



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 6, 2017 – 10:40 AM EST

PDB ID : 4FXT
Title : Crystal structure of a DUF3823 family protein (BACOVA_02663) from *Bacteroides ovatus* ATCC 8483 at 2.77 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : unknown
Resolution : 2.77 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

<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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

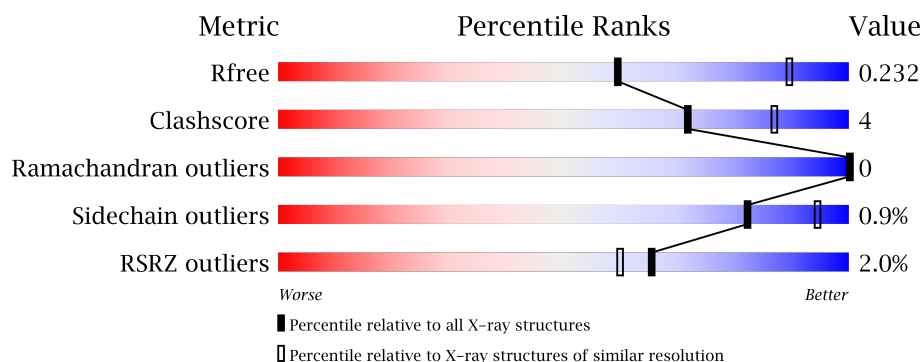
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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}	100719	3276 (2.80-2.76)
Clashscore	112137	3771 (2.80-2.76)
Ramachandran outliers	110173	3707 (2.80-2.76)
Sidechain outliers	110143	3709 (2.80-2.76)
RSRZ outliers	101464	3307 (2.80-2.76)

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	202	
1	B	202	
1	C	202	
1	D	202	
1	E	202	

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Mol	Chain	Length	Quality of chain
1	F	202	
1	G	202	
1	H	202	
1	I	202	
1	J	202	
1	K	202	
1	L	202	
1	M	202	
1	N	202	

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
2	EDO	B	302	-	-	-	X
2	EDO	G	301	-	-	-	X
2	EDO	H	301	-	-	-	X
2	EDO	M	301	-	-	-	X
2	EDO	N	301	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 21144 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	Se	0	0	0
			1475	932	244	296	1	2			
1	B	193	Total	C	N	O	S	Se	0	1	0
			1489	941	248	297	1	2			
1	C	191	Total	C	N	O	S	Se	0	0	0
			1460	923	242	292	1	2			
1	D	192	Total	C	N	O	S	Se	0	0	0
			1470	927	246	294	1	2			
1	E	202	Total	C	N	O	S	Se	0	0	0
			1545	979	257	306	1	2			
1	F	197	Total	C	N	O	S	Se	0	0	0
			1515	958	252	302	1	2			
1	G	196	Total	C	N	O	S	Se	0	0	0
			1505	952	251	300	1	1			
1	H	196	Total	C	N	O	S	Se	0	0	0
			1502	951	248	300	1	2			
1	I	191	Total	C	N	O	S	Se	0	0	0
			1468	928	243	294	1	2			
1	J	197	Total	C	N	O	S	Se	0	0	0
			1509	953	252	301	1	2			
1	K	193	Total	C	N	O	S	Se	0	0	0
			1485	939	248	295	1	2			
1	L	191	Total	C	N	O	S	Se	0	1	0
			1473	929	247	294	1	2			
1	M	195	Total	C	N	O	S	Se	0	0	0
			1506	953	250	300	1	2			
1	N	188	Total	C	N	O	S	Se	0	1	0
			1447	915	240	289	1	2			

There are 14 discrepancies between the modelled and reference sequences:

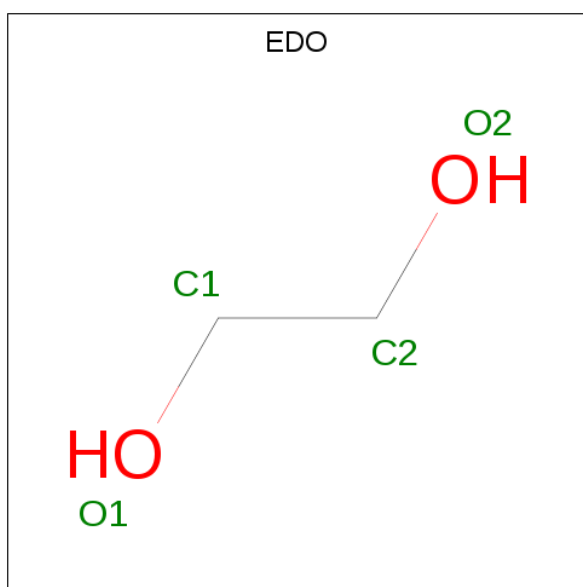
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP A7LXU7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	leader sequence	UNP A7LXU7
C	0	GLY	-	leader sequence	UNP A7LXU7
D	0	GLY	-	leader sequence	UNP A7LXU7
E	0	GLY	-	leader sequence	UNP A7LXU7
F	0	GLY	-	leader sequence	UNP A7LXU7
G	0	GLY	-	leader sequence	UNP A7LXU7
H	0	GLY	-	leader sequence	UNP A7LXU7
I	0	GLY	-	leader sequence	UNP A7LXU7
J	0	GLY	-	leader sequence	UNP A7LXU7
K	0	GLY	-	leader sequence	UNP A7LXU7
L	0	GLY	-	leader sequence	UNP A7LXU7
M	0	GLY	-	leader sequence	UNP A7LXU7
N	0	GLY	-	leader sequence	UNP A7LXU7

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



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	I	1	Total C O 4 2 2	0	0
2	J	1	Total C O 4 2 2	0	0
2	K	1	Total C O 4 2 2	0	0
2	L	1	Total C O 4 2 2	0	0
2	M	1	Total C O 4 2 2	0	0
2	N	1	Total C O 4 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	24	Total O 24 24	0	0
3	B	31	Total O 31 31	0	0
3	C	13	Total O 13 13	0	0
3	D	21	Total O 21 21	0	0
3	E	15	Total O 15 15	0	0
3	F	19	Total O 19 19	0	0
3	G	21	Total O 21 21	0	0
3	H	16	Total O 16 16	0	0

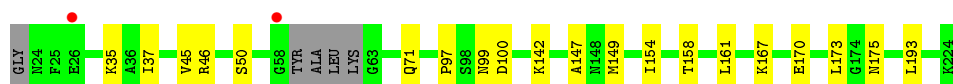
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	I	17	Total 17	O 17	0	0
3	J	12	Total 12	O 12	0	0
3	K	14	Total 14	O 14	0	0
3	L	8	Total 8	O 8	0	0
3	M	10	Total 10	O 10	0	0
3	N	10	Total 10	O 10	0	0

- Molecule 1: Uncharacterized protein





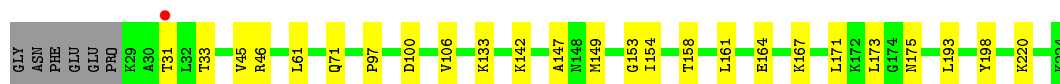
- Molecule 1: Uncharacterized protein

Chain G: 87% 10% .



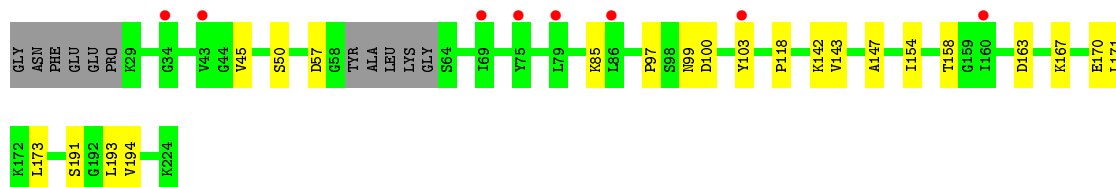
- Molecule 1: Uncharacterized protein

Chain H: 85% 12% .



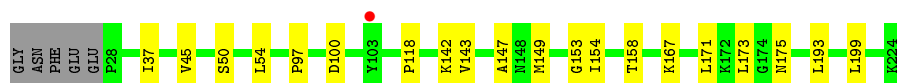
- Molecule 1: Uncharacterized protein

Chain I: 4% 84% 11% 5%



- Molecule 1: Uncharacterized protein

Chain J: 88% 10% .



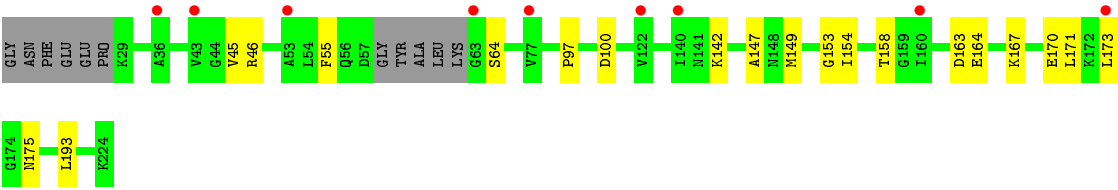
- Molecule 1: Uncharacterized protein

Chain K: 83% 12% .

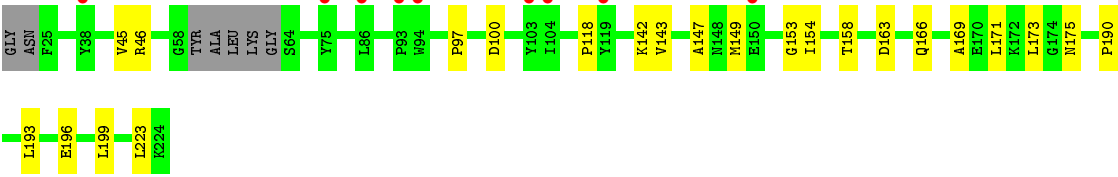
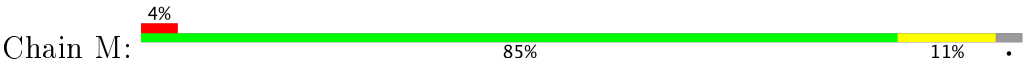


- Molecule 1: Uncharacterized protein

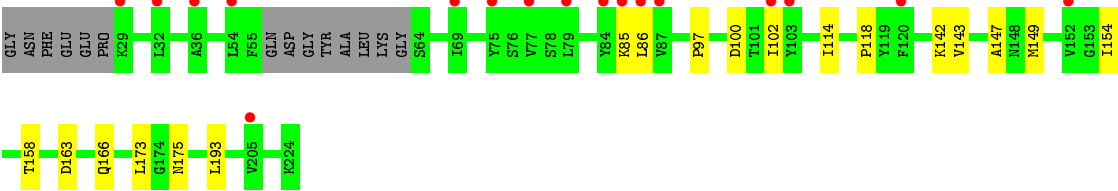
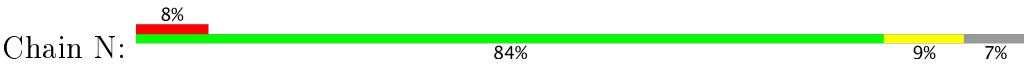
Chain L: 4% 85% 10% 5%



• Molecule 1: Uncharacterized protein



• Molecule 1: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.81Å 351.00Å 90.34Å 90.00° 93.42° 90.00°	Depositor
Resolution (Å)	49.08 – 2.77 48.06 – 2.77	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.08-2.77) 99.8 (48.06-2.77)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.77Å)	Xtriage
Refinement program	BUSTER-TNT 2.10.0, BUSTER 2.10.0	Depositor
R, R_{free}	0.183 , 0.217 0.200 , 0.232	Depositor DCC
R_{free} test set	4158 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	58.9	Xtriage
Anisotropy	0.606	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21144	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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: 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	0.57	0/1498	0.75	0/2025
1	B	0.56	0/1515	0.78	0/2046
1	C	0.51	0/1483	0.73	0/2005
1	D	0.50	0/1492	0.73	0/2016
1	E	0.51	0/1571	0.74	0/2125
1	F	0.51	0/1539	0.74	0/2079
1	G	0.50	0/1529	0.75	0/2068
1	H	0.57	0/1526	0.76	0/2064
1	I	0.50	0/1490	0.72	0/2013
1	J	0.52	0/1534	0.76	0/2074
1	K	0.51	0/1508	0.74	0/2038
1	L	0.49	0/1497	0.73	0/2021
1	M	0.51	0/1530	0.74	0/2067
1	N	0.51	0/1471	0.74	0/1988
All	All	0.52	0/21183	0.74	0/28629

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	1475	0	1448	14	0
1	B	1489	0	1471	11	0
1	C	1460	0	1432	14	0
1	D	1470	0	1449	12	0
1	E	1545	0	1518	17	0
1	F	1515	0	1488	13	0
1	G	1505	0	1483	13	0
1	H	1502	0	1479	14	0
1	I	1468	0	1448	11	0
1	J	1509	0	1480	10	0
1	K	1485	0	1465	11	0
1	L	1473	0	1460	12	0
1	M	1506	0	1483	12	0
1	N	1447	0	1434	9	0
2	A	4	0	6	0	0
2	B	12	0	18	2	0
2	C	4	0	6	1	0
2	D	4	0	6	0	0
2	E	4	0	6	0	0
2	F	4	0	6	2	0
2	G	4	0	6	0	0
2	H	4	0	6	0	0
2	I	4	0	6	0	0
2	J	4	0	6	0	0
2	K	4	0	6	0	0
2	L	4	0	6	0	0
2	M	4	0	6	1	0
2	N	4	0	6	1	0
3	A	24	0	0	0	0
3	B	31	0	0	2	0
3	C	13	0	0	0	0
3	D	21	0	0	1	0
3	E	15	0	0	2	0
3	F	19	0	0	0	0
3	G	21	0	0	0	0
3	H	16	0	0	0	0
3	I	17	0	0	1	0
3	J	12	0	0	0	0
3	K	14	0	0	0	0
3	L	8	0	0	0	0
3	M	10	0	0	0	0
3	N	10	0	0	0	0
All	All	21144	0	20634	164	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

The worst 5 of 164 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:E:169:ALA:HB2	1:E:190:PRO:HG3	1.64	0.79
1:H:164:GLU:O	1:H:167:LYS:HE3	1.83	0.78
1:A:168:GLU:HG2	1:A:193:LEU:HD13	1.72	0.70
1:F:158:THR:H	2:F:301:EDO:H21	1.57	0.70
1:D:161:LEU:HD11	1:G:171:LEU:HD11	1.73	0.69

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	189/202 (94%)	186 (98%)	3 (2%)	0	100	100
1	B	190/202 (94%)	184 (97%)	6 (3%)	0	100	100
1	C	187/202 (93%)	183 (98%)	4 (2%)	0	100	100
1	D	188/202 (93%)	185 (98%)	3 (2%)	0	100	100
1	E	200/202 (99%)	193 (96%)	7 (4%)	0	100	100
1	F	193/202 (96%)	189 (98%)	4 (2%)	0	100	100
1	G	194/202 (96%)	189 (97%)	5 (3%)	0	100	100
1	H	194/202 (96%)	189 (97%)	5 (3%)	0	100	100
1	I	187/202 (93%)	184 (98%)	3 (2%)	0	100	100
1	J	195/202 (96%)	192 (98%)	3 (2%)	0	100	100
1	K	189/202 (94%)	185 (98%)	4 (2%)	0	100	100
1	L	188/202 (93%)	185 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	M	191/202 (95%)	185 (97%)	6 (3%)	0	100	100
1	N	185/202 (92%)	182 (98%)	3 (2%)	0	100	100
All	All	2670/2828 (94%)	2611 (98%)	59 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	160/167 (96%)	159 (99%)	1 (1%)	89	96
1	B	162/167 (97%)	160 (99%)	2 (1%)	75	92
1	C	158/167 (95%)	156 (99%)	2 (1%)	73	92
1	D	160/167 (96%)	159 (99%)	1 (1%)	89	96
1	E	165/167 (99%)	164 (99%)	1 (1%)	89	96
1	F	164/167 (98%)	163 (99%)	1 (1%)	89	96
1	G	162/167 (97%)	162 (100%)	0	100	100
1	H	162/167 (97%)	161 (99%)	1 (1%)	89	96
1	I	160/167 (96%)	158 (99%)	2 (1%)	73	92
1	J	162/167 (97%)	159 (98%)	3 (2%)	62	87
1	K	161/167 (96%)	158 (98%)	3 (2%)	62	87
1	L	161/167 (96%)	160 (99%)	1 (1%)	89	96
1	M	164/167 (98%)	163 (99%)	1 (1%)	89	96
1	N	158/167 (95%)	157 (99%)	1 (1%)	89	96
All	All	2259/2338 (97%)	2239 (99%)	20 (1%)	82	94

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	57	ASP

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Mol	Chain	Res	Type
1	I	163	ASP
1	K	163	ASP
1	F	99	ASN
1	H	106	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	GLN
1	C	71	GLN
1	D	71	GLN
1	F	99	ASN
1	J	112	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

16 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
2	EDO	A	301	-	3,3,3	0.76	0	2,2,2	0.25	0
2	EDO	B	301	-	3,3,3	0.65	0	2,2,2	0.09	0
2	EDO	B	302	-	3,3,3	1.08	0	2,2,2	0.08	0
2	EDO	B	303	-	3,3,3	0.69	0	2,2,2	0.20	0
2	EDO	C	301	-	3,3,3	0.67	0	2,2,2	0.10	0
2	EDO	D	301	-	3,3,3	0.65	0	2,2,2	0.10	0
2	EDO	E	301	-	3,3,3	0.69	0	2,2,2	0.09	0
2	EDO	F	301	-	3,3,3	0.64	0	2,2,2	0.37	0
2	EDO	G	301	-	3,3,3	0.82	0	2,2,2	0.20	0
2	EDO	H	301	-	3,3,3	0.85	0	2,2,2	0.29	0
2	EDO	I	301	-	3,3,3	0.67	0	2,2,2	0.24	0
2	EDO	J	301	-	3,3,3	0.70	0	2,2,2	0.11	0
2	EDO	K	301	-	3,3,3	0.68	0	2,2,2	0.14	0
2	EDO	L	301	-	3,3,3	0.76	0	2,2,2	0.27	0
2	EDO	M	301	-	3,3,3	0.63	0	2,2,2	0.34	0
2	EDO	N	301	-	3,3,3	0.73	0	2,2,2	0.06	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	EDO	A	301	-	-	0/1/1/1	0/0/0/0
2	EDO	B	301	-	-	0/1/1/1	0/0/0/0
2	EDO	B	302	-	-	0/1/1/1	0/0/0/0
2	EDO	B	303	-	-	0/1/1/1	0/0/0/0
2	EDO	C	301	-	-	0/1/1/1	0/0/0/0
2	EDO	D	301	-	-	0/1/1/1	0/0/0/0
2	EDO	E	301	-	-	0/1/1/1	0/0/0/0
2	EDO	F	301	-	-	0/1/1/1	0/0/0/0
2	EDO	G	301	-	-	0/1/1/1	0/0/0/0
2	EDO	H	301	-	-	0/1/1/1	0/0/0/0
2	EDO	I	301	-	-	0/1/1/1	0/0/0/0
2	EDO	J	301	-	-	0/1/1/1	0/0/0/0
2	EDO	K	301	-	-	0/1/1/1	0/0/0/0
2	EDO	L	301	-	-	0/1/1/1	0/0/0/0
2	EDO	M	301	-	-	0/1/1/1	0/0/0/0
2	EDO	N	301	-	-	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.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	EDO	1	0
2	B	302	EDO	1	0
2	C	301	EDO	1	0
2	F	301	EDO	2	0
2	M	301	EDO	1	0
2	N	301	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	191/202 (94%)	-0.02	0 100 100	38, 57, 102, 142	0
1	B	191/202 (94%)	0.08	2 (1%) 82 79	44, 61, 87, 108	0
1	C	189/202 (93%)	0.07	0 100 100	41, 67, 104, 124	0
1	D	190/202 (94%)	0.28	2 (1%) 80 77	46, 68, 108, 140	0
1	E	200/202 (99%)	0.13	1 (0%) 90 89	47, 68, 104, 132	0
1	F	195/202 (96%)	0.08	2 (1%) 82 79	43, 66, 108, 132	0
1	G	194/202 (96%)	0.09	0 100 100	43, 68, 94, 117	0
1	H	194/202 (96%)	0.06	1 (0%) 90 89	34, 61, 102, 122	0
1	I	189/202 (93%)	0.32	8 (4%) 37 30	50, 70, 117, 164	0
1	J	195/202 (96%)	0.14	1 (0%) 90 89	43, 64, 102, 132	0
1	K	191/202 (94%)	0.22	3 (1%) 72 68	58, 75, 104, 123	0
1	L	189/202 (93%)	0.37	9 (4%) 31 24	61, 82, 108, 126	0
1	M	193/202 (95%)	0.43	9 (4%) 32 25	62, 88, 141, 157	0
1	N	186/202 (92%)	0.51	17 (9%) 10 6	56, 83, 152, 170	0
All	All	2687/2828 (95%)	0.20	55 (2%) 65 60	34, 70, 113, 170	0

The worst 5 of 55 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	77	VAL	6.8
1	N	86	LEU	6.0
1	N	79	LEU	5.9
1	N	87	VAL	4.5
1	N	103	TYR	3.9

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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
2	EDO	B	302	4/4	0.75	0.47	8.24	63,64,67,68	0
2	EDO	H	301	4/4	0.90	0.40	5.44	56,59,64,66	0
2	EDO	M	301	4/4	0.93	0.31	4.55	55,55,57,57	0
2	EDO	G	301	4/4	0.90	0.30	3.81	61,63,63,64	0
2	EDO	N	301	4/4	0.92	0.30	3.04	57,58,60,60	0
2	EDO	A	301	4/4	0.94	0.27	2.50	67,67,68,68	0
2	EDO	J	301	4/4	0.91	0.27	1.48	64,66,69,71	0
2	EDO	B	303	4/4	0.89	0.21	0.88	77,79,80,81	0
2	EDO	D	301	4/4	0.94	0.23	0.76	53,54,54,54	0
2	EDO	C	301	4/4	0.98	0.19	0.74	66,66,68,69	0
2	EDO	F	301	4/4	0.92	0.21	0.41	66,68,69,71	0
2	EDO	I	301	4/4	0.91	0.19	0.08	56,58,64,70	0
2	EDO	B	301	4/4	0.96	0.19	-0.31	52,54,57,61	0
2	EDO	K	301	4/4	0.93	0.18	-0.53	57,58,58,59	0
2	EDO	E	301	4/4	0.96	0.17	-0.59	68,70,72,73	0
2	EDO	L	301	4/4	0.90	0.16	-0.83	54,58,61,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.