



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 21, 2016 – 08:40 PM EDT

PDB ID : 4XGW
Title : Crystal structure of Escherichia coli Flavin trafficking protein, an FMN transferase, E169K mutant
Authors : Tomchick, D.R.; Brautigam, C.A.; Deka, R.K.; Norgard, M.V.
Deposited on : 2015-01-02
Resolution : 1.75 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

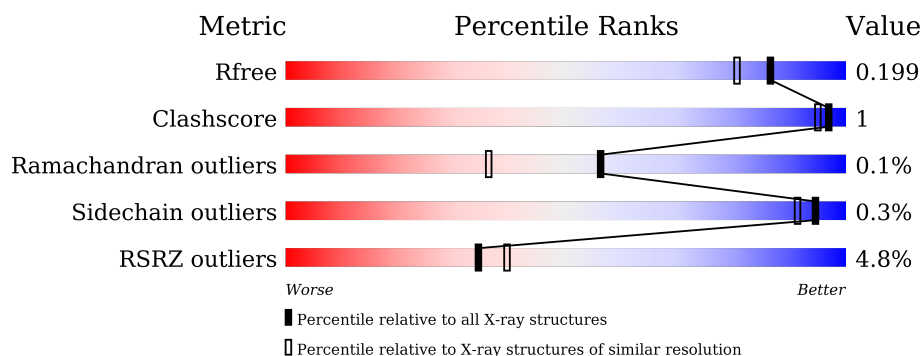
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	91344	2417 (1.76-1.72)
Clashscore	102246	2570 (1.76-1.72)
Ramachandran outliers	100387	2544 (1.76-1.72)
Sidechain outliers	100360	2544 (1.76-1.72)
RSRZ outliers	91569	2420 (1.76-1.72)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	340	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>•</div> <div>11%</div> </div> </div>
1	B	340	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>•</div> <div>9%</div> </div> </div>
1	C	340	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>•</div> <div>10%</div> </div> </div>
1	D	340	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>•</div> <div>11%</div> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NA	A	502	-	-	-	X
3	NA	C	502	-	-	-	X
3	NA	D	502	-	-	-	X
4	EDO	D	503	-	-	-	X
4	EDO	D	504	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19641 atoms, of which 9481 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 FAD:protein FMN transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	304	Total	C	H	N	O	S	0	3	0
			4704	1484	2347	404	459	10			
1	B	308	Total	C	H	N	O	S	0	0	0
			4757	1501	2377	409	460	10			
1	C	307	Total	C	H	N	O	S	0	2	0
			4766	1501	2386	409	460	10			
1	D	303	Total	C	H	N	O	S	0	3	0
			4688	1473	2347	404	454	10			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AB85
A	169	LYS	GLU	engineered mutation	UNP P0AB85
A	333	LEU	-	expression tag	UNP P0AB85
A	334	GLU	-	expression tag	UNP P0AB85
A	335	HIS	-	expression tag	UNP P0AB85
A	336	HIS	-	expression tag	UNP P0AB85
A	337	HIS	-	expression tag	UNP P0AB85
A	338	HIS	-	expression tag	UNP P0AB85
A	339	HIS	-	expression tag	UNP P0AB85
A	340	HIS	-	expression tag	UNP P0AB85
B	1	MET	-	initiating methionine	UNP P0AB85
B	169	LYS	GLU	engineered mutation	UNP P0AB85
B	333	LEU	-	expression tag	UNP P0AB85
B	334	GLU	-	expression tag	UNP P0AB85
B	335	HIS	-	expression tag	UNP P0AB85
B	336	HIS	-	expression tag	UNP P0AB85
B	337	HIS	-	expression tag	UNP P0AB85
B	338	HIS	-	expression tag	UNP P0AB85
B	339	HIS	-	expression tag	UNP P0AB85
B	340	HIS	-	expression tag	UNP P0AB85
C	1	MET	-	initiating methionine	UNP P0AB85

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	169	LYS	GLU	engineered mutation	UNP P0AB85
C	333	LEU	-	expression tag	UNP P0AB85
C	334	GLU	-	expression tag	UNP P0AB85
C	335	HIS	-	expression tag	UNP P0AB85
C	336	HIS	-	expression tag	UNP P0AB85
C	337	HIS	-	expression tag	UNP P0AB85
C	338	HIS	-	expression tag	UNP P0AB85
C	339	HIS	-	expression tag	UNP P0AB85
C	340	HIS	-	expression tag	UNP P0AB85
D	1	MET	-	initiating methionine	UNP P0AB85
D	169	LYS	GLU	engineered mutation	UNP P0AB85
D	333	LEU	-	expression tag	UNP P0AB85
D	334	GLU	-	expression tag	UNP P0AB85
D	335	HIS	-	expression tag	UNP P0AB85
D	336	HIS	-	expression tag	UNP P0AB85
D	337	HIS	-	expression tag	UNP P0AB85
D	338	HIS	-	expression tag	UNP P0AB85
D	339	HIS	-	expression tag	UNP P0AB85
D	340	HIS	-	expression tag	UNP P0AB85

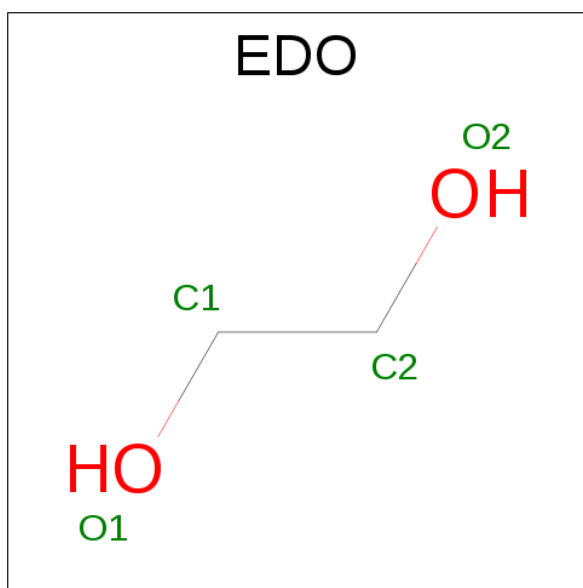
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0
3	D	1	Total Na 1 1	0	0
3	C	1	Total Na 1 1	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	0
			10	2	6	2		
4	C	1	Total	C	H	O	0	0
			10	2	6	2		
4	D	1	Total	C	H	O	0	0
			10	2	6	2		
4	D	1	Total	C	H	O	0	0
			10	2	6	2		

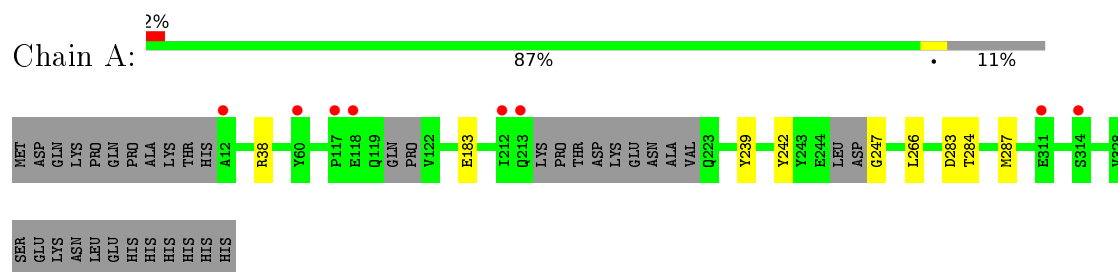
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	181	Total	O	0	0
			181	181		
5	B	161	Total	O	0	0
			161	161		
5	C	147	Total	O	0	0
			147	147		
5	D	190	Total	O	0	0
			190	190		

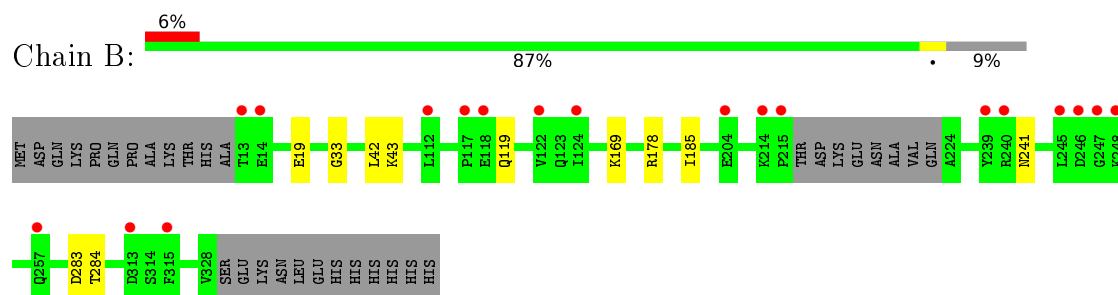
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

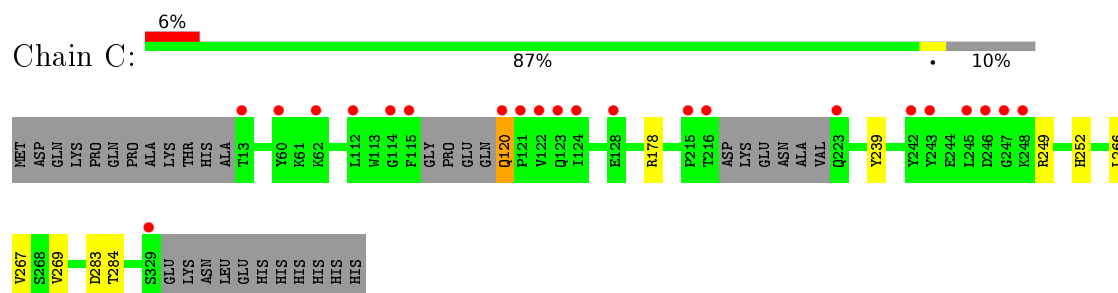
- Molecule 1: FAD:protein FMN transferase



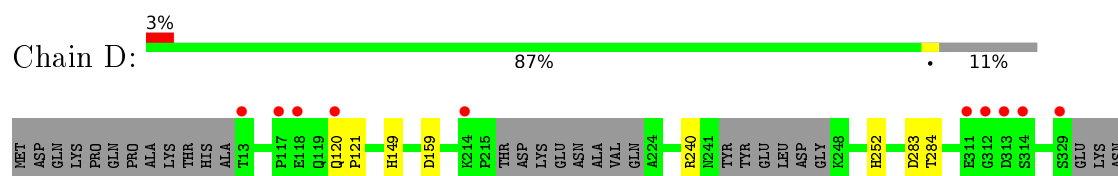
- Molecule 1: FAD:protein FMN transferase



- Molecule 1: FAD:protein FMN transferase



- Molecule 1: FAD:protein FMN transferase



LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.22Å 70.65Å 86.28Å 75.72° 71.91° 69.72°	Depositor
Resolution (Å)	33.92 – 1.75 33.92 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.5 (33.92-1.75) 91.7 (33.92-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 1.75Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.172 , 0.199 0.172 , 0.199	Depositor DCC
R_{free} test set	5874 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 41.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 116591 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19641	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, CA, 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	0.24	0/2411	0.46	0/3264
1	B	0.24	0/2426	0.45	0/3289
1	C	0.24	0/2432	0.46	0/3294
1	D	0.24	0/2396	0.46	0/3247
All	All	0.24	0/9665	0.46	0/13094

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 [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	2357	2347	2340	6	0
1	B	2380	2377	2376	7	0
1	C	2380	2386	2381	6	0
1	D	2341	2347	2340	6	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	C	8	12	12	0	0
4	D	8	12	12	0	0
5	A	181	0	0	1	0
5	B	161	0	0	3	0
5	C	147	0	0	2	0
5	D	190	0	0	0	0
All	All	10160	9481	9461	25	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 1.

All (25) 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:249:ARG:O	5:C:601:HOH:O	2.14	0.66
1:A:266:LEU:O	1:A:287:MET:HE3	1.96	0.66
1:B:19:GLU:O	1:B:43:LYS:NZ	2.37	0.58
1:B:178:ARG:NH2	5:B:602:HOH:O	2.38	0.56
1:B:283:ASP:OD1	1:B:284:THR:N	2.40	0.55
1:C:283:ASP:OD1	1:C:284:THR:N	2.44	0.51
1:D:283:ASP:OD1	1:D:284:THR:N	2.45	0.49
1:C:178:ARG:NH2	5:C:605:HOH:O	2.44	0.49
1:A:38:ARG:NH2	1:A:183:GLU:O	2.46	0.48
1:C:239:TYR:HA	1:C:267:VAL:CG1	2.44	0.48
1:D:159:ASP:OD1	1:D:159:ASP:N	2.41	0.48
1:D:120:GLN:HB3	1:D:121:PRO:HD2	1.96	0.48
1:A:283:ASP:OD1	1:A:284:THR:N	2.48	0.47
1:C:120:GLN:HE21	1:C:120:GLN:N	2.12	0.47
1:A:247:GLY:N	5:A:606:HOH:O	2.48	0.45
1:B:119:GLN:NE2	1:B:241:ASN:O	2.50	0.44
1:D:240:ARG:HD2	1:D:252:HIS:CG	2.53	0.43
1:A:239:TYR:HA	1:A:242:TYR:HB2	2.00	0.43
1:B:42:LEU:HD11	1:B:185:ILE:HD11	2.00	0.43
1:B:178:ARG:CZ	5:B:602:HOH:O	2.69	0.41
1:B:169:LYS:HE3	5:B:708:HOH:O	2.20	0.41
1:D:240:ARG:HD2	1:D:252:HIS:CD2	2.55	0.41
1:C:266:LEU:HD21	1:C:269:VAL:CG2	2.51	0.41
1:A:38:ARG:NH2	1:A:183:GLU:HB3	2.36	0.40
1:D:120:GLN:HB3	1:D:121:PRO:CD	2.51	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	299/340 (88%)	292 (98%)	7 (2%)	0	100	100
1	B	304/340 (89%)	297 (98%)	6 (2%)	1 (0%)	46	26
1	C	303/340 (89%)	300 (99%)	3 (1%)	0	100	100
1	D	300/340 (88%)	293 (98%)	7 (2%)	0	100	100
All	All	1206/1360 (89%)	1182 (98%)	23 (2%)	1 (0%)	56	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	33	GLY

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	257/288 (89%)	257 (100%)	0	100	100
1	B	259/288 (90%)	259 (100%)	0	100	100
1	C	261/288 (91%)	259 (99%)	2 (1%)	86	77
1	D	258/288 (90%)	257 (100%)	1 (0%)	93	90
All	All	1035/1152 (90%)	1032 (100%)	3 (0%)	94	91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	120	GLN
1	C	252	HIS
1	D	149	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 7 are monoatomic - leaving 4 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	$\# Z > 2$
4	EDO	C	503	-	3,3,3	0.44	0	2,2,2	0.44	0
4	EDO	C	504	-	3,3,3	0.44	0	2,2,2	0.36	0
4	EDO	D	503	-	3,3,3	0.46	0	2,2,2	0.38	0
4	EDO	D	504	-	3,3,3	0.44	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
4	EDO	C	503	-	-	0/1/1/1	0/0/0/0
4	EDO	C	504	-	-	0/1/1/1	0/0/0/0
4	EDO	D	503	-	-	0/1/1/1	0/0/0/0
4	EDO	D	504	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	304/340 (89%)	-0.06	8 (2%) 59 66	13, 23, 55, 91	0
1	B	308/340 (90%)	0.23	19 (6%) 24 28	16, 26, 63, 85	0
1	C	307/340 (90%)	0.29	22 (7%) 18 22	15, 26, 67, 89	0
1	D	303/340 (89%)	-0.04	10 (3%) 50 56	13, 21, 57, 90	0
All	All	1222/1360 (89%)	0.10	59 (4%) 34 39	13, 24, 61, 91	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	122	VAL	9.2
1	B	245	LEU	6.9
1	C	60	TYR	6.6
1	C	13	THR	6.4
1	C	124	ILE	6.0
1	B	246	ASP	5.6
1	B	247	GLY	5.6
1	C	114	GLY	5.2
1	D	311	GLU	5.2
1	C	247	GLY	5.1
1	C	120	GLN	4.6
1	D	313	ASP	4.5
1	C	216	THR	4.4
1	C	115	PHE	4.4
1	C	245	LEU	4.3
1	B	313	ASP	4.1
1	D	120	GLN	4.0
1	B	13	THR	4.0
1	C	62	LYS	4.0
1	B	14	GLU	4.0
1	A	213	GLN	4.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	215	PRO	3.9
1	B	214	LYS	3.9
1	C	243	TYR	3.8
1	B	122	VAL	3.7
1	A	12	ALA	3.6
1	C	121	PRO	3.6
1	A	60	TYR	3.5
1	C	246	ASP	3.5
1	C	123	GLN	3.5
1	C	242	TYR	3.5
1	C	248	LYS	3.4
1	A	311	GLU	3.3
1	D	118	GLU	3.3
1	A	118	GLU	3.2
1	D	329	SER	3.2
1	B	117	PRO	3.2
1	C	329	SER	3.2
1	B	248	LYS	3.1
1	D	314	SER	3.1
1	B	124	ILE	3.0
1	C	128	GLU	2.8
1	A	212	ILE	2.8
1	C	112	LEU	2.7
1	D	13	THR	2.6
1	A	117	PRO	2.6
1	A	314	SER	2.6
1	D	117	PRO	2.5
1	D	312	GLY	2.5
1	B	118	GLU	2.5
1	B	315	PHE	2.5
1	B	239	TYR	2.4
1	B	240	ARG	2.3
1	B	204	GLU	2.3
1	B	112	LEU	2.2
1	C	215	PRO	2.2
1	D	214	LYS	2.2
1	B	257	GLN	2.1
1	C	223	GLN	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	EDO	D	503	4/4	0.64	0.21	14.23	41,53,63,63	0
4	EDO	D	504	4/4	0.72	0.15	4.33	40,49,53,54	0
3	NA	D	502	1/1	0.91	0.21	4.17	31,31,31,31	0
3	NA	C	502	1/1	0.84	0.32	4.09	40,40,40,40	0
3	NA	A	502	1/1	0.96	0.17	2.49	29,29,29,29	0
4	EDO	C	503	4/4	0.89	0.15	0.44	29,38,46,46	0
4	EDO	C	504	4/4	0.73	0.15	-0.69	63,76,80,84	0
2	CA	C	501	1/1	0.97	0.05	-2.98	26,26,26,26	0
2	CA	B	501	1/1	0.99	0.03	-3.98	27,27,27,27	0
2	CA	A	501	1/1	0.99	0.04	-4.45	24,24,24,24	0
2	CA	D	501	1/1	0.98	0.05	-5.51	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.