
1	B	257	ASN	2.3
1	B	28	SER	2.3
1	B	219	ARG	2.3
1	A	5	GLU	2.2
1	A	84	GLU	2.1
1	B	216	GLU	2.0
1	A	19	ARG	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	A	305	4/4	0.75	0.40	123,123,126,127	0
3	EDO	A	307	4/4	0.77	0.32	89,89,90,92	0
2	PO4	B	301	5/5	0.85	0.13	129,129,133,138	0
3	EDO	A	308	4/4	0.85	0.34	89,95,98,104	0
4	EPE	A	310	15/15	0.86	0.25	60,67,96,102	15
3	EDO	A	309	4/4	0.87	0.23	85,87,91,91	0
3	EDO	A	304	4/4	0.87	0.26	83,89,90,93	0
2	PO4	A	301	5/5	0.88	0.12	87,97,109,113	0
2	PO4	A	302	5/5	0.96	0.09	70,77,83,85	0
3	EDO	A	303	4/4	0.97	0.15	56,64,66,75	0
3	EDO	A	306	4/4	0.97	0.11	51,52,53,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	B	302	5/5	0.98	0.08	67,67,73,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.