Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 10:15 AM EDT

PDB ID	:	4H3P
Title	:	Crystal structure of human ERK2 complexed with a MAPK docking peptide
Authors	:	Gogl, G.; Toeroe, I.; Remenyi, A.
Deposited on	:	2012-09-14
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 (i) 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 (i)) were used in the production of this report:

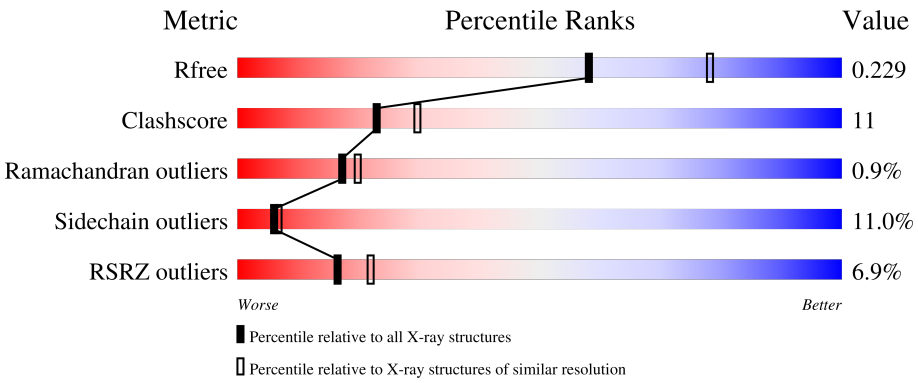
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 <=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	362	4%, 71%, 21%, •, 6%
1	D	362	5%, 66%, 24%, • •, 7%
2	B	24	29%, 29%, 29%, 8%, 33%
2	E	24	33%, 33%, 29%, 8%, 29%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5926 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 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	342	2741	1764	465	497	15	0	0	0
1	D	338	2707	1740	460	493	14	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P28482
A	0	SER	-	expression tag	UNP P28482
A	77	ALA	ARG	engineered mutation	UNP P28482
A	314	ALA	GLU	engineered mutation	UNP P28482
D	-1	GLY	-	expression tag	UNP P28482
D	0	SER	-	expression tag	UNP P28482
D	77	ALA	ARG	engineered mutation	UNP P28482
D	314	ALA	GLU	engineered mutation	UNP P28482

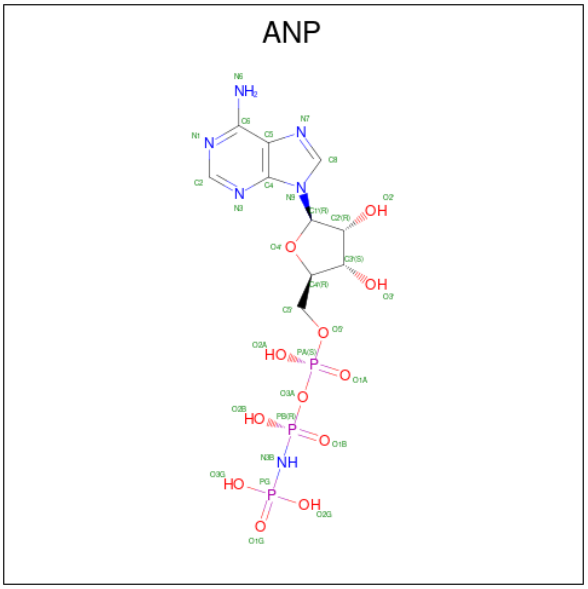
- Molecule 2 is a protein called Ribosomal protein S6 kinase alpha-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	16	117	74	23	20	0	0	0
2	E	17	122	78	24	20	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	719	ALA	SER	engineered mutation	UNP Q15418
B	724	ALA	GLN	engineered mutation	UNP Q15418
E	719	ALA	SER	engineered mutation	UNP Q15418
E	724	ALA	GLN	engineered mutation	UNP Q15418

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



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

- Molecule 4 is water.

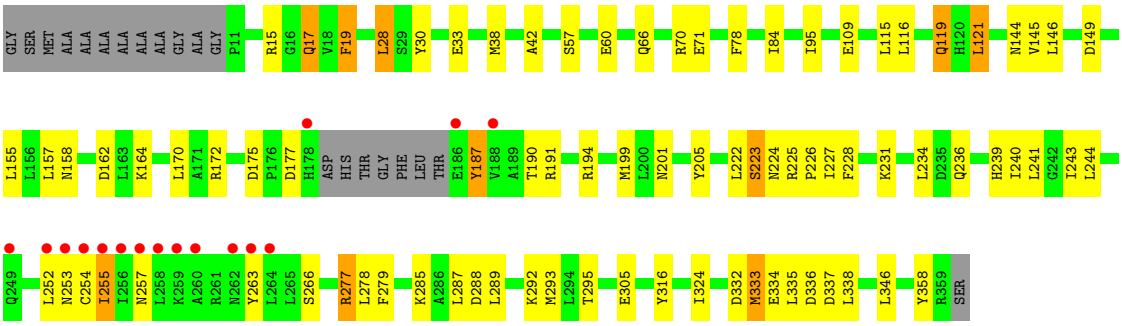
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	85	85	85	0	0
			Total	O		
4	D	92	92	92	0	0

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.

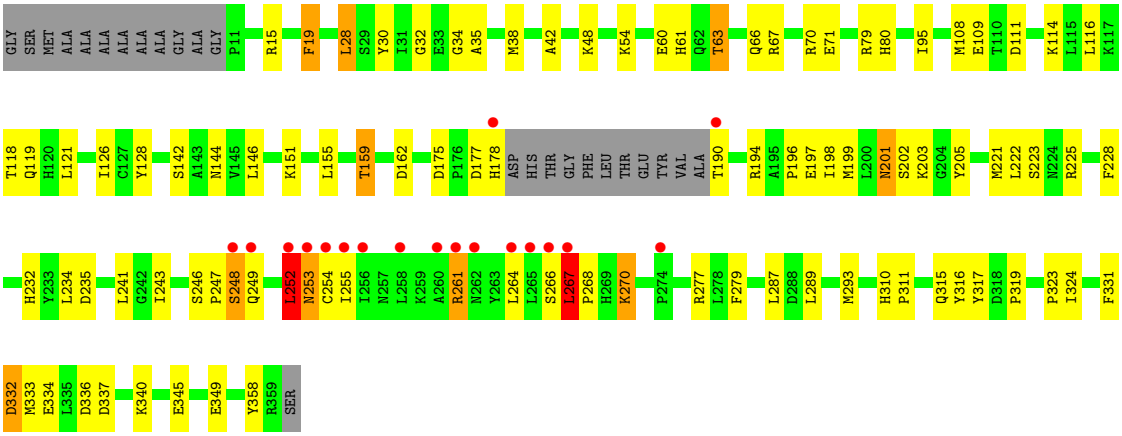
• Molecule 1: Mitogen-activated protein kinase 1

Chain A:



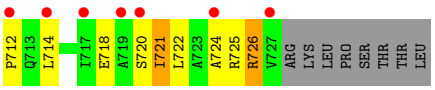
• Molecule 1: Mitogen-activated protein kinase 1

Chain D:



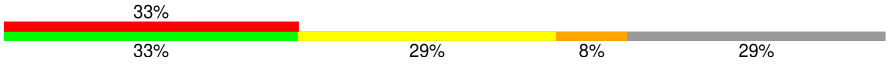
• Molecule 2: Ribosomal protein S6 kinase alpha-1

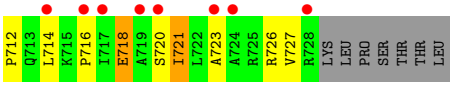
Chain B:



• Molecule 2: Ribosomal protein S6 kinase alpha-1

Chain E:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	41.48Å 58.77Å 79.18Å 100.93° 98.96° 90.01°	Depositor
Resolution (Å)	42.37 – 2.30 42.37 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.9 (42.37-2.30) 89.1 (42.37-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.77 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.178 , 0.224 0.182 , 0.229	Depositor DCC
R_{free} test set	1543 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	32.1	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 57.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5926	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	0.46	0/2807	0.65	0/3809
1	D	0.49	0/2772	0.66	2/3761 (0.1%)
2	B	0.32	0/118	0.84	1/159 (0.6%)
2	E	0.33	0/123	0.57	0/165
All	All	0.47	0/5820	0.66	3/7894 (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
1	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	267	LEU	CA-CB-CG	6.21	129.59	115.30
1	D	252	LEU	CB-CG-CD2	5.55	120.43	111.00
2	B	722	LEU	CA-CB-CG	-5.39	102.91	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	187	TYR	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	2741	0	2706	55	0
1	D	2707	0	2665	72	0
2	B	117	0	122	6	0
2	E	122	0	129	10	0
3	A	31	0	13	0	0
3	D	31	0	13	3	0
4	A	85	0	0	9	1
4	D	92	0	0	9	1
All	All	5926	0	5648	130	1

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 11.

All (130) 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:288:ASP:OD2	1:A:292:LYS:NZ	1.88	1.06
1:A:33:GLU:OE1	4:A:574:HOH:O	1.84	0.94
1:D:109:GLU:OE1	4:D:556:HOH:O	1.86	0.93
1:D:177:ASP:OD2	4:D:577:HOH:O	1.85	0.93
1:A:316:TYR:O	2:B:726:ARG:NH1	2.10	0.84
1:A:175:ASP:O	1:A:177:ASP:HA	1.79	0.82
1:D:119:GLN:NE2	2:E:712:PRO:O	2.13	0.81
1:A:19:PHE:HE1	1:A:38:MET:HE3	1.46	0.80
1:D:316:TYR:O	2:E:726:ARG:NH1	2.15	0.79
1:A:119:GLN:NE2	2:B:712:PRO:O	2.16	0.79
1:D:337:ASP:N	4:D:563:HOH:O	2.06	0.72
1:D:175:ASP:O	1:D:177:ASP:HA	1.90	0.71
2:E:716:PRO:HB2	2:E:718:GLU:HG2	1.75	0.69
1:A:277:ARG:CG	1:A:277:ARG:HH11	2.06	0.69
1:A:19:PHE:CE1	1:A:38:MET:HE3	2.27	0.68
1:D:34:GLY:O	4:D:573:HOH:O	2.11	0.67
1:A:289:LEU:O	1:A:293:MET:HG3	1.94	0.67
1:D:255:ILE:H	1:D:261:ARG:NH1	1.93	0.66
1:D:255:ILE:H	1:D:261:ARG:NH2	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ARG:HH11	1:A:277:ARG:HG2	1.62	0.65
1:A:334:GLU:HG2	1:D:61:HIS:CE1	2.33	0.63
1:D:255:ILE:H	1:D:261:ARG:CZ	2.12	0.63
1:D:177:ASP:O	1:D:178:HIS:CB	2.46	0.62
1:D:316:TYR:HA	2:E:726:ARG:HD2	1.81	0.62
1:D:128:TYR:CZ	2:E:723:ALA:HB2	2.35	0.61
1:D:126:ILE:HG12	1:D:221:MET:HB3	1.83	0.61
1:D:255:ILE:H	1:D:261:ARG:HH12	1.48	0.61
1:D:111:ASP:OD2	1:D:114:LYS:HE2	2.00	0.59
1:D:162:ASP:OD2	2:E:720:SER:OG	2.21	0.59
1:D:255:ILE:H	1:D:261:ARG:HH22	1.49	0.59
1:A:201:ASN:HD22	1:A:255:ILE:HA	1.68	0.58
1:A:277:ARG:HH11	1:A:277:ARG:CB	2.17	0.58
1:D:349:GLU:OE2	4:D:535:HOH:O	2.17	0.57
1:A:277:ARG:HG2	1:A:277:ARG:NH1	2.18	0.57
1:A:78:PHE:HZ	1:A:145:VAL:HG21	1.69	0.57
2:B:721:ILE:H	2:B:721:ILE:HD12	1.69	0.56
1:A:162:ASP:OD1	4:A:566:HOH:O	2.17	0.56
1:D:19:PHE:HE1	1:D:38:MET:HE3	1.71	0.56
1:A:243:ILE:HD12	1:A:278:LEU:HD11	1.88	0.56
1:A:334:GLU:OE2	1:A:336:ASP:HB2	2.06	0.56
1:D:334:GLU:HG3	4:D:563:HOH:O	2.07	0.55
1:D:142:SER:HB2	1:D:324:ILE:HG23	1.89	0.55
1:A:115:LEU:HD21	2:B:714:LEU:HD11	1.90	0.54
1:A:223:SER:HB3	1:A:279:PHE:CE1	2.43	0.54
1:D:95:ILE:HD11	1:D:345:GLU:HA	1.89	0.53
1:D:248:SER:O	1:D:252:LEU:N	2.32	0.53
1:A:15:ARG:O	1:A:17:GLN:NE2	2.37	0.53
1:D:63:THR:OG1	1:D:67:ARG:NH1	2.42	0.53
3:D:401:ANP:O1B	3:D:401:ANP:O1A	2.27	0.53
1:D:28:LEU:HD12	1:D:42:ALA:HB2	1.91	0.52
1:D:128:TYR:CE1	2:E:723:ALA:HB2	2.45	0.52
1:D:315:GLN:NE2	4:D:580:HOH:O	1.88	0.52
1:A:95:ILE:HG21	1:D:232:HIS:CD2	2.44	0.51
1:D:255:ILE:N	1:D:261:ARG:HH12	2.07	0.51
1:D:109:GLU:CD	1:D:159:THR:HG1	2.13	0.51
1:A:177:ASP:OD2	1:A:177:ASP:O	2.29	0.51
1:A:201:ASN:ND2	1:A:255:ILE:HA	2.26	0.51
1:D:66:GLN:O	1:D:70:ARG:HG3	2.11	0.51
1:A:239:HIS:HE1	4:A:533:HOH:O	1.94	0.50
1:D:109:GLU:OE1	1:D:159:THR:OG1	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:253:ASN:HA	1:D:261:ARG:NH1	2.26	0.50
1:A:227:ILE:HG23	1:A:228:PHE:CD2	2.46	0.50
1:A:191:ARG:HA	1:A:194:ARG:HG3	1.93	0.50
1:D:194:ARG:HD2	1:D:198:ILE:HG21	1.93	0.50
2:E:721:ILE:H	2:E:721:ILE:HD12	1.75	0.50
1:A:253:ASN:HA	1:A:255:ILE:HG22	1.93	0.50
1:A:66:GLN:O	1:A:70:ARG:HG3	2.12	0.50
1:D:177:ASP:OD2	1:D:177:ASP:O	2.29	0.50
1:A:305:GLU:OE1	4:A:525:HOH:O	2.18	0.49
1:D:316:TYR:HD1	2:E:726:ARG:HG3	1.75	0.49
1:A:277:ARG:NH1	4:A:541:HOH:O	2.45	0.49
1:A:28:LEU:HD12	1:A:42:ALA:HB2	1.95	0.48
1:D:255:ILE:N	1:D:261:ARG:HH22	2.09	0.48
1:A:223:SER:HB3	1:A:279:PHE:HE1	1.79	0.48
1:D:228:PHE:HE1	1:D:243:ILE:HG13	1.78	0.48
1:A:337:ASP:HB3	1:D:35:ALA:HB1	1.96	0.47
1:D:254:CYS:H	1:D:261:ARG:NH2	2.12	0.47
1:A:116:LEU:O	1:A:224:ASN:ND2	2.46	0.47
1:D:317:TYR:CZ	1:D:319:PRO:HD3	2.49	0.47
1:A:28:LEU:HD12	1:A:42:ALA:CB	2.45	0.47
1:A:84:ILE:HD12	1:A:164:LYS:HD3	1.97	0.47
1:D:80:HIS:CD2	1:D:323:PRO:HG2	2.49	0.46
1:D:196:PRO:HA	1:D:199:MET:HE2	1.96	0.46
1:D:201:ASN:C	1:D:201:ASN:HD22	2.19	0.46
1:D:228:PHE:CE1	1:D:243:ILE:HG13	2.50	0.46
1:D:254:CYS:H	1:D:261:ARG:HH22	1.64	0.46
1:D:289:LEU:O	1:D:293:MET:HG3	2.16	0.46
1:A:162:ASP:OD2	2:B:720:SER:OG	2.23	0.46
1:A:226:PRO:HA	4:A:508:HOH:O	2.15	0.45
1:D:66:GLN:NE2	1:D:336:ASP:HA	2.32	0.45
1:A:234:LEU:HD22	1:A:263:TYR:CZ	2.52	0.45
1:D:249:GLN:HA	1:D:252:LEU:HB2	1.99	0.45
1:D:151:LYS:NZ	1:D:190:THR:HG21	2.31	0.45
1:D:232:HIS:CE1	1:D:235:ASP:HB2	2.52	0.45
1:D:267:LEU:HB3	1:D:268:PRO:HD2	1.97	0.45
2:B:724:ALA:C	2:B:726:ARG:H	2.20	0.45
1:A:333:MET:HG2	1:A:338:LEU:HD11	1.98	0.45
1:A:17:GLN:HG3	1:A:38:MET:CE	2.47	0.44
1:D:247:PRO:HB2	1:D:252:LEU:HD22	1.98	0.44
1:D:340:LYS:HE2	4:D:583:HOH:O	2.17	0.44
1:A:191:ARG:HB3	1:A:236:GLN:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:SER:HA	1:D:247:PRO:HD3	1.80	0.44
1:D:267:LEU:CB	1:D:268:PRO:HD2	2.47	0.44
1:A:30:TYR:OH	4:A:539:HOH:O	1.99	0.43
1:A:255:ILE:HG12	1:A:257:ASN:O	2.18	0.43
1:D:268:PRO:HG2	1:D:270:LYS:HE3	2.00	0.43
1:A:17:GLN:OE1	4:A:535:HOH:O	2.22	0.43
1:A:228:PHE:CE1	1:A:240:ILE:HA	2.53	0.43
1:D:336:ASP:N	4:D:563:HOH:O	2.51	0.43
1:D:146:LEU:HD11	1:D:205:TYR:C	2.38	0.43
1:D:63:THR:O	1:D:67:ARG:HD2	2.19	0.43
1:D:108:MET:O	3:D:401:ANP:H2	2.19	0.42
1:D:197:GLU:O	1:D:202:SER:N	2.52	0.42
1:D:15:ARG:HD2	1:D:30:TYR:CD1	2.54	0.42
1:D:48:LYS:HZ2	1:D:48:LYS:HG2	1.57	0.42
1:A:146:LEU:HD11	1:A:205:TYR:C	2.41	0.42
1:A:17:GLN:HB3	4:A:535:HOH:O	2.20	0.41
1:A:109:GLU:HG3	1:A:158:ASN:OD1	2.21	0.41
1:D:151:LYS:HZ3	1:D:190:THR:HG21	1.85	0.41
1:D:223:SER:HB3	1:D:279:PHE:CE1	2.55	0.41
1:D:116:LEU:HD23	1:D:116:LEU:HA	1.86	0.41
1:D:310:HIS:CG	1:D:311:PRO:HD2	2.55	0.41
1:A:121:LEU:HD12	1:A:121:LEU:HA	1.77	0.41
1:A:335:LEU:HD23	1:A:335:LEU:HA	1.84	0.41
1:A:244:LEU:HD23	1:A:244:LEU:HA	1.90	0.40
1:D:32:GLY:HA3	3:D:401:ANP:H4'	2.02	0.40
1:A:149:ASP:HB2	1:A:170:LEU:HD12	2.03	0.40
1:D:331:PHE:O	1:D:332:ASP:HB2	2.22	0.40
1:A:253:ASN:C	1:A:255:ILE:H	2.15	0.40
2:E:723:ALA:O	2:E:726:ARG:HB2	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:581:HOH:O	4:D:590:HOH:O[1_655]	2.15	0.05

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	338/362 (93%)	322 (95%)	15 (4%)	1 (0%)	41	50
1	D	334/362 (92%)	320 (96%)	11 (3%)	3 (1%)	17	20
2	B	14/24 (58%)	11 (79%)	2 (14%)	1 (7%)	1	0
2	E	15/24 (62%)	14 (93%)	0	1 (7%)	1	0
All	All	701/772 (91%)	667 (95%)	28 (4%)	6 (1%)	17	20

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	253	ASN
1	D	332	ASP
1	A	332	ASP
1	D	252	LEU
2	B	725	ARG
2	E	727	VAL

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	295/317 (93%)	263 (89%)	32 (11%)	6	7
1	D	291/317 (92%)	262 (90%)	29 (10%)	7	9
2	B	11/21 (52%)	8 (73%)	3 (27%)	0	0
2	E	11/21 (52%)	8 (73%)	3 (27%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	608/676 (90%)	541 (89%)	67 (11%)	6 7

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	19	PHE
1	A	28	LEU
1	A	57	SER
1	A	60	GLU
1	A	71	GLU
1	A	119	GLN
1	A	121	LEU
1	A	144	ASN
1	A	155	LEU
1	A	157	LEU
1	A	172	ARG
1	A	187	TYR
1	A	190	THR
1	A	199	MET
1	A	222	LEU
1	A	223	SER
1	A	225	ARG
1	A	231	LYS
1	A	241	LEU
1	A	252	LEU
1	A	254	CYS
1	A	255	ILE
1	A	266	SER
1	A	277	ARG
1	A	285	LYS
1	A	287	LEU
1	A	295	THR
1	A	324	ILE
1	A	333	MET
1	A	346	LEU
1	A	358	TYR
1	D	19	PHE
1	D	28	LEU
1	D	54	LYS
1	D	60	GLU
1	D	63	THR

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Mol	Chain	Res	Type
1	D	71	GLU
1	D	79	ARG
1	D	118	THR
1	D	121	LEU
1	D	144	ASN
1	D	155	LEU
1	D	159	THR
1	D	201	ASN
1	D	203	LYS
1	D	222	LEU
1	D	225	ARG
1	D	234	LEU
1	D	241	LEU
1	D	248	SER
1	D	252	LEU
1	D	261	ARG
1	D	264	LEU
1	D	266	SER
1	D	267	LEU
1	D	270	LYS
1	D	277	ARG
1	D	287	LEU
1	D	333	MET
1	D	358	TYR
2	B	718	GLU
2	B	721	ILE
2	B	726	ARG
2	E	714	LEU
2	E	718	GLU
2	E	721	ILE

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	119	GLN
1	A	201	ASN
1	D	61	HIS
1	D	66	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
3	ANP	A	401	-	29,33,33	2.55	7 (24%)	31,52,52	2.71	6 (19%)
3	ANP	D	401	-	29,33,33	2.55	7 (24%)	31,52,52	2.64	9 (29%)

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	401	-	-	8/14/38/38	0/3/3/3
3	ANP	D	401	-	-	8/14/38/38	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	401	ANP	PB-O1B	9.46	1.60	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	ANP	PB-O1B	9.46	1.60	1.46
3	D	401	ANP	PB-N3B	4.79	1.75	1.63
3	A	401	ANP	PB-N3B	4.72	1.75	1.63
3	D	401	ANP	PG-O1G	3.90	1.52	1.46
3	A	401	ANP	PG-O1G	3.89	1.52	1.46
3	A	401	ANP	C2'-C3'	-3.69	1.43	1.53
3	D	401	ANP	C2'-C3'	-3.63	1.43	1.53
3	D	401	ANP	C6-N6	3.10	1.45	1.34
3	A	401	ANP	O4'-C4'	-2.95	1.38	1.45
3	A	401	ANP	C6-N6	2.89	1.44	1.34
3	D	401	ANP	O4'-C4'	-2.71	1.39	1.45
3	D	401	ANP	PA-O3A	-2.37	1.56	1.59
3	A	401	ANP	C2-N3	2.19	1.35	1.32

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	401	ANP	O1B-PB-N3B	-11.29	95.14	111.77
3	D	401	ANP	O1B-PB-N3B	-10.86	95.78	111.77
3	D	401	ANP	O2B-PB-O1B	4.77	120.11	109.87
3	A	401	ANP	O2B-PB-O1B	4.75	120.06	109.87
3	A	401	ANP	N3-C2-N1	-4.70	122.30	128.67
3	A	401	ANP	O2B-PB-O3A	4.55	119.82	104.64
3	D	401	ANP	N3-C2-N1	-4.12	123.09	128.67
3	D	401	ANP	O2B-PB-O3A	4.07	118.22	104.64
3	D	401	ANP	O5'-C5'-C4'	2.98	119.15	108.99
3	A	401	ANP	O5'-C5'-C4'	2.96	119.06	108.99
3	A	401	ANP	O3A-PB-N3B	-2.61	99.36	106.59
3	D	401	ANP	C4-C5-N7	-2.56	106.64	109.34
3	D	401	ANP	O3A-PB-N3B	-2.41	99.91	106.59
3	D	401	ANP	C2'-C3'-C4'	2.29	107.04	102.61
3	D	401	ANP	O4'-C1'-N9	2.24	111.71	108.75

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ANP	PB-N3B-PG-O1G
3	A	401	ANP	PG-N3B-PB-O3A
3	A	401	ANP	C5'-O5'-PA-O1A
3	A	401	ANP	C5'-O5'-PA-O2A
3	A	401	ANP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
3	A	401	ANP	O4'-C4'-C5'-O5'
3	A	401	ANP	C3'-C4'-C5'-O5'
3	D	401	ANP	PB-N3B-PG-O1G
3	D	401	ANP	PG-N3B-PB-O3A
3	D	401	ANP	PA-O3A-PB-O1B
3	D	401	ANP	C5'-O5'-PA-O2A
3	D	401	ANP	O4'-C4'-C5'-O5'
3	D	401	ANP	C3'-C4'-C5'-O5'
3	D	401	ANP	C5'-O5'-PA-O1A
3	D	401	ANP	C5'-O5'-PA-O3A
3	A	401	ANP	PA-O3A-PB-O1B

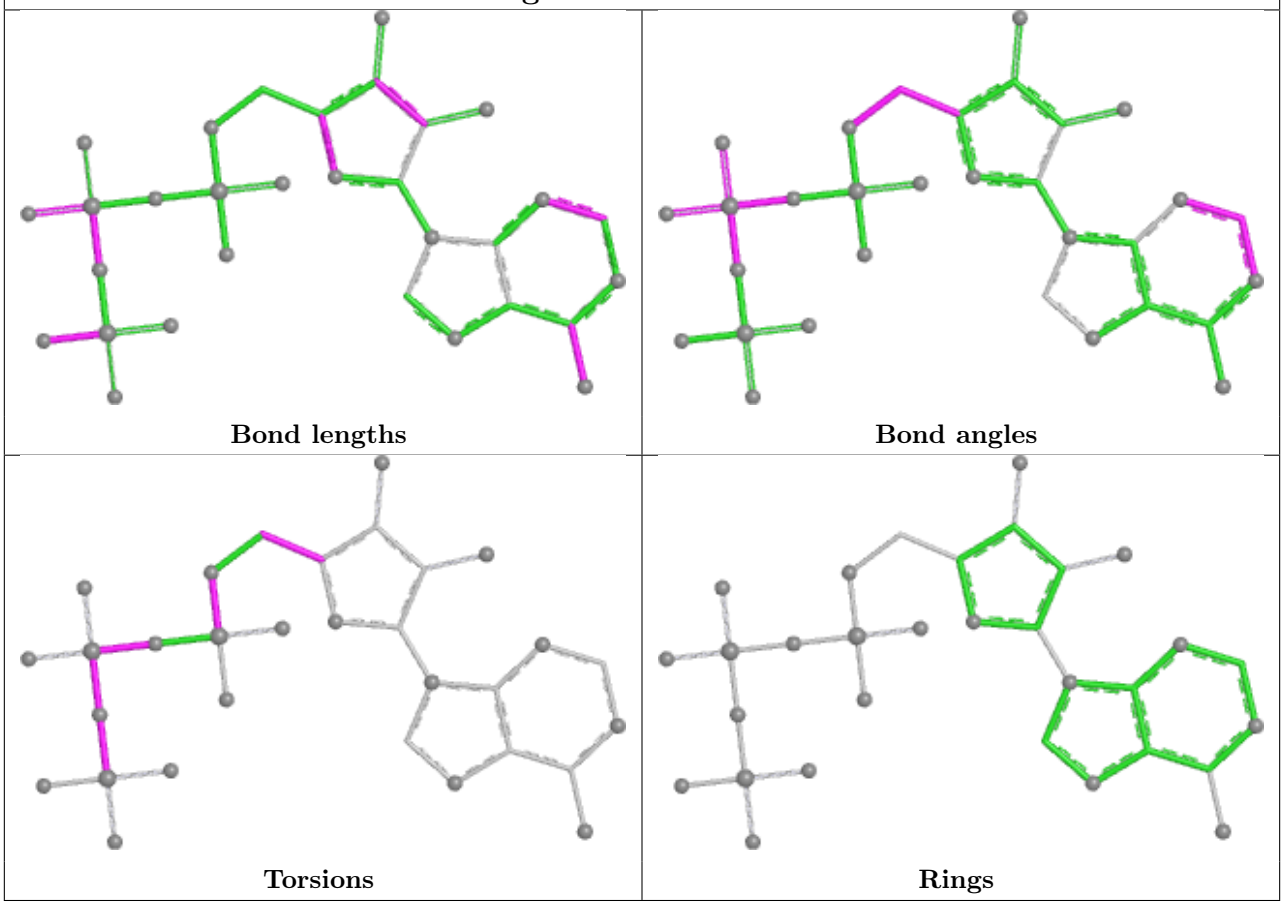
There are no ring outliers.

1 monomer is involved in 3 short contacts:

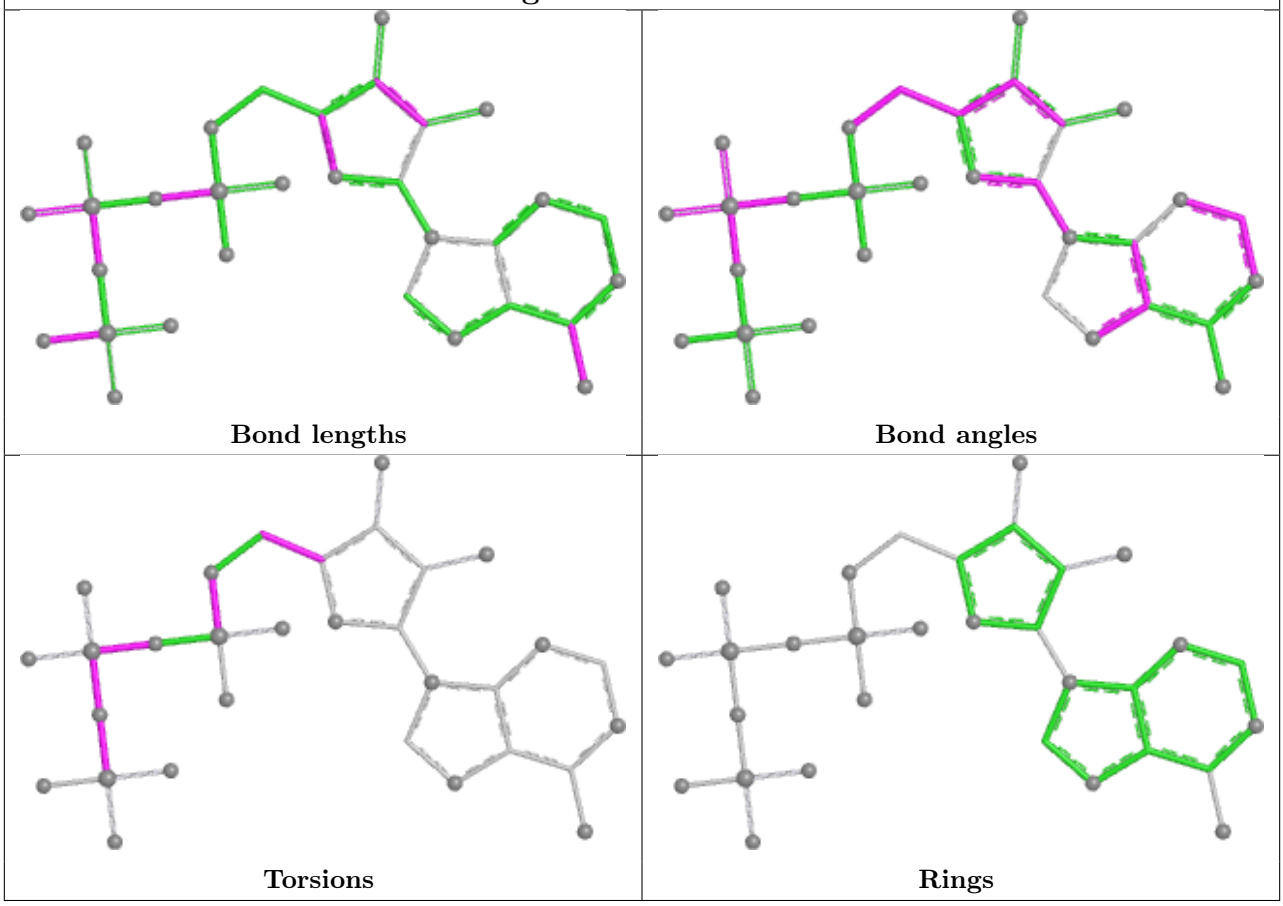
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	401	ANP	3	0

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 then 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.

Ligand ANP A 401



Ligand ANP D 401



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	342/362 (94%)	0.07	16 (4%)	31	38	19, 40, 80, 124	0
1	D	338/362 (93%)	0.08	18 (5%)	26	33	19, 39, 77, 151	0
2	B	16/24 (66%)	1.87	7 (43%)	0	0	65, 81, 105, 106	0
2	E	17/24 (70%)	2.02	8 (47%)	0	0	67, 78, 97, 118	0
All	All	713/772 (92%)	0.16	49 (6%)	16	22	19, 40, 88, 151	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	256	ILE	9.9
1	A	256	ILE	8.3
1	D	254	CYS	5.7
1	A	254	CYS	5.6
2	B	719	ALA	5.4
2	E	719	ALA	5.3
1	D	260	ALA	4.9
2	B	724	ALA	4.8
1	D	261	ARG	4.7
2	E	724	ALA	4.7
1	D	253	ASN	4.6
1	A	260	ALA	4.5
2	E	728	ARG	4.4
2	B	727	VAL	4.4
1	D	264	LEU	4.3
1	A	258	LEU	4.0
1	D	258	LEU	3.8
1	A	252	LEU	3.6
1	A	257	ASN	3.3
2	B	712	PRO	3.3
1	D	255	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	252	LEU	2.9
1	D	249	GLN	2.8
1	D	266	SER	2.8
1	A	186	GLU	2.7
1	A	253	ASN	2.6
1	D	178	HIS	2.6
1	D	267	LEU	2.6
2	E	717	ILE	2.6
1	A	264	LEU	2.5
1	A	262	ASN	2.4
1	D	190	THR	2.4
1	A	263	TYR	2.4
2	B	720	SER	2.4
1	A	249	GLN	2.4
1	A	259	LYS	2.3
2	E	720	SER	2.4
2	B	717	ILE	2.3
2	E	714	LEU	2.3
1	A	188	VAL	2.3
2	E	723	ALA	2.2
1	A	178	HIS	2.2
1	D	274	PRO	2.2
2	E	716	PRO	2.1
1	A	255	ILE	2.0
1	D	265	LEU	2.0
1	D	262	ASN	2.0
1	D	248	SER	2.0
2	B	714	LEU	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 monosaccharides 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,

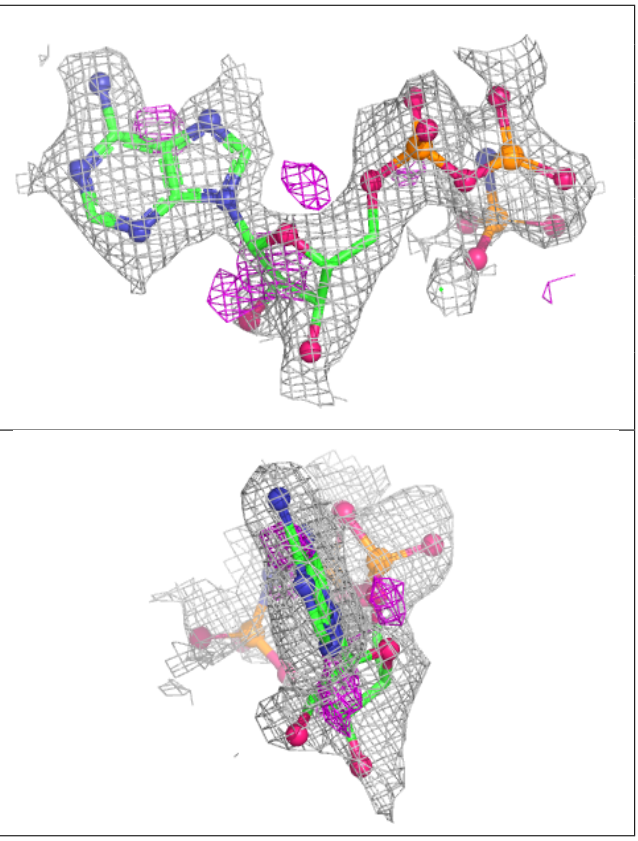
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ANP	D	401	31/31	0.75	0.24	26,77,172,249	0
3	ANP	A	401	31/31	0.82	0.24	22,77,202,212	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.

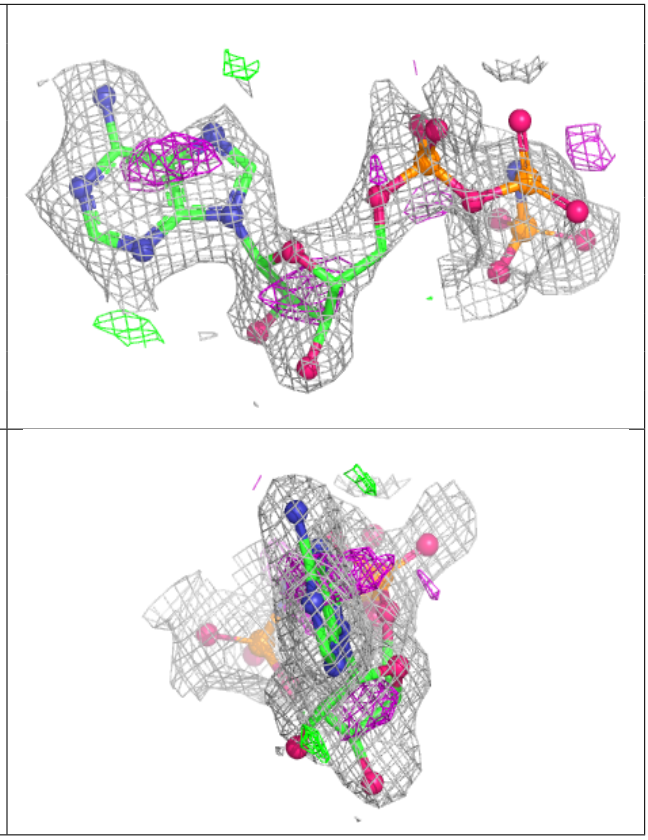
Electron density around ANP D 401:

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 401:

$2mF_o$-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.