



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:13 PM UTC

PDB ID : 2F2C / pdb_00002f2c
Title : X-ray structure of human CDK6-Vcyclinwith the inhibitor aminopurvalanol
Authors : Schulze-Gahmen, U.; Lu, H.
Deposited on : 2005-11-16
Resolution : 2.80 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
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

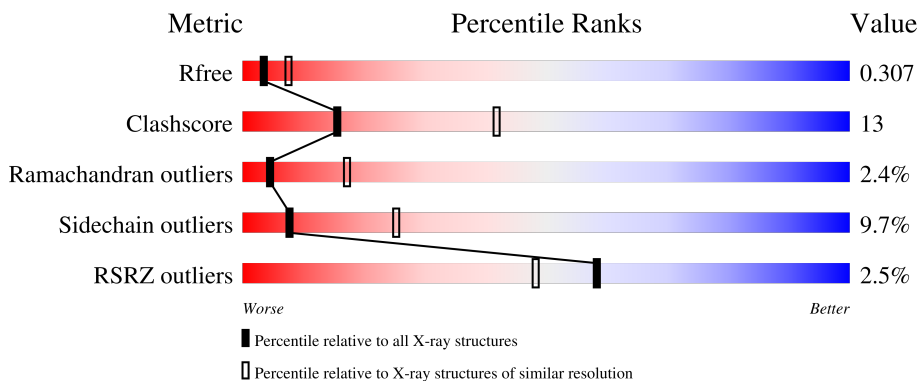
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	254	
2	B	308	

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
3	SO4	B	402	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4059 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.

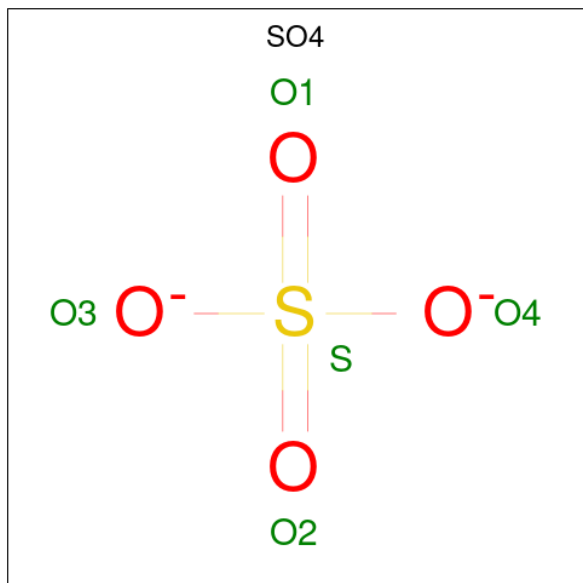
- Molecule 1 is a protein called Cyclin homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	244	1903	1228	306	358	11	0	0	0

- Molecule 2 is a protein called Cell division protein kinase 6.

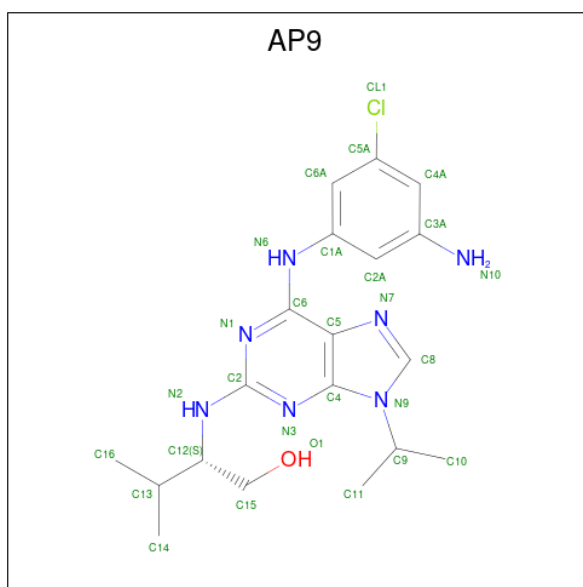
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	280	2119	1370	351	391	7	0	0	0

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



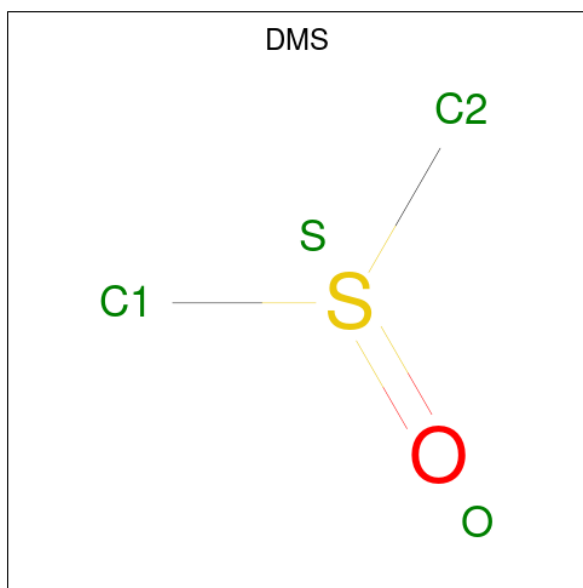
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
3	B	1	5	4	1	0	0

- Molecule 4 is (2S)-2-({6-[(3-AMINO-5-CHLOROPHENYL)AMINO]-9-ISOPROPYL-9H-PURIN-2-YL}AMINO)-3-METHYLBUTAN-1-OL (CCD ID: AP9) (formula: C₁₉H₂₆ClN₇O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Cl	N			O
4	B	1	28	19	1	7	1	0	0

- Molecule 5 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).

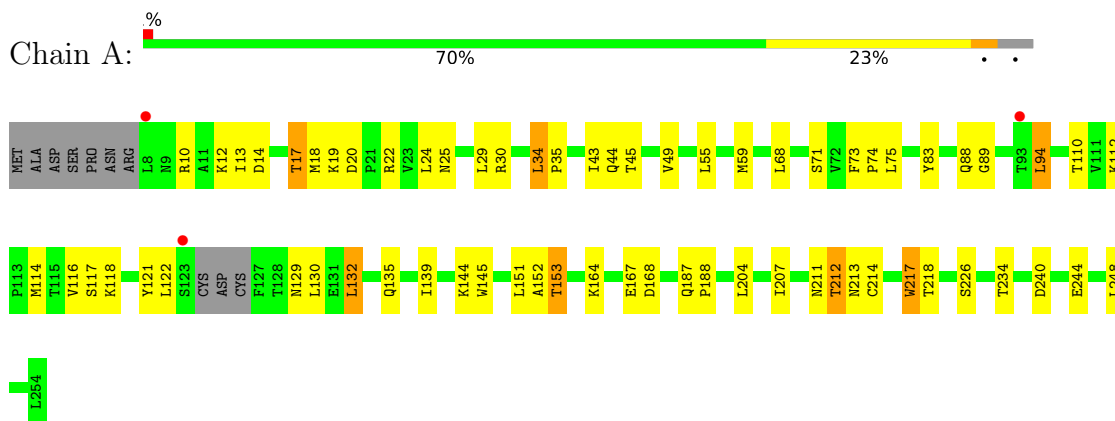


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	O			S
5	B	1	4	2	1	1	0	0

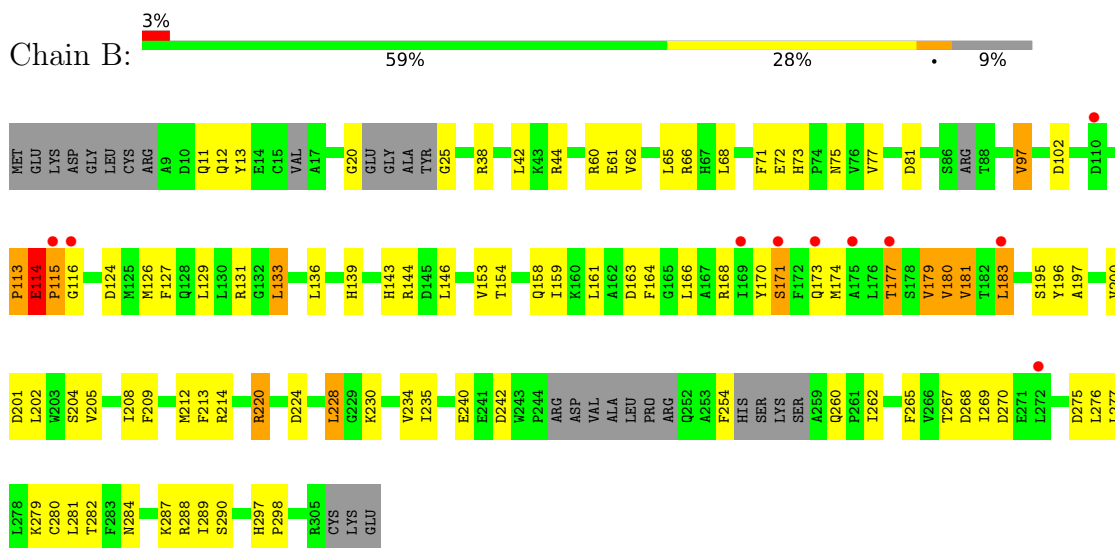
3 Residue-property plots [i](#)

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: Cyclin homolog



- Molecule 2: Cell division protein kinase 6



4 Data and refinement statistics i

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	71.52Å 71.52Å 449.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.9 (20.00-2.80) 96.5 (20.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.77Å)	Xtrriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.238 , 0.301 (Not available) , 0.307	Depositor DCC
R_{free} test set	903 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	87.0	Xtrriage
Anisotropy	0.212	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 92.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4059	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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: AP9, SO4, DMS

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	0.69	0/1934	1.00	0/2632
2	B	0.69	2/2162 (0.1%)	0.96	1/2943 (0.0%)
All	All	0.69	2/4096 (0.0%)	0.98	1/5575 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	114	GLU	CD-OE2	10.05	1.44	1.25
2	B	114	GLU	CD-OE1	7.87	1.40	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	115	PRO	N-CD-CG	-5.25	95.32	103.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	114	GLU	Peptide

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	1903	0	1961	45	0
2	B	2119	0	2028	67	0
3	B	5	0	0	2	0
4	B	28	0	26	3	0
5	B	4	0	6	2	0
All	All	4059	0	4021	108	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 13.

All (108) 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:168:ARG:HE	2:B:174:MET:HE1	1.36	0.90
2:B:202:LEU:HD13	2:B:289:ILE:O	1.71	0.88
1:A:55:LEU:HD22	1:A:94:LEU:HD22	1.57	0.85
2:B:60:ARG:NH1	2:B:168:ARG:HD2	1.97	0.80
1:A:68:LEU:HD11	1:A:114:MET:CE	2.17	0.75
1:A:240:ASP:O	1:A:244:GLU:HG2	1.87	0.74
2:B:168:ARG:NE	2:B:174:MET:HE1	2.03	0.73
2:B:114:GLU:O	2:B:116:GLY:N	2.22	0.73
1:A:116:VAL:HG23	1:A:132:LEU:HD12	1.68	0.73
2:B:60:ARG:HH12	2:B:168:ARG:HD2	1.55	0.72
1:A:68:LEU:HD11	1:A:114:MET:HE2	1.73	0.71
1:A:12:LYS:HD2	1:A:13:ILE:H	1.58	0.67
2:B:126:MET:HE1	2:B:205:VAL:HG13	1.76	0.67
2:B:267:THR:O	2:B:268:ASP:HB2	1.94	0.67
2:B:235:ILE:HD13	2:B:281:LEU:O	1.96	0.66
2:B:114:GLU:C	2:B:116:GLY:H	2.04	0.66
1:A:30:ARG:HD3	2:B:139:HIS:ND1	2.13	0.64
1:A:44:GLN:HE22	1:A:89:GLY:H	1.46	0.64
1:A:34:LEU:HB2	1:A:35:PRO:CD	2.28	0.63
2:B:20:GLY:HA3	4:B:401:AP9:H141	1.81	0.62
1:A:12:LYS:CD	1:A:13:ILE:H	2.11	0.62
1:A:207:ILE:O	1:A:211:ASN:HA	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:144:ARG:NH1	3:B:402:SO4:O3	2.36	0.59
2:B:114:GLU:C	2:B:116:GLY:N	2.60	0.59
2:B:65:LEU:HA	5:B:403:DMS:H11	1.85	0.58
2:B:129:LEU:HD12	2:B:161:LEU:HD21	1.86	0.58
1:A:34:LEU:HB2	1:A:35:PRO:HD2	1.86	0.58
1:A:116:VAL:HG23	1:A:132:LEU:CD1	2.32	0.58
1:A:217:TRP:C	1:A:217:TRP:CD1	2.82	0.57
2:B:214:ARG:HD2	2:B:265:PHE:O	2.03	0.57
1:A:17:THR:HG22	1:A:18:MET:HG3	1.87	0.57
1:A:19:LYS:HE2	1:A:167:GLU:OE2	2.05	0.56
2:B:242:ASP:HB3	2:B:284:ASN:HB2	1.88	0.56
1:A:12:LYS:HD2	1:A:13:ILE:N	2.21	0.55
2:B:127:PHE:CE2	2:B:131:ARG:HD3	2.42	0.54
1:A:49:VAL:CG1	1:A:188:PRO:HB3	2.38	0.54
1:A:45:THR:H	1:A:88:GLN:NE2	2.07	0.53
2:B:143:HIS:O	2:B:201:ASP:OD1	2.25	0.53
2:B:171:SER:O	2:B:174:MET:HG2	2.09	0.53
2:B:228:LEU:HD12	2:B:254:PHE:CE2	2.43	0.53
1:A:43:ILE:HG22	1:A:43:ILE:O	2.10	0.52
1:A:211:ASN:O	1:A:214:CYS:HB3	2.10	0.52
2:B:275:ASP:HB3	2:B:297:HIS:HE2	1.73	0.52
2:B:235:ILE:CD1	2:B:281:LEU:O	2.57	0.52
1:A:151:LEU:O	1:A:152:ALA:C	2.53	0.52
2:B:262:ILE:HG12	2:B:277:LEU:HD23	1.91	0.52
1:A:43:ILE:O	1:A:43:ILE:CG2	2.59	0.51
2:B:68:LEU:HB2	5:B:403:DMS:H13	1.91	0.51
2:B:183:LEU:H	2:B:183:LEU:HD23	1.76	0.51
2:B:153:VAL:HA	2:B:158:GLN:O	2.12	0.49
2:B:177:THR:N	3:B:402:SO4:O2	2.38	0.49
1:A:116:VAL:C	1:A:118:LYS:H	2.21	0.49
1:A:121:TYR:HD2	1:A:122:LEU:HD12	1.77	0.49
2:B:126:MET:HG3	2:B:209:PHE:CD1	2.48	0.49
2:B:279:LYS:HG2	2:B:289:ILE:HD13	1.95	0.48
2:B:204:SER:O	2:B:208:ILE:HD12	2.14	0.48
1:A:20:ASP:OD1	1:A:22:ARG:HG2	2.13	0.48
2:B:25:GLY:HA3	2:B:44:ARG:O	2.13	0.48
2:B:209:PHE:CE2	2:B:276:LEU:HD23	2.49	0.48
2:B:77:VAL:HG21	4:B:401:AP9:H102	1.96	0.48
2:B:133:LEU:HD11	2:B:146:LEU:HD11	1.94	0.47
2:B:280:CYS:O	2:B:288:ARG:HD3	2.13	0.47
2:B:220:ARG:O	2:B:230:LYS:NZ	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:213:PHE:CE1	2:B:269:ILE:HA	2.50	0.47
1:A:240:ASP:O	1:A:244:GLU:CG	2.61	0.46
2:B:154:THR:OG1	2:B:158:GLN:HB3	2.15	0.46
1:A:145:TRP:HB3	2:B:62:VAL:HG11	1.98	0.46
1:A:217:TRP:O	1:A:218:THR:C	2.60	0.45
1:A:19:LYS:CE	1:A:167:GLU:OE2	2.64	0.45
2:B:42:LEU:CD2	2:B:97:VAL:HG13	2.46	0.45
2:B:202:LEU:CD1	2:B:290:SER:HA	2.46	0.45
1:A:151:LEU:C	1:A:153:THR:N	2.69	0.45
1:A:10:ARG:HD3	2:B:173:GLN:O	2.16	0.45
1:A:121:TYR:CD2	1:A:122:LEU:HD12	2.52	0.45
2:B:196:TYR:HB2	2:B:200:VAL:HG21	2.00	0.44
1:A:68:LEU:HD11	1:A:114:MET:HE1	1.97	0.44
2:B:61:GLU:HG3	2:B:164:PHE:HB2	1.99	0.44
2:B:71:PHE:CZ	2:B:139:HIS:HE1	2.36	0.44
2:B:113:PRO:O	2:B:114:GLU:O	2.35	0.44
2:B:177:THR:HB	2:B:179:VAL:HG23	1.99	0.43
2:B:153:VAL:HG12	2:B:159:ILE:HD13	2.00	0.43
2:B:212:MET:HB3	2:B:212:MET:HE2	1.78	0.43
2:B:180:VAL:HG13	2:B:181:VAL:HG23	1.99	0.43
1:A:55:LEU:O	1:A:59:MET:HG3	2.18	0.42
1:A:121:TYR:HD2	1:A:122:LEU:CD1	2.32	0.42
2:B:196:TYR:HB2	2:B:200:VAL:CG2	2.49	0.42
1:A:12:LYS:CG	1:A:13:ILE:H	2.32	0.42
2:B:73:HIS:HE1	2:B:75:ASN:HB2	1.84	0.42
1:A:73:PHE:HB3	1:A:74:PRO:HD3	2.02	0.42
2:B:42:LEU:HD21	2:B:97:VAL:HG13	2.00	0.42
2:B:77:VAL:HG21	4:B:401:AP9:C10	2.50	0.42
2:B:202:LEU:HD11	2:B:290:SER:HA	2.02	0.42
2:B:81:ASP:HB2	2:B:97:VAL:CG2	2.50	0.41
1:A:45:THR:H	1:A:88:GLN:HE21	1.66	0.41
1:A:83:TYR:CE1	1:A:139:ILE:HA	2.55	0.41
2:B:81:ASP:HB2	2:B:97:VAL:HG23	2.02	0.41
1:A:135:GLN:HE21	1:A:139:ILE:HG13	1.85	0.41
1:A:204:LEU:HD22	1:A:217:TRP:CD1	2.55	0.41
1:A:14:ASP:OD2	2:B:197:ALA:HB1	2.21	0.41
2:B:73:HIS:CE1	2:B:75:ASN:HB2	2.56	0.41
2:B:170:TYR:CE2	2:B:197:ALA:HA	2.56	0.41
2:B:42:LEU:HD23	2:B:42:LEU:HA	1.91	0.41
2:B:275:ASP:O	2:B:279:LYS:HB2	2.21	0.41
2:B:114:GLU:HA	2:B:114:GLU:OE2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:ARG:HA	1:A:25:ASN:HB2	2.02	0.40
1:A:212:THR:O	1:A:213:ASN:C	2.65	0.40
2:B:262:ILE:HG12	2:B:277:LEU:CD2	2.51	0.40
2:B:11:GLN:C	2:B:13:TYR:H	2.29	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	240/254 (94%)	210 (88%)	28 (12%)	2 (1%)	16	44
2	B	268/308 (87%)	229 (85%)	29 (11%)	10 (4%)	2	9
All	All	508/562 (90%)	439 (86%)	57 (11%)	12 (2%)	4	17

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	72	GLU
2	B	114	GLU
2	B	115	PRO
1	A	117	SER
1	A	212	THR
2	B	113	PRO
2	B	181	VAL
2	B	12	GLN
2	B	38	ARG
2	B	163	ASP
2	B	298	PRO
2	B	260	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	216/234 (92%)	195 (90%)	21 (10%)	8	25
2	B	217/272 (80%)	196 (90%)	21 (10%)	8	25
All	All	433/506 (86%)	391 (90%)	42 (10%)	8	25

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	24	LEU
1	A	29	LEU
1	A	34	LEU
1	A	71	SER
1	A	75	LEU
1	A	94	LEU
1	A	110	THR
1	A	112	LYS
1	A	129	ASN
1	A	130	LEU
1	A	132	LEU
1	A	144	LYS
1	A	153	THR
1	A	164	LYS
1	A	168	ASP
1	A	187	GLN
1	A	217	TRP
1	A	226	SER
1	A	234	THR
1	A	248	LEU
2	B	66	ARG
2	B	97	VAL
2	B	102	ASP
2	B	124	ASP
2	B	133	LEU
2	B	136	LEU

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Mol	Chain	Res	Type
2	B	166	LEU
2	B	171	SER
2	B	177	THR
2	B	179	VAL
2	B	180	VAL
2	B	183	LEU
2	B	195	SER
2	B	220	ARG
2	B	224	ASP
2	B	228	LEU
2	B	234	VAL
2	B	240	GLU
2	B	270	ASP
2	B	282	THR
2	B	287	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	88	GLN
1	A	95	GLN
1	A	135	GLN
1	A	187	GLN
1	A	233	ASN
2	B	67	HIS
2	B	103	GLN
2	B	158	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](#)

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)

3 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
5	DMS	B	403	-	3,3,3	2.92	1 (33%)	3,3,3	1.14	0
4	AP9	B	401	-	30,30,30	2.40	6 (20%)	43,43,43	2.42	13 (30%)
3	SO4	B	402	-	4,4,4	0.37	0	6,6,6	0.52	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
4	AP9	B	401	-	-	10/18/18/18	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	401	AP9	C5A-CL1	-10.22	1.50	1.74
5	B	403	DMS	O-S	4.92	1.82	1.50
4	B	401	AP9	C5-C4	4.71	1.47	1.39
4	B	401	AP9	C5-C6	3.31	1.49	1.41
4	B	401	AP9	C1A-N6	-2.44	1.35	1.40
4	B	401	AP9	C8-N7	2.34	1.36	1.31
4	B	401	AP9	C5-N7	-2.23	1.35	1.39

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	AP9	C5-C4-N3	-8.94	117.77	127.18
4	B	401	AP9	N3-C4-N9	6.37	135.08	126.99
4	B	401	AP9	C4-C5-N7	-4.14	105.85	110.58
4	B	401	AP9	N2-C2-N1	3.91	123.24	117.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	AP9	C9-N9-C8	-3.28	124.15	128.05
4	B	401	AP9	C6-C5-N7	3.13	135.84	132.43
4	B	401	AP9	N6-C6-N1	3.12	123.69	119.08
4	B	401	AP9	C5-N7-C8	2.90	108.00	103.45
4	B	401	AP9	C2-N3-C4	2.73	122.75	113.99
4	B	401	AP9	N1-C2-N3	-2.58	122.04	126.26
4	B	401	AP9	C2-N2-C12	2.48	128.18	124.32
4	B	401	AP9	C2-N1-C6	2.22	121.82	116.36
4	B	401	AP9	C11-C9-C10	-2.21	106.45	112.46

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	401	AP9	C13-C12-C15-O1
4	B	401	AP9	N2-C12-C15-O1
4	B	401	AP9	C15-C12-N2-C2
4	B	401	AP9	C11-C9-N9-C4
4	B	401	AP9	C11-C9-N9-C8
4	B	401	AP9	C5-C6-N6-C1A
4	B	401	AP9	N1-C6-N6-C1A
4	B	401	AP9	C13-C12-N2-C2
4	B	401	AP9	C2A-C1A-N6-C6
4	B	401	AP9	C6A-C1A-N6-C6

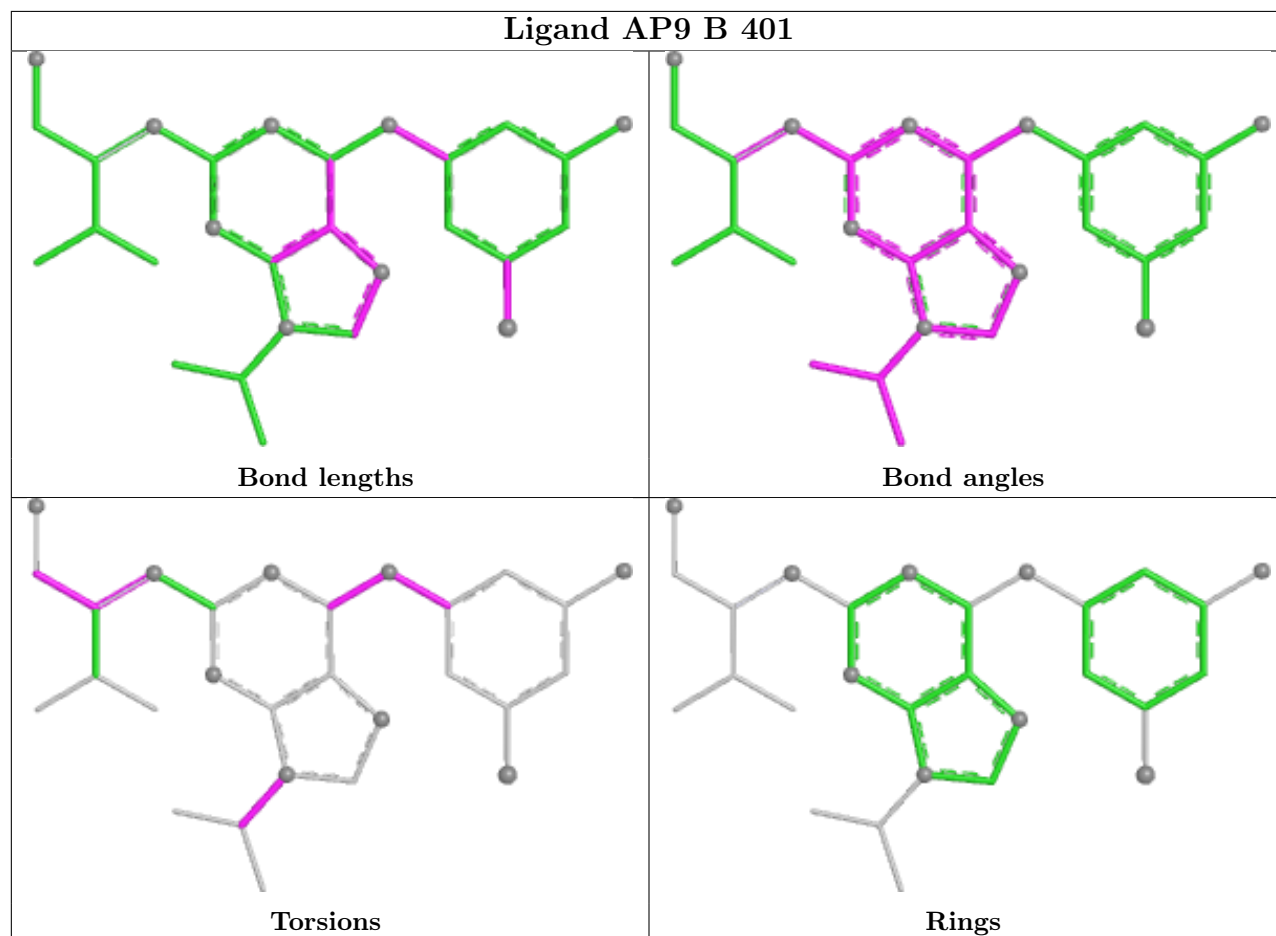
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	403	DMS	2	0
4	B	401	AP9	3	0
3	B	402	SO4	2	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	244/254 (96%)	-0.02	3 (1%) 76 68	62, 72, 86, 95	0
2	B	280/308 (90%)	0.24	10 (3%) 46 37	65, 72, 87, 102	0
All	All	524/562 (93%)	0.12	13 (2%) 58 48	62, 72, 87, 102	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	110	ASP	4.7
1	A	8	LEU	4.1
2	B	169	ILE	3.3
1	A	93	THR	3.2
2	B	175	ALA	2.9
2	B	173	GLN	2.5
2	B	171	SER	2.5
2	B	116	GLY	2.4
1	A	123	SER	2.4
2	B	183	LEU	2.3
2	B	115	PRO	2.3
2	B	177	THR	2.2
2	B	272	LEU	2.2

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](#)

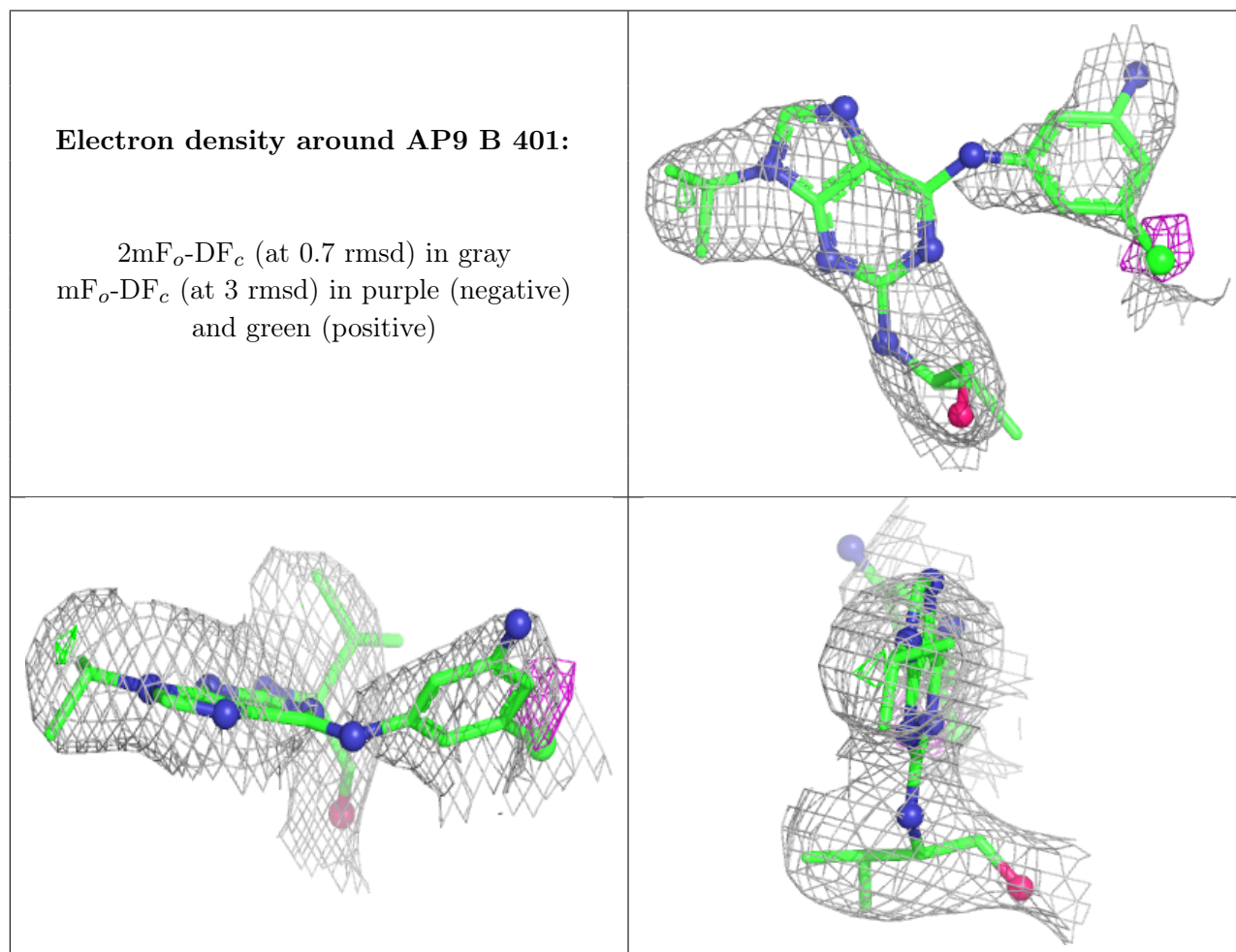
There are no oligosaccharides in this entry.

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
4	AP9	B	401	28/28	0.74	0.14	78,80,81,81	0
3	SO4	B	402	5/5	0.80	0.29	94,95,97,97	0
5	DMS	B	403	4/4	0.91	0.12	87,87,87,88	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.