



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 08:12 AM UTC

PDB ID : 4KAG / pdb_00004kag
Title : Crystal structure analysis of a single amino acid deletion mutation in EGFP
Authors : Arpino, J.A.J.; Rizkallah, P.J.
Deposited on : 2013-04-22
Resolution : 1.12 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

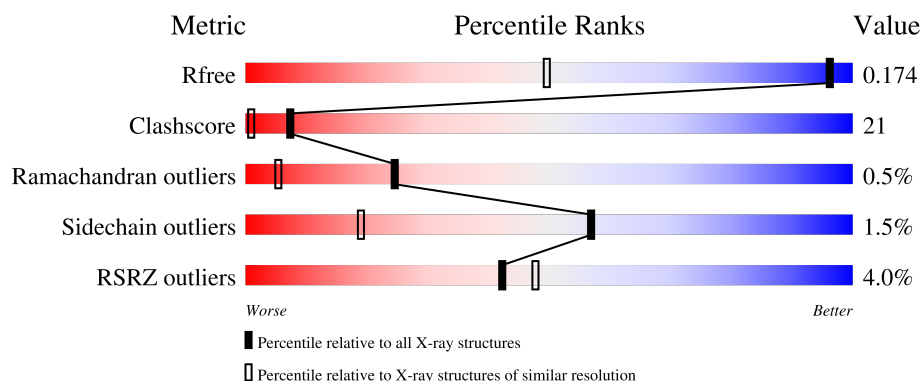
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.12 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	1985 (1.14-1.10)
Clashscore	190562	2021 (1.14-1.10)
Ramachandran outliers	187476	1978 (1.14-1.10)
Sidechain outliers	187428	1974 (1.14-1.10)
RSRZ outliers	180081	1984 (1.14-1.10)

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	236	

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
2	EDO	A	301	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	A	302	-	-	X	-
2	EDO	A	303	-	-	X	-
2	EDO	A	307	-	-	X	-
2	EDO	A	321	-	X	X	-
4	TRS	A	318	-	X	-	-
4	TRS	A	319	-	-	X	-
4	TRS	A	320	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2444 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.

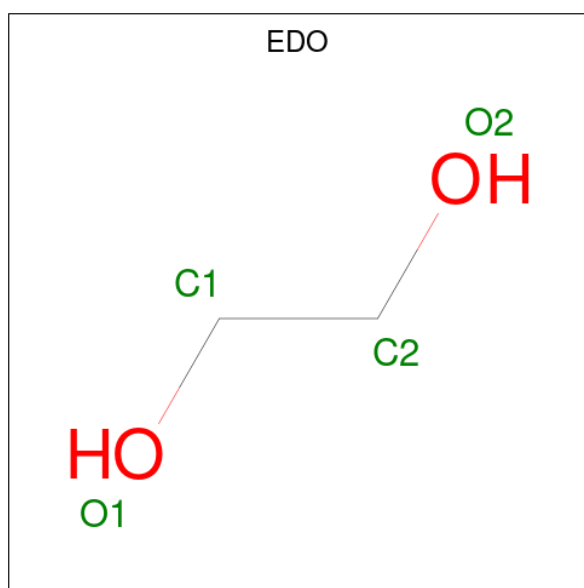
- Molecule 1 is a protein called Green fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	225	2066	1308	348	402	8	0	31	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	VAL	-	insertion	UNP P42212
A	64	LEU	PHE	engineered mutation	UNP P42212
A	66	CRO	SER	chromophore	UNP P42212
A	66	CRO	TYR	chromophore	UNP P42212
A	66	CRO	GLY	chromophore	UNP P42212
A	?	-	ASP	deletion	UNP P42212
A	231	LEU	HIS	conflict	UNP P42212

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



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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		
4	A	1	Total	C	N	O	0	0
			8	4	1	3		

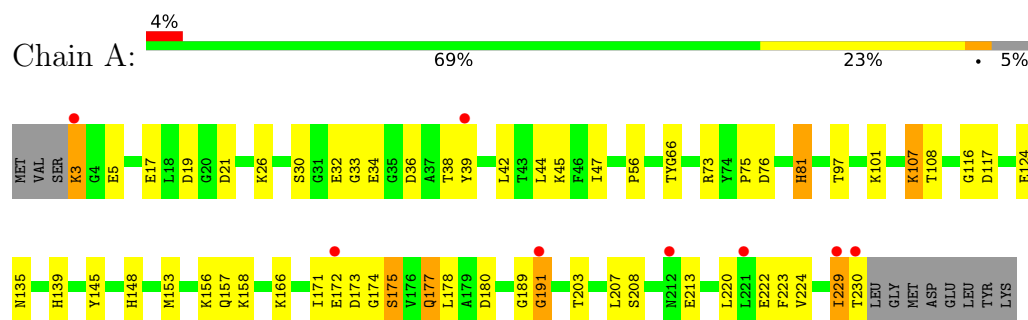
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	280	Total	O	0	0
			280	280		

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: Green fluorescent protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	57.14Å 57.14Å 135.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.94 – 1.12 39.94 – 1.12	Depositor EDS
% Data completeness (in resolution range)	96.1 (39.94-1.12) 96.1 (39.94-1.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 1.12Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.144 , 0.161 0.153 , 0.174	Depositor DCC
R_{free} test set	4723 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2444	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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: SO4, CRO, TRS, 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	1.55	14/2086 (0.7%)	1.17	3/2817 (0.1%)

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	177	GLN	CD-NE2	-7.00	1.18	1.33
1	A	81[A]	HIS	N-CA	-6.27	1.39	1.46
1	A	81[B]	HIS	N-CA	-6.27	1.39	1.46
1	A	191	GLY	C-O	6.15	1.32	1.24
1	A	39	TYR	C-N	-5.78	1.26	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	229	ILE	CB-CA-C	6.08	118.45	110.91
1	A	229	ILE	CA-CB-CG1	-5.93	100.32	110.40
1	A	175	SER	N-CA-CB	-5.83	102.33	110.38

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	2066	0	1994	84	1
2	A	64	0	93	42	0
3	A	10	0	0	1	0
4	A	24	0	36	14	0
5	A	280	0	0	14	2
All	All	2444	0	2123	89	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 21.

The worst 5 of 89 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:157:GLN:OE1	4:A:320:TRS:O1	1.59	1.17
1:A:81[B]:HIS:CE1	2:A:302:EDO:H22	1.86	1.11
1:A:180[B]:ASP:OD1	2:A:312:EDO:C2	1.99	1.09
1:A:180[B]:ASP:OD1	2:A:312:EDO:H21	1.54	1.07
1:A:81[B]:HIS:HE1	2:A:302:EDO:H11	1.22	1.03

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)
5:A:450:HOH:O	5:A:634:HOH:O[4_465]	1.68	0.52
1:A:223[B]:PHE:CE2	5:A:598:HOH:O[4_465]	1.81	0.39

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	254/236 (108%)	249 (98%)	3 (1%)	2 (1%)	16 3

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174[A]	GLY
1	A	174[B]	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	229/206 (111%)	226 (99%)	3 (1%)	61 23

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

Mol	Chain	Res	Type
1	A	3	LYS
1	A	32	GLU
1	A	107	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	177	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](#)

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CRO	A	66	1	22,23,24	2.22	5 (22%)	30,32,34	1.85	8 (26%)

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
1	CRO	A	66	1	-	0/12/31/32	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CRO	CB2-CA2	6.44	1.41	1.35
1	A	66	CRO	CA2-C2	-6.38	1.41	1.48
1	A	66	CRO	C2-N3	-2.86	1.33	1.40
1	A	66	CRO	C1-N3	-2.64	1.32	1.37
1	A	66	CRO	CG2-CB2	-2.25	1.42	1.46

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRO	C3-CA3-N3	4.14	121.84	112.43
1	A	66	CRO	C2-CA2-N2	-3.77	106.25	108.95
1	A	66	CRO	CA2-N2-C1	3.51	108.54	105.80
1	A	66	CRO	CA2-C2-N3	3.22	106.21	103.50
1	A	66	CRO	O2-C2-CA2	-3.12	129.03	131.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

21 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	$\# Z > 2$
4	TRS	A	318	-	7,7,7	0.96	0	9,9,9	1.64	2 (22%)
2	EDO	A	315	-	3,3,3	0.87	0	2,2,2	1.20	0
2	EDO	A	308	-	3,3,3	1.44	0	2,2,2	0.12	0
2	EDO	A	312	-	3,3,3	1.07	0	2,2,2	0.34	0
4	TRS	A	319	-	7,7,7	0.69	0	9,9,9	1.36	1 (11%)
2	EDO	A	314	-	3,3,3	1.76	1 (33%)	2,2,2	1.39	0
2	EDO	A	303	-	3,3,3	0.41	0	2,2,2	0.96	0
4	TRS	A	320	-	7,7,7	1.81	2 (28%)	9,9,9	2.99	5 (55%)
2	EDO	A	313	-	3,3,3	1.17	0	2,2,2	1.55	1 (50%)
2	EDO	A	310	-	3,3,3	0.69	0	2,2,2	0.53	0
2	EDO	A	321	-	3,3,3	3.25	2 (66%)	2,2,2	2.11	1 (50%)
3	SO4	A	316	-	4,4,4	1.72	1 (25%)	6,6,6	0.94	0
3	SO4	A	317	-	4,4,4	1.10	0	6,6,6	0.74	0
2	EDO	A	301	-	3,3,3	1.16	0	2,2,2	1.75	1 (50%)
2	EDO	A	305	-	3,3,3	0.49	0	2,2,2	0.69	0
2	EDO	A	311	-	3,3,3	0.54	0	2,2,2	0.39	0
2	EDO	A	309	-	3,3,3	0.75	0	2,2,2	1.15	0
2	EDO	A	307	-	3,3,3	0.39	0	2,2,2	0.81	0
2	EDO	A	304	-	3,3,3	1.24	0	2,2,2	1.67	1 (50%)
2	EDO	A	306	-	3,3,3	0.96	0	2,2,2	0.49	0
2	EDO	A	302	-	3,3,3	0.96	0	2,2,2	0.35	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	TRS	A	318	-	-	8/9/9/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	315	-	-	1/1/1/1	-
2	EDO	A	308	-	-	1/1/1/1	-
2	EDO	A	312	-	-	0/1/1/1	-
4	TRS	A	319	-	-	6/9/9/9	-
2	EDO	A	314	-	-	0/1/1/1	-
2	EDO	A	303	-	-	1/1/1/1	-
4	TRS	A	320	-	-	9/9/9/9	-
2	EDO	A	313	-	-	1/1/1/1	-
2	EDO	A	310	-	-	1/1/1/1	-
2	EDO	A	321	-	-	0/1/1/1	-
2	EDO	A	301	-	-	0/1/1/1	-
2	EDO	A	305	-	-	1/1/1/1	-
2	EDO	A	311	-	-	1/1/1/1	-
2	EDO	A	309	-	-	1/1/1/1	-
2	EDO	A	307	-	-	1/1/1/1	-
2	EDO	A	304	-	-	1/1/1/1	-
2	EDO	A	306	-	-	0/1/1/1	-
2	EDO	A	302	-	-	1/1/1/1	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	321	EDO	O1-C1	-4.98	1.16	1.42
4	A	320	TRS	C3-C	-3.63	1.43	1.53
3	A	316	SO4	O1-S	-2.64	1.29	1.44
4	A	320	TRS	C-N	-2.59	1.41	1.49
2	A	314	EDO	O2-C2	2.27	1.53	1.42

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	320	TRS	C1-C-N	4.77	120.33	108.17
4	A	320	TRS	O1-C1-C	4.48	123.36	110.88
4	A	320	TRS	C2-C-N	-3.99	97.99	108.17
4	A	318	TRS	O2-C2-C	-3.42	101.33	110.88
4	A	320	TRS	O3-C3-C	-3.26	101.80	110.88

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	318	TRS	N-C-C3-O3
4	A	319	TRS	C2-C-C1-O1
4	A	319	TRS	C3-C-C1-O1
4	A	319	TRS	N-C-C1-O1
4	A	319	TRS	N-C-C2-O2

There are no ring outliers.

18 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	318	TRS	3	0
2	A	315	EDO	1	0
2	A	308	EDO	2	0
2	A	312	EDO	3	0
4	A	319	TRS	9	0
2	A	314	EDO	1	0
2	A	303	EDO	8	0
4	A	320	TRS	2	0
2	A	313	EDO	1	0
2	A	310	EDO	1	0
2	A	321	EDO	4	0
3	A	316	SO4	1	0
2	A	301	EDO	4	0
2	A	305	EDO	2	0
2	A	307	EDO	5	0
2	A	304	EDO	2	0
2	A	306	EDO	1	0
2	A	302	EDO	7	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	224/236 (94%)	0.32	9 (4%)	42 48	3, 11, 21, 72	31 (13%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	230	THR	5.5
1	A	229	ILE	3.8
1	A	39	TYR	2.9
1	A	133	ASP	2.8
1	A	3	LYS	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CRO	A	66	22/23	0.95	0.07	8,9,12,13	0

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
4	TRS	A	319	8/8	0.64	0.38	22,25,29,34	8
2	EDO	A	311	4/4	0.74	0.21	50,53,53,55	0
2	EDO	A	308	4/4	0.76	0.18	22,24,31,40	0
4	TRS	A	318	8/8	0.77	0.21	31,33,44,44	8
2	EDO	A	302	4/4	0.78	0.17	36,38,40,45	0
4	TRS	A	320	8/8	0.79	0.30	23,25,35,38	8
2	EDO	A	314	4/4	0.80	0.16	31,31,31,35	0
2	EDO	A	313	4/4	0.80	0.15	24,30,33,36	0
3	SO4	A	317	5/5	0.82	0.17	26,72,84,92	0
2	EDO	A	315	4/4	0.84	0.14	27,30,40,51	0
2	EDO	A	312	4/4	0.85	0.14	26,30,32,35	0
2	EDO	A	304	4/4	0.86	0.13	27,32,37,43	0
2	EDO	A	301	4/4	0.86	0.14	15,29,32,35	0
2	EDO	A	309	4/4	0.86	0.14	25,27,27,40	0
2	EDO	A	310	4/4	0.87	0.13	27,34,41,41	0
2	EDO	A	307	4/4	0.87	0.12	21,34,38,42	0
2	EDO	A	305	4/4	0.89	0.12	22,30,31,34	0
2	EDO	A	303	4/4	0.89	0.11	30,33,34,38	0
3	SO4	A	316	5/5	0.90	0.14	13,13,20,23	5
2	EDO	A	306	4/4	0.92	0.10	23,23,24,26	0
2	EDO	A	321	4/4	0.95	0.17	10,15,15,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.