



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 06:22 PM UTC

PDB ID : 5G3O / pdb_00005g3o
Title : Bacillus cereus formamidase (BceAmiF) inhibited with urea.
Authors : Gavira, J.A.; Martinez-Rodriguez, S.; Conejero-Muriel, M.
Deposited on : 2016-04-29
Resolution : 2.15 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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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

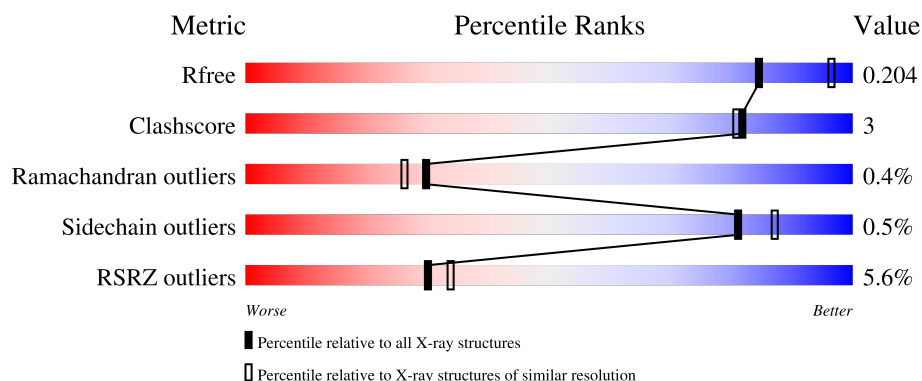
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	346	5% 75% 5% 20%
1	B	346	4% 73% 6% 21%
1	C	346	3% 75% • 21%
1	D	346	4% 75% • 20%
1	E	346	5% 77% • 20%

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Mol	Chain	Length	Quality of chain
1	F	346	 A horizontal bar chart showing the quality of chain 1. The bar is divided into four segments: a small red segment at the beginning labeled '5%', a large green segment labeled '74%', a small yellow segment labeled '6%', and a grey segment at the end labeled '20%'.

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	PEG	D	1311	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14672 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 FORMAMIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	277	Total	C	N	O	S	0	18	0
			2336	1498	381	444	13			
1	B	275	Total	C	N	O	S	0	19	0
			2325	1490	378	444	13			
1	C	274	Total	C	N	O	S	0	17	0
			2296	1473	375	435	13			
1	D	278	Total	C	N	O	S	0	21	0
			2354	1507	384	450	13			
1	E	278	Total	C	N	O	S	0	22	0
			2376	1522	392	450	12			
1	F	277	Total	C	N	O	S	0	20	0
			2351	1504	384	450	13			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	333	TRP	-	expression tag	UNP E5LR94
A	334	PHE	-	expression tag	UNP E5LR94
A	335	ARG	-	expression tag	UNP E5LR94
A	336	VAL	-	expression tag	UNP E5LR94
A	337	ASP	-	expression tag	UNP E5LR94
A	338	PRO	-	expression tag	UNP E5LR94
A	339	LEU	-	expression tag	UNP E5LR94
A	340	GLU	-	expression tag	UNP E5LR94
A	341	HIS	-	expression tag	UNP E5LR94
A	342	HIS	-	expression tag	UNP E5LR94
A	343	HIS	-	expression tag	UNP E5LR94
A	344	HIS	-	expression tag	UNP E5LR94
A	345	HIS	-	expression tag	UNP E5LR94
A	346	HIS	-	expression tag	UNP E5LR94
B	333	TRP	-	expression tag	UNP E5LR94
B	334	PHE	-	expression tag	UNP E5LR94
B	335	ARG	-	expression tag	UNP E5LR94

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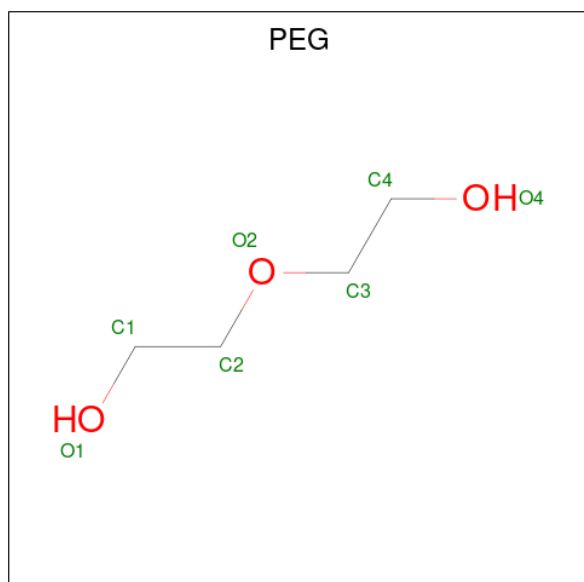
Chain	Residue	Modelled	Actual	Comment	Reference
B	336	VAL	-	expression tag	UNP E5LR94
B	337	ASP	-	expression tag	UNP E5LR94
B	338	PRO	-	expression tag	UNP E5LR94
B	339	LEU	-	expression tag	UNP E5LR94
B	340	GLU	-	expression tag	UNP E5LR94
B	341	HIS	-	expression tag	UNP E5LR94
B	342	HIS	-	expression tag	UNP E5LR94
B	343	HIS	-	expression tag	UNP E5LR94
B	344	HIS	-	expression tag	UNP E5LR94
B	345	HIS	-	expression tag	UNP E5LR94
B	346	HIS	-	expression tag	UNP E5LR94
C	333	TRP	-	expression tag	UNP E5LR94
C	334	PHE	-	expression tag	UNP E5LR94
C	335	ARG	-	expression tag	UNP E5LR94
C	336	VAL	-	expression tag	UNP E5LR94
C	337	ASP	-	expression tag	UNP E5LR94
C	338	PRO	-	expression tag	UNP E5LR94
C	339	LEU	-	expression tag	UNP E5LR94
C	340	GLU	-	expression tag	UNP E5LR94
C	341	HIS	-	expression tag	UNP E5LR94
C	342	HIS	-	expression tag	UNP E5LR94
C	343	HIS	-	expression tag	UNP E5LR94
C	344	HIS	-	expression tag	UNP E5LR94
C	345	HIS	-	expression tag	UNP E5LR94
C	346	HIS	-	expression tag	UNP E5LR94
D	333	TRP	-	expression tag	UNP E5LR94
D	334	PHE	-	expression tag	UNP E5LR94
D	335	ARG	-	expression tag	UNP E5LR94
D	336	VAL	-	expression tag	UNP E5LR94
D	337	ASP	-	expression tag	UNP E5LR94
D	338	PRO	-	expression tag	UNP E5LR94
D	339	LEU	-	expression tag	UNP E5LR94
D	340	GLU	-	expression tag	UNP E5LR94
D	341	HIS	-	expression tag	UNP E5LR94
D	342	HIS	-	expression tag	UNP E5LR94
D	343	HIS	-	expression tag	UNP E5LR94
D	344	HIS	-	expression tag	UNP E5LR94
D	345	HIS	-	expression tag	UNP E5LR94
D	346	HIS	-	expression tag	UNP E5LR94
E	333	TRP	-	expression tag	UNP E5LR94
E	334	PHE	-	expression tag	UNP E5LR94
E	335	ARG	-	expression tag	UNP E5LR94

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Chain	Residue	Modelled	Actual	Comment	Reference
E	336	VAL	-	expression tag	UNP E5LR94
E	337	ASP	-	expression tag	UNP E5LR94
E	338	PRO	-	expression tag	UNP E5LR94
E	339	LEU	-	expression tag	UNP E5LR94
E	340	GLU	-	expression tag	UNP E5LR94
E	341	HIS	-	expression tag	UNP E5LR94
E	342	HIS	-	expression tag	UNP E5LR94
E	343	HIS	-	expression tag	UNP E5LR94
E	344	HIS	-	expression tag	UNP E5LR94
E	345	HIS	-	expression tag	UNP E5LR94
E	346	HIS	-	expression tag	UNP E5LR94
F	333	TRP	-	expression tag	UNP E5LR94
F	334	PHE	-	expression tag	UNP E5LR94
F	335	ARG	-	expression tag	UNP E5LR94
F	336	VAL	-	expression tag	UNP E5LR94
F	337	ASP	-	expression tag	UNP E5LR94
F	338	PRO	-	expression tag	UNP E5LR94
F	339	LEU	-	expression tag	UNP E5LR94
F	340	GLU	-	expression tag	UNP E5LR94
F	341	HIS	-	expression tag	UNP E5LR94
F	342	HIS	-	expression tag	UNP E5LR94
F	343	HIS	-	expression tag	UNP E5LR94
F	344	HIS	-	expression tag	UNP E5LR94
F	345	HIS	-	expression tag	UNP E5LR94
F	346	HIS	-	expression tag	UNP E5LR94

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



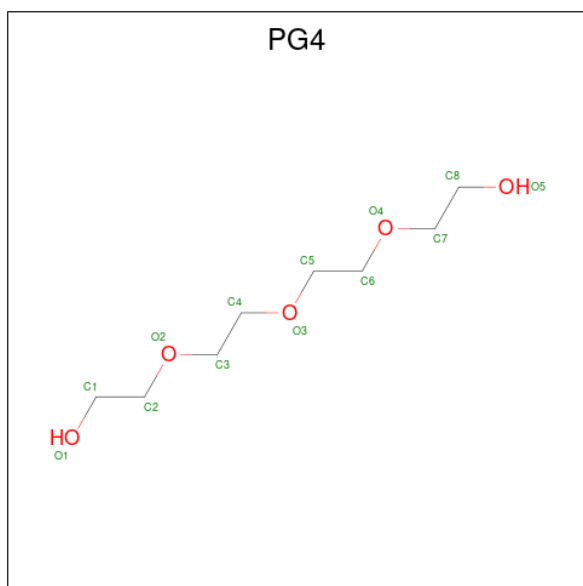
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	F	1	Total C O 7 4 3	0	0
2	F	1	Total C O 7 4 3	0	0
2	F	1	Total C O 7 4 3	0	0

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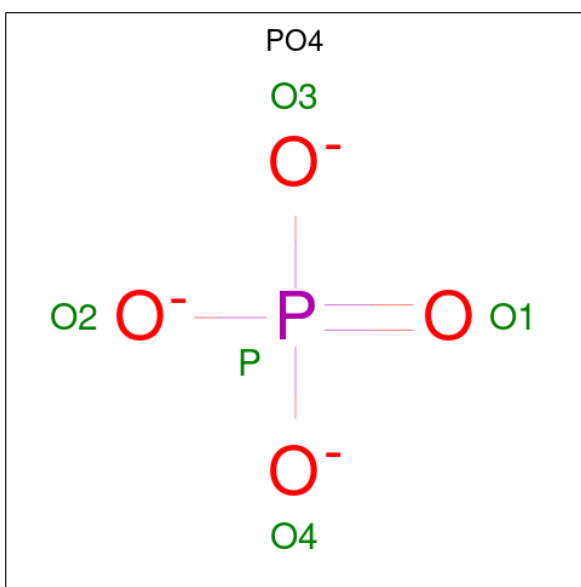
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			7	4	3		

- Molecule 3 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



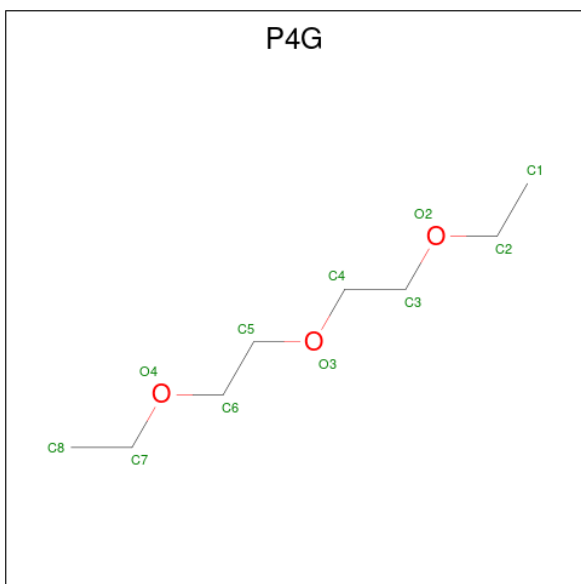
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	8	5		
3	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

- Molecule 5 is 1-ETHOXY-2-(2-ETHOXYETHOXY)ETHANE (CCD ID: P4G) (formula: $C_8H_{18}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			11	8	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	C	O	0	0
			11	8	3		

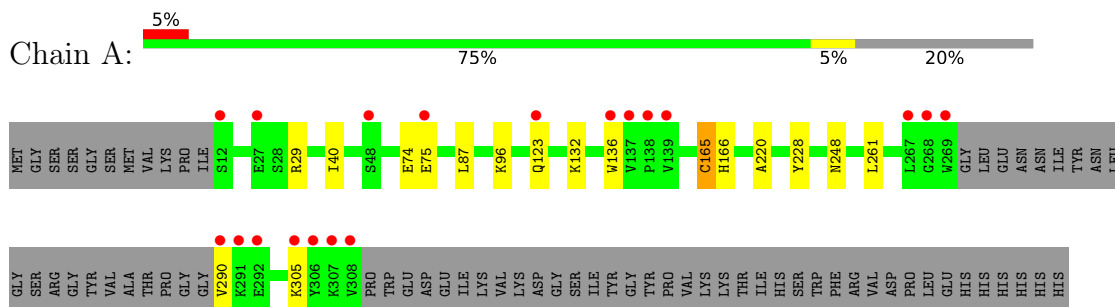
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	95	Total	O	0	0
			95	95		
6	B	74	Total	O	0	0
			74	74		
6	C	70	Total	O	0	0
			70	70		
6	D	69	Total	O	0	0
			69	69		
6	E	61	Total	O	0	0
			61	61		
6	F	46	Total	O	0	0
			46	46		

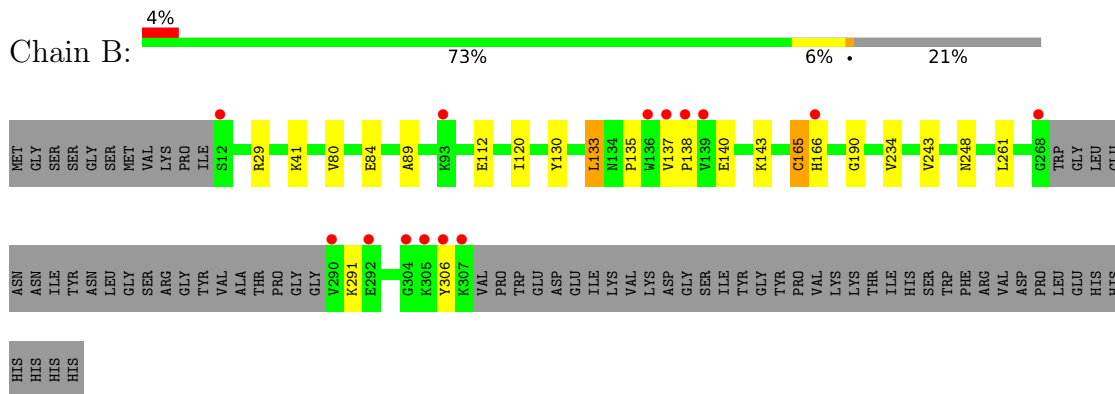
3 Residue-property plots [i](#)

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.

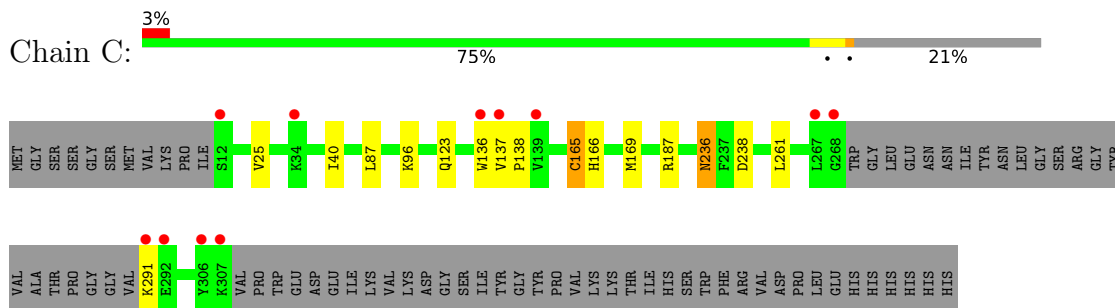
• Molecule 1: FORMAMIDASE



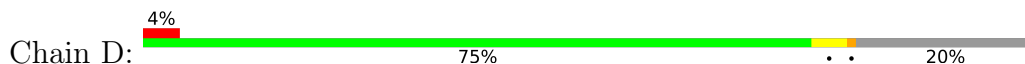
• Molecule 1: FORMAMIDASE

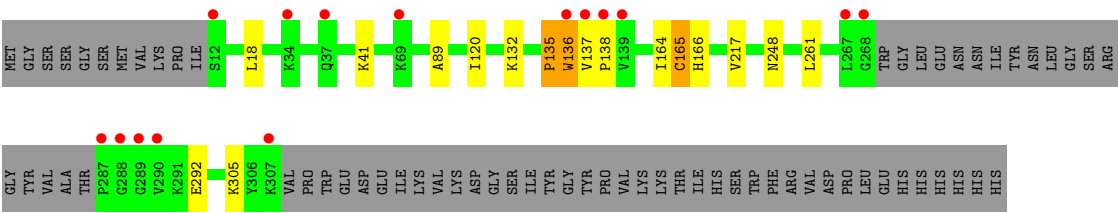


• Molecule 1: FORMAMIDASE

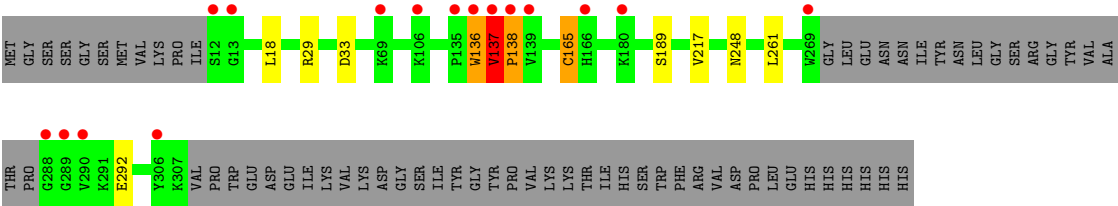
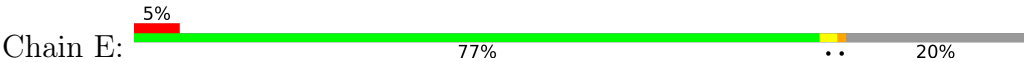


• Molecule 1: FORMAMIDASE

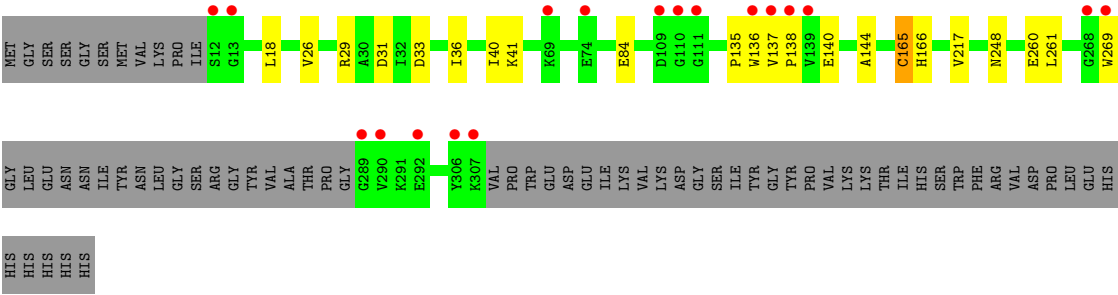




● Molecule 1: FORMAMIDASE



● Molecule 1: FORMAMIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	142.69Å 149.75Å 97.22Å 90.00° 98.68° 90.00°	Depositor
Resolution (Å)	47.06 – 2.15 47.06 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.06-2.15) 99.2 (47.06-2.15)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.14Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.179 , 0.202 0.183 , 0.204	Depositor DCC
R_{free} test set	5444 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14672	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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, PO4, PG4, P4G, QCS

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	0.13	0/2380	0.32	0/3237
1	B	0.12	0/2367	0.31	0/3217
1	C	0.13	0/2338	0.31	0/3176
1	D	0.12	0/2397	0.31	0/3254
1	E	0.13	0/2420	0.31	0/3284
1	F	0.12	0/2395	0.30	0/3258
All	All	0.13	0/14297	0.31	0/19426

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	E	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	136[A]	TRP	Peptide
1	E	137[A]	VAL	Peptide

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	2336	0	2253	11	0
1	B	2325	0	2248	16	0
1	C	2296	0	2228	11	0
1	D	2354	0	2273	15	0
1	E	2376	0	2300	10	0
1	F	2351	0	2255	13	0
2	A	21	0	30	0	0
2	B	28	0	40	0	0
2	C	21	0	30	0	0
2	D	28	0	40	1	0
2	E	35	0	50	0	0
2	F	28	0	40	0	0
3	A	26	0	36	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	C	11	0	18	1	0
5	F	11	0	18	1	0
6	A	95	0	0	2	0
6	B	74	0	0	4	0
6	C	70	0	0	0	0
6	D	69	0	0	0	0
6	E	61	0	0	0	0
6	F	46	0	0	0	0
All	All	14672	0	13859	71	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 3.

The worst 5 of 71 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:D:165[B]:QCS:SG	1:D:166[B]:HIS:N	2.43	0.91
1:D:41:LYS:NZ	2:D:1311:PEG:O4	2.10	0.83
1:F:29[A]:ARG:NH1	1:F:33:ASP:OD1	2.14	0.81
1:C:165[B]:QCS:SG	1:C:166[B]:HIS:N	2.54	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:29[B]:ARG:NH1	1:E:33:ASP:OD1	2.16	0.79

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	289/346 (84%)	278 (96%)	9 (3%)	2 (1%)	18	13
1	B	288/346 (83%)	279 (97%)	9 (3%)	0	100	100
1	C	285/346 (82%)	270 (95%)	13 (5%)	2 (1%)	18	13
1	D	293/346 (85%)	275 (94%)	14 (5%)	4 (1%)	9	4
1	E	294/346 (85%)	281 (96%)	11 (4%)	2 (1%)	18	13
1	F	291/346 (84%)	280 (96%)	9 (3%)	2 (1%)	18	13
All	All	1740/2076 (84%)	1663 (96%)	65 (4%)	12 (1%)	30	13

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	135[A]	PRO
1	D	135[B]	PRO
1	D	136[A]	TRP
1	D	136[B]	TRP
1	F	136[A]	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	251/293 (86%)	251 (100%)	0	100	100
1	B	250/293 (85%)	248 (99%)	2 (1%)	73	79
1	C	246/293 (84%)	245 (100%)	1 (0%)	84	89
1	D	252/293 (86%)	250 (99%)	2 (1%)	73	79
1	E	254/293 (87%)	251 (99%)	3 (1%)	63	70
1	F	252/293 (86%)	250 (99%)	2 (1%)	73	79
All	All	1505/1758 (86%)	1495 (99%)	10 (1%)	81	82

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	292	GLU
1	F	260[A]	GLU
1	F	260[B]	GLU
1	D	292[A]	GLU
1	D	292[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	244	GLN
1	F	194	GLN
1	F	244	GLN
1	F	248	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 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	QCS	A	165[A]	1	6,8,9	1.91	1 (16%)	4,9,11	2.00	1 (25%)
1	QCS	A	165[B]	1	6,8,9	1.93	1 (16%)	4,9,11	1.67	1 (25%)
1	QCS	C	165[A]	1	6,8,9	1.89	1 (16%)	4,9,11	1.99	1 (25%)
1	QCS	B	165[A]	1	6,8,9	1.89	1 (16%)	4,9,11	1.53	1 (25%)
1	QCS	C	165[B]	1	6,8,9	1.85	1 (16%)	4,9,11	1.36	1 (25%)
1	QCS	F	165[A]	1	6,8,9	1.86	1 (16%)	4,9,11	1.30	1 (25%)
1	QCS	B	165[B]	1	6,8,9	1.88	1 (16%)	4,9,11	2.04	1 (25%)
1	QCS	D	165[A]	1	6,8,9	1.92	1 (16%)	4,9,11	1.87	1 (25%)
1	QCS	F	165[B]	1	6,8,9	1.93	1 (16%)	4,9,11	1.98	1 (25%)
1	QCS	E	165[B]	1	6,8,9	1.90	1 (16%)	4,9,11	1.92	1 (25%)
1	QCS	D	165[B]	1	6,8,9	1.84	1 (16%)	4,9,11	1.45	1 (25%)
1	QCS	E	165[A]	1	6,8,9	1.87	1 (16%)	4,9,11	1.48	1 (25%)

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	QCS	A	165[A]	1	-	3/4/7/9	-
1	QCS	A	165[B]	1	-	1/4/7/9	-
1	QCS	C	165[A]	1	-	1/4/7/9	-
1	QCS	B	165[A]	1	-	0/4/7/9	-
1	QCS	C	165[B]	1	-	1/4/7/9	-
1	QCS	F	165[A]	1	-	1/4/7/9	-
1	QCS	B	165[B]	1	-	2/4/7/9	-
1	QCS	D	165[A]	1	-	2/4/7/9	-
1	QCS	F	165[B]	1	-	1/4/7/9	-
1	QCS	E	165[B]	1	-	1/4/7/9	-
1	QCS	D	165[B]	1	-	3/4/7/9	-
1	QCS	E	165[A]	1	-	1/4/7/9	-

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	165[B]	QCS	CD-NE2	3.75	1.42	1.33
1	A	165[A]	QCS	CD-NE2	3.70	1.42	1.33
1	E	165[B]	QCS	CD-NE2	3.70	1.42	1.33
1	C	165[A]	QCS	CD-NE2	3.69	1.42	1.33
1	D	165[A]	QCS	CD-NE2	3.67	1.42	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	165[B]	QCS	OE1-CD-NE2	-3.74	119.78	125.59
1	C	165[A]	QCS	OE1-CD-NE2	-3.68	119.86	125.59
1	F	165[B]	QCS	OE1-CD-NE2	-3.67	119.88	125.59
1	E	165[B]	QCS	OE1-CD-NE2	-3.57	120.04	125.59
1	A	165[A]	QCS	OE1-CD-NE2	-3.55	120.07	125.59

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	165[A]	QCS	CA-CB-SG-CD
1	A	165[A]	QCS	OE1-CD-SG-CB
1	A	165[B]	QCS	OE1-CD-SG-CB
1	B	165[B]	QCS	OE1-CD-SG-CB
1	C	165[B]	QCS	OE1-CD-SG-CB

There are no ring outliers.

9 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	165[A]	QCS	1	0
1	A	165[B]	QCS	1	0
1	B	165[A]	QCS	2	0
1	C	165[B]	QCS	2	0
1	F	165[A]	QCS	1	0
1	B	165[B]	QCS	1	0
1	D	165[A]	QCS	1	0
1	D	165[B]	QCS	3	0
1	E	165[A]	QCS	1	0

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

29 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	PO4	C	1312	-	4,4,4	0.91	0	6,6,6	0.52	0
2	PEG	C	1310	-	6,6,6	0.50	0	5,5,5	0.38	0
2	PEG	F	1309	-	6,6,6	0.50	0	5,5,5	0.30	0
2	PEG	B	1311	-	6,6,6	0.50	0	5,5,5	0.40	0
2	PEG	F	1310	-	6,6,6	0.50	0	5,5,5	0.28	0
2	PEG	C	1309	-	6,6,6	0.48	0	5,5,5	0.35	0
2	PEG	E	1311	-	6,6,6	0.50	0	5,5,5	0.31	0
2	PEG	F	1308	-	6,6,6	0.48	0	5,5,5	0.46	0
2	PEG	B	1308	-	6,6,6	0.48	0	5,5,5	0.40	0
2	PEG	D	1309	-	6,6,6	0.49	0	5,5,5	0.39	0
2	PEG	F	1311	-	6,6,6	0.48	0	5,5,5	0.52	0
2	PEG	E	1309	-	6,6,6	0.51	0	5,5,5	0.30	0
3	PG4	A	1312	-	12,12,12	0.52	0	11,11,11	0.39	0
2	PEG	E	1310	-	6,6,6	0.50	0	5,5,5	0.26	0
2	PEG	D	1308	-	6,6,6	0.48	0	5,5,5	0.42	0
2	PEG	A	1309	-	6,6,6	0.49	0	5,5,5	0.32	0
2	PEG	C	1308	-	6,6,6	0.49	0	5,5,5	0.28	0
2	PEG	D	1311	-	6,6,6	0.62	0	5,5,5	1.40	1 (20%)
2	PEG	A	1310	-	6,6,6	0.49	0	5,5,5	0.31	0
2	PEG	E	1312	-	6,6,6	0.50	0	5,5,5	0.34	0
2	PEG	E	1308	-	6,6,6	0.48	0	5,5,5	0.45	0
2	PEG	A	1311	-	6,6,6	0.50	0	5,5,5	0.30	0
3	PG4	A	1313	-	12,12,12	0.53	0	11,11,11	0.32	0
2	PEG	B	1309	-	6,6,6	0.50	0	5,5,5	0.33	0
2	PEG	B	1310	-	6,6,6	0.51	0	5,5,5	0.30	0
2	PEG	D	1310	-	6,6,6	0.48	0	5,5,5	0.39	0
4	PO4	B	1312	-	4,4,4	0.91	0	6,6,6	0.46	0
5	P4G	F	1312	-	10,10,10	0.56	0	9,9,9	0.44	0
5	P4G	C	1311	-	10,10,10	0.56	0	9,9,9	0.52	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
2	PEG	C	1310	-	-	2/4/4/4	-
2	PEG	F	1309	-	-	3/4/4/4	-
2	PEG	B	1311	-	-	1/4/4/4	-
2	PEG	F	1310	-	-	1/4/4/4	-
2	PEG	C	1309	-	-	2/4/4/4	-
2	PEG	E	1311	-	-	1/4/4/4	-
2	PEG	F	1308	-	-	2/4/4/4	-
2	PEG	B	1308	-	-	1/4/4/4	-
2	PEG	D	1309	-	-	3/4/4/4	-
2	PEG	F	1311	-	-	4/4/4/4	-
2	PEG	E	1309	-	-	1/4/4/4	-
3	PG4	A	1312	-	-	7/10/10/10	-
2	PEG	E	1310	-	-	1/4/4/4	-
2	PEG	D	1308	-	-	2/4/4/4	-
2	PEG	A	1309	-	-	2/4/4/4	-
2	PEG	C	1308	-	-	2/4/4/4	-
2	PEG	D	1311	-	-	1/4/4/4	-
2	PEG	A	1310	-	-	1/4/4/4	-
2	PEG	E	1312	-	-	2/4/4/4	-
2	PEG	E	1308	-	-	0/4/4/4	-
2	PEG	A	1311	-	-	2/4/4/4	-
3	PG4	A	1313	-	-	3/10/10/10	-
2	PEG	B	1309	-	-	2/4/4/4	-
2	PEG	B	1310	-	-	1/4/4/4	-
2	PEG	D	1310	-	-	3/4/4/4	-
5	P4G	F	1312	-	-	4/8/8/8	-
5	P4G	C	1311	-	-	3/8/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1311	PEG	C3-O2-C2	2.87	125.83	113.26

There are no chirality outliers.

5 of 57 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1312	PG4	O3-C5-C6-O4
2	B	1310	PEG	O1-C1-C2-O2
2	D	1308	PEG	O2-C3-C4-O4
2	F	1309	PEG	O2-C3-C4-O4
2	D	1309	PEG	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1311	PEG	1	0
5	F	1312	P4G	1	0
5	C	1311	P4G	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 [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	276/346 (79%)	0.12	19 (6%) 23 26	16, 37, 58, 107	17 (6%)
1	B	274/346 (79%)	0.32	14 (5%) 33 37	19, 43, 73, 103	18 (6%)
1	C	273/346 (78%)	0.03	11 (4%) 42 47	17, 38, 55, 103	16 (5%)
1	D	277/346 (80%)	0.10	15 (5%) 31 35	17, 40, 59, 117	20 (7%)
1	E	277/346 (80%)	0.19	16 (5%) 29 32	17, 40, 68, 95	21 (7%)
1	F	276/346 (79%)	0.25	18 (6%) 25 28	19, 41, 67, 100	19 (6%)
All	All	1653/2076 (79%)	0.17	93 (5%) 30 34	16, 40, 65, 117	111 (6%)

The worst 5 of 93 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	136[A]	TRP	11.4
1	B	136[A]	TRP	9.3
1	A	290	VAL	8.0
1	F	136[A]	TRP	7.7
1	A	269	TRP	5.8

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	QCS	D	165[A]	9/10	0.88	0.13	46,47,52,53	9
1	QCS	D	165[B]	9/10	0.88	0.13	46,47,51,52	9
1	QCS	E	165[A]	9/10	0.89	0.17	47,49,52,53	9
1	QCS	E	165[B]	9/10	0.89	0.17	47,49,53,54	9

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	QCS	F	165[A]	9/10	0.89	0.17	46,47,52,53	9
1	QCS	F	165[B]	9/10	0.89	0.17	46,47,53,53	9
1	QCS	A	165[A]	9/10	0.90	0.12	44,45,50,51	9
1	QCS	A	165[B]	9/10	0.90	0.12	44,45,49,50	9
1	QCS	C	165[A]	9/10	0.92	0.12	40,41,46,47	9
1	QCS	C	165[B]	9/10	0.92	0.12	40,41,45,47	9
1	QCS	B	165[A]	9/10	0.92	0.14	46,47,50,50	9
1	QCS	B	165[B]	9/10	0.92	0.14	46,46,51,54	9

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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($\text{\AA}^2$)	Q<0.9
5	P4G	F	1312	11/11	0.60	0.26	63,64,65,65	0
2	PEG	F	1310	7/7	0.72	0.28	85,86,87,87	0
2	PEG	B	1311	7/7	0.75	0.20	86,87,88,88	0
2	PEG	D	1311	7/7	0.78	0.44	131,132,134,134	0
2	PEG	C	1308	7/7	0.79	0.18	73,73,74,74	0
3	PG4	A	1313	13/13	0.80	0.31	109,112,114,114	0
2	PEG	B	1308	7/7	0.80	0.21	66,66,67,67	0
2	PEG	B	1309	7/7	0.81	0.23	76,77,79,79	0
2	PEG	F	1309	7/7	0.82	0.17	79,79,80,80	0
2	PEG	D	1309	7/7	0.84	0.17	72,72,73,73	0
2	PEG	D	1308	7/7	0.84	0.17	64,64,64,64	0
2	PEG	E	1308	7/7	0.84	0.23	69,69,71,71	0
2	PEG	E	1309	7/7	0.84	0.19	76,77,79,79	0
5	P4G	C	1311	11/11	0.85	0.17	56,57,58,58	0
2	PEG	A	1310	7/7	0.85	0.21	84,85,87,87	0
2	PEG	E	1311	7/7	0.86	0.20	88,88,89,89	0
2	PEG	E	1310	7/7	0.86	0.16	73,73,74,74	0
3	PG4	A	1312	13/13	0.87	0.23	85,88,92,92	0
2	PEG	F	1308	7/7	0.87	0.17	65,65,65,66	0
2	PEG	D	1310	7/7	0.87	0.18	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PEG	A	1311	7/7	0.87	0.14	68,68,69,69	0
2	PEG	C	1309	7/7	0.88	0.20	74,74,75,75	0
2	PEG	A	1309	7/7	0.89	0.15	70,71,72,73	0
2	PEG	F	1311	7/7	0.89	0.15	73,73,73,73	0
2	PEG	B	1310	7/7	0.90	0.23	109,110,110,110	0
2	PEG	E	1312	7/7	0.90	0.14	74,75,76,76	0
2	PEG	C	1310	7/7	0.92	0.14	70,70,72,73	0
4	PO4	C	1312	5/5	0.97	0.07	28,30,31,31	0
4	PO4	B	1312	5/5	0.99	0.06	31,32,33,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.