



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 04:21 AM UTC

PDB ID : 6SEN / pdb_00006sen
Title : TEAD4 bound to a FAM181A peptide
Authors : Scheufler, C.; Villard, F.
Deposited on : 2019-07-30
Resolution : 1.65 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (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

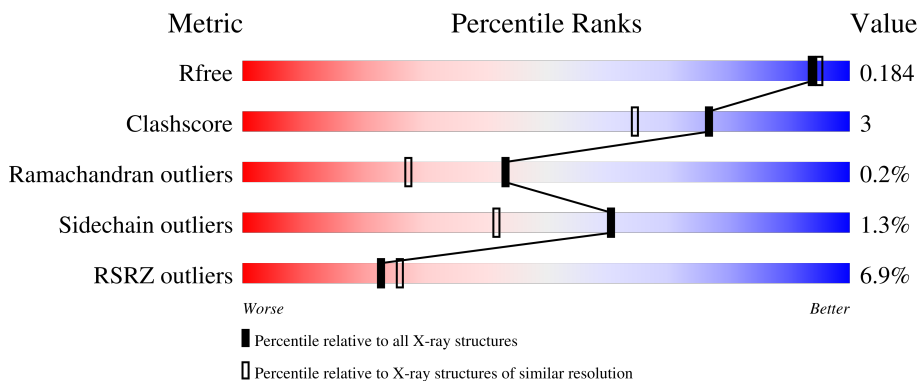
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	220	 7% 88% 5% 7%
1	B	220	 6% 90% 7% 7%
2	L	18	 6% 100%
2	M	18	 6% 94% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4070 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 Transcriptional enhancer factor TEF-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	205	Total	C	N	O	S	0	7	0
			1728	1110	287	322	9			
1	B	213	Total	C	N	O	S	0	7	0
			1789	1146	294	338	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	215	GLY	-	expression tag	UNP Q15561
A	216	PRO	-	expression tag	UNP Q15561
B	215	GLY	-	expression tag	UNP Q15561
B	216	PRO	-	expression tag	UNP Q15561

- Molecule 2 is a protein called Protein FAM181A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	18	Total	C	N	O	S	0	0	1
			138	89	25	23	1			
2	M	18	Total	C	N	O	S	0	1	1
			151	97	28	25	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	189	ACE	-	acetylation	UNP Q8N9Y4
L	206	NH2	-	amidation	UNP Q8N9Y4
M	189	ACE	-	acetylation	UNP Q8N9Y4
M	206	NH2	-	amidation	UNP Q8N9Y4

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	L	1	Total	O	S	0	0
			5	4	1		

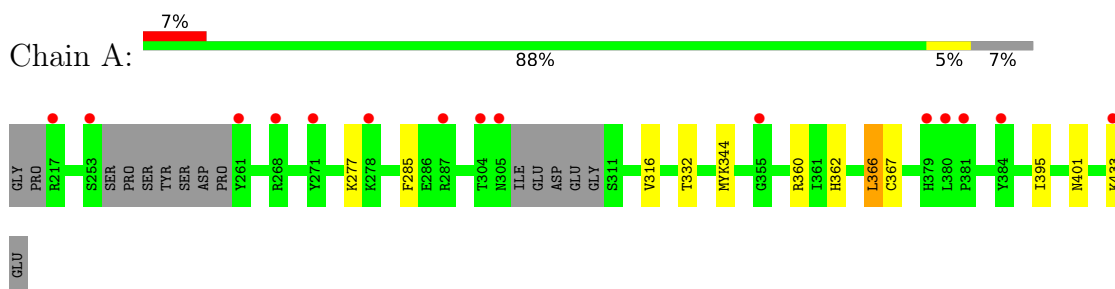
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	108	Total	O	0	0
			108	108		
4	B	103	Total	O	0	0
			103	103		
4	L	17	Total	O	0	0
			17	17		
4	M	11	Total	O	0	0
			11	11		

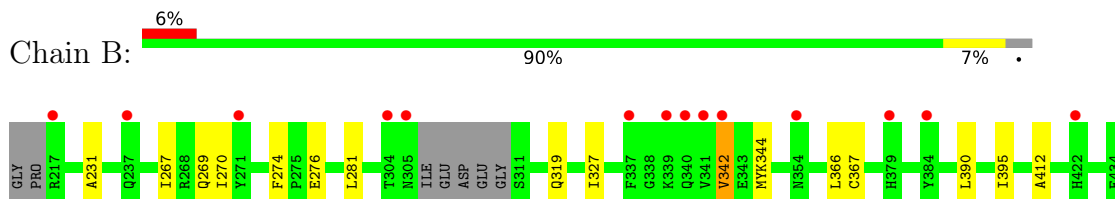
3 Residue-property plots [i](#)

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.

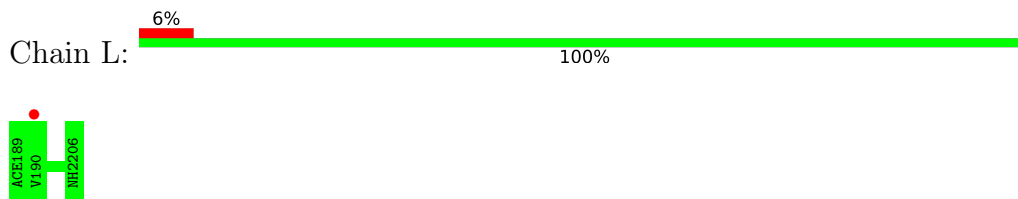
- Molecule 1: Transcriptional enhancer factor TEF-3



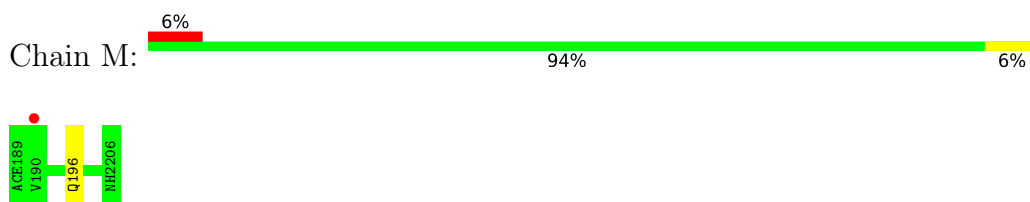
- Molecule 1: Transcriptional enhancer factor TEF-3



- Molecule 2: Protein FAM181A



- Molecule 2: Protein FAM181A



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	66.49Å 132.07Å 62.02Å 90.00° 115.87° 90.00°	Depositor
Resolution (Å)	23.75 – 1.65 23.75 – 1.65	Depositor EDS
% Data completeness (in resolution range)	81.8 (23.75-1.65) 81.7 (23.75-1.65)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 1.65Å)	Xtrriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.190 , 0.219 0.193 , 0.184	Depositor DCC
R_{free} test set	2479 reflections (3.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtrriage
Anisotropy	0.016	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4070	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% 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: SO4, MYK, NH2, ACE

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	0.77	0/1744	1.09	1/2358 (0.0%)
1	B	0.81	0/1810	1.12	2/2450 (0.1%)
2	L	0.75	0/140	1.09	0/191
2	M	0.75	0/153	0.88	0/207
All	All	0.79	0/3847	1.10	3/5206 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	401	ASN	N-CA-C	-5.15	101.37	109.25
1	B	276	GLU	CB-CG-CD	5.04	121.17	112.60
1	B	327	ILE	N-CA-C	-5.03	100.64	107.99

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	1728	0	1647	13	0
1	B	1789	0	1697	9	0
2	L	138	0	130	0	0
2	M	151	0	148	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	L	5	0	0	0	0
4	A	108	0	0	0	0
4	B	103	0	0	0	0
4	L	17	0	0	0	0
4	M	11	0	0	0	0
All	All	4070	0	3622	22	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 3.

All (22) 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:A:332:THR:HG21	1:A:344:MYK:HT	1.70	0.72
1:B:342:VAL:HG11	1:B:367[A]:CYS:SG	2.43	0.58
1:A:360:ARG:HD3	1:A:362:HIS:HB2	1.90	0.53
1:A:316[A]:VAL:CG2	1:A:366:LEU:HD12	2.39	0.53
1:A:344:MYK:HD	1:A:367:CYS:HB3	1.91	0.52
1:A:316[A]:VAL:CG2	1:A:366:LEU:CD1	2.88	0.51
1:A:344:MYK:HLA	1:A:395:ILE:HD13	1.93	0.51
1:A:344:MYK:HP	1:A:395:ILE:CD1	2.43	0.49
1:A:285:PHE:HE2	1:A:433:LYS:HB3	1.79	0.48
1:A:316[A]:VAL:HG21	1:A:344:MYK:HKA	1.96	0.48
1:A:316[A]:VAL:HG21	1:A:344:MYK:HYA	1.96	0.47
1:B:344:MYK:HI	1:B:395:ILE:HD13	1.96	0.47
1:A:332:THR:HG21	1:A:344:MYK:CT	2.41	0.47
1:B:231:ALA:HB1	1:B:344:MYK:HPA	1.97	0.46
1:A:316[A]:VAL:HG23	1:A:366:LEU:HD12	1.99	0.45
1:B:231:ALA:CB	1:B:344:MYK:HPA	2.47	0.45
1:A:344:MYK:HI	1:A:344:MYK:OX	2.18	0.44
1:B:269:GLN:HG3	2:M:196[A]:GLN:O	2.19	0.43
1:B:344:MYK:HV	1:B:390:LEU:HD21	2.01	0.42
1:B:274:PHE:CZ	1:B:412:ALA:HB1	2.54	0.42
1:B:344:MYK:HU	1:B:344:MYK:HW	1.93	0.42
1:B:267:ILE:HD11	1:B:281:LEU:HG	2.02	0.41

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	205/220 (93%)	203 (99%)	1 (0%)	1 (0%)	24	10
1	B	215/220 (98%)	207 (96%)	8 (4%)	0	100	100
2	L	16/18 (89%)	16 (100%)	0	0	100	100
2	M	17/18 (94%)	17 (100%)	0	0	100	100
All	All	453/476 (95%)	443 (98%)	9 (2%)	1 (0%)	43	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	LYS

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	186/200 (93%)	185 (100%)	1 (0%)	81	72
1	B	196/200 (98%)	192 (98%)	4 (2%)	48	26
2	L	14/15 (93%)	14 (100%)	0	100	100
2	M	16/15 (107%)	16 (100%)	0	100	100
All	All	412/430 (96%)	407 (99%)	5 (1%)	61	45

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	366	LEU
1	B	270	ILE
1	B	319	GLN
1	B	342	VAL
1	B	366	LEU

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

Mol	Chain	Res	Type
1	A	236	GLN
1	A	422	HIS
1	B	243	ASN
1	B	291	ASN
1	B	305	ASN
1	B	340	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](#)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
1	MYK	B	344	1	22,23,24	0.79	0	19,24,26	0.75	0
1	MYK	A	344	1	22,23,24	0.72	0	19,24,26	0.77	0

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
1	MYK	B	344	1	-	10/22/23/25	-
1	MYK	A	344	1	-	7/22/23/25	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	344	MYK	CI-CT-CY-CX
1	A	344	MYK	CG-CD-CE-NZ
1	B	344	MYK	CT-CI-CK-CL
1	B	344	MYK	CM-CP-CR-CU
1	A	344	MYK	CP-CR-CU-CQ
1	B	344	MYK	CK-CI-CT-CY
1	B	344	MYK	CE-CD-CG-CB
1	A	344	MYK	CM-CP-CR-CU
1	A	344	MYK	CU-CQ-CS-CW
1	B	344	MYK	CK-CL-CM-CP
1	B	344	MYK	CQ-CS-CW-CV
1	A	344	MYK	CK-CI-CT-CY
1	B	344	MYK	C-CA-CB-CG
1	B	344	MYK	CP-CR-CU-CQ
1	B	344	MYK	CU-CQ-CS-CW
1	A	344	MYK	CQ-CS-CW-CV
1	A	344	MYK	CD-CE-NZ-CX

There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	344	MYK	5	0
1	A	344	MYK	8	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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	$\# Z > 2$
3	SO4	B	502	-	4,4,4	0.29	0	6,6,6	0.58	0
3	SO4	A	501	-	4,4,4	0.39	0	6,6,6	0.39	0
3	SO4	B	501	-	4,4,4	0.42	0	6,6,6	0.18	0
3	SO4	L	301	-	4,4,4	0.29	0	6,6,6	0.19	0
3	SO4	A	502	-	4,4,4	0.28	0	6,6,6	0.37	0

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	204/220 (92%)	0.33	15 (7%) 20 23	12, 32, 59, 86	7 (3%)
1	B	212/220 (96%)	0.37	14 (6%) 24 27	11, 35, 57, 84	7 (3%)
2	L	16/18 (88%)	0.46	1 (6%) 26 29	26, 33, 63, 71	0
2	M	16/18 (88%)	0.20	1 (6%) 26 29	21, 32, 59, 60	1 (6%)
All	All	448/476 (94%)	0.35	31 (6%) 23 26	11, 33, 60, 86	15 (3%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	384	TYR	4.2
2	L	190	VAL	4.2
1	A	217	ARG	3.9
1	A	305	ASN	3.7
1	B	339	LYS	3.3
1	A	261	TYR	3.2
1	A	271	TYR	3.2
1	B	217	ARG	3.2
1	B	341	VAL	3.1
1	A	380	LEU	3.1
1	B	305	ASN	3.0
1	B	337	PHE	3.0
1	A	287[A]	ARG	2.9
1	A	304	THR	2.8
1	A	433	LYS	2.7
1	A	253	SER	2.6
1	A	355	GLY	2.6
1	B	271	TYR	2.5
1	B	342	VAL	2.5
1	B	422	HIS	2.4
1	A	381	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	379	HIS	2.3
1	A	278	LYS	2.3
1	B	237	GLN	2.3
1	B	340	GLN	2.2
1	A	379	HIS	2.2
1	B	304	THR	2.2
1	B	384	TYR	2.1
2	M	190	VAL	2.1
1	B	354	ASN	2.1
1	A	268	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	MYK	A	344	24/25	0.85	0.16	34,51,59,59	0
1	MYK	B	344	24/25	0.90	0.14	30,47,53,57	0

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(Å ²)	Q<0.9
3	SO4	L	301	5/5	0.82	0.12	60,60,63,63	0
3	SO4	A	501	5/5	0.97	0.06	38,38,39,40	0
3	SO4	B	502	5/5	0.98	0.05	36,37,38,39	0
3	SO4	A	502	5/5	0.98	0.06	29,30,33,36	0
3	SO4	B	501	5/5	0.99	0.04	27,30,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.