



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 15, 2018 – 08:16 PM EST

PDB ID : 5UKA
Title : Salmonella typhimurium AhpC E49Q mutant
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Deposited on : 2017-01-20
Resolution : 1.90 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

<http://wwpdb.org/validation/2016/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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030736

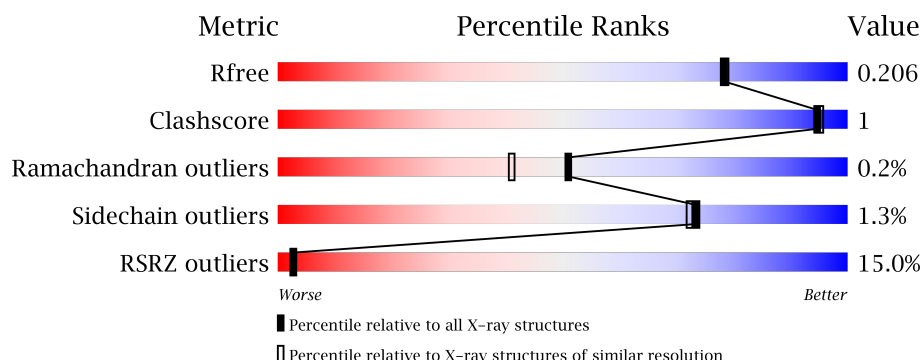
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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. 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	186	<div> <div>12%</div> <div> <div></div> <div>84%</div> <div>11%</div> </div> </div>
1	B	186	<div> <div>9%</div> <div> <div></div> <div>83%</div> <div>12%</div> </div> </div>
1	C	186	<div> <div>13%</div> <div> <div></div> <div>82%</div> <div>12%</div> </div> </div>
1	D	186	<div> <div>10%</div> <div> <div></div> <div>85%</div> <div>12%</div> </div> </div>
1	E	186	<div> <div>23%</div> <div> <div></div> <div>84%</div> <div>13%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	K	A	202	-	-	-	X
3	K	D	202	-	-	-	X
4	NA	B	203	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13493 atoms, of which 6347 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 Alkyl hydroperoxide reductase subunit C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	165	Total	C	H	N	O	S	0	6	0
			2603	840	1279	222	258	4			
1	B	164	Total	C	H	N	O	S	0	13	0
			2693	873	1322	227	267	4			
1	C	163	Total	C	H	N	O	S	0	7	0
			2582	836	1262	219	262	3			
1	D	164	Total	C	H	N	O	S	0	4	0
			2569	831	1260	219	256	3			
1	E	162	Total	C	H	N	O	S	0	0	0
			2495	807	1224	214	247	3			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	49	GLN	GLU	engineered mutation	UNP P0A251
B	49	GLN	GLU	engineered mutation	UNP P0A251
C	49	GLN	GLU	engineered mutation	UNP P0A251
D	49	GLN	GLU	engineered mutation	UNP P0A251
E	49	GLN	GLU	engineered mutation	UNP P0A251

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total K 1 1	0	0
3	A	1	Total K 1 1	0	0
3	D	1	Total K 1 1	0	0

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Na 1 1	0	0

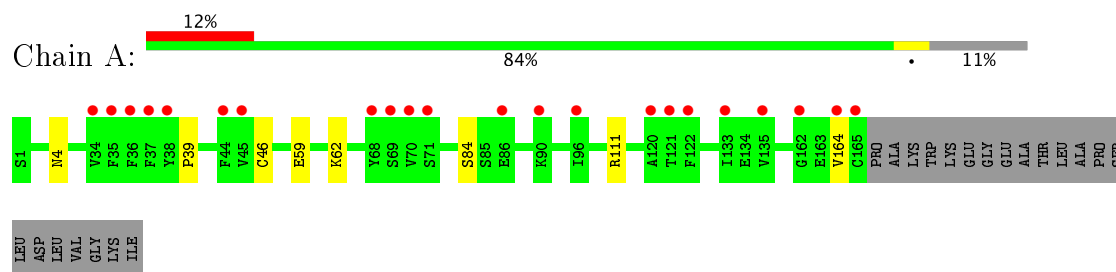
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	127	Total O 128 128	0	1
5	B	137	Total O 139 139	0	2
5	C	125	Total O 127 127	0	2
5	D	118	Total O 118 118	0	0
5	E	30	Total O 30 30	0	0

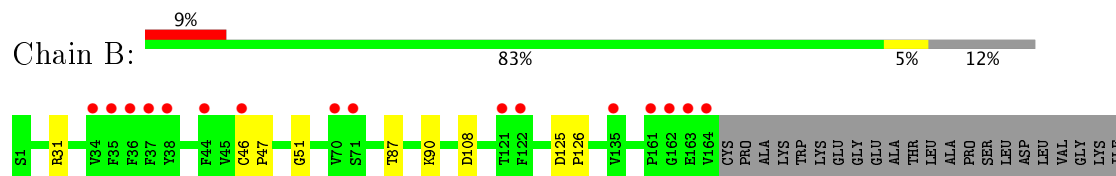
3 Residue-property plots

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.

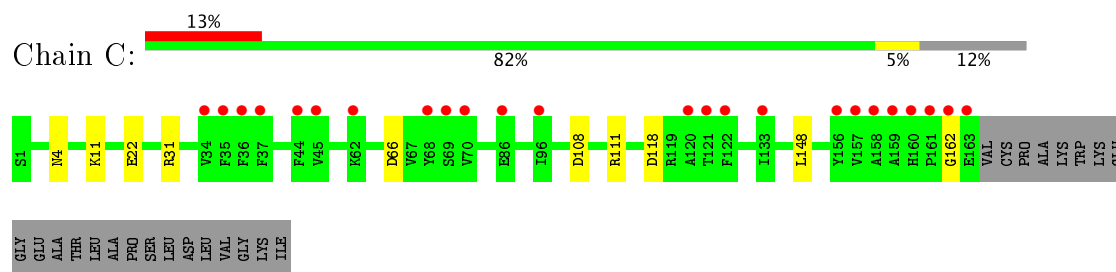
- Molecule 1: Alkyl hydroperoxide reductase subunit C



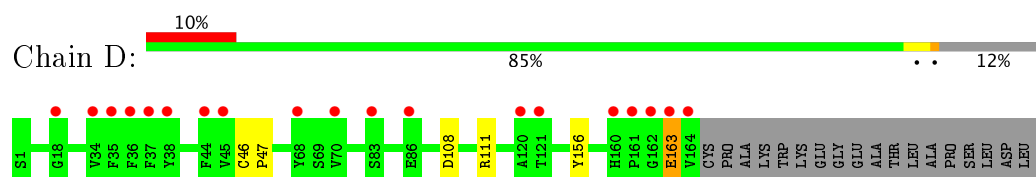
- Molecule 1: Alkyl hydroperoxide reductase subunit C



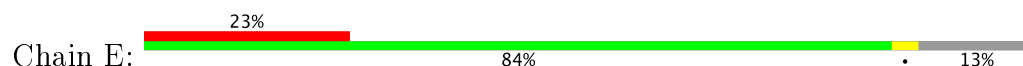
- Molecule 1: Alkyl hydroperoxide reductase subunit C

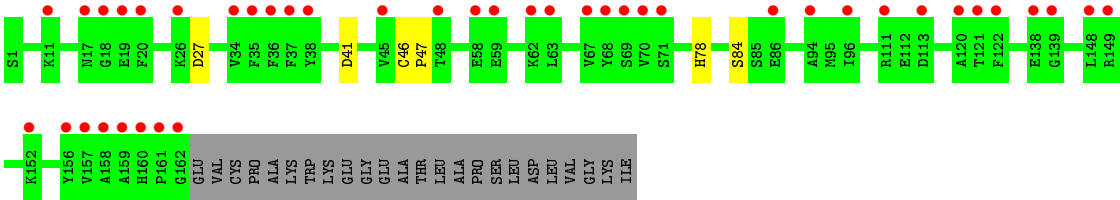


- Molecule 1: Alkyl hydroperoxide reductase subunit C



- Molecule 1: Alkyl hydroperoxide reductase subunit C





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	126.90Å 171.95Å 135.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.26 – 1.90 27.97 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (27.26-1.90) 99.9 (27.97-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 1.91Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.175 , 0.209 0.172 , 0.206	Depositor DCC
R_{free} test set	5792 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 60.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13493	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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: NA, K, CL

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	0.68	0/1377	0.67	0/1866
1	B	0.65	0/1428	0.69	0/1938
1	C	0.59	0/1372	0.66	0/1860
1	D	0.54	0/1354	0.61	0/1836
1	E	0.40	0/1300	0.56	0/1762
All	All	0.58	0/6831	0.64	0/9262

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	1324	1279	1254	3	0
1	B	1371	1322	1298	7	0
1	C	1320	1262	1238	4	0
1	D	1309	1260	1247	3	0
1	E	1271	1224	1228	2	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
4	B	1	0	0	0	0
5	A	128	0	0	1	0
5	B	139	0	0	3	0
5	C	127	0	0	2	0
5	D	118	0	0	1	0
5	E	30	0	0	0	0
All	All	7146	6347	6265	19	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 1.

All (19) 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:46[B]:CYS:SG	5:B:327:HOH:O	2.16	1.01
1:D:108[A]:ASP:OD2	5:D:301:HOH:O	1.91	0.88
1:C:108[A]:ASP:OD2	5:C:401:HOH:O	2.03	0.75
1:A:4[B]:ASN:ND2	5:A:301:HOH:O	2.23	0.71
1:B:31:ARG:NH2	5:B:307:HOH:O	2.42	0.52
1:C:11:LYS:NZ	5:C:409:HOH:O	2.44	0.49
1:C:31:ARG:HD3	1:C:66:ASP:OD2	2.13	0.48
1:A:59:GLU:O	1:A:62[B]:LYS:HG2	2.14	0.47
1:B:46[B]:CYS:N	1:B:47[B]:PRO:HD2	2.30	0.47
1:C:111:ARG:HG2	1:C:118:ASP:OD1	2.16	0.46
1:E:41:ASP:OD2	1:E:78:HIS:ND1	2.46	0.43
1:B:47[B]:PRO:O	1:B:51:GLY:N	2.51	0.42
1:A:39:PRO:HD2	1:A:46:CYS:SG	2.60	0.41
1:B:108[A]:ASP:OD2	5:B:301:HOH:O	2.22	0.41
1:D:46:CYS:N	1:D:47:PRO:CD	2.83	0.41
1:B:87:THR:O	1:B:90:LYS:HG3	2.21	0.41
1:D:156:TYR:OH	1:D:163:GLU:OE1	2.37	0.41
1:E:46:CYS:N	1:E:47:PRO:CD	2.84	0.40
1:B:125:ASP:HB2	1:B:126:PRO:CD	2.51	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	169/186 (91%)	168 (99%)	1 (1%)	0	100	100
1	B	175/186 (94%)	173 (99%)	2 (1%)	0	100	100
1	C	168/186 (90%)	165 (98%)	2 (1%)	1 (1%)	28	16
1	D	166/186 (89%)	163 (98%)	2 (1%)	1 (1%)	28	16
1	E	160/186 (86%)	156 (98%)	4 (2%)	0	100	100
All	All	838/930 (90%)	825 (98%)	11 (1%)	2 (0%)	51	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	163	GLU
1	C	162	GLY

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	144/154 (94%)	141 (98%)	3 (2%)	59	53
1	B	150/154 (97%)	150 (100%)	0	100	100
1	C	143/154 (93%)	139 (97%)	4 (3%)	49	40
1	D	141/154 (92%)	140 (99%)	1 (1%)	87	87
1	E	134/154 (87%)	132 (98%)	2 (2%)	70	67
All	All	712/770 (92%)	702 (99%)	10 (1%)	73	69

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	SER
1	A	111	ARG
1	A	164	VAL
1	C	4	ASN
1	C	22[A]	GLU
1	C	22[B]	GLU
1	C	148	LEU
1	D	111	ARG
1	E	27	ASP
1	E	84	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	C	56	HIS
1	C	160	HIS
1	D	49	GLN
1	D	160	HIS

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 9 ligands modelled in this entry, 9 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	165/186 (88%)	0.40	22 (13%) 4 4	26, 42, 77, 136	0
1	B	164/186 (88%)	0.32	16 (9%) 8 9	27, 40, 70, 145	0
1	C	163/186 (87%)	0.38	24 (14%) 3 3	29, 44, 74, 137	0
1	D	164/186 (88%)	0.45	19 (11%) 5 6	36, 48, 83, 139	0
1	E	162/186 (87%)	1.26	42 (25%) 1 1	47, 74, 110, 122	0
All	All	818/930 (87%)	0.56	123 (15%) 3 3	26, 49, 96, 145	0

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	VAL	9.9
1	D	164	VAL	9.7
1	E	159	ALA	7.7
1	E	156	TYR	6.8
1	E	45	VAL	6.4
1	E	36	PHE	6.1
1	C	162	GLY	5.9
1	E	70	VAL	5.9
1	B	162	GLY	5.9
1	D	163	GLU	5.9
1	E	161	PRO	5.3
1	D	162	GLY	5.2
1	E	35	PHE	5.2
1	E	160	HIS	5.0
1	E	157	VAL	4.8
1	E	34	VAL	4.8
1	B	163	GLU	4.7
1	E	37	PHE	4.5
1	A	164	VAL	4.4
1	B	46[A]	CYS	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	37	PHE	4.3
1	D	44	PHE	4.3
1	C	44	PHE	4.3
1	E	162	GLY	4.3
1	C	45	VAL	4.2
1	A	70	VAL	4.2
1	C	161	PRO	4.1
1	A	122	PHE	4.1
1	E	38	TYR	4.0
1	E	58	GLU	4.0
1	E	62	LYS	4.0
1	B	44[A]	PHE	3.9
1	E	122	PHE	3.9
1	E	121	THR	3.9
1	C	163	GLU	3.8
1	D	161	PRO	3.7
1	E	148	LEU	3.7
1	E	69	SER	3.7
1	A	120	ALA	3.6
1	C	157	VAL	3.6
1	A	44	PHE	3.6
1	E	68	TYR	3.5
1	E	67	VAL	3.5
1	C	37	PHE	3.5
1	A	35	PHE	3.5
1	E	120	ALA	3.4
1	A	36	PHE	3.4
1	E	18	GLY	3.4
1	C	156	TYR	3.4
1	E	63	LEU	3.4
1	D	86	GLU	3.3
1	D	70	VAL	3.3
1	C	121	THR	3.2
1	D	45	VAL	3.2
1	E	26	LYS	3.2
1	C	122	PHE	3.2
1	E	152	LYS	3.1
1	D	35	PHE	3.1
1	A	96	ILE	3.1
1	B	37	PHE	3.1
1	D	36	PHE	3.1
1	B	70	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	34	VAL	3.0
1	E	71	SER	3.0
1	A	121	THR	3.0
1	A	69	SER	2.9
1	C	70	VAL	2.9
1	C	35	PHE	2.9
1	E	20	PHE	2.9
1	C	159	ALA	2.9
1	E	96	ILE	2.8
1	B	35	PHE	2.8
1	D	37	PHE	2.8
1	E	138	GLU	2.7
1	C	160	HIS	2.7
1	D	34	VAL	2.7
1	D	83	SER	2.7
1	C	86[A]	GLU	2.7
1	A	133	ILE	2.7
1	D	160	HIS	2.7
1	E	111	ARG	2.6
1	D	18	GLY	2.6
1	B	38	TYR	2.6
1	A	34	VAL	2.6
1	A	71	SER	2.5
1	C	62	LYS	2.5
1	E	59	GLU	2.5
1	E	86	GLU	2.5
1	C	69	SER	2.5
1	E	17	ASN	2.5
1	A	86	GLU	2.4
1	D	121	THR	2.4
1	A	38	TYR	2.4
1	B	161	PRO	2.4
1	D	68	TYR	2.4
1	C	36	PHE	2.4
1	B	122	PHE	2.3
1	E	139	GLY	2.3
1	E	158	ALA	2.3
1	A	68	TYR	2.3
1	C	96	ILE	2.3
1	A	90	LYS	2.3
1	A	165	CYS	2.3
1	E	149	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	94	ALA	2.2
1	E	11	LYS	2.2
1	C	133	ILE	2.2
1	D	120	ALA	2.2
1	B	121	THR	2.2
1	E	48	THR	2.2
1	C	120	ALA	2.2
1	B	71	SER	2.1
1	B	34	VAL	2.1
1	B	135	VAL	2.1
1	E	19	GLU	2.1
1	D	38	TYR	2.1
1	B	36	PHE	2.1
1	C	158	ALA	2.1
1	A	45	VAL	2.0
1	E	113	ASP	2.0
1	C	68	TYR	2.0
1	A	135	VAL	2.0
1	A	162	GLY	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 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, 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	LLDF	B-factors(Å ²)	Q<0.9
4	NA	B	203	1/1	0.53	0.36	13.88	89,89,89,89	0
3	K	D	202	1/1	0.74	0.39	6.84	99,99,99,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	K	A	202	1/1	0.93	0.26	2.28	66,66,66,66	1
3	K	B	202	1/1	0.95	0.22	1.75	72,72,72,72	0
2	CL	E	201	1/1	0.95	0.08	-0.56	73,73,73,73	0
2	CL	A	201	1/1	0.99	0.08	-0.65	42,42,42,42	0
2	CL	C	201	1/1	0.99	0.06	-1.10	48,48,48,48	0
2	CL	B	201	1/1	1.00	0.06	-1.13	43,43,43,43	0
2	CL	D	201	1/1	0.98	0.05	-1.26	51,51,51,51	0

6.5 Other polymers

There are no such residues in this entry.