



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2022 – 11:22 AM EDT

PDB ID : 8DPD
Title : superfolder GFP Tyr74pCNPhe mutant
Authors : Phillips-Piro, C.M.; Papoutsis, B.; Piacentini, J.
Deposited on : 2022-07-15
Resolution : 1.51 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.29
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

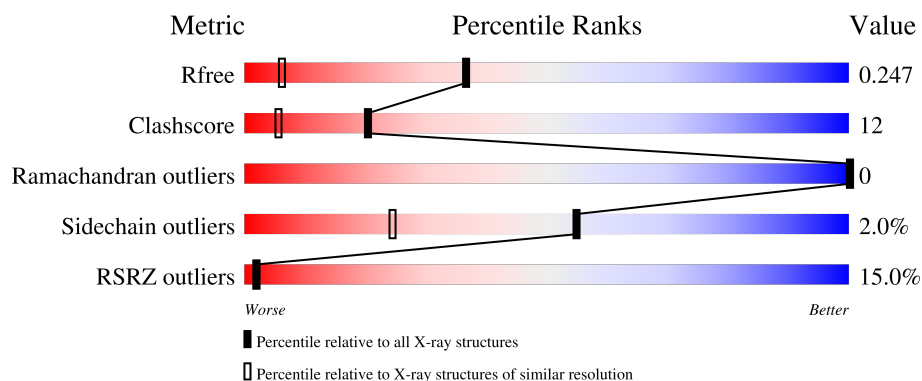
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.51 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}	130704	4009 (1.54-1.50)
Clashscore	141614	4249 (1.54-1.50)
Ramachandran outliers	138981	4148 (1.54-1.50)
Sidechain outliers	138945	4146 (1.54-1.50)
RSRZ outliers	127900	3943 (1.54-1.50)

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	237	<div> <div>14%</div> <div>74%</div> <div>20%</div> <div>• •</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4073 atoms, of which 1896 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.

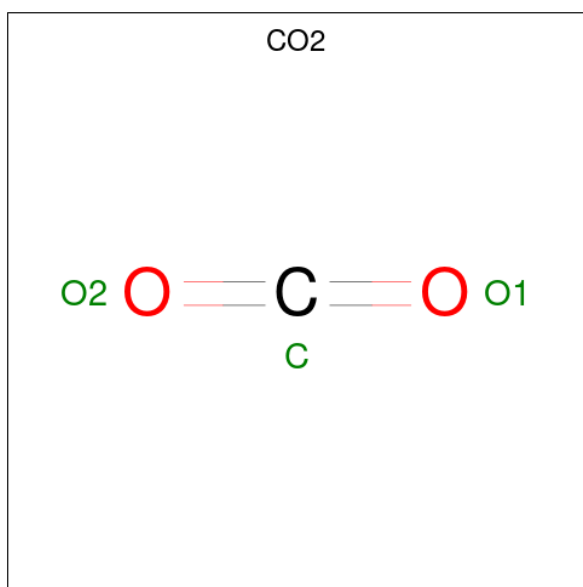
- Molecule 1 is a protein called Superfolder green fluorescent protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	228	Total	C	H	N	O	S	0	13	0
			3782	1214	1862	333	368	5			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P42212
A	1	VAL	-	expression tag	UNP P42212
A	30	ARG	SER	engineered mutation	UNP P42212
A	39	ASN	TYR	engineered mutation	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	74	4CF	TYR	engineered mutation	UNP P42212
A	80	ARG	GLN	engineered mutation	UNP P42212
A	99	SER	PHE	engineered mutation	UNP P42212
A	105	THR	ASN	engineered mutation	UNP P42212
A	145	PHE	TYR	engineered mutation	UNP P42212
A	153	THR	MET	engineered mutation	UNP P42212
A	163	ALA	VAL	engineered mutation	UNP P42212
A	171	VAL	ILE	engineered mutation	UNP P42212
A	206	VAL	ALA	engineered mutation	UNP P42212

- Molecule 2 is CARBON DIOXIDE (three-letter code: CO2) (formula: CO₂).



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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	Mg	0	0
			4	4		

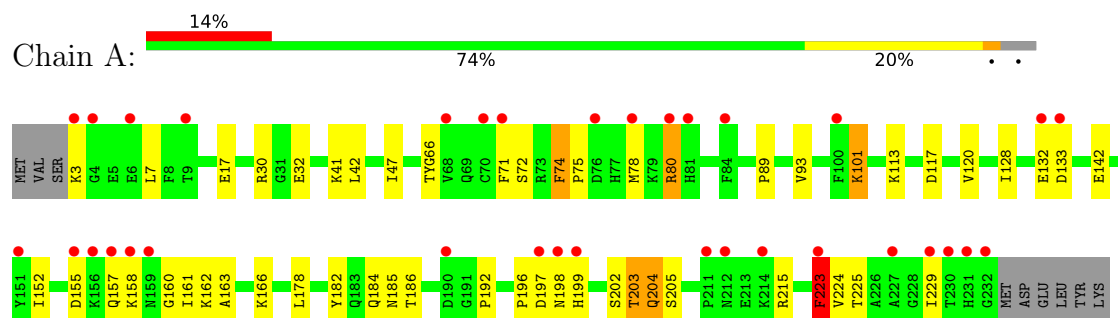
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	227	Total	O	0	0
			227	227		

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: Superfolder green fluorescent protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	47.22Å 47.22Å 346.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.90 – 1.51 19.90 – 1.51	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.90-1.51) 98.8 (19.90-1.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 1.51Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.208 , 0.252 0.205 , 0.247	Depositor DCC
R_{free} test set	2000 reflections (5.36%)	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.8	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	0.97	EDS
Total number of atoms	4073	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.75% 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: PEG, 4CF, MG, CO2, EDO, CRO

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	0.63	3/1934 (0.2%)	0.92	13/2607 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	223[A]	PHE	C-O	6.30	1.35	1.23
1	A	223[B]	PHE	C-O	6.30	1.35	1.23
1	A	203	THR	C-N	-5.34	1.21	1.34

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223[A]	PHE	N-CA-CB	8.90	126.61	110.60
1	A	223[B]	PHE	N-CA-CB	8.90	126.61	110.60
1	A	223[A]	PHE	O-C-N	6.28	132.75	122.70
1	A	223[B]	PHE	O-C-N	6.28	132.75	122.70
1	A	223[A]	PHE	CB-CG-CD2	-6.11	116.53	120.80
1	A	223[B]	PHE	CB-CG-CD2	-6.11	116.53	120.80
1	A	223[A]	PHE	CA-C-O	-6.10	107.30	120.10
1	A	223[B]	PHE	CA-C-O	-6.10	107.30	120.10
1	A	41	LYS	CA-CB-CG	5.52	125.55	113.40
1	A	223[A]	PHE	CB-CG-CD1	5.48	124.64	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223[B]	PHE	CB-CG-CD1	5.48	124.64	120.80
1	A	223[A]	PHE	CA-CB-CG	5.04	126.00	113.90
1	A	223[B]	PHE	CA-CB-CG	5.04	126.00	113.90

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	204[A]	GLN	Mainchain
1	A	204[B]	GLN	Mainchain
1	A	223[A]	PHE	Mainchain

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	1920	1862	1839	47	0
2	A	3	0	0	1	0
3	A	16	24	24	0	0
4	A	7	10	10	0	0
5	A	4	0	0	0	0
6	A	227	0	0	15	1
All	All	2177	1896	1873	48	1

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 12.

All (48) 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:204[B]:GLN:N	1:A:204[B]:GLN:CA	1.69	1.50
1:A:101:LYS:NZ	6:A:405:HOH:O	1.95	0.99
1:A:117[B]:ASP:OD1	6:A:401:HOH:O	1.88	0.91
1:A:133:ASP:OD1	6:A:402:HOH:O	1.90	0.89
1:A:155:ASP:O	6:A:403:HOH:O	1.91	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLU:OE1	6:A:404:HOH:O	1.93	0.87
1:A:182:TYR:OH	1:A:184:GLN:NE2	2.11	0.84
1:A:166:LYS:NZ	6:A:405:HOH:O	1.95	0.76
1:A:160:GLY:N	6:A:403:HOH:O	2.19	0.75
1:A:142[A]:GLU:OE1	6:A:406:HOH:O	2.06	0.72
2:A:301:CO2:O2	6:A:407:HOH:O	2.14	0.65
1:A:203:THR:C	1:A:204[B]:GLN:CA	2.65	0.63
1:A:113:LYS:O	1:A:120[B]:VAL:HG22	1.97	0.63
1:A:17:GLU:OE2	1:A:30[A]:ARG:NH2	2.35	0.60
1:A:30[A]:ARG:NH2	6:A:415:HOH:O	2.35	0.58
1:A:158:LYS:HE2	1:A:186:THR:HG22	1.90	0.53
1:A:155:ASP:OD1	1:A:157:GLN:OE1	2.26	0.53
1:A:162:LYS:HE2	1:A:184:GLN:NE2	2.26	0.51
1:A:223[A]:PHE:CZ	1:A:225:THR:CG2	2.94	0.51
1:A:75:PRO:HG2	1:A:78:MET:CE	2.42	0.50
1:A:158:LYS:HE2	1:A:186:THR:CG2	2.42	0.49
1:A:229:ILE:HG23	1:A:229:ILE:O	2.12	0.48
1:A:3:LYS:HE3	1:A:89:PRO:HG3	1.94	0.47
1:A:42:LEU:HD21	1:A:71:PHE:CB	2.44	0.47
1:A:192:PRO:O	6:A:408:HOH:O	2.20	0.46
1:A:101:LYS:HD3	6:A:495:HOH:O	2.16	0.45
1:A:197:ASP:O	1:A:198:ASN:C	2.54	0.45
1:A:204[B]:GLN:HG3	6:A:472:HOH:O	2.17	0.45
1:A:152:ILE:HD13	1:A:163:ALA:HB2	2.00	0.44
1:A:120[B]:VAL:HG23	6:A:458:HOH:O	2.18	0.44
1:A:204[B]:GLN:CG	1:A:205:SER:N	2.80	0.44
1:A:3:LYS:CE	1:A:89:PRO:HG3	2.48	0.44
1:A:42:LEU:HD21	1:A:71:PHE:HB2	2.00	0.43
1:A:196:PRO:HB3	1:A:199[A]:HIS:CE1	2.54	0.42
1:A:32:GLU:OE2	6:A:409:HOH:O	2.22	0.42
1:A:196:PRO:HB2	1:A:199[A]:HIS:ND1	2.35	0.42
1:A:3:LYS:HE3	1:A:7:LEU:HD11	2.01	0.42
1:A:74:4CF:HD2	1:A:78:MET:HB3	2.02	0.42
1:A:101:LYS:HE3	1:A:178:LEU:HD12	2.01	0.42
1:A:93:VAL:O	1:A:185:ASN:HA	2.20	0.41
1:A:223[A]:PHE:CZ	1:A:225:THR:HG21	2.55	0.41
1:A:47:ILE:HD13	1:A:215:ARG:CZ	2.51	0.41
1:A:155:ASP:OD2	1:A:162:LYS:HE3	2.20	0.41
1:A:80:ARG:H	1:A:80:ARG:HG3	1.70	0.41
1:A:75:PRO:HG2	1:A:78:MET:HE2	2.03	0.41
1:A:72[A]:SER:HA	1:A:224[A]:VAL:HG13	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:GLU:OE2	1:A:30[A]:ARG:CZ	2.69	0.40
1:A:161:ILE:C	1:A:161:ILE:HD12	2.41	0.40

All (1) 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)
6:A:548:HOH:O	6:A:585:HOH:O[8_665]	2.02	0.18

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	235/237 (99%)	230 (98%)	5 (2%)	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	A	210/206 (102%)	205 (98%)	5 (2%)	49	19

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	ARG
1	A	101	LYS
1	A	128	ILE
1	A	202[A]	SER
1	A	202[B]	SER

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	157	GLN
1	A	184	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	CRO	A	66	1	23,23,24	2.40	6 (26%)	30,32,34	3.06	18 (60%)
1	4CF	A	74	1	12,13,14	0.80	0	13,16,18	1.34	1 (7%)

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	-	1/12/31/32	0/2/2/2
1	4CF	A	74	1	-	2/7/8/10	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CRO	CA2-C2	9.39	1.57	1.48
1	A	66	CRO	C2-N3	3.26	1.47	1.39
1	A	66	CRO	CB2-CA2	3.09	1.37	1.35
1	A	66	CRO	CG2-CB2	2.75	1.52	1.46
1	A	66	CRO	OH-CZ	2.12	1.42	1.37
1	A	66	CRO	CA1-N1	-2.02	1.41	1.47

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRO	CA2-N2-C1	9.32	112.64	105.77
1	A	66	CRO	O2-C2-CA2	-4.76	128.29	130.96
1	A	66	CRO	C2-CA2-N2	-4.70	105.64	108.93
1	A	66	CRO	CA2-C2-N3	-4.50	101.25	103.37
1	A	66	CRO	CB2-CA2-N2	4.23	134.70	128.83
1	A	74	4CF	CB-CA-C	-4.08	103.81	111.47
1	A	66	CRO	CD2-CG2-CD1	3.82	123.30	117.64
1	A	66	CRO	O3-C3-CA3	-3.68	115.28	126.39
1	A	66	CRO	O2-C2-N3	3.05	130.41	124.35
1	A	66	CRO	CG2-CB2-CA2	-2.99	126.28	129.94
1	A	66	CRO	C2-N3-C1	2.62	109.29	107.97
1	A	66	CRO	C1-CA1-N1	-2.58	105.78	109.96
1	A	66	CRO	CE2-CZ-CE1	2.43	123.86	119.77
1	A	66	CRO	CD2-CG2-CB2	-2.28	113.44	121.22
1	A	66	CRO	OG1-CB1-CA1	2.24	113.84	109.04
1	A	66	CRO	CB2-CA2-C2	-2.21	119.64	122.28
1	A	66	CRO	CE2-CD2-CG2	-2.19	118.39	121.25
1	A	66	CRO	CG1-CB1-CA1	-2.15	107.10	112.16
1	A	66	CRO	CE1-CD1-CG2	-2.14	118.45	121.25

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	74	4CF	C-CA-CB-CG
1	A	74	4CF	N-CA-CB-CG
1	A	66	CRO	C3-CA3-N3-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	74	4CF	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	EDO	A	304	-	3,3,3	0.49	0	2,2,2	0.33	0
3	EDO	A	302	-	3,3,3	0.38	0	2,2,2	0.90	0
4	PEG	A	306	-	6,6,6	0.18	0	5,5,5	0.10	0
3	EDO	A	305	-	3,3,3	0.44	0	2,2,2	0.44	0
3	EDO	A	303	-	3,3,3	0.54	0	2,2,2	0.12	0
2	CO2	A	301	-	2,2,2	1.06	0	1,1,1	0.28	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	304	-	-	1/1/1/1	-
3	EDO	A	302	-	-	0/1/1/1	-
4	PEG	A	306	-	-	3/4/4/4	-
3	EDO	A	305	-	-	1/1/1/1	-
3	EDO	A	303	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	306	PEG	O1-C1-C2-O2
4	A	306	PEG	O2-C3-C4-O4
4	A	306	PEG	C4-C3-O2-C2
3	A	305	EDO	O1-C1-C2-O2
3	A	304	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	CO2	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	A	226/237 (95%)	0.86	34 (15%) 2 2	16, 28, 75, 122	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	HIS	12.3
1	A	230	THR	8.9
1	A	232	GLY	7.9
1	A	157	GLN	6.5
1	A	3	LYS	6.3
1	A	229	ILE	5.3
1	A	155	ASP	5.2
1	A	159	ASN	4.0
1	A	214	LYS	3.9
1	A	223[A]	PHE	3.9
1	A	198	ASN	3.8
1	A	158	LYS	3.7
1	A	76	ASP	3.6
1	A	212	ASN	3.5
1	A	80	ARG	3.5
1	A	78	MET	3.0
1	A	9	THR	3.0
1	A	68	VAL	2.9
1	A	70	CYS	2.8
1	A	4	GLY	2.8
1	A	199[A]	HIS	2.8
1	A	190	ASP	2.6
1	A	156	LYS	2.5
1	A	197	ASP	2.5
1	A	100	PHE	2.4
1	A	6	GLU	2.4
1	A	71	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	84	PHE	2.3
1	A	132	GLU	2.3
1	A	81	HIS	2.3
1	A	211	PRO	2.2
1	A	133	ASP	2.2
1	A	151	TYR	2.1
1	A	227	ALA	2.1

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	4CF	A	74	13/14	0.91	0.14	33,43,68,69	0
1	CRO	A	66	22/23	0.95	0.11	14,20,43,43	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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	PEG	A	306	7/7	0.73	0.24	52,66,80,85	0
3	EDO	A	303	4/4	0.84	0.29	40,57,60,72	0
3	EDO	A	304	4/4	0.87	0.21	37,45,58,69	0
3	EDO	A	305	4/4	0.88	0.14	44,53,57,60	0
3	EDO	A	302	4/4	0.90	0.13	36,46,56,56	0
2	CO2	A	301	3/3	0.94	0.22	21,21,24,26	3
5	MG	A	307	1/1	0.95	0.15	25,25,25,25	0
5	MG	A	308	1/1	0.95	0.10	23,23,23,23	0
5	MG	A	310	1/1	0.96	0.09	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	MG	A	309	1/1	0.97	0.09	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.