



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:55 PM UTC

PDB ID : 4U65 / pdb_00004u65
Title : Structure of the periplasmic output domain of the Legionella pneumophila LapD ortholog CdgS9 in complex with Pseudomonas fluorescens LapG
Authors : Chatterjee, D.; Cooley, R.B.; Boyd, C.D.; Mehl, R.A.; O'Toole, G.A.; Sondermann, H.S.
Deposited on : 2014-07-27
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

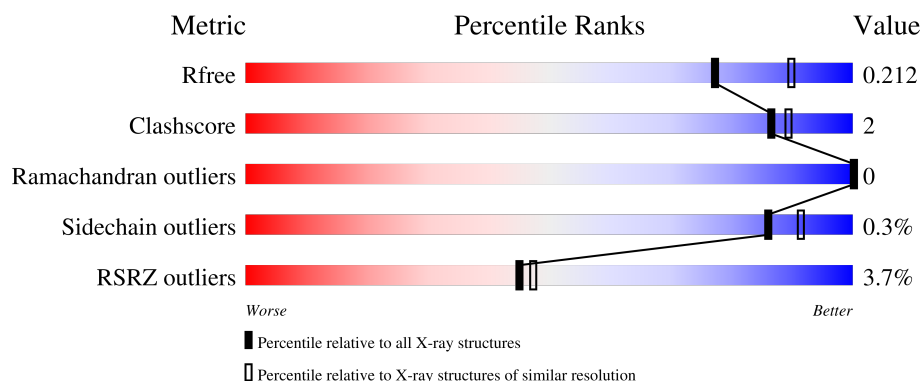
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



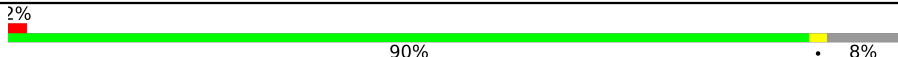
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	131	5% 95% 5%
1	B	131	5% 89% 11% .
1	C	131	% 96% .
1	D	131	9% 95% 5% .
2	E	197	2% 88% 8% ..

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Mol	Chain	Length	Quality of chain
2	F	197	 A horizontal bar chart showing the quality of chain F. The bar is green, indicating a high quality score. The bar is labeled with '2%' at the start, '90%' in the middle, and '8%' at the end. The bar is green for the first 90% and yellow for the last 8%.

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7826 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 Two component histidine kinase, GGDEF domain protein/EAL domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	2	0
			1056	678	174	198	6			
1	B	131	Total	C	N	O	S	0	2	0
			1054	676	172	200	6			
1	C	131	Total	C	N	O	S	0	0	0
			1042	670	170	196	6			
1	D	131	Total	C	N	O	S	0	2	0
			1054	676	172	200	6			

- Molecule 2 is a protein called Putative cystine protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	181	Total	C	N	O	S	0	2	0
			1474	948	248	272	6			
2	F	181	Total	C	N	O	S	0	0	0
			1465	945	246	268	6			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total	Ca	0	0
			2	2		
3	F	2	Total	Ca	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	113	Total	O	0	0
			113	113		

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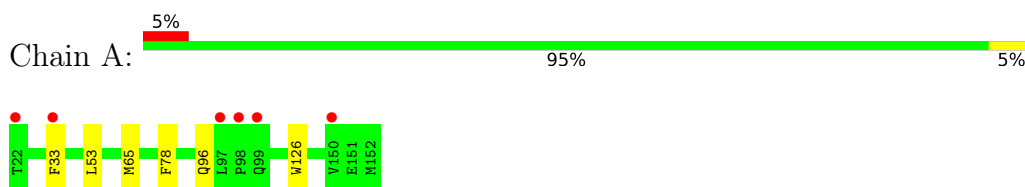
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	70	Total 70	O 70	0	0
4	C	120	Total 120	O 120	0	0
4	D	68	Total 68	O 68	0	0
4	E	163	Total 163	O 163	0	0
4	F	143	Total 143	O 143	0	0

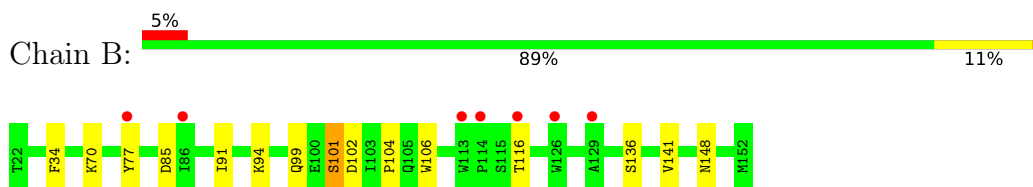
3 Residue-property plots [i](#)

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.

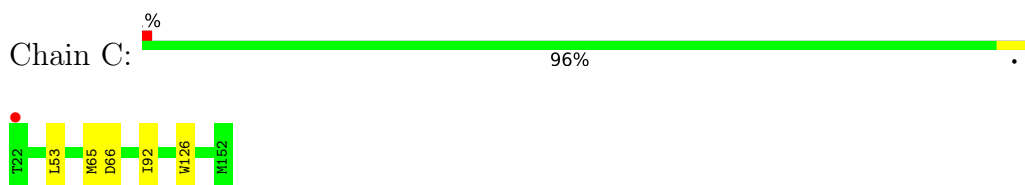
- Molecule 1: Two component histidine kinase, GGDEF domain protein/EAL domain protein



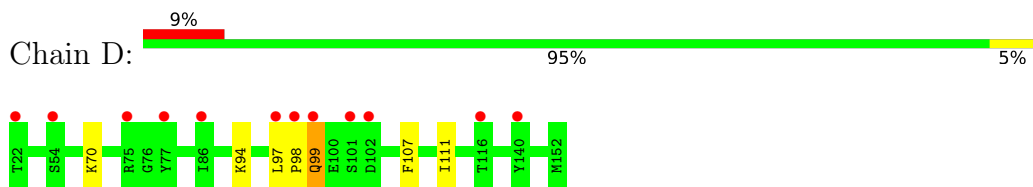
- Molecule 1: Two component histidine kinase, GGDEF domain protein/EAL domain protein



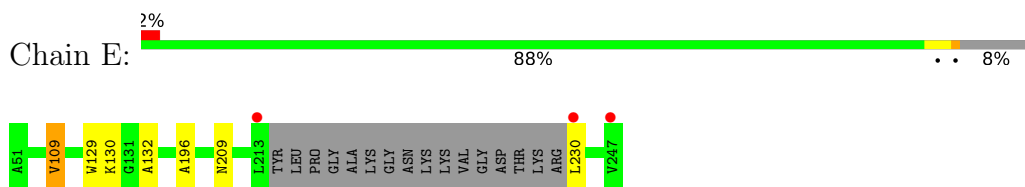
- Molecule 1: Two component histidine kinase, GGDEF domain protein/EAL domain protein



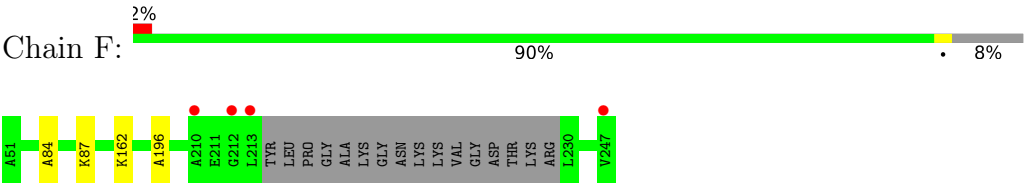
- Molecule 1: Two component histidine kinase, GGDEF domain protein/EAL domain protein



- Molecule 2: Putative cystine protease



- Molecule 2: Putative cystine protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.85Å 73.67Å 88.22Å 90.00° 92.88° 90.00°	Depositor
Resolution (Å)	44.06 – 2.10 44.06 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.6 (44.06-2.10) 93.1 (44.06-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.170 , 0.211 0.172 , 0.212	Depositor DCC
R_{free} test set	2001 reflections (3.50%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l 0.016 for -k,-h,-l 0.116 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7826	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.16	0/1082	0.40	0/1467
1	B	0.15	0/1077	0.49	2/1461 (0.1%)
1	C	0.15	0/1065	0.41	0/1445
1	D	0.18	0/1077	0.42	0/1461
2	E	0.11	0/1510	0.31	0/2048
2	F	0.11	0/1501	0.31	0/2036
All	All	0.14	0/7312	0.39	2/9918 (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	D	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	101	SER	CA-C-N	6.73	133.82	121.70
1	B	101	SER	C-N-CA	6.73	133.82	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	98	PRO	Peptide
1	D	99	GLN	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	1056	0	1057	6	0
1	B	1054	0	1048	12	0
1	C	1042	0	1040	4	0
1	D	1054	0	1048	6	0
2	E	1474	0	1448	5	0
2	F	1465	0	1449	3	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	113	0	0	0	0
4	B	70	0	0	0	0
4	C	120	0	0	0	0
4	D	68	0	0	0	0
4	E	163	0	0	0	0
4	F	143	0	0	1	0
All	All	7826	0	7090	30	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 2.

All (30) 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:70:LYS:HG3	1:B:94:LYS:HE3	1.73	0.71
1:A:33:PHE:HZ	1:B:148:ASN:HD22	1.41	0.67
1:D:97:LEU:HB3	1:D:99:GLN:OE1	1.99	0.63
1:B:116:THR:HG21	1:B:136:SER:HB2	1.85	0.59
1:D:70:LYS:HG3	1:D:94:LYS:HE2	1.85	0.58
1:B:85:ASP:HB3	1:B:91:ILE:HD11	1.85	0.57
1:A:33:PHE:HZ	1:B:148:ASN:ND2	2.03	0.56
1:A:126:TRP:CD2	2:E:196:ALA:HB1	2.42	0.54
1:C:126:TRP:CD2	2:F:196:ALA:HB1	2.42	0.54
1:B:101:SER:HA	1:B:102:ASP:CB	2.39	0.53
1:C:53:LEU:HD22	1:C:65:MET:HG3	1.91	0.52
1:B:101:SER:HA	1:B:102:ASP:HB3	1.92	0.51
1:B:104:PRO:HB2	1:B:106:TRP:CD1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:109:VAL:HG13	2:E:132:ALA:HB1	1.93	0.50
1:C:65:MET:HG2	1:C:92:ILE:HD12	1.94	0.49
1:B:77:TYR:HE2	2:E:230:LEU:HB3	1.78	0.49
2:F:162:LYS:NZ	4:F:402:HOH:O	2.47	0.48
1:D:99:GLN:HG2	1:D:99:GLN:O	2.14	0.47
1:D:99:GLN:H	1:D:99:GLN:CD	2.23	0.46
1:B:99:GLN:N	1:B:99:GLN:OE1	2.49	0.45
1:C:66:ASP:HA	1:C:92:ILE:HD13	1.99	0.44
1:B:34:PHE:CD1	1:B:141:VAL:HG22	2.53	0.44
1:D:99:GLN:OE1	1:D:99:GLN:N	2.43	0.43
1:A:53:LEU:HD22	1:A:65:MET:HG3	1.99	0.43
1:D:107:PHE:CE2	1:D:111:ILE:HD13	2.54	0.43
1:A:78:PHE:O	1:A:96:GLN:NE2	2.44	0.41
1:A:33:PHE:CZ	1:B:148:ASN:ND2	2.83	0.41
2:F:84:ALA:HA	2:F:87:LYS:HE3	2.02	0.41
2:E:129:TRP:CZ3	2:E:130:LYS:HE2	2.56	0.40
2:E:230:LEU:HD12	2:E:230:LEU:HA	1.92	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	131/131 (100%)	129 (98%)	2 (2%)	0	100	100
1	B	131/131 (100%)	128 (98%)	3 (2%)	0	100	100
1	C	129/131 (98%)	128 (99%)	1 (1%)	0	100	100
1	D	131/131 (100%)	129 (98%)	2 (2%)	0	100	100
2	E	179/197 (91%)	176 (98%)	3 (2%)	0	100	100
2	F	177/197 (90%)	172 (97%)	5 (3%)	0	100	100
All	All	878/918 (96%)	862 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	120/118 (102%)	120 (100%)	0	100	100
1	B	120/118 (102%)	120 (100%)	0	100	100
1	C	118/118 (100%)	118 (100%)	0	100	100
1	D	120/118 (102%)	120 (100%)	0	100	100
2	E	155/166 (93%)	153 (99%)	2 (1%)	61	69
2	F	154/166 (93%)	154 (100%)	0	100	100
All	All	787/804 (98%)	785 (100%)	2 (0%)	86	91

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	109	VAL
2	E	209	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	28	ASN
1	B	84	GLN
1	B	105	GLN
1	D	148	ASN
2	E	86	GLN
2	E	209	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	131/131 (100%)	-0.01	6 (4%)	37	39	13, 28, 57, 87	2 (1%)
1	B	131/131 (100%)	0.44	7 (5%)	32	34	13, 40, 70, 86	2 (1%)
1	C	131/131 (100%)	-0.12	1 (0%)	82	84	18, 28, 49, 70	0
1	D	131/131 (100%)	0.49	12 (9%)	14	15	12, 40, 75, 98	2 (1%)
2	E	181/197 (91%)	-0.33	3 (1%)	69	71	11, 26, 51, 74	2 (1%)
2	F	181/197 (91%)	-0.28	4 (2%)	62	65	16, 26, 49, 78	0
All	All	886/918 (96%)	-0.01	33 (3%)	45	47	11, 29, 62, 98	8 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	77	TYR	5.0
1	B	116	THR	4.7
1	D	116	THR	4.6
2	F	247	VAL	4.2
1	A	33	PHE	3.9
2	F	212	GLY	3.7
1	D	140	TYR	3.7
2	F	213	LEU	3.7
2	E	230	LEU	3.3
1	B	113	TRP	3.3
1	D	99	GLN	3.2
1	A	98	PRO	3.2
1	D	97	LEU	3.0
2	F	210	ALA	3.0
1	D	101	SER	2.8
1	D	77	TYR	2.8
1	D	75	ARG	2.8
1	D	98	PRO	2.8
1	B	86	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	97	LEU	2.6
1	D	102	ASP	2.6
1	B	129	ALA	2.6
1	D	86	ILE	2.5
2	E	247	VAL	2.5
2	E	213	LEU	2.3
1	D	22	THR	2.3
1	A	22	THR	2.3
1	B	114	PRO	2.1
1	A	150	VAL	2.1
1	A	99	GLN	2.1
1	D	54	SER	2.1
1	C	22	THR	2.1
1	B	126	TRP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	F	302	1/1	0.98	0.05	35,35,35,35	0
3	CA	E	302	1/1	0.99	0.03	36,36,36,36	0
3	CA	F	301	1/1	0.99	0.06	36,36,36,36	0
3	CA	E	301	1/1	0.99	0.05	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.