



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:28 PM GMT

PDB ID : 4IFK
Title : Arginines 51 and 239* from a Neighboring Subunit are Essential for Catalysis
in a Zinc-dependent Decarboxylase
Authors : Huo, L.; Davis, I.; Chen, L.; Liu, A.
Deposited on : 2012-12-14
Resolution : 2.01 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

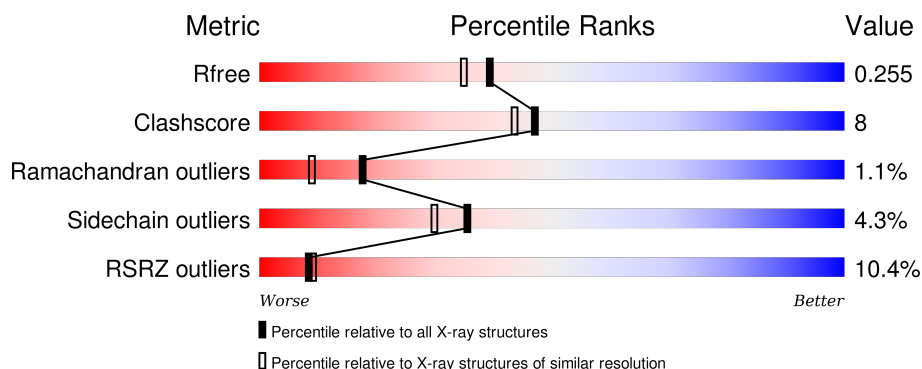
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

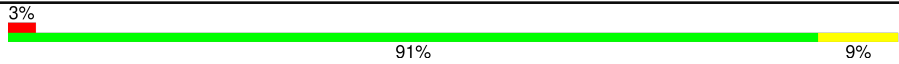

The reported resolution of this entry is 2.01 Å.

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}	91344	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	331	
2	B	331	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5416 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 2-amino-3-carboxymuconate 6-semialdehyde decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	4	0
			2588	1652	447	471	18			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	239	ALA	ARG	engineered mutation	UNP Q83V25

- Molecule 2 is a protein called 2-amino-3-carboxymuconate 6-semialdehyde decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	331	Total	C	N	O	S	0	3	0
			2596	1654	454	470	18			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	51	ALA	ARG	engineered mutation	UNP Q83V25

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		

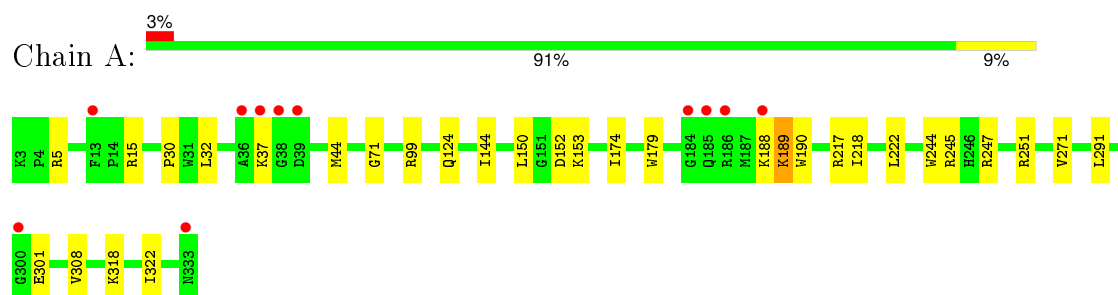
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	161	Total 161	O 161	0	0
5	B	68	Total 68	O 68	0	0

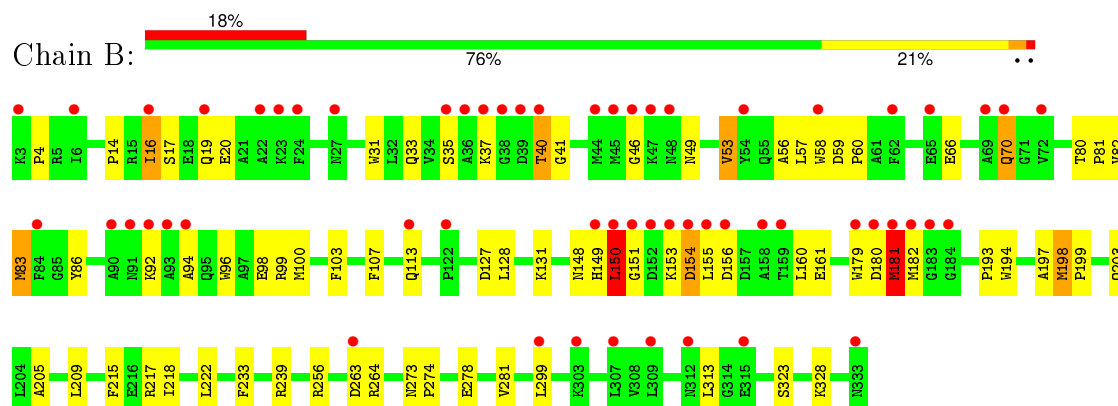
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 2-amino-3-carboxymuconate 6-semialdehyde decarboxylase



- Molecule 2: 2-amino-3-carboxymuconate 6-semialdehyde decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.85Å 49.06Å 110.55Å 90.00° 127.10° 90.00°	Depositor
Resolution (Å)	22.92 – 2.01 22.92 – 2.01	Depositor EDS
% Data completeness (in resolution range)	96.8 (22.92-2.01) 91.9 (22.92-2.01)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 2.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.209 , 0.263 0.203 , 0.255	Depositor DCC
R_{free} test set	2017 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 42741 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5416	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.41	0/2678	0.56	0/3630
2	B	0.35	0/2678	0.55	0/3628
All	All	0.38	0/5356	0.55	0/7258

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

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	2588	0	2528	22	0
2	B	2596	0	2540	58	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	161	0	0	7	0
5	B	68	0	0	7	0
All	All	5416	0	5068	77	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 8.

All (77) 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)
2:B:86:TYR:HB3	2:B:150:LEU:HD22	1.35	1.08
2:B:151:GLY:H	2:B:179:TRP:HZ3	1.21	0.89
2:B:181:MET:HG2	2:B:193:PRO:HB3	1.57	0.86
1:A:37:LYS:NZ	5:A:625:HOH:O	2.16	0.79
2:B:86:TYR:CB	2:B:150:LEU:HD22	2.15	0.75
2:B:66:GLU:O	5:B:561:HOH:O	2.03	0.75
1:A:152:ASP:OD1	5:A:649:HOH:O	2.03	0.74
2:B:239[A]:ARG:NE	5:B:515:HOH:O	2.24	0.70
1:A:144[B]:ILE:HD11	1:A:174:ILE:HG12	1.74	0.70
2:B:198:MET:HG3	2:B:199:PRO:HD3	1.75	0.67
1:A:189:LYS:HE3	1:A:190:TRP:H	1.60	0.67
1:A:37:LYS:NZ	5:A:634:HOH:O	2.29	0.66
2:B:181:MET:N	2:B:181:MET:SD	2.68	0.66
2:B:263:ASP:OD2	5:B:553:HOH:O	2.15	0.64
2:B:218:ILE:HG21	2:B:222:LEU:HD22	1.81	0.62
2:B:151:GLY:O	2:B:180:ASP:OD2	2.18	0.61
2:B:53:VAL:HG21	2:B:57:LEU:HD22	1.85	0.59
2:B:70:GLN:N	5:B:561:HOH:O	2.09	0.57
1:A:15:ARG:HD2	1:A:32:LEU:HD23	1.86	0.57
2:B:160:LEU:HD11	5:B:565:HOH:O	2.06	0.55
2:B:153:LYS:HG2	2:B:154:ASP:N	2.22	0.54
2:B:156:ASP:OD1	5:B:565:HOH:O	2.17	0.54
2:B:149:HIS:NE2	2:B:153:LYS:O	2.40	0.53
2:B:182:MET:O	5:B:527:HOH:O	2.19	0.53
5:A:647:HOH:O	2:B:239[B]:ARG:HG3	2.08	0.52
2:B:4:PRO:HG2	2:B:323:SER:OG	2.09	0.52
2:B:14:PRO:HG3	2:B:100:MET:HE3	1.92	0.52
1:A:124:GLN:HA	1:A:150:LEU:HD21	1.92	0.51
1:A:271:VAL:HG12	2:B:239[B]:ARG:HH21	1.75	0.51
2:B:80:THR:O	2:B:83:MET:HG2	2.10	0.51
2:B:127:ASP:O	2:B:131:LYS:HG2	2.12	0.50
1:A:291:LEU:HD22	1:A:322:ILE:HD12	1.93	0.50
2:B:86:TYR:CG	2:B:150:LEU:HD22	2.47	0.50
2:B:150:LEU:HB2	2:B:153:LYS:HE3	1.93	0.50
1:A:144[B]:ILE:CD1	1:A:174:ILE:HG12	2.40	0.50
2:B:151:GLY:O	2:B:180:ASP:CG	2.51	0.49
1:A:217:ARG:NH1	5:A:581:HOH:O	2.30	0.49
2:B:96:TRP:O	2:B:100:MET:HG2	2.13	0.48
2:B:81:PRO:HB3	2:B:179:TRP:CD1	2.49	0.48
2:B:86:TYR:CG	2:B:150:LEU:CD2	2.97	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:181:MET:HG3	2:B:197:ALA:HB3	1.96	0.47
2:B:80:THR:HG22	2:B:82:VAL:HG23	1.97	0.47
1:A:30:PRO:HA	1:A:44:MET:O	2.15	0.46
2:B:281:VAL:HG21	2:B:313:LEU:HD11	1.98	0.46
1:A:318:LYS:O	1:A:322:ILE:HG12	2.16	0.46
2:B:86:TYR:HB3	2:B:150:LEU:CD2	2.25	0.46
2:B:328:LYS:HE3	2:B:328:LYS:HB2	1.59	0.46
2:B:203:GLN:HG3	2:B:233:PHE:HA	1.97	0.46
2:B:58:TRP:CD1	2:B:59:ASP:HB2	2.51	0.46
2:B:60:PRO:HG3	2:B:107:PHE:CD1	2.51	0.45
2:B:274:PRO:O	2:B:278:GLU:HG3	2.16	0.45
1:A:99:ARG:NH1	5:A:588:HOH:O	2.31	0.45
1:A:247:ARG:HG2	2:B:299:LEU:HD21	1.98	0.44
1:A:5:ARG:HD3	1:A:71:GLY:O	2.17	0.44
2:B:131:LYS:HA	2:B:131:LYS:HD2	1.72	0.44
1:A:308:VAL:HG23	5:A:520:HOH:O	2.16	0.44
2:B:17:SER:OG	2:B:19:GLN:HG2	2.18	0.44
2:B:273:ASN:HA	2:B:274:PRO:HD3	1.89	0.43
2:B:14:PRO:HG2	2:B:16:ILE:HG12	1.99	0.43
2:B:40:THR:HG23	2:B:41:GLY:H	1.83	0.43
2:B:150:LEU:HB3	2:B:153:LYS:CG	2.49	0.43
1:A:244:TRP:O	1:A:251:ARG:HG3	2.19	0.43
2:B:161:GLU:HG3	2:B:218:ILE:HD11	2.02	0.42
2:B:150:LEU:HB2	2:B:153:LYS:CE	2.50	0.42
1:A:189:LYS:HA	1:A:189:LYS:HD2	1.82	0.42
2:B:20:GLU:OE2	2:B:99[B]:ARG:NH2	2.53	0.42
2:B:16:ILE:HD12	2:B:31:TRP:HA	2.02	0.42
2:B:81:PRO:HB3	2:B:179:TRP:NE1	2.35	0.42
1:A:218:ILE:HG21	1:A:222:LEU:HD22	2.02	0.42
2:B:150:LEU:HB3	2:B:153:LYS:CB	2.51	0.41
2:B:209:LEU:HA	2:B:209:LEU:HD23	1.88	0.41
1:A:271:VAL:HG12	2:B:239[B]:ARG:NH2	2.36	0.41
2:B:215:PHE:O	2:B:264:ARG:NH1	2.49	0.41
1:A:150:LEU:O	1:A:153:LYS:HB2	2.21	0.41
2:B:148:ASN:HB2	2:B:205:ALA:HB2	2.03	0.40
2:B:94:ALA:HB2	2:B:128:LEU:HD13	2.03	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	333/331 (101%)	325 (98%)	7 (2%)	1 (0%)	46	40
2	B	332/331 (100%)	308 (93%)	18 (5%)	6 (2%)	11	4
All	All	665/662 (100%)	633 (95%)	25 (4%)	7 (1%)	17	9

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	GLU
2	B	35	SER
2	B	181	MET
2	B	56	ALA
2	B	70	GLN
2	B	150	LEU
2	B	46	GLY

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	274/270 (102%)	270 (98%)	4 (2%)	72	74
2	B	273/270 (101%)	254 (93%)	19 (7%)	19	12
All	All	547/540 (101%)	524 (96%)	23 (4%)	35	31

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

Mol	Chain	Res	Type
1	A	179	TRP
1	A	188	LYS
1	A	189	LYS
1	A	245	ARG
2	B	16	ILE
2	B	33	GLN
2	B	37	LYS
2	B	40	THR
2	B	49	ASN
2	B	53	VAL
2	B	83	MET
2	B	92	LYS
2	B	98	GLU
2	B	103	PHE
2	B	113	GLN
2	B	150	LEU
2	B	154	ASP
2	B	155	LEU
2	B	181	MET
2	B	194	TRP
2	B	198	MET
2	B	217	ARG
2	B	256	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 3 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	331/331 (100%)	0.31	11 (3%)	50	52	17, 30, 49, 72	0
2	B	331/331 (100%)	1.11	58 (17%)	2	2	24, 45, 73, 90	0
All	All	662/662 (100%)	0.71	69 (10%)	8	9	17, 37, 66, 90	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	151	GLY	8.4
2	B	150	LEU	7.7
2	B	38	GLY	6.9
2	B	152	ASP	6.8
2	B	155	LEU	6.7
2	B	35	SER	6.0
1	A	36	ALA	5.8
2	B	54	TYR	5.7
2	B	181	MET	5.2
2	B	39	ASP	5.1
2	B	48	ASN	5.1
2	B	154	ASP	4.9
2	B	156	ASP	4.8
2	B	3	LYS	4.5
2	B	27	ASN	4.4
2	B	149	HIS	4.4
2	B	36	ALA	4.3
1	A	38	GLY	3.9
2	B	179	TRP	3.8
2	B	184	GLY	3.5
2	B	69	ALA	3.4
2	B	47	LYS	3.4
2	B	183	GLY	3.3
2	B	312	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	333	ASN	3.1
2	B	309	LEU	3.0
2	B	45	MET	2.9
2	B	91	ASN	2.9
2	B	22	ALA	2.9
2	B	46	GLY	2.9
2	B	16	ILE	2.8
2	B	23	LYS	2.8
2	B	62	PHE	2.8
2	B	153	LYS	2.7
2	B	65	GLU	2.7
2	B	180	ASP	2.7
2	B	159	THR	2.7
2	B	24	PHE	2.7
2	B	70	GLN	2.6
1	A	300	GLY	2.6
1	A	37	LYS	2.6
2	B	90	ALA	2.6
2	B	58	TRP	2.6
2	B	315	GLU	2.6
1	A	184	GLY	2.5
1	A	185	GLN	2.5
2	B	72	VAL	2.5
2	B	40	THR	2.5
1	A	186	ARG	2.4
2	B	113	GLN	2.4
2	B	93	ALA	2.4
2	B	19	GLN	2.4
1	A	13	PHE	2.3
1	A	39	ASP	2.3
2	B	6	ILE	2.3
2	B	158	ALA	2.3
2	B	92	LYS	2.2
2	B	307	LEU	2.2
1	A	188	LYS	2.1
2	B	122	PRO	2.1
2	B	182	MET	2.1
2	B	37	LYS	2.1
2	B	263	ASP	2.1
2	B	303	LYS	2.0
1	A	333	ASN	2.0
2	B	299	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
2	B	84	PHE	2.0
2	B	44	MET	2.0
2	B	94	ALA	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	ZN	B	401	1/1	0.99	0.06	-5.78	38,38,38,38	0
3	MG	A	401	1/1	0.61	0.16	-	60,60,60,60	0
4	ZN	A	402	1/1	0.99	0.10	-	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.