



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 07:34 am BST

PDB ID : 2BW1
Title : Iron-bound crystal structure of Dps-like peroxide resistance protein (Dpr) from *Streptococcus suis*.
Authors : Kauko, A.; Pulliainen, A.; Haataja, S.; Finne, J.; Papageorgiou, A.C.
Deposited on : 2005-07-07
Resolution : 1.81 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

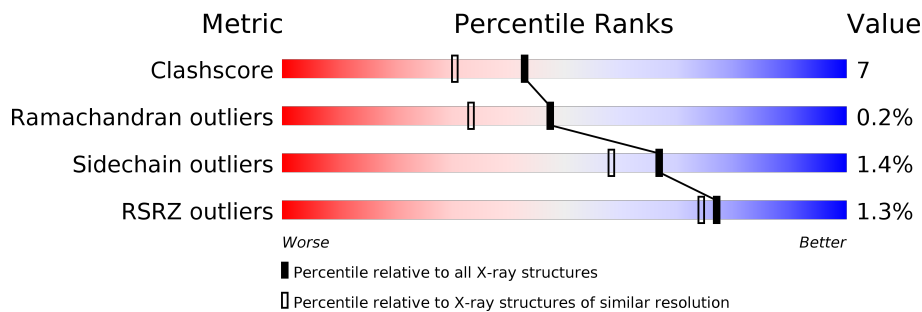
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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))
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	165	
1	B	165	
1	C	165	
1	D	165	
1	E	165	
1	F	165	
1	G	165	

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Mol	Chain	Length	Quality of chain
1	H	165	<p>% 84% 8% 8%</p>
1	I	165	<p>% 85% 6% 8%</p>
1	J	165	<p>% 79% 10% 8%</p>
1	K	165	<p>% 83% 9% 7%</p>
1	L	165	<p>% 82% 9% 8%</p>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16600 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.

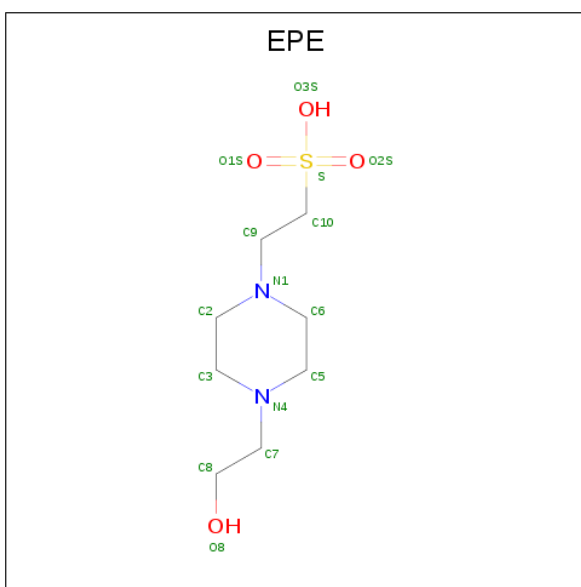
- Molecule 1 is a protein called DPS-LIKE PEROXIDE RESISTANCE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	151	1258	800	205	246	7	0	9	0
1	B	151	1250	796	204	243	7	0	8	0
1	C	151	1255	800	202	246	7	0	8	0
1	D	152	1256	796	206	247	7	0	8	0
1	E	151	1241	789	201	245	6	0	6	0
1	F	151	1249	792	204	246	7	0	8	0
1	G	165	1322	840	215	260	7	0	6	0
1	H	151	1231	783	200	242	6	0	5	0
1	I	151	1229	781	200	241	7	0	5	0
1	J	151	1255	797	205	247	6	0	9	0
1	K	153	1246	792	206	242	6	0	5	0
1	L	151	1233	784	202	241	6	0	6	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	8	GLY	GLN	engineered mutation	UNP Q9F5J9

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Fe	0	0
			1	1		
3	J	1	Total	Fe	0	0
			1	1		
3	D	1	Total	Fe	0	0
			1	1		
3	K	1	Total	Fe	0	0
			1	1		
3	E	1	Total	Fe	0	0
			1	1		
3	H	1	Total	Fe	0	0
			1	1		
3	B	1	Total	Fe	0	0
			1	1		
3	I	1	Total	Fe	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total 1	Fe 1	0	0
3	A	1	Total 1	Fe 1	0	0
3	L	1	Total 1	Fe 1	0	0
3	F	1	Total 1	Fe 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	K	1	Total 1	Ca 1	0	0

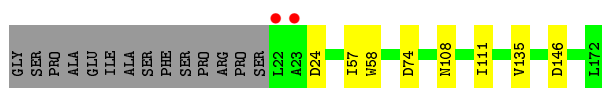
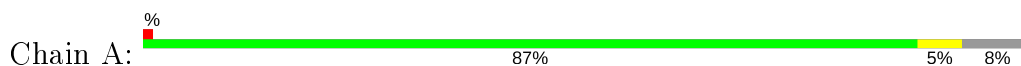
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	145	Total 145	O 145	0	0
5	B	128	Total 128	O 128	0	0
5	C	124	Total 124	O 124	0	0
5	D	151	Total 151	O 151	0	0
5	E	126	Total 126	O 126	0	0
5	F	98	Total 98	O 98	0	0
5	G	146	Total 146	O 146	0	0
5	H	99	Total 99	O 99	0	0
5	I	124	Total 124	O 124	0	0
5	J	93	Total 93	O 93	0	0
5	K	147	Total 147	O 147	0	0
5	L	121	Total 121	O 121	0	0

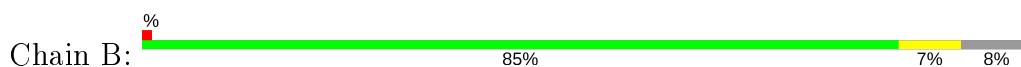
3 Residue-property plots [i](#)

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

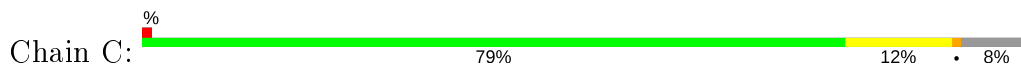
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



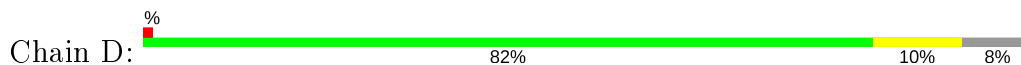
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



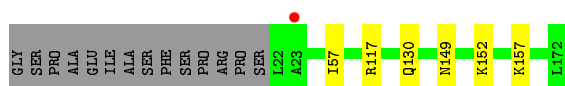
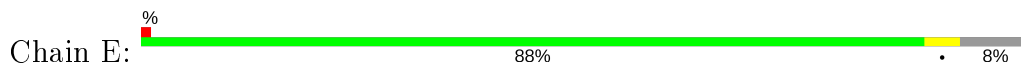
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



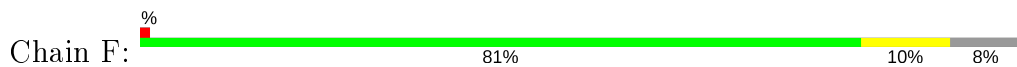
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



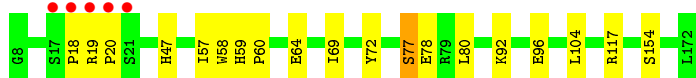
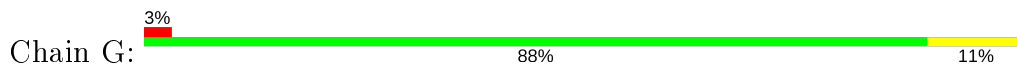
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



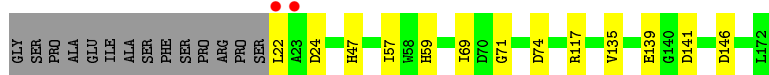
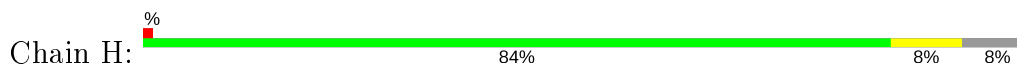
- Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



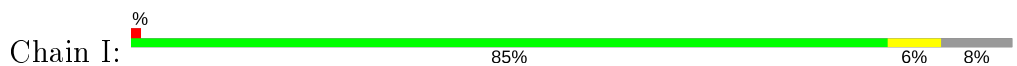
● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



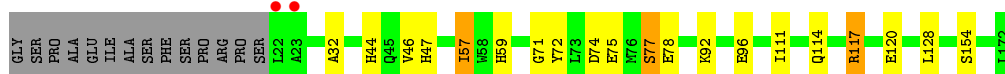
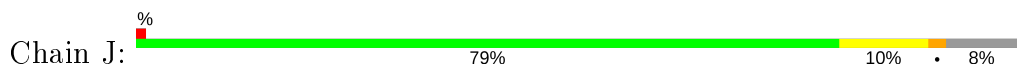
● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



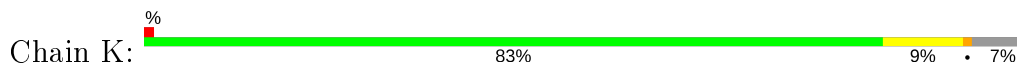
● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



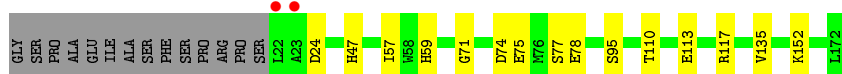
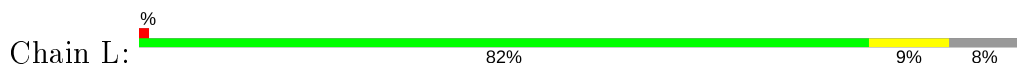
● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



● Molecule 1: DPS-LIKE PEROXIDE RESISTANCE PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.13Å 138.05Å 142.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.53 – 1.81 19.53 – 1.81	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.53-1.81) 99.3 (19.53-1.81)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 1.81Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.187 , 0.230 0.187 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtrriage
Anisotropy	0.043	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.024 for -h,l,k	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16600	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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: CA, EPE, FE

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	0.66	0/1303	0.64	0/1759
1	B	0.66	0/1291	0.68	0/1743
1	C	1.24	2/1297 (0.2%)	0.68	0/1752
1	D	0.72	0/1297	0.68	0/1751
1	E	0.62	0/1275	0.65	0/1725
1	F	0.86	2/1290 (0.2%)	0.70	2/1743 (0.1%)
1	G	1.12	2/1361 (0.1%)	0.69	0/1839
1	H	0.64	0/1264	0.64	0/1710
1	I	0.70	0/1262	0.65	0/1706
1	J	1.03	2/1300 (0.2%)	0.59	0/1756
1	K	1.14	2/1279 (0.2%)	0.69	1/1728 (0.1%)
1	L	0.83	2/1264 (0.2%)	0.68	2/1707 (0.1%)
All	All	0.88	12/15483 (0.1%)	0.66	5/20919 (0.0%)

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	K	0	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	77[A]	SER	C-N	26.91	1.96	1.34
1	C	77[B]	SER	C-N	26.91	1.96	1.34
1	G	77[A]	SER	C-N	23.63	1.88	1.34
1	G	77[B]	SER	C-N	23.63	1.88	1.34
1	K	77[A]	SER	C-N	22.62	1.86	1.34
1	K	77[B]	SER	C-N	22.62	1.86	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	77[A]	SER	C-N	21.00	1.82	1.34
1	J	77[B]	SER	C-N	21.00	1.82	1.34
1	F	77[A]	SER	C-N	14.17	1.66	1.34
1	F	77[B]	SER	C-N	14.17	1.66	1.34
1	L	77[A]	SER	C-N	12.52	1.62	1.34
1	L	77[B]	SER	C-N	12.52	1.62	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	77[A]	SER	O-C-N	-6.61	112.12	122.70
1	F	77[B]	SER	O-C-N	-6.61	112.12	122.70
1	L	77[A]	SER	O-C-N	-5.40	114.06	122.70
1	L	77[B]	SER	O-C-N	-5.40	114.06	122.70
1	K	51	ARG	NE-CZ-NH2	-5.12	117.74	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	K	20	PRO	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1258	0	1203	32	0
1	B	1250	0	1201	16	0
1	C	1255	0	1198	34	0
1	D	1256	0	1198	12	0
1	E	1241	0	1172	16	0
1	F	1249	0	1181	23	0
1	G	1322	0	1249	31	0
1	H	1231	0	1169	24	0
1	I	1229	0	1168	15	0
1	J	1255	0	1195	11	0
1	K	1246	0	1193	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1233	0	1169	29	0
2	A	15	0	17	0	0
2	B	15	0	17	1	0
2	C	15	0	17	0	0
2	D	15	0	17	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	K	1	0	0	0	0
5	A	145	0	0	2	0
5	B	128	0	0	3	0
5	C	124	0	0	1	0
5	D	151	0	0	3	0
5	E	126	0	0	4	0
5	F	98	0	0	3	0
5	G	146	0	0	3	0
5	H	99	0	0	0	0
5	I	124	0	0	0	0
5	J	93	0	0	0	0
5	K	147	0	0	0	0
5	L	121	0	0	4	0
All	All	16600	0	14364	201	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 (201) 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:C:57[B]:ILE:HD11	1:L:57[B]:ILE:CG1	1.26	1.62
1:A:57[B]:ILE:CD1	1:H:57[B]:ILE:HD11	1.24	1.56
1:C:77[B]:SER:C	1:C:78[B]:GLU:CA	1.86	1.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57[B]:ILE:HD11	1:F:57[B]:ILE:CD1	1.47	1.42
1:C:57[B]:ILE:CD1	1:L:57[B]:ILE:HG13	1.47	1.41
1:B:57[B]:ILE:CD1	1:F:57[B]:ILE:HD11	1.52	1.39
1:G:77[B]:SER:C	1:G:78[B]:GLU:CA	1.95	1.33
1:J:77[A]:SER:C	1:J:78[A]:GLU:N	1.82	1.31
1:A:57[B]:ILE:CD1	1:K:57[B]:ILE:HD11	1.59	1.31
1:K:77[B]:SER:C	1:K:78[B]:GLU:CA	1.97	1.30
1:C:57[B]:ILE:HD11	1:L:57[B]:ILE:CD1	1.58	1.30
1:A:57[B]:ILE:CG1	1:K:57[B]:ILE:HD11	1.60	1.30
1:C:77[B]:SER:CA	1:C:78[B]:GLU:N	1.92	1.29
1:K:77[A]:SER:C	1:K:78[A]:GLU:N	1.86	1.29
1:G:77[A]:SER:C	1:G:78[A]:GLU:N	1.88	1.25
1:G:77[B]:SER:CA	1:G:78[B]:GLU:N	2.01	1.24
1:A:57[B]:ILE:HD11	1:H:57[B]:ILE:CD1	1.68	1.23
1:A:57[B]:ILE:CD1	1:K:57[B]:ILE:CG1	2.18	1.22
1:A:57[B]:ILE:CD1	1:H:57[B]:ILE:CD1	2.17	1.21
1:K:77[B]:SER:CA	1:K:78[B]:GLU:N	2.04	1.20
1:A:57[B]:ILE:HD11	1:K:57[B]:ILE:CG1	1.71	1.19
1:A:57[B]:ILE:CD1	1:K:57[B]:ILE:CD1	2.20	1.19
1:C:77[A]:SER:C	1:C:78[A]:GLU:N	1.95	1.18
1:C:57[B]:ILE:CD1	1:L:57[B]:ILE:CG1	2.09	1.15
1:A:57[B]:ILE:HG12	1:K:57[B]:ILE:HD11	1.24	1.14
1:A:57[B]:ILE:HD11	1:K:57[B]:ILE:CD1	1.78	1.14
1:K:77[B]:SER:O	1:K:78[B]:GLU:N	1.79	1.13
1:C:77[B]:SER:O	1:C:78[B]:GLU:N	1.79	1.12
1:C:57[B]:ILE:CD1	1:L:57[B]:ILE:CD1	2.25	1.11
1:E:57:ILE:HD11	1:L:57[B]:ILE:HD11	1.14	1.10
1:G:77[B]:SER:O	1:G:78[B]:GLU:N	1.83	1.09
1:B:57[B]:ILE:CD1	1:F:57[B]:ILE:CD1	2.17	1.09
1:A:24:ASP:HB3	1:A:135:VAL:HG11	1.36	1.07
1:A:57[B]:ILE:HD13	1:K:57[B]:ILE:HG12	1.33	1.06
1:G:57[B]:ILE:CG2	1:I:57[B]:ILE:HD11	1.84	1.06
1:B:57[B]:ILE:HG13	1:F:57[B]:ILE:HD13	1.34	1.05
1:A:57[B]:ILE:HD11	1:K:57[B]:ILE:HG13	1.39	1.03
1:A:57[B]:ILE:CG1	1:H:57[B]:ILE:HD11	1.89	1.01
1:A:57[B]:ILE:HG13	1:H:57[B]:ILE:HD13	1.43	1.00
1:E:57:ILE:CD1	1:L:57[B]:ILE:HD11	1.93	0.99
1:G:57[B]:ILE:HG21	1:I:57[B]:ILE:HD11	1.45	0.97
1:C:77[B]:SER:O	1:C:78[B]:GLU:CA	2.08	0.97
1:B:57[B]:ILE:CG1	1:F:57[B]:ILE:HD13	1.95	0.97
1:G:57[B]:ILE:CG2	1:I:57[B]:ILE:CD1	2.43	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57[B]:ILE:CG1	1:H:57[B]:ILE:CD1	2.43	0.96
1:C:57[B]:ILE:HD11	1:L:57[B]:ILE:HG13	1.03	0.96
1:C:57[B]:ILE:HD13	1:L:57[B]:ILE:HG13	1.45	0.95
1:C:57[B]:ILE:CD1	1:L:57[B]:ILE:HD11	1.93	0.95
1:C:57[B]:ILE:HD11	1:L:57[B]:ILE:HD11	1.45	0.95
1:A:57[B]:ILE:HD11	1:K:57[B]:ILE:HD11	1.41	0.94
1:A:57[B]:ILE:HG13	1:H:57[B]:ILE:CD1	1.97	0.94
1:L:74[A]:ASP:C	1:L:75[A]:GLU:N	2.19	0.94
1:B:57[B]:ILE:CG1	1:F:57[B]:ILE:CD1	2.44	0.94
1:A:57[B]:ILE:HD13	1:K:57[B]:ILE:CG1	1.90	0.93
1:A:24:ASP:HB3	1:A:135:VAL:CG1	1.99	0.92
1:K:77[B]:SER:O	1:K:78[B]:GLU:CA	2.15	0.92
1:G:77[B]:SER:C	1:G:78[B]:GLU:N	0.81	0.85
1:K:77[B]:SER:C	1:K:78[B]:GLU:N	0.81	0.85
1:G:77[B]:SER:O	1:G:78[B]:GLU:CA	2.20	0.85
1:F:57[B]:ILE:HG13	1:J:57:ILE:HD13	1.58	0.85
1:C:57[B]:ILE:HG13	1:E:57:ILE:HD11	1.58	0.83
1:A:57[B]:ILE:HG12	1:K:57[B]:ILE:CD1	2.08	0.82
1:C:57[B]:ILE:HG13	1:E:57:ILE:CD1	2.10	0.82
1:C:57[B]:ILE:HD11	1:L:57[B]:ILE:HG12	1.56	0.81
1:C:77[B]:SER:C	1:C:78[B]:GLU:N	0.73	0.79
1:E:157:LYS:NZ	5:E:2116:HOH:O	2.17	0.78
1:C:77[B]:SER:O	1:C:78[B]:GLU:HA	1.82	0.78
1:H:57[B]:ILE:HD11	1:K:57[B]:ILE:HD11	1.67	0.77
1:J:72:TYR:HH	1:J:154[B]:SER:HG	1.29	0.77
1:D:63[A]:ASP:OD2	5:D:2045:HOH:O	2.01	0.76
1:A:57[B]:ILE:HD11	1:H:57[B]:ILE:HD11	0.76	0.75
2:B:200:EPE:H51	5:B:2127:HOH:O	1.85	0.75
1:G:57[B]:ILE:HG22	1:I:57[B]:ILE:CD1	2.15	0.75
1:C:77[B]:SER:C	1:C:78[B]:GLU:HA	2.05	0.75
1:F:152:LYS:HE3	5:F:2074:HOH:O	1.87	0.74
1:A:57[B]:ILE:CD1	1:K:57[B]:ILE:HG12	1.98	0.72
1:F:152:LYS:CE	5:F:2074:HOH:O	2.37	0.72
1:C:57[B]:ILE:HD12	1:L:57[B]:ILE:CD1	2.18	0.71
1:L:24:ASP:HB3	1:L:135:VAL:CG1	2.19	0.71
1:H:22:LEU:HD21	1:H:141:ASP:HB2	1.74	0.70
1:A:57[B]:ILE:HD12	1:H:57[B]:ILE:HD11	1.59	0.70
1:F:57[B]:ILE:CG1	1:J:57:ILE:HD13	2.23	0.69
1:K:77[B]:SER:O	1:K:78[B]:GLU:HA	1.91	0.69
1:D:146:ASP:OD1	5:D:2127:HOH:O	2.11	0.68
1:A:57[B]:ILE:HD13	1:K:57[B]:ILE:CD1	2.18	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:57[B]:ILE:CG2	1:I:57[B]:ILE:HD13	2.24	0.68
1:C:40:HIS:CE1	1:C:66[B]:MET:SD	2.88	0.67
1:D:57:ILE:HD11	1:G:57[A]:ILE:HD13	1.77	0.67
1:A:108:ASN:HB2	5:A:2094:HOH:O	1.93	0.67
1:D:57:ILE:CD1	1:G:57[A]:ILE:HD13	2.24	0.66
1:H:57[B]:ILE:HD11	1:K:57[B]:ILE:CD1	2.26	0.66
1:B:57[B]:ILE:CG1	1:F:57[B]:ILE:HD11	2.14	0.65
1:E:57:ILE:HD11	1:L:57[B]:ILE:CD1	2.09	0.65
1:D:72:TYR:HH	1:D:154[B]:SER:HG	1.39	0.64
1:L:117:ARG:HG3	5:L:2089:HOH:O	1.96	0.64
1:G:77[B]:SER:O	1:G:78[B]:GLU:HA	1.95	0.63
1:B:57[B]:ILE:HD11	1:F:57[B]:ILE:HD11	0.69	0.63
1:B:71:GLY:O	1:B:74[B]:ASP:HB2	1.98	0.62
1:C:57[B]:ILE:CG1	1:E:57:ILE:HD11	2.28	0.62
1:F:24:ASP:HB3	1:F:135:VAL:HG11	1.81	0.62
1:B:24:ASP:HB2	1:B:139:GLU:OE2	2.00	0.62
1:L:24:ASP:HB3	1:L:135:VAL:HG13	1.82	0.61
1:H:22:LEU:CD2	1:H:141:ASP:HB2	2.30	0.61
1:H:24:ASP:HB3	1:H:135:VAL:CG1	2.31	0.61
1:L:152:LYS:HE3	5:L:2096:HOH:O	1.99	0.61
1:B:127:ALA:O	1:B:131:LYS:HG3	2.01	0.60
1:C:130:GLN:OE1	1:C:152:LYS:HE3	2.02	0.60
1:H:24:ASP:HB3	1:H:135:VAL:HG11	1.83	0.60
1:G:92:LYS:HD3	5:G:2033:HOH:O	2.01	0.59
1:J:117[B]:ARG:NH1	1:J:120:GLU:HB2	2.17	0.59
1:I:40:HIS:CE1	1:I:66[B]:MET:HG2	2.39	0.57
1:L:24:ASP:HB3	1:L:135:VAL:HG11	1.84	0.57
1:C:68:GLU:OE2	1:C:72[B]:TYR:OH	2.13	0.57
1:F:24:ASP:HB3	1:F:135:VAL:CG1	2.37	0.55
1:H:57[B]:ILE:CD1	1:K:57[B]:ILE:HD11	2.36	0.55
1:A:57[B]:ILE:CD1	1:K:57[B]:ILE:HG13	2.10	0.55
1:C:92:LYS:O	1:C:96:GLU:HG3	2.07	0.54
1:J:92:LYS:NZ	1:J:96:GLU:OE1	2.41	0.54
1:C:135:VAL:O	1:C:139:GLU:HG3	2.07	0.54
1:A:24:ASP:CB	1:A:135:VAL:CG1	2.81	0.53
1:E:57:ILE:CD1	1:L:57[B]:ILE:CD1	2.77	0.53
1:G:92:LYS:NZ	1:G:96:GLU:OE1	2.41	0.53
1:C:152:LYS:CE	5:C:2100:HOH:O	2.56	0.53
1:I:24:ASP:HB3	1:I:135:VAL:CG1	2.38	0.53
1:G:59:HIS:HB3	1:G:60:PRO:HD3	1.92	0.52
1:C:77[B]:SER:N	1:C:78[B]:GLU:N	2.54	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:57[B]:ILE:HG22	1:I:57[B]:ILE:HD13	1.86	0.51
1:L:71:GLY:O	1:L:74[B]:ASP:HB2	2.10	0.51
1:L:75[A]:GLU:OE2	5:L:2053:HOH:O	2.19	0.50
1:C:57[B]:ILE:CD1	1:E:57:ILE:HD11	2.42	0.50
1:D:40:HIS:CD2	1:D:70:ASP:OD2	2.65	0.50
1:K:28:VAL:HG11	1:K:132:GLY:HA2	1.94	0.50
1:H:57[B]:ILE:CG1	1:K:57[B]:ILE:HD12	2.42	0.50
1:G:117:ARG:HG3	5:G:2106:HOH:O	2.12	0.50
1:J:71:GLY:O	1:J:74[B]:ASP:HB2	2.12	0.49
1:B:40:HIS:HE1	5:B:2015:HOH:O	1.94	0.49
1:D:71:GLY:O	1:D:74[B]:ASP:HB2	2.12	0.49
1:G:57[B]:ILE:HG23	1:I:57[B]:ILE:CD1	2.40	0.49
1:G:77[B]:SER:N	1:G:78[B]:GLU:N	2.59	0.49
1:L:152:LYS:CE	5:L:2096:HOH:O	2.60	0.49
1:H:71:GLY:O	1:H:74[B]:ASP:HB2	2.13	0.49
1:D:57:ILE:HD13	1:G:57[A]:ILE:HD13	1.95	0.49
1:J:117[B]:ARG:HH12	1:J:120:GLU:HB2	1.77	0.48
1:H:57[B]:ILE:CG1	1:K:57[B]:ILE:CD1	2.90	0.48
1:G:47:HIS:CE1	1:G:59:HIS:CE1	3.01	0.48
1:H:57[B]:ILE:HG12	1:K:57[B]:ILE:HD12	1.95	0.48
1:B:160:TRP:CH2	1:L:78[A]:GLU:HG2	2.48	0.48
1:B:57[B]:ILE:CD1	1:F:57[B]:ILE:HD13	2.19	0.48
1:E:117:ARG:HD3	5:E:2097:HOH:O	2.12	0.48
1:I:40:HIS:ND1	1:I:66[B]:MET:HG2	2.29	0.48
1:H:47:HIS:CE1	1:H:59:HIS:CE1	3.02	0.47
1:I:135:VAL:O	1:I:139:GLU:HG3	2.14	0.47
1:C:71:GLY:O	1:C:74[B]:ASP:HB2	2.13	0.47
1:K:46:VAL:O	1:K:50:MET:HB2	2.15	0.46
1:L:74[B]:ASP:O	1:L:78[B]:GLU:HB2	2.14	0.46
1:A:57[B]:ILE:HD12	1:H:57[B]:ILE:CD1	2.29	0.46
1:G:77[B]:SER:C	1:G:78[B]:GLU:HA	2.14	0.46
1:F:74[B]:ASP:O	1:F:78[B]:GLU:HB2	2.15	0.45
1:G:57[B]:ILE:HD11	1:G:58:TRP:CE2	2.51	0.45
1:I:24:ASP:HB3	1:I:135:VAL:HG11	1.99	0.45
1:E:130:GLN:OE1	1:E:152:LYS:HE3	2.16	0.45
1:G:64:GLU:OE2	5:G:2052:HOH:O	2.20	0.45
1:I:57[A]:ILE:HG23	1:I:58:TRP:CD1	2.52	0.45
1:A:74[B]:ASP:HB2	5:A:2057:HOH:O	2.16	0.44
1:L:110:THR:OG1	1:L:113:GLU:HG3	2.17	0.44
1:D:56:MET:CG	1:G:57[B]:ILE:HD13	2.48	0.44
2:D:200:EPE:H51	5:D:2149:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:32:ALA:HA	1:J:128:LEU:HD21	2.00	0.44
1:B:40:HIS:ND1	5:B:2011:HOH:O	1.93	0.43
1:E:152:LYS:CE	5:E:2104:HOH:O	2.65	0.43
1:C:57[B]:ILE:HG13	1:E:57:ILE:HD13	1.97	0.43
1:D:72:TYR:OH	1:D:154[B]:SER:OG	2.16	0.43
1:G:72:TYR:HH	1:G:154[B]:SER:HB3	1.82	0.43
1:A:146:ASP:OD1	1:E:149:ASN:ND2	2.52	0.42
1:L:47:HIS:CE1	1:L:59:HIS:CE1	3.07	0.42
1:F:82:THR:HG22	1:K:172:LEU:HD11	2.01	0.42
1:G:59:HIS:HB3	1:G:60:PRO:CD	2.49	0.42
1:F:28:VAL:HG11	1:F:132:GLY:HA2	2.00	0.42
1:C:57[A]:ILE:HG23	1:C:58:TRP:CD1	2.54	0.42
1:D:62:MET:O	1:D:66[B]:MET:HG3	2.19	0.42
1:F:143:VAL:HG23	1:K:152:LYS:HE3	2.02	0.42
1:K:47:HIS:CE1	1:K:59:HIS:CE1	3.08	0.42
1:G:77[A]:SER:O	1:G:80:LEU:HB3	2.20	0.41
1:I:40:HIS:CE1	1:I:66[B]:MET:SD	3.13	0.41
1:D:130:GLN:OE1	1:D:152:LYS:HE3	2.20	0.41
1:F:64:GLU:OE2	5:F:2032:HOH:O	2.22	0.41
1:F:71:GLY:O	1:F:74[B]:ASP:HB2	2.20	0.41
1:F:40:HIS:CE1	1:F:66[B]:MET:SD	3.14	0.41
1:I:48:TRP:HB3	1:K:81:ILE:HD11	2.03	0.41
1:B:46:VAL:HG13	1:B:114:GLN:HB3	2.04	0.40
1:J:47:HIS:CE1	1:J:59:HIS:CE1	3.09	0.40
1:H:57[B]:ILE:CD1	1:K:57[B]:ILE:CD1	2.95	0.40
1:E:152:LYS:HE3	5:E:2104:HOH:O	2.22	0.40
1:H:22:LEU:HD23	1:H:139:GLU:HB3	2.02	0.40
1:J:46:VAL:HG13	1:J:114:GLN:HB3	2.04	0.40
1:E:57:ILE:CG1	1:L:57[B]:ILE:HD11	2.47	0.40
1:F:117[B]:ARG:HE	1:F:117[B]:ARG:HB2	1.71	0.40
1:C:47:HIS:CE1	1:C:59:HIS:CE1	3.10	0.40
1:A:57[A]:ILE:HG23	1:A:58:TRP:CD1	2.57	0.40
1:K:22:LEU:HD22	1:K:141:ASP:HB2	2.04	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	158/165 (96%)	157 (99%)	1 (1%)	0	100	100
1	B	157/165 (95%)	156 (99%)	1 (1%)	0	100	100
1	C	157/165 (95%)	154 (98%)	3 (2%)	0	100	100
1	D	158/165 (96%)	157 (99%)	1 (1%)	0	100	100
1	E	155/165 (94%)	154 (99%)	1 (1%)	0	100	100
1	F	157/165 (95%)	156 (99%)	1 (1%)	0	100	100
1	G	169/165 (102%)	162 (96%)	4 (2%)	3 (2%)	8	1
1	H	154/165 (93%)	153 (99%)	1 (1%)	0	100	100
1	I	154/165 (93%)	150 (97%)	4 (3%)	0	100	100
1	J	158/165 (96%)	156 (99%)	2 (1%)	0	100	100
1	K	156/165 (94%)	155 (99%)	1 (1%)	0	100	100
1	L	155/165 (94%)	154 (99%)	1 (1%)	0	100	100
All	All	1888/1980 (95%)	1864 (99%)	21 (1%)	3 (0%)	47	33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	18	PRO
1	G	19	ARG
1	G	20	PRO

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	136/140 (97%)	135 (99%)	1 (1%)	84	80
1	B	135/140 (96%)	135 (100%)	0	100	100
1	C	136/140 (97%)	130 (96%)	6 (4%)	28	13
1	D	136/140 (97%)	134 (98%)	2 (2%)	65	55
1	E	132/140 (94%)	132 (100%)	0	100	100
1	F	134/140 (96%)	133 (99%)	1 (1%)	84	80
1	G	140/140 (100%)	138 (99%)	2 (1%)	67	58
1	H	131/140 (94%)	128 (98%)	3 (2%)	50	37
1	I	131/140 (94%)	129 (98%)	2 (2%)	65	55
1	J	136/140 (97%)	128 (94%)	8 (6%)	19	7
1	K	133/140 (95%)	133 (100%)	0	100	100
1	L	131/140 (94%)	130 (99%)	1 (1%)	81	77
All	All	1611/1680 (96%)	1585 (98%)	26 (2%)	67	53

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	ILE
1	C	74[A]	ASP
1	C	74[B]	ASP
1	C	75[A]	GLU
1	C	75[B]	GLU
1	C	101	LYS
1	C	160	TRP
1	D	99	GLN
1	D	111	ILE
1	F	69	ILE
1	G	69	ILE
1	G	104	LEU
1	H	69	ILE
1	H	117	ARG
1	H	146	ASP
1	I	44	HIS
1	I	69	ILE
1	J	44	HIS
1	J	57	ILE
1	J	75[A]	GLU
1	J	75[B]	GLU
1	J	111[A]	ILE

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Mol	Chain	Res	Type
1	J	111[B]	ILE
1	J	117[A]	ARG
1	J	117[B]	ARG
1	L	95	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	GLN
1	C	99	GLN
1	F	99	GLN
1	I	99	GLN
1	J	99	GLN
1	J	145	ASN
1	K	130	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 13 are monoatomic - leaving 4 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	EPE	B	200	-	15,15,15	0.92	1 (6%)	18,20,20	3.03	7 (38%)
2	EPE	C	200	-	15,15,15	0.73	1 (6%)	18,20,20	2.36	8 (44%)
2	EPE	A	200	-	15,15,15	1.07	1 (6%)	18,20,20	2.35	6 (33%)
2	EPE	D	200	-	15,15,15	1.07	1 (6%)	18,20,20	2.41	8 (44%)

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	EPE	B	200	-	-	2/9/19/19	0/1/1/1
2	EPE	C	200	-	-	3/9/19/19	0/1/1/1
2	EPE	A	200	-	-	5/9/19/19	0/1/1/1
2	EPE	D	200	-	-	5/9/19/19	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	200	EPE	C10-S	3.68	1.82	1.77
2	D	200	EPE	C10-S	3.33	1.82	1.77
2	B	200	EPE	C10-S	2.72	1.81	1.77
2	C	200	EPE	C10-S	2.24	1.80	1.77

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	200	EPE	O1S-S-C10	8.51	117.17	106.92
2	A	200	EPE	O1S-S-C10	5.41	113.43	106.92
2	D	200	EPE	O3S-S-C10	5.09	114.01	105.77
2	D	200	EPE	C5-N4-C3	4.93	119.92	108.83
2	B	200	EPE	O2S-S-C10	-4.69	101.27	106.92
2	C	200	EPE	O3S-S-C10	4.68	113.33	105.77
2	B	200	EPE	C5-N4-C3	4.63	119.25	108.83
2	A	200	EPE	C5-N4-C3	4.56	119.08	108.83
2	C	200	EPE	O1S-S-C10	4.04	111.78	106.92
2	C	200	EPE	C5-N4-C3	4.04	117.91	108.83
2	B	200	EPE	C7-N4-C3	3.76	120.84	111.23
2	C	200	EPE	O2S-S-C10	-3.64	102.53	106.92
2	A	200	EPE	O2S-S-C10	-3.60	102.58	106.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	200	EPE	C7-N4-C3	3.59	120.42	111.23
2	D	200	EPE	C7-N4-C5	3.48	120.13	111.23
2	A	200	EPE	C7-N4-C5	3.29	119.66	111.23
2	C	200	EPE	C7-N4-C3	3.17	119.33	111.23
2	A	200	EPE	C6-N1-C2	3.06	115.71	108.83
2	D	200	EPE	C5-C6-N1	-2.94	104.60	110.64
2	D	200	EPE	C7-N4-C3	2.93	118.73	111.23
2	B	200	EPE	C6-N1-C2	2.90	115.36	108.83
2	B	200	EPE	C7-N4-C5	2.69	118.11	111.23
2	B	200	EPE	C9-N1-C2	-2.60	104.57	111.23
2	D	200	EPE	C3-C2-N1	2.41	115.59	110.64
2	D	200	EPE	C2-C3-N4	2.41	115.58	110.64
2	C	200	EPE	O3S-S-O1S	-2.36	105.51	111.27
2	D	200	EPE	C6-N1-C2	2.33	114.07	108.83
2	C	200	EPE	C6-N1-C2	2.20	113.79	108.83
2	C	200	EPE	C2-C3-N4	2.02	114.79	110.64

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	200	EPE	C9-C10-S-O1S
2	D	200	EPE	C9-C10-S-O2S
2	C	200	EPE	C10-C9-N1-C2
2	C	200	EPE	C10-C9-N1-C6
2	C	200	EPE	C8-C7-N4-C3
2	D	200	EPE	C9-C10-S-O3S
2	A	200	EPE	C9-C10-S-O3S
2	B	200	EPE	N4-C7-C8-O8
2	A	200	EPE	N4-C7-C8-O8
2	B	200	EPE	C8-C7-N4-C5
2	A	200	EPE	C9-C10-S-O1S
2	A	200	EPE	C9-C10-S-O2S
2	A	200	EPE	C8-C7-N4-C5
2	D	200	EPE	C8-C7-N4-C3
2	D	200	EPE	C8-C7-N4-C5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	200	EPE	1	0
2	D	200	EPE	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	L	4
1	G	2
1	J	2
1	K	2
1	C	2
1	F	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	L	74[A]:ASP	C	75[A]:GLU	N	2.19
1	C	77[A]:SER	C	78[A]:GLU	N	1.95
1	G	77[A]:SER	C	78[A]:GLU	N	1.88
1	K	77[A]:SER	C	78[A]:GLU	N	1.86
1	J	77[A]:SER	C	78[A]:GLU	N	1.82
1	F	77[A]:SER	C	78[A]:GLU	N	1.66
1	L	77[A]:SER	C	78[A]:GLU	N	1.62
1	L	77[B]:SER	C	78[B]:GLU	N	1.12
1	F	77[B]:SER	C	78[B]:GLU	N	1.09
1	J	77[B]:SER	C	78[B]:GLU	N	0.87
1	G	77[B]:SER	C	78[B]:GLU	N	0.81
1	K	77[B]:SER	C	78[B]:GLU	N	0.81
1	C	77[B]:SER	C	78[B]:GLU	N	0.73
1	L	74[B]:ASP	C	75[B]:GLU	N	0.47

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	151/165 (91%)	-0.38	2 (1%) 77 74	16, 20, 29, 39	0
1	B	151/165 (91%)	-0.32	2 (1%) 77 74	16, 21, 29, 43	0
1	C	151/165 (91%)	-0.28	2 (1%) 77 74	18, 22, 32, 43	0
1	D	152/165 (92%)	-0.33	1 (0%) 87 86	16, 20, 29, 42	0
1	E	151/165 (91%)	-0.27	1 (0%) 87 86	19, 23, 32, 42	0
1	F	151/165 (91%)	-0.14	2 (1%) 77 74	20, 24, 35, 43	0
1	G	165/165 (100%)	-0.17	5 (3%) 50 44	17, 23, 35, 49	0
1	H	151/165 (91%)	-0.08	2 (1%) 77 74	19, 26, 37, 50	0
1	I	151/165 (91%)	-0.19	2 (1%) 77 74	17, 22, 31, 39	0
1	J	151/165 (91%)	-0.05	2 (1%) 77 74	20, 25, 34, 42	0
1	K	153/165 (92%)	-0.24	1 (0%) 87 86	17, 22, 31, 41	0
1	L	151/165 (91%)	-0.22	2 (1%) 77 74	20, 24, 31, 37	0
All	All	1829/1980 (92%)	-0.22	24 (1%) 77 74	16, 23, 33, 50	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	22	LEU	5.5
1	D	21	SER	5.3
1	G	18	PRO	5.1
1	A	22	LEU	4.5
1	J	22	LEU	4.3
1	K	20	PRO	4.1
1	C	23	ALA	3.9
1	H	23	ALA	3.8
1	G	19	ARG	3.8
1	I	23	ALA	3.7
1	L	22	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	172	LEU	3.5
1	L	23	ALA	3.4
1	F	23	ALA	3.3
1	B	23	ALA	3.1
1	G	20	PRO	2.9
1	A	23	ALA	2.8
1	H	22	LEU	2.7
1	G	17	SER	2.7
1	E	23	ALA	2.6
1	F	22	LEU	2.5
1	G	21	SER	2.2
1	C	22	LEU	2.2
1	J	23	ALA	2.2

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. 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(\AA^2)	Q<0.9
3	FE	L	2000	1/1	0.94	0.06	31,31,31,31	1
2	EPE	A	200	15/15	0.94	0.16	26,39,47,50	0
2	EPE	D	200	15/15	0.95	0.15	25,32,37,37	0
3	FE	G	2000	1/1	0.95	0.06	26,26,26,26	1
2	EPE	C	200	15/15	0.96	0.13	28,35,43,45	0
3	FE	C	2000	1/1	0.97	0.06	26,26,26,26	1
3	FE	F	2000	1/1	0.97	0.07	27,27,27,27	1
3	FE	B	2000	1/1	0.97	0.05	24,24,24,24	1
2	EPE	B	200	15/15	0.97	0.11	22,29,38,40	0
3	FE	A	2000	1/1	0.97	0.04	25,25,25,25	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	FE	H	2000	1/1	0.98	0.04	26,26,26,26	1
3	FE	E	2000	1/1	0.98	0.05	26,26,26,26	1
3	FE	K	2000	1/1	0.98	0.03	24,24,24,24	1
3	FE	D	2000	1/1	0.98	0.03	26,26,26,26	1
3	FE	I	2000	1/1	0.98	0.04	26,26,26,26	1
4	CA	K	1501	1/1	0.98	0.06	25,25,25,25	0
3	FE	J	2000	1/1	0.99	0.04	23,23,23,23	1

6.5 Other polymers [i](#)

There are no such residues in this entry.