



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:17 AM UTC

PDB ID : 4FXT / pdb_00004fxt
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 : 2012-07-03
Resolution : 2.77 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 : 4-5-2 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

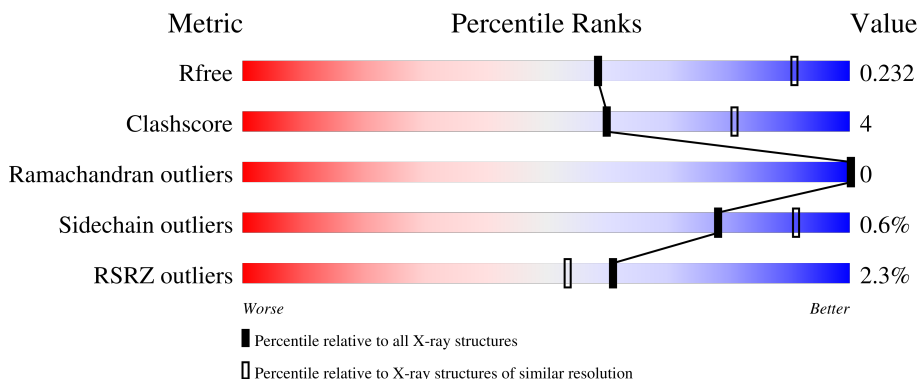
1 Overall quality at a glance i

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}	180053	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (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 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	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	<p>87% 11%</p>
1	G	202	<p>84% 13%</p>
1	H	202	<p>84% 13%</p>
1	I	202	<p>84% 10% 5%</p>
1	J	202	<p>84% 13%</p>
1	K	202	<p>82% 13%</p>
1	L	202	<p>84% 10% 5%</p>
1	M	202	<p>84% 12%</p>
1	N	202	<p>84% 9% 7%</p>

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

There are 14 discrepancies between the modelled and reference sequences:

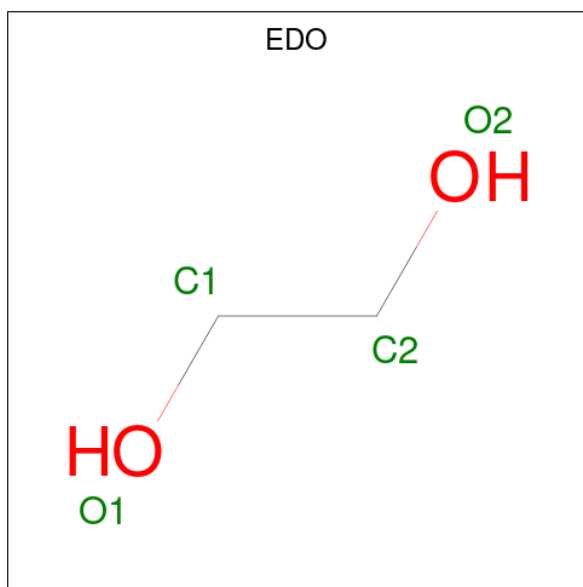
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP A7LXU7

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

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: 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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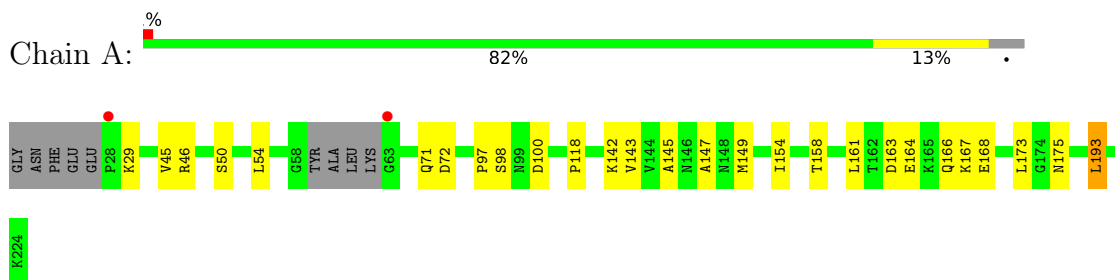
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	I	17	Total O 17 17	0	0
3	J	12	Total O 12 12	0	0
3	K	14	Total O 14 14	0	0
3	L	8	Total O 8 8	0	0
3	M	10	Total O 10 10	0	0
3	N	10	Total O 10 10	0	0

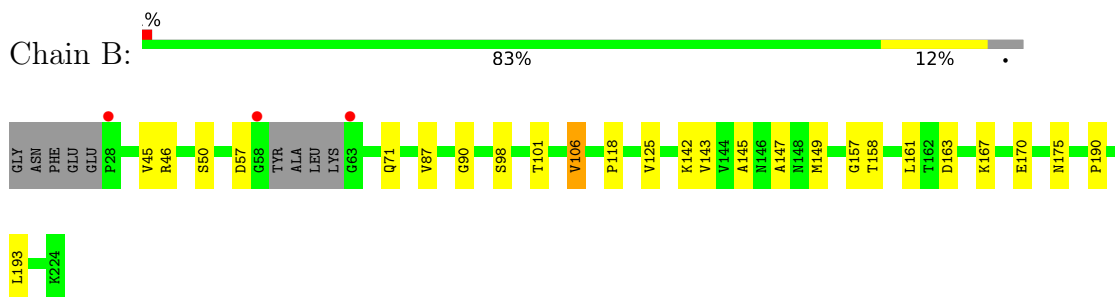
3 Residue-property plots

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.

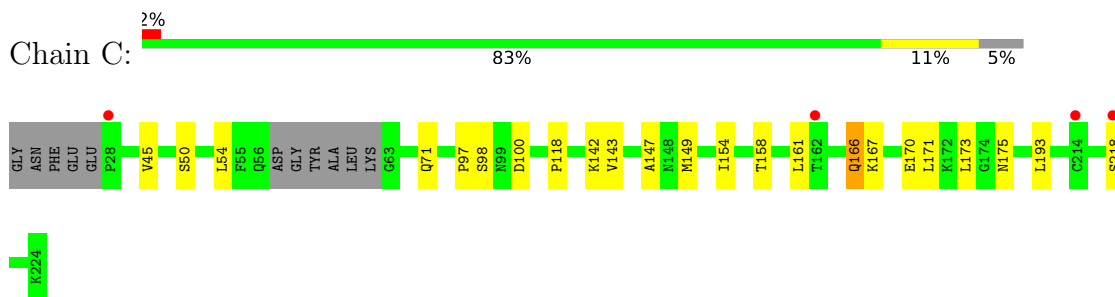
- Molecule 1: Uncharacterized protein



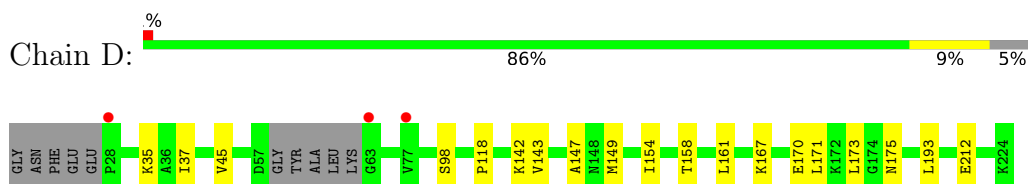
- Molecule 1: Uncharacterized protein



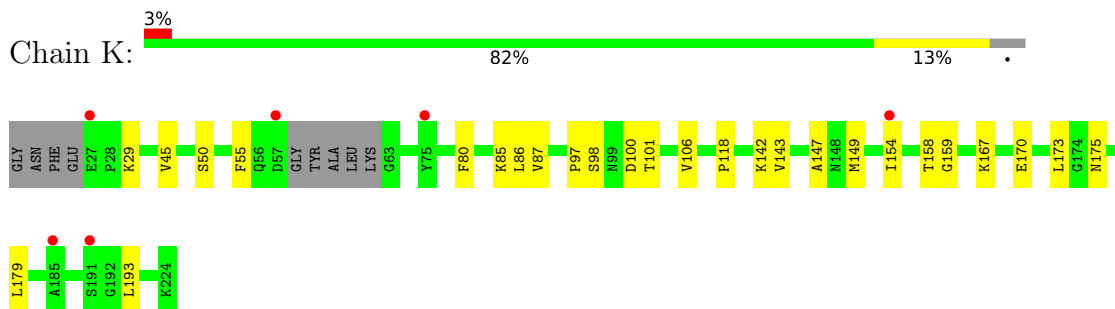
- Molecule 1: Uncharacterized protein



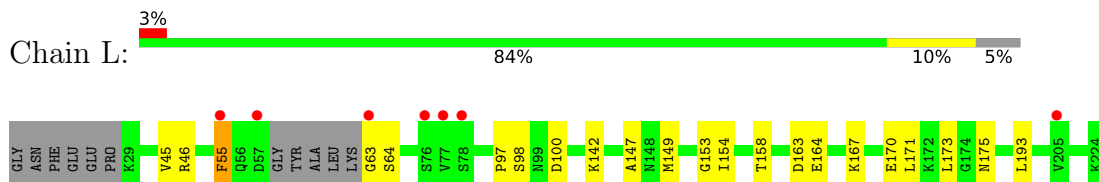
- Molecule 1: Uncharacterized protein



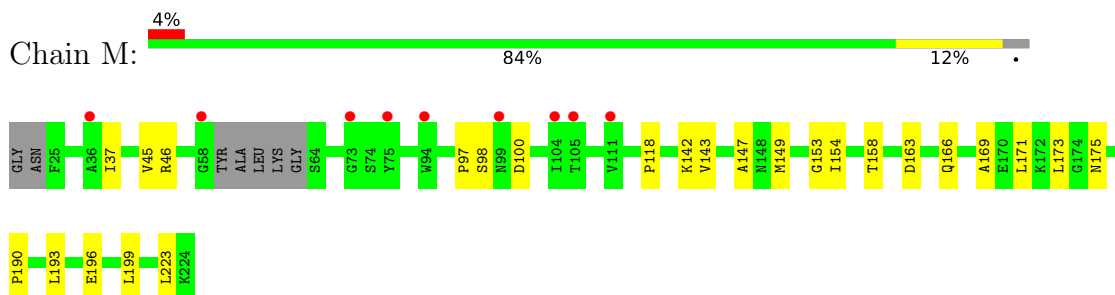
- Molecule 1: Uncharacterized protein



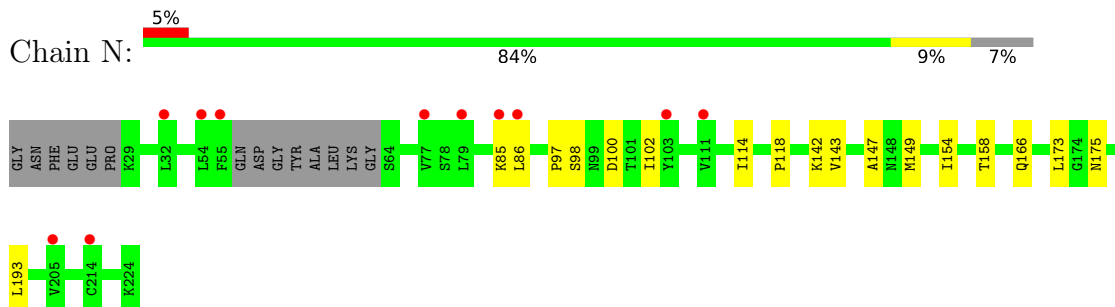
- Molecule 1: Uncharacterized protein



- 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 49.08 – 2.77	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.08-2.77) 99.8 (49.08-2.77)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.77Å)	Xtrriage
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 (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	58.9	Xtrriage
Anisotropy	0.607	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 80.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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	0.92	0/1498	1.21	7/2025 (0.3%)
1	B	0.93	0/1515	1.23	9/2046 (0.4%)
1	C	0.84	0/1483	1.16	3/2005 (0.1%)
1	D	0.83	0/1492	1.18	2/2016 (0.1%)
1	E	0.85	0/1571	1.21	13/2125 (0.6%)
1	F	0.84	0/1539	1.18	4/2079 (0.2%)
1	G	0.83	0/1529	1.22	11/2068 (0.5%)
1	H	0.92	1/1526 (0.1%)	1.21	4/2064 (0.2%)
1	I	0.84	1/1490 (0.1%)	1.14	2/2013 (0.1%)
1	J	0.86	0/1534	1.24	9/2074 (0.4%)
1	K	0.85	0/1508	1.20	4/2038 (0.2%)
1	L	0.84	1/1497 (0.1%)	1.18	3/2021 (0.1%)
1	M	0.86	0/1530	1.17	2/2067 (0.1%)
1	N	0.86	0/1471	1.18	2/1988 (0.1%)
All	All	0.86	3/21183 (0.0%)	1.20	75/28629 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	45	VAL	CA-C	5.47	1.57	1.52
1	L	63	GLY	C-N	5.38	1.41	1.33
1	I	45	VAL	CA-C	5.19	1.58	1.52

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	130	ASN	CA-CB-CG	7.62	120.22	112.60
1	G	98	SER	CA-C-N	7.50	130.33	120.28
1	G	98	SER	C-N-CA	7.50	130.33	120.28
1	A	98	SER	CA-C-N	7.29	130.05	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	SER	C-N-CA	7.29	130.05	120.28
1	J	98	SER	CA-C-N	7.01	129.67	120.28
1	J	98	SER	C-N-CA	7.01	129.67	120.28
1	E	98	SER	CA-C-N	6.82	129.41	120.28
1	E	98	SER	C-N-CA	6.82	129.41	120.28
1	D	98	SER	CA-C-N	6.81	129.40	120.28
1	D	98	SER	C-N-CA	6.81	129.40	120.28
1	L	98	SER	CA-C-N	6.79	129.38	120.28
1	L	98	SER	C-N-CA	6.79	129.38	120.28
1	H	98	SER	CA-C-N	6.76	129.34	120.28
1	H	98	SER	C-N-CA	6.76	129.34	120.28
1	C	98	SER	CA-C-N	6.72	129.29	120.28
1	C	98	SER	C-N-CA	6.72	129.29	120.28
1	J	29	LYS	CA-C-N	6.67	130.63	120.82
1	J	29	LYS	C-N-CA	6.67	130.63	120.82
1	M	98	SER	CA-C-N	6.60	129.12	120.28
1	M	98	SER	C-N-CA	6.60	129.12	120.28
1	B	57	ASP	CA-CB-CG	6.57	119.17	112.60
1	N	98	SER	CA-C-N	6.54	129.05	120.28
1	N	98	SER	C-N-CA	6.54	129.05	120.28
1	J	59	TYR	CA-C-N	6.50	128.99	120.28
1	J	59	TYR	C-N-CA	6.50	128.99	120.28
1	A	145	ALA	CA-C-N	6.42	128.88	120.28
1	A	145	ALA	C-N-CA	6.42	128.88	120.28
1	B	98	SER	CA-C-N	6.41	129.52	120.28
1	B	98	SER	C-N-CA	6.41	129.52	120.28
1	I	98	SER	CA-C-N	6.19	128.57	120.28
1	I	98	SER	C-N-CA	6.19	128.57	120.28
1	G	60	ALA	CA-C-N	6.17	129.68	120.31
1	G	60	ALA	C-N-CA	6.17	129.68	120.31
1	K	98	SER	CA-C-N	6.15	128.52	120.28
1	K	98	SER	C-N-CA	6.15	128.52	120.28
1	J	72	ASP	CA-CB-CG	6.03	118.63	112.60
1	F	98	SER	CA-C-N	5.99	128.31	120.28
1	F	98	SER	C-N-CA	5.99	128.31	120.28
1	L	55	PHE	CA-CB-CG	5.91	119.70	113.80
1	B	106	VAL	CA-C-N	5.86	132.25	123.23
1	B	106	VAL	C-N-CA	5.86	132.25	123.23
1	A	29	LYS	CA-C-N	5.82	131.30	120.95
1	A	29	LYS	C-N-CA	5.82	131.30	120.95
1	B	145	ALA	CA-C-N	5.79	128.04	120.28
1	B	145	ALA	C-N-CA	5.79	128.04	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	190	PRO	CA-C-N	5.77	129.47	120.82
1	B	190	PRO	C-N-CA	5.77	129.47	120.82
1	K	159	GLY	CA-C-N	5.71	128.59	120.53
1	K	159	GLY	C-N-CA	5.71	128.59	120.53
1	E	72	ASP	CA-CB-CG	5.62	118.22	112.60
1	H	145	ALA	CA-C-N	5.54	128.16	120.29
1	H	145	ALA	C-N-CA	5.54	128.16	120.29
1	E	106	VAL	CA-C-N	5.50	130.87	122.95
1	E	106	VAL	C-N-CA	5.50	130.87	122.95
1	E	60	ALA	CA-C-N	5.48	127.89	120.38
1	E	60	ALA	C-N-CA	5.48	127.89	120.38
1	G	72	ASP	CA-CB-CG	5.47	118.07	112.60
1	E	159	GLY	CA-C-N	5.43	127.60	120.60
1	E	159	GLY	C-N-CA	5.43	127.60	120.60
1	A	72	ASP	CA-CB-CG	5.41	118.01	112.60
1	C	166	GLN	CB-CG-CD	-5.34	103.52	112.60
1	G	64	SER	CA-C-N	5.32	131.98	122.13
1	G	64	SER	C-N-CA	5.32	131.98	122.13
1	E	145	ALA	CA-C-N	5.27	127.34	120.28
1	E	145	ALA	C-N-CA	5.27	127.34	120.28
1	E	123	ARG	CA-C-N	5.08	130.15	122.68
1	E	123	ARG	C-N-CA	5.08	130.15	122.68
1	G	145	ALA	CA-C-N	5.08	127.51	120.29
1	G	145	ALA	C-N-CA	5.08	127.51	120.29
1	G	190	PRO	CA-C-N	5.08	127.34	120.38
1	G	190	PRO	C-N-CA	5.08	127.34	120.38
1	F	145	ALA	CA-C-N	5.04	127.44	120.29
1	F	145	ALA	C-N-CA	5.04	127.44	120.29
1	J	166	GLN	CB-CG-CD	-5.03	104.06	112.60

There are no chirality outliers.

There are no planarity outliers.

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	1475	0	1448	15	0
1	B	1489	0	1471	13	0
1	C	1460	0	1432	15	0
1	D	1470	0	1449	12	0
1	E	1545	0	1518	19	0
1	F	1515	0	1488	15	0
1	G	1505	0	1483	15	0
1	H	1502	0	1479	15	0
1	I	1468	0	1448	11	0
1	J	1509	0	1480	10	0
1	K	1485	0	1465	13	0
1	L	1473	0	1460	13	0
1	M	1506	0	1483	14	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	180	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.

All (180) 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
1:C:161:LEU:HD11	1:D:171:LEU:HD11	1.79	0.65
1:B:161:LEU:HD11	1:C:171:LEU:HD11	1.80	0.63
1:K:167:LYS:HD2	1:K:170:GLU:HB3	1.80	0.63
1:B:157:GLY:HA3	2:B:301:EDO:H12	1.81	0.62
1:L:164:GLU:O	1:L:167:LYS:HE3	2.00	0.61
1:A:158:THR:HG22	1:A:193:LEU:HD11	1.84	0.59
1:G:158:THR:HG22	1:G:193:LEU:HD13	1.85	0.58
1:M:169:ALA:HB2	1:M:190:PRO:HG3	1.85	0.58
1:K:55:PHE:HB2	1:K:85:LYS:HG3	1.86	0.57
1:I:142:LYS:HE3	1:I:147:ALA:O	2.05	0.57
1:I:158:THR:HG22	1:I:193:LEU:HD13	1.87	0.57
1:M:142:LYS:HE3	1:M:147:ALA:O	2.05	0.56
1:L:158:THR:HG22	1:L:193:LEU:HD13	1.86	0.56
1:C:142:LYS:HE3	1:C:147:ALA:O	2.06	0.56
1:K:142:LYS:HE3	1:K:147:ALA:O	2.06	0.56
1:E:158:THR:HG22	1:E:193:LEU:HD13	1.89	0.55
1:G:97:PRO:HG2	1:G:100:ASP:HB3	1.87	0.55
1:K:87:VAL:CG1	1:K:101:THR:HG22	2.37	0.55
1:M:158:THR:HG22	1:M:193:LEU:HD13	1.88	0.55
1:B:142:LYS:HE3	1:B:147:ALA:O	2.07	0.54
1:H:142:LYS:HE3	1:H:147:ALA:O	2.07	0.54
1:L:142:LYS:HE3	1:L:147:ALA:O	2.08	0.54
1:J:142:LYS:HE3	1:J:147:ALA:O	2.07	0.54
1:A:142:LYS:HE3	1:A:147:ALA:O	2.08	0.54
1:J:158:THR:HG22	1:J:193:LEU:HD13	1.90	0.54
1:D:142:LYS:HE3	1:D:147:ALA:O	2.06	0.54
1:E:142:LYS:HE3	1:E:147:ALA:O	2.07	0.54
1:B:167:LYS:HD2	1:B:170:GLU:HB3	1.91	0.53
1:K:118:PRO:O	1:K:143:VAL:HB	2.08	0.53
1:A:164:GLU:O	1:A:167:LYS:HE3	2.08	0.53
1:F:142:LYS:HE3	1:F:147:ALA:O	2.09	0.53
1:L:154:ILE:HD11	1:L:173:LEU:HD11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:GLY:HA3	3:B:430:HOH:O	2.10	0.52
1:I:154:ILE:HD11	1:I:173:LEU:HD11	1.91	0.52
1:D:161:LEU:CD1	1:G:171:LEU:HD11	2.39	0.52
1:L:167:LYS:HD2	1:L:170:GLU:HB3	1.90	0.52
1:N:142:LYS:HE3	1:N:147:ALA:O	2.09	0.52
1:G:142:LYS:HE3	1:G:147:ALA:O	2.10	0.51
1:K:154:ILE:HD11	1:K:173:LEU:HD11	1.92	0.51
1:B:71:GLN:NE2	3:B:407:HOH:O	2.43	0.51
1:F:154:ILE:HD11	1:F:173:LEU:HD11	1.93	0.51
1:I:191:SER:O	1:I:194:VAL:HG12	2.11	0.51
1:D:158:THR:HG22	1:D:193:LEU:HD13	1.92	0.51
1:B:158:THR:HG22	1:B:193:LEU:HD13	1.91	0.51
1:A:154:ILE:HD11	1:A:173:LEU:HD11	1.92	0.50
1:M:199:LEU:HD12	1:M:199:LEU:C	2.36	0.50
1:D:118:PRO:O	1:D:143:VAL:HB	2.12	0.50
1:E:149:MSE:CE	3:E:408:HOH:O	2.59	0.50
1:C:71:GLN:HE22	1:C:218:SER:HB3	1.77	0.50
1:J:199:LEU:C	1:J:199:LEU:HD12	2.37	0.49
1:K:158:THR:HG22	1:K:193:LEU:HD21	1.95	0.49
1:J:154:ILE:HD11	1:J:173:LEU:HD11	1.95	0.49
1:I:99:ASN:ND2	3:I:406:HOH:O	2.46	0.49
1:H:154:ILE:HD11	1:H:173:LEU:HD11	1.95	0.48
1:M:154:ILE:HD11	1:M:173:LEU:HD11	1.94	0.48
1:N:158:THR:HG22	1:N:193:LEU:HD13	1.95	0.48
1:D:154:ILE:HD11	1:D:173:LEU:HD11	1.94	0.48
1:F:158:THR:HG22	1:F:193:LEU:HD13	1.95	0.48
1:K:29:LYS:HE2	1:K:80:PHE:CE2	2.49	0.48
1:C:149:MSE:HE3	1:C:175:ASN:O	2.14	0.48
1:F:161:LEU:HD11	1:I:171:LEU:HD11	1.95	0.48
1:F:158:THR:H	2:F:301:EDO:C2	2.25	0.47
1:N:85:LYS:HA	1:N:102:ILE:O	2.14	0.47
1:N:154:ILE:HD11	1:N:173:LEU:HD11	1.97	0.47
1:F:35:LYS:HE3	1:F:37:ILE:HD11	1.96	0.47
1:F:97:PRO:HG2	1:F:100:ASP:HB3	1.96	0.47
1:I:167:LYS:HD2	1:I:170:GLU:HB3	1.96	0.47
1:E:154:ILE:HD11	1:E:173:LEU:HD11	1.97	0.47
1:C:154:ILE:HD11	1:C:173:LEU:HD11	1.97	0.47
1:F:167:LYS:HD2	1:F:170:GLU:HB3	1.96	0.47
1:H:149:MSE:HE3	1:H:175:ASN:O	2.14	0.47
1:M:166:GLN:HB3	2:M:301:EDO:H11	1.96	0.47
1:N:118:PRO:O	1:N:143:VAL:HB	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:97:PRO:HG2	1:J:100:ASP:HB3	1.95	0.47
1:I:45:VAL:HG21	1:I:50:SER:OG	2.15	0.47
1:K:149:MSE:HE3	1:K:175:ASN:O	2.14	0.47
1:F:149:MSE:HE3	1:F:175:ASN:O	2.14	0.46
1:I:85:LYS:HE2	1:I:103:TYR:HE2	1.79	0.46
1:H:31:THR:HG22	1:H:33:THR:HG23	1.97	0.46
1:J:118:PRO:O	1:J:143:VAL:HB	2.14	0.46
1:E:35:LYS:HE3	1:E:37:ILE:HD11	1.98	0.45
1:E:199:LEU:HD12	1:E:199:LEU:C	2.41	0.45
1:D:167:LYS:HD2	1:D:170:GLU:HB3	1.98	0.45
1:H:61:LEU:HD23	1:H:61:LEU:H	1.81	0.45
1:C:71:GLN:HE22	1:C:218:SER:CB	2.30	0.45
1:L:149:MSE:HE3	1:L:175:ASN:O	2.17	0.45
1:L:158:THR:HG22	1:L:193:LEU:CD1	2.47	0.45
1:C:166:GLN:HB3	2:C:301:EDO:H11	1.99	0.45
1:E:97:PRO:HG2	1:E:100:ASP:HB3	1.98	0.45
1:C:161:LEU:CD1	1:D:171:LEU:HD11	2.46	0.45
1:N:97:PRO:HG2	1:N:100:ASP:HB3	1.98	0.45
1:E:149:MSE:HE1	3:E:408:HOH:O	2.17	0.45
1:M:97:PRO:HG2	1:M:100:ASP:HB3	1.99	0.45
1:A:161:LEU:HD11	1:H:171:LEU:HD11	1.98	0.44
1:A:97:PRO:HG2	1:A:100:ASP:HB3	1.98	0.44
1:A:149:MSE:HE3	1:A:175:ASN:O	2.17	0.44
1:G:149:MSE:HE3	1:G:175:ASN:C	2.42	0.44
1:H:97:PRO:HG2	1:H:100:ASP:HB3	2.00	0.44
1:I:97:PRO:HG2	1:I:100:ASP:HB3	2.00	0.44
1:A:158:THR:HG22	1:A:193:LEU:CD1	2.46	0.44
1:G:154:ILE:HD11	1:G:173:LEU:HD11	2.00	0.44
1:M:149:MSE:HE3	1:M:175:ASN:O	2.17	0.44
1:J:45:VAL:HG21	1:J:50:SER:OG	2.18	0.44
1:L:97:PRO:HG2	1:L:100:ASP:HB3	1.99	0.44
1:M:118:PRO:O	1:M:143:VAL:HB	2.18	0.44
1:M:149:MSE:HE3	1:M:175:ASN:C	2.43	0.43
1:B:149:MSE:HE3	1:B:175:ASN:C	2.44	0.43
1:J:149:MSE:HE3	1:J:175:ASN:O	2.18	0.43
1:C:158:THR:HG22	1:C:193:LEU:HD13	2.00	0.43
1:A:45:VAL:HG21	1:A:50:SER:OG	2.19	0.43
1:H:71:GLN:OE1	1:H:161:LEU:HA	2.19	0.43
1:N:166[A]:GLN:HB3	2:N:301:EDO:H22	2.00	0.43
1:B:125:VAL:H	2:B:302:EDO:H11	1.84	0.43
1:C:97:PRO:HG2	1:C:100:ASP:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:53:ALA:HB3	1:E:87:VAL:HG23	1.99	0.43
1:H:153:GLY:HA2	1:H:171:LEU:O	2.18	0.43
1:I:118:PRO:O	1:I:143:VAL:HB	2.18	0.43
1:K:97:PRO:HG2	1:K:100:ASP:HB3	2.00	0.43
1:E:167:LYS:HD2	1:E:170:GLU:HB3	2.01	0.43
1:M:45:VAL:HG22	1:M:46:ARG:H	1.83	0.43
1:D:35:LYS:HE3	1:D:37:ILE:HD11	2.00	0.42
1:H:149:MSE:HE3	1:H:175:ASN:C	2.44	0.42
1:E:125:VAL:HG21	1:E:216:SER:CB	2.48	0.42
1:H:45:VAL:HG22	1:H:46:ARG:H	1.83	0.42
1:F:71:GLN:OE1	1:F:161:LEU:HA	2.19	0.42
1:G:45:VAL:HG21	1:G:50:SER:OG	2.20	0.42
1:E:85:LYS:NZ	1:E:103:TYR:HE2	2.17	0.42
1:E:149:MSE:HE3	1:E:175:ASN:C	2.44	0.42
1:J:153:GLY:HA2	1:J:171:LEU:O	2.19	0.42
1:H:158:THR:HG22	1:H:193:LEU:HD13	2.00	0.42
1:C:45:VAL:HG21	1:C:50:SER:OG	2.19	0.42
1:E:45:VAL:HG21	1:E:50:SER:OG	2.20	0.42
1:K:45:VAL:HG21	1:K:50:SER:OG	2.20	0.42
1:B:87:VAL:CG1	1:B:101:THR:HG22	2.49	0.42
1:K:86:LEU:C	1:K:86:LEU:HD23	2.43	0.42
1:B:118:PRO:O	1:B:143:VAL:HB	2.20	0.42
1:A:54:LEU:HD12	1:A:54:LEU:HA	1.92	0.42
1:E:54:LEU:HD21	1:E:79:LEU:HD11	2.02	0.42
1:L:55:PHE:HB3	1:L:64:SER:HA	2.02	0.42
1:N:149:MSE:HE3	1:N:175:ASN:O	2.20	0.42
1:D:212:GLU:HG3	3:D:401:HOH:O	2.19	0.41
1:E:25:PHE:HB2	1:L:64:SER:O	2.20	0.41
1:A:45:VAL:HG22	1:A:46:ARG:H	1.85	0.41
1:F:45:VAL:HG21	1:F:50:SER:OG	2.20	0.41
1:G:45:VAL:HG22	1:G:46:ARG:H	1.85	0.41
1:C:149:MSE:HE3	1:C:175:ASN:C	2.45	0.41
1:D:149:MSE:HE3	1:D:175:ASN:O	2.20	0.41
1:M:46:ARG:HG3	1:M:163:ASP:CG	2.45	0.41
1:B:46:ARG:HG3	1:B:163:ASP:CG	2.45	0.41
1:G:167:LYS:HD2	1:G:170:GLU:HB3	2.03	0.41
1:F:46:ARG:HG3	1:F:163:ASP:CG	2.45	0.41
1:K:149:MSE:HE3	1:K:175:ASN:C	2.46	0.41
1:A:46:ARG:HG3	1:A:163:ASP:CG	2.46	0.41
1:A:118:PRO:O	1:A:143:VAL:HB	2.20	0.41
1:B:45:VAL:HG21	1:B:50:SER:OG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:PRO:O	1:C:143:VAL:HB	2.20	0.41
1:L:153:GLY:HA2	1:L:171:LEU:O	2.21	0.41
1:M:153:GLY:HA2	1:M:171:LEU:O	2.21	0.40
1:N:86:LEU:HD13	1:N:114:ILE:HG13	2.03	0.40
1:C:167:LYS:HD2	1:C:170:GLU:HB3	2.03	0.40
1:F:149:MSE:HE3	1:F:175:ASN:C	2.45	0.40
1:G:54:LEU:CD2	1:G:79:LEU:HD11	2.51	0.40
1:G:149:MSE:HE3	1:G:175:ASN:O	2.21	0.40
1:E:125:VAL:HG21	1:E:216:SER:HB2	2.03	0.40
1:G:61:LEU:N	1:G:61:LEU:HD12	2.36	0.40
1:H:198:TYR:CG	1:H:220:LYS:HE3	2.56	0.40
1:E:198:TYR:CG	1:E:220:LYS:HE3	2.56	0.40
1:G:153:GLY:HA2	1:G:171:LEU:O	2.21	0.40
1:L:45:VAL:HG22	1:L:46:ARG:H	1.86	0.40
1:L:46:ARG:HG3	1:L:163:ASP:CG	2.47	0.40
1:A:166:GLN:HE21	1:H:133:LYS:NZ	2.20	0.40
1:F:45:VAL:HG22	1:F:46:ARG:H	1.86	0.40
1:G:46:ARG:HG3	1:G:163:ASP:CG	2.46	0.40
1:J:158:THR:HG22	1:J:193:LEU:CD1	2.52	0.40
1:M:196:GLU:O	1:M:223:LEU:HB2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
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
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 [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	160/167 (96%)	158 (99%)	2 (1%)	61	83
1	B	162/167 (97%)	161 (99%)	1 (1%)	78	91
1	C	158/167 (95%)	157 (99%)	1 (1%)	78	91
1	D	160/167 (96%)	159 (99%)	1 (1%)	78	91
1	E	165/167 (99%)	164 (99%)	1 (1%)	78	91
1	F	164/167 (98%)	164 (100%)	0	100	100
1	G	162/167 (97%)	162 (100%)	0	100	100
1	H	162/167 (97%)	161 (99%)	1 (1%)	78	91
1	I	160/167 (96%)	160 (100%)	0	100	100
1	J	162/167 (97%)	158 (98%)	4 (2%)	42	73
1	K	161/167 (96%)	159 (99%)	2 (1%)	63	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	161/167 (96%)	161 (100%)	0	100	100
1	M	164/167 (98%)	163 (99%)	1 (1%)	78	91
1	N	158/167 (95%)	158 (100%)	0	100	100
All	All	2259/2338 (97%)	2245 (99%)	14 (1%)	78	91

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	193	LEU
1	B	106	VAL
1	C	54	LEU
1	D	45	VAL
1	E	179	LEU
1	H	106	VAL
1	J	37	ILE
1	J	54	LEU
1	J	112	GLN
1	J	167	LYS
1	K	106	VAL
1	K	179	LEU
1	M	37	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	148	ASN
1	A	166	GLN
1	A	217	GLN
1	B	91	ASN
1	B	99	ASN
1	B	148	ASN
1	B	151	ASN
1	C	71	GLN
1	C	148	ASN
1	C	151	ASN
1	C	166	GLN
1	D	71	GLN
1	D	132	ASN

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Mol	Chain	Res	Type
1	D	151	ASN
1	E	132	ASN
1	E	148	ASN
1	E	151	ASN
1	F	91	ASN
1	F	99	ASN
1	F	148	ASN
1	F	183	ASN
1	F	217	GLN
1	G	183	ASN
1	G	217	GLN
1	H	132	ASN
1	H	146	ASN
1	I	148	ASN
1	J	112	GLN
1	J	166	GLN
1	J	217	GLN
1	K	132	ASN
1	K	151	ASN
1	K	166	GLN
1	L	217	GLN
1	M	91	ASN
1	M	151	ASN
1	M	166	GLN
1	M	217	GLN
1	N	71	GLN
1	N	109	ASN
1	N	151	ASN
1	N	183	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	K	301	-	3,3,3	0.61	0	2,2,2	0.16	0
2	EDO	I	301	-	3,3,3	0.63	0	2,2,2	0.26	0
2	EDO	A	301	-	3,3,3	0.70	0	2,2,2	0.20	0
2	EDO	B	302	-	3,3,3	1.02	0	2,2,2	0.06	0
2	EDO	H	301	-	3,3,3	0.79	0	2,2,2	0.24	0
2	EDO	B	301	-	3,3,3	0.59	0	2,2,2	0.08	0
2	EDO	G	301	-	3,3,3	0.74	0	2,2,2	0.15	0
2	EDO	F	301	-	3,3,3	0.58	0	2,2,2	0.37	0
2	EDO	L	301	-	3,3,3	0.69	0	2,2,2	0.23	0
2	EDO	D	301	-	3,3,3	0.59	0	2,2,2	0.11	0
2	EDO	M	301	-	3,3,3	0.58	0	2,2,2	0.35	0
2	EDO	E	301	-	3,3,3	0.63	0	2,2,2	0.07	0
2	EDO	J	301	-	3,3,3	0.64	0	2,2,2	0.13	0
2	EDO	N	301	-	3,3,3	0.66	0	2,2,2	0.09	0
2	EDO	C	301	-	3,3,3	0.61	0	2,2,2	0.13	0
2	EDO	B	303	-	3,3,3	0.63	0	2,2,2	0.15	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	K	301	-	-	0/1/1/1	-
2	EDO	I	301	-	-	0/1/1/1	-
2	EDO	A	301	-	-	0/1/1/1	-
2	EDO	B	302	-	-	0/1/1/1	-
2	EDO	H	301	-	-	0/1/1/1	-
2	EDO	B	301	-	-	0/1/1/1	-
2	EDO	G	301	-	-	0/1/1/1	-
2	EDO	F	301	-	-	1/1/1/1	-
2	EDO	L	301	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	D	301	-	-	0/1/1/1	-
2	EDO	M	301	-	-	0/1/1/1	-
2	EDO	E	301	-	-	0/1/1/1	-
2	EDO	J	301	-	-	0/1/1/1	-
2	EDO	N	301	-	-	0/1/1/1	-
2	EDO	C	301	-	-	0/1/1/1	-
2	EDO	B	303	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	301	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	EDO	1	0
2	B	301	EDO	1	0
2	F	301	EDO	2	0
2	M	301	EDO	1	0
2	N	301	EDO	1	0
2	C	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%)	-0.06	2 (1%) 79 74	38, 57, 102, 142	0
1	B	191/202 (94%)	-0.03	3 (1%) 70 63	32, 61, 87, 108	1 (0%)
1	C	189/202 (93%)	0.03	4 (2%) 63 56	41, 67, 104, 124	0
1	D	190/202 (94%)	0.01	3 (1%) 70 63	46, 68, 108, 140	0
1	E	200/202 (99%)	-0.01	3 (1%) 72 65	47, 68, 104, 132	0
1	F	195/202 (96%)	-0.02	2 (1%) 79 74	43, 66, 108, 132	0
1	G	194/202 (96%)	-0.03	0 100 100	43, 68, 94, 117	0
1	H	194/202 (96%)	0.05	2 (1%) 79 74	34, 61, 102, 122	0
1	I	189/202 (93%)	0.32	5 (2%) 57 49	50, 70, 117, 164	0
1	J	195/202 (96%)	-0.01	4 (2%) 63 56	43, 64, 102, 132	0
1	K	191/202 (94%)	0.26	6 (3%) 51 44	58, 75, 104, 123	0
1	L	189/202 (93%)	0.25	7 (3%) 45 38	42, 82, 108, 126	1 (0%)
1	M	193/202 (95%)	0.50	9 (4%) 36 31	62, 88, 141, 157	0
1	N	186/202 (92%)	0.43	11 (5%) 28 23	35, 83, 152, 170	1 (0%)
All	All	2687/2828 (95%)	0.12	61 (2%) 61 53	32, 70, 113, 170	3 (0%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	58	GLY	5.1
1	N	77	VAL	4.7
1	B	58	GLY	4.3
1	L	63	GLY	4.2
1	D	63	GLY	4.2
1	M	58	GLY	4.1
1	F	58	GLY	3.9
1	L	76	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	N	79	LEU	3.6
1	I	43	VAL	3.5
1	H	61	LEU	3.4
1	K	27	GLU	3.4
1	C	28	PRO	3.3
1	M	75	TYR	3.3
1	D	28	PRO	3.1
1	A	28	PRO	3.1
1	M	36	ALA	3.0
1	K	57	ASP	3.0
1	B	63	GLY	3.0
1	B	28	PRO	2.9
1	L	78	SER	2.8
1	H	60	ALA	2.8
1	L	77	VAL	2.8
1	A	63	GLY	2.7
1	N	86	LEU	2.7
1	N	205	VAL	2.6
1	J	28	PRO	2.5
1	C	214	CYS	2.5
1	E	26	GLU	2.5
1	M	73	GLY	2.5
1	C	218	SER	2.4
1	K	154	ILE	2.4
1	C	162	THR	2.4
1	K	191	SER	2.4
1	M	94	TRP	2.3
1	J	63	GLY	2.3
1	N	32	LEU	2.3
1	N	111	VAL	2.3
1	D	77	VAL	2.2
1	N	103	TYR	2.2
1	M	105	THR	2.2
1	L	205	VAL	2.2
1	E	160	ILE	2.2
1	K	185	ALA	2.2
1	I	103	TYR	2.2
1	N	85	LYS	2.2
1	I	44	GLY	2.2
1	M	104	ILE	2.1
1	N	55	PHE	2.1
1	J	60	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	J	59	TYR	2.1
1	F	24	ASN	2.1
1	N	214	CYS	2.1
1	L	57	ASP	2.1
1	N	54	LEU	2.1
1	E	24	ASN	2.1
1	I	101	THR	2.0
1	M	111	VAL	2.0
1	L	55	PHE	2.0
1	M	99	ASN	2.0
1	K	75	TYR	2.0

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 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
2	EDO	B	303	4/4	0.77	0.14	77,79,80,81	0
2	EDO	B	302	4/4	0.82	0.18	63,64,67,68	0
2	EDO	A	301	4/4	0.88	0.11	67,67,68,68	0
2	EDO	F	301	4/4	0.88	0.12	66,68,69,71	0
2	EDO	G	301	4/4	0.91	0.12	61,63,63,64	0
2	EDO	N	301	4/4	0.91	0.12	57,58,60,60	0
2	EDO	I	301	4/4	0.92	0.09	56,58,64,70	0
2	EDO	L	301	4/4	0.92	0.08	54,58,61,62	0
2	EDO	M	301	4/4	0.92	0.10	55,55,57,57	0
2	EDO	D	301	4/4	0.92	0.11	53,54,54,54	0
2	EDO	H	301	4/4	0.93	0.12	56,59,64,66	0
2	EDO	K	301	4/4	0.94	0.11	57,58,58,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	EDO	E	301	4/4	0.94	0.09	68,70,72,73	0
2	EDO	C	301	4/4	0.96	0.08	66,66,68,69	0
2	EDO	J	301	4/4	0.96	0.08	64,66,69,71	0
2	EDO	B	301	4/4	0.97	0.06	52,54,57,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.