



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:08 PM UTC

PDB ID : 6MYN / pdb_00006myn
Title : Crystal structure of murine NF-kappaB inducing kinase (NIK) bound to inhibitor R7
Authors : Harris, S.F.; Smith, M.; Barker, J.
Deposited on : 2018-11-01
Resolution : 2.74 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)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

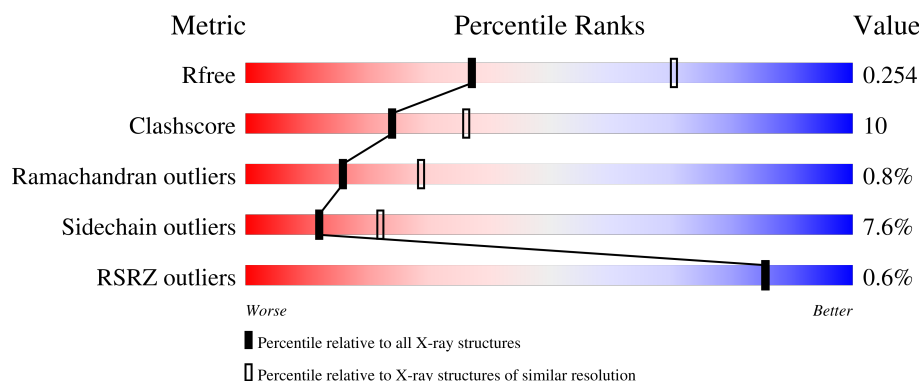
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	347	 68% 21% • 7%
1	B	347	 69% 22% • 5%

2 Entry composition [i](#)

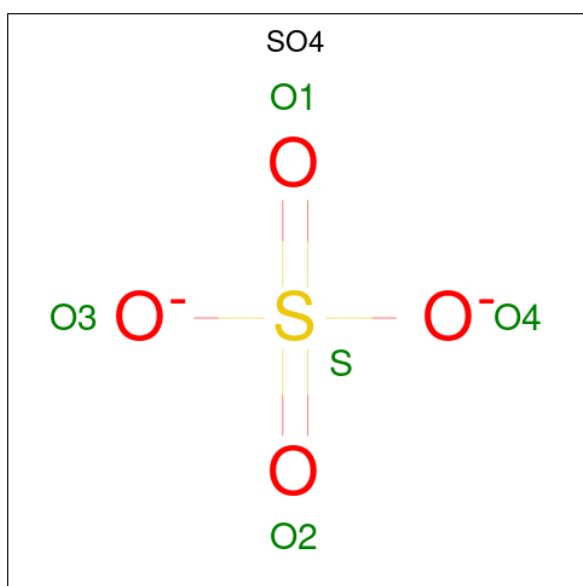
There are 4 unique types of molecules in this entry. The entry contains 5280 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Mitogen-activated protein kinase kinase kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	2	0
			2525	1598	449	457	21			
1	B	328	Total	C	N	O	S	0	1	0
			2557	1618	454	464	21			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



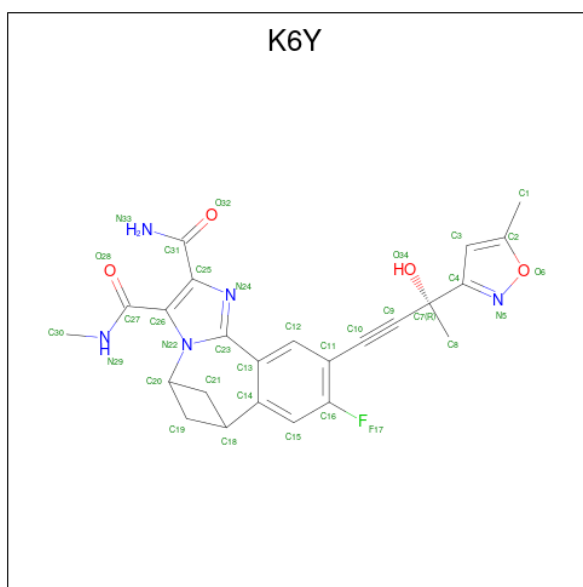
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is (5s,7s)-9-fluoro-10-[(3R)-3-hydroxy-3-(5-methyl-1,2-oxazol-3-yl)but-1-yn-1-yl]-N 3 -methyl-6,7-dihydro-5H-5,7-methanoimidazo[2,1-a][2]benzazepine-2,3-dicarboxamide (CCD ID: K6Y) (formula: C₂₄H₂₂FN₅O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			34	24	1	5	4		
3	B	1	Total	C	F	N	O	0	0
			34	24	1	5	4		

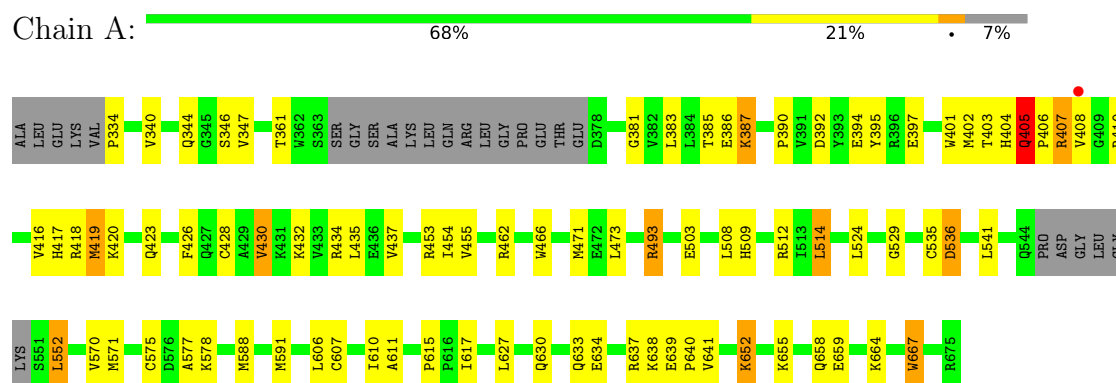
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		
4	B	46	Total	O	0	0
			46	46		

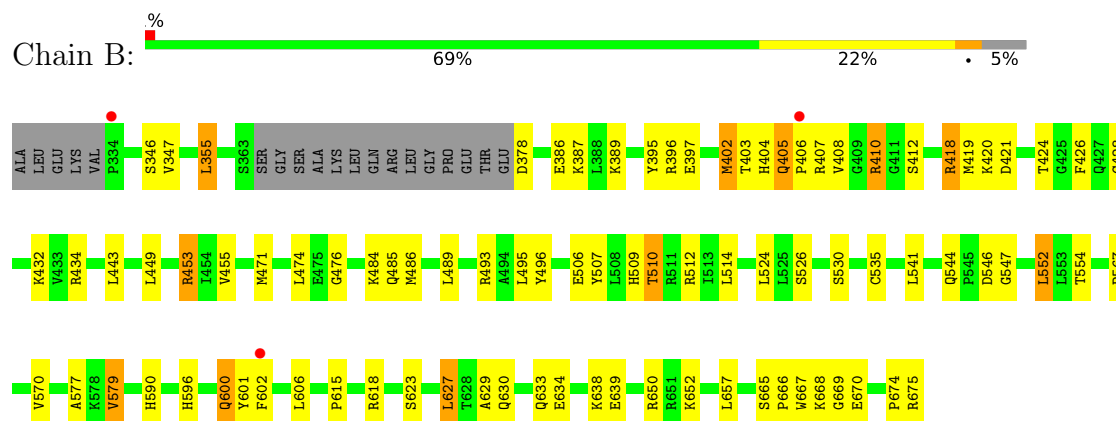
3 Residue-property plots

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: Mitogen-activated protein kinase kinase kinase 14



- Molecule 1: Mitogen-activated protein kinase kinase kinase 14



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	143.22Å 143.22Å 46.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.19 – 2.74 42.19 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.19-2.74) 99.8 (42.19-2.74)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.12-2829_final	Depositor
R, R_{free}	0.179 , 0.255 0.180 , 0.254	Depositor DCC
R_{free} test set	2604 reflections (10.34%)	wwPDB-VP
Wilson B-factor (Å ²)	58.3	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5280	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: SO4, K6Y

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	0.45	0/2584	0.64	0/3493
1	B	0.44	0/2618	0.65	0/3540
All	All	0.44	0/5202	0.64	0/7033

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	2525	0	2519	52	0
1	B	2557	0	2557	54	0
2	A	10	0	0	1	0
2	B	15	0	0	1	0
3	A	34	0	0	3	0
3	B	34	0	0	1	0
4	A	59	0	0	4	0
4	B	46	0	0	1	0
All	All	5280	0	5076	103	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 10.

All (103) 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:392:ASP:OD2	1:B:407:ARG:NH1	1.86	1.08
1:A:664:LYS:NZ	4:A:801:HOH:O	2.05	0.90
1:B:615:PRO:HG3	1:B:638:LYS:HG3	1.59	0.83
1:A:536:ASP:OD2	3:A:703:K6Y:F17	1.94	0.75
1:A:634:GLU:OE1	1:A:652:LYS:NZ	2.19	0.73
1:B:403:THR:HG21	1:B:407:ARG:HH21	1.53	0.72
1:B:405:GLN:HB2	1:B:406:PRO:HD3	1.73	0.70
1:A:453:ARG:HG3	1:A:503:GLU:HB3	1.74	0.69
1:A:615:PRO:HG3	1:A:638:LYS:HG3	1.75	0.68
1:A:571:MET:HE3	1:A:611:ALA:HB1	1.78	0.66
1:B:600:GLN:HG3	1:B:675:ARG:HH21	1.62	0.64
1:A:658[A]:GLN:NE2	4:A:802:HOH:O	2.15	0.62
1:B:405:GLN:HB2	1:B:406:PRO:CD	2.29	0.62
1:B:506:GLU:O	1:B:510:THR:HG23	1.98	0.62
1:A:405:GLN:HB3	1:A:406:PRO:HD2	1.83	0.61
1:B:509:HIS:CE1	1:B:577:ALA:HB1	2.37	0.60
1:B:453:ARG:HH11	1:B:453:ARG:HA	1.67	0.59
1:B:666:PRO:HG2	1:B:669:GLY:HA2	1.85	0.59
1:A:347:VAL:HG11	1:A:395:TYR:HB2	1.85	0.58
1:B:489:LEU:O	4:B:801:HOH:O	2.17	0.57
1:A:404:HIS:CD2	1:A:405:GLN:HB2	2.40	0.57
1:B:455:VAL:HG21	1:B:535:CYS:HB3	1.86	0.57
1:A:655:LYS:NZ	1:A:659:GLU:OE2	2.36	0.56
1:A:432:LYS:HD3	1:A:466:TRP:CD2	2.41	0.56
1:B:386:GLU:HG3	1:B:389:LYS:HE2	1.87	0.55
1:B:410:ARG:NH1	1:B:418:ARG:NH2	2.54	0.55
1:A:455:VAL:HG21	1:A:535:CYS:HB3	1.88	0.55
1:B:600:GLN:HG3	1:B:675:ARG:NH2	2.21	0.55
1:B:404:HIS:HB2	1:B:418:ARG:HG2	1.89	0.55
1:B:397:GLU:OE1	1:B:432:LYS:NZ	2.41	0.53
1:A:428:CYS:HB2	1:A:471:MET:O	2.09	0.52
1:A:435:LEU:HD22	1:A:462:ARG:HD2	1.92	0.52
1:B:541:LEU:HD22	1:B:552:LEU:HD11	1.91	0.52
1:A:346:SER:HB3	1:B:434:ARG:NH1	2.25	0.52
1:B:453:ARG:NH2	2:B:702:SO4:O3	2.43	0.52
1:A:552:LEU:HD13	1:A:575:CYS:HB2	1.92	0.51
1:A:514:LEU:HD23	1:A:541:LEU:HB3	1.91	0.51
1:B:514:LEU:HD11	1:B:579:VAL:HG22	1.92	0.51
1:B:601:TYR:HB2	1:B:602:PHE:CD2	2.46	0.51
1:B:449:LEU:HD22	1:B:507:TYR:HE2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:LYS:HB2	1:A:426:PHE:CG	2.47	0.49
1:B:493:ARG:NH1	1:B:665:SER:O	2.45	0.49
1:A:386:GLU:HG2	1:B:484:LYS:HD3	1.94	0.49
1:B:378:ASP:CG	1:B:396:ARG:HH22	2.20	0.49
1:A:397:GLU:OE2	1:A:407:ARG:NH2	2.46	0.48
1:A:529:GLY:HA3	1:A:667:TRP:CD2	2.48	0.48
1:A:401:TRP:CZ3	1:A:430:VAL:HG21	2.48	0.48
1:A:637:ARG:NH1	2:A:701:SO4:O2	2.43	0.48
1:B:486:MET:HE3	1:B:668:LYS:HA	1.94	0.47
1:B:544:GLN:O	1:B:547:GLY:N	2.31	0.47
1:A:405:GLN:HB3	1:A:406:PRO:CD	2.44	0.47
1:A:509:HIS:CE1	1:A:577:ALA:HB1	2.50	0.47
1:A:434:ARG:HB2	1:A:437:VAL:HG22	1.96	0.47
1:A:512:ARG:NH1	4:A:812:HOH:O	2.48	0.47
1:A:578:LYS:NZ	1:A:640:PRO:O	2.47	0.46
1:A:410:ARG:HD2	3:A:703:K6Y:O28	2.15	0.46
1:B:667:TRP:CZ2	1:B:668:LYS:HE2	2.50	0.46
1:A:340:VAL:O	1:A:344:GLN:HG3	2.17	0.45
1:A:655:LYS:O	1:A:659:GLU:HG3	2.17	0.45
1:B:476:GLY:HA3	1:B:526:SER:O	2.17	0.45
1:B:514:LEU:HD13	1:B:577:ALA:HA	1.99	0.45
1:B:485:GLN:HB3	1:B:486:MET:HE2	1.99	0.45
1:B:402:MET:HB3	1:B:402:MET:HE3	1.69	0.45
1:B:552:LEU:HD12	1:B:552:LEU:HA	1.64	0.45
1:A:617:ILE:HG21	1:A:633:GLN:HG2	1.98	0.45
1:A:347:VAL:HA	1:A:381:GLY:O	2.17	0.45
1:A:334:PRO:HB2	1:B:570:VAL:O	2.18	0.44
1:B:634:GLU:HG3	1:B:652:LYS:NZ	2.33	0.44
1:A:383:LEU:HD12	1:A:390:PRO:HD3	2.00	0.43
1:A:588:MET:HA	1:A:591:MET:HE3	2.00	0.43
1:A:418:ARG:NH1	1:A:473:LEU:HD21	2.32	0.43
1:B:567:PRO:HG2	1:B:638:LYS:HG2	1.99	0.43
1:A:536:ASP:HA	3:A:703:K6Y:C9	2.48	0.43
1:B:590:HIS:CE1	1:B:596:HIS:CD2	3.07	0.43
1:A:394:GLU:HG2	4:A:805:HOH:O	2.19	0.42
1:B:509:HIS:CG	1:B:577:ALA:HB1	2.53	0.42
1:B:590:HIS:CE1	1:B:596:HIS:HD2	2.37	0.42
1:B:627:LEU:HD12	1:B:627:LEU:HA	1.56	0.42
1:B:495:LEU:HB2	1:B:657:LEU:HD13	2.01	0.42
3:B:704:K6Y:O32	3:B:704:K6Y:N29	2.53	0.42
1:B:355:LEU:HD11	1:B:443:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:419:MET:HE3	1:B:420:LYS:C	2.45	0.42
1:B:421:ASP:OD2	1:B:424:THR:HG23	2.20	0.42
1:B:509:HIS:CD2	1:B:577:ALA:HB1	2.55	0.42
1:A:386:GLU:H	1:A:386:GLU:CD	2.26	0.42
1:A:407:ARG:HH11	1:A:417:HIS:CE1	2.38	0.42
1:B:355:LEU:HD12	1:B:355:LEU:O	2.19	0.42
1:B:387:LYS:HB3	1:B:426:PHE:CG	2.54	0.42
1:B:474:LEU:HD13	1:B:524:LEU:HB2	2.02	0.42
1:A:454:ILE:HD11	1:A:508:LEU:HG	2.02	0.41
1:B:347:VAL:HG11	1:B:395:TYR:HB2	2.01	0.41
1:A:627:LEU:HD11	1:A:652:LYS:HD2	2.01	0.41
1:A:630:GLN:O	1:A:634:GLU:HG3	2.20	0.41
1:B:428:CYS:HB2	1:B:471:MET:O	2.20	0.41
1:A:493:ARG:NH1	1:A:667:TRP:HA	2.36	0.41
1:B:544:GLN:O	1:B:546:ASP:N	2.54	0.41
1:A:419:MET:HE3	1:A:420:LYS:C	2.46	0.41
1:A:419:MET:HG2	1:A:430:VAL:HG22	2.03	0.41
1:A:570:VAL:HG11	1:A:607:CYS:O	2.21	0.41
1:B:496:TYR:OH	1:B:530:SER:O	2.36	0.40
1:A:416:VAL:HA	1:A:430:VAL:O	2.22	0.40
1:B:629:ALA:O	1:B:633:GLN:HG3	2.22	0.40
1:A:455:VAL:HG21	1:A:524:LEU:HD12	2.04	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	318/347 (92%)	298 (94%)	18 (6%)	2 (1%)	21	36
1	B	325/347 (94%)	303 (93%)	19 (6%)	3 (1%)	14	26
All	All	643/694 (93%)	601 (94%)	37 (6%)	5 (1%)	16	29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	405	GLN
1	A	405	GLN
1	B	412	SER
1	B	674	PRO
1	A	536	ASP

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	271/289 (94%)	251 (93%)	20 (7%)	13	23
1	B	275/289 (95%)	254 (92%)	21 (8%)	12	22
All	All	546/578 (94%)	505 (92%)	41 (8%)	12	23

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	361	THR
1	A	385	THR
1	A	387	LYS
1	A	402	MET
1	A	403	THR
1	A	405	GLN
1	A	407	ARG
1	A	408	VAL
1	A	419	MET
1	A	423	GLN
1	A	430	VAL
1	A	493	ARG
1	A	514	LEU
1	A	552	LEU
1	A	606	LEU
1	A	610	ILE
1	A	639	GLU
1	A	641	VAL

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Mol	Chain	Res	Type
1	A	652	LYS
1	A	667	TRP
1	B	346	SER
1	B	355	LEU
1	B	402	MET
1	B	408	VAL
1	B	410	ARG
1	B	418	ARG
1	B	453	ARG
1	B	510	THR
1	B	512	ARG
1	B	552	LEU
1	B	554	THR
1	B	579	VAL
1	B	600	GLN
1	B	606	LEU
1	B	618	ARG
1	B	623	SER
1	B	627	LEU
1	B	630	GLN
1	B	639	GLU
1	B	650	ARG
1	B	670	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	351	GLN
1	A	405	GLN
1	A	423	GLN
1	B	596	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	SO4	A	702	-	4,4,4	0.25	0	6,6,6	0.28	0
2	SO4	B	702	-	4,4,4	0.30	0	6,6,6	0.25	0
2	SO4	A	701	-	4,4,4	0.38	0	6,6,6	0.37	0
3	K6Y	A	703	-	36,38,38	0.58	1 (2%)	41,59,59	1.50	9 (21%)
2	SO4	B	701	-	4,4,4	0.24	0	6,6,6	0.37	0
3	K6Y	B	704	-	36,38,38	0.80	2 (5%)	41,59,59	1.43	7 (17%)
2	SO4	B	703	-	4,4,4	0.38	0	6,6,6	0.41	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
3	K6Y	A	703	-	-	1/18/42/42	0/4/5/5
3	K6Y	B	704	-	-	2/18/42/42	0/4/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	704	K6Y	O34-C7	-3.32	1.40	1.43
3	A	703	K6Y	C21-C20	-2.32	1.50	1.53
3	B	704	K6Y	C8-C7	-2.06	1.49	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	703	K6Y	C21-C18-C14	-4.22	109.95	122.04
3	B	704	K6Y	C14-C18-C19	-3.88	110.92	122.04
3	B	704	K6Y	C21-C18-C14	-3.70	111.45	122.04
3	B	704	K6Y	C27-C26-N22	-3.66	118.30	122.33
3	A	703	K6Y	C27-C26-N22	-3.55	118.42	122.33
3	A	703	K6Y	C18-C21-C20	3.32	92.43	83.51
3	A	703	K6Y	C14-C18-C19	-3.13	113.08	122.04
3	B	704	K6Y	C18-C21-C20	2.83	91.11	83.51
3	B	704	K6Y	O34-C7-C8	-2.73	99.88	107.53
3	A	703	K6Y	C12-C13-C23	-2.69	111.22	119.55
3	A	703	K6Y	C18-C19-C20	2.68	90.71	83.51
3	B	704	K6Y	C18-C19-C20	2.50	90.24	83.51
3	A	703	K6Y	C16-C11-C10	2.47	126.20	120.75
3	A	703	K6Y	C14-C13-C23	2.11	128.01	123.28
3	A	703	K6Y	C13-C23-N24	-2.05	118.60	122.58
3	B	704	K6Y	C12-C13-C23	-2.04	113.24	119.55

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	704	K6Y	C9-C10-C11-C16
3	A	703	K6Y	C11-C10-C9-C7
3	B	704	K6Y	C9-C10-C11-C12

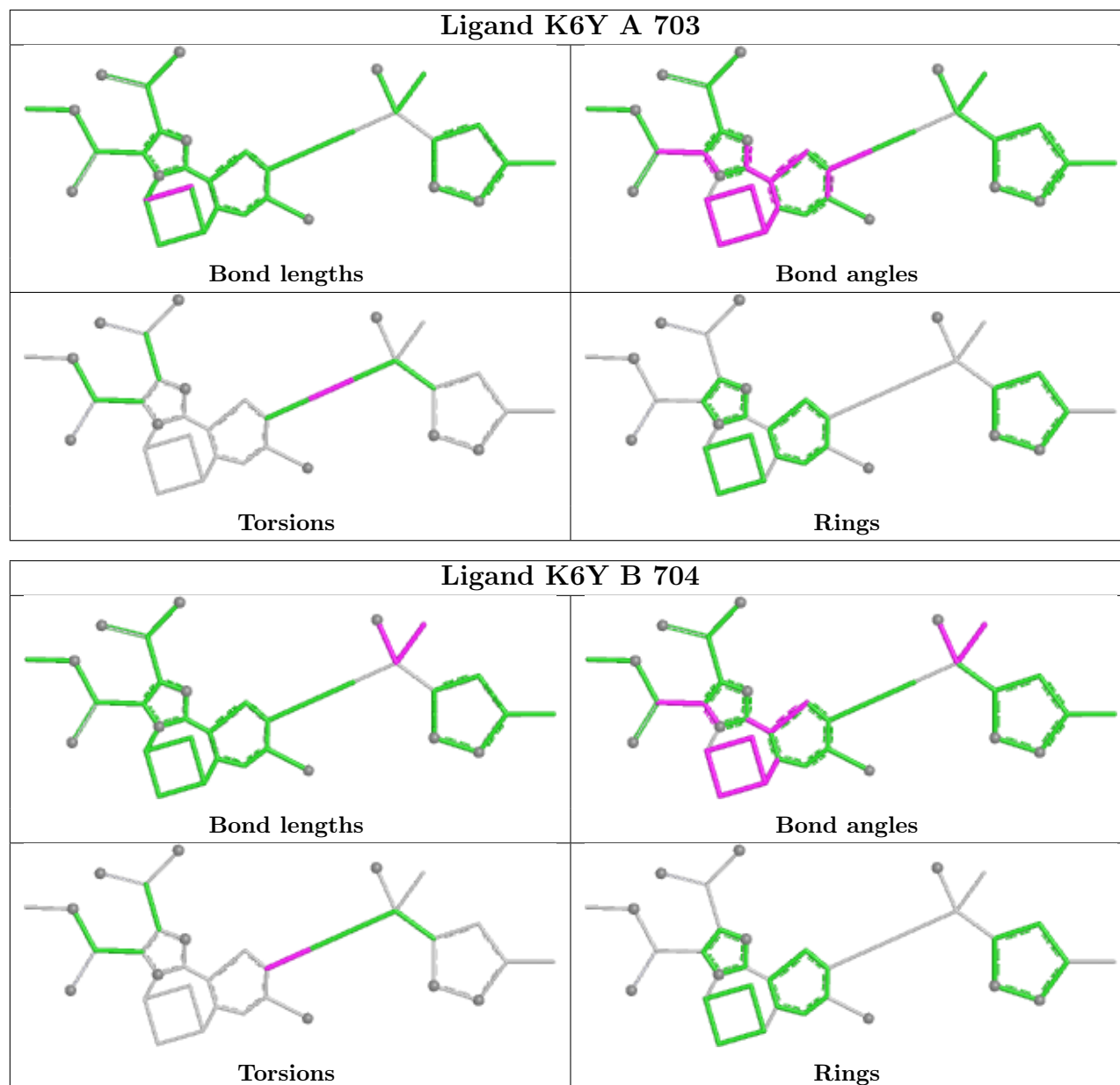
There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	702	SO4	1	0
2	A	701	SO4	1	0
3	A	703	K6Y	3	0
3	B	704	K6Y	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	322/347 (92%)	-0.35	1 (0%) 90 91	23, 48, 83, 112	2 (0%)
1	B	328/347 (94%)	-0.18	3 (0%) 81 81	25, 53, 89, 103	1 (0%)
All	All	650/694 (93%)	-0.27	4 (0%) 85 85	23, 50, 88, 112	3 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	408	VAL	3.7
1	B	334	PRO	3.1
1	B	406	PRO	2.6
1	B	602	PHE	2.6

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 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.

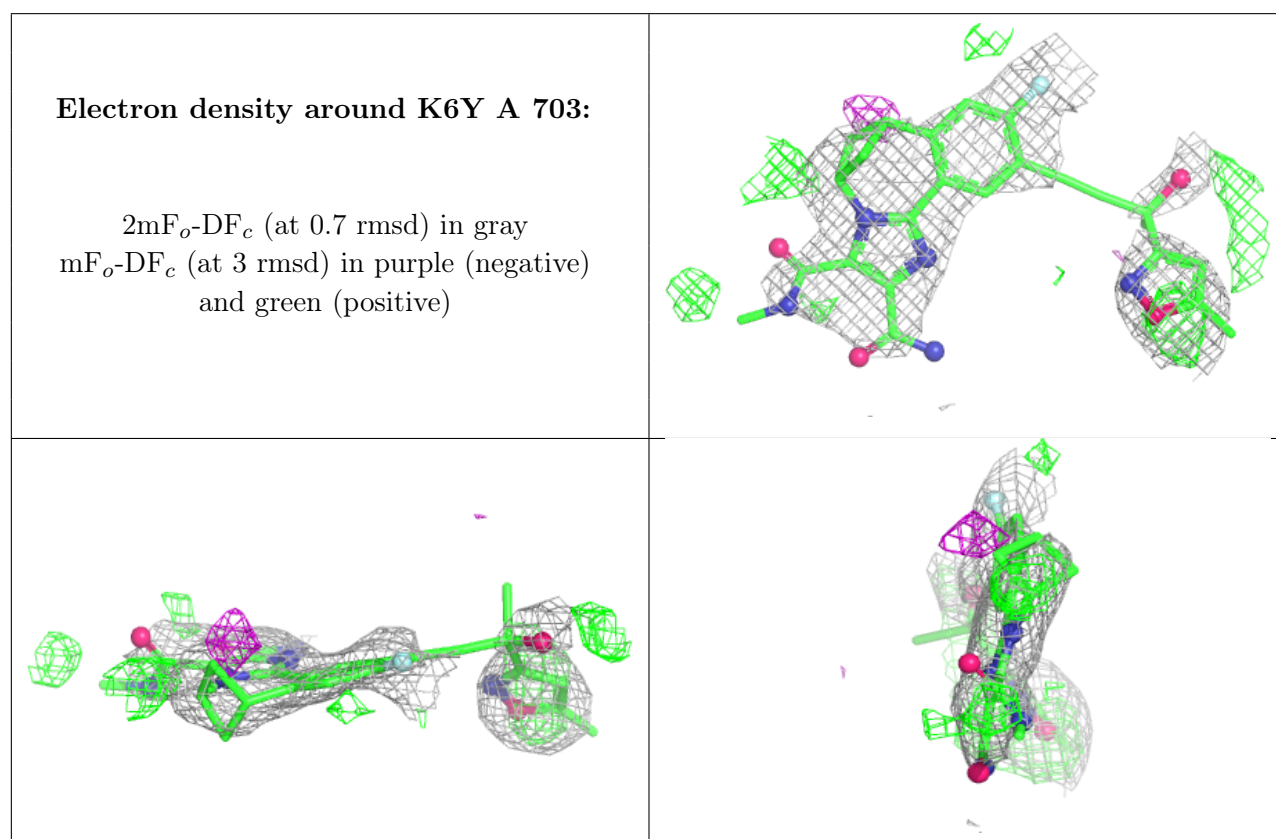
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	B	702	5/5	0.69	0.11	97,100,129,134	0

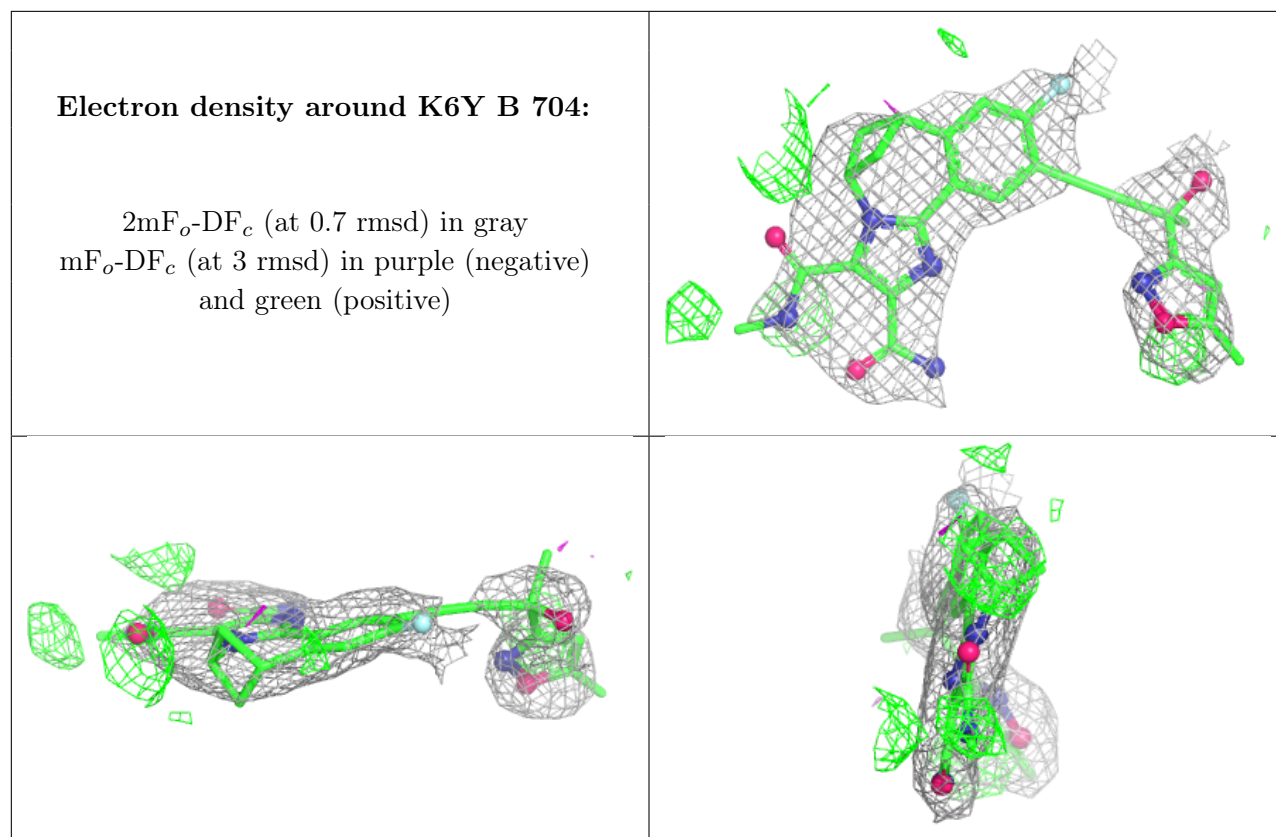
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	701	5/5	0.87	0.09	67,71,90,97	0
3	K6Y	A	703	34/34	0.88	0.21	34,48,58,68	34
2	SO4	B	703	5/5	0.90	0.20	62,67,73,93	0
3	K6Y	B	704	34/34	0.90	0.17	34,50,59,69	34
2	SO4	A	702	5/5	0.91	0.11	66,78,89,90	0
2	SO4	B	701	5/5	0.94	0.07	74,74,87,95	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.