



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:51 AM GMT

PDB ID : 2YJK  
Title : STRUCTURE OF DPS FROM MICROBACTERIUM ARBORESCENS IN  
THE HIGH IRON FORM  
Authors : Zeth, K.; Buecheler, R.; Boland, W.  
Deposited on : 2011-05-19  
Resolution : 2.00 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

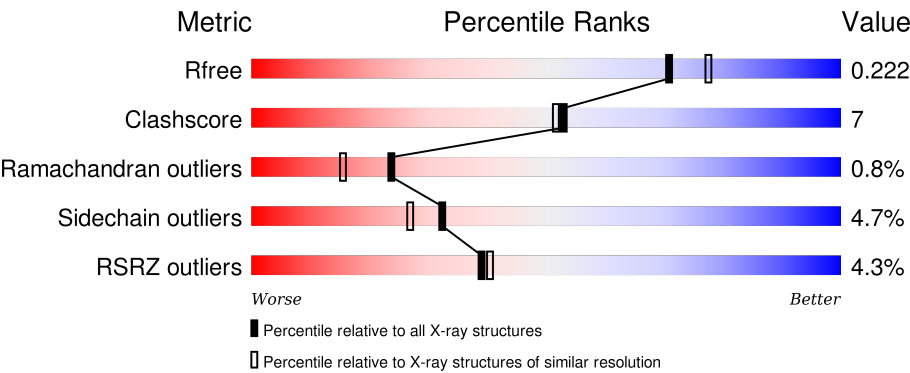
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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 <sub>free</sub>	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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  $\geq 3$ , 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	161	<div><div>76%14%7%</div></div>
1	B	161	<div><div>78%14%6%</div></div>
1	C	161	<div><div>78%16%...</div></div>
1	D	161	<div><div>81%10%6%</div></div>
1	E	161	<div><div>80%12%6%</div></div>

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Mol	Chain	Length	Quality of chain
1	F	161	<div><div></div><div>4%</div><div>83%</div><div>10%</div><div>•</div><div>6%</div></div>
1	G	161	<div><div></div><div>3%</div><div>80%</div><div>11%</div><div>•</div><div>•</div><div>5%</div></div>
1	H	161	<div><div></div><div>4%</div><div>81%</div><div>9%</div><div>•</div><div>•</div><div>6%</div></div>
1	I	161	<div><div></div><div>4%</div><div>81%</div><div>9%</div><div>•</div><div>•</div><div>8%</div></div>
1	J	161	<div><div></div><div>2%</div><div>80%</div><div>10%</div><div>•</div><div>8%</div></div>
1	K	161	<div><div></div><div>6%</div><div>84%</div><div>9%</div><div>•</div><div>6%</div></div>
1	L	161	<div><div></div><div>2%</div><div>75%</div><div>16%</div><div>•</div><div>6%</div></div>

## 2 Entry composition

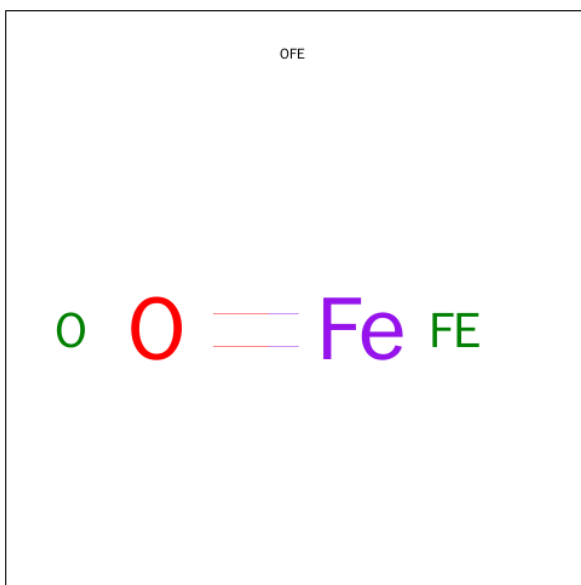
There are 4 unique types of molecules in this entry. The entry contains 14492 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AFP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	1	0
			1134	717	195	220	2			
1	B	152	Total	C	N	O	S	0	0	0
			1148	725	197	224	2			
1	C	157	Total	C	N	O	S	0	0	0
			1173	740	202	229	2			
1	D	152	Total	C	N	O	S	0	3	0
			1158	732	197	227	2			
1	E	152	Total	C	N	O	S	0	0	0
			1146	724	197	223	2			
1	F	152	Total	C	N	O	S	0	0	0
			1146	724	197	223	2			
1	G	153	Total	C	N	O	S	0	0	0
			1153	728	198	225	2			
1	H	152	Total	C	N	O	S	0	1	0
			1151	727	197	225	2			
1	I	148	Total	C	N	O	S	0	0	0
			1123	709	193	219	2			
1	J	148	Total	C	N	O	S	0	0	0
			1123	709	193	219	2			
1	K	152	Total	C	N	O	S	0	1	0
			1151	727	197	225	2			
1	L	152	Total	C	N	O	S	0	1	0
			1153	728	197	226	2			

- Molecule 2 is IRON(II) OXIDE (three-letter code: OFE) (formula: FeO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	O	0	0
			2	1	1		
2	B	1	Total	Fe	O	0	0
			2	1	1		
2	C	1	Total	Fe	O	0	0
			2	1	1		
2	D	1	Total	Fe	O	0	0
			2	1	1		
2	E	1	Total	Fe	O	0	0
			2	1	1		
2	F	1	Total	Fe	O	0	0
			2	1	1		
2	G	1	Total	Fe	O	0	0
			2	1	1		
2	H	1	Total	Fe	O	0	0
			2	1	1		
2	J	1	Total	Fe	O	0	0
			2	1	1		
2	K	1	Total	Fe	O	0	0
			2	1	1		
2	L	1	Total	Fe	O	0	0
			2	1	1		

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	J	1	Total	Fe	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Fe 1	0	0
3	B	1	Total 1	Fe 1	0	0
3	I	1	Total 1	Fe 1	0	0
3	C	1	Total 1	Fe 1	0	0
3	L	1	Total 1	Fe 1	0	0

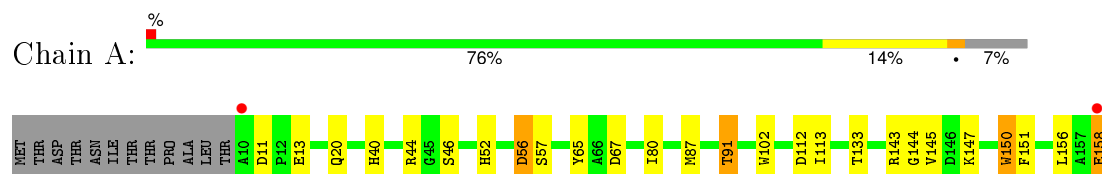
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total 59	O 59	0	0
4	B	58	Total 58	O 58	0	0
4	C	43	Total 43	O 43	0	0
4	D	65	Total 65	O 65	0	0
4	E	54	Total 54	O 54	0	0
4	F	62	Total 62	O 62	0	0
4	G	93	Total 93	O 93	0	0
4	H	70	Total 70	O 70	0	0
4	I	39	Total 39	O 39	0	0
4	J	56	Total 56	O 56	0	0
4	K	50	Total 50	O 50	0	0
4	L	56	Total 56	O 56	0	0

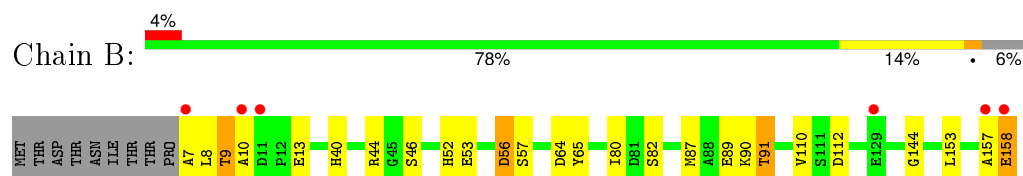
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

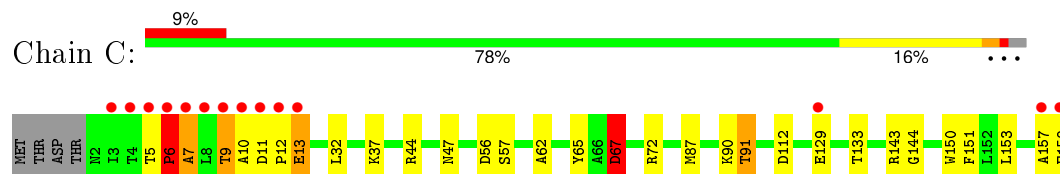
- Molecule 1: AFP



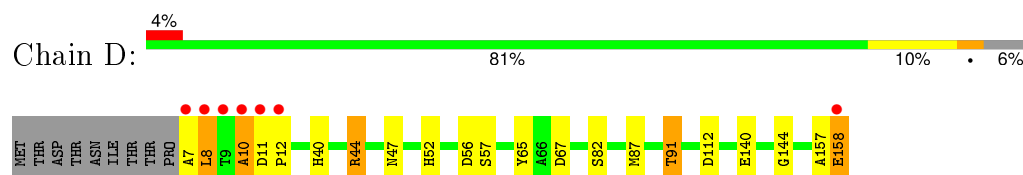
- Molecule 1: AFP



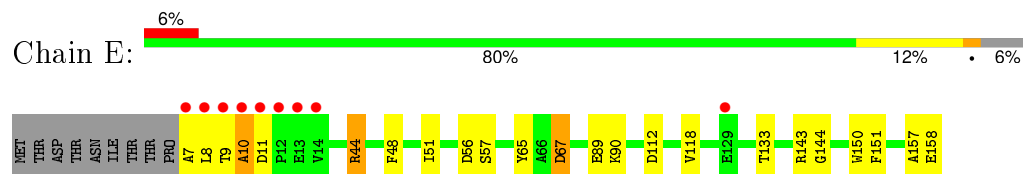
- Molecule 1: AFP



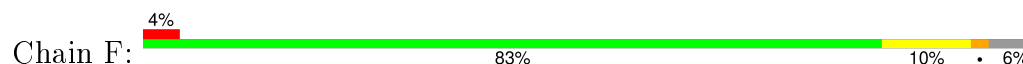
- Molecule 1: AFP

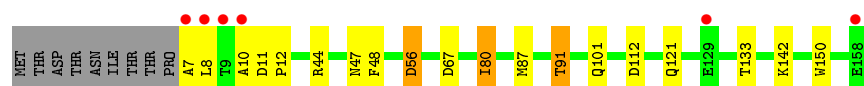


- Molecule 1: AFP

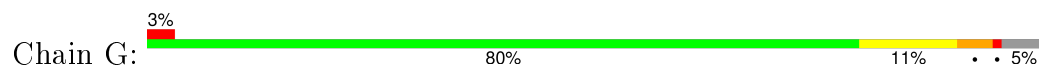


- Molecule 1: AFP

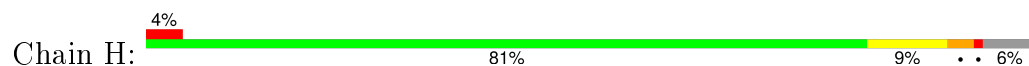




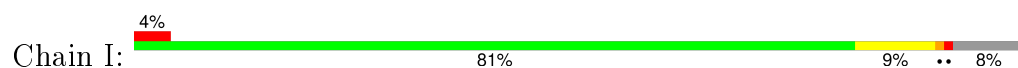
- Molecule 1: AFP



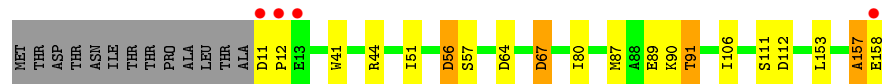
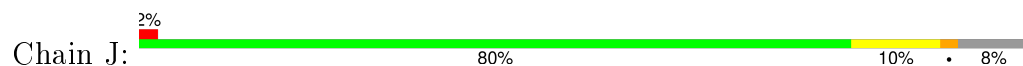
- Molecule 1: AFP



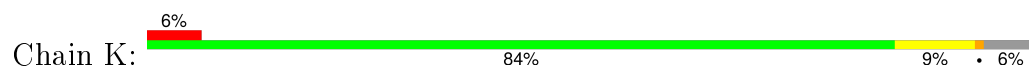
- Molecule 1: AFP



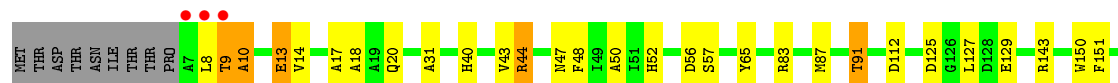
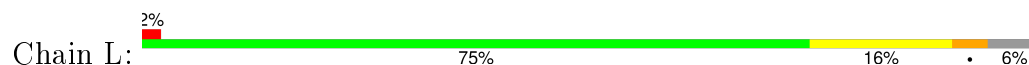
- Molecule 1: AFP



- Molecule 1: AFP



- Molecule 1: AFP





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.58Å 92.05Å 128.58Å 90.00° 96.04° 90.00°	Depositor
Resolution (Å)	28.98 – 2.00 28.98 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (28.98-2.00) 99.3 (28.98-2.00)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.170 , 0.216 0.178 , 0.222	Depositor DCC
$R_{free}$ test set	6820 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.3	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 42.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 136426 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	14492	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: FE, OFE

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	1.13	3/1156 (0.3%)	0.90	2/1576 (0.1%)
1	B	1.25	4/1167 (0.3%)	0.90	2/1593 (0.1%)
1	C	1.11	1/1192 (0.1%)	0.93	4/1628 (0.2%)
1	D	1.14	2/1186 (0.2%)	0.92	2/1620 (0.1%)
1	E	1.11	3/1165 (0.3%)	0.91	3/1590 (0.2%)
1	F	1.17	2/1165 (0.2%)	0.92	2/1590 (0.1%)
1	G	1.22	2/1172 (0.2%)	1.00	6/1600 (0.4%)
1	H	1.30	4/1173 (0.3%)	1.03	5/1601 (0.3%)
1	I	1.08	1/1142 (0.1%)	0.94	2/1558 (0.1%)
1	J	1.19	3/1142 (0.3%)	0.91	2/1558 (0.1%)
1	K	1.13	3/1173 (0.3%)	0.86	1/1601 (0.1%)
1	L	1.14	1/1175 (0.1%)	0.93	6/1604 (0.4%)
All	All	1.17	29/14008 (0.2%)	0.93	37/19119 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
1	D	0	1
1	F	0	1
1	G	0	1
1	H	0	1
1	J	0	1
1	L	0	1
All	All	0	8

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	112	ASP	CB-CG	9.04	1.70	1.51
1	J	112	ASP	CB-CG	7.78	1.68	1.51
1	I	112	ASP	CB-CG	7.78	1.68	1.51
1	F	112	ASP	CB-CG	7.51	1.67	1.51
1	B	112	ASP	CB-CG	7.38	1.67	1.51
1	E	89	GLU	CG-CD	7.30	1.62	1.51
1	D	112	ASP	CB-CG	7.27	1.67	1.51
1	G	112	ASP	CB-CG	6.91	1.66	1.51
1	E	112	ASP	CB-CG	6.91	1.66	1.51
1	J	67	ASP	CB-CG	-6.83	1.37	1.51
1	K	140	GLU	CG-CD	6.74	1.62	1.51
1	L	112	ASP	CB-CG	6.44	1.65	1.51
1	C	112	ASP	CB-CG	6.40	1.65	1.51
1	A	112	ASP	CB-CG	6.17	1.64	1.51
1	J	89	GLU	CG-CD	5.90	1.60	1.51
1	B	13	GLU	CG-CD	5.87	1.60	1.51
1	K	112	ASP	CB-CG	5.86	1.64	1.51
1	B	89	GLU	CG-CD	5.85	1.60	1.51
1	A	150	TRP	CB-CG	5.64	1.60	1.50
1	K	63	GLN	CG-CD	5.64	1.64	1.51
1	B	53	GLU	CB-CG	5.57	1.62	1.52
1	H	89	GLU	CG-CD	5.51	1.60	1.51
1	E	48	PHE	CE2-CZ	5.49	1.47	1.37
1	H	43	VAL	CB-CG1	5.46	1.64	1.52
1	H	100	ALA	CA-CB	5.36	1.63	1.52
1	D	140	GLU	CG-CD	5.17	1.59	1.51
1	F	150	TRP	CB-CG	5.03	1.59	1.50
1	G	140	GLU	CG-CD	5.03	1.59	1.51
1	A	102	TRP	CE3-CZ3	5.01	1.47	1.38

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	44	ARG	NE-CZ-NH2	-10.69	114.96	120.30
1	C	67	ASP	CB-CG-OD1	-10.06	109.25	118.30
1	L	112	ASP	CB-CG-OD1	9.68	127.01	118.30
1	I	67	ASP	CB-CG-OD2	-8.79	110.39	118.30
1	B	112	ASP	CB-CG-OD1	7.88	125.39	118.30
1	H	112	ASP	CB-CG-OD1	7.85	125.37	118.30
1	J	67	ASP	CB-CG-OD1	-7.79	111.29	118.30
1	G	112	ASP	CB-CG-OD1	7.59	125.13	118.30
1	D	44	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	A	112	ASP	CB-CG-OD1	7.32	124.89	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	112	ASP	CB-CG-OD1	6.92	124.53	118.30
1	C	72	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	C	6	PRO	N-CA-CB	6.60	111.22	103.30
1	A	11	ASP	CB-CG-OD1	6.33	123.99	118.30
1	G	6	PRO	N-CA-CB	5.96	110.45	103.30
1	E	112	ASP	CB-CG-OD1	5.89	123.60	118.30
1	C	112	ASP	CB-CG-OD1	5.87	123.58	118.30
1	E	67	ASP	CB-CG-OD2	-5.76	113.11	118.30
1	H	44	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	G	81	ASP	CB-CG-OD1	5.43	123.19	118.30
1	B	112	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	I	72	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	L	44	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	J	112	ASP	CB-CG-OD1	5.35	123.11	118.30
1	G	44	ARG	NE-CZ-NH2	-5.25	117.68	120.30
1	F	80	ILE	CG1-CB-CG2	-5.20	99.97	111.40
1	L	112	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	G	57	SER	CB-CA-C	5.14	119.86	110.10
1	H	91	THR	OG1-CB-CG2	5.13	121.81	110.00
1	G	72	ARG	NE-CZ-NH1	-5.12	117.74	120.30
1	H	8	LEU	CA-CB-CG	5.12	127.08	115.30
1	L	125	ASP	CB-CG-OD1	5.08	122.88	118.30
1	L	83	ARG	NE-CZ-NH2	-5.07	117.76	120.30
1	K	112	ASP	CB-CG-OD1	5.04	122.84	118.30
1	E	44	ARG	NE-CZ-NH2	-5.02	117.79	120.30
1	L	143	ARG	NE-CZ-NH1	-5.01	117.79	120.30
1	D	112	ASP	CB-CG-OD1	5.01	122.81	118.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	157	ALA	Peptide
1	C	157	ALA	Peptide
1	D	10	ALA	Peptide
1	F	10	ALA	Peptide
1	G	157	ALA	Peptide
1	H	157	ALA	Peptide
1	J	157	ALA	Peptide
1	L	157	ALA	Peptide

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1134	0	1128	20	0
1	B	1148	0	1138	20	0
1	C	1173	0	1147	23	0
1	D	1158	0	1149	15	0
1	E	1146	0	1133	15	0
1	F	1146	0	1133	11	0
1	G	1153	0	1140	17	0
1	H	1151	0	1137	15	0
1	I	1123	0	1110	13	0
1	J	1123	0	1110	19	0
1	K	1151	0	1137	13	0
1	L	1153	0	1142	23	0
2	A	2	0	0	1	0
2	B	2	0	0	1	0
2	C	2	0	0	0	0
2	D	2	0	0	1	0
2	E	2	0	0	1	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	1	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	L	1	0	0	0	0
4	A	59	0	0	5	0
4	B	58	0	0	6	0
4	C	43	0	0	1	0
4	D	65	0	0	2	0
4	E	54	0	0	7	0
4	F	62	0	0	1	0
4	G	93	0	0	9	0
4	H	70	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	39	0	0	1	0
4	J	56	0	0	8	0
4	K	50	0	0	3	0
4	L	56	0	0	13	0
All	All	14492	0	13604	182	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 7.

All (182) 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)
4:E:2013:HOH:O	1:F:67:ASP:HB2	1.25	1.30
1:B:64:ASP:HB3	4:B:2018:HOH:O	1.09	1.25
1:G:67:ASP:HB2	4:G:2035:HOH:O	1.45	1.16
1:A:20:GLN:NE2	4:A:2002:HOH:O	1.76	1.12
1:D:67[B]:ASP:OD2	4:D:2024:HOH:O	1.67	1.11
1:H:13:GLU:HG2	4:H:2004:HOH:O	1.51	1.08
1:J:90:LYS:HE2	4:J:2033:HOH:O	1.59	1.02
1:A:67:ASP:HB2	4:A:2024:HOH:O	1.59	1.01
1:L:47:ASN:OD1	4:L:2012:HOH:O	1.85	0.95
1:E:9:THR:HA	1:E:10:ALA:HB3	1.46	0.95
1:J:91:THR:HG23	4:J:2008:HOH:O	1.68	0.94
1:H:9:THR:CB	1:H:10:ALA:HA	1.99	0.93
1:J:157:ALA:O	1:J:158:GLU:HG3	1.67	0.92
1:G:91:THR:HG23	4:G:2013:HOH:O	1.72	0.90
1:J:90:LYS:CE	4:J:2033:HOH:O	2.19	0.90
1:K:11:ASP:O	4:K:2001:HOH:O	1.95	0.85
1:B:87:MET:O	1:B:91:THR:HB	1.77	0.84
1:D:91:THR:HG23	4:D:2009:HOH:O	1.78	0.84
1:J:90:LYS:HE3	4:J:2032:HOH:O	1.78	0.83
1:B:91:THR:HG23	4:B:2011:HOH:O	1.78	0.83
1:B:8:LEU:HB3	1:B:9:THR:HA	1.61	0.82
1:L:9:THR:HG23	1:L:10:ALA:H	1.43	0.82
1:A:91:THR:HG23	4:A:2011:HOH:O	1.79	0.81
1:I:91:THR:HG23	4:I:2008:HOH:O	1.80	0.81
1:K:91:THR:HG23	4:K:2008:HOH:O	1.81	0.80
1:B:90:LYS:HE3	4:B:2026:HOH:O	1.83	0.79
1:B:7:ALA:N	1:B:8:LEU:HA	1.97	0.79
1:L:50:ALA:HB3	4:L:2012:HOH:O	1.82	0.78
1:H:91:THR:HG23	4:H:2015:HOH:O	1.82	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:90:LYS:CE	4:J:2032:HOH:O	2.32	0.76
1:E:9:THR:HA	1:E:10:ALA:CB	2.15	0.76
1:L:129:GLU:HB2	4:L:2044:HOH:O	1.87	0.75
1:D:11:ASP:OD1	1:D:12:PRO:HD2	1.88	0.73
1:F:91:THR:HG23	4:F:2012:HOH:O	1.89	0.73
1:L:48:PHE:HA	4:L:2010:HOH:O	1.87	0.72
1:A:87:MET:O	1:A:91:THR:HB	1.89	0.72
1:A:44:ARG:HH22	1:B:7:ALA:HB3	1.54	0.71
1:D:158:GLU:HA	1:D:158:GLU:OE1	1.90	0.70
1:L:43:VAL:HG22	4:L:2010:HOH:O	1.90	0.70
1:C:10:ALA:H	1:C:11:ASP:HA	1.57	0.68
1:J:157:ALA:O	1:J:158:GLU:CG	2.39	0.68
1:J:64:ASP:HB3	4:J:2022:HOH:O	1.93	0.67
1:E:150:TRP:CZ3	1:E:151:PHE:HE1	2.13	0.66
1:L:87:MET:O	1:L:91:THR:HB	1.96	0.65
1:H:9:THR:CB	1:H:10:ALA:CA	2.73	0.64
1:J:90:LYS:NZ	4:J:2033:HOH:O	2.30	0.64
1:H:157:ALA:O	1:H:158:GLU:HB2	1.98	0.64
1:F:7:ALA:HB3	4:H:2069:HOH:O	1.98	0.64
1:E:11:ASP:HB2	4:E:2001:HOH:O	1.97	0.64
1:H:10:ALA:O	1:H:11:ASP:C	2.35	0.63
1:L:65:TYR:HE2	4:L:2017:HOH:O	1.80	0.63
1:D:8:LEU:HD22	4:L:2039:HOH:O	1.99	0.63
1:C:87:MET:O	1:C:91:THR:HB	1.99	0.63
1:J:87:MET:O	1:J:91:THR:HB	1.99	0.62
1:H:9:THR:H	1:H:10:ALA:HB2	1.63	0.62
1:E:90:LYS:HE3	4:E:2036:HOH:O	1.99	0.62
1:G:67:ASP:CB	4:G:2035:HOH:O	2.22	0.62
1:E:157:ALA:O	1:E:158:GLU:C	2.37	0.62
4:A:2017:HOH:O	1:B:7:ALA:HB3	1.99	0.62
1:I:157:ALA:O	1:I:158:GLU:HB2	1.98	0.61
1:D:82[A]:SER:HG	1:J:41:TRP:HZ3	1.48	0.61
1:C:91:THR:HG23	4:C:2010:HOH:O	1.99	0.60
1:I:87:MET:O	1:I:91:THR:HB	2.00	0.60
1:K:11:ASP:OD1	1:K:12:PRO:HD2	2.02	0.60
1:L:91:THR:CG2	4:L:2008:HOH:O	2.48	0.60
1:C:129:GLU:OE1	1:C:129:GLU:HA	2.02	0.60
1:L:65:TYR:CE2	4:L:2017:HOH:O	2.51	0.60
1:D:157:ALA:O	1:D:158:GLU:HB2	2.02	0.60
1:C:10:ALA:N	1:C:11:ASP:HA	2.17	0.60
1:K:11:ASP:HB3	4:K:2002:HOH:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:ASP:OD1	1:C:12:PRO:HD2	2.04	0.58
1:E:9:THR:CA	1:E:10:ALA:HB3	2.28	0.57
1:A:147[B]:LYS:HG2	4:E:2028:HOH:O	2.04	0.57
2:E:1159:OFE:O	1:F:56:ASP:OD2	2.21	0.57
1:D:87:MET:O	1:D:91:THR:HB	2.04	0.57
1:H:87:MET:O	1:H:91:THR:HB	2.05	0.57
1:K:87:MET:O	1:K:91:THR:HB	2.05	0.57
1:H:9:THR:N	1:H:10:ALA:HB2	2.20	0.56
4:G:2022:HOH:O	1:J:158:GLU:O	2.18	0.56
1:C:7:ALA:H	1:E:118:VAL:HG22	1.68	0.56
1:B:40:HIS:CE1	1:B:52:HIS:CE1	2.94	0.56
1:L:9:THR:HG23	1:L:10:ALA:N	2.16	0.55
1:K:56:ASP:OD2	2:L:1159:OFE:O	2.24	0.55
1:I:40:HIS:CE1	1:I:52:HIS:CE1	2.95	0.55
1:L:150:TRP:CZ3	1:L:151:PHE:HE1	2.25	0.55
1:F:87:MET:O	1:F:91:THR:HB	2.06	0.54
4:E:2013:HOH:O	1:F:67:ASP:CB	2.08	0.54
1:L:91:THR:HG23	4:L:2008:HOH:O	2.05	0.54
1:B:65:TYR:OH	1:B:144:GLY:HA3	2.08	0.54
1:C:150:TRP:CZ3	1:C:151:PHE:HE1	2.25	0.54
1:J:90:LYS:HE2	4:J:2032:HOH:O	2.05	0.53
1:G:90:LYS:HE2	4:G:2052:HOH:O	2.08	0.53
1:L:129:GLU:CB	4:L:2044:HOH:O	2.51	0.53
1:J:51:ILE:HD12	1:J:106:ILE:HD11	1.90	0.53
1:C:6:PRO:O	1:C:7:ALA:HB3	2.09	0.53
1:H:65:TYR:OH	1:H:144:GLY:HA3	2.09	0.52
1:G:6:PRO:CB	4:G:2001:HOH:O	2.58	0.52
4:B:2013:HOH:O	1:C:158:GLU:C	2.48	0.52
1:D:65:TYR:OH	1:D:144:GLY:HA3	2.09	0.52
2:A:1159:OFE:O	1:B:56:ASP:OD2	2.28	0.52
1:B:8:LEU:HB3	1:B:9:THR:CA	2.38	0.51
1:E:150:TRP:CZ3	1:E:151:PHE:CE1	2.96	0.51
1:K:40:HIS:CE1	1:K:52:HIS:CE1	2.98	0.51
1:A:44:ARG:NH2	1:B:7:ALA:HB3	2.24	0.51
1:K:157:ALA:O	1:K:158:GLU:HG2	2.10	0.51
1:G:44:ARG:HD2	1:G:99:PHE:HB3	1.93	0.51
4:G:2014:HOH:O	1:H:67:ASP:HB2	2.11	0.50
1:A:65:TYR:OH	1:A:144:GLY:HA3	2.11	0.50
1:A:143:ARG:HG3	1:E:133:THR:HG23	1.94	0.50
1:L:40:HIS:CE1	1:L:52:HIS:CE1	3.00	0.49
1:A:143:ARG:O	1:A:147[B]:LYS:HE3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:GLU:OE1	1:C:13:GLU:HA	2.13	0.49
1:H:7:ALA:N	1:H:8:LEU:HA	2.28	0.49
1:G:87:MET:O	1:G:91:THR:HB	2.13	0.49
1:G:157:ALA:O	1:G:158:GLU:HB2	2.13	0.49
1:D:82[A]:SER:OG	1:J:41:TRP:HZ3	1.95	0.49
1:C:7:ALA:HB2	4:E:2047:HOH:O	2.12	0.48
1:C:37:LYS:NZ	1:I:67:ASP:OD2	2.40	0.48
1:B:91:THR:CG2	4:B:2011:HOH:O	2.49	0.48
1:A:56:ASP:OD2	2:B:1159:OFE:O	2.32	0.48
1:L:47:ASN:CG	4:L:2012:HOH:O	2.43	0.48
1:F:11:ASP:HA	1:F:12:PRO:HD2	1.79	0.48
1:C:32:LEU:HD23	1:C:62:ALA:HB2	1.95	0.47
2:D:1159:OFE:O	1:J:56:ASP:OD2	2.32	0.47
1:L:150:TRP:CZ3	1:L:151:PHE:CE1	3.02	0.47
1:G:37:LYS:HE3	4:G:2014:HOH:O	2.13	0.47
1:I:129:GLU:HA	1:I:129:GLU:OE1	2.15	0.47
1:E:7:ALA:HA	1:E:8:LEU:C	2.35	0.47
1:G:158:GLU:OE1	1:G:158:GLU:HA	2.15	0.46
1:C:7:ALA:CB	4:E:2047:HOH:O	2.63	0.46
1:G:8:LEU:HD13	1:I:153:LEU:HG	1.98	0.46
1:B:8:LEU:CB	1:B:9:THR:HA	2.39	0.45
1:A:40:HIS:CE1	1:A:52:HIS:CE1	3.03	0.45
1:I:11:ASP:HA	1:I:12:PRO:HD2	1.70	0.45
1:L:13:GLU:HG3	1:L:14:VAL:N	2.31	0.45
1:C:133:THR:HG23	1:E:143:ARG:HG3	1.97	0.45
1:B:46:SER:O	1:C:47:ASN:ND2	2.48	0.45
1:H:157:ALA:O	1:H:158:GLU:CB	2.64	0.45
1:A:150:TRP:CZ3	1:A:151:PHE:HE1	2.35	0.45
1:F:47:ASN:O	1:F:48:PHE:C	2.55	0.45
1:B:158:GLU:HA	1:B:158:GLU:OE1	2.17	0.44
1:G:44:ARG:HB3	1:J:157:ALA:HB3	1.99	0.44
1:G:7:ALA:O	1:G:8:LEU:CB	2.66	0.44
1:D:7:ALA:O	1:D:8:LEU:HB2	2.17	0.44
1:A:156:LEU:O	1:A:158:GLU:HG3	2.18	0.43
1:G:28:LYS:NZ	1:G:119:ASP:OD2	2.36	0.43
1:A:67:ASP:CB	4:A:2024:HOH:O	2.40	0.43
1:G:158:GLU:C	4:G:2093:HOH:O	2.57	0.43
1:F:133:THR:HG23	1:H:143:ARG:HG3	2.01	0.43
1:B:158:GLU:HB3	1:K:101:GLN:NE2	2.33	0.43
1:D:40:HIS:CE1	1:D:52:HIS:CE1	3.07	0.43
1:I:133:THR:CG2	1:K:143:ARG:HG3	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:HB3	1:D:157:ALA:HB3	2.01	0.42
1:G:6:PRO:HA	1:I:153:LEU:HD11	2.01	0.42
1:I:65:TYR:OH	1:I:144:GLY:HA3	2.19	0.42
1:C:6:PRO:O	1:C:7:ALA:CB	2.66	0.42
1:A:158:GLU:HA	1:F:101:GLN:OE1	2.20	0.42
1:B:110:VAL:HG11	1:B:153:LEU:HD13	2.02	0.42
1:C:90:LYS:HD3	1:C:90:LYS:HA	1.88	0.42
1:A:46:SER:O	1:D:47:ASN:ND2	2.53	0.42
1:B:82:SER:OG	4:B:2023:HOH:O	2.11	0.41
1:L:31:ALA:HB2	1:L:91:THR:HG21	2.02	0.41
1:I:133:THR:HG23	1:K:143:ARG:HG3	2.01	0.41
1:A:113:ILE:HG23	1:A:145:VAL:CG1	2.50	0.41
1:L:129:GLU:CG	4:L:2044:HOH:O	2.68	0.41
1:E:51:ILE:HG23	1:E:51:ILE:HD12	1.84	0.41
1:G:65:TYR:OH	1:G:144:GLY:HA3	2.20	0.41
1:F:121:GLN:OE1	1:F:142:LYS:HE3	2.20	0.41
1:C:9:THR:HA	1:C:10:ALA:HA	1.74	0.41
1:C:65:TYR:OH	1:C:144:GLY:HA3	2.20	0.41
1:H:8:LEU:HB2	1:J:153:LEU:HD21	2.03	0.41
1:L:17:ALA:O	1:L:20:GLN:HG2	2.20	0.41
1:D:11:ASP:HA	1:D:12:PRO:HD3	2.02	0.41
1:K:157:ALA:O	1:K:158:GLU:CG	2.68	0.41
1:A:133:THR:HG23	1:C:143:ARG:HG3	2.01	0.41
1:E:65:TYR:OH	1:E:144:GLY:HA3	2.19	0.41
1:J:11:ASP:HA	1:J:12:PRO:HD2	1.73	0.41
1:E:9:THR:CA	1:E:10:ALA:CB	2.92	0.40
1:L:157:ALA:O	1:L:158:GLU:HB2	2.21	0.40
1:L:18:ALA:HA	1:L:127:LEU:HD21	2.02	0.40
1:C:67:ASP:OD1	1:I:37:LYS:HE3	2.22	0.40

There are no symmetry-related clashes.

## 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	148/161 (92%)	147 (99%)	1 (1%)	0	100	100
1	B	150/161 (93%)	146 (97%)	3 (2%)	1 (1%)	26	19
1	C	155/161 (96%)	150 (97%)	2 (1%)	3 (2%)	10	4
1	D	153/161 (95%)	149 (97%)	2 (1%)	2 (1%)	15	7
1	E	150/161 (93%)	146 (97%)	3 (2%)	1 (1%)	26	19
1	F	150/161 (93%)	144 (96%)	6 (4%)	0	100	100
1	G	151/161 (94%)	149 (99%)	1 (1%)	1 (1%)	26	19
1	H	151/161 (94%)	147 (97%)	1 (1%)	3 (2%)	9	3
1	I	146/161 (91%)	145 (99%)	1 (1%)	0	100	100
1	J	146/161 (91%)	146 (100%)	0	0	100	100
1	K	151/161 (94%)	145 (96%)	4 (3%)	2 (1%)	15	7
1	L	151/161 (94%)	148 (98%)	1 (1%)	2 (1%)	15	7
All	All	1802/1932 (93%)	1762 (98%)	25 (1%)	15 (1%)	24	15

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	8	LEU
1	B	10	ALA
1	C	7	ALA
1	D	8	LEU
1	D	10	ALA
1	K	10	ALA
1	L	10	ALA
1	C	6	PRO
1	E	10	ALA
1	L	9	THR
1	C	5	THR
1	H	8	LEU
1	H	11	ASP
1	H	10	ALA
1	K	11	ASP

### 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	117/127 (92%)	111 (95%)	6 (5%)	29	23
1	B	118/127 (93%)	111 (94%)	7 (6%)	24	18
1	C	118/127 (93%)	110 (93%)	8 (7%)	20	13
1	D	120/127 (94%)	115 (96%)	5 (4%)	36	31
1	E	117/127 (92%)	113 (97%)	4 (3%)	44	41
1	F	117/127 (92%)	112 (96%)	5 (4%)	35	30
1	G	118/127 (93%)	111 (94%)	7 (6%)	24	18
1	H	118/127 (93%)	113 (96%)	5 (4%)	36	31
1	I	116/127 (91%)	112 (97%)	4 (3%)	44	41
1	J	116/127 (91%)	109 (94%)	7 (6%)	24	17
1	K	118/127 (93%)	116 (98%)	2 (2%)	68	71
1	L	119/127 (94%)	113 (95%)	6 (5%)	30	24
All	All	1412/1524 (93%)	1346 (95%)	66 (5%)	32	27

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	56	ASP
1	A	57	SER
1	A	80	ILE
1	A	91	THR
1	A	158	GLU
1	B	9	THR
1	B	44	ARG
1	B	56	ASP
1	B	57	SER
1	B	80	ILE
1	B	91	THR
1	B	158	GLU
1	C	9	THR
1	C	13	GLU
1	C	44	ARG
1	C	56	ASP
1	C	57	SER
1	C	67	ASP

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Mol	Chain	Res	Type
1	C	91	THR
1	C	153	LEU
1	D	44	ARG
1	D	56	ASP
1	D	57	SER
1	D	91	THR
1	D	158	GLU
1	E	44	ARG
1	E	56	ASP
1	E	57	SER
1	E	67	ASP
1	F	8	LEU
1	F	44	ARG
1	F	56	ASP
1	F	80	ILE
1	F	91	THR
1	G	13	GLU
1	G	44	ARG
1	G	56	ASP
1	G	57	SER
1	G	80	ILE
1	G	91	THR
1	G	158	GLU
1	H	13	GLU
1	H	44	ARG
1	H	56	ASP
1	H	80	ILE
1	H	91	THR
1	I	11	ASP
1	I	44	ARG
1	I	67	ASP
1	I	91	THR
1	J	44	ARG
1	J	56	ASP
1	J	57	SER
1	J	67	ASP
1	J	80	ILE
1	J	91	THR
1	J	111	SER
1	K	44	ARG
1	K	91	THR
1	L	8	LEU

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Mol	Chain	Res	Type
1	L	13	GLU
1	L	44	ARG
1	L	56	ASP
1	L	57	SER
1	L	91	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	63	GLN
1	A	101	GLN
1	B	20	GLN
1	C	20	GLN
1	C	101	GLN
1	D	20	GLN
1	D	30	GLN
1	D	63	GLN
1	D	101	GLN
1	E	101	GLN
1	F	20	GLN
1	G	30	GLN
1	G	63	GLN
1	G	101	GLN
1	H	30	GLN
1	H	63	GLN
1	I	30	GLN
1	I	63	GLN
1	J	20	GLN
1	J	101	GLN
1	K	20	GLN
1	K	30	GLN
1	K	63	GLN
1	K	101	GLN
1	L	30	GLN
1	L	63	GLN
1	L	101	GLN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 6 are monoatomic - leaving 11 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	OFE	A	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	B	1159	1	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	C	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	D	1159	1	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	E	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	F	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	G	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	H	1159	1,4	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	J	1159	1	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	K	1159	1	0,1,1	0.00	-	0,0,0	0.00	-
2	OFE	L	1159	1	0,1,1	0.00	-	0,0,0	0.00	-

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	OFE	A	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	B	1159	1	-	0/0/0/0	0/0/0/0
2	OFE	C	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	D	1159	1	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OFE	E	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	F	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	G	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	H	1159	1,4	-	0/0/0/0	0/0/0/0
2	OFE	J	1159	1	-	0/0/0/0	0/0/0/0
2	OFE	K	1159	1	-	0/0/0/0	0/0/0/0
2	OFE	L	1159	1	-	0/0/0/0	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.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1159	OFE	1	0
2	B	1159	OFE	1	0
2	D	1159	OFE	1	0
2	E	1159	OFE	1	0
2	L	1159	OFE	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	149/161 (92%)	-0.50	2 (1%) 79 80	20, 29, 43, 68	0
1	B	152/161 (94%)	-0.31	6 (3%) 43 45	18, 27, 47, 72	0
1	C	157/161 (97%)	0.43	14 (8%) 12 13	24, 33, 88, 115	0
1	D	152/161 (94%)	-0.14	7 (4%) 36 38	20, 29, 54, 91	0
1	E	152/161 (94%)	-0.02	9 (5%) 26 27	20, 31, 60, 96	0
1	F	152/161 (94%)	-0.19	6 (3%) 43 45	19, 28, 46, 83	0
1	G	153/161 (95%)	-0.36	5 (3%) 50 51	18, 26, 52, 67	0
1	H	152/161 (94%)	-0.17	6 (3%) 43 45	16, 25, 50, 76	0
1	I	148/161 (91%)	-0.24	7 (4%) 35 37	21, 31, 47, 72	0
1	J	148/161 (91%)	-0.33	4 (2%) 58 58	19, 28, 41, 68	0
1	K	152/161 (94%)	0.01	9 (5%) 26 27	21, 31, 58, 92	0
1	L	152/161 (94%)	-0.30	4 (2%) 59 60	20, 28, 47, 93	0
All	All	1819/1932 (94%)	-0.17	79 (4%) 39 40	16, 29, 53, 115	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	3	ILE	18.7
1	C	5	THR	15.9
1	C	4	THR	14.7
1	F	8	LEU	13.2
1	C	8	LEU	11.9
1	C	7	ALA	11.9
1	C	10	ALA	11.0
1	H	9	THR	11.0
1	K	8	LEU	10.6
1	C	9	THR	10.5
1	L	8	LEU	10.4

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Mol	Chain	Res	Type	RSRZ
1	E	8	LEU	9.9
1	D	8	LEU	9.3
1	E	7	ALA	9.1
1	E	9	THR	9.1
1	K	7	ALA	8.5
1	E	10	ALA	8.3
1	F	7	ALA	8.2
1	D	10	ALA	7.9
1	K	10	ALA	7.8
1	F	9	THR	7.6
1	D	9	THR	7.5
1	B	158	GLU	7.4
1	K	9	THR	7.1
1	C	6	PRO	7.0
1	B	7	ALA	7.0
1	J	12	PRO	6.8
1	L	7	ALA	6.8
1	C	12	PRO	6.2
1	C	11	ASP	6.2
1	J	11	ASP	6.1
1	H	7	ALA	6.0
1	E	11	ASP	5.9
1	K	11	ASP	5.8
1	G	158	GLU	5.8
1	K	158	GLU	5.5
1	D	7	ALA	5.4
1	L	9	THR	5.2
1	I	13	GLU	5.1
1	I	11	ASP	5.0
1	F	158	GLU	4.8
1	E	14	VAL	4.8
1	K	12	PRO	4.7
1	I	12	PRO	4.6
1	H	10	ALA	4.5
1	B	157	ALA	4.4
1	C	13	GLU	4.4
1	G	7	ALA	4.4
1	C	158	GLU	4.4
1	H	8	LEU	4.4
1	D	11	ASP	4.2
1	G	8	LEU	3.9
1	E	12	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
1	J	13	GLU	3.8
1	I	158	GLU	3.6
1	K	13	GLU	3.6
1	E	13	GLU	3.5
1	H	158	GLU	3.4
1	L	158	GLU	3.4
1	I	15	ALA	3.4
1	F	129	GLU	3.3
1	I	14	VAL	3.2
1	D	158	GLU	3.0
1	A	158	GLU	2.8
1	C	129	GLU	2.8
1	G	129	GLU	2.7
1	E	129	GLU	2.7
1	K	129	GLU	2.7
1	J	158	GLU	2.7
1	F	10	ALA	2.6
1	A	10	ALA	2.6
1	G	157	ALA	2.5
1	C	157	ALA	2.5
1	D	12	PRO	2.5
1	B	10	ALA	2.5
1	I	16	ALA	2.4
1	H	11	ASP	2.3
1	B	11	ASP	2.2
1	B	129	GLU	2.1

## 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	OFE	H	1159	2/2	0.99	0.10	0.59	26,26,26,44	1
2	OFE	F	1159	2/2	0.99	0.10	0.13	27,27,27,35	1
2	OFE	L	1159	2/2	0.99	0.05	-2.01	18,18,18,46	1
2	OFE	J	1159	2/2	0.99	0.05	-2.48	19,19,19,39	1
2	OFE	D	1159	2/2	0.99	0.07	-3.26	19,19,19,41	1
2	OFE	K	1159	2/2	0.99	0.06	-3.26	23,23,23,40	1
2	OFE	B	1159	2/2	0.99	0.04	-3.45	38,38,38,42	0
2	OFE	A	1159	2/2	0.99	0.04	-3.79	24,24,24,42	1
2	OFE	C	1159	2/2	1.00	0.03	-3.87	15,15,15,41	1
2	OFE	E	1159	2/2	1.00	0.04	-4.24	19,19,19,41	1
3	FE	I	1159	1/1	0.99	0.03	-4.32	34,34,34,34	0
2	OFE	G	1159	2/2	1.00	0.04	-5.81	17,17,17,34	1
3	FE	D	1160	1/1	0.91	0.07	-	14,14,14,14	1
3	FE	C	1160	1/1	0.97	0.39	-	19,19,19,19	1
3	FE	L	1160	1/1	0.98	0.12	-	5,5,5,5	1
3	FE	B	1160	1/1	0.92	0.07	-	15,15,15,15	1
3	FE	J	1160	1/1	0.95	0.07	-	6,6,6,6	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.