



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:23 PM UTC

PDB ID : 7Y8X / pdb_00007y8x
Title : Crystal structure of AlbEF homolog from Quasibacillus thermotolerans in complex with Ni(II)
Authors : Ishida, K.; Nakamura, A.; Kojima, S.
Deposited on : 2022-06-24
Resolution : 2.25 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

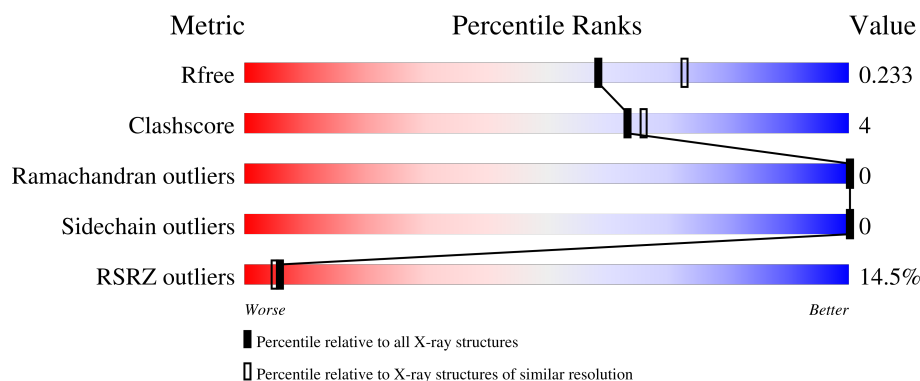
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	E	395	16% 78% 15% 6%
2	F	366	11% 88% 7% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6059 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 AlbE homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	370	Total	C	N	O	S	0	0	0
			3021	1939	505	566	11			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-13	MET	-	initiating methionine	UNP A0A837GIQ1
E	-12	GLY	-	expression tag	UNP A0A837GIQ1
E	-11	SER	-	expression tag	UNP A0A837GIQ1
E	-10	SER	-	expression tag	UNP A0A837GIQ1
E	-9	HIS	-	expression tag	UNP A0A837GIQ1
E	-8	HIS	-	expression tag	UNP A0A837GIQ1
E	-7	HIS	-	expression tag	UNP A0A837GIQ1
E	-6	HIS	-	expression tag	UNP A0A837GIQ1
E	-5	HIS	-	expression tag	UNP A0A837GIQ1
E	-4	HIS	-	expression tag	UNP A0A837GIQ1
E	-3	SER	-	expression tag	UNP A0A837GIQ1
E	-2	GLN	-	expression tag	UNP A0A837GIQ1
E	-1	ASP	-	expression tag	UNP A0A837GIQ1
E	0	PRO	-	expression tag	UNP A0A837GIQ1

- Molecule 2 is a protein called AlbF homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	347	Total	C	N	O	S	0	3	0
			2879	1846	484	538	11			

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	1	Total	Ni	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		

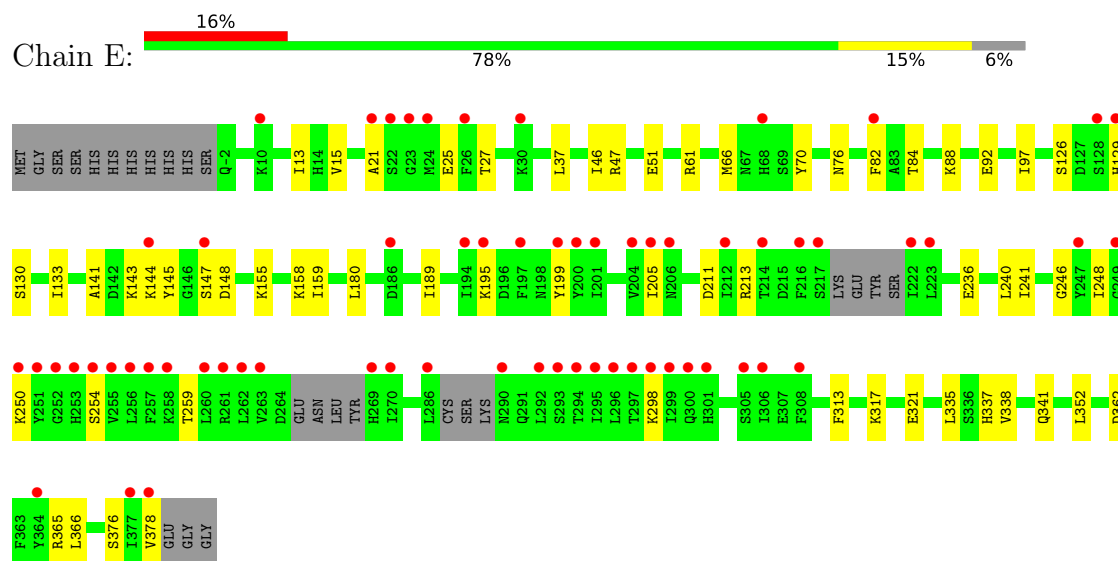
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	E	10	Total	O	0	0
			10	10		
6	F	70	Total	O	0	0
			70	70		

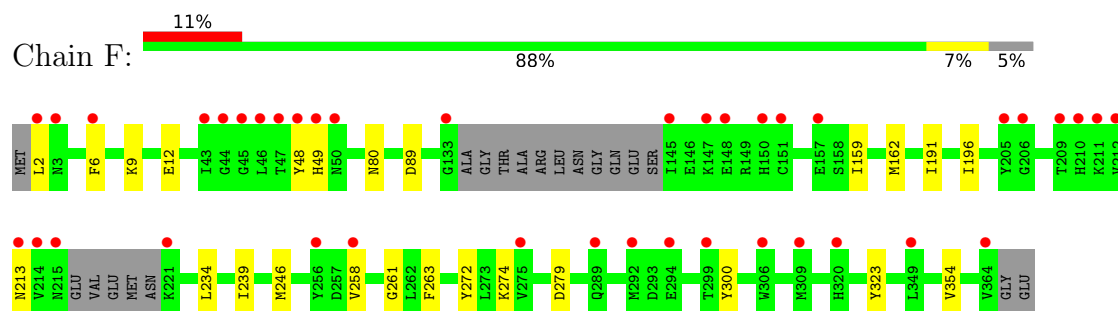
3 Residue-property plots [i](#)

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: AlbE homolog



• Molecule 2: AlbF homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.38Å 136.38Å 118.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.71 – 2.25 44.71 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.71-2.25) 99.8 (44.71-2.25)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.85 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.200 , 0.233 0.200 , 0.233	Depositor DCC
R_{free} test set	2882 reflections (3.83%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6059	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: EDO, NI, PO4

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	E	0.09	0/3082	0.24	0/4153
2	F	0.09	0/2946	0.24	0/3977
All	All	0.09	0/6028	0.24	0/8130

There are no bond length outliers.

There are no bond angle outliers.

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	E	3021	0	3005	34	0
2	F	2879	0	2850	15	0
3	E	20	0	0	1	0
3	F	30	0	0	0	0
4	F	1	0	0	0	0
5	F	28	0	42	2	0
6	E	10	0	0	0	0
6	F	70	0	0	1	0
All	All	6059	0	5897	47	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 (47) 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)
2:F:49:HIS:NE2	2:F:213:ASN:O	2.21	0.73
1:E:236:GLU:HA	1:E:240:LEU:HB2	1.72	0.71
1:E:246:GLY:HA2	1:E:250:LYS:HB3	1.71	0.71
1:E:213:ARG:HD3	1:E:378:VAL:HG22	1.79	0.64
1:E:241:ILE:HG12	1:E:366:LEU:HD21	1.84	0.60
2:F:234:LEU:HD12	5:F:408:EDO:H21	1.83	0.60
1:E:236:GLU:OE2	1:E:365:ARG:NH1	2.35	0.59
1:E:126:SER:O	1:E:130:SER:N	2.35	0.58
2:F:239:ILE:HG21	2:F:354:VAL:HG21	1.85	0.57
1:E:84:THR:HA	2:F:2:LEU:HD13	1.88	0.55
1:E:143:LYS:HG2	1:E:147:SER:HB2	1.89	0.55
1:E:259:THR:HG23	1:E:298:LYS:HE3	1.89	0.55
1:E:321:GLU:HG3	1:E:352:LEU:HD22	1.88	0.55
1:E:337:HIS:CE1	1:E:341:GLN:HG3	2.43	0.54
2:F:263:PHE:HB3	2:F:272:TYR:HB2	1.90	0.53
1:E:155:LYS:HA	1:E:158:LYS:HE2	1.91	0.52
1:E:211:ASP:HB3	1:E:376:SER:HA	1.90	0.52
2:F:246:MET:HE2	2:F:300:TYR:HB3	1.90	0.52
1:E:46:ILE:HD13	1:E:97:ILE:HG12	1.91	0.51
2:F:89:ASP:OD2	2:F:323:TYR:OH	2.24	0.50
1:E:144:LYS:HA	1:E:148:ASP:HB2	1.94	0.49
1:E:21:ALA:HA	1:E:70:TYR:CD1	2.49	0.48
1:E:25:GLU:OE1	1:E:27:THR:OG1	2.30	0.48
2:F:261:GLY:N	2:F:274:LYS:O	2.44	0.47
1:E:61:ARG:NE	3:E:402:PO4:O2	2.47	0.47
1:E:180:LEU:HD11	1:E:195:LYS:HB3	1.96	0.47
1:E:13:ILE:HD12	1:E:76:ASN:HB3	1.97	0.46
2:F:159:ILE:HA	2:F:162:MET:HE2	1.97	0.46
1:E:82:PHE:HB3	2:F:6:PHE:HB2	1.98	0.45
2:F:191:ILE:HG23	2:F:196:ILE:HD13	1.98	0.45
1:E:37:LEU:HD23	1:E:159:ILE:HG21	1.99	0.44
1:E:129:HIS:O	1:E:133:ILE:HG12	2.17	0.44
1:E:25:GLU:OE2	1:E:145:TYR:OH	2.34	0.44
2:F:48:TYR:O	6:F:501:HOH:O	2.21	0.44
1:E:189:ILE:HD13	1:E:338:VAL:HG21	2.00	0.44
2:F:258:VAL:HG13	2:F:279:ASP:HB3	1.99	0.44
1:E:248:ILE:C	1:E:254:SER:HB2	2.44	0.43
1:E:313:PHE:CE2	1:E:317:LYS:HD2	2.54	0.43
1:E:88:LYS:HB3	1:E:92:GLU:HB2	2.01	0.43
2:F:9:LYS:O	2:F:12:GLU:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:ALA:HB3	1:E:205:ILE:HD12	2.00	0.42
1:E:15:VAL:HG11	1:E:335:LEU:HD22	2.02	0.42
1:E:66:MET:HG3	1:E:199:TYR:CD1	2.55	0.42
1:E:47:ARG:HG3	1:E:51:GLU:CD	2.45	0.41
1:E:21:ALA:HA	1:E:70:TYR:HD1	1.85	0.41
1:E:240:LEU:HD13	1:E:362:ASP:HB3	2.01	0.41
2:F:80[B]:ASN:HA	5:F:412:EDO:H11	2.02	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	E	362/395 (92%)	348 (96%)	14 (4%)	0	100	100
2	F	344/366 (94%)	331 (96%)	13 (4%)	0	100	100
All	All	706/761 (93%)	679 (96%)	27 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	E	340/362 (94%)	340 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	321/332 (97%)	321 (100%)	0	100	100
All	All	661/694 (95%)	661 (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	E	28	GLN
1	E	95	GLN
1	E	121	GLN
1	E	172	ASN
1	E	291	GLN
2	F	23	ASN
2	F	179	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 18 ligands modelled in this entry, 1 is monoatomic - leaving 17 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
3	PO4	F	406	-	4,4,4	0.95	0	6,6,6	0.46	0
5	EDO	F	408	-	3,3,3	0.44	0	2,2,2	0.37	0
3	PO4	E	403	-	4,4,4	0.94	0	6,6,6	0.48	0
3	PO4	F	401	4	4,4,4	0.96	0	6,6,6	0.45	0
5	EDO	F	410	-	3,3,3	0.44	0	2,2,2	0.31	0
3	PO4	F	402	-	4,4,4	0.94	0	6,6,6	0.52	0
5	EDO	F	411	-	3,3,3	0.45	0	2,2,2	0.28	0
5	EDO	F	414	-	3,3,3	0.42	0	2,2,2	0.36	0
5	EDO	F	413	-	3,3,3	0.45	0	2,2,2	0.33	0
3	PO4	E	404	-	4,4,4	0.97	0	6,6,6	0.47	0
3	PO4	E	402	-	4,4,4	0.95	0	6,6,6	0.46	0
3	PO4	E	401	-	4,4,4	0.95	0	6,6,6	0.49	0
5	EDO	F	409	-	3,3,3	0.44	0	2,2,2	0.35	0
5	EDO	F	412	-	3,3,3	0.44	0	2,2,2	0.37	0
3	PO4	F	405	-	4,4,4	0.92	0	6,6,6	0.47	0
3	PO4	F	403	-	4,4,4	0.96	0	6,6,6	0.46	0
3	PO4	F	404	-	4,4,4	0.97	0	6,6,6	0.43	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
5	EDO	F	408	-	-	1/1/1/1	-
5	EDO	F	411	-	-	0/1/1/1	-
5	EDO	F	410	-	-	0/1/1/1	-
5	EDO	F	414	-	-	0/1/1/1	-
5	EDO	F	413	-	-	0/1/1/1	-
5	EDO	F	409	-	-	0/1/1/1	-
5	EDO	F	412	-	-	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
5	F	408	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	408	EDO	1	0
3	E	402	PO4	1	0
5	F	412	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	E	370/395 (93%)	1.04	64 (17%)	4 3	41, 79, 151, 198	0
2	F	347/366 (94%)	0.56	40 (11%)	9 9	29, 58, 120, 158	4 (1%)
All	All	717/761 (94%)	0.81	104 (14%)	6 5	29, 69, 147, 198	4 (0%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	197	PHE	8.3
2	F	48	TYR	7.4
2	F	145	ILE	6.6
1	E	292	LEU	6.6
1	E	212	ILE	5.9
2	F	2	LEU	5.9
2	F	215	ASN	5.8
2	F	50	ASN	5.6
2	F	256	TYR	5.6
2	F	221	LYS	5.0
1	E	251	TYR	4.9
1	E	21	ALA	4.8
1	E	378	VAL	4.8
2	F	364	VAL	4.8
1	E	262	LEU	4.8
2	F	211	LYS	4.8
2	F	147	LYS	4.4
1	E	253	HIS	4.3
2	F	210	HIS	4.2
1	E	22	SER	4.2
2	F	214	VAL	4.2
2	F	45	GLY	4.1
1	E	216	PHE	4.1
1	E	286	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	E	364	TYR	3.9
1	E	296	LEU	3.8
2	F	49	HIS	3.8
1	E	257	PHE	3.8
1	E	23	GLY	3.8
2	F	46	LEU	3.8
2	F	306	TRP	3.7
2	F	47	THR	3.6
1	E	255	VAL	3.6
2	F	209	THR	3.6
1	E	269	HIS	3.4
1	E	205	ILE	3.4
1	E	214	THR	3.3
1	E	194	ILE	3.3
1	E	299	ILE	3.3
1	E	217	SER	3.3
1	E	263	VAL	3.3
1	E	294	THR	3.2
1	E	295	ILE	3.2
1	E	252	GLY	3.2
1	E	222	ILE	3.1
2	F	213	ASN	3.1
1	E	250	LYS	3.1
1	E	270	ILE	3.0
1	E	129	HIS	3.0
1	E	308	PHE	3.0
2	F	258	VAL	3.0
1	E	128	SER	3.0
1	E	30	LYS	2.9
1	E	195	LYS	2.9
1	E	26	PHE	2.9
1	E	260	LEU	2.9
2	F	275	VAL	2.9
1	E	199	TYR	2.9
1	E	256	LEU	2.9
1	E	200	TYR	2.9
2	F	3	ASN	2.9
2	F	150	HIS	2.8
1	E	24	MET	2.8
1	E	206	ASN	2.8
1	E	82	PHE	2.8
2	F	44	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	298	LYS	2.7
2	F	320[A]	HIS	2.7
2	F	212	VAL	2.7
1	E	290	ASN	2.7
2	F	148	GLU	2.7
1	E	297	THR	2.6
2	F	151	CYS	2.6
1	E	144	LYS	2.6
2	F	309	MET	2.6
1	E	204	VAL	2.6
1	E	10	LYS	2.6
2	F	157	GLU	2.5
2	F	43	ILE	2.5
2	F	289	GLN	2.5
1	E	223	LEU	2.5
1	E	249	GLY	2.5
1	E	301	HIS	2.5
1	E	300	GLN	2.4
2	F	206	GLY	2.3
1	E	147	SER	2.3
2	F	133	GLY	2.3
1	E	306	ILE	2.3
1	E	377	ILE	2.3
1	E	186	ASP	2.2
2	F	292	MET	2.2
1	E	293	SER	2.2
1	E	305	SER	2.2
2	F	294	GLU	2.1
1	E	68	HIS	2.1
2	F	205	TYR	2.1
1	E	201	ILE	2.1
2	F	349	LEU	2.1
1	E	258	LYS	2.1
2	F	6	PHE	2.1
1	E	254	SER	2.0
2	F	299	THR	2.0
1	E	247	TYR	2.0
1	E	261	ARG	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 [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	PO4	F	404	5/5	0.60	0.16	84,97,115,123	0
5	EDO	F	412	4/4	0.68	0.24	66,74,80,83	0
5	EDO	F	408	4/4	0.76	0.25	68,70,78,83	0
3	PO4	E	403	5/5	0.77	0.13	88,89,104,105	0
3	PO4	F	406	5/5	0.79	0.12	58,90,96,110	0
3	PO4	E	404	5/5	0.80	0.15	89,92,101,140	0
5	EDO	F	411	4/4	0.80	0.20	69,77,77,77	0
3	PO4	F	403	5/5	0.80	0.11	71,88,99,100	0
5	EDO	F	409	4/4	0.85	0.20	65,66,67,67	0
5	EDO	F	413	4/4	0.85	0.22	55,69,78,84	0
3	PO4	E	402	5/5	0.88	0.08	73,80,99,102	0
3	PO4	F	401	5/5	0.89	0.09	62,79,82,97	0
3	PO4	F	402	5/5	0.90	0.12	68,68,88,104	0
3	PO4	F	405	5/5	0.93	0.12	51,61,76,78	0
5	EDO	F	414	4/4	0.93	0.19	64,64,72,74	0
5	EDO	F	410	4/4	0.94	0.12	57,62,67,70	0
4	NI	F	407	1/1	0.96	0.13	123,123,123,123	0
3	PO4	E	401	5/5	0.96	0.06	63,70,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.