



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:53 AM UTC

PDB ID : 3SLS / pdb_00003sls
Title : Crystal Structure of human MEK-1 kinase in complex with UCB1353770 and AMPPNP
Authors : Meier, C.; Ceska, T.A.
Deposited on : 2011-06-25
Resolution : 2.30 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

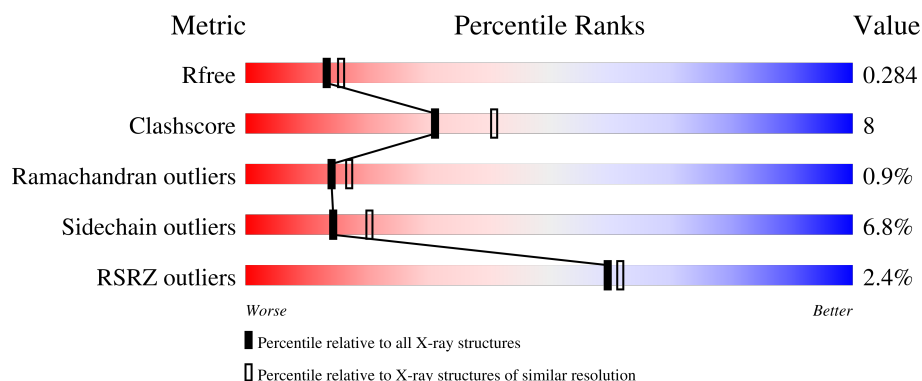
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

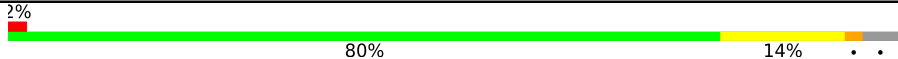
The reported resolution of this entry is 2.30 Å.

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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	304	
1	B	304	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5030 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 Dual specificity mitogen-activated protein kinase kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	3	0
			2264	1441	392	416	15			
1	B	295	Total	C	N	O	S	0	1	1
			2280	1458	386	420	16			

There are 106 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	MET	-	expression tag	UNP Q02750
A	43	THR	-	expression tag	UNP Q02750
A	44	LEU	-	expression tag	UNP Q02750
A	264	GLY	-	insertion	UNP Q02750
A	265	SER	-	insertion	UNP Q02750
A	266	GLY	-	insertion	UNP Q02750
A	267	SER	-	insertion	UNP Q02750
A	268	GLY	-	insertion	UNP Q02750
A	269	SER	-	insertion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	ASP	deletion	UNP Q02750
A	?	-	ALA	deletion	UNP Q02750
A	?	-	LYS	deletion	UNP Q02750
A	?	-	GLU	deletion	UNP Q02750
A	?	-	LEU	deletion	UNP Q02750
A	?	-	GLU	deletion	UNP Q02750
A	?	-	LEU	deletion	UNP Q02750
A	?	-	MET	deletion	UNP Q02750
A	?	-	PHE	deletion	UNP Q02750
A	?	-	GLY	deletion	UNP Q02750
A	?	-	CYS	deletion	UNP Q02750
A	?	-	GLN	deletion	UNP Q02750
A	?	-	VAL	deletion	UNP Q02750

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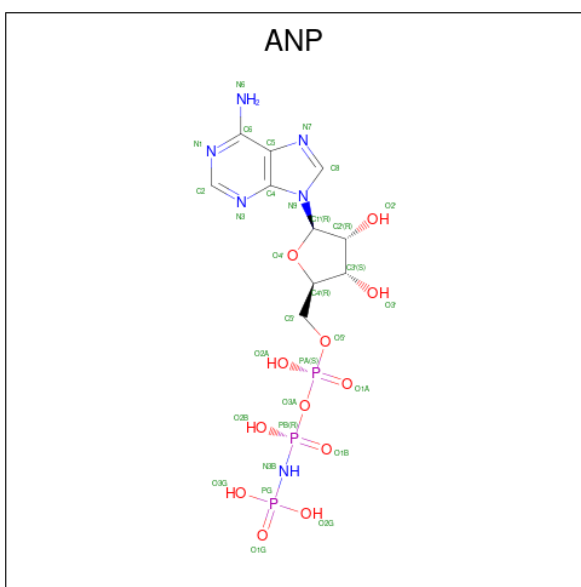
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	deletion	UNP Q02750
A	?	-	GLY	deletion	UNP Q02750
A	?	-	ASP	deletion	UNP Q02750
A	?	-	ALA	deletion	UNP Q02750
A	?	-	ALA	deletion	UNP Q02750
A	?	-	GLU	deletion	UNP Q02750
A	?	-	THR	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	ARG	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	ARG	deletion	UNP Q02750
A	?	-	THR	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	GLY	deletion	UNP Q02750
A	?	-	ARG	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	LEU	deletion	UNP Q02750
A	?	-	SER	deletion	UNP Q02750
A	?	-	SER	deletion	UNP Q02750
A	?	-	TYR	deletion	UNP Q02750
A	?	-	GLY	deletion	UNP Q02750
A	?	-	MET	deletion	UNP Q02750
A	?	-	ASP	deletion	UNP Q02750
A	?	-	SER	deletion	UNP Q02750
A	?	-	ARG	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
A	?	-	PRO	deletion	UNP Q02750
B	42	MET	-	expression tag	UNP Q02750
B	43	THR	-	expression tag	UNP Q02750
B	44	LEU	-	expression tag	UNP Q02750
B	264	GLY	-	insertion	UNP Q02750
B	265	SER	-	insertion	UNP Q02750
B	266	GLY	-	insertion	UNP Q02750
B	267	SER	-	insertion	UNP Q02750
B	268	GLY	-	insertion	UNP Q02750
B	269	SER	-	insertion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	ASP	deletion	UNP Q02750
B	?	-	ALA	deletion	UNP Q02750

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LYS	deletion	UNP Q02750
B	?	-	GLU	deletion	UNP Q02750
B	?	-	LEU	deletion	UNP Q02750
B	?	-	GLU	deletion	UNP Q02750
B	?	-	LEU	deletion	UNP Q02750
B	?	-	MET	deletion	UNP Q02750
B	?	-	PHE	deletion	UNP Q02750
B	?	-	GLY	deletion	UNP Q02750
B	?	-	CYS	deletion	UNP Q02750
B	?	-	GLN	deletion	UNP Q02750
B	?	-	VAL	deletion	UNP Q02750
B	?	-	GLU	deletion	UNP Q02750
B	?	-	GLY	deletion	UNP Q02750
B	?	-	ASP	deletion	UNP Q02750
B	?	-	ALA	deletion	UNP Q02750
B	?	-	ALA	deletion	UNP Q02750
B	?	-	GLU	deletion	UNP Q02750
B	?	-	THR	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	ARG	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	ARG	deletion	UNP Q02750
B	?	-	THR	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	GLY	deletion	UNP Q02750
B	?	-	ARG	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	LEU	deletion	UNP Q02750
B	?	-	SER	deletion	UNP Q02750
B	?	-	SER	deletion	UNP Q02750
B	?	-	TYR	deletion	UNP Q02750
B	?	-	GLY	deletion	UNP Q02750
B	?	-	MET	deletion	UNP Q02750
B	?	-	ASP	deletion	UNP Q02750
B	?	-	SER	deletion	UNP Q02750
B	?	-	ARG	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750
B	?	-	PRO	deletion	UNP Q02750

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).

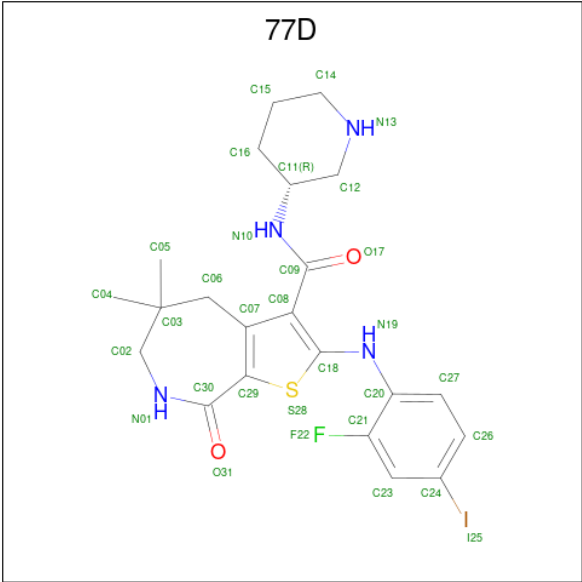


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	B	1	Total 31	C 10	N 6	O 12	P 3	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0

- Molecule 4 is 2-[(2-fluoro-4-iodophenyl)amino]-5,5-dimethyl-8-oxo-N-[(3R)-piperidin-3-yl]-5,6,7,8-tetrahydro-4H-thieno[2,3-c]azepine-3-carboxamide (CCD ID: 77D) (formula: $C_{22}H_{26}FIN_4O_2S$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
4	A	1	Total	C	F	I	N	O	S	0	0
			31	22	1	1	4	2	1		
4	B	1	Total	C	F	I	N	O	S	0	0
			31	22	1	1	4	2	1		

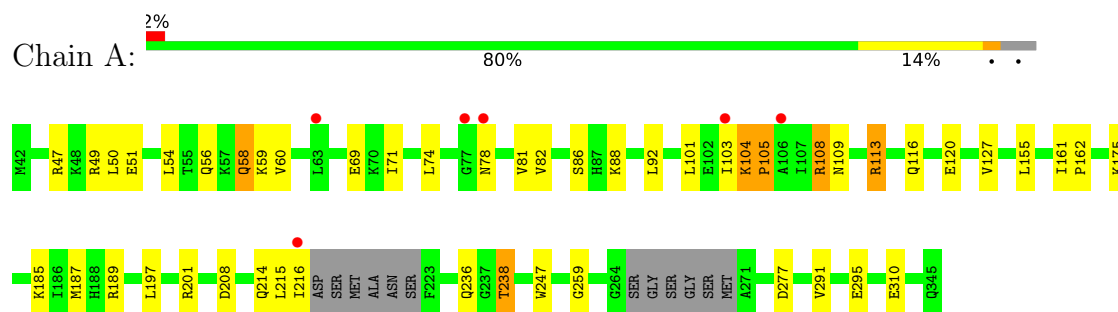
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	184	Total	O	0	0
			184	184		
5	B	176	Total	O	0	0
			176	176		

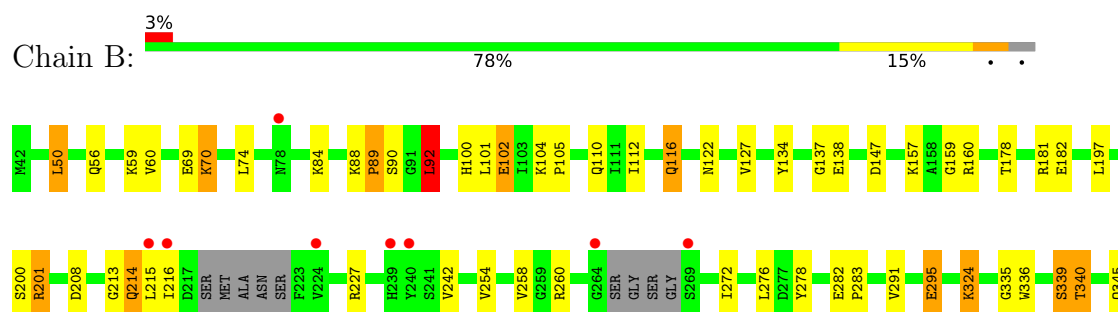
3 Residue-property plots [i](#)

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: Dual specificity mitogen-activated protein kinase kinase 1



- Molecule 1: Dual specificity mitogen-activated protein kinase kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.50Å 153.89Å 48.21Å 90.00° 101.48° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.30) 99.2 (30.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.206 , 0.287 0.204 , 0.284	Depositor DCC
R_{free} test set	1508 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.037 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5030	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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: 77D, MG, ANP

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	0.58	0/2314	0.88	0/3114
1	B	0.60	1/2322 (0.0%)	0.90	3/3125 (0.1%)
All	All	0.59	1/4636 (0.0%)	0.89	3/6239 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	283	PRO	CA-C	5.28	1.54	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	104	LYS	CA-C-N	6.45	127.91	119.84
1	B	104	LYS	C-N-CA	6.45	127.91	119.84
1	B	92	LEU	N-CA-C	6.15	119.18	109.96

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	2264	0	2253	31	0
1	B	2280	0	2288	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	31	0	13	1	0
2	B	31	0	13	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	31	0	26	5	0
4	B	31	0	26	2	0
5	A	184	0	0	3	0
5	B	176	0	0	10	0
All	All	5030	0	4619	76	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 8.

All (76) 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:104:LYS:HB3	1:A:105:PRO:HA	1.28	1.10
1:B:336:TRP:O	1:B:340:THR:HG23	1.69	0.93
1:A:109:ASN:O	1:A:113:ARG:HG2	1.74	0.87
1:A:49:ARG:HG2	1:A:201:ARG:HH21	1.40	0.82
1:A:104:LYS:HB3	1:A:105:PRO:CA	2.11	0.79
1:B:50:LEU:HD11	1:B:122:ASN:HB3	1.64	0.79
1:B:335:GLY:O	1:B:339:SER:HB2	1.86	0.76
1:B:213:GLY:HA3	1:B:214:GLN:CB	2.16	0.75
1:B:213:GLY:CA	1:B:214:GLN:HB2	2.17	0.74
1:A:49:ARG:HG2	1:A:201:ARG:NH2	2.04	0.72
1:A:104:LYS:CB	1:A:105:PRO:HA	2.15	0.72
1:B:213:GLY:CA	1:B:214:GLN:CB	2.71	0.67
1:B:147:ASP:CG	1:B:201:ARG:HH22	2.02	0.66
1:B:116:GLN:HG2	5:B:495:HOH:O	1.96	0.65
1:B:56:GLN:CG	1:B:92:LEU:HD11	2.27	0.65
1:A:277:ASP:OD1	5:A:367:HOH:O	2.15	0.64
1:B:260:ARG:NH1	5:B:356:HOH:O	2.25	0.61
4:A:500:77D:H06	4:A:500:77D:N10	2.15	0.61
1:A:214:GLN:HA	1:A:216:ILE:N	2.17	0.59
1:A:127:VAL:HG21	1:A:197:LEU:HD12	1.86	0.58
1:A:69:GLU:HB2	1:A:88:LYS:HD2	1.86	0.58
1:A:104:LYS:CB	1:A:105:PRO:CA	2.81	0.57
1:A:47:ARG:O	1:A:51:GLU:HG2	2.06	0.56
1:B:213:GLY:HA2	1:B:214:GLN:HB2	1.88	0.56
1:B:127:VAL:HG21	1:B:197:LEU:HD12	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LYS:HE3	1:B:340:THR:O	2.07	0.55
1:B:213:GLY:HA3	1:B:214:GLN:HB3	1.89	0.55
1:B:69:GLU:HB3	5:B:392:HOH:O	2.07	0.54
1:A:78:ASN:OD1	2:A:400:ANP:N3B	2.28	0.54
1:A:109:ASN:O	1:A:113:ARG:CG	2.52	0.53
1:B:208:ASP:HB3	5:B:476:HOH:O	2.09	0.52
1:B:88:LYS:C	1:B:90:SER:H	2.17	0.52
1:B:159:GLY:O	5:B:502:HOH:O	2.19	0.52
1:B:102:GLU:HG3	5:B:444:HOH:O	2.10	0.51
1:A:155:LEU:HD11	1:A:259:GLY:HA2	1.92	0.51
1:A:105:PRO:O	1:A:108:ARG:HG3	2.11	0.51
1:A:49:ARG:NH2	1:A:201:ARG:HB3	2.27	0.50
1:A:78:ASN:HB2	4:A:500:77D:H14A	1.93	0.49
1:B:147:ASP:CG	1:B:201:ARG:NH2	2.68	0.49
1:B:160:ARG:HD3	5:B:379:HOH:O	2.13	0.49
1:A:56:GLN:HA	1:A:59:LYS:HD3	1.95	0.48
4:A:500:77D:H06	4:A:500:77D:HN10	1.76	0.48
1:B:112:ILE:O	1:B:116:GLN:NE2	2.46	0.48
1:A:214:GLN:HA	1:A:215:LEU:C	2.38	0.48
1:B:295:GLU:H	1:B:295:GLU:CD	2.22	0.48
1:B:215:LEU:HA	1:B:216:ILE:HA	1.63	0.47
1:A:238:THR:HA	5:A:484:HOH:O	2.15	0.47
1:B:100:HIS:HA	1:B:138:GLU:HG2	1.97	0.47
1:B:74:LEU:HD21	1:B:84:LYS:HB2	1.97	0.46
1:B:134:TYR:OH	1:B:137:GLY:HA2	2.16	0.46
4:B:500:77D:H06	4:B:500:77D:N10	2.31	0.46
1:A:247:TRP:CD1	1:A:247:TRP:C	2.94	0.45
1:B:324:LYS:HE3	5:B:351:HOH:O	2.16	0.45
1:A:187:MET:HE2	1:A:189:ARG:HG2	1.98	0.44
4:A:500:77D:HN10	4:A:500:77D:C06	2.30	0.44
1:B:88:LYS:HB2	1:B:89:PRO:HD3	1.99	0.44
1:B:178:THR:O	1:B:182:GLU:HG2	2.17	0.44
1:B:116:GLN:CG	5:B:495:HOH:O	2.60	0.43
1:B:254:VAL:O	1:B:258:VAL:HG22	2.18	0.43
1:A:105:PRO:O	1:A:109:ASN:ND2	2.51	0.43
1:B:227:ARG:NH2	5:B:398:HOH:O	2.49	0.43
1:B:70:LYS:H	1:B:70:LYS:CD	2.32	0.43
1:A:161:ILE:HA	1:A:162:PRO:HD3	1.96	0.42
1:B:200:SER:OG	1:B:201:ARG:NH2	2.31	0.42
1:A:208:ASP:HA	4:A:500:77D:C20	2.50	0.42
1:B:208:ASP:HA	4:B:500:77D:C20	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:ARG:HA	1:B:181:ARG:HD2	1.75	0.42
1:A:74:LEU:HD12	1:A:82:VAL:HG12	2.03	0.41
1:B:272:ILE:O	1:B:276:LEU:HG	2.20	0.41
1:B:336:TRP:O	1:B:340:THR:CG2	2.53	0.41
1:B:147:ASP:OD2	1:B:201:ARG:NH2	2.54	0.40
1:A:71:ILE:HD11	1:A:86:SER:HB2	2.02	0.40
1:A:175:LYS:NZ	5:A:369:HOH:O	2.53	0.40
1:A:54:LEU:O	1:A:58[A]:GLN:NE2	2.54	0.40
1:B:278:TYR:CE2	1:B:282:GLU:HG3	2.57	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	289/304 (95%)	275 (95%)	11 (4%)	3 (1%)	12	15
1	B	289/304 (95%)	277 (96%)	10 (4%)	2 (1%)	18	23
All	All	578/608 (95%)	552 (96%)	21 (4%)	5 (1%)	14	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	LYS
1	A	103	ILE
1	A	105	PRO
1	B	105	PRO
1	B	89	PRO

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	242/262 (92%)	225 (93%)	17 (7%)	14	19
1	B	246/262 (94%)	228 (93%)	18 (7%)	13	18
All	All	488/524 (93%)	453 (93%)	35 (7%)	14	18

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU
1	A	58[A]	GLN
1	A	58[B]	GLN
1	A	81	VAL
1	A	92	LEU
1	A	101	LEU
1	A	108	ARG
1	A	113	ARG
1	A	116	GLN
1	A	120	GLU
1	A	185	LYS
1	A	236	GLN
1	A	238	THR
1	A	291	VAL
1	A	295	GLU
1	A	310[A]	GLU
1	A	310[B]	GLU
1	B	50	LEU
1	B	59	LYS
1	B	60	VAL
1	B	70	LYS
1	B	92	LEU
1	B	101	LEU
1	B	102	GLU
1	B	110	GLN
1	B	116	GLN
1	B	201	ARG

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Mol	Chain	Res	Type
1	B	214	GLN
1	B	242	VAL
1	B	291	VAL
1	B	295	GLU
1	B	324	LYS
1	B	339	SER
1	B	340	THR
1	B	345	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	145	HIS
1	A	164	GLN
1	A	239	HIS
1	A	316	GLN
1	B	78	ASN
1	B	109	ASN
1	B	116	GLN
1	B	214	GLN
1	B	345	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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
4	77D	A	500	-	30,34,34	0.95	1 (3%)	38,50,50	1.28	5 (13%)
2	ANP	A	400	3	33,33,33	2.08	10 (30%)	45,52,52	1.99	14 (31%)
2	ANP	B	400	3	33,33,33	2.09	13 (39%)	45,52,52	1.97	13 (28%)
4	77D	B	500	-	30,34,34	0.96	1 (3%)	38,50,50	1.35	5 (13%)

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	77D	A	500	-	-	0/12/35/35	0/4/4/4
2	ANP	A	400	3	-	3/18/38/38	0/3/3/3
2	ANP	B	400	3	-	5/18/38/38	0/3/3/3
4	77D	B	500	-	-	0/12/35/35	0/4/4/4

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	400	ANP	PG-N3B	5.06	1.76	1.63
2	B	400	ANP	PG-N3B	4.96	1.76	1.63
2	A	400	ANP	C5-C4	4.88	1.47	1.39
2	B	400	ANP	C5-C4	4.73	1.47	1.39
2	A	400	ANP	PB-N3B	4.62	1.75	1.63
2	B	400	ANP	PB-N3B	4.45	1.75	1.63
2	A	400	ANP	PA-O3A	3.15	1.62	1.59
2	B	400	ANP	C5-C6	3.03	1.49	1.41
2	B	400	ANP	PG-O1G	3.02	1.50	1.46
2	A	400	ANP	PG-O1G	2.92	1.50	1.46
4	A	500	77D	C30-N01	2.89	1.37	1.33
2	A	400	ANP	PB-O1B	2.87	1.50	1.46
2	B	400	ANP	PB-O1B	2.85	1.50	1.46
2	A	400	ANP	C5-C6	2.82	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	500	77D	C30-N01	2.67	1.37	1.33
2	B	400	ANP	PA-O3A	2.59	1.62	1.59
2	B	400	ANP	PB-O2B	-2.51	1.50	1.56
2	A	400	ANP	PB-O2B	-2.48	1.50	1.56
2	B	400	ANP	PB-O3A	2.47	1.62	1.59
2	B	400	ANP	C5-N7	-2.45	1.34	1.39
2	B	400	ANP	C8-N7	2.34	1.36	1.31
2	A	400	ANP	C8-N7	2.30	1.36	1.31
2	B	400	ANP	PG-O2G	-2.19	1.51	1.56
2	A	400	ANP	PG-O3G	-2.10	1.51	1.56
2	B	400	ANP	PG-O3G	-2.01	1.51	1.56

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	ANP	C5-C4-N3	-5.91	118.58	126.72
2	A	400	ANP	C5-C4-N3	-5.28	119.45	126.72
2	B	400	ANP	N3-C4-N9	4.73	135.22	127.17
2	A	400	ANP	N3-C4-N9	4.31	134.50	127.17
2	A	400	ANP	N3-C2-N1	-3.93	122.63	128.58
2	A	400	ANP	O1G-PG-N3B	-3.87	106.06	111.77
2	A	400	ANP	C2-N3-C4	3.80	121.11	111.83
2	B	400	ANP	C2-N3-C4	3.77	121.05	111.83
2	B	400	ANP	O2B-PB-O1B	3.73	117.86	109.87
4	A	500	77D	C30-C29-S28	3.59	119.53	110.60
2	B	400	ANP	C4-C5-N7	-3.48	106.61	110.58
2	A	400	ANP	C4-C5-N7	-3.29	106.82	110.58
4	B	500	77D	C30-C29-S28	3.26	118.72	110.60
4	A	500	77D	C18-C08-C07	3.26	113.43	111.75
2	A	400	ANP	O2B-PB-O1B	3.22	116.77	109.87
4	B	500	77D	C18-C08-C07	3.12	113.36	111.75
2	A	400	ANP	C4-N9-C8	2.97	108.86	105.74
2	B	400	ANP	N3-C2-N1	-2.97	124.08	128.58
4	A	500	77D	O31-C30-C29	2.86	125.04	116.81
2	B	400	ANP	C5-N7-C8	2.77	107.81	103.45
2	B	400	ANP	O2G-PG-O3G	2.70	114.86	107.59
2	B	400	ANP	C4-N9-C8	2.62	108.48	105.74
2	A	400	ANP	C6-C5-N7	2.51	136.93	132.09
2	A	400	ANP	C5-N7-C8	2.47	107.34	103.45
4	B	500	77D	C02-N01-C30	-2.45	119.93	125.18
4	B	500	77D	O31-C30-C29	2.38	123.66	116.81
2	A	400	ANP	O2A-PA-O3A	2.35	113.62	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	ANP	C2-N1-C6	2.28	122.48	118.73
2	B	400	ANP	O4'-C1'-N9	2.21	112.34	108.09
2	B	400	ANP	C6-C5-N7	2.19	136.31	132.09
4	B	500	77D	C29-S28-C18	2.17	93.19	91.30
2	B	400	ANP	N9-C8-N7	-2.14	110.89	113.94
2	A	400	ANP	N9-C8-N7	-2.12	110.93	113.94
2	B	400	ANP	C3'-C2'-C1'	2.10	105.44	101.46
4	A	500	77D	C05-C03-C06	2.10	112.74	109.09
4	A	500	77D	C23-C21-C20	-2.01	121.37	123.43
2	A	400	ANP	O1B-PB-N3B	-2.00	108.82	111.77

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	400	ANP	PG-N3B-PB-O3A
2	B	400	ANP	C5'-O5'-PA-O1A
2	A	400	ANP	PG-N3B-PB-O3A
2	B	400	ANP	C5'-O5'-PA-O3A
2	A	400	ANP	PB-O3A-PA-O2A
2	B	400	ANP	PB-O3A-PA-O2A
2	A	400	ANP	PB-O3A-PA-O1A
2	B	400	ANP	PB-O3A-PA-O1A

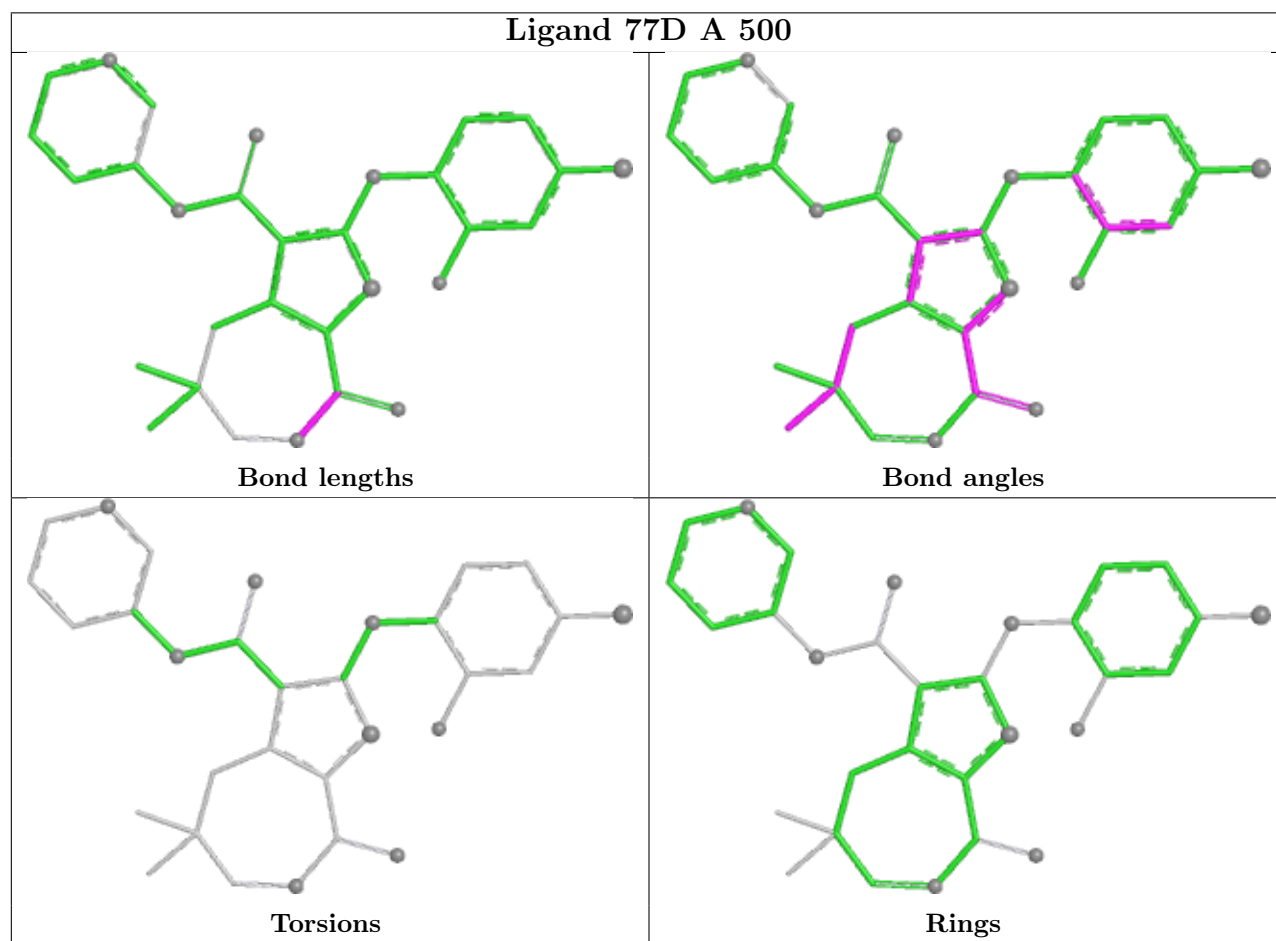
There are no ring outliers.

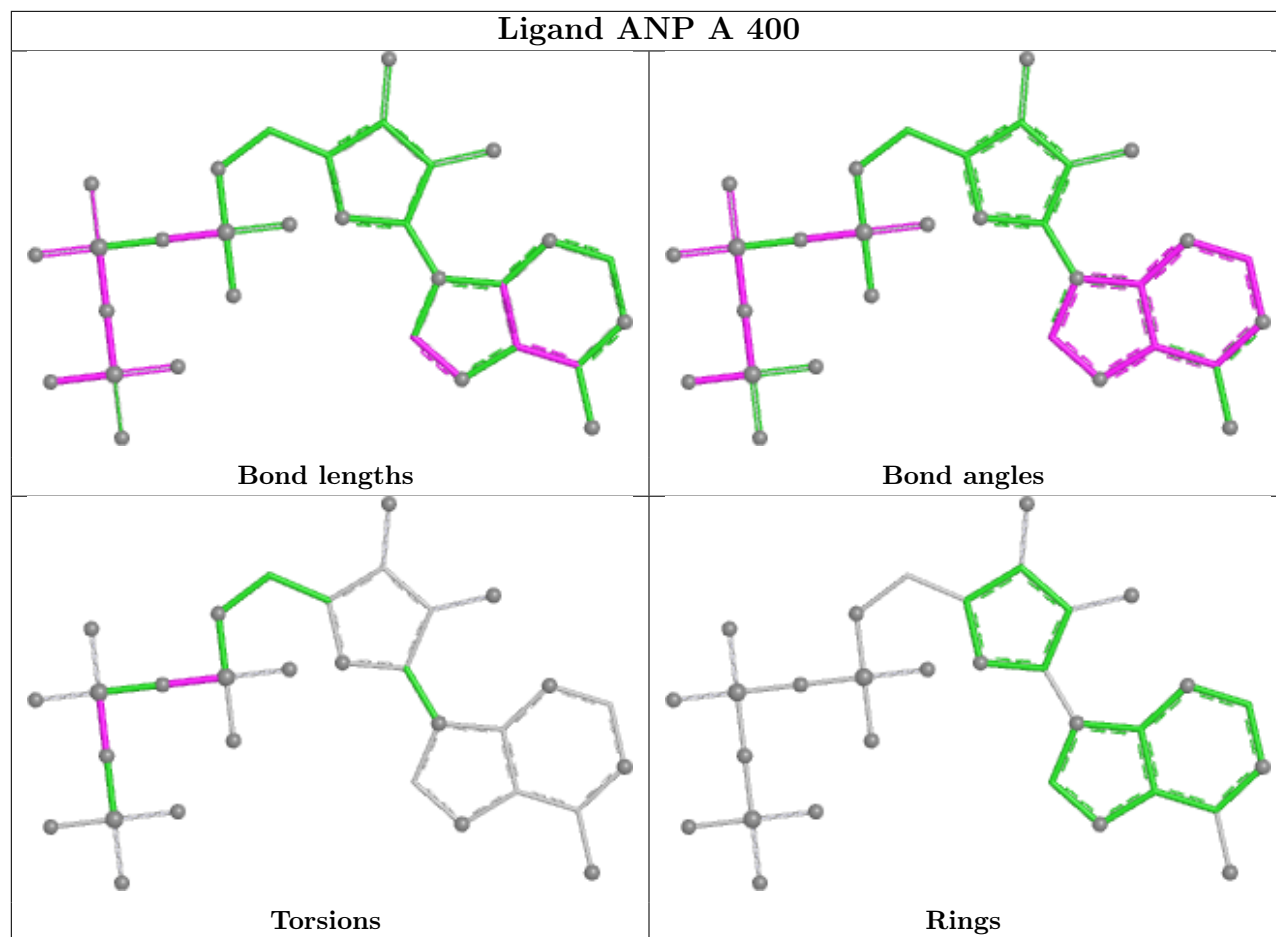
3 monomers are involved in 8 short contacts:

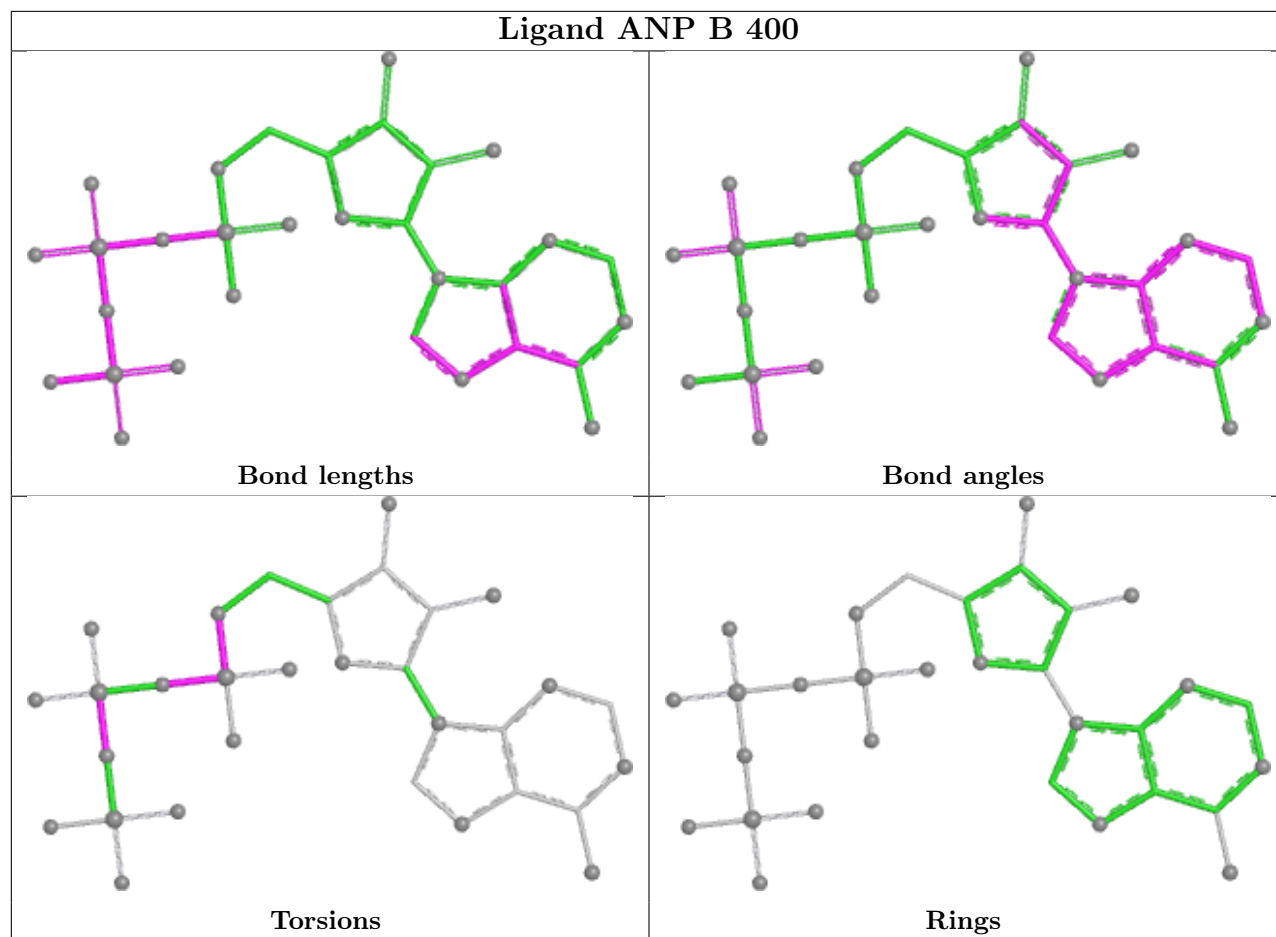
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	500	77D	5	0
2	A	400	ANP	1	0
4	B	500	77D	2	0

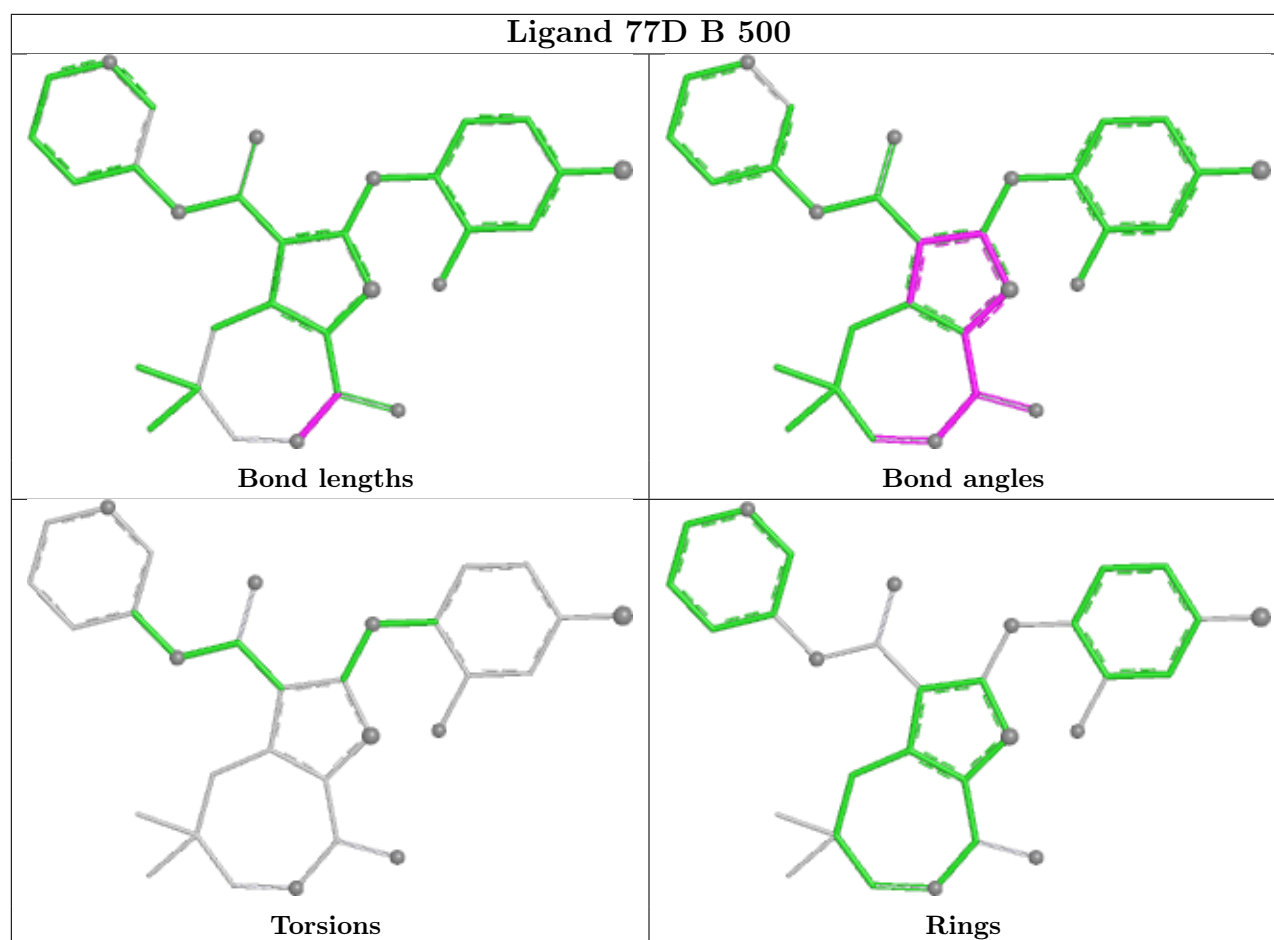
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	292/304 (96%)	0.21	6 (2%) 63 65	19, 42, 65, 72	11 (3%)
1	B	295/304 (97%)	0.24	8 (2%) 56 58	22, 42, 65, 73	11 (3%)
All	All	587/608 (96%)	0.22	14 (2%) 59 62	19, 42, 65, 73	22 (3%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	264	GLY	3.9
1	A	106	ALA	3.7
1	B	269	SER	3.1
1	B	224	VAL	3.0
1	B	239	HIS	3.0
1	A	63	LEU	2.8
1	B	216	ILE	2.8
1	A	216	ILE	2.7
1	B	240	TYR	2.3
1	A	103	ILE	2.2
1	B	215	LEU	2.2
1	A	78	ASN	2.2
1	B	78	ASN	2.1
1	A	77	GLY	2.0

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](#)

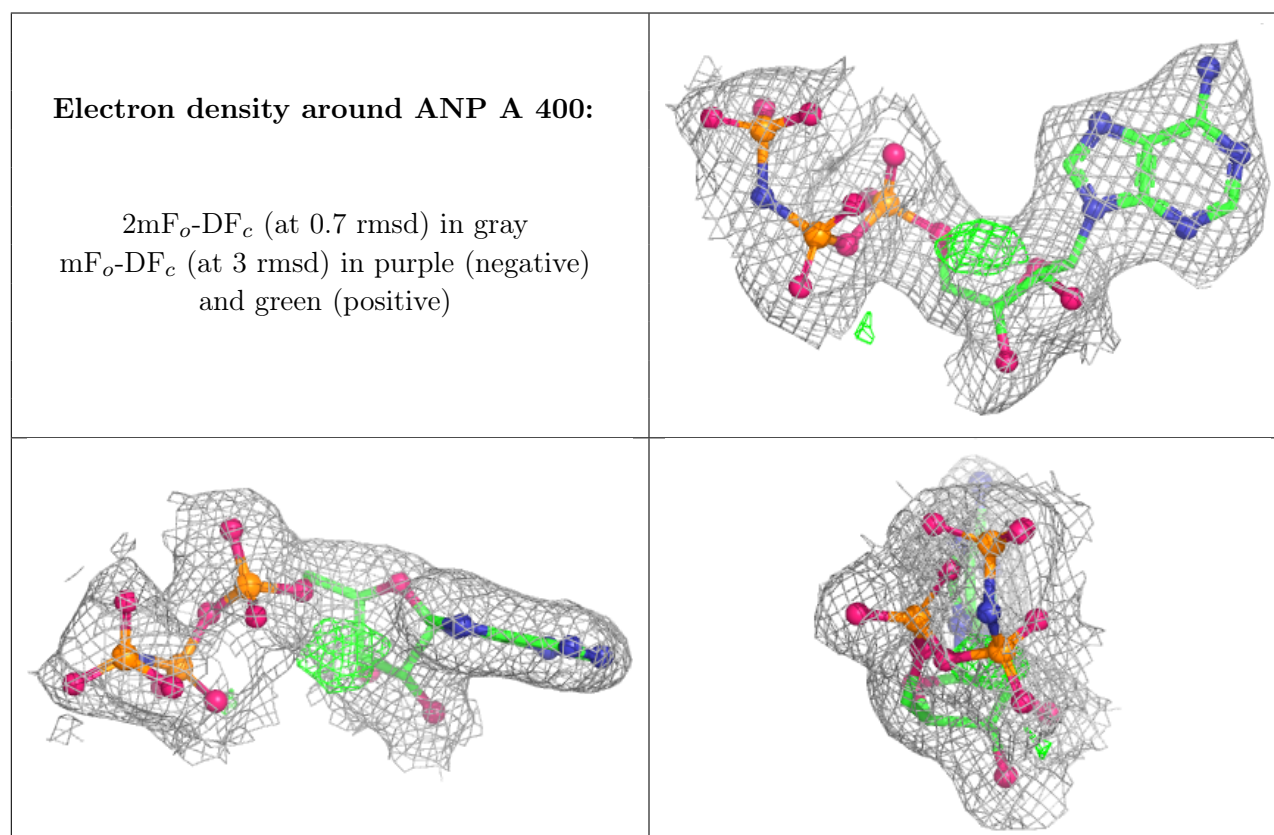
There are no oligosaccharides in this entry.

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. 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.

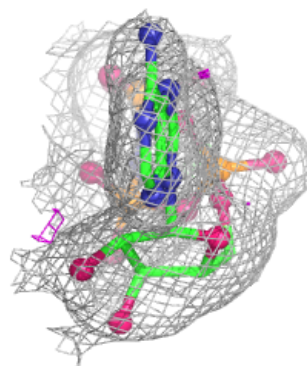
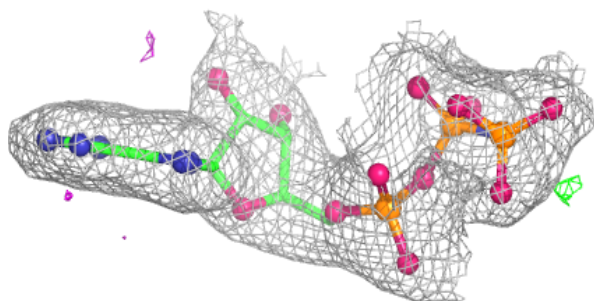
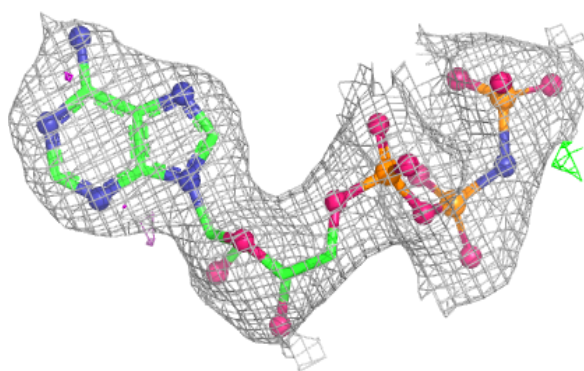
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ANP	A	400	31/31	0.96	0.06	34,39,42,43	0
2	ANP	B	400	31/31	0.98	0.05	30,37,40,43	0
3	MG	B	401	1/1	0.98	0.02	32,32,32,32	0
4	77D	B	500	31/31	0.98	0.07	35,40,43,45	0
4	77D	A	500	31/31	0.99	0.07	37,41,42,44	0
3	MG	A	401	1/1	1.00	0.01	40,40,40,40	0

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.

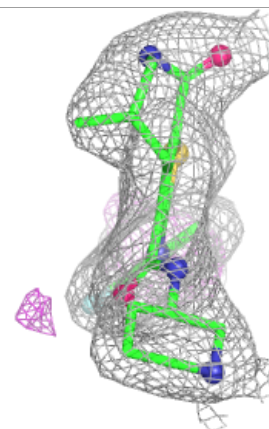
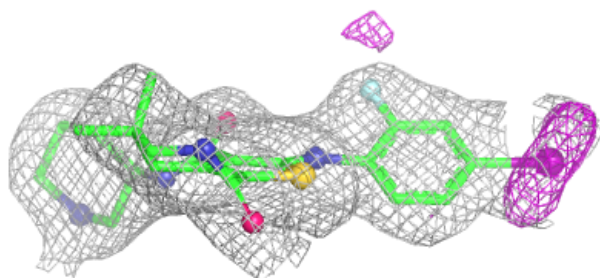
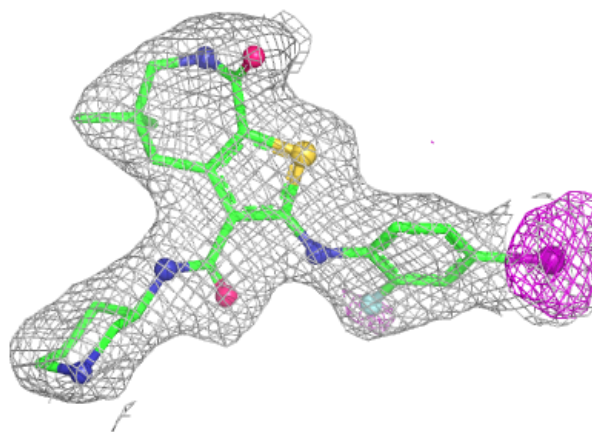


Electron density around ANP B 400:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

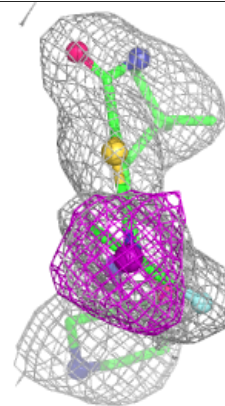
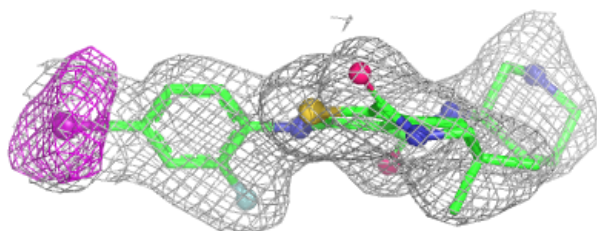
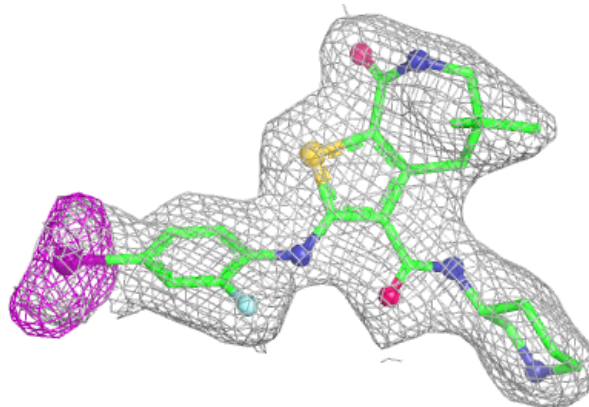
**Electron density around 77D B 500:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 77D A 500:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.