



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:15 AM UTC

PDB ID : 7JV0 / pdb_00007jv0
Title : Crystal Structure of KSR1:MEK1 in complex with AMP-PNP, and allosteric MEK inhibitor PD0325901
Authors : Khan, Z.M.; Dar, A.C.
Deposited on : 2020-08-20
Resolution : 3.63 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

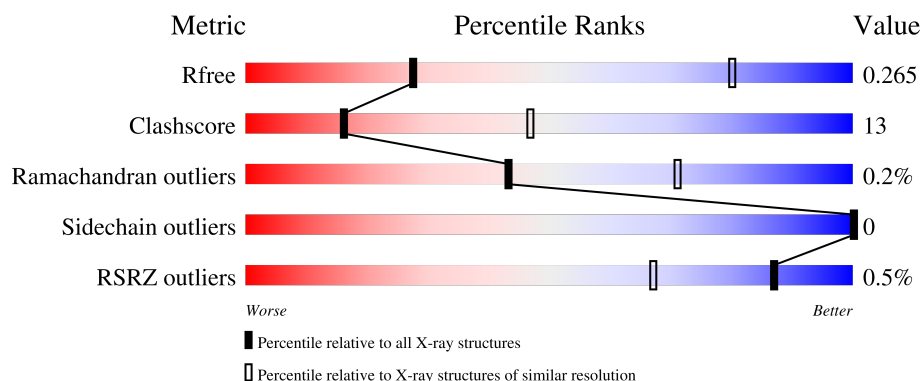
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.63 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	1132 (3.76-3.52)
Clashscore	190562	1014 (3.74-3.54)
Ramachandran outliers	187476	1123 (3.76-3.52)
Sidechain outliers	187428	1121 (3.76-3.52)
RSRZ outliers	180081	1130 (3.76-3.52)

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	334	 63% 19% 17%
2	C	384	 54% 26% 20%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4741 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 Kinase suppressor of Ras 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2221	1422	387	399	13			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	566	MET	-	initiating methionine	UNP Q8IVT5
A	567	SER	-	expression tag	UNP Q8IVT5
A	568	TYR	-	expression tag	UNP Q8IVT5
A	569	TYR	-	expression tag	UNP Q8IVT5
A	570	HIS	-	expression tag	UNP Q8IVT5
A	571	HIS	-	expression tag	UNP Q8IVT5
A	572	HIS	-	expression tag	UNP Q8IVT5
A	573	HIS	-	expression tag	UNP Q8IVT5
A	574	HIS	-	expression tag	UNP Q8IVT5
A	575	HIS	-	expression tag	UNP Q8IVT5
A	576	ASP	-	expression tag	UNP Q8IVT5
A	577	TYR	-	expression tag	UNP Q8IVT5
A	578	ASP	-	expression tag	UNP Q8IVT5
A	579	ILE	-	expression tag	UNP Q8IVT5
A	580	PRO	-	expression tag	UNP Q8IVT5
A	581	THR	-	expression tag	UNP Q8IVT5
A	582	THR	-	expression tag	UNP Q8IVT5
A	583	GLU	-	expression tag	UNP Q8IVT5
A	584	ASN	-	expression tag	UNP Q8IVT5
A	585	LEU	-	expression tag	UNP Q8IVT5
A	586	TYR	-	expression tag	UNP Q8IVT5
A	587	PHE	-	expression tag	UNP Q8IVT5
A	588	GLN	-	expression tag	UNP Q8IVT5
A	589	GLY	-	expression tag	UNP Q8IVT5
A	590	ALA	-	expression tag	UNP Q8IVT5
A	898	GLU	ASP	conflict	UNP Q8IVT5

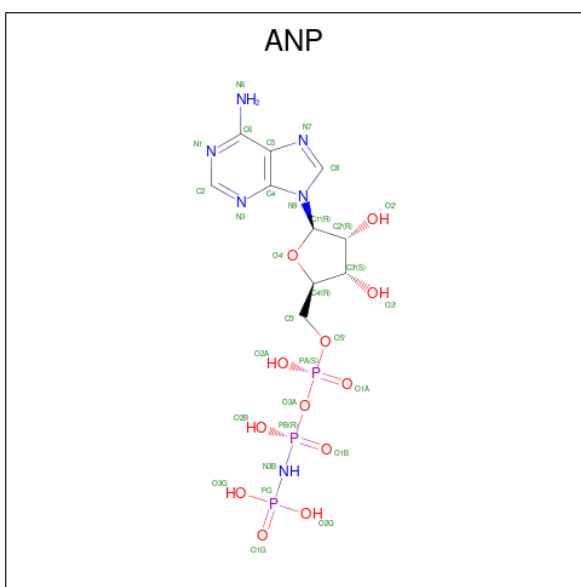
- Molecule 2 is a protein called Dual specificity mitogen-activated protein kinase kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	308	Total	C	N	O	S	0	0	0
			2431	1554	413	448	16			

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	10	MET	-	initiating methionine	UNP P29678
C	11	SER	-	expression tag	UNP P29678
C	12	TYR	-	expression tag	UNP P29678
C	13	TYR	-	expression tag	UNP P29678
C	14	HIS	-	expression tag	UNP P29678
C	15	HIS	-	expression tag	UNP P29678
C	16	HIS	-	expression tag	UNP P29678
C	17	HIS	-	expression tag	UNP P29678
C	18	HIS	-	expression tag	UNP P29678
C	19	HIS	-	expression tag	UNP P29678
C	20	ASP	-	expression tag	UNP P29678
C	21	TYR	-	expression tag	UNP P29678
C	22	ASP	-	expression tag	UNP P29678
C	23	ILE	-	expression tag	UNP P29678
C	24	PRO	-	expression tag	UNP P29678
C	25	THR	-	expression tag	UNP P29678
C	26	THR	-	expression tag	UNP P29678
C	27	GLU	-	expression tag	UNP P29678
C	28	ASN	-	expression tag	UNP P29678
C	29	LEU	-	expression tag	UNP P29678
C	30	TYR	-	expression tag	UNP P29678
C	31	PHE	-	expression tag	UNP P29678
C	32	GLN	-	expression tag	UNP P29678
C	33	GLY	-	expression tag	UNP P29678
C	34	ALA	-	expression tag	UNP P29678

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).

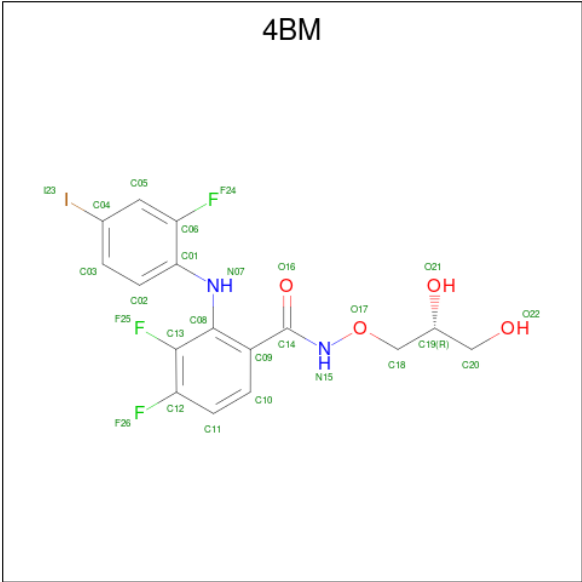


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	C	1	Total 31	C 10	N 6	O 12	P 3	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 5 is N-([(2R)-2,3-dihydroxypropyl]oxy)-3,4-difluoro-2-[(2-fluoro-4-iodophenyl)amino]benzamide (CCD ID: 4BM) (formula: $C_{16}H_{14}F_3IN_2O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	F	I	N	O		
5	C	1	26	16	3	1	2	4	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	137.66Å 137.66Å 220.89Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 3.63 40.00 – 3.63	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-3.63) 99.8 (40.00-3.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 3.66Å)	Xtriage
Refinement program	PHENIX (1.14rc1_3177)	Depositor
R, R_{free}	0.246 , 0.260 0.247 , 0.265	Depositor DCC
R_{free} test set	727 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	156.8	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 99.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4741	wwPDB-VP
Average B, all atoms (Å ²)	161.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: 4BM, 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	0.17	0/2274	0.47	0/3068
2	C	0.19	0/2477	0.55	0/3334
All	All	0.18	0/4751	0.51	0/6402

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
1	A	0	1
2	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	844	THR	Peptide
2	C	62	GLU	Peptide

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	2221	0	2199	42	0
2	C	2431	0	2473	82	0
3	A	31	0	13	1	0
3	C	31	0	13	3	0
4	A	1	0	0	0	0
5	C	26	0	14	5	0
All	All	4741	0	4712	124	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 (124) 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:C:63:LEU:HD21	2:C:87:HIS:NE2	2.00	0.77
1:A:644:ASP:HB2	1:A:759:VAL:HG22	1.67	0.76
1:A:646:HIS:O	1:A:646:HIS:ND1	2.18	0.76
2:C:64:LYS:HG3	2:C:65:ASP:H	1.52	0.72
2:C:180:LEU:HD22	2:C:186:ILE:HD11	1.73	0.70
2:C:143:MET:SD	3:C:401:ANP:N6	2.64	0.70
2:C:97:LYS:NZ	3:C:401:ANP:O2A	2.25	0.70
2:C:181:ARG:HH11	2:C:181:ARG:HB3	1.56	0.70
2:C:43:ASP:HB3	2:C:46:GLN:HG3	1.75	0.68
1:A:639:ARG:NH2	1:A:750:ASP:OD1	2.28	0.66
2:C:115:LEU:HD22	2:C:132:ALA:HB2	1.78	0.65
2:C:266:PRO:HB2	2:C:271:LEU:HG	1.79	0.65
2:C:271:LEU:HA	2:C:274:MET:HE2	1.78	0.65
2:C:151:LEU:HA	2:C:154:VAL:HB	1.79	0.64
2:C:168:LYS:HE3	2:C:369:VAL:HG12	1.81	0.62
1:A:811:TRP:CE2	1:A:855:ILE:HD11	2.35	0.61
1:A:715:GLN:NE2	1:A:874:MET:SD	2.73	0.61
2:C:215:LEU:HD11	2:C:219:MET:HE2	1.82	0.61
1:A:855:ILE:HG22	1:A:876:MET:HE2	1.82	0.60
1:A:700:ASP:O	1:A:703:THR:HG22	2.01	0.60
2:C:100:HIS:NE2	2:C:102:GLU:OE1	2.35	0.60
1:A:807:PHE:HA	1:A:810:VAL:HG12	1.86	0.58
2:C:92:LEU:HD23	2:C:93:VAL:N	2.19	0.58
2:C:181:ARG:HB3	2:C:181:ARG:NH1	2.19	0.58
2:C:234:ARG:HG3	2:C:240:TYR:CD1	2.39	0.57
2:C:127:VAL:HA	2:C:205:LYS:HE2	1.85	0.57
1:A:716:GLU:OE1	1:A:719:LYS:NZ	2.29	0.56
2:C:87:HIS:CE1	2:C:89:PRO:HB2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:108:ARG:HH12	2:C:137:GLY:H	1.52	0.56
2:C:46:GLN:HA	2:C:49:ARG:HH21	1.70	0.56
2:C:155:LEU:HD12	2:C:256:MET:HA	1.87	0.56
1:A:662:ARG:O	1:A:662:ARG:HD2	2.06	0.56
1:A:784:VAL:HA	1:A:787:MET:HG3	1.87	0.55
2:C:108:ARG:HH11	2:C:134:TYR:HB2	1.71	0.55
2:C:59:LYS:NZ	2:C:90:SER:HB2	2.21	0.55
2:C:164:GLN:NE2	2:C:367:GLU:OE2	2.40	0.54
1:A:791:LYS:HE3	1:A:795:GLN:HB3	1.89	0.53
2:C:110:GLN:HG3	2:C:113:ARG:NH2	2.23	0.53
2:C:363:ARG:O	2:C:367:GLU:HG3	2.09	0.52
2:C:49:ARG:HG2	2:C:201:ARG:NH2	2.24	0.52
1:A:706:ASP:O	1:A:710:THR:HG23	2.09	0.52
1:A:839:MET:HA	1:A:842:VAL:HG12	1.92	0.51
2:C:177:LEU:HB2	2:C:352:LEU:HD21	1.92	0.51
2:C:87:HIS:HE1	2:C:89:PRO:HB2	1.75	0.51
1:A:654:PHE:O	1:A:658:VAL:HG22	2.10	0.51
2:C:56:GLN:HG3	2:C:59:LYS:HE3	1.93	0.51
2:C:230:MET:HE3	2:C:234:ARG:HB3	1.92	0.51
1:A:769:LEU:HB3	2:C:224:VAL:HG13	1.93	0.51
2:C:188:HIS:CD2	2:C:209:PHE:HB3	2.46	0.51
2:C:44:GLU:OE2	2:C:47:ARG:NH1	2.45	0.50
2:C:308:MET:HE2	2:C:312:GLU:HB3	1.92	0.50
2:C:165:ILE:HD11	2:C:371:PHE:CD1	2.47	0.50
2:C:170:SER:O	2:C:174:ILE:HG13	2.12	0.50
2:C:379:ILE:HD11	2:C:381:LEU:HB2	1.94	0.49
2:C:101:LEU:HD11	2:C:219:MET:SD	2.52	0.49
2:C:123:SER:HB3	2:C:126:ILE:HD13	1.94	0.49
1:A:852:VAL:O	1:A:855:ILE:HG12	2.13	0.49
1:A:610:PHE:CE2	1:A:613:VAL:HG13	2.47	0.49
1:A:840:LYS:O	1:A:844:THR:HG23	2.12	0.48
1:A:766:GLU:N	1:A:766:GLU:OE1	2.47	0.48
2:C:127:VAL:HG11	2:C:197:LEU:HD12	1.94	0.48
2:C:63:LEU:HA	2:C:63:LEU:HD23	1.28	0.48
1:A:841:ARG:O	1:A:844:THR:OG1	2.26	0.48
2:C:270:GLU:O	2:C:274:MET:HG3	2.14	0.48
1:A:673:MET:HE3	1:A:687:SER:HB2	1.96	0.47
5:C:402:4BM:F25	5:C:402:4BM:H02	2.05	0.47
1:A:698:VAL:HG12	1:A:705:LEU:HD11	1.96	0.47
2:C:43:ASP:OD1	2:C:44:GLU:N	2.46	0.47
1:A:782:GLU:CD	1:A:867:ARG:HH22	2.22	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:801:ALA:O	1:A:804:VAL:HG22	2.15	0.47
2:C:100:HIS:C	2:C:101:LEU:HD12	2.38	0.47
2:C:115:LEU:HD22	2:C:132:ALA:CB	2.44	0.47
1:A:643:MET:HG3	1:A:682:LEU:HD11	1.97	0.46
2:C:64:LYS:HG3	2:C:65:ASP:N	2.26	0.46
2:C:63:LEU:HD13	2:C:68:PHE:CZ	2.50	0.46
1:A:622:GLY:HA3	3:A:901:ANP:O3G	2.15	0.46
1:A:807:PHE:O	1:A:810:VAL:HG12	2.16	0.46
2:C:255:GLU:HG3	2:C:261:TYR:HA	1.98	0.46
1:A:711:ARG:O	1:A:715:GLN:HG3	2.16	0.45
2:C:63:LEU:O	2:C:64:LYS:HB3	2.16	0.45
2:C:143:MET:HE2	5:C:402:4BM:H05	1.97	0.45
2:C:49:ARG:HE	2:C:49:ARG:HB2	1.67	0.45
2:C:127:VAL:HG23	5:C:402:4BM:I23	2.86	0.45
2:C:155:LEU:CD1	2:C:256:MET:HA	2.47	0.45
2:C:321:PRO:HA	2:C:322:PRO:HD3	1.85	0.45
1:A:831:TRP:CZ2	2:C:318:VAL:HG11	2.52	0.45
2:C:103:ILE:HD13	2:C:107:ILE:HG13	1.99	0.44
2:C:136:ASP:CG	2:C:136:ASP:O	2.60	0.44
2:C:82:VAL:HG11	3:C:401:ANP:C8	2.47	0.44
2:C:121:CYS:HG	2:C:184:HIS:CD2	2.35	0.44
2:C:125:TYR:CE2	2:C:175:LYS:HD3	2.52	0.44
2:C:226:THR:HG23	2:C:227:ARG:HG3	1.99	0.44
2:C:165:ILE:HD11	2:C:371:PHE:HD1	1.81	0.44
2:C:178:THR:HG23	2:C:352:LEU:HD22	2.00	0.44
1:A:659:MET:O	1:A:662:ARG:HB3	2.19	0.43
1:A:707:ILE:HG12	1:A:881:PRO:HB2	2.00	0.43
1:A:610:PHE:HB2	1:A:632:TRP:CZ2	2.54	0.43
1:A:833:ILE:HD13	1:A:839:MET:HE2	1.99	0.43
2:C:143:MET:CE	5:C:402:4BM:H05	2.49	0.43
1:A:625:GLY:HA3	1:A:640:LEU:O	2.19	0.42
1:A:648:GLN:O	1:A:652:LYS:HG3	2.19	0.42
1:A:785:ARG:NH1	1:A:834:GLY:O	2.51	0.42
2:C:71:ILE:HB	2:C:84:LYS:HB3	2.02	0.42
2:C:230:MET:HE2	2:C:235:LEU:HD23	2.02	0.42
2:C:265:PRO:HA	2:C:266:PRO:HD3	1.92	0.42
5:C:402:4BM:O16	5:C:402:4BM:N07	2.53	0.42
2:C:381:LEU:HA	2:C:381:LEU:HD23	1.78	0.42
2:C:230:MET:HE3	2:C:234:ARG:HD3	2.01	0.42
2:C:260:ARG:HD3	2:C:266:PRO:HG3	2.01	0.41
1:A:610:PHE:CD1	1:A:610:PHE:C	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:708:ASN:OD1	1:A:711:ARG:NH2	2.42	0.41
2:C:104:LYS:O	2:C:107:ILE:HG12	2.20	0.41
2:C:121:CYS:HG	2:C:184:HIS:CE1	2.38	0.41
1:A:602:TYR:C	1:A:604:GLN:H	2.28	0.41
1:A:776:LEU:HA	1:A:779:LEU:HD12	2.03	0.41
1:A:800:LYS:O	1:A:804:VAL:HG13	2.20	0.41
2:C:40:LEU:O	2:C:40:LEU:HD13	2.20	0.41
2:C:340:LYS:HA	2:C:343:ILE:HD12	2.01	0.41
2:C:140:SER:O	2:C:141:ILE:HD13	2.20	0.41
2:C:179:TYR:CZ	2:C:183:LYS:HD2	2.56	0.41
2:C:97:LYS:O	2:C:140:SER:HA	2.21	0.41
2:C:181:ARG:O	2:C:185:LYS:HD3	2.21	0.40
2:C:375:LEU:O	2:C:378:THR:OG1	2.39	0.40
2:C:186:ILE:HD12	2:C:186:ILE:C	2.46	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	270/334 (81%)	253 (94%)	17 (6%)	0	100	100
2	C	302/384 (79%)	289 (96%)	12 (4%)	1 (0%)	36	64
All	All	572/718 (80%)	542 (95%)	29 (5%)	1 (0%)	43	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	64	LYS

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	238/291 (82%)	238 (100%)	0	100	100
2	C	270/333 (81%)	270 (100%)	0	100	100
All	All	508/624 (81%)	508 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	824	GLN
1	A	832	GLN
2	C	46	GLN
2	C	145	HIS

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](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ANP	A	901	4	33,33,33	0.94	3 (9%)	45,52,52	0.60	0
3	ANP	C	401	-	33,33,33	0.89	3 (9%)	45,52,52	0.72	1 (2%)
5	4BM	C	402	-	26,27,27	1.95	6 (23%)	31,37,37	1.03	1 (3%)

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	ANP	A	901	4	-	3/18/38/38	0/3/3/3
3	ANP	C	401	-	-	3/18/38/38	0/3/3/3
5	4BM	C	402	-	-	3/16/16/16	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	402	4BM	C14-N15	5.93	1.43	1.34
5	C	402	4BM	C01-N07	4.41	1.52	1.39
3	A	901	ANP	PG-O1G	2.65	1.50	1.46
5	C	402	4BM	C09-C14	2.58	1.55	1.50
3	A	901	ANP	PG-N3B	2.57	1.70	1.63
5	C	402	4BM	C08-N07	2.55	1.48	1.41
3	A	901	ANP	PB-O1B	2.51	1.50	1.46
5	C	402	4BM	C09-C08	-2.46	1.37	1.41
3	C	401	ANP	PG-N3B	2.41	1.69	1.63
3	C	401	ANP	PG-O1G	2.38	1.49	1.46
3	C	401	ANP	PB-O1B	2.34	1.49	1.46
5	C	402	4BM	O16-C14	-2.11	1.18	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	402	4BM	C05-C06-C01	-3.20	120.16	123.43
3	C	401	ANP	O1G-PG-N3B	-2.49	108.10	111.77

There are no chirality outliers.

All (9) torsion outliers are listed below:

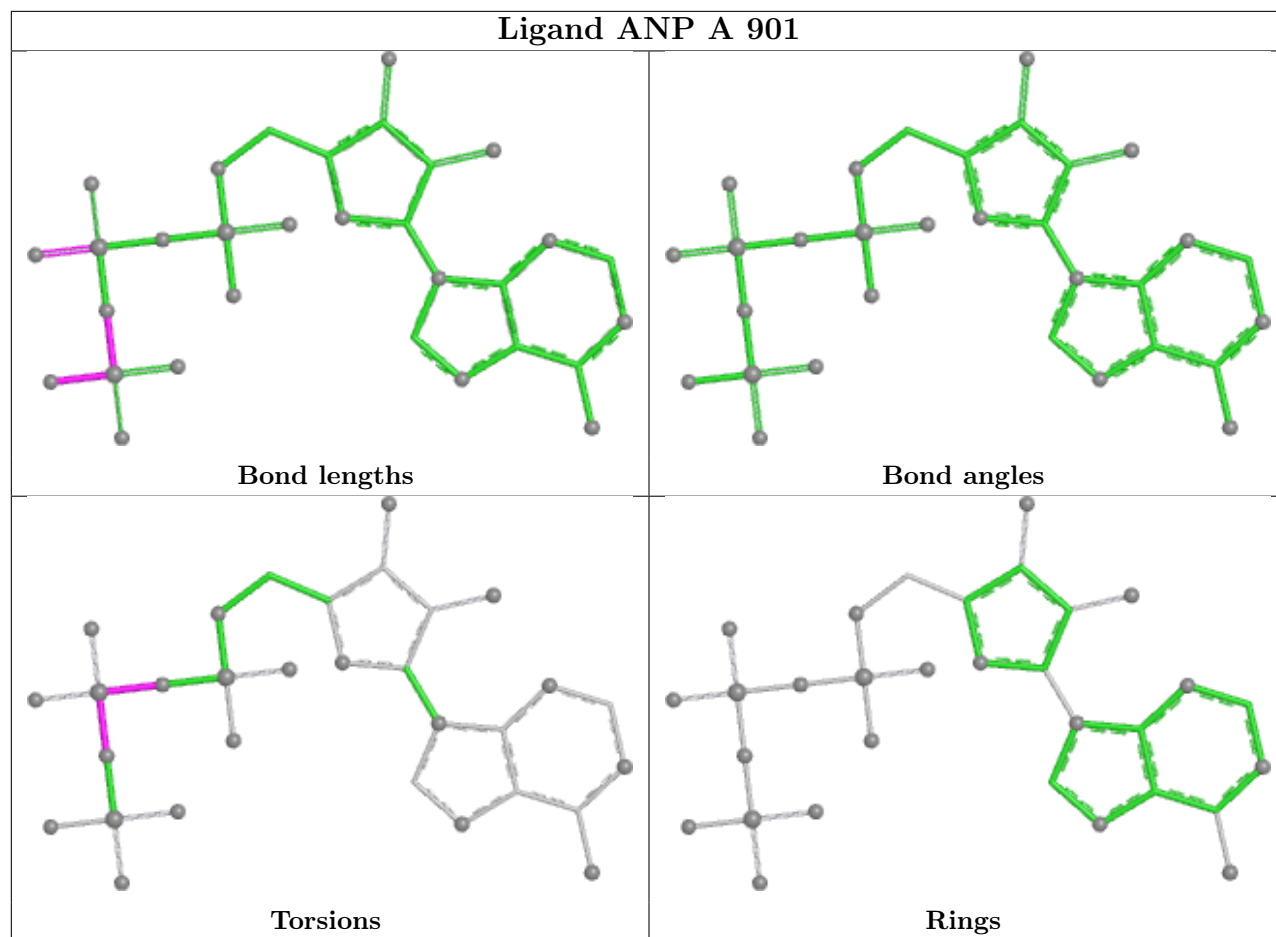
Mol	Chain	Res	Type	Atoms
3	A	901	ANP	PG-N3B-PB-O1B
3	A	901	ANP	PA-O3A-PB-O2B
3	C	401	ANP	PB-N3B-PG-O1G
3	C	401	ANP	PG-N3B-PB-O1B
3	C	401	ANP	PA-O3A-PB-O1B
5	C	402	4BM	C14-N15-O17-C18
5	C	402	4BM	C19-C18-O17-N15
3	A	901	ANP	PA-O3A-PB-O1B
5	C	402	4BM	C10-C09-C14-N15

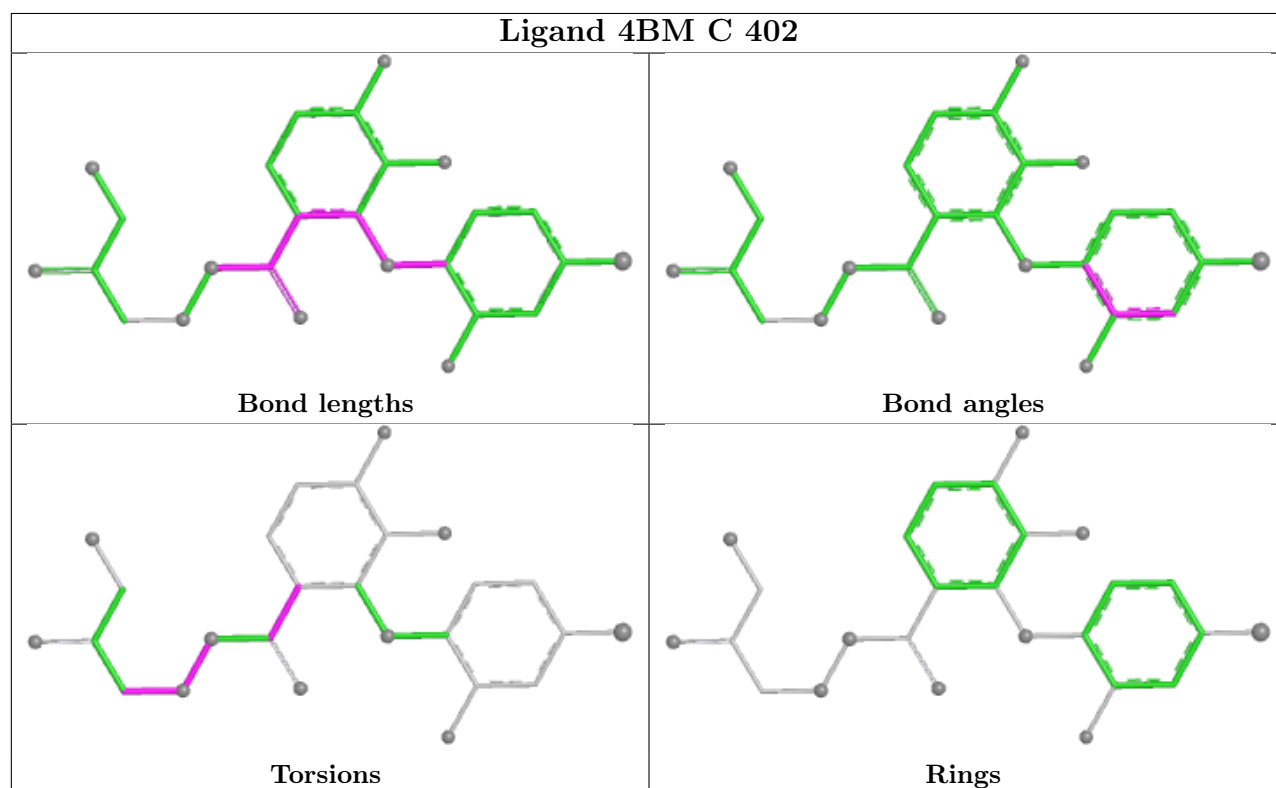
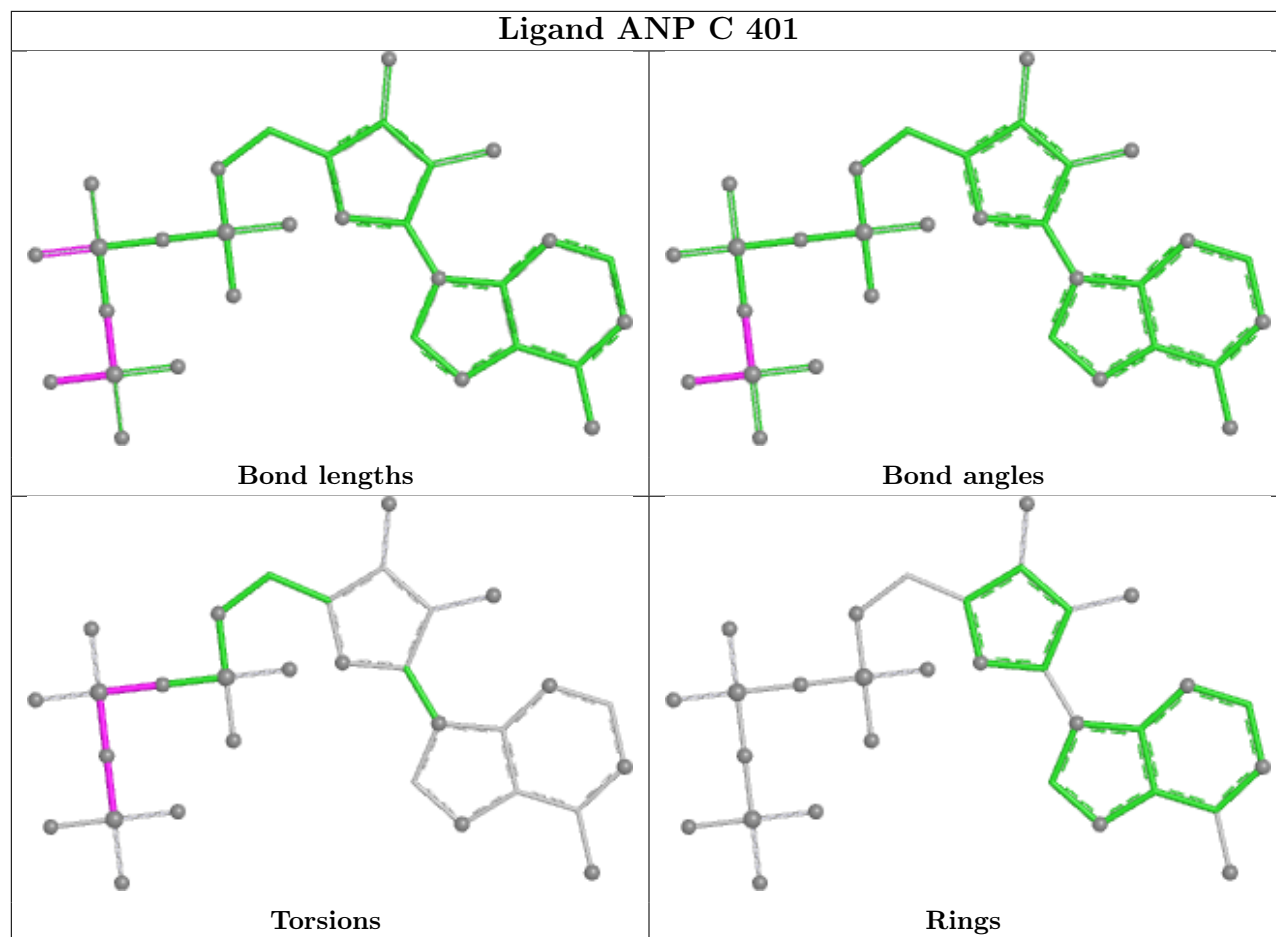
There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	ANP	1	0
3	C	401	ANP	3	0
5	C	402	4BM	5	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	632:TRP	C	633:HIS	N	5.41

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	276/334 (82%)	-0.22	1 (0%)	88 71	94, 131, 174, 229	0
2	C	308/384 (80%)	-0.00	2 (0%)	85 65	115, 188, 233, 250	0
All	All	584/718 (81%)	-0.11	3 (0%)	87 67	94, 153, 229, 250	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	846	VAL	2.9
2	C	219	MET	2.7
2	C	139	ILE	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.

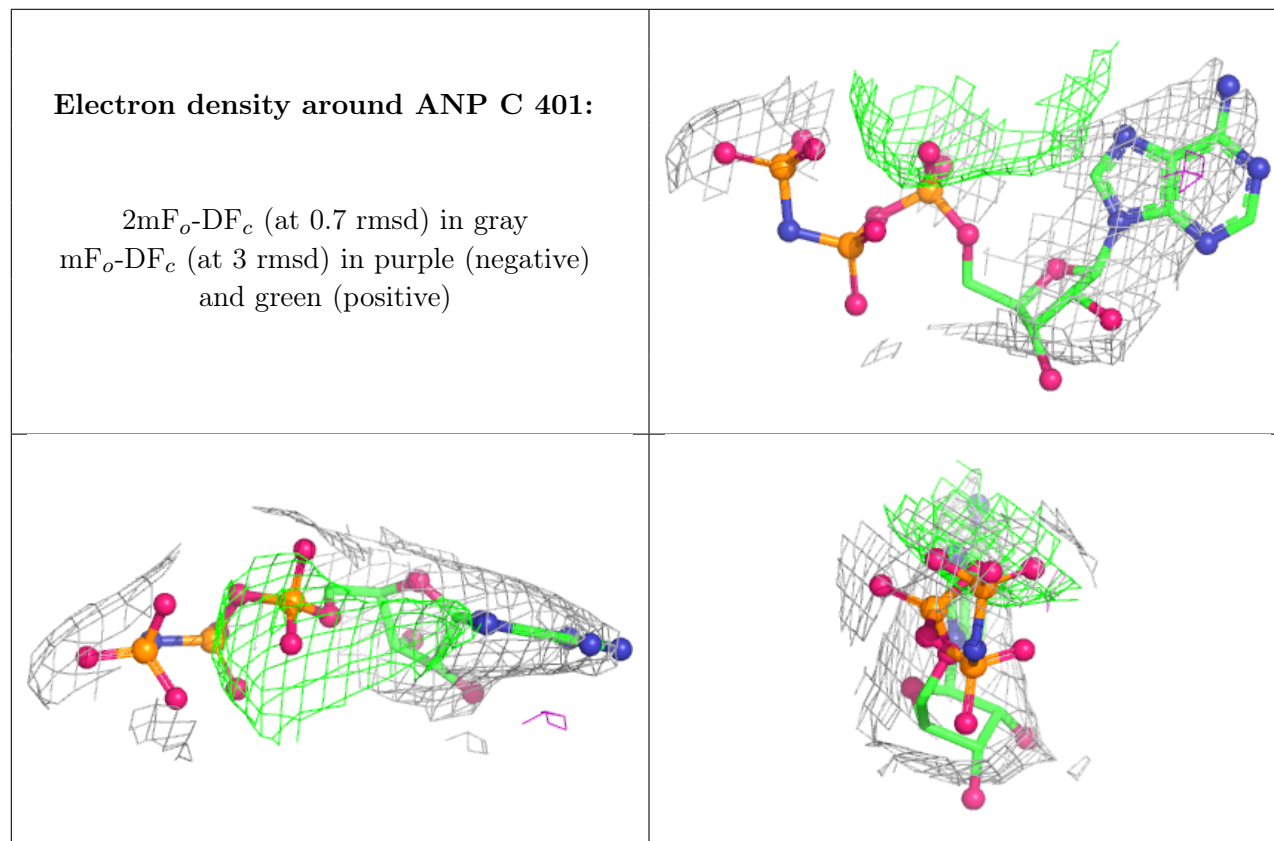
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
3	ANP	C	401	31/31	0.84	0.11	185,200,211,212	0
5	4BM	C	402	26/26	0.90	0.12	170,188,199,226	0

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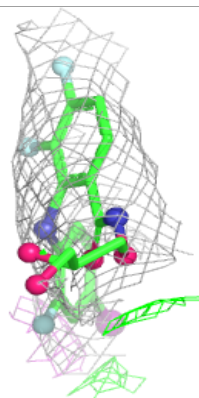
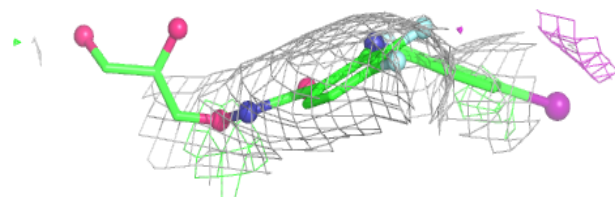
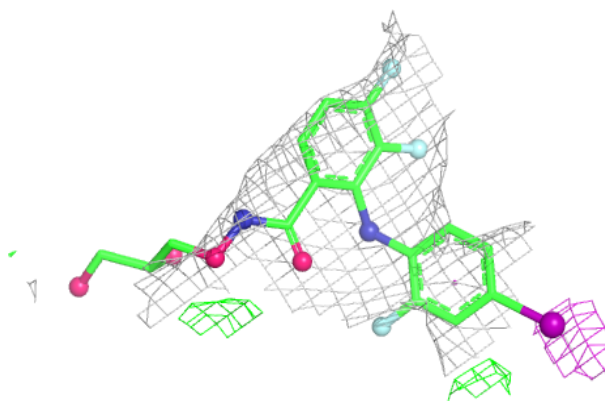
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	902	1/1	0.91	0.13	119,119,119,119	0
3	ANP	A	901	31/31	0.91	0.08	94,130,163,171	0

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.



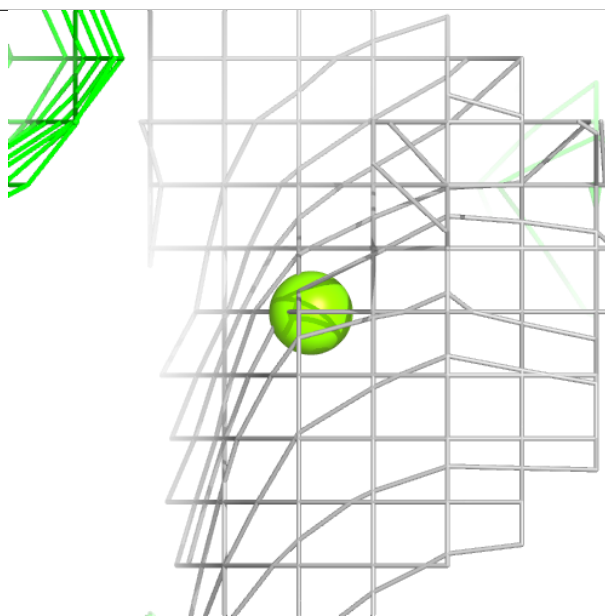
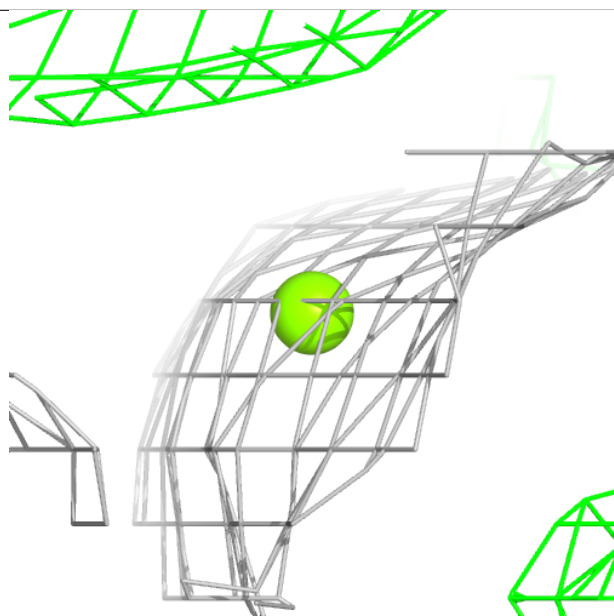
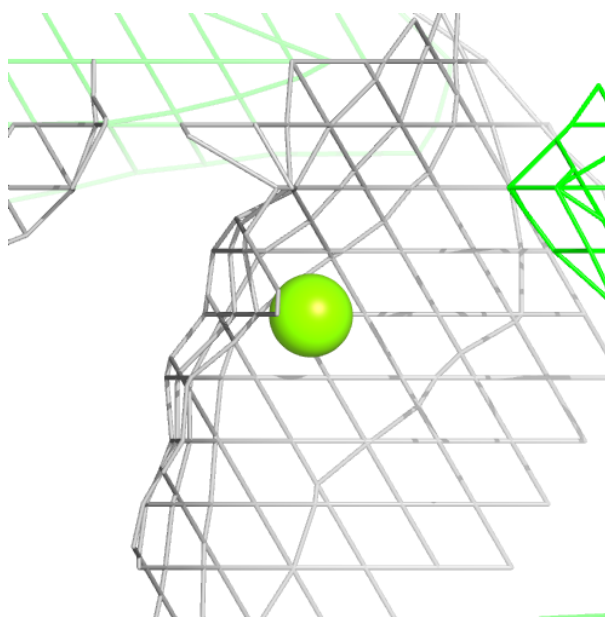
Electron density around 4BM C 402:

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



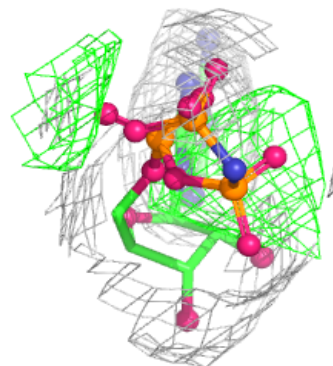
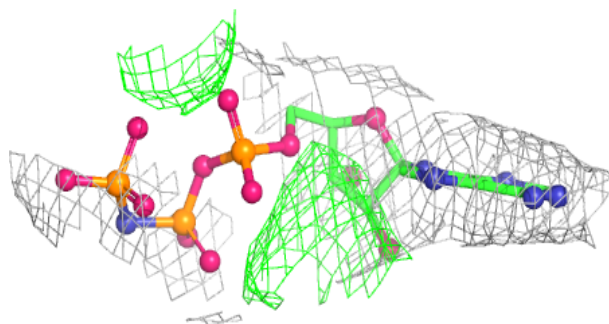
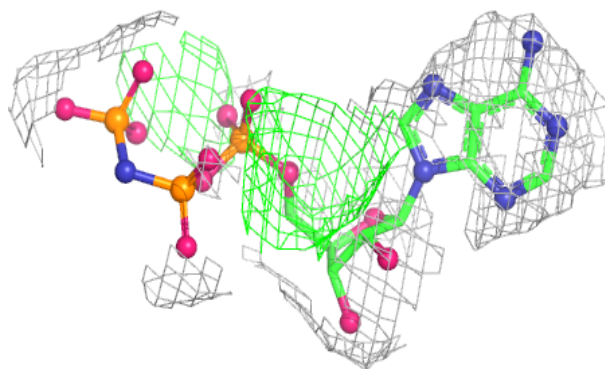
Electron density around MG A 902:

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



Electron density around ANP A 901:

$2mF_o-DF_c$ (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.