



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2026 – 12:11 AM UTC

PDB ID : 8QTT / pdb_00008qtt
Title : Crystal structure of a C-terminally truncated version of Arabidopsis thaliana 14-3-3 omega in complex with a phosphopeptide from the inhibitor protein BKI1.
Authors : Hothorn, M.; Obergfell, E.
Deposited on : 2023-10-13
Resolution : 2.35 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

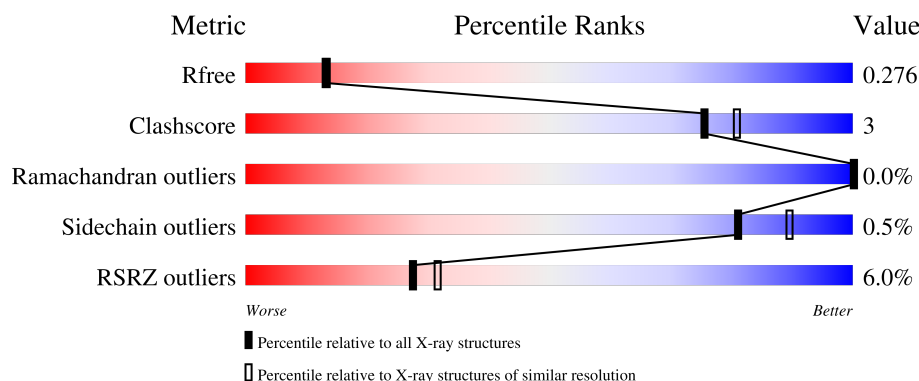
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	241	 5% 88% 9% .
1	B	241	 7% 88% 9% .
1	C	241	 7% 90% 6% .
1	D	241	 6% 91% 7% .
1	E	241	 4% 93% . .

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Mol	Chain	Length	Quality of chain
1	F	241	 5% 92% 6% •
1	G	241	 7% 90% 7% ••
1	H	241	 7% 91% • 6%
1	I	241	 3% 93% • •
1	J	241	 5% 91% 6% •
2	K	6	 17% 83% 17%
2	L	6	 33% 67% 17% 17%
2	M	6	 17% 83% 17%
2	N	6	 33% 67% 17% 17%
2	O	6	 83% 17%
2	P	6	 33% 83% 17%
2	Q	6	 33% 83% 17%
2	R	6	 83% 17%
2	S	6	 83% 17%
2	T	6	 17% 67% 17% 17%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 38323 atoms, of which 18751 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 14-3-3-like protein GF14 omega.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	234	Total	C	H	N	O	S	0	2	0
			3702	1172	1832	319	372	7			
1	B	234	Total	C	H	N	O	S	0	0	0
			3684	1167	1823	317	370	7			
1	C	232	Total	C	H	N	O	S	0	0	0
			3661	1160	1814	315	365	7			
1	D	236	Total	C	H	N	O	S	0	3	0
			3741	1183	1854	321	375	8			
1	E	234	Total	C	H	N	O	S	0	2	0
			3704	1173	1834	317	373	7			
1	F	237	Total	C	H	N	O	S	0	3	0
			3762	1188	1865	325	376	8			
1	G	236	Total	C	H	N	O	S	0	1	0
			3726	1178	1846	322	373	7			
1	H	227	Total	C	H	N	O	S	0	2	0
			3609	1143	1790	312	357	7			
1	I	234	Total	C	H	N	O	S	0	1	0
			3705	1172	1836	320	370	7			
1	J	234	Total	C	H	N	O	S	0	3	0
			3723	1177	1845	322	372	7			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	expression tag	UNP Q01525
A	0	GLY	-	expression tag	UNP Q01525
A	238	ARG	-	expression tag	UNP Q01525
A	239	GLY	-	expression tag	UNP Q01525
B	-1	ALA	-	expression tag	UNP Q01525
B	0	GLY	-	expression tag	UNP Q01525
B	238	ARG	-	expression tag	UNP Q01525
B	239	GLY	-	expression tag	UNP Q01525
C	-1	ALA	-	expression tag	UNP Q01525

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP Q01525
C	238	ARG	-	expression tag	UNP Q01525
C	239	GLY	-	expression tag	UNP Q01525
D	-1	ALA	-	expression tag	UNP Q01525
D	0	GLY	-	expression tag	UNP Q01525
D	238	ARG	-	expression tag	UNP Q01525
D	239	GLY	-	expression tag	UNP Q01525
E	-1	ALA	-	expression tag	UNP Q01525
E	0	GLY	-	expression tag	UNP Q01525
E	238	ARG	-	expression tag	UNP Q01525
E	239	GLY	-	expression tag	UNP Q01525
F	-1	ALA	-	expression tag	UNP Q01525
F	0	GLY	-	expression tag	UNP Q01525
F	238	ARG	-	expression tag	UNP Q01525
F	239	GLY	-	expression tag	UNP Q01525
G	-1	ALA	-	expression tag	UNP Q01525
G	0	GLY	-	expression tag	UNP Q01525
G	238	ARG	-	expression tag	UNP Q01525
G	239	GLY	-	expression tag	UNP Q01525
H	-1	ALA	-	expression tag	UNP Q01525
H	0	GLY	-	expression tag	UNP Q01525
H	238	ARG	-	expression tag	UNP Q01525
H	239	GLY	-	expression tag	UNP Q01525
I	-1	ALA	-	expression tag	UNP Q01525
I	0	GLY	-	expression tag	UNP Q01525
I	238	ARG	-	expression tag	UNP Q01525
I	239	GLY	-	expression tag	UNP Q01525
J	-1	ALA	-	expression tag	UNP Q01525
J	0	GLY	-	expression tag	UNP Q01525
J	238	ARG	-	expression tag	UNP Q01525
J	239	GLY	-	expression tag	UNP Q01525

- Molecule 2 is a protein called BRI1 kinase inhibitor 1.

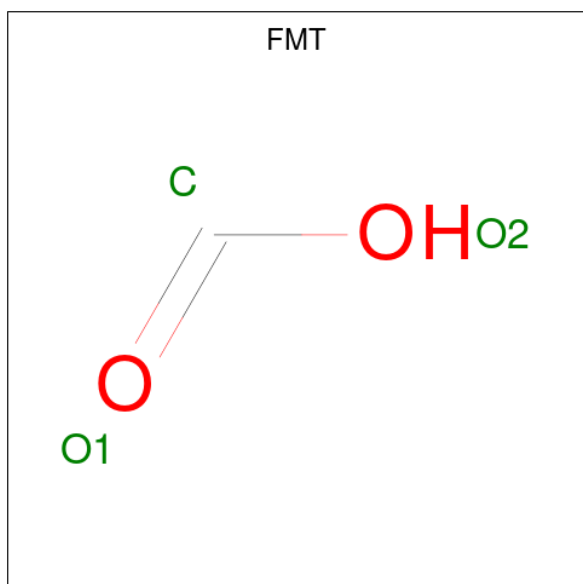
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	K	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	L	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	M	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	N	5	Total	C	H	N	O	P	0	0	0
			75	26	34	5	9	1			

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	O	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	P	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	Q	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	R	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	S	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			
2	T	6	Total	C	H	N	O	P	0	0	0
			90	31	40	6	12	1			

- Molecule 3 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



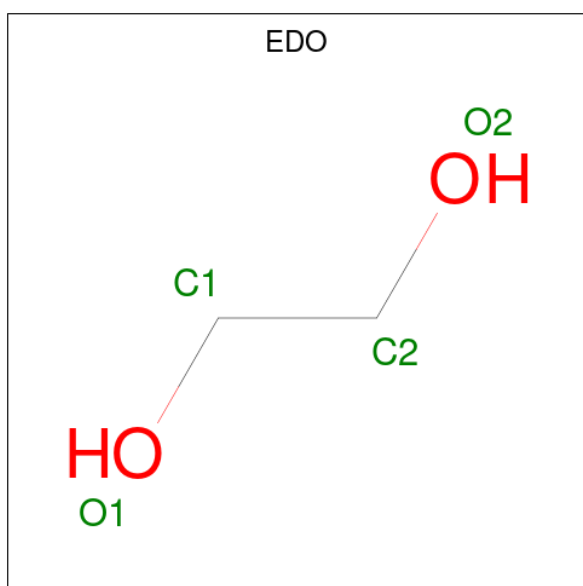
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	K	1	Total	C	H	O	0	0
			4	1	1	2		
3	B	1	Total	C	H	O	0	0
			4	1	1	2		
3	C	1	Total	C	H	O	0	0
			4	1	1	2		
3	D	1	Total	C	H	O	0	0
			4	1	1	2		
3	E	1	Total	C	H	O	0	0
			4	1	1	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	Q	1	Total	C	H	O	0	0
			4	1	1	2		
3	H	1	Total	C	H	O	0	0
			4	1	1	2		
3	S	1	Total	C	H	O	0	0
			4	1	1	2		
3	J	1	Total	C	H	O	0	0
			5	1	2	2		
3	J	1	Total	C	H	O	0	0
			5	1	2	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	E	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	35	Total	O	0	0
			35	35		
5	K	1	Total	O	0	0
			1	1		
5	B	35	Total	O	0	0
			35	35		

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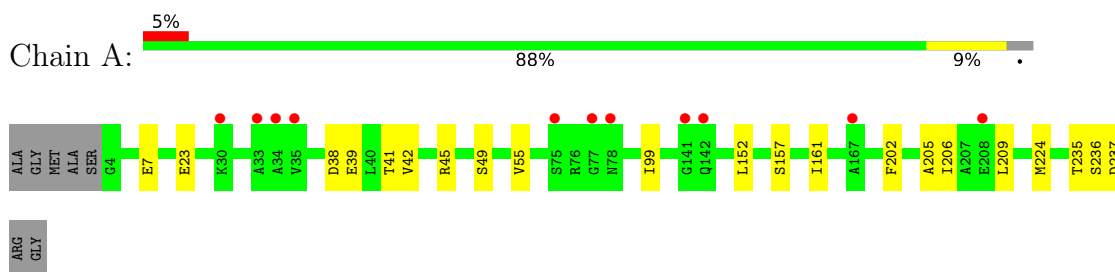
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	L	3	Total O 3 3	0	0
5	C	43	Total O 43 43	0	0
5	D	31	Total O 31 31	0	0
5	N	1	Total O 1 1	0	0
5	E	35	Total O 35 35	0	0
5	O	2	Total O 2 2	0	0
5	F	29	Total O 29 29	0	0
5	P	5	Total O 5 5	0	0
5	G	24	Total O 24 24	0	0
5	Q	1	Total O 1 1	0	0
5	H	33	Total O 33 33	0	0
5	R	3	Total O 3 3	0	0
5	I	44	Total O 44 44	0	0
5	S	3	Total O 3 3	0	0
5	J	38	Total O 38 38	0	0
5	T	3	Total O 3 3	0	0

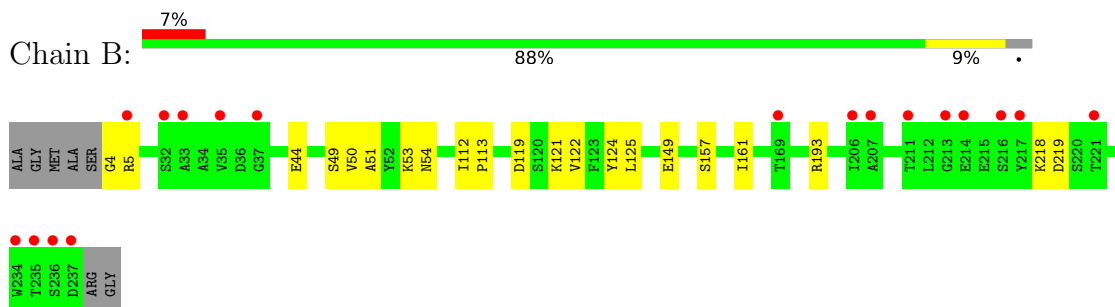
3 Residue-property plots [i](#)

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.

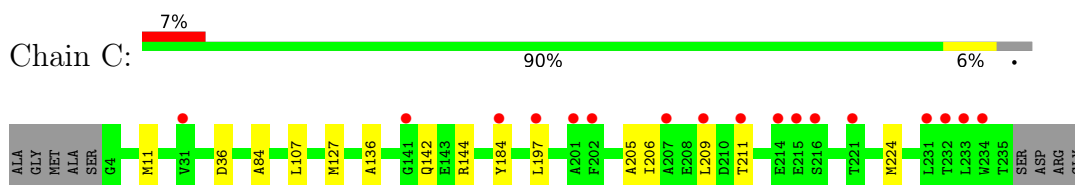
- Molecule 1: 14-3-3-like protein GF14 omega



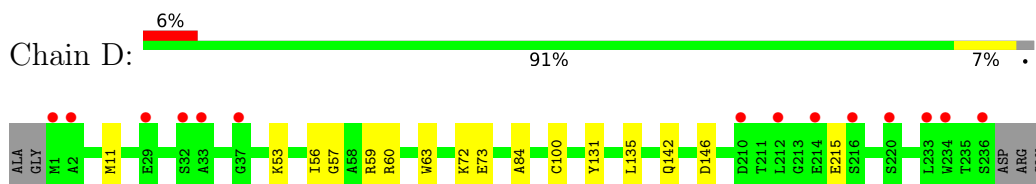
- Molecule 1: 14-3-3-like protein GF14 omega



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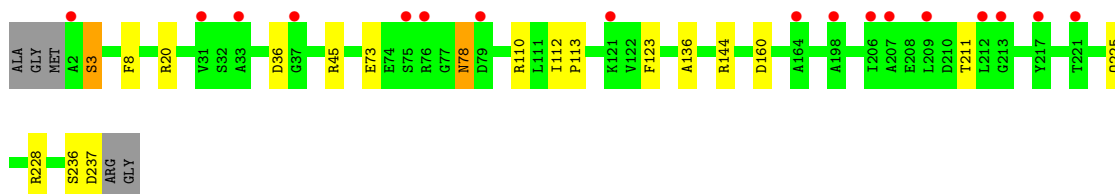
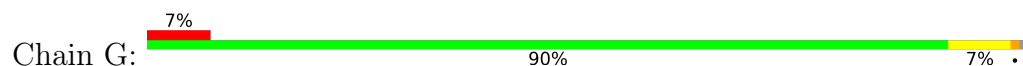




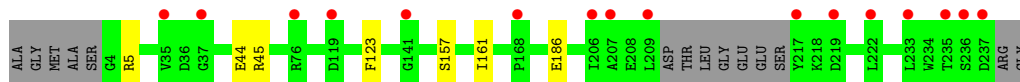
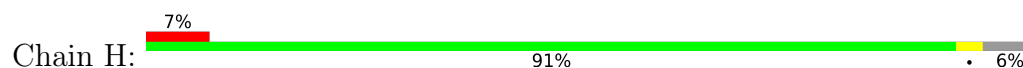
- Molecule 1: 14-3-3-like protein GF14 omega



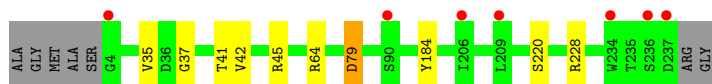
- Molecule 1: 14-3-3-like protein GF14 omega



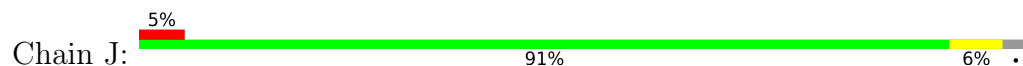
- Molecule 1: 14-3-3-like protein GF14 omega



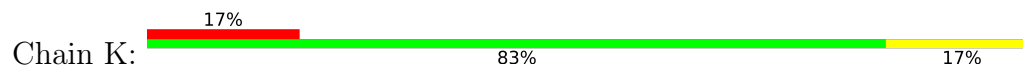
- Molecule 1: 14-3-3-like protein GF14 omega

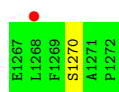


- Molecule 1: 14-3-3-like protein GF14 omega

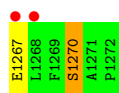


- Molecule 2: BRI1 kinase inhibitor 1

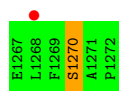
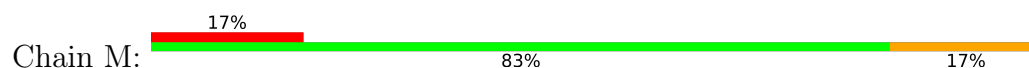




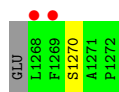
- Molecule 2: BRI1 kinase inhibitor 1



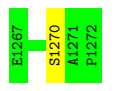
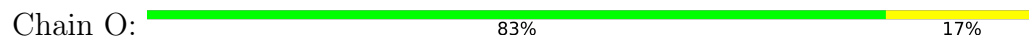
- Molecule 2: BRI1 kinase inhibitor 1



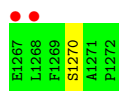
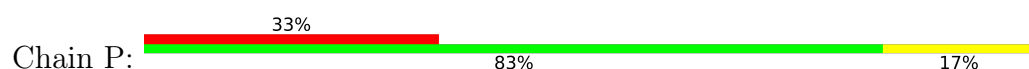
- Molecule 2: BRI1 kinase inhibitor 1



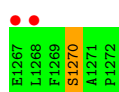
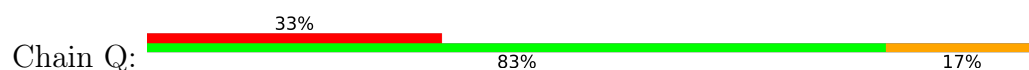
- Molecule 2: BRI1 kinase inhibitor 1



- Molecule 2: BRI1 kinase inhibitor 1

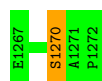


- Molecule 2: BRI1 kinase inhibitor 1



- Molecule 2: BRI1 kinase inhibitor 1

Chain R:  83% 17%



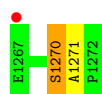
- Molecule 2: BRI1 kinase inhibitor 1

Chain S:  83% 17%



- Molecule 2: BRI1 kinase inhibitor 1

Chain T:  17% 67% 17% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.22Å 71.08Å 150.99Å 100.47° 95.53° 88.74°	Depositor
Resolution (Å)	50.33 – 2.35 50.33 – 2.35	Depositor EDS
% Data completeness (in resolution range)	97.9 (50.33-2.35) 98.0 (50.33-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.234 , 0.276 0.235 , 0.276	Depositor DCC
R_{free} test set	5917 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.3	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	0.94	EDS
Total number of atoms	38323	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.03% 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: SEP, FMT, 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	0.09	0/1908	0.23	0/2572
1	B	0.08	0/1890	0.21	0/2548
1	C	0.08	0/1876	0.21	0/2529
1	D	0.08	0/1928	0.21	0/2598
1	E	0.08	0/1905	0.24	0/2568
1	F	0.08	0/1938	0.21	0/2611
1	G	0.08	0/1912	0.22	0/2577
1	H	0.08	0/1856	0.21	0/2500
1	I	0.08	0/1901	0.20	0/2562
1	J	0.08	0/1919	0.21	0/2586
2	K	0.09	0/40	0.22	0/52
2	L	0.07	0/40	0.21	0/52
2	M	0.09	0/40	0.22	0/52
2	N	0.12	0/31	0.23	0/40
2	O	0.10	0/40	0.22	0/52
2	P	0.09	0/40	0.22	0/52
2	Q	0.11	0/40	0.26	0/52
2	R	0.10	0/40	0.23	0/52
2	S	0.12	0/40	0.18	0/52
2	T	0.11	0/40	0.28	0/52
All	All	0.08	0/19424	0.22	0/26159

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 ⓘ

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	1870	1832	1829	21	0
1	B	1861	1823	1827	13	1
1	C	1847	1814	1818	10	1
1	D	1887	1854	1853	16	0
1	E	1870	1834	1838	9	1
1	F	1897	1865	1864	10	1
1	G	1880	1846	1850	13	0
1	H	1819	1790	1786	4	0
1	I	1869	1836	1840	6	2
1	J	1878	1845	1842	11	0
2	K	50	40	40	0	0
2	L	50	40	41	2	0
2	M	50	40	40	1	0
2	N	41	34	34	0	0
2	O	50	40	40	0	0
2	P	50	40	41	0	0
2	Q	50	40	41	1	0
2	R	50	40	40	1	0
2	S	50	40	40	0	0
2	T	50	40	40	2	0
3	B	3	1	1	0	0
3	C	3	1	1	1	0
3	D	3	1	1	0	0
3	E	3	1	1	0	0
3	H	3	1	1	1	0
3	J	6	4	2	0	0
3	K	3	1	1	0	0
3	Q	3	1	1	0	0
3	S	3	1	1	0	0
4	E	4	6	6	0	0
5	A	35	0	0	4	0
5	B	35	0	0	9	0
5	C	43	0	0	1	0
5	D	31	0	0	7	0
5	E	35	0	0	1	0
5	F	29	0	0	1	0
5	G	24	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	H	33	0	0	0	0
5	I	44	0	0	5	0
5	J	38	0	0	6	0
5	K	1	0	0	0	0
5	L	3	0	0	2	0
5	N	1	0	0	0	0
5	O	2	0	0	0	0
5	P	5	0	0	0	0
5	Q	1	0	0	1	0
5	R	3	0	0	1	0
5	S	3	0	0	0	0
5	T	3	0	0	1	0
All	All	19572	18751	18760	112	3

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.

All (112) 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:45:ARG:O	5:A:1401:HOH:O	1.75	1.03
1:D:56:ILE:N	5:D:1401:HOH:O	1.99	0.96
1:D:63:TRP:N	5:D:1402:HOH:O	1.99	0.95
1:B:125:LEU:N	5:B:1402:HOH:O	1.98	0.95
1:B:50:VAL:O	5:B:1401:HOH:O	1.89	0.89
1:B:121:LYS:O	5:B:1402:HOH:O	1.93	0.86
2:L:1267:GLU:O	5:L:1301:HOH:O	1.95	0.84
1:D:59:ARG:O	5:D:1402:HOH:O	1.99	0.80
1:J:152:LEU:N	5:J:1402:HOH:O	2.15	0.80
2:R:1270:SEP:O1P	5:R:1401:HOH:O	2.03	0.76
2:L:1270:SEP:O1P	5:L:1302:HOH:O	2.04	0.75
1:I:220:SER:OG	5:I:1401:HOH:O	2.05	0.74
1:B:54:ASN:N	5:B:1401:HOH:O	2.05	0.72
1:B:53:LYS:N	5:B:1401:HOH:O	2.21	0.72
1:G:225:GLN:OE1	1:G:228:ARG:NH2	2.24	0.71
2:Q:1270:SEP:O1P	5:Q:1401:HOH:O	2.09	0.71
1:H:5:ARG:NH2	1:H:44:GLU:OE2	2.24	0.71
1:I:35:VAL:O	5:I:1402:HOH:O	2.10	0.69
1:J:140:THR:O	5:J:1401:HOH:O	2.10	0.69
1:A:209:LEU:HD11	1:A:224:MET:HE1	1.75	0.67
1:C:205:ALA:HB1	1:C:224:MET:HE2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:235:THR:N	5:E:1402:HOH:O	2.30	0.64
1:B:4:GLY:N	5:B:1403:HOH:O	2.33	0.62
1:J:60:ARG:NH2	2:T:1270:SEP:O3P	2.34	0.60
1:F:131:TYR:O	1:F:135:LEU:HD12	2.02	0.60
1:A:206:ILE:HA	1:A:209:LEU:HG	1.83	0.59
1:G:20:ARG:NE	5:G:301:HOH:O	2.38	0.57
3:C:1301:FMT:O2	2:M:1270:SEP:O2P	2.24	0.55
1:D:131:TYR:O	1:D:135:LEU:HD12	2.07	0.55
1:F:209:LEU:HD13	1:F:224:MET:HE1	1.89	0.54
1:G:136:ALA:O	1:G:144:ARG:NH1	2.40	0.54
1:F:145:LYS:NZ	5:F:1302:HOH:O	2.39	0.54
1:G:36:ASP:OD1	1:G:110:ARG:NH2	2.41	0.53
1:C:211:THR:HG23	1:G:160:ASP:HB3	1.89	0.53
1:B:5:ARG:NH2	1:B:44:GLU:OE2	2.42	0.53
1:I:228:ARG:NH2	5:I:1406:HOH:O	2.41	0.52
1:A:23:GLU:OE2	5:A:1402:HOH:O	2.19	0.52
1:F:205:ALA:HB1	1:F:224:MET:HE2	1.92	0.51
1:A:205:ALA:HB3	1:A:224:MET:HE3	1.92	0.51
1:C:11:MET:CE	1:D:84:ALA:HB1	2.41	0.51
1:D:56:ILE:HG22	5:D:1401:HOH:O	2.10	0.51
1:C:107:LEU:HD21	1:C:127:MET:HE2	1.94	0.50
1:C:136:ALA:O	1:C:144:ARG:NH1	2.44	0.50
1:I:64:ARG:NH1	5:I:1407:HOH:O	2.44	0.50
1:J:126:LYS:NZ	5:J:1407:HOH:O	2.44	0.50
1:D:53:LYS:C	5:D:1401:HOH:O	2.53	0.50
1:J:149:GLU:C	5:J:1402:HOH:O	2.55	0.50
1:B:124:TYR:N	5:B:1402:HOH:O	2.45	0.49
1:D:142[B]:GLN:NE2	1:D:146:ASP:OD2	2.45	0.49
1:J:142[B]:GLN:NE2	1:J:146:ASP:OD2	2.45	0.49
1:C:206:ILE:HA	1:C:209:LEU:HB2	1.93	0.49
1:A:152:LEU:HG	1:E:207:ALA:HB1	1.94	0.49
1:G:236:SER:O	1:G:237:ASP:HB2	2.13	0.48
1:G:3:SER:HB3	1:G:8:PHE:CE2	2.48	0.48
1:F:100:CYS:HB3	1:F:135:LEU:HD11	1.95	0.48
1:E:72:LYS:HD3	1:E:76:ARG:CZ	2.44	0.48
1:B:122:VAL:C	5:B:1402:HOH:O	2.58	0.47
1:F:142[B]:GLN:NE2	1:F:146:ASP:OD2	2.48	0.46
1:J:73:GLU:OE2	1:J:76:ARG:NH1	2.48	0.46
1:D:60:ARG:C	5:D:1402:HOH:O	2.58	0.46
1:D:72:LYS:NZ	1:D:73:GLU:OE2	2.48	0.46
1:D:100:CYS:CB	1:D:135:LEU:HD11	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:37:GLY:N	5:J:1403:HOH:O	2.38	0.46
1:I:37:GLY:N	5:I:1402:HOH:O	2.46	0.46
1:A:205:ALA:O	1:A:209:LEU:HG	2.16	0.46
1:A:209:LEU:HD21	1:A:224:MET:SD	2.55	0.46
1:C:84:ALA:HB1	1:D:11:MET:CE	2.46	0.46
1:A:236:SER:O	1:A:237:ASP:HB2	2.16	0.45
1:A:38:ASP:OD1	1:A:39:GLU:N	2.47	0.45
1:D:100:CYS:HB3	1:D:135:LEU:HD11	1.96	0.45
1:G:78:ASN:N	1:G:78:ASN:OD1	2.49	0.45
1:C:209:LEU:HD13	1:C:224:MET:HE1	1.99	0.45
1:D:215:GLU:OE1	1:D:215:GLU:N	2.50	0.44
1:J:153:ALA:N	5:J:1402:HOH:O	2.23	0.44
1:A:49:SER:CB	5:A:1401:HOH:O	2.65	0.44
1:G:73:GLU:OE1	1:G:78:ASN:ND2	2.50	0.44
1:E:84:ALA:HB1	1:F:11:MET:CE	2.48	0.44
1:A:236:SER:O	1:A:237:ASP:CB	2.66	0.43
1:A:206:ILE:HA	1:A:209:LEU:CG	2.48	0.43
1:E:209:LEU:HD11	1:E:217:TYR:CE2	2.53	0.43
2:T:1271:ALA:O	5:T:1301:HOH:O	2.20	0.43
1:B:112:ILE:HB	1:B:113:PRO:HD3	2.00	0.43
1:H:186:GLU:OE2	3:H:1301:FMT:O2	2.36	0.43
1:A:55:VAL:HG12	1:A:99:ILE:HD13	2.00	0.43
1:I:41:THR:HG22	1:I:42:VAL:N	2.33	0.43
1:F:112:ILE:HB	1:F:113:PRO:HD3	1.99	0.43
1:A:41:THR:HG22	1:A:42:VAL:N	2.34	0.42
1:J:112:ILE:HB	1:J:113:PRO:HD3	2.01	0.42
1:B:51:ALA:C	5:B:1401:HOH:O	2.62	0.42
1:J:211:THR:O	1:J:211:THR:HG22	2.19	0.42
1:C:11:MET:HE1	1:D:84:ALA:HB1	2.02	0.42
1:G:45:ARG:HD2	1:G:123:PHE:CD1	2.55	0.42
1:C:36:ASP:OD1	1:C:36:ASP:N	2.52	0.42
1:G:236:SER:O	1:G:237:ASP:CB	2.67	0.42
1:D:57:GLY:N	5:D:1401:HOH:O	2.23	0.41
1:A:202:PHE:CE1	1:A:224:MET:HE2	2.55	0.41
1:E:209:LEU:HD13	1:E:209:LEU:O	2.20	0.41
1:A:7:GLU:OE1	1:A:7:GLU:N	2.48	0.41
1:A:49:SER:N	5:A:1401:HOH:O	2.20	0.41
1:A:157:SER:O	1:A:161:ILE:HD12	2.21	0.41
1:E:235:THR:O	1:E:235:THR:HG22	2.20	0.41
1:G:211:THR:O	1:G:211:THR:HG22	2.19	0.41
1:H:157:SER:O	1:H:161:ILE:HD12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:THR:HG21	5:C:1403:HOH:O	2.19	0.41
1:A:152:LEU:HD23	1:E:208[B]:GLU:HG2	2.03	0.41
1:B:157:SER:O	1:B:161:ILE:HD12	2.21	0.41
1:B:218:LYS:HG3	1:B:219:ASP:N	2.36	0.41
1:E:209:LEU:HD11	1:E:217:TYR:CZ	2.56	0.41
1:F:215:GLU:OE1	1:F:215:GLU:N	2.53	0.40
1:G:112:ILE:HB	1:G:113:PRO:HD3	2.03	0.40
1:H:45:ARG:HD2	1:H:123:PHE:CD1	2.57	0.40
1:F:222:LEU:C	1:F:222:LEU:HD23	2.46	0.40

All (3) 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)
1:F:149:GLU:OE2	1:I:184:TYR:OH[1_556]	2.15	0.05
1:B:149:GLU:OE2	1:C:184:TYR:OH[1_665]	2.16	0.04
1:E:143:GLU:H	1:I:79:ASP:OD2[1_666]	1.60	0.00

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	234/241 (97%)	230 (98%)	4 (2%)	0	100	100
1	B	232/241 (96%)	226 (97%)	6 (3%)	0	100	100
1	C	230/241 (95%)	226 (98%)	4 (2%)	0	100	100
1	D	237/241 (98%)	233 (98%)	4 (2%)	0	100	100
1	E	234/241 (97%)	229 (98%)	5 (2%)	0	100	100
1	F	238/241 (99%)	234 (98%)	4 (2%)	0	100	100
1	G	235/241 (98%)	229 (97%)	5 (2%)	1 (0%)	30	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	225/241 (93%)	222 (99%)	3 (1%)	0	100	100
1	I	233/241 (97%)	228 (98%)	5 (2%)	0	100	100
1	J	235/241 (98%)	231 (98%)	4 (2%)	0	100	100
2	K	3/6 (50%)	3 (100%)	0	0	100	100
2	L	3/6 (50%)	3 (100%)	0	0	100	100
2	M	3/6 (50%)	3 (100%)	0	0	100	100
2	N	2/6 (33%)	2 (100%)	0	0	100	100
2	O	3/6 (50%)	3 (100%)	0	0	100	100
2	P	3/6 (50%)	3 (100%)	0	0	100	100
2	Q	3/6 (50%)	3 (100%)	0	0	100	100
2	R	3/6 (50%)	3 (100%)	0	0	100	100
2	S	3/6 (50%)	3 (100%)	0	0	100	100
2	T	3/6 (50%)	3 (100%)	0	0	100	100
All	All	2362/2470 (96%)	2317 (98%)	44 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	3	SER

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	198/199 (100%)	198 (100%)	0	100	100
1	B	196/199 (98%)	193 (98%)	3 (2%)	57	72
1	C	194/199 (98%)	192 (99%)	2 (1%)	68	81
1	D	200/199 (100%)	200 (100%)	0	100	100
1	E	198/199 (100%)	197 (100%)	1 (0%)	81	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	201/199 (101%)	201 (100%)	0	100	100
1	G	198/199 (100%)	197 (100%)	1 (0%)	81	89
1	H	192/199 (96%)	192 (100%)	0	100	100
1	I	197/199 (99%)	194 (98%)	3 (2%)	57	72
1	J	199/199 (100%)	198 (100%)	1 (0%)	81	89
2	K	4/4 (100%)	4 (100%)	0	100	100
2	L	4/4 (100%)	4 (100%)	0	100	100
2	M	4/4 (100%)	4 (100%)	0	100	100
2	N	3/4 (75%)	3 (100%)	0	100	100
2	O	4/4 (100%)	4 (100%)	0	100	100
2	P	4/4 (100%)	4 (100%)	0	100	100
2	Q	4/4 (100%)	4 (100%)	0	100	100
2	R	4/4 (100%)	4 (100%)	0	100	100
2	S	4/4 (100%)	4 (100%)	0	100	100
2	T	4/4 (100%)	4 (100%)	0	100	100
All	All	2012/2030 (99%)	2001 (100%)	11 (0%)	81	89

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	49	SER
1	B	119	ASP
1	B	193	ARG
1	C	142	GLN
1	C	197	LEU
1	E	199	LYS
1	G	78	ASN
1	I	45[A]	ARG
1	I	45[B]	ARG
1	I	79	ASP
1	J	45	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	HIS

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Mol	Chain	Res	Type
1	B	81	HIS
1	C	78	ASN
1	D	81	HIS
1	E	78	ASN
1	E	81	HIS
1	F	196	ASN
1	H	163	ASN
1	J	196	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 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$
2	SEP	L	1270	2	8,9,10	1.63	1 (12%)	7,12,14	1.28	1 (14%)
2	SEP	K	1270	2	8,9,10	1.62	1 (12%)	7,12,14	1.53	1 (14%)
2	SEP	S	1270	2	8,9,10	1.64	1 (12%)	7,12,14	1.53	1 (14%)
2	SEP	R	1270	2	8,9,10	1.61	1 (12%)	7,12,14	1.42	1 (14%)
2	SEP	T	1270	2	8,9,10	1.60	1 (12%)	7,12,14	1.68	1 (14%)
2	SEP	Q	1270	2	8,9,10	1.64	1 (12%)	7,12,14	1.60	1 (14%)
2	SEP	O	1270	2	8,9,10	1.62	1 (12%)	7,12,14	1.30	1 (14%)
2	SEP	N	1270	2	8,9,10	1.61	1 (12%)	7,12,14	1.70	1 (14%)
2	SEP	P	1270	2	8,9,10	1.64	1 (12%)	7,12,14	1.54	1 (14%)
2	SEP	M	1270	2	8,9,10	1.61	1 (12%)	7,12,14	1.55	1 (14%)

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	SEP	L	1270	2	-	0/6/8/10	-
2	SEP	K	1270	2	-	0/6/8/10	-
2	SEP	S	1270	2	-	0/6/8/10	-
2	SEP	R	1270	2	-	0/6/8/10	-
2	SEP	T	1270	2	-	0/6/8/10	-
2	SEP	Q	1270	2	-	0/6/8/10	-
2	SEP	O	1270	2	-	0/6/8/10	-
2	SEP	N	1270	2	-	0/6/8/10	-
2	SEP	P	1270	2	-	0/6/8/10	-
2	SEP	M	1270	2	-	0/6/8/10	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	S	1270	SEP	P-O1P	3.57	1.61	1.50
2	P	1270	SEP	P-O1P	3.52	1.61	1.50
2	Q	1270	SEP	P-O1P	3.52	1.61	1.50
2	O	1270	SEP	P-O1P	3.52	1.61	1.50
2	L	1270	SEP	P-O1P	3.52	1.61	1.50
2	N	1270	SEP	P-O1P	3.51	1.61	1.50
2	K	1270	SEP	P-O1P	3.50	1.61	1.50
2	R	1270	SEP	P-O1P	3.49	1.61	1.50
2	M	1270	SEP	P-O1P	3.49	1.61	1.50
2	T	1270	SEP	P-O1P	3.46	1.61	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	1270	SEP	OG-CB-CA	3.99	112.03	108.14
2	T	1270	SEP	OG-CB-CA	3.89	111.93	108.14
2	Q	1270	SEP	OG-CB-CA	3.67	111.72	108.14
2	M	1270	SEP	OG-CB-CA	3.60	111.65	108.14
2	S	1270	SEP	OG-CB-CA	3.59	111.64	108.14
2	K	1270	SEP	OG-CB-CA	3.50	111.55	108.14
2	P	1270	SEP	OG-CB-CA	3.42	111.47	108.14
2	R	1270	SEP	OG-CB-CA	3.20	111.26	108.14
2	L	1270	SEP	OG-CB-CA	2.91	110.97	108.14
2	O	1270	SEP	OG-CB-CA	2.74	110.82	108.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	1270	SEP	1	0
2	R	1270	SEP	1	0
2	T	1270	SEP	1	0
2	Q	1270	SEP	1	0
2	M	1270	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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	$\# Z > 2$
3	FMT	B	1301	-	2,2,2	0.71	0	1,1,1	0.46	0
3	FMT	C	1301	-	2,2,2	0.73	0	1,1,1	0.45	0
3	FMT	H	1301	-	2,2,2	0.75	0	1,1,1	0.46	0
3	FMT	Q	1301	-	2,2,2	0.70	0	1,1,1	0.47	0
3	FMT	J	1301	-	2,2,2	0.72	0	1,1,1	0.43	0
3	FMT	E	1302	-	2,2,2	0.74	0	1,1,1	0.46	0
3	FMT	J	1302	-	2,2,2	0.73	0	1,1,1	0.45	0
3	FMT	K	1301	-	2,2,2	0.70	0	1,1,1	0.46	0
3	FMT	S	1301	-	2,2,2	0.71	0	1,1,1	0.46	0
4	EDO	E	1301	-	3,3,3	0.42	0	2,2,2	0.56	0
3	FMT	D	1301	-	2,2,2	0.75	0	1,1,1	0.45	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	E	1301	-	-	1/1/1/1	-

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
4	E	1301	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1301	FMT	1	0
3	H	1301	FMT	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	234/241 (97%)	0.49	11 (4%)	36	42	35, 62, 115, 168	1 (0%)
1	B	234/241 (97%)	0.41	18 (7%)	19	22	33, 63, 132, 182	0
1	C	232/241 (96%)	0.44	17 (7%)	21	24	38, 65, 113, 137	0
1	D	236/241 (97%)	0.42	14 (5%)	28	32	37, 67, 132, 183	2 (0%)
1	E	234/241 (97%)	0.30	9 (3%)	44	50	35, 58, 110, 138	2 (0%)
1	F	237/241 (98%)	0.40	12 (5%)	33	39	36, 64, 129, 163	2 (0%)
1	G	236/241 (97%)	0.63	17 (7%)	21	24	40, 70, 124, 160	1 (0%)
1	H	227/241 (94%)	0.49	16 (7%)	22	25	34, 64, 113, 171	1 (0%)
1	I	234/241 (97%)	0.32	7 (2%)	52	59	34, 63, 116, 157	1 (0%)
1	J	234/241 (97%)	0.35	11 (4%)	36	42	35, 61, 120, 160	2 (0%)
2	K	5/6 (83%)	1.15	1 (20%)	3	2	55, 60, 80, 89	0
2	L	5/6 (83%)	1.49	2 (40%)	1	0	49, 59, 85, 118	0
2	M	5/6 (83%)	1.26	1 (20%)	3	2	54, 63, 92, 106	0
2	N	4/6 (66%)	1.57	2 (50%)	0	0	61, 67, 85, 102	0
2	O	5/6 (83%)	1.20	0	100	100	51, 60, 85, 121	0
2	P	5/6 (83%)	1.44	2 (40%)	1	0	58, 63, 106, 112	0
2	Q	5/6 (83%)	1.32	2 (40%)	1	0	62, 67, 77, 99	0
2	R	5/6 (83%)	1.31	0	100	100	65, 79, 88, 98	0
2	S	5/6 (83%)	0.85	0	100	100	60, 68, 87, 108	0
2	T	5/6 (83%)	1.53	1 (20%)	3	2	68, 69, 97, 122	0
All	All	2387/2470 (96%)	0.44	143 (5%)	27	31	33, 64, 118, 183	12 (0%)

All (143) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	142[A]	GLN	6.8
1	G	2	ALA	6.6
1	F	234	TRP	5.0
1	B	37	GLY	4.3
1	H	217	TYR	4.2
1	H	209	LEU	3.9
1	D	234	TRP	3.8
1	H	236	SER	3.8
1	H	235	THR	3.6
1	B	217	TYR	3.6
1	I	237	ASP	3.5
1	B	216	SER	3.4
1	C	209	LEU	3.4
1	C	211	THR	3.3
1	I	209	LEU	3.3
1	A	77	GLY	3.2
1	A	78	ASN	3.2
1	E	216	SER	3.2
1	B	235	THR	3.2
1	H	37	GLY	3.2
1	J	209	LEU	3.1
2	N	1268	LEU	3.1
1	D	37	GLY	3.1
1	H	206	ILE	3.0
2	P	1268	LEU	3.0
1	F	230	ASN	3.0
1	H	233	LEU	3.0
1	J	231	LEU	3.0
1	J	206	ILE	3.0
1	F	64[A]	ARG	3.0
1	C	202	PHE	2.9
2	L	1268	LEU	2.9
1	E	235	THR	2.9
1	A	33	ALA	2.9
1	A	34	ALA	2.9
1	G	212	LEU	2.9
1	E	35	VAL	2.9
1	B	236	SER	2.9
1	F	237	ASP	2.9
1	B	33	ALA	2.8
1	G	207	ALA	2.8
1	F	1	MET	2.8
1	D	220	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	34	ALA	2.8
1	D	33	ALA	2.7
1	B	213	GLY	2.7
1	D	1	MET	2.7
1	G	209	LEU	2.7
1	H	237	ASP	2.7
1	G	213	GLY	2.7
1	D	212	LEU	2.7
1	A	35	VAL	2.6
1	J	36	ASP	2.6
1	C	184	TYR	2.6
1	G	76	ARG	2.6
1	E	37	GLY	2.6
1	E	213	GLY	2.6
2	N	1269	PHE	2.6
1	F	232	THR	2.6
1	H	35	VAL	2.6
1	G	217	TYR	2.6
1	H	219	ASP	2.5
1	B	234	TRP	2.5
1	B	35	VAL	2.5
1	H	76	ARG	2.5
1	F	227	LEU	2.5
1	B	207	ALA	2.5
1	G	79	ASP	2.5
1	J	216	SER	2.5
1	C	197	LEU	2.5
1	B	211	THR	2.5
1	G	206	ILE	2.5
1	J	35	VAL	2.5
1	C	233	LEU	2.5
1	E	209	LEU	2.5
1	A	167	ALA	2.5
1	G	33	ALA	2.5
1	G	221	THR	2.4
1	D	2	ALA	2.4
1	J	234	TRP	2.4
1	A	141	GLY	2.4
1	J	182	VAL	2.4
1	C	234	TRP	2.4
1	I	90	SER	2.4
1	H	168	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	207	ALA	2.3
1	G	75	SER	2.3
1	I	236	SER	2.3
1	B	206	ILE	2.3
1	G	198	ALA	2.3
1	I	206	ILE	2.3
1	B	237	ASP	2.3
1	J	219	ASP	2.3
1	G	37	GLY	2.3
1	G	164	ALA	2.3
1	C	232	THR	2.3
1	C	231	LEU	2.3
1	F	37	GLY	2.3
1	C	216	SER	2.3
2	Q	1267	GLU	2.3
2	T	1267	GLU	2.3
2	K	1268	LEU	2.2
1	B	5	ARG	2.2
1	H	141	GLY	2.2
1	C	201	ALA	2.2
1	D	214	GLU	2.2
1	D	32	SER	2.2
1	I	4	GLY	2.2
1	D	29	GLU	2.2
2	L	1267	GLU	2.2
1	F	212	LEU	2.2
1	C	141	GLY	2.2
1	B	214	GLU	2.2
1	C	215	GLU	2.2
1	H	207	ALA	2.2
1	J	207	ALA	2.2
1	B	32	SER	2.2
1	A	30	LYS	2.2
1	D	236	SER	2.2
1	E	207	ALA	2.2
1	F	217	TYR	2.1
1	E	231	LEU	2.1
1	H	222	LEU	2.1
2	Q	1268	LEU	2.1
1	D	216	SER	2.1
1	F	235	THR	2.1
1	G	121	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
2	P	1267	GLU	2.1
1	C	31	VAL	2.1
1	I	234	TRP	2.1
1	A	75	SER	2.1
1	C	221	THR	2.1
1	J	235	THR	2.1
1	E	212	LEU	2.1
1	D	210	ASP	2.1
1	H	119	ASP	2.1
2	M	1268	LEU	2.1
1	G	31	VAL	2.1
1	B	221	THR	2.0
1	A	208	GLU	2.0
1	B	169	THR	2.0
1	C	214	GLU	2.0
1	D	233	LEU	2.0

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
2	SEP	M	1270	10/11	0.95	0.08	42,48,58,59	0
2	SEP	S	1270	10/11	0.95	0.07	44,49,57,67	0
2	SEP	K	1270	10/11	0.96	0.06	37,45,53,57	0
2	SEP	N	1270	10/11	0.96	0.08	47,48,65,72	0
2	SEP	O	1270	10/11	0.96	0.07	39,43,54,54	0
2	SEP	Q	1270	10/11	0.96	0.07	40,46,56,58	0
2	SEP	R	1270	10/11	0.96	0.09	54,57,67,68	0
2	SEP	L	1270	10/11	0.96	0.07	39,43,47,49	0
2	SEP	T	1270	10/11	0.96	0.08	45,53,61,63	0
2	SEP	P	1270	10/11	0.97	0.08	38,47,59,66	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FMT	J	1301	3/3	0.57	0.20	92,94,111,113	0
3	FMT	S	1301	3/3	0.79	0.21	73,76,76,87	0
3	FMT	Q	1301	3/3	0.80	0.24	65,69,71,86	0
3	FMT	J	1302	3/3	0.83	0.35	75,91,110,123	0
4	EDO	E	1301	4/4	0.84	0.19	51,83,103,103	0
3	FMT	D	1301	3/3	0.86	0.14	58,60,60,72	0
3	FMT	B	1301	3/3	0.86	0.15	54,57,59,71	0
3	FMT	C	1301	3/3	0.88	0.14	58,60,79,95	0
3	FMT	K	1301	3/3	0.89	0.17	50,54,62,65	0
3	FMT	E	1302	3/3	0.91	0.12	43,53,57,64	0
3	FMT	H	1301	3/3	0.91	0.11	50,53,53,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.