



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 9, 2017 – 10:49 am GMT

PDB ID : 5K0K
Title : Crystal structure of the catalytic domain of the proto-oncogene tyrosine-protein kinase MER in complex with inhibitor UNC2434
Authors : Wang, X.; Liu, J.; Zhang, W.; Stashko, M.A.; Nichols, J.; DeRyckere, D.; Miley, M.J.; Norris-Drouin, J.; Chen, Z.; Machius, M.; Wood, E.; Graham, D.K.; Earp, H.S.; Graham, K.; Kireev, D.; Frye, S.V.
Deposited on : 2016-05-17
Resolution : 2.54 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : recalc28906
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) : recalc28906

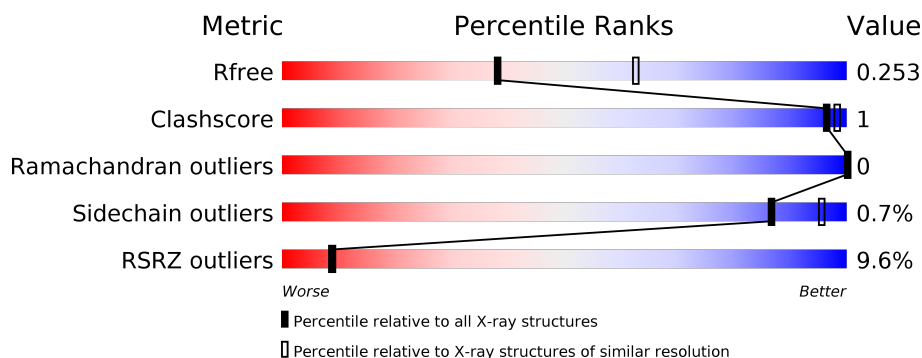
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	4993 (2.58-2.50)
Clashscore	112137	5755 (2.58-2.50)
Ramachandran outliers	110173	5652 (2.58-2.50)
Sidechain outliers	110143	5654 (2.58-2.50)
RSRZ outliers	101464	5026 (2.58-2.50)

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	313	<div> <div>7%</div> <div>81%</div> <div>17%</div> </div>
1	B	313	<div> <div>8%</div> <div>77%</div> <div>21%</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
2	CL	A	901	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8443 atoms, of which 4226 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 Tyrosine-protein kinase Mer.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	260	Total	C	H	N	O	S	0	2	0
			4239	1349	2128	354	389	19			
1	B	248	Total	C	H	N	O	S	0	3	0
			4031	1282	2028	337	365	19			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	552	MET	-	initiating methionine	UNP Q12866
A	553	GLY	-	expression tag	UNP Q12866
A	554	SER	-	expression tag	UNP Q12866
A	555	SER	-	expression tag	UNP Q12866
A	556	HIS	-	expression tag	UNP Q12866
A	557	HIS	-	expression tag	UNP Q12866
A	558	HIS	-	expression tag	UNP Q12866
A	559	HIS	-	expression tag	UNP Q12866
A	560	HIS	-	expression tag	UNP Q12866
A	561	HIS	-	expression tag	UNP Q12866
A	562	SER	-	expression tag	UNP Q12866
A	563	SER	-	expression tag	UNP Q12866
A	564	GLY	-	expression tag	UNP Q12866
A	565	LEU	-	expression tag	UNP Q12866
A	566	VAL	-	expression tag	UNP Q12866
A	567	PRO	-	expression tag	UNP Q12866
A	568	ARG	-	expression tag	UNP Q12866
A	569	GLY	-	expression tag	UNP Q12866
B	552	MET	-	initiating methionine	UNP Q12866
B	553	GLY	-	expression tag	UNP Q12866
B	554	SER	-	expression tag	UNP Q12866
B	555	SER	-	expression tag	UNP Q12866
B	556	HIS	-	expression tag	UNP Q12866
B	557	HIS	-	expression tag	UNP Q12866
B	558	HIS	-	expression tag	UNP Q12866

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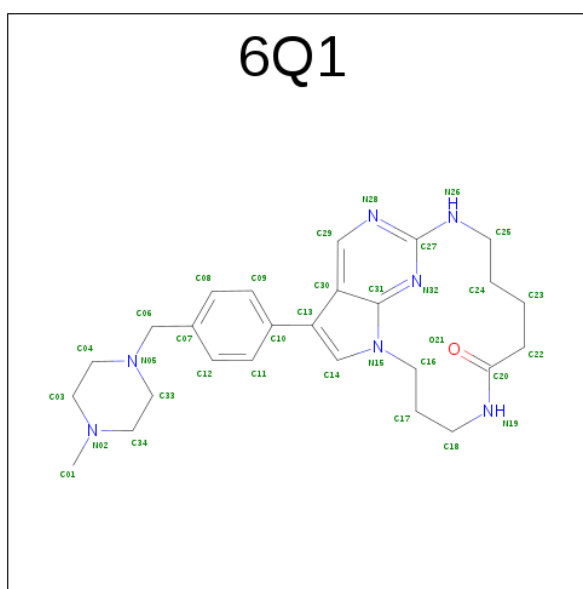
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Chain	Residue	Modelled	Actual	Comment	Reference
B	559	HIS	-	expression tag	UNP Q12866
B	560	HIS	-	expression tag	UNP Q12866
B	561	HIS	-	expression tag	UNP Q12866
B	562	SER	-	expression tag	UNP Q12866
B	563	SER	-	expression tag	UNP Q12866
B	564	GLY	-	expression tag	UNP Q12866
B	565	LEU	-	expression tag	UNP Q12866
B	566	VAL	-	expression tag	UNP Q12866
B	567	PRO	-	expression tag	UNP Q12866
B	568	ARG	-	expression tag	UNP Q12866
B	569	GLY	-	expression tag	UNP Q12866

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Cl 2 2	0	0
2	A	4	Total Cl 4 4	0	0

- Molecule 3 is 15-{4-[(4-methylpiperazin-1-yl)methyl]phenyl}-4,5,6,7,9,10,11,12-octahydro-2,16-(azenometheno)pyrrolo[2,1-d][1,3,5,9]tetraazacyclotetradecin-8(3H)-one (three-letter code: 6Q1) (formula: C₂₆H₃₅N₇O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			69	26	35	7	1		
3	B	1	Total	C	H	N	O	0	0
			69	26	35	7	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

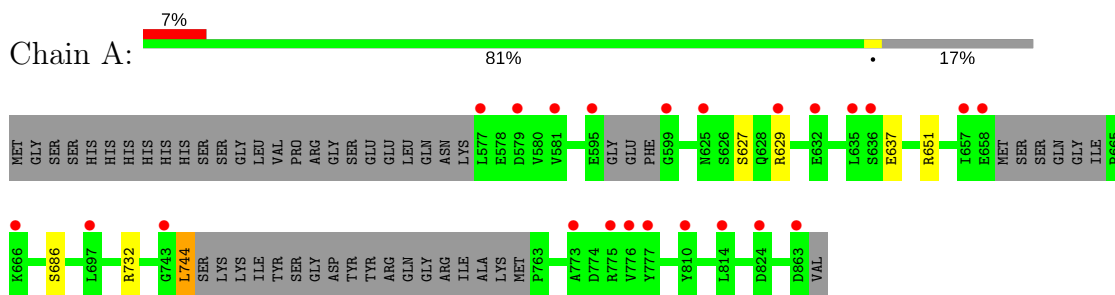
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	12	Total	O	0	0
			12	12		
5	B	16	Total	O	0	0
			16	16		

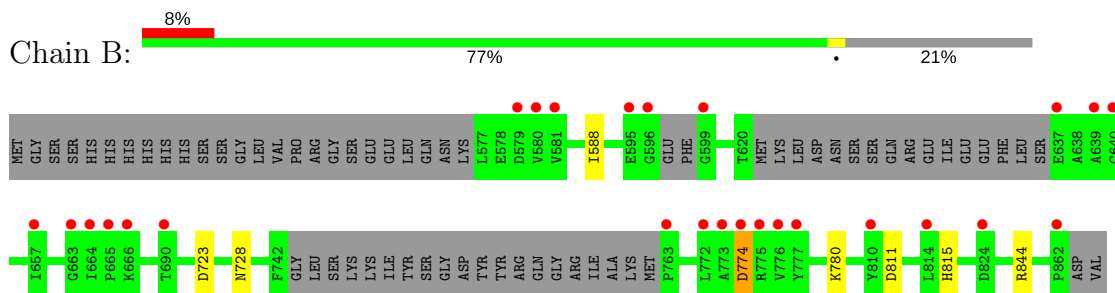
3 Residue-property plots [i](#)

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: Tyrosine-protein kinase Mer



• Molecule 1: Tyrosine-protein kinase Mer



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.07Å 91.31Å 69.20Å 90.00° 100.42° 90.00°	Depositor
Resolution (Å)	34.55 – 2.54 34.55 – 2.54	Depositor EDS
% Data completeness (in resolution range)	86.0 (34.55-2.54) 82.4 (34.55-2.54)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.54Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.198 , 0.254 0.192 , 0.253	Depositor DCC
R_{free} test set	1772 reflections (10.01%)	DCC
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8443	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.24% 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: MG, 6Q1, 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	0.24	0/2153	0.41	0/2907
1	B	0.24	0/2052	0.39	0/2773
All	All	0.24	0/4205	0.40	0/5680

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	2111	2128	2120	5	0
1	B	2003	2028	2004	5	0
2	A	4	0	0	3	0
2	B	2	0	0	1	0
3	A	34	35	0	0	0
3	B	34	35	0	0	0
4	B	1	0	0	0	0
5	A	12	0	0	0	0
5	B	16	0	0	0	0
All	All	4217	4226	4124	10	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 (10) 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:723:ASP:O	1:B:728[B]:ASN:ND2	2.20	0.74
1:B:780:LYS:NZ	1:B:844:ARG:O	2.29	0.65
1:B:588:ILE:HA	2:B:901:CL:CL	2.48	0.51
1:B:774:ASP:OD1	1:B:774:ASP:N	2.45	0.50
1:A:637:GLU:OE1	1:A:744:LEU:N	2.43	0.50
1:A:651:ARG:HB2	2:A:901:CL:CL	2.49	0.49
1:A:732:ARG:NH1	2:A:901:CL:CL	2.77	0.48
1:B:811:ASP:O	1:B:815:HIS:ND1	2.46	0.46
1:A:627:SER:OG	1:A:629:ARG:NH2	2.49	0.44
1:A:651:ARG:CB	2:A:901:CL:CL	3.04	0.43

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	254/313 (81%)	249 (98%)	5 (2%)	0	100	100
1	B	243/313 (78%)	235 (97%)	8 (3%)	0	100	100
All	All	497/626 (79%)	484 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	237/280 (85%)	235 (99%)	2 (1%)	85	94
1	B	225/280 (80%)	224 (100%)	1 (0%)	93	98
All	All	462/560 (82%)	459 (99%)	3 (1%)	87	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	686	SER
1	A	744	LEU
1	B	774	ASP

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

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, 7 are monoatomic - leaving 2 for Mogul analysis.

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	6Q1	A	904	-	35,38,38	2.73	8 (22%)	42,52,52	2.30	8 (19%)
3	6Q1	B	904	-	35,38,38	2.69	8 (22%)	42,52,52	2.44	11 (26%)

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
3	6Q1	A	904	-	-	0/23/33/33	0/3/5/5
3	6Q1	B	904	-	-	0/23/33/33	0/3/5/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	904	6Q1	C13-C10	-3.45	1.43	1.49
3	A	904	6Q1	C13-C10	-3.44	1.43	1.49
3	A	904	6Q1	C16-N15	-2.03	1.44	1.49
3	B	904	6Q1	C25-N26	-2.01	1.41	1.46
3	B	904	6Q1	C14-C13	2.20	1.41	1.38
3	B	904	6Q1	C31-N32	2.36	1.39	1.35
3	B	904	6Q1	C11-C12	2.44	1.43	1.38
3	A	904	6Q1	C11-C12	2.51	1.43	1.38
3	A	904	6Q1	C14-C13	2.53	1.41	1.38
3	A	904	6Q1	C31-N32	2.54	1.39	1.35
3	B	904	6Q1	C20-N19	4.57	1.44	1.33
3	A	904	6Q1	C20-N19	4.82	1.44	1.33
3	B	904	6Q1	C27-N28	4.82	1.40	1.34
3	A	904	6Q1	C27-N28	5.13	1.40	1.34
3	A	904	6Q1	C27-N26	11.78	1.53	1.34
3	B	904	6Q1	C27-N26	11.80	1.53	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	904	6Q1	N28-C27-N32	-5.66	121.00	126.65
3	A	904	6Q1	N28-C27-N32	-5.48	121.18	126.65
3	B	904	6Q1	O21-C20-N19	-2.41	118.38	122.97
3	B	904	6Q1	C25-N26-C27	-2.05	119.94	123.68
3	B	904	6Q1	C24-C25-N26	2.12	117.11	111.46
3	B	904	6Q1	C03-N02-C34	2.24	112.46	109.47
3	B	904	6Q1	C29-N28-C27	2.31	119.64	115.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	904	6Q1	C29-N28-C27	2.49	119.94	115.89
3	A	904	6Q1	C24-C25-N26	2.96	119.33	111.46
3	A	904	6Q1	C06-N05-C33	3.23	117.91	111.08
3	B	904	6Q1	C22-C20-N19	3.34	122.25	116.49
3	B	904	6Q1	C01-N02-C34	4.35	117.23	110.67
3	A	904	6Q1	C06-N05-C04	4.61	120.85	111.08
3	A	904	6Q1	C01-N02-C34	4.70	117.76	110.67
3	A	904	6Q1	C27-N32-C31	5.14	120.98	115.11
3	B	904	6Q1	C06-N05-C04	5.63	123.01	111.08
3	B	904	6Q1	C27-N32-C31	5.76	121.68	115.11
3	A	904	6Q1	C01-N02-C03	7.82	122.46	110.67
3	B	904	6Q1	C01-N02-C03	8.14	122.94	110.67

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	260/313 (83%)	0.49	23 (8%) 11 11	26, 46, 91, 147	0
1	B	248/313 (79%)	0.54	26 (10%) 7 7	29, 49, 87, 118	0
All	All	508/626 (81%)	0.52	49 (9%) 9 9	26, 47, 90, 147	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	657	ILE	5.4
1	B	637	GLU	5.2
1	A	577	LEU	4.6
1	B	639	ALA	4.4
1	A	777	TYR	4.0
1	B	777	TYR	3.9
1	A	773	ALA	3.8
1	B	774	ASP	3.8
1	B	776	VAL	3.7
1	A	579	ASP	3.6
1	A	635	LEU	3.6
1	A	658	GLU	3.5
1	A	810	TYR	3.5
1	B	773	ALA	3.5
1	A	775	ARG	3.4
1	B	862	PRO	3.4
1	B	580	VAL	3.3
1	A	629	ARG	3.3
1	B	665	PRO	3.2
1	A	824	ASP	3.2
1	B	772	LEU	3.1
1	B	824	ASP	3.1
1	B	579	ASP	3.0
1	A	595	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	666	LYS	3.0
1	B	775	ARG	2.9
1	B	664	ILE	2.9
1	B	814	LEU	2.9
1	A	657	ILE	2.9
1	B	663	GLY	2.8
1	B	810	TYR	2.8
1	A	666	LYS	2.7
1	A	743	GLY	2.7
1	A	599	GLY	2.4
1	B	690	THR	2.4
1	A	636	SER	2.3
1	B	596	GLY	2.3
1	B	595	GLU	2.3
1	B	763	PRO	2.3
1	A	632	GLU	2.3
1	A	581	VAL	2.2
1	B	640	CYS	2.2
1	A	814	LEU	2.1
1	A	776	VAL	2.1
1	A	625	ASN	2.1
1	A	697	LEU	2.1
1	B	599	GLY	2.0
1	B	581	VAL	2.0
1	A	863	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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(\AA^2)	Q<0.9
3	6Q1	A	904	34/34	0.91	0.21	0.51	46,56,67,69	0
3	6Q1	B	904	34/34	0.90	0.19	0.40	38,57,71,72	0
2	CL	A	905	1/1	0.98	0.17	0.35	48,48,48,48	0
2	CL	A	902	1/1	0.99	0.14	-0.98	40,40,40,40	0
2	CL	B	901	1/1	0.97	0.11	-1.05	44,44,44,44	0
2	CL	B	902	1/1	0.99	0.13	-1.24	37,37,37,37	0
2	CL	A	903	1/1	0.96	0.14	-	64,64,64,64	0
2	CL	A	901	1/1	0.94	0.09	-	46,46,46,46	0
4	MG	B	903	1/1	0.94	0.15	-	49,49,49,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.