



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 07:16 AM GMT

PDB ID : 3A2C
Title : Crystal structure of a pyrazolopyrimidine inhibitor complex bound to MAP-KAP Kinase-2 (MK2)
Authors : Fujino, A.; Takimoto-Kamimura, M.
Deposited on : 2009-05-12
Resolution : 2.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

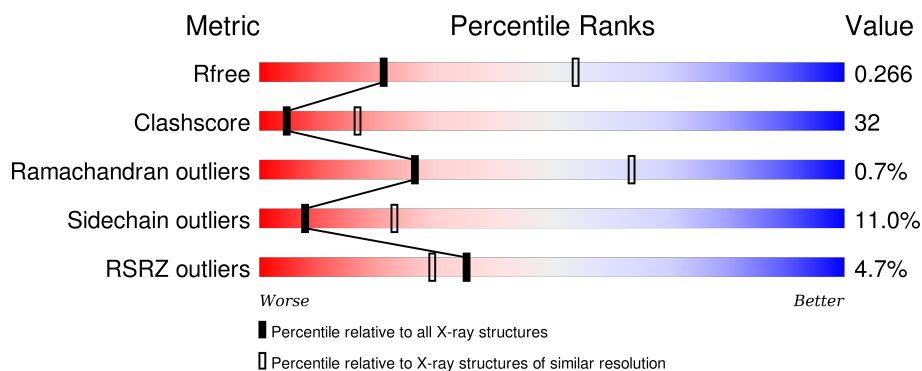
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	324	<div> <div>3%</div> <div> <div>44%</div> <div>35%</div> <div>5%</div> <div>15%</div> </div> </div>
1	B	324	<div> <div>2%</div> <div> <div>40%</div> <div>40%</div> <div>5%</div> <div>15%</div> </div> </div>
1	C	324	<div> <div>4%</div> <div> <div>40%</div> <div>35%</div> <div>9%</div> <div>16%</div> </div> </div>
1	D	324	<div> <div>3%</div> <div> <div>46%</div> <div>33%</div> <div>6%</div> <div>15%</div> </div> </div>
1	E	324	<div> <div>3%</div> <div> <div>47%</div> <div>32%</div> <div>6%</div> <div>15%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	324	
1	G	324	
1	H	324	
1	I	324	
1	J	324	
1	K	324	
1	L	324	

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	PDY	A	2	-	-	-	X
2	PDY	C	2	-	-	X	-
2	PDY	D	2	-	-	-	X
2	PDY	E	1	-	-	X	X
2	PDY	E	2	-	-	X	X
2	PDY	F	2	-	-	X	-
2	PDY	G	1	-	-	X	-
2	PDY	G	2	-	-	X	X
2	PDY	H	1	-	-	X	-
2	PDY	H	2	-	-	X	-
2	PDY	I	1	-	-	X	-
2	PDY	I	2	-	-	X	X
2	PDY	J	1	-	-	X	-
2	PDY	K	1	-	-	X	-
2	PDY	K	2	-	-	X	-
2	PDY	L	1	-	-	X	-
2	PDY	L	2	-	-	X	X

2 Entry composition

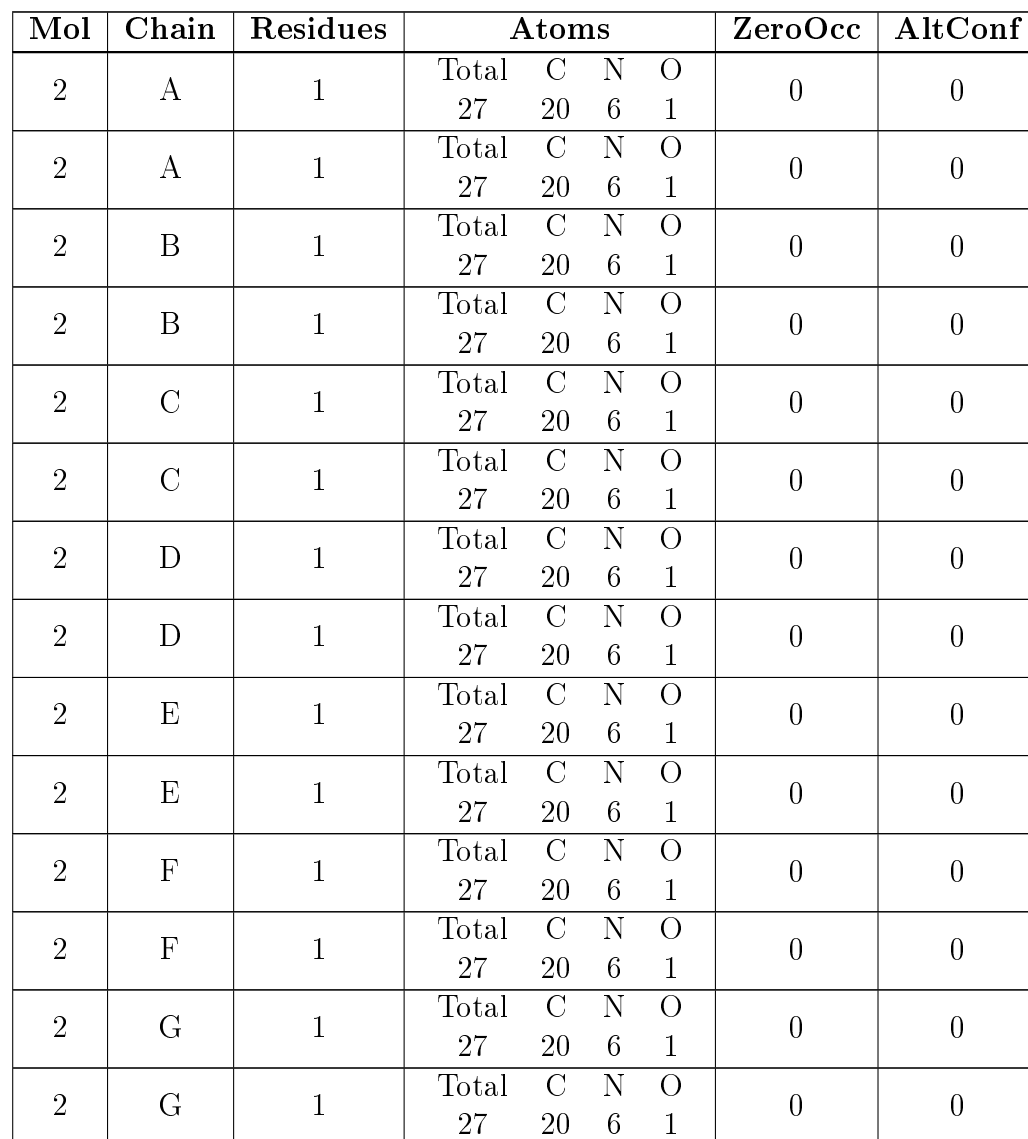
There are 3 unique types of molecules in this entry. The entry contains 27440 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MAP kinase-activated protein kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2218	1415	384	402	17			
1	B	277	Total	C	N	O	S	0	0	0
			2240	1426	388	409	17			
1	C	273	Total	C	N	O	S	0	0	0
			2207	1409	380	401	17			
1	D	277	Total	C	N	O	S	0	0	0
			2244	1430	389	408	17			
1	E	276	Total	C	N	O	S	0	0	0
			2234	1425	385	407	17			
1	F	287	Total	C	N	O	S	0	0	0
			2316	1476	401	422	17			
1	G	273	Total	C	N	O	S	0	0	0
			2210	1411	381	401	17			
1	H	275	Total	C	N	O	S	0	0	0
			2226	1421	385	403	17			
1	I	275	Total	C	N	O	S	0	0	0
			2225	1420	384	404	17			
1	J	275	Total	C	N	O	S	0	0	0
			2225	1420	384	404	17			
1	K	273	Total	C	N	O	S	0	0	0
			2208	1409	381	401	17			
1	L	276	Total	C	N	O	S	0	0	0
			2234	1425	385	407	17			

- Molecule 2 is N 7 -(4-ETHOXYPHENYL)-6-METHYL-N 5 -[(3S)-PIPERIDIN-3-YL]PYRAZOLO[1,5-A]PYRIMIDINE-5,7-DIAMINE (three-letter code: PDY) (formula: C₂₀H₂₆N₆O).

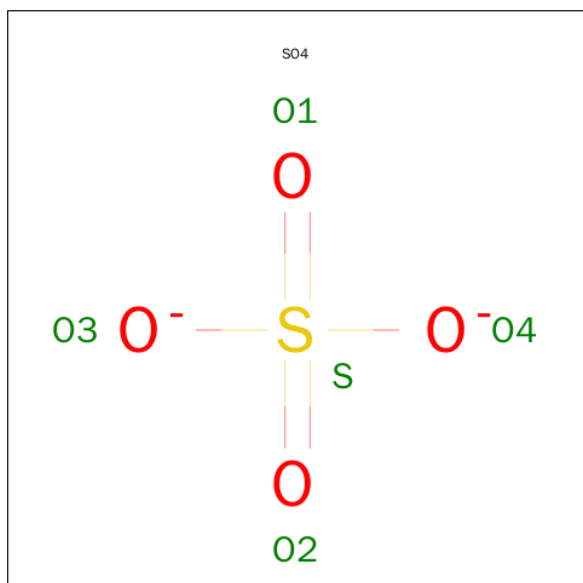


WORLDWIDE
PDB
PROTEIN DATA BANK

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	H	1	Total	C	N	O	0	0
			27	20	6	1		
2	H	1	Total	C	N	O	0	0
			27	20	6	1		
2	I	1	Total	C	N	O	0	0
			27	20	6	1		
2	I	1	Total	C	N	O	0	0
			27	20	6	1		
2	J	1	Total	C	N	O	0	0
			27	20	6	1		
2	J	1	Total	C	N	O	0	0
			27	20	6	1		
2	K	1	Total	C	N	O	0	0
			27	20	6	1		
2	K	1	Total	C	N	O	0	0
			27	20	6	1		
2	L	1	Total	C	N	O	0	0
			27	20	6	1		
2	L	1	Total	C	N	O	0	0
			27	20	6	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

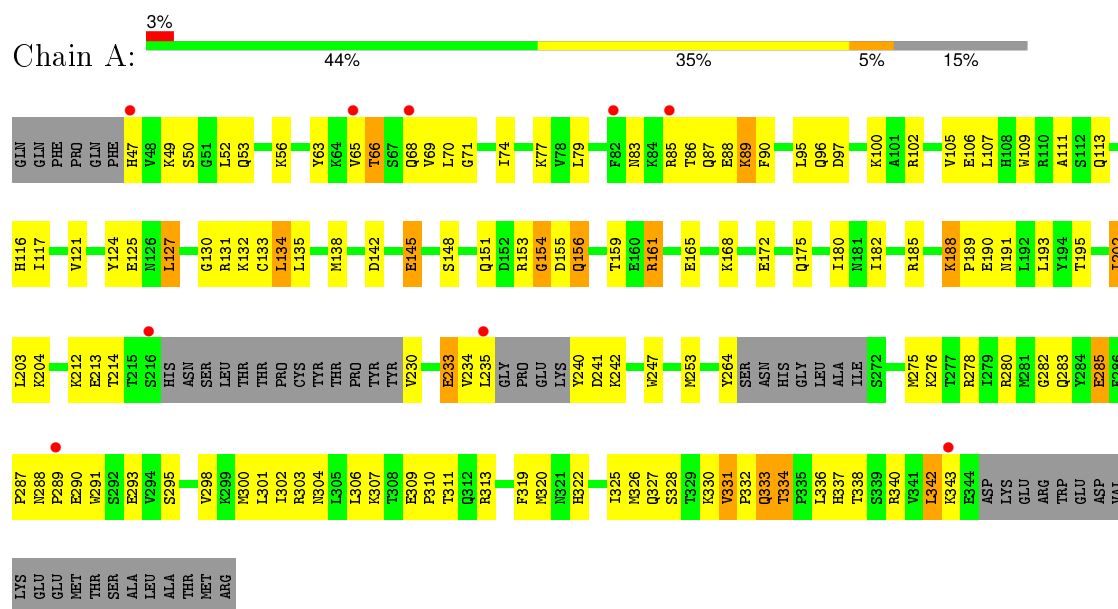


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	O	S	0	0
			5	4	1		

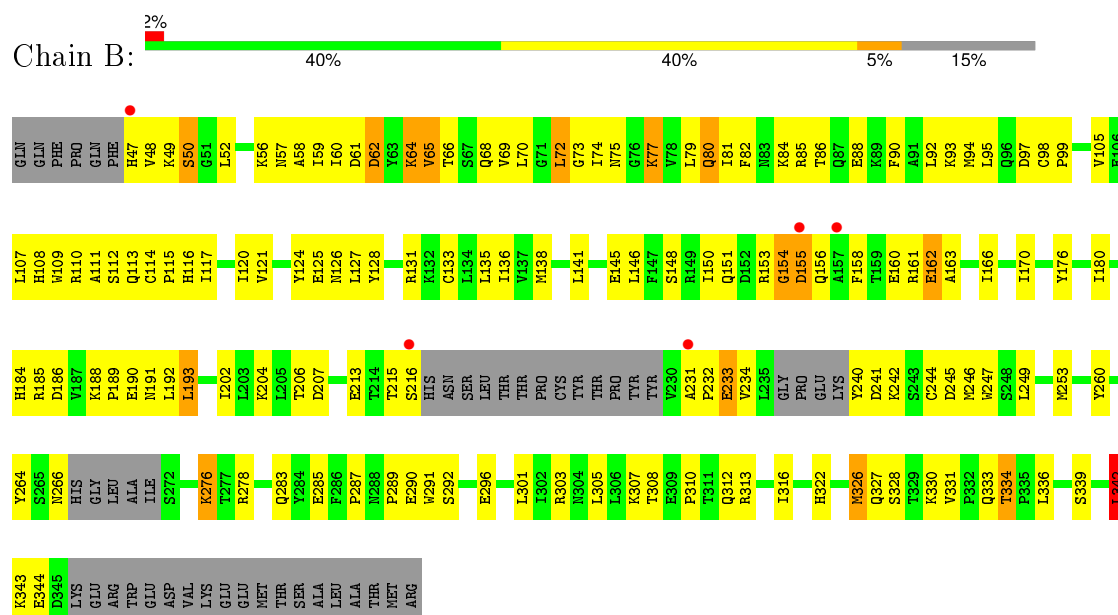
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

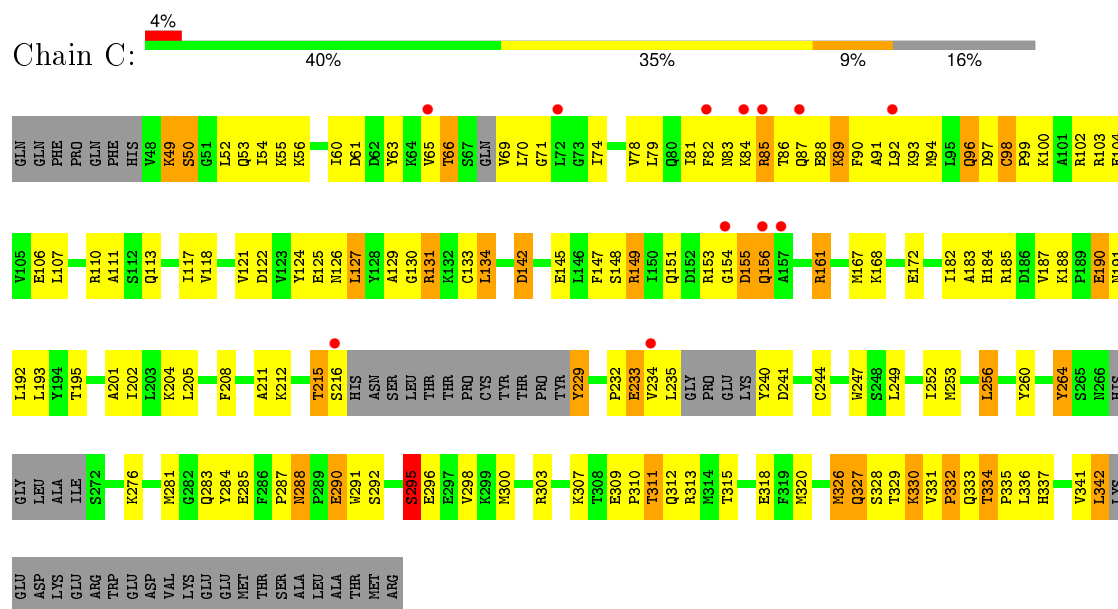
- Molecule 1: MAP kinase-activated protein kinase 2



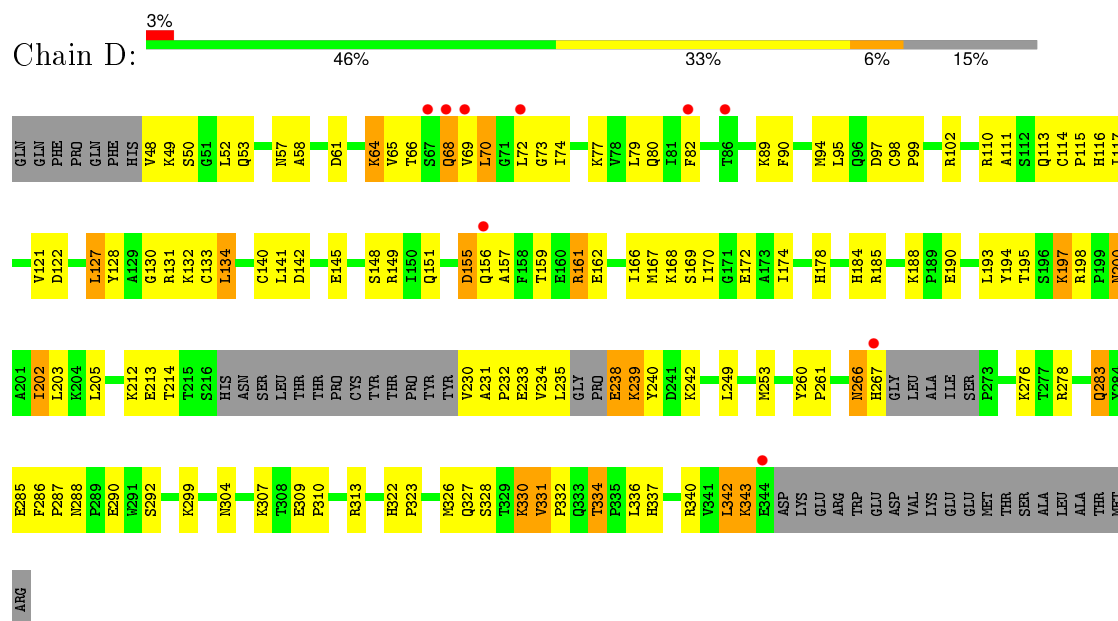
- Molecule 1: MAP kinase-activated protein kinase 2



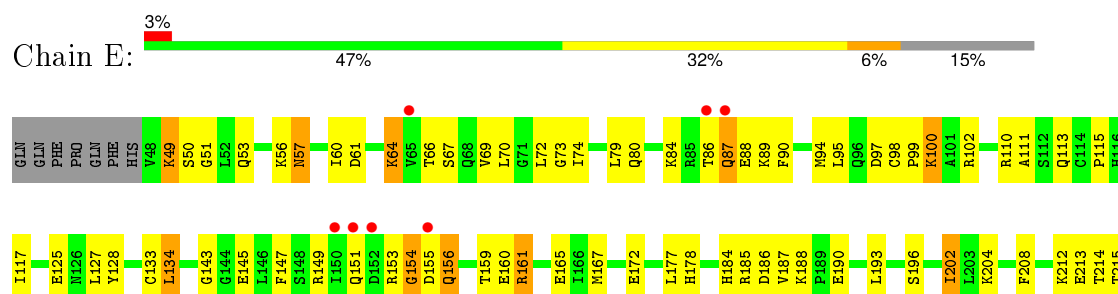
- Molecule 1: MAP kinase-activated protein kinase 2

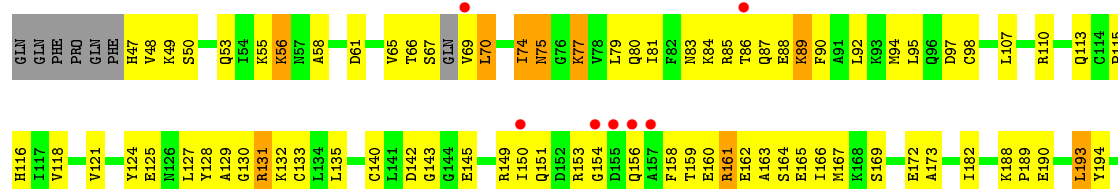


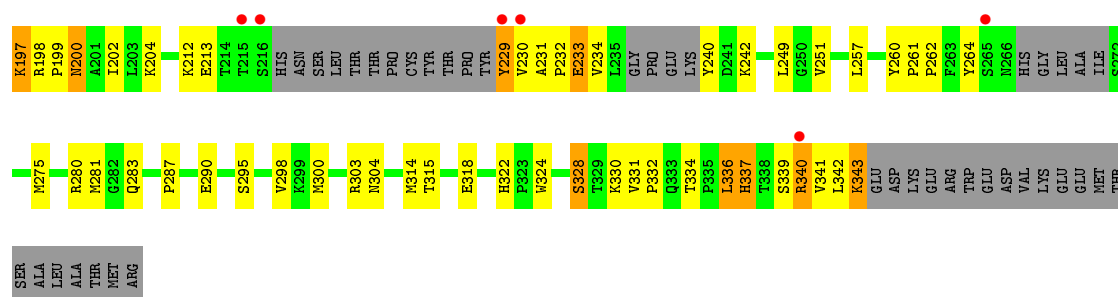
- Molecule 1: MAP kinase-activated protein kinase 2



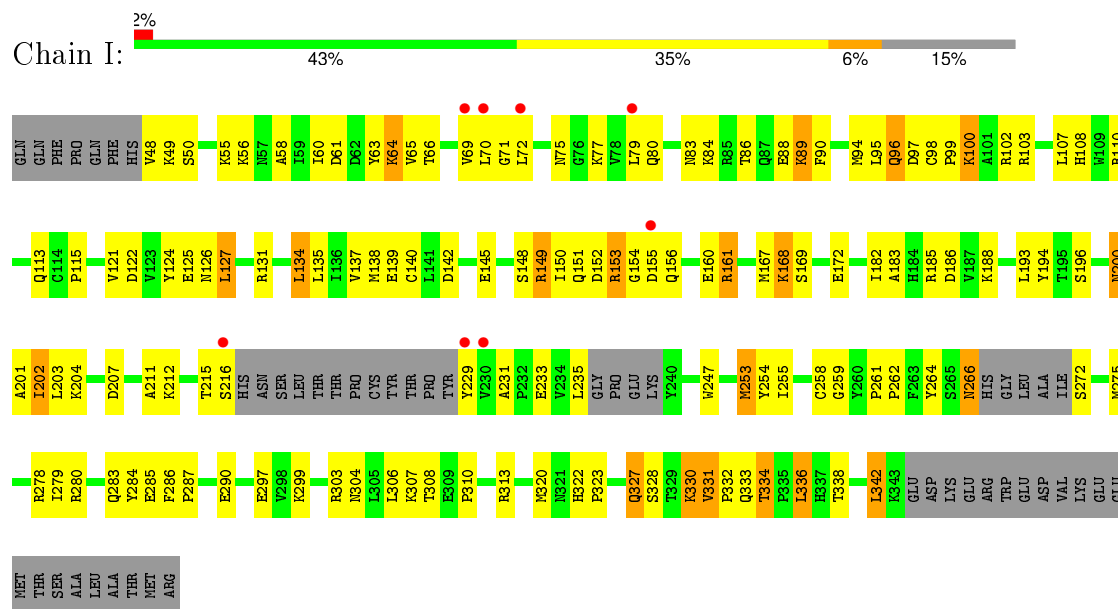
- Molecule 1: MAP kinase-activated protein kinase 2



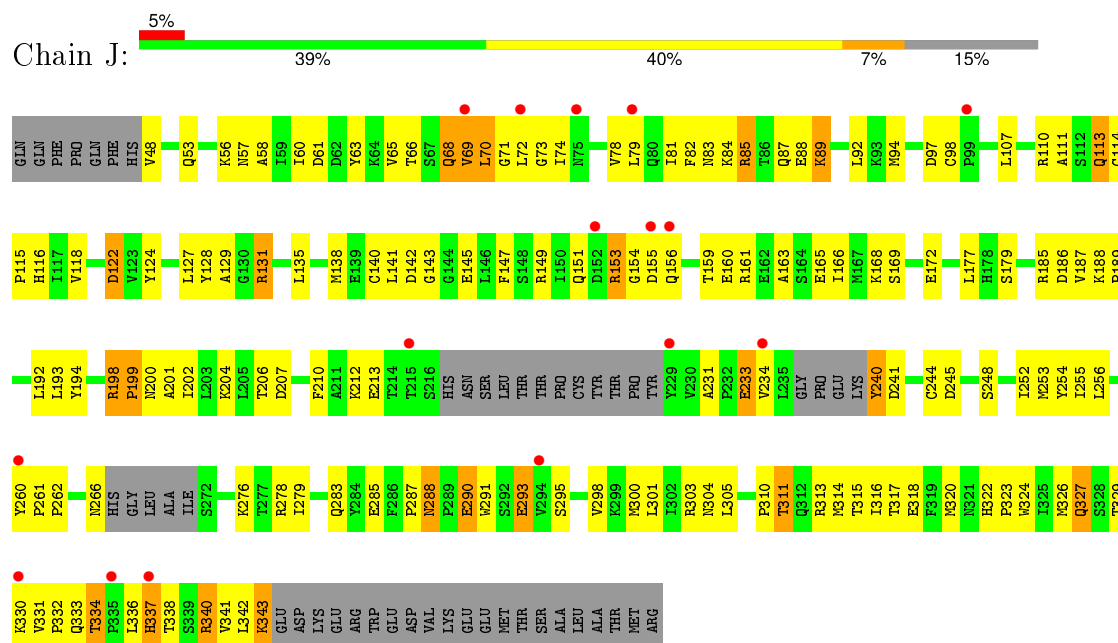




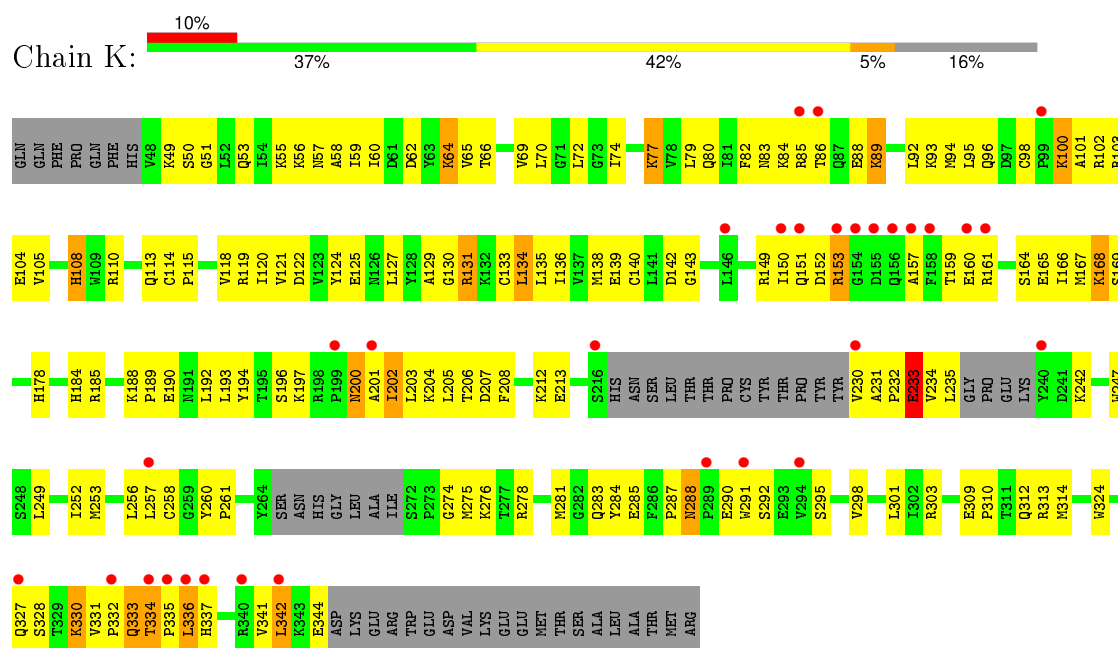
• Molecule 1: MAP kinase-activated protein kinase 2



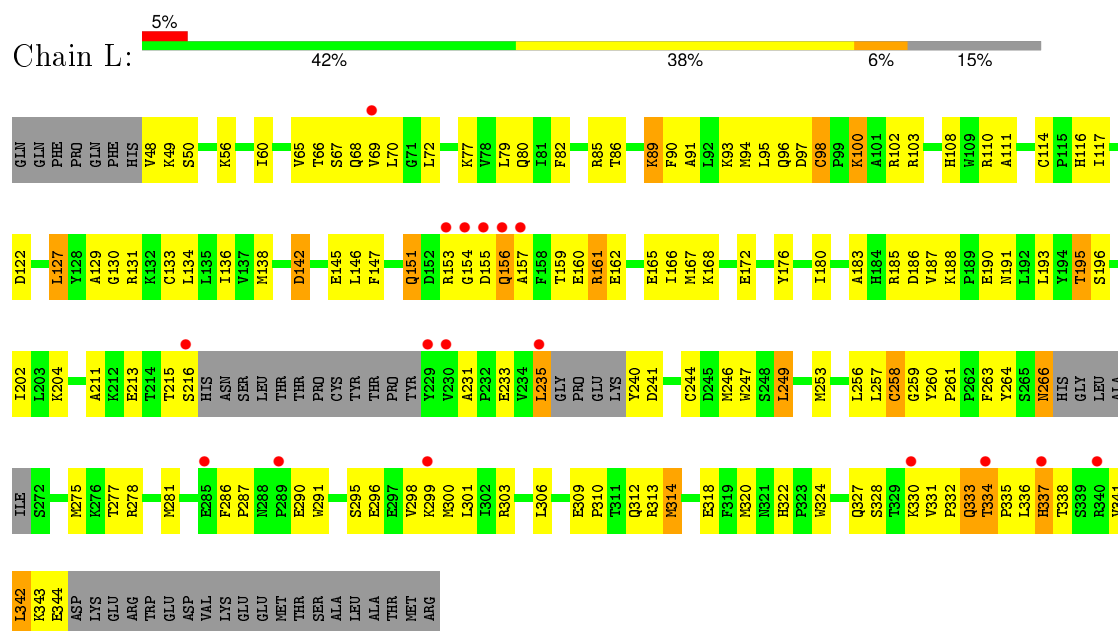
• Molecule 1: MAP kinase-activated protein kinase 2



• Molecule 1: MAP kinase-activated protein kinase 2



• Molecule 1: MAP kinase-activated protein kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	139.16Å 180.96Å 216.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.90 49.26 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.99-2.90) 99.6 (49.26-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.84 (at 2.91Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.288 , 0.335 0.212 , 0.266	Depositor DCC
R_{free} test set	6069 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 71.0	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	2 of 121020 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	27440	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 2.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, PDY

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	0.42	0/2262	0.58	0/3048
1	B	0.40	0/2284	0.57	1/3078 (0.0%)
1	C	0.36	0/2250	0.53	0/3032
1	D	0.40	0/2288	0.55	0/3081
1	E	0.42	0/2278	0.57	0/3070
1	F	0.44	0/2365	0.59	0/3191
1	G	0.36	0/2254	0.54	0/3038
1	H	0.38	0/2270	0.53	0/3058
1	I	0.41	0/2269	0.55	0/3058
1	J	0.31	0/2269	0.48	0/3058
1	K	0.33	0/2251	0.51	0/3033
1	L	0.33	0/2278	0.51	0/3070
All	All	0.38	0/27318	0.54	1/36815 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	342	LEU	CA-CB-CG	5.02	126.85	115.30

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	2218	0	2255	134	0
1	B	2240	0	2270	153	0
1	C	2207	0	2240	156	0
1	D	2244	0	2281	113	0
1	E	2234	0	2268	127	0
1	F	2316	0	2347	123	0
1	G	2210	0	2245	119	0
1	H	2226	0	2260	160	0
1	I	2225	0	2262	147	0
1	J	2225	0	2262	173	0
1	K	2208	0	2248	160	0
1	L	2234	0	2268	149	0
2	A	54	0	52	7	0
2	B	54	0	52	16	0
2	C	54	0	52	16	0
2	D	54	0	52	13	0
2	E	54	0	52	19	0
2	F	54	0	52	18	0
2	G	54	0	52	18	0
2	H	54	0	51	39	0
2	I	54	0	52	34	0
2	J	54	0	52	18	0
2	K	54	0	52	24	0
2	L	54	0	52	20	0
3	F	5	0	0	1	0
All	All	27440	0	27829	1767	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 32.

The worst 5 of 1767 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:I:2:PDY:C18	2:I:2:PDY:H14	1.44	1.48
2:L:2:PDY:C14	2:L:2:PDY:H18	1.53	1.38
1:H:264:TYR:CZ	2:H:2:PDY:H20	1.61	1.32
1:I:264:TYR:CG	2:I:2:PDY:H17	1.65	1.32
2:L:2:PDY:C18	2:L:2:PDY:H14	1.58	1.32

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	266/324 (82%)	251 (94%)	14 (5%)	1 (0%)	39	74
1	B	269/324 (83%)	252 (94%)	15 (6%)	2 (1%)	26	63
1	C	263/324 (81%)	237 (90%)	21 (8%)	5 (2%)	10	35
1	D	269/324 (83%)	249 (93%)	19 (7%)	1 (0%)	39	74
1	E	268/324 (83%)	249 (93%)	16 (6%)	3 (1%)	17	51
1	F	283/324 (87%)	262 (93%)	19 (7%)	2 (1%)	26	63
1	G	265/324 (82%)	240 (91%)	24 (9%)	1 (0%)	39	74
1	H	265/324 (82%)	242 (91%)	23 (9%)	0	100	100
1	I	267/324 (82%)	251 (94%)	16 (6%)	0	100	100
1	J	267/324 (82%)	231 (86%)	34 (13%)	2 (1%)	26	63
1	K	265/324 (82%)	238 (90%)	24 (9%)	3 (1%)	17	51
1	L	268/324 (83%)	241 (90%)	24 (9%)	3 (1%)	17	51
All	All	3215/3888 (83%)	2943 (92%)	249 (8%)	23 (1%)	26	63

5 of 23 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	154	GLY
1	A	154	GLY
1	C	328	SER
1	G	295	SER
1	L	156	GLN

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	248/293 (85%)	217 (88%)	31 (12%)	6	17
1	B	251/293 (86%)	228 (91%)	23 (9%)	11	33
1	C	247/293 (84%)	212 (86%)	35 (14%)	4	12
1	D	251/293 (86%)	223 (89%)	28 (11%)	7	22
1	E	250/293 (85%)	223 (89%)	27 (11%)	8	24
1	F	258/293 (88%)	225 (87%)	33 (13%)	5	16
1	G	247/293 (84%)	228 (92%)	19 (8%)	16	42
1	H	249/293 (85%)	224 (90%)	25 (10%)	9	28
1	I	249/293 (85%)	222 (89%)	27 (11%)	8	24
1	J	249/293 (85%)	219 (88%)	30 (12%)	6	18
1	K	247/293 (84%)	219 (89%)	28 (11%)	7	22
1	L	250/293 (85%)	226 (90%)	24 (10%)	10	31
All	All	2996/3516 (85%)	2666 (89%)	330 (11%)	8	23

5 of 330 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	127	LEU
1	G	161	ARG
1	K	336	LEU
1	F	138	MET
1	F	295	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 113 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	108	HIS
1	G	116	HIS
1	K	337	HIS
1	F	116	HIS
1	F	327	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

25 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$
2	PDY	A	1	-	23,30,30	1.79	3 (13%)	24,41,41	2.57	4 (16%)
2	PDY	A	2	-	23,30,30	1.75	3 (13%)	24,41,41	2.42	6 (25%)
2	PDY	B	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.49	2 (8%)
2	PDY	B	2	-	23,30,30	1.76	3 (13%)	24,41,41	2.42	2 (8%)
2	PDY	C	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.42	2 (8%)
2	PDY	C	2	-	23,30,30	1.75	3 (13%)	24,41,41	2.48	3 (12%)
2	PDY	D	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.57	4 (16%)
2	PDY	D	2	-	23,30,30	1.73	3 (13%)	24,41,41	2.27	4 (16%)
2	PDY	E	1	-	23,30,30	1.79	3 (13%)	24,41,41	2.43	4 (16%)
2	PDY	E	2	-	23,30,30	1.75	3 (13%)	24,41,41	2.64	5 (20%)
2	PDY	F	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.62	3 (12%)
2	PDY	F	2	-	23,30,30	1.77	3 (13%)	24,41,41	2.44	3 (12%)
3	SO4	F	3	-	4,4,4	0.20	0	6,6,6	0.21	0
2	PDY	G	1	-	23,30,30	1.78	3 (13%)	24,41,41	2.52	5 (20%)
2	PDY	G	2	-	23,30,30	1.76	3 (13%)	24,41,41	2.34	4 (16%)
2	PDY	H	1	-	23,30,30	1.79	3 (13%)	24,41,41	2.66	5 (20%)
2	PDY	H	2	-	23,30,30	1.80	3 (13%)	24,41,41	2.71	10 (41%)
2	PDY	I	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.46	4 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PDY	I	2	-	23,30,30	1.74	3 (13%)	24,41,41	2.16	8 (33%)
2	PDY	J	1	-	23,30,30	1.75	3 (13%)	24,41,41	2.29	3 (12%)
2	PDY	J	2	-	23,30,30	1.77	3 (13%)	24,41,41	2.38	3 (12%)
2	PDY	K	1	-	23,30,30	1.77	3 (13%)	24,41,41	2.41	5 (20%)
2	PDY	K	2	-	23,30,30	1.76	3 (13%)	24,41,41	2.50	5 (20%)
2	PDY	L	1	-	23,30,30	1.79	3 (13%)	24,41,41	2.46	4 (16%)
2	PDY	L	2	-	23,30,30	1.74	3 (13%)	24,41,41	2.03	4 (16%)

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
2	PDY	A	1	-	-	1/11/19/19	0/4/4/4
2	PDY	A	2	-	-	1/11/19/19	0/4/4/4
2	PDY	B	1	-	-	1/11/19/19	0/4/4/4
2	PDY	B	2	-	-	1/11/19/19	0/4/4/4
2	PDY	C	1	-	-	1/11/19/19	0/4/4/4
2	PDY	C	2	-	-	1/11/19/19	0/4/4/4
2	PDY	D	1	-	-	1/11/19/19	0/4/4/4
2	PDY	D	2	-	-	1/11/19/19	0/4/4/4
2	PDY	E	1	-	-	1/11/19/19	0/4/4/4
2	PDY	E	2	-	-	1/11/19/19	0/4/4/4
2	PDY	F	1	-	-	1/11/19/19	0/4/4/4
2	PDY	F	2	-	-	1/11/19/19	0/4/4/4
3	SO4	F	3	-	-	0/0/0/0	0/0/0/0
2	PDY	G	1	-	-	1/11/19/19	0/4/4/4
2	PDY	G	2	-	-	1/11/19/19	0/4/4/4
2	PDY	H	1	-	-	1/11/19/19	0/4/4/4
2	PDY	H	2	-	-	1/11/19/19	0/4/4/4
2	PDY	I	1	-	-	1/11/19/19	0/4/4/4
2	PDY	I	2	-	-	1/11/19/19	0/4/4/4
2	PDY	J	1	-	-	1/11/19/19	0/4/4/4
2	PDY	J	2	-	-	1/11/19/19	0/4/4/4
2	PDY	K	1	-	-	1/11/19/19	0/4/4/4
2	PDY	K	2	-	-	1/11/19/19	0/4/4/4
2	PDY	L	1	-	-	1/11/19/19	0/4/4/4
2	PDY	L	2	-	-	1/11/19/19	0/4/4/4

The worst 5 of 72 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1	PDY	C13-N9	-3.76	1.32	1.40
2	F	1	PDY	C13-N9	-3.66	1.32	1.40
2	K	1	PDY	C13-N9	-3.64	1.33	1.40
2	E	1	PDY	C13-N9	-3.63	1.33	1.40
2	A	1	PDY	C13-N9	-3.60	1.33	1.40

The worst 5 of 102 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	PDY	C24-O21-C16	-6.23	105.36	117.92
2	F	1	PDY	C24-O21-C16	-4.70	108.44	117.92
2	F	1	PDY	C8-C11-N7	-4.66	100.08	111.38
2	B	1	PDY	C8-C11-N7	-4.62	100.18	111.38
2	L	1	PDY	C8-C11-N7	-4.59	100.25	111.38

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	2	PDY	C3-C1-N9-C13
2	C	2	PDY	C3-C1-N9-C13
2	A	1	PDY	C3-C1-N9-C13
2	B	1	PDY	C3-C1-N9-C13
2	G	2	PDY	C3-C1-N9-C13

There are no ring outliers.

25 monomers are involved in 243 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	PDY	4	0
2	A	2	PDY	3	0
2	B	1	PDY	8	0
2	B	2	PDY	8	0
2	C	1	PDY	6	0
2	C	2	PDY	10	0
2	D	1	PDY	8	0
2	D	2	PDY	5	0
2	E	1	PDY	10	0
2	E	2	PDY	9	0
2	F	1	PDY	6	0
2	F	2	PDY	12	0
3	F	3	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	PDY	9	0
2	G	2	PDY	9	0
2	H	1	PDY	13	0
2	H	2	PDY	26	0
2	I	1	PDY	9	0
2	I	2	PDY	25	0
2	J	1	PDY	10	0
2	J	2	PDY	8	0
2	K	1	PDY	14	0
2	K	2	PDY	10	0
2	L	1	PDY	9	0
2	L	2	PDY	11	0

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	274/324 (84%)	0.16	9 (3%) 50 42	8, 24, 56, 81	0
1	B	277/324 (85%)	0.22	5 (1%) 71 68	7, 29, 71, 80	0
1	C	273/324 (84%)	0.31	12 (4%) 38 32	15, 40, 73, 83	0
1	D	277/324 (85%)	0.28	9 (3%) 51 43	8, 32, 70, 81	0
1	E	276/324 (85%)	0.23	9 (3%) 50 42	10, 28, 58, 76	0
1	F	287/324 (88%)	0.17	14 (4%) 33 27	4, 25, 70, 86	0
1	G	273/324 (84%)	0.36	11 (4%) 42 35	16, 38, 67, 79	0
1	H	275/324 (84%)	0.29	13 (4%) 35 29	13, 39, 68, 84	0
1	I	275/324 (84%)	0.19	8 (2%) 55 49	14, 30, 66, 97	0
1	J	275/324 (84%)	0.53	16 (5%) 26 20	27, 52, 81, 100	0
1	K	273/324 (84%)	0.61	31 (11%) 7 4	25, 49, 76, 98	0
1	L	276/324 (85%)	0.45	17 (6%) 24 17	15, 46, 83, 110	0
All	All	3311/3888 (85%)	0.32	154 (4%) 35 29	4, 37, 73, 110	0

The worst 5 of 154 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	216	SER	6.9
1	J	155	ASP	6.8
1	I	229	TYR	6.0
1	K	157	ALA	6.0
1	H	155	ASP	6.0

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
2	PDY	D	2	27/27	0.69	0.42	7.79	65,74,80,81	0
2	PDY	L	2	27/27	0.82	0.43	5.01	71,75,77,77	0
2	PDY	E	2	27/27	0.82	0.43	3.98	60,67,69,69	0
2	PDY	G	2	27/27	0.90	0.37	2.95	52,56,64,64	0
2	PDY	A	2	27/27	0.82	0.37	2.61	60,65,69,71	0
2	PDY	I	2	27/27	0.80	0.38	2.24	53,61,74,75	0
2	PDY	E	1	27/27	0.97	0.28	2.24	7,23,37,38	0
3	SO4	F	3	5/5	0.87	0.26	1.30	61,62,64,65	0
2	PDY	K	2	27/27	0.85	0.40	1.24	61,67,79,80	0
2	PDY	C	2	27/27	0.91	0.25	0.89	48,50,62,63	0
2	PDY	H	2	27/27	0.90	0.33	0.83	52,55,60,60	0
2	PDY	J	2	27/27	0.90	0.33	0.79	59,62,67,68	0
2	PDY	K	1	27/27	0.94	0.24	0.74	27,44,54,55	0
2	PDY	B	2	27/27	0.96	0.21	0.66	20,29,34,35	0
2	PDY	C	1	27/27	0.92	0.23	0.49	18,32,56,58	0
2	PDY	A	1	27/27	0.97	0.22	0.46	2,11,31,32	0
2	PDY	H	1	27/27	0.96	0.22	0.43	22,31,54,55	0
2	PDY	J	1	27/27	0.96	0.23	0.38	31,39,54,55	0
2	PDY	D	1	27/27	0.95	0.23	0.17	16,23,43,44	0
2	PDY	L	1	27/27	0.95	0.21	0.07	30,34,47,50	0
2	PDY	F	2	27/27	0.96	0.17	-0.37	20,24,28,29	0
2	PDY	I	1	27/27	0.96	0.19	-0.39	5,19,43,45	0
2	PDY	B	1	27/27	0.96	0.21	-0.40	6,16,37,38	0
2	PDY	F	1	27/27	0.96	0.18	-0.45	5,15,32,34	0
2	PDY	G	1	27/27	0.96	0.16	-0.58	4,20,38,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.