Full wwPDB X-ray Structure Validation Report (i)

Sep 24, 2023 – 03:05 AM EDT

PDB ID	:	5KSR
Title	:	Stationary phase survival protein E (SurE) from Xylella fastidiosa - XFSurE-TB (Tetramer Bigger).
Authors	:	Machado, A.T.P.; Fonseca, E.M.B.; Dos Reis, M.A.; Saraiva, A.M.; Dos Santos, C.A.; De Toledo, M.A.; Polikarpov, I.; De Souza, A.P.; De Aparicio, R.; Iulek, J.
Deposited on	:	2016-07-09
Resolution	:	1.96 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 (i) 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 (1)) were used in the production of this report:

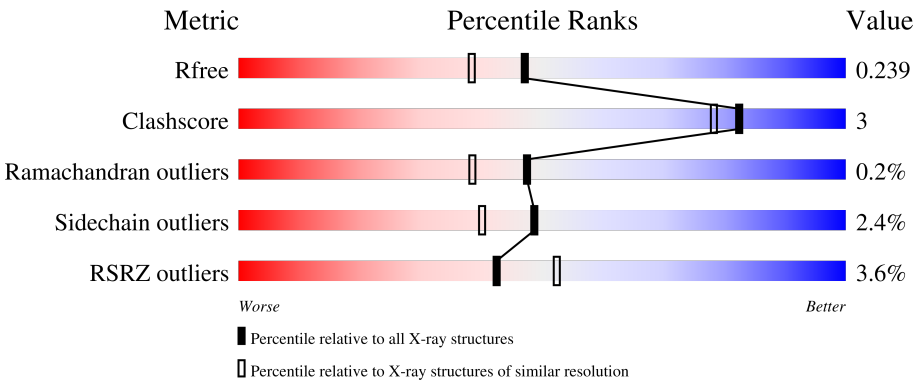
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 1.96 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}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 <=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	270	3% 90% • • 6%
1	B	270	2% 89% 5% • 6%
1	C	270	2% 88% 7% • •
1	D	270	6% 90% 5% •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 8554 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 5'-nucleotidase SurE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	1956	1219	351	378	8	10	3	0
1	B	254	1943	1212	346	377	8	24	3	0
1	C	258	1999	1253	354	383	9	43	6	0
1	D	259	1981	1233	354	386	8	41	3	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	263	LEU	-	expression tag	UNP Q9PF20
A	264	GLU	-	expression tag	UNP Q9PF20
A	265	HIS	-	expression tag	UNP Q9PF20
A	266	HIS	-	expression tag	UNP Q9PF20
A	267	HIS	-	expression tag	UNP Q9PF20
A	268	HIS	-	expression tag	UNP Q9PF20
A	269	HIS	-	expression tag	UNP Q9PF20
A	270	HIS	-	expression tag	UNP Q9PF20
B	263	LEU	-	expression tag	UNP Q9PF20
B	264	GLU	-	expression tag	UNP Q9PF20
B	265	HIS	-	expression tag	UNP Q9PF20
B	266	HIS	-	expression tag	UNP Q9PF20
B	267	HIS	-	expression tag	UNP Q9PF20
B	268	HIS	-	expression tag	UNP Q9PF20
B	269	HIS	-	expression tag	UNP Q9PF20
B	270	HIS	-	expression tag	UNP Q9PF20
C	263	LEU	-	expression tag	UNP Q9PF20
C	264	GLU	-	expression tag	UNP Q9PF20
C	265	HIS	-	expression tag	UNP Q9PF20
C	266	HIS	-	expression tag	UNP Q9PF20
C	267	HIS	-	expression tag	UNP Q9PF20

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Chain	Residue	Modelled	Actual	Comment	Reference
C	268	HIS	-	expression tag	UNP Q9PF20
C	269	HIS	-	expression tag	UNP Q9PF20
C	270	HIS	-	expression tag	UNP Q9PF20
D	263	LEU	-	expression tag	UNP Q9PF20
D	264	GLU	-	expression tag	UNP Q9PF20
D	265	HIS	-	expression tag	UNP Q9PF20
D	266	HIS	-	expression tag	UNP Q9PF20
D	267	HIS	-	expression tag	UNP Q9PF20
D	268	HIS	-	expression tag	UNP Q9PF20
D	269	HIS	-	expression tag	UNP Q9PF20
D	270	HIS	-	expression tag	UNP Q9PF20

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mn 1 1	0	0
2	B	1	Total Mn 1 1	0	0
2	C	1	Total Mn 1 1	0	0
2	D	1	Total Mn 1 1	0	0

- Molecule 3 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total I 2 2	0	0
3	B	1	Total I 1 1	0	0
3	C	2	Total I 2 2	0	0
3	D	1	Total I 1 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
4	B	1	1	1	0	0
			Total	Cl		
4	C	1	1	1	0	0
			Total	Cl		
4	D	1	1	1	0	0

- Molecule 5 is water.

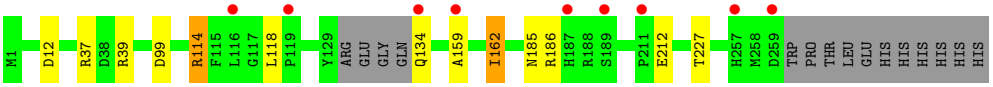
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	171	171	171	0	0
			Total	O		
5	B	165	165	165	0	0
			Total	O		
5	C	168	168	168	0	0
			Total	O		
5	D	157	157	157	0	0

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.

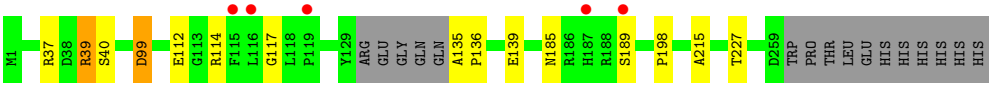
- Molecule 1: 5'-nucleotidase SurE

Chain A:



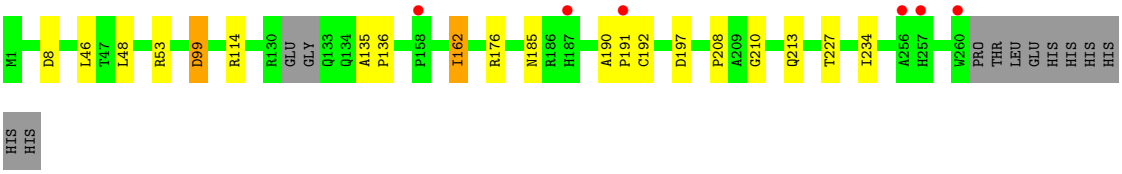
- Molecule 1: 5'-nucleotidase SurE

Chain B:



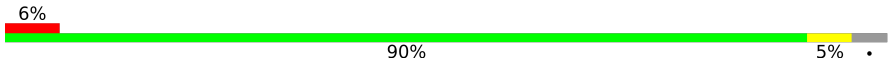
- Molecule 1: 5'-nucleotidase SurE

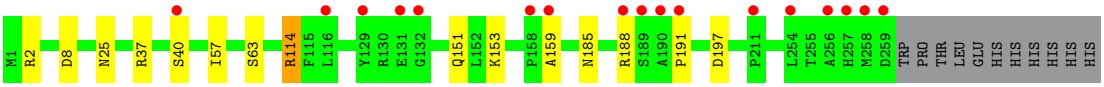
Chain C:



- Molecule 1: 5'-nucleotidase SurE

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	172.34Å 84.17Å 87.31Å 90.00° 96.62° 90.00°	Depositor
Resolution (Å)	43.25 – 1.96 75.53 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.25-1.96) 99.5 (75.53-1.96)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.42 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.196 , 0.238 0.198 , 0.239	Depositor DCC
R_{free} test set	4432 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	29.8	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 79.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8554	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.41% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: IOD, CL, MN

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	0.89	0/2004	0.74	1/2742 (0.0%)
1	B	0.86	0/1991	0.74	0/2726
1	C	0.81	0/2058	0.71	1/2817 (0.0%)
1	D	0.76	0/2024	0.71	0/2771
All	All	0.83	0/8077	0.72	2/11056 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	46	LEU	CA-CB-CG	5.06	126.95	115.30
1	A	118	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1956	0	1935	10	0
1	B	1943	0	1921	8	0
1	C	1999	0	1988	11	0
1	D	1981	0	1950	9	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	1	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	171	0	0	5	0
5	B	165	0	0	6	0
5	C	168	0	0	4	1
5	D	157	0	0	6	1
All	All	8554	0	7794	39	1

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 (39) 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:B:99[B]:ASP:OD2	1:B:185:ASN:ND2	2.07	0.88
1:A:99[A]:ASP:OD2	1:A:185:ASN:ND2	2.12	0.83
4:B:503:CL:CL	5:B:692:HOH:O	2.41	0.74
1:A:114:ARG:NH1	1:A:159:ALA:HB1	2.09	0.68
1:B:139:GLU:OE2	5:B:601:HOH:O	2.13	0.67
1:D:185:ASN:ND2	5:D:605:HOH:O	2.28	0.67
1:A:39:ARG:NH1	5:A:604:HOH:O	2.28	0.66
1:A:134:GLN:N	5:A:606:HOH:O	2.31	0.62
1:A:114:ARG:HH12	1:A:159:ALA:HB1	1.63	0.62
1:B:215:ALA:O	5:B:602:HOH:O	2.15	0.61
1:A:186[A]:ARG:NH1	5:A:610:HOH:O	2.35	0.60
1:D:188:ARG:O	5:D:601:HOH:O	2.16	0.60
1:A:12:ASP:OD2	5:A:601:HOH:O	2.16	0.59
1:B:39:ARG:NH2	5:B:606:HOH:O	2.36	0.58
1:D:153:LYS:NZ	5:D:612:HOH:O	2.37	0.56
1:B:114:ARG:NH2	1:B:117:GLY:HA2	2.21	0.56
1:A:114:ARG:HG3	1:A:162:ILE:HG22	1.88	0.56
1:C:210:GLY:O	1:C:213:GLN:NE2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99[B]:ASP:OD2	1:C:185[B]:ASN:ND2	2.41	0.54
1:C:176:ARG:NH2	5:C:611:HOH:O	2.40	0.54
1:D:40:SER:HB3	5:D:675:HOH:O	2.07	0.54
1:D:25:ASN:ND2	5:D:614:HOH:O	2.44	0.50
1:D:114[A]:ARG:NH1	1:D:159:ALA:HB1	2.28	0.49
1:D:57[B]:ILE:HD11	1:D:63:SER:HB2	1.94	0.48
1:C:48:LEU:O	5:C:601:HOH:O	2.20	0.47
1:B:40:SER:HB3	5:B:604:HOH:O	2.14	0.46
1:A:212:GLU:H	1:A:212:GLU:CD	2.20	0.45
1:D:2:ARG:NH1	5:D:607:HOH:O	2.29	0.45
1:C:208:PRO:O	5:C:602:HOH:O	2.21	0.45
1:B:112:GLU:OE2	5:B:603:HOH:O	2.21	0.45
1:A:39:ARG:NH2	5:A:611:HOH:O	2.37	0.44
1:C:53:ARG:NE	5:C:609:HOH:O	2.38	0.44
1:B:135:ALA:HA	1:B:136:PRO:HD3	1.86	0.44
1:D:114[A]:ARG:CZ	1:D:159:ALA:HB1	2.50	0.42
1:C:162[B]:ILE:HD11	1:C:234:ILE:HD11	2.02	0.41
1:C:114:ARG:HB2	1:C:162[A]:ILE:HD11	2.01	0.41
1:C:190:ALA:HA	1:C:191:PRO:HD3	1.82	0.41
1:C:135:ALA:HA	1:C:136:PRO:HD3	1.82	0.40

All (1) 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)
5:C:759:HOH:O	5:D:726:HOH:O[2_555]	2.19	0.01

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	254/270 (94%)	247 (97%)	7 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	253/270 (94%)	244 (96%)	8 (3%)	1 (0%)	34	22
1	C	260/270 (96%)	251 (96%)	9 (4%)	0	100	100
1	D	259/270 (96%)	249 (96%)	9 (4%)	1 (0%)	34	22
All	All	1026/1080 (95%)	991 (97%)	33 (3%)	2 (0%)	47	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	191	PRO
1	B	198	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	213/224 (95%)	209 (98%)	4 (2%)	57	50
1	B	212/224 (95%)	206 (97%)	6 (3%)	43	33
1	C	219/224 (98%)	211 (96%)	8 (4%)	34	22
1	D	215/224 (96%)	210 (98%)	5 (2%)	50	42
All	All	859/896 (96%)	836 (97%)	23 (3%)	49	34

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	114	ARG
1	A	162	ILE
1	A	227	THR
1	B	37	ARG
1	B	39	ARG
1	B	99[A]	ASP
1	B	99[B]	ASP
1	B	189	SER

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Mol	Chain	Res	Type
1	B	227	THR
1	C	8	ASP
1	C	99[A]	ASP
1	C	99[B]	ASP
1	C	162[A]	ILE
1	C	162[B]	ILE
1	C	192	CYS
1	C	197	ASP
1	C	227	THR
1	D	8	ASP
1	D	37	ARG
1	D	114[A]	ARG
1	D	151	GLN
1	D	197	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	ASN
1	D	173	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 14 ligands modelled in this entry, 14 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 (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	255/270 (94%)	0.11	9 (3%)	44	53	15, 29, 57, 90	3 (1%)
1	B	254/270 (94%)	0.00	5 (1%)	65	73	15, 32, 57, 78	7 (2%)
1	C	258/270 (95%)	0.05	6 (2%)	60	69	17, 31, 64, 90	10 (3%)
1	D	259/270 (95%)	0.20	17 (6%)	18	26	16, 34, 71, 98	11 (4%)
All	All	1026/1080 (95%)	0.09	37 (3%)	42	52	15, 32, 62, 98	31 (3%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	158	PRO	6.1
1	D	257	HIS	5.7
1	B	116	LEU	5.3
1	C	260	TRP	4.6
1	D	254	LEU	4.6
1	D	131	GLU	4.3
1	A	134	GLN	4.1
1	A	116	LEU	3.9
1	A	259	ASP	3.8
1	D	191	PRO	3.7
1	D	132	GLY	3.6
1	D	259	ASP	3.5
1	D	40	SER	3.3
1	D	129	TYR	3.1
1	C	256	ALA	3.1
1	C	187	HIS	3.0
1	D	258	MET	3.0
1	D	159	ALA	2.9
1	D	116	LEU	2.9
1	A	159	ALA	2.8
1	D	256	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	190	ALA	2.7
1	B	115	PHE	2.7
1	C	257	HIS	2.6
1	A	189	SER	2.6
1	D	189	SER	2.5
1	D	188	ARG	2.5
1	D	158	PRO	2.4
1	A	187	HIS	2.4
1	A	257	HIS	2.4
1	C	191	PRO	2.3
1	B	189	SER	2.2
1	A	119	PRO	2.1
1	A	211	PRO	2.1
1	B	119	PRO	2.1
1	D	211	PRO	2.0
1	B	187	HIS	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	CL	C	504	1/1	0.92	0.09	37,37,37,37	0
2	MN	B	501	1/1	0.93	0.05	43,43,43,43	0
4	CL	B	503	1/1	0.94	0.08	31,31,31,31	0
4	CL	A	504	1/1	0.94	0.14	31,31,31,31	0
4	CL	D	503	1/1	0.94	0.08	36,36,36,36	0
2	MN	D	501	1/1	0.98	0.08	37,37,37,37	0
3	IOD	A	503	1/1	0.98	0.11	39,39,39,39	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MN	A	501	1/1	0.98	0.04	41,41,41,41	0
2	MN	C	501	1/1	0.99	0.05	35,35,35,35	0
3	IOD	B	502	1/1	0.99	0.09	40,40,40,40	1
3	IOD	C	503	1/1	0.99	0.09	32,32,32,32	1
3	IOD	D	502	1/1	0.99	0.11	31,31,31,31	1
3	IOD	A	502	1/1	1.00	0.11	20,20,20,20	0
3	IOD	C	502	1/1	1.00	0.11	25,25,25,25	1

6.5 Other polymers (i)

There are no such residues in this entry.