



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:22 AM UTC

PDB ID : 5ZCT / pdb_00005zct
Title : The crystal structure of the poly-alpha-L-glutamate peptides synthetase RimK at 2.05 angstrom resolution.
Authors : Arimura, Y.; Kono, T.; Kino, K.; Kurumizaka, H.
Deposited on : 2018-02-20
Resolution : 2.05 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

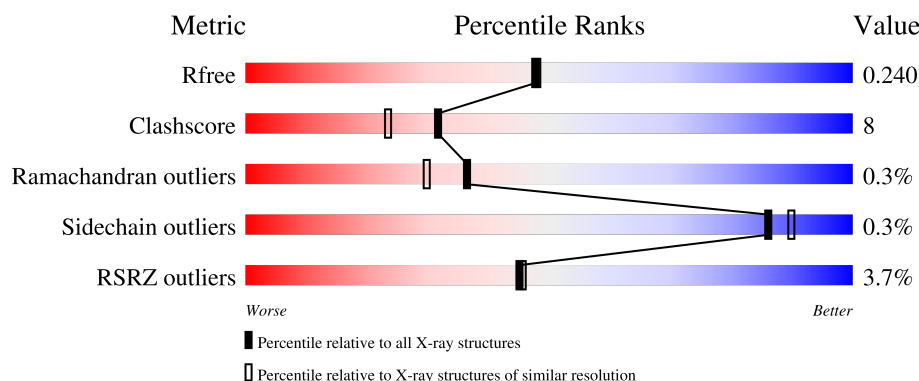
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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.



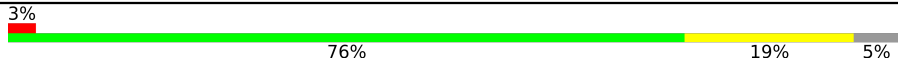
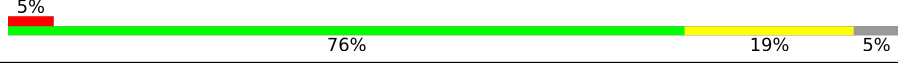
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	306	0% 86% 9% 5%
1	B	306	12% 72% 23% 5%
1	C	306	2% 80% 14% 5%
1	D	306	3% 76% 19% 5%
1	E	306	2% 82% 13% 5%

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Mol	Chain	Length	Quality of chain
1	F	306	
1	G	306	
1	H	306	

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
4	SO4	F	403	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 19202 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 Ribosomal protein S6–L-glutamate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	B	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	C	290	Total	C	N	O	S	0	0	0
			2199	1376	404	407	12			
1	D	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	E	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	F	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	G	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			
1	H	292	Total	C	N	O	S	0	0	0
			2213	1384	406	411	12			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	293	GLU	-	expression tag	UNP P0C0U4
A	294	TYR	-	expression tag	UNP P0C0U4
A	295	CYS	-	expression tag	UNP P0C0U4
A	296	LEU	-	expression tag	UNP P0C0U4
A	297	LYS	-	expression tag	UNP P0C0U4
A	298	THR	-	expression tag	UNP P0C0U4
A	299	GLY	-	expression tag	UNP P0C0U4
A	300	GLY	-	expression tag	UNP P0C0U4
A	301	GLY	-	expression tag	UNP P0C0U4
A	302	THR	-	expression tag	UNP P0C0U4
A	303	LEU	-	expression tag	UNP P0C0U4
A	304	VAL	-	expression tag	UNP P0C0U4
A	305	PRO	-	expression tag	UNP P0C0U4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	306	ARG	-	expression tag	UNP P0C0U4
B	293	GLU	-	expression tag	UNP P0C0U4
B	294	TYR	-	expression tag	UNP P0C0U4
B	295	CYS	-	expression tag	UNP P0C0U4
B	296	LEU	-	expression tag	UNP P0C0U4
B	297	LYS	-	expression tag	UNP P0C0U4
B	298	THR	-	expression tag	UNP P0C0U4
B	299	GLY	-	expression tag	UNP P0C0U4
B	300	GLY	-	expression tag	UNP P0C0U4
B	301	GLY	-	expression tag	UNP P0C0U4
B	302	THR	-	expression tag	UNP P0C0U4
B	303	LEU	-	expression tag	UNP P0C0U4
B	304	VAL	-	expression tag	UNP P0C0U4
B	305	PRO	-	expression tag	UNP P0C0U4
B	306	ARG	-	expression tag	UNP P0C0U4
C	293	GLU	-	expression tag	UNP P0C0U4
C	294	TYR	-	expression tag	UNP P0C0U4
C	295	CYS	-	expression tag	UNP P0C0U4
C	296	LEU	-	expression tag	UNP P0C0U4
C	297	LYS	-	expression tag	UNP P0C0U4
C	298	THR	-	expression tag	UNP P0C0U4
C	299	GLY	-	expression tag	UNP P0C0U4
C	300	GLY	-	expression tag	UNP P0C0U4
C	301	GLY	-	expression tag	UNP P0C0U4
C	302	THR	-	expression tag	UNP P0C0U4
C	303	LEU	-	expression tag	UNP P0C0U4
C	304	VAL	-	expression tag	UNP P0C0U4
C	305	PRO	-	expression tag	UNP P0C0U4
C	306	ARG	-	expression tag	UNP P0C0U4
D	293	GLU	-	expression tag	UNP P0C0U4
D	294	TYR	-	expression tag	UNP P0C0U4
D	295	CYS	-	expression tag	UNP P0C0U4
D	296	LEU	-	expression tag	UNP P0C0U4
D	297	LYS	-	expression tag	UNP P0C0U4
D	298	THR	-	expression tag	UNP P0C0U4
D	299	GLY	-	expression tag	UNP P0C0U4
D	300	GLY	-	expression tag	UNP P0C0U4
D	301	GLY	-	expression tag	UNP P0C0U4
D	302	THR	-	expression tag	UNP P0C0U4
D	303	LEU	-	expression tag	UNP P0C0U4
D	304	VAL	-	expression tag	UNP P0C0U4
D	305	PRO	-	expression tag	UNP P0C0U4

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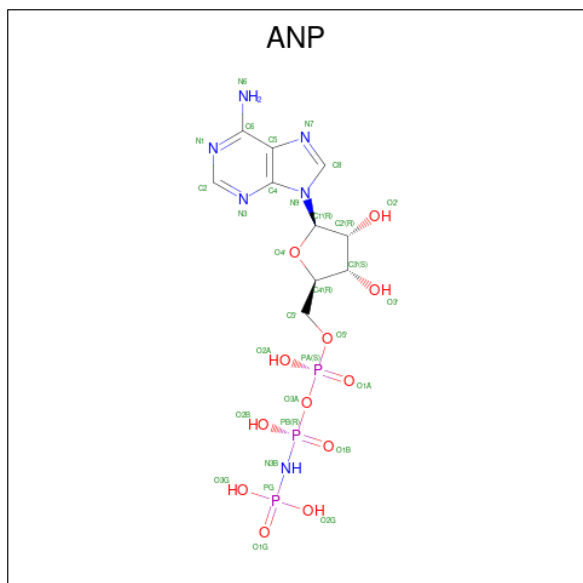
Chain	Residue	Modelled	Actual	Comment	Reference
D	306	ARG	-	expression tag	UNP P0C0U4
E	293	GLU	-	expression tag	UNP P0C0U4
E	294	TYR	-	expression tag	UNP P0C0U4
E	295	CYS	-	expression tag	UNP P0C0U4
E	296	LEU	-	expression tag	UNP P0C0U4
E	297	LYS	-	expression tag	UNP P0C0U4
E	298	THR	-	expression tag	UNP P0C0U4
E	299	GLY	-	expression tag	UNP P0C0U4
E	300	GLY	-	expression tag	UNP P0C0U4
E	301	GLY	-	expression tag	UNP P0C0U4
E	302	THR	-	expression tag	UNP P0C0U4
E	303	LEU	-	expression tag	UNP P0C0U4
E	304	VAL	-	expression tag	UNP P0C0U4
E	305	PRO	-	expression tag	UNP P0C0U4
E	306	ARG	-	expression tag	UNP P0C0U4
F	293	GLU	-	expression tag	UNP P0C0U4
F	294	TYR	-	expression tag	UNP P0C0U4
F	295	CYS	-	expression tag	UNP P0C0U4
F	296	LEU	-	expression tag	UNP P0C0U4
F	297	LYS	-	expression tag	UNP P0C0U4
F	298	THR	-	expression tag	UNP P0C0U4
F	299	GLY	-	expression tag	UNP P0C0U4
F	300	GLY	-	expression tag	UNP P0C0U4
F	301	GLY	-	expression tag	UNP P0C0U4
F	302	THR	-	expression tag	UNP P0C0U4
F	303	LEU	-	expression tag	UNP P0C0U4
F	304	VAL	-	expression tag	UNP P0C0U4
F	305	PRO	-	expression tag	UNP P0C0U4
F	306	ARG	-	expression tag	UNP P0C0U4
G	293	GLU	-	expression tag	UNP P0C0U4
G	294	TYR	-	expression tag	UNP P0C0U4
G	295	CYS	-	expression tag	UNP P0C0U4
G	296	LEU	-	expression tag	UNP P0C0U4
G	297	LYS	-	expression tag	UNP P0C0U4
G	298	THR	-	expression tag	UNP P0C0U4
G	299	GLY	-	expression tag	UNP P0C0U4
G	300	GLY	-	expression tag	UNP P0C0U4
G	301	GLY	-	expression tag	UNP P0C0U4
G	302	THR	-	expression tag	UNP P0C0U4
G	303	LEU	-	expression tag	UNP P0C0U4
G	304	VAL	-	expression tag	UNP P0C0U4
G	305	PRO	-	expression tag	UNP P0C0U4

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Chain	Residue	Modelled	Actual	Comment	Reference
G	306	ARG	-	expression tag	UNP P0C0U4
H	293	GLU	-	expression tag	UNP P0C0U4
H	294	TYR	-	expression tag	UNP P0C0U4
H	295	CYS	-	expression tag	UNP P0C0U4
H	296	LEU	-	expression tag	UNP P0C0U4
H	297	LYS	-	expression tag	UNP P0C0U4
H	298	THR	-	expression tag	UNP P0C0U4
H	299	GLY	-	expression tag	UNP P0C0U4
H	300	GLY	-	expression tag	UNP P0C0U4
H	301	GLY	-	expression tag	UNP P0C0U4
H	302	THR	-	expression tag	UNP P0C0U4
H	303	LEU	-	expression tag	UNP P0C0U4
H	304	VAL	-	expression tag	UNP P0C0U4
H	305	PRO	-	expression tag	UNP P0C0U4
H	306	ARG	-	expression tag	UNP P0C0U4

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



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

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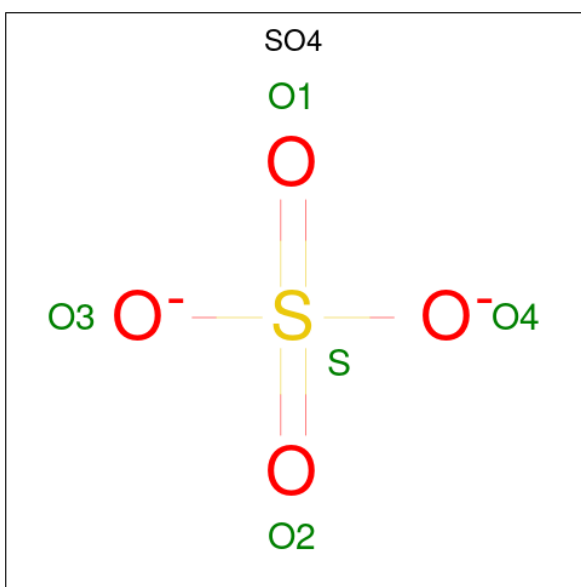
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	E	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	F	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	G	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	H	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		
3	G	1	Total	Mg	0	0
			1	1		
3	H	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		

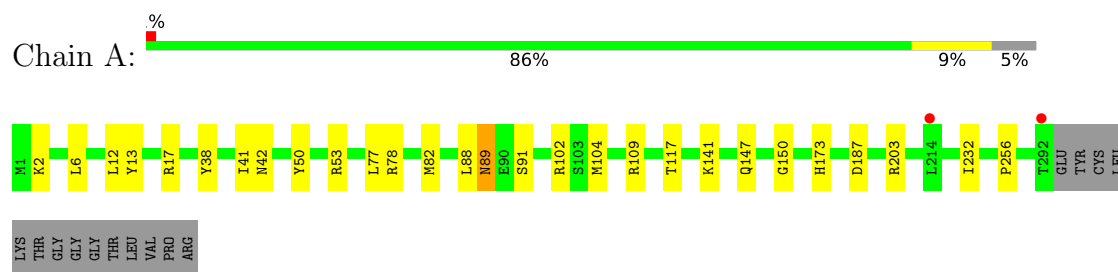
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	237	Total 237	O 237	0	0
5	B	72	Total 72	O 72	0	0
5	C	166	Total 166	O 166	0	0
5	D	94	Total 94	O 94	0	0
5	E	167	Total 167	O 167	0	0
5	F	108	Total 108	O 108	0	0
5	G	243	Total 243	O 243	0	0
5	H	99	Total 99	O 99	0	0

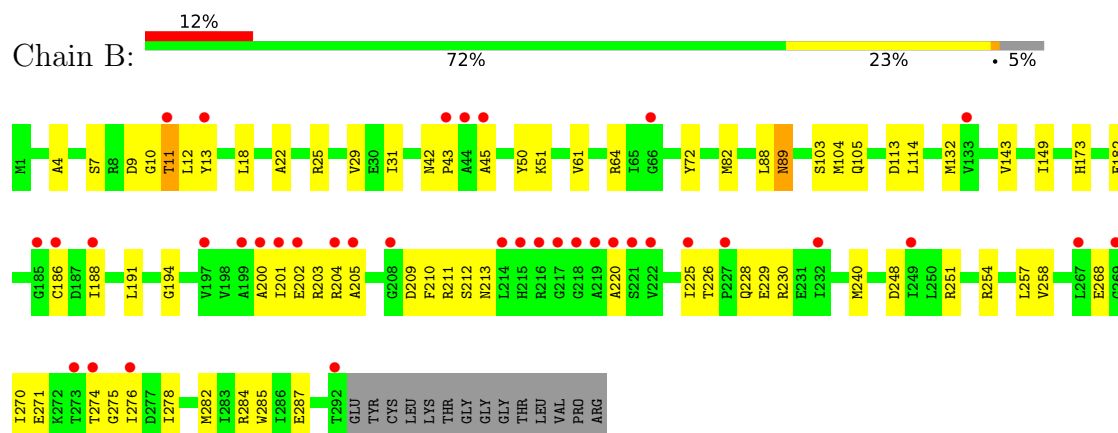
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

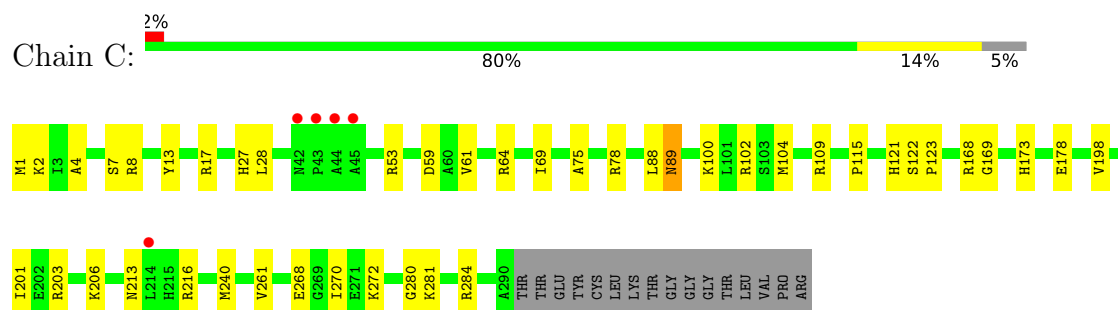
- Molecule 1: Ribosomal protein S6–L-glutamate ligase



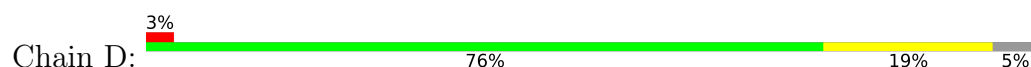
- Molecule 1: Ribosomal protein S6–L-glutamate ligase

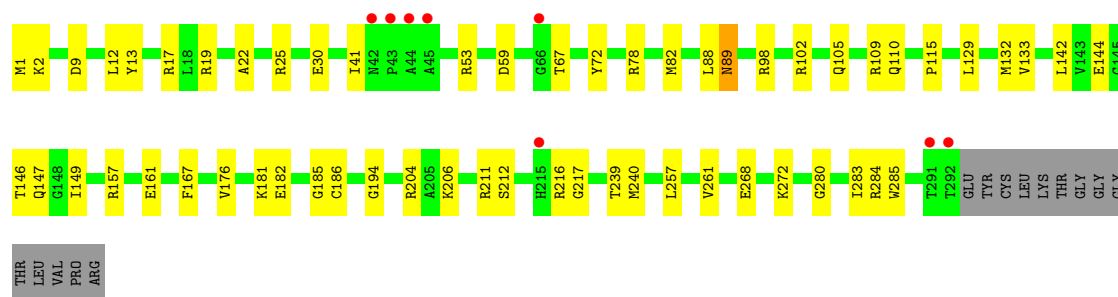


- Molecule 1: Ribosomal protein S6–L-glutamate ligase

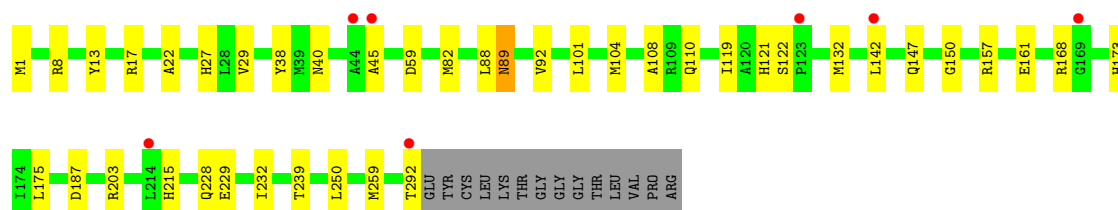
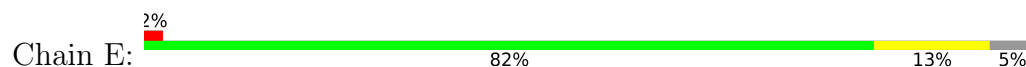


- Molecule 1: Ribosomal protein S6–L-glutamate ligase

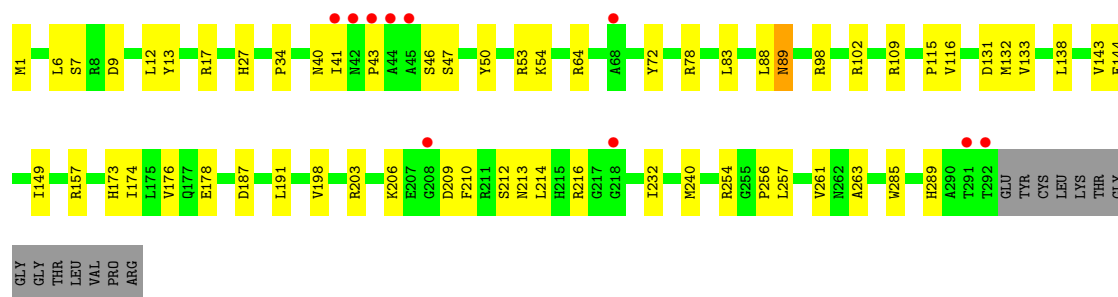
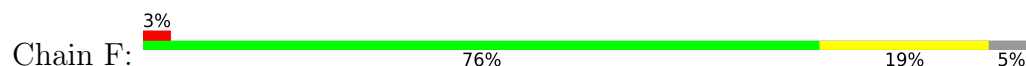




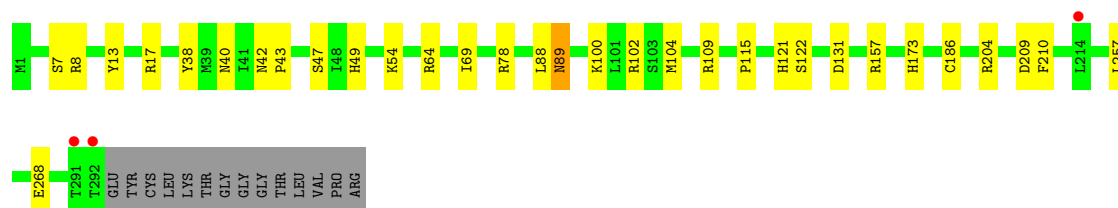
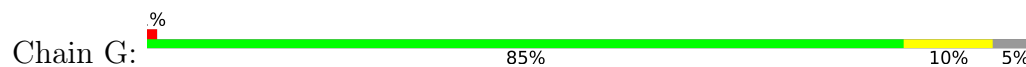
- Molecule 1: Ribosomal protein S6-L-glutamate ligase



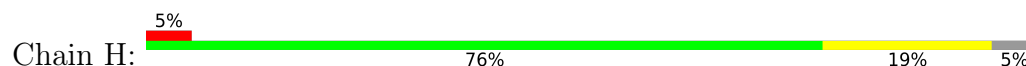
- Molecule 1: Ribosomal protein S6-L-glutamate ligase

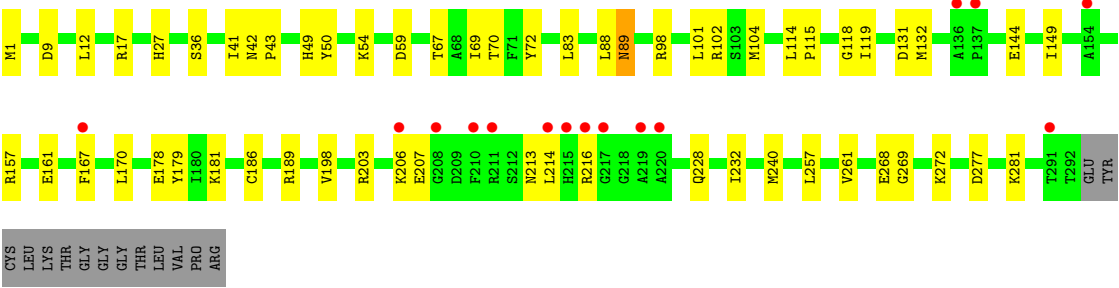


- Molecule 1: Ribosomal protein S6-L-glutamate ligase



- Molecule 1: Ribosomal protein S6-L-glutamate ligase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	212.62Å 121.63Å 119.74Å 90.00° 118.74° 90.00°	Depositor
Resolution (Å)	48.21 – 2.05 48.21 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.2 (48.21-2.05) 98.2 (48.21-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.05Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.196 , 0.240 0.197 , 0.240	Depositor DCC
R_{free} test set	8478 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19202	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: SO4, 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	0.42	0/2243	0.67	2/3032 (0.1%)
1	B	0.34	0/2243	0.64	0/3032
1	C	0.37	0/2229	0.61	0/3012
1	D	0.34	0/2243	0.60	0/3032
1	E	0.40	0/2243	0.62	0/3032
1	F	0.35	0/2243	0.60	0/3032
1	G	0.43	0/2243	0.64	0/3032
1	H	0.36	0/2243	0.61	0/3032
All	All	0.38	0/17930	0.62	2/24236 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	GLN	CA-C-N	-6.71	108.50	121.78
1	A	147	GLN	C-N-CA	-6.71	108.50	121.78

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	2213	0	2278	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2213	0	2279	61	0
1	C	2199	0	2267	29	0
1	D	2213	0	2281	45	0
1	E	2213	0	2278	34	0
1	F	2213	0	2281	48	0
1	G	2213	0	2279	24	0
1	H	2213	0	2280	54	0
2	A	31	0	13	1	0
2	B	31	0	13	1	0
2	C	31	0	13	0	0
2	D	31	0	13	0	0
2	E	31	0	13	0	0
2	F	31	0	13	0	0
2	G	31	0	13	0	0
2	H	31	0	13	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	15	0	0	1	0
4	B	5	0	0	0	0
4	C	15	0	0	2	0
4	D	5	0	0	1	0
4	E	10	0	0	0	0
4	F	5	0	0	2	0
4	G	10	0	0	1	0
4	H	5	0	0	1	0
5	A	237	0	0	2	0
5	B	72	0	0	0	0
5	C	166	0	0	0	0
5	D	94	0	0	1	0
5	E	167	0	0	2	0
5	F	108	0	0	4	0
5	G	243	0	0	1	0
5	H	99	0	0	2	0
All	All	19202	0	18327	293	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 (293) 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:C:206:LYS:HE2	1:C:216:ARG:O	1.67	0.94
1:D:206:LYS:NZ	1:D:211:ARG:O	2.01	0.93
1:F:9:ASP:HB3	1:F:12:LEU:HD23	1.60	0.83
1:F:144:GLU:HG3	1:F:149:ILE:HD11	1.62	0.81
1:D:206:LYS:HE2	1:D:216:ARG:C	2.06	0.79
1:G:104:MET:HE3	1:H:41:ILE:HD11	1.63	0.78
1:C:104:MET:HE3	1:D:41:ILE:HD11	1.64	0.78
1:B:104:MET:HE3	1:B:114:LEU:HD13	1.64	0.78
1:B:186:CYS:HB3	1:B:204:ARG:HG3	1.63	0.77
1:E:8:ARG:HG2	1:E:8:ARG:HH11	1.50	0.76
1:H:189:ARG:HH21	1:H:203:ARG:HH12	1.32	0.74
1:H:189:ARG:HH21	1:H:203:ARG:NH1	1.87	0.73
1:F:102:ARG:NH2	4:F:403:SO4:O4	2.23	0.72
1:F:43:PRO:HA	1:F:46:SER:HB3	1.71	0.72
1:B:274:THR:HB	1:B:276:ILE:HG13	1.72	0.71
1:B:209:ASP:OD1	1:B:210:PHE:N	2.23	0.71
1:C:203:ARG:NH1	4:C:403:SO4:O4	2.25	0.70
1:F:213:ASN:OD1	1:F:216:ARG:NH1	2.24	0.70
1:B:225:ILE:HD11	1:B:230:ARG:HG3	1.72	0.70
1:D:211:ARG:NH1	1:D:216:ARG:HE	1.91	0.68
1:F:203:ARG:HD3	1:F:214:LEU:HG	1.74	0.67
1:A:232:ILE:HD12	1:A:256:PRO:HG2	1.76	0.67
1:D:144:GLU:HG3	1:D:149:ILE:HD11	1.75	0.67
1:C:1:MET:HG3	1:C:59:ASP:HB2	1.78	0.66
1:B:113:ASP:OD2	1:B:257:LEU:HD23	1.95	0.66
1:E:108:ALA:HB1	1:F:43:PRO:HD3	1.78	0.65
1:D:280:GLY:O	1:D:284:ARG:HG3	1.96	0.64
1:A:102:ARG:NH2	4:A:405:SO4:O3	2.27	0.64
1:E:1:MET:HG3	1:E:59:ASP:HB2	1.78	0.64
1:H:189:ARG:HD2	1:H:203:ARG:HH11	1.63	0.64
1:H:131:ASP:OD1	1:H:157:ARG:NH2	2.31	0.64
1:B:209:ASP:OD1	1:B:211:ARG:N	2.23	0.64
1:B:202:GLU:O	1:B:220:ALA:HA	1.98	0.63
1:B:186:CYS:HB2	1:B:202:GLU:OE1	1.98	0.63
1:B:45:ALA:HB2	1:B:82:MET:HE2	1.80	0.63
1:B:4:ALA:HB3	1:B:61:VAL:HG22	1.80	0.63
1:B:50:TYR:CE2	1:B:51:LYS:HD2	2.33	0.63
1:D:78:ARG:NH2	5:D:501:HOH:O	2.28	0.62
1:C:173:HIS:CE1	1:D:72:TYR:HB2	2.34	0.62
1:C:123:PRO:O	1:C:168:ARG:NH1	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:198:VAL:HB	1:H:281:LYS:HD2	1.81	0.62
1:E:119:ILE:HG22	1:E:175:LEU:HD23	1.82	0.62
1:G:49:HIS:NE2	1:G:54:LYS:NZ	2.48	0.62
1:G:186:CYS:SG	1:G:204:ARG:HG2	2.41	0.61
1:G:100:LYS:O	1:G:104:MET:HG3	2.01	0.61
1:E:173:HIS:CE1	1:F:72:TYR:HB2	2.36	0.60
1:D:157:ARG:O	1:D:161:GLU:HG3	2.03	0.59
1:D:102:ARG:NH2	4:D:403:SO4:O1	2.33	0.59
1:F:131:ASP:OD1	1:F:157:ARG:NH2	2.36	0.59
1:H:115:PRO:HD3	1:H:257:LEU:HB3	1.85	0.59
1:F:78:ARG:NH2	5:F:501:HOH:O	2.30	0.58
1:F:178:GLU:CD	1:F:254:ARG:HH12	2.09	0.58
1:H:228:GLN:HB2	5:H:588:HOH:O	2.03	0.58
1:B:226:THR:HG22	1:B:228:GLN:N	2.18	0.58
1:F:232:ILE:HD12	1:F:256:PRO:HG2	1.86	0.58
2:B:401:ANP:O1B	2:B:401:ANP:O1G	2.23	0.57
1:G:204:ARG:HG3	1:G:204:ARG:HH11	1.70	0.56
1:D:1:MET:HG2	1:D:59:ASP:HB2	1.86	0.56
1:D:206:LYS:HE2	1:D:217:GLY:N	2.21	0.56
1:D:19:ARG:HG2	1:D:19:ARG:HH11	1.70	0.56
1:C:1:MET:HB3	1:C:27:HIS:ND1	2.21	0.55
1:E:119:ILE:HG12	1:F:41:ILE:HD11	1.88	0.55
1:H:206:LYS:NZ	1:H:207:GLU:HB2	2.21	0.55
1:H:240:MET:HE2	1:H:261:VAL:HB	1.88	0.55
1:D:206:LYS:H	1:D:206:LYS:CD	2.21	0.55
1:E:119:ILE:HG23	1:F:41:ILE:CD1	2.38	0.54
1:F:254:ARG:HD2	5:F:592:HOH:O	2.08	0.54
1:H:214:LEU:HD21	1:H:269:GLY:HA3	1.87	0.54
1:G:42:ASN:HA	1:G:43:PRO:C	2.32	0.54
1:D:240:MET:HE2	1:D:261:VAL:HB	1.88	0.54
1:E:38:TYR:CE2	1:F:132:MET:HE1	2.41	0.54
1:H:189:ARG:HD2	1:H:203:ARG:NH1	2.22	0.54
1:C:280:GLY:O	1:C:284:ARG:HG3	2.08	0.54
1:D:115:PRO:HD3	1:D:257:LEU:HB3	1.91	0.53
1:D:67:THR:HB	1:D:147:GLN:NE2	2.23	0.53
1:A:38:TYR:CE2	1:B:132:MET:HE1	2.43	0.53
1:C:78:ARG:NH1	1:D:105:GLN:OE1	2.36	0.53
1:H:206:LYS:CD	1:H:207:GLU:H	2.22	0.53
1:B:211:ARG:C	1:B:213:ASN:H	2.17	0.53
1:C:100:LYS:O	1:C:104:MET:HG3	2.08	0.53
1:C:115:PRO:HA	1:C:178:GLU:OE2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:LYS:NZ	1:D:216:ARG:HB2	2.24	0.53
1:H:54:LYS:HD3	1:H:54:LYS:N	2.23	0.53
1:B:201:ILE:HD13	1:B:270:ILE:HG13	1.90	0.53
1:H:101:LEU:HD22	1:H:119:ILE:HG21	1.90	0.53
1:D:211:ARG:CZ	1:D:216:ARG:HE	2.21	0.52
1:E:101:LEU:HD22	1:E:119:ILE:HG21	1.91	0.52
1:C:7:SER:HB2	1:C:64:ARG:HB2	1.91	0.52
1:D:129:LEU:O	1:D:132:MET:HB2	2.09	0.52
1:G:38:TYR:CD2	1:H:132:MET:HE1	2.44	0.52
1:B:25:ARG:NH1	1:B:284:ARG:CZ	2.73	0.52
1:C:2:LYS:HG3	1:C:28:LEU:HD22	1.91	0.52
1:E:119:ILE:HG22	1:E:175:LEU:CD2	2.40	0.52
1:H:157:ARG:O	1:H:161:GLU:HG3	2.09	0.52
1:E:119:ILE:HG23	1:F:41:ILE:HD11	1.92	0.52
1:C:109:ARG:HB2	1:D:82:MET:HE3	1.90	0.52
1:H:144:GLU:HG3	1:H:149:ILE:HD11	1.90	0.52
1:B:7:SER:HB2	1:B:64:ARG:HB2	1.92	0.52
1:F:47:SER:HB3	1:F:54:LYS:HG3	1.90	0.52
1:C:13:TYR:CE2	1:C:268:GLU:HA	2.45	0.51
1:B:13:TYR:CE2	1:B:268:GLU:HA	2.46	0.51
1:D:206:LYS:H	1:D:206:LYS:HD2	1.75	0.51
1:F:98:ARG:HG2	1:F:98:ARG:HH21	1.75	0.51
1:E:13:TYR:CE1	1:E:17:ARG:HG3	2.45	0.51
1:G:38:TYR:CE2	1:H:132:MET:HE1	2.45	0.51
1:H:70:THR:HG21	1:H:98:ARG:HD3	1.92	0.51
1:B:143:VAL:HB	1:B:173:HIS:CE1	2.44	0.51
1:A:104:MET:HE1	1:A:117:THR:HG21	1.92	0.51
1:E:157:ARG:O	1:E:161:GLU:HG3	2.11	0.51
1:H:102:ARG:NH2	4:H:403:SO4:O2	2.28	0.51
1:A:187:ASP:OD2	1:A:203:ARG:NH1	2.43	0.50
1:H:214:LEU:HD11	1:H:269:GLY:HA3	1.93	0.50
1:D:186:CYS:SG	1:D:204:ARG:HG3	2.52	0.49
1:G:173:HIS:CE1	1:H:72:TYR:HB2	2.47	0.49
1:H:67:THR:HG22	5:H:577:HOH:O	2.12	0.49
1:A:88:LEU:O	1:A:89:ASN:HB2	2.11	0.49
1:E:1:MET:HE2	1:E:292:THR:CG2	2.43	0.49
1:H:213:ASN:HB3	1:H:216:ARG:NH2	2.27	0.49
1:B:200:ALA:C	1:B:201:ILE:HD12	2.37	0.49
1:D:22:ALA:HB2	1:D:283:ILE:HD12	1.93	0.49
1:B:103:SER:OG	1:B:240:MET:HE1	2.13	0.49
1:H:49:HIS:NE2	1:H:54:LYS:HG3	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:ILE:CG2	1:B:220:ALA:HB1	2.43	0.49
1:H:189:ARG:NH2	1:H:203:ARG:HH12	2.05	0.49
1:F:178:GLU:OE1	1:F:254:ARG:NH1	2.45	0.49
1:H:167:PHE:HA	1:H:170:LEU:HD12	1.95	0.49
1:A:109:ARG:HG3	1:B:82:MET:HE3	1.95	0.49
1:B:271:GLU:O	1:B:275:GLY:N	2.46	0.49
1:C:75:ALA:HA	1:C:78:ARG:NH1	2.28	0.49
1:F:116:VAL:HG23	1:F:178:GLU:OE2	2.12	0.49
1:F:7:SER:HB2	1:F:64:ARG:HB2	1.94	0.48
1:G:78:ARG:NH2	1:H:102:ARG:HD2	2.27	0.48
1:D:109:ARG:NH2	1:D:110:GLN:OE1	2.45	0.48
1:E:187:ASP:OD1	1:E:203:ARG:HB2	2.13	0.48
1:H:206:LYS:HD3	1:H:207:GLU:H	1.78	0.48
1:F:41:ILE:HD12	1:F:41:ILE:N	2.28	0.48
1:F:240:MET:HG3	1:F:261:VAL:HG11	1.94	0.48
1:A:173:HIS:CE1	1:B:72:TYR:HB2	2.49	0.48
1:H:1:MET:HB3	1:H:27:HIS:ND1	2.28	0.48
1:D:67:THR:HB	1:D:147:GLN:HE21	1.78	0.48
1:E:147:GLN:HG2	1:E:215:HIS:CD2	2.48	0.48
1:F:191:LEU:HD23	1:F:198:VAL:HG22	1.94	0.48
1:H:69:ILE:HD12	1:H:69:ILE:O	2.14	0.48
1:D:9:ASP:HB3	1:D:12:LEU:HG	1.96	0.48
1:G:8:ARG:HD2	1:G:69:ILE:HD13	1.95	0.48
1:H:9:ASP:HB3	1:H:12:LEU:HG	1.95	0.48
1:C:4:ALA:HB3	1:C:61:VAL:HG22	1.95	0.47
1:C:203:ARG:HH11	1:C:213:ASN:HA	1.79	0.47
1:E:121:HIS:HD2	1:E:122:SER:OG	1.97	0.47
1:B:114:LEU:CD2	1:B:258:VAL:O	2.63	0.47
1:F:50:TYR:O	1:F:53:ARG:HG2	2.15	0.47
1:F:98:ARG:HH12	1:F:263:ALA:H	1.63	0.47
1:F:216:ARG:O	1:F:216:ARG:HG2	2.15	0.47
2:H:401:ANP:O2G	2:H:401:ANP:O1A	2.33	0.47
1:F:46:SER:HB2	1:F:83:LEU:HD11	1.97	0.47
1:D:181:LYS:HD3	1:D:182:GLU:OE1	2.15	0.47
1:H:206:LYS:HD3	1:H:207:GLU:N	2.30	0.47
1:H:268:GLU:O	1:H:272:LYS:HG3	2.15	0.47
1:A:50:TYR:O	1:A:53:ARG:HG2	2.15	0.46
1:G:204:ARG:HG3	1:G:204:ARG:NH1	2.29	0.46
1:H:49:HIS:CE1	1:H:54:LYS:HG3	2.50	0.46
1:H:104:MET:HE2	1:H:104:MET:HB3	1.80	0.46
1:H:104:MET:HE3	1:H:114:LEU:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:206:LYS:CG	1:H:207:GLU:H	2.28	0.46
1:H:213:ASN:HD22	1:H:216:ARG:NH1	2.14	0.46
1:E:187:ASP:OD2	1:E:203:ARG:NH1	2.49	0.46
1:F:285:TRP:CE2	1:F:289:HIS:CE1	3.04	0.46
1:D:25:ARG:NH1	1:D:284:ARG:NH1	2.64	0.46
1:B:10:GLY:HA3	1:B:31:ILE:HG21	1.98	0.46
1:D:133:VAL:HG21	1:D:176:VAL:HG12	1.96	0.46
1:E:142:LEU:HD12	1:E:142:LEU:N	2.31	0.46
1:G:131:ASP:CG	1:G:157:ARG:HH22	2.24	0.46
1:H:42:ASN:HA	1:H:43:PRO:C	2.39	0.46
1:B:42:ASN:HA	1:B:43:PRO:C	2.41	0.46
1:F:102:ARG:NE	4:F:403:SO4:O2	2.49	0.46
1:D:98:ARG:NH1	1:D:146:THR:O	2.47	0.46
1:B:132:MET:HE3	1:B:132:MET:HB2	1.75	0.45
1:D:30:GLU:OE2	1:D:53:ARG:NH2	2.32	0.45
1:F:13:TYR:CE2	1:F:17:ARG:HG3	2.51	0.45
1:B:9:ASP:C	1:B:11:THR:H	2.24	0.45
1:D:206:LYS:HZ3	1:D:216:ARG:HB2	1.81	0.45
1:C:88:LEU:O	1:C:89:ASN:HB2	2.16	0.45
1:C:201:ILE:HG23	1:C:270:ILE:HG13	1.98	0.45
1:F:13:TYR:CZ	1:F:17:ARG:HG3	2.52	0.45
1:H:1:MET:HG3	1:H:59:ASP:HB2	1.97	0.45
1:D:13:TYR:CE1	1:D:268:GLU:HA	2.51	0.45
1:D:181:LYS:CD	1:D:181:LYS:H	2.30	0.45
1:F:115:PRO:HD3	1:F:257:LEU:HB3	1.98	0.45
2:H:401:ANP:O2G	2:H:401:ANP:O1B	2.35	0.45
1:B:9:ASP:HB3	1:B:12:LEU:HG	1.97	0.45
1:B:25:ARG:NH2	1:B:287:GLU:OE1	2.43	0.45
1:B:25:ARG:NH1	1:B:284:ARG:NH2	2.65	0.45
1:D:13:TYR:CE1	1:D:17:ARG:HG3	2.50	0.45
1:G:40:ASN:HA	1:H:118:GLY:HA2	1.98	0.45
1:A:13:TYR:CE1	1:A:17:ARG:HG3	2.52	0.45
1:E:22:ALA:HB3	1:E:29:VAL:HG21	1.99	0.45
1:E:92:VAL:HG12	5:E:646:HOH:O	2.16	0.45
1:B:188:ILE:HA	1:B:201:ILE:O	2.16	0.44
1:B:25:ARG:HH12	1:B:284:ARG:NH2	2.15	0.44
1:B:205:ALA:HB2	1:B:212:SER:HA	1.99	0.44
1:D:185:GLY:O	1:D:212:SER:HB3	2.16	0.44
1:D:206:LYS:HE3	1:D:217:GLY:C	2.42	0.44
1:E:229:GLU:HA	1:E:232:ILE:HG22	1.99	0.44
1:B:201:ILE:CD1	1:B:270:ILE:HG13	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:228:GLN:HB2	5:E:639:HOH:O	2.18	0.44
1:B:88:LEU:O	1:B:89:ASN:HB2	2.17	0.44
1:D:194:GLY:HA3	1:D:285:TRP:CH2	2.53	0.44
1:E:40:ASN:HB3	1:E:45:ALA:O	2.16	0.44
1:G:7:SER:HB2	1:G:64:ARG:HB2	1.99	0.44
1:D:142:LEU:HD22	1:D:167:PHE:CG	2.52	0.44
1:E:110:GLN:HG3	1:E:239:THR:HG23	2.00	0.44
1:F:88:LEU:O	1:F:89:ASN:HB2	2.17	0.44
1:H:36:SER:HB2	1:H:50:TYR:CE1	2.52	0.44
1:H:178:GLU:O	1:H:178:GLU:HG3	2.16	0.44
1:A:2:LYS:HE3	1:A:2:LYS:HB2	1.43	0.44
1:B:188:ILE:O	1:B:248:ASP:HA	2.18	0.44
1:F:1:MET:HB3	1:F:27:HIS:ND1	2.33	0.44
1:D:268:GLU:O	1:D:272:LYS:HG3	2.18	0.44
1:E:82:MET:HE3	1:E:82:MET:HB3	1.65	0.44
1:E:142:LEU:HD13	1:E:150:GLY:HA3	2.00	0.43
1:F:138:LEU:HD13	1:F:176:VAL:CG1	2.49	0.43
1:G:88:LEU:O	1:G:89:ASN:HB2	2.17	0.43
1:B:194:GLY:HA3	1:B:285:TRP:CH2	2.53	0.43
1:G:102:ARG:NE	4:G:404:SO4:O3	2.41	0.43
1:H:70:THR:HG21	1:H:98:ARG:CD	2.48	0.43
1:G:47:SER:HB3	1:G:54:LYS:HG3	2.00	0.43
1:H:228:GLN:O	1:H:232:ILE:HD13	2.17	0.43
1:A:2:LYS:NZ	5:A:501:HOH:O	2.30	0.43
1:F:6:LEU:HD22	1:F:34:PRO:HG3	2.01	0.43
1:H:206:LYS:HZ2	1:H:207:GLU:HB2	1.82	0.43
1:B:114:LEU:HD22	1:B:258:VAL:O	2.19	0.43
1:E:122:SER:HB3	1:E:168:ARG:CZ	2.48	0.43
1:B:226:THR:CG2	1:B:228:GLN:HB3	2.49	0.43
1:E:104:MET:HE1	1:E:175:LEU:HD21	2.01	0.43
1:H:1:MET:CG	1:H:59:ASP:HB2	2.49	0.43
1:C:240:MET:HG3	1:C:261:VAL:HG11	2.00	0.43
1:G:13:TYR:CE1	1:G:17:ARG:HG3	2.53	0.43
1:H:17:ARG:NH2	1:H:277:ASP:OD1	2.44	0.43
2:A:401:ANP:O5'	2:A:401:ANP:H8	2.19	0.42
1:B:254:ARG:NH1	1:B:257:LEU:HD21	2.34	0.42
1:A:141:LYS:HA	1:A:150:GLY:O	2.20	0.42
1:G:121:HIS:HD2	1:G:122:SER:OG	2.02	0.42
1:B:22:ALA:HB3	1:B:29:VAL:HG21	2.01	0.42
1:B:229:GLU:OE1	1:B:251:ARG:NH2	2.48	0.42
1:C:121:HIS:HD2	1:C:122:SER:OG	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:ARG:HD2	1:B:105:GLN:OE1	2.19	0.42
1:B:113:ASP:OD2	1:B:257:LEU:CD2	2.66	0.42
1:H:186:CYS:HA	1:H:203:ARG:O	2.19	0.42
1:B:114:LEU:CD2	1:B:258:VAL:HB	2.49	0.42
1:B:104:MET:HE2	1:B:104:MET:HB3	1.87	0.42
1:C:268:GLU:OE2	1:C:272:LYS:NZ	2.48	0.42
1:A:12:LEU:HB2	5:A:637:HOH:O	2.19	0.42
1:A:41:ILE:HG13	1:A:42:ASN:H	1.85	0.42
1:D:88:LEU:O	1:D:89:ASN:HB2	2.20	0.42
1:F:109:ARG:NH2	5:F:502:HOH:O	2.38	0.42
1:F:232:ILE:HD13	5:F:542:HOH:O	2.19	0.42
1:H:206:LYS:HZ2	1:H:207:GLU:H	1.68	0.42
1:A:77:LEU:HD23	1:A:91:SER:HA	2.02	0.41
1:E:250:LEU:HG	1:E:259:MET:HG3	2.01	0.41
1:F:133:VAL:HG21	1:F:176:VAL:HG12	2.02	0.41
1:B:203:ARG:NH2	1:B:213:ASN:HA	2.35	0.41
1:C:13:TYR:CZ	1:C:17:ARG:HG3	2.55	0.41
1:G:13:TYR:CE1	1:G:268:GLU:HA	2.55	0.41
1:F:187:ASP:OD1	1:F:212:SER:OG	2.33	0.41
1:B:18:LEU:HD21	1:B:282:MET:HE2	2.03	0.41
1:B:191:LEU:HD13	1:B:278:ILE:HG21	2.02	0.41
1:F:143:VAL:HB	1:F:173:HIS:CE1	2.55	0.41
1:B:9:ASP:C	1:B:11:THR:N	2.78	0.41
1:F:209:ASP:OD1	1:F:210:PHE:N	2.54	0.41
1:G:109:ARG:NH2	5:G:512:HOH:O	2.52	0.41
1:G:209:ASP:OD1	1:G:210:PHE:N	2.53	0.41
1:C:102:ARG:NH2	4:C:404:SO4:O2	2.34	0.41
1:F:173:HIS:C	1:F:174:ILE:HG13	2.46	0.41
1:B:211:ARG:C	1:B:213:ASN:N	2.77	0.41
1:E:132:MET:HE3	1:F:40:ASN:OD1	2.21	0.41
1:F:206:LYS:HB2	1:F:206:LYS:HE3	1.94	0.41
1:H:179:TYR:CE2	1:H:181:LYS:HG3	2.56	0.41
1:B:182:GLU:OE1	1:B:182:GLU:N	2.48	0.41
1:B:201:ILE:HG21	1:B:220:ALA:HB1	2.03	0.40
1:C:198:VAL:HG13	1:C:281:LYS:HD3	2.02	0.40
1:F:138:LEU:HD23	1:F:138:LEU:HA	1.66	0.40
1:H:88:LEU:O	1:H:89:ASN:HB2	2.22	0.40
1:B:114:LEU:HD23	1:B:258:VAL:HB	2.04	0.40
1:B:149:ILE:O	1:C:169:GLY:HA3	2.21	0.40
1:C:104:MET:HE3	1:D:41:ILE:CD1	2.44	0.40
1:E:88:LEU:O	1:E:89:ASN:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:110:GLN:HG3	1:D:239:THR:HG23	2.03	0.40
1:E:104:MET:HE1	1:E:175:LEU:CD2	2.51	0.40
1:G:115:PRO:HD3	1:G:257:LEU:HB3	2.02	0.40
1:A:38:TYR:HE2	1:B:132:MET:HE1	1.83	0.40
1:A:82:MET:HE1	1:B:105:GLN:HG2	2.02	0.40
1:C:8:ARG:HD2	1:C:69:ILE:HD13	2.02	0.40
1:E:1:MET:HB3	1:E:27:HIS:ND1	2.36	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	290/306 (95%)	283 (98%)	6 (2%)	1 (0%)	36	30
1	B	290/306 (95%)	275 (95%)	14 (5%)	1 (0%)	36	30
1	C	288/306 (94%)	279 (97%)	8 (3%)	1 (0%)	36	30
1	D	290/306 (95%)	278 (96%)	11 (4%)	1 (0%)	36	30
1	E	290/306 (95%)	283 (98%)	6 (2%)	1 (0%)	36	30
1	F	290/306 (95%)	280 (97%)	9 (3%)	1 (0%)	36	30
1	G	290/306 (95%)	284 (98%)	5 (2%)	1 (0%)	36	30
1	H	290/306 (95%)	277 (96%)	12 (4%)	1 (0%)	36	30
All	All	2318/2448 (95%)	2239 (97%)	71 (3%)	8 (0%)	36	30

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	ASN
1	C	89	ASN
1	B	89	ASN

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Mol	Chain	Res	Type
1	E	89	ASN
1	F	89	ASN
1	G	89	ASN
1	H	89	ASN
1	D	89	ASN

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	231/242 (96%)	230 (100%)	1 (0%)	84	88
1	B	231/242 (96%)	230 (100%)	1 (0%)	84	88
1	C	229/242 (95%)	228 (100%)	1 (0%)	84	88
1	D	231/242 (96%)	230 (100%)	1 (0%)	84	88
1	E	231/242 (96%)	231 (100%)	0	100	100
1	F	231/242 (96%)	231 (100%)	0	100	100
1	G	231/242 (96%)	231 (100%)	0	100	100
1	H	231/242 (96%)	230 (100%)	1 (0%)	84	88
All	All	1846/1936 (95%)	1841 (100%)	5 (0%)	86	90

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	B	11	THR
1	C	53	ARG
1	D	2	LYS
1	H	83	LEU

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

Mol	Chain	Res	Type
1	B	228	GLN
1	C	42	ASN
1	D	147	GLN
1	E	110	GLN
1	E	215	HIS
1	E	289	HIS
1	G	289	HIS
1	H	42	ASN
1	H	184	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](#)

Of 30 ligands modelled in this entry, 8 are monoatomic - leaving 22 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	$\# Z > 2$
2	ANP	A	401	3	33,33,33	2.35	7 (21%)	45,52,52	1.12	3 (6%)
2	ANP	H	401	3	33,33,33	2.14	5 (15%)	45,52,52	0.67	2 (4%)
4	SO4	C	404	-	4,4,4	0.25	0	6,6,6	0.14	0
4	SO4	A	403	-	4,4,4	0.27	0	6,6,6	0.52	0
4	SO4	H	403	-	4,4,4	0.30	0	6,6,6	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	G	404	-	4,4,4	0.22	0	6,6,6	0.23	0
2	ANP	C	401	3	33,33,33	2.13	6 (18%)	45,52,52	0.89	4 (8%)
2	ANP	F	401	3	33,33,33	1.03	4 (12%)	45,52,52	0.83	3 (6%)
2	ANP	B	401	3	33,33,33	2.23	5 (15%)	45,52,52	0.86	3 (6%)
4	SO4	A	405	-	4,4,4	0.32	0	6,6,6	0.21	0
2	ANP	D	401	3	33,33,33	2.14	6 (18%)	45,52,52	0.77	2 (4%)
4	SO4	B	403	-	4,4,4	0.25	0	6,6,6	0.20	0
4	SO4	E	404	-	4,4,4	0.25	0	6,6,6	0.32	0
2	ANP	G	401	3	33,33,33	1.63	4 (12%)	45,52,52	0.70	2 (4%)
4	SO4	D	403	-	4,4,4	0.26	0	6,6,6	0.19	0
4	SO4	A	404	-	4,4,4	0.35	0	6,6,6	0.21	0
4	SO4	F	403	-	4,4,4	0.19	0	6,6,6	0.21	0
2	ANP	E	402	3	33,33,33	2.79	7 (21%)	45,52,52	0.93	3 (6%)
4	SO4	C	405	-	4,4,4	0.23	0	6,6,6	0.49	0
4	SO4	G	403	-	4,4,4	0.30	0	6,6,6	0.35	0
4	SO4	E	401	-	4,4,4	0.28	0	6,6,6	0.39	0
4	SO4	C	403	-	4,4,4	0.34	0	6,6,6	0.35	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
2	ANP	A	401	3	-	3/18/38/38	0/3/3/3
2	ANP	B	401	3	-	3/18/38/38	0/3/3/3
2	ANP	H	401	3	-	6/18/38/38	0/3/3/3
2	ANP	C	401	3	-	5/18/38/38	0/3/3/3
2	ANP	D	401	3	-	1/18/38/38	0/3/3/3
2	ANP	G	401	3	-	4/18/38/38	0/3/3/3
2	ANP	F	401	3	-	3/18/38/38	0/3/3/3
2	ANP	E	402	3	-	5/18/38/38	0/3/3/3

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	402	ANP	PB-O1B	11.58	1.63	1.46
2	C	401	ANP	PG-O1G	8.86	1.59	1.46
2	B	401	ANP	PG-O1G	8.74	1.59	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	ANP	PG-O1G	8.47	1.59	1.46
2	H	401	ANP	PG-O1G	8.47	1.59	1.46
2	E	402	ANP	PG-O1G	7.98	1.58	1.46
2	D	401	ANP	PG-O1G	7.86	1.58	1.46
2	D	401	ANP	PB-O1B	7.43	1.57	1.46
2	G	401	ANP	PG-O1G	7.37	1.57	1.46
2	B	401	ANP	PB-O1B	7.36	1.57	1.46
2	H	401	ANP	PB-O1B	7.09	1.56	1.46
2	C	401	ANP	PB-O1B	6.35	1.55	1.46
2	A	401	ANP	PA-O3A	5.88	1.65	1.59
2	A	401	ANP	PG-N3B	5.03	1.76	1.63
2	A	401	ANP	PB-O2B	4.70	1.69	1.56
2	E	402	ANP	PA-O3A	4.64	1.64	1.59
2	E	402	ANP	PG-N3B	3.99	1.73	1.63
2	A	401	ANP	PB-N3B	3.52	1.72	1.63
2	D	401	ANP	PA-O3A	3.01	1.62	1.59
2	B	401	ANP	PG-N3B	2.96	1.71	1.63
2	C	401	ANP	PB-O2B	-2.82	1.49	1.56
2	G	401	ANP	PA-O3A	2.80	1.62	1.59
2	G	401	ANP	PG-N3B	2.80	1.70	1.63
2	F	401	ANP	PG-N3B	2.78	1.70	1.63
2	F	401	ANP	PB-O1B	2.74	1.50	1.46
2	H	401	ANP	PG-O3G	-2.72	1.49	1.56
2	G	401	ANP	PG-O3G	-2.71	1.49	1.56
2	E	402	ANP	PB-N3B	2.69	1.70	1.63
2	F	401	ANP	PG-O1G	2.69	1.50	1.46
2	E	402	ANP	PB-O2B	-2.63	1.49	1.56
2	H	401	ANP	PG-N3B	2.61	1.70	1.63
2	C	401	ANP	PG-N3B	2.60	1.70	1.63
2	D	401	ANP	PB-O2B	-2.59	1.49	1.56
2	A	401	ANP	PG-O3G	-2.59	1.49	1.56
2	H	401	ANP	PB-O2B	-2.50	1.50	1.56
2	B	401	ANP	PB-O2B	-2.46	1.50	1.56
2	D	401	ANP	PG-O2G	-2.43	1.50	1.56
2	E	402	ANP	PG-O2G	-2.43	1.50	1.56
2	D	401	ANP	PG-N3B	2.38	1.69	1.63
2	A	401	ANP	PB-O1B	2.31	1.49	1.46
2	B	401	ANP	PG-O2G	2.28	1.62	1.56
2	C	401	ANP	PG-O2G	-2.10	1.51	1.56
2	C	401	ANP	PB-N3B	2.03	1.68	1.63
2	F	401	ANP	PB-O3A	-2.02	1.56	1.59

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ANP	O1B-PB-N3B	-5.05	104.34	111.77
2	E	402	ANP	O2G-PG-O1G	-3.45	104.79	113.45
2	E	402	ANP	O1B-PB-N3B	3.13	116.38	111.77
2	A	401	ANP	O3G-PG-O1G	-3.01	105.90	113.45
2	C	401	ANP	O1B-PB-N3B	-2.90	107.50	111.77
2	A	401	ANP	O2B-PB-O3A	2.88	114.24	104.64
2	D	401	ANP	O2B-PB-O3A	2.74	113.80	104.64
2	B	401	ANP	O2B-PB-O3A	2.69	113.61	104.64
2	E	402	ANP	O3A-PA-O1A	2.58	118.47	110.70
2	C	401	ANP	O2G-PG-O1G	-2.50	107.18	113.45
2	F	401	ANP	O2A-PA-O3A	2.39	113.74	107.27
2	C	401	ANP	O2B-PB-O3A	2.39	112.61	104.64
2	H	401	ANP	O3G-PG-O1G	-2.27	107.77	113.45
2	G	401	ANP	O3G-PG-O1G	-2.25	107.80	113.45
2	B	401	ANP	O3G-PG-O1G	-2.23	107.86	113.45
2	H	401	ANP	O2B-PB-O3A	2.16	111.87	104.64
2	C	401	ANP	O2B-PB-O1B	-2.15	105.26	109.87
2	G	401	ANP	O2A-PA-O3A	2.14	113.05	107.27
2	D	401	ANP	O2G-PG-O1G	-2.13	108.11	113.45
2	F	401	ANP	O2G-PG-O1G	-2.06	108.29	113.45
2	B	401	ANP	O2B-PB-O1B	-2.05	105.46	109.87
2	F	401	ANP	O2B-PB-O3A	2.00	111.33	104.64

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ANP	PA-O3A-PB-O2B
2	C	401	ANP	PB-N3B-PG-O1G
2	C	401	ANP	PG-N3B-PB-O3A
2	C	401	ANP	PA-O3A-PB-O2B
2	D	401	ANP	PA-O3A-PB-O2B
2	E	402	ANP	PB-N3B-PG-O1G
2	E	402	ANP	PG-N3B-PB-O1B
2	E	402	ANP	PG-N3B-PB-O3A
2	E	402	ANP	PA-O3A-PB-O2B
2	F	401	ANP	PG-N3B-PB-O1B
2	G	401	ANP	PA-O3A-PB-O2B
2	H	401	ANP	PB-N3B-PG-O1G
2	H	401	ANP	C5'-O5'-PA-O1A
2	H	401	ANP	C5'-O5'-PA-O3A
2	G	401	ANP	PB-O3A-PA-O1A
2	H	401	ANP	C5'-O5'-PA-O2A

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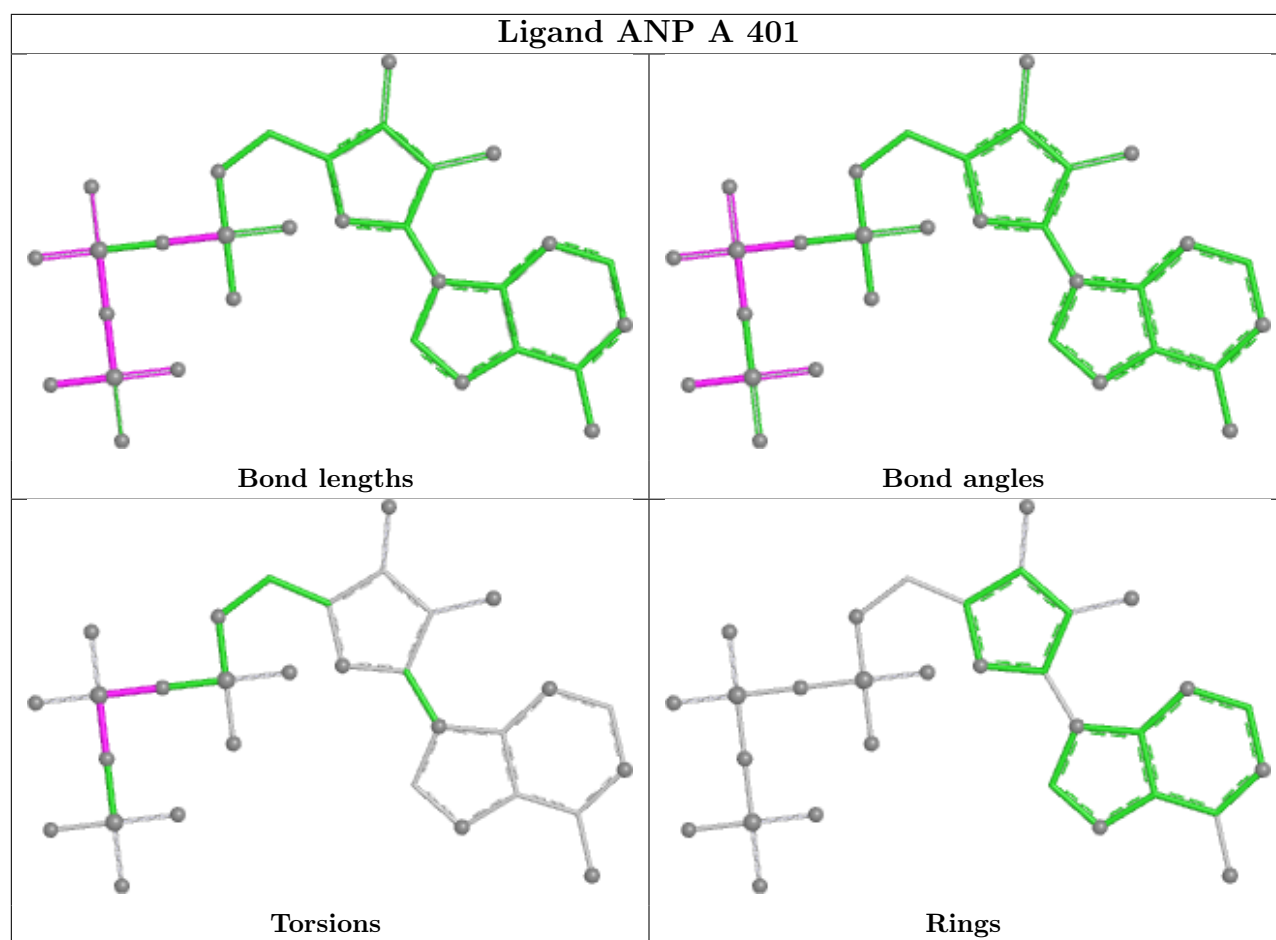
Mol	Chain	Res	Type	Atoms
2	F	401	ANP	PB-O3A-PA-O2A
2	B	401	ANP	PB-O3A-PA-O2A
2	C	401	ANP	PG-N3B-PB-O1B
2	B	401	ANP	PA-O3A-PB-O2B
2	H	401	ANP	C4'-C5'-O5'-PA
2	G	401	ANP	PB-O3A-PA-O2A
2	A	401	ANP	PA-O3A-PB-O1B
2	C	401	ANP	PA-O3A-PB-O1B
2	E	402	ANP	PA-O3A-PB-O1B
2	F	401	ANP	PA-O3A-PB-O1B
2	G	401	ANP	PA-O3A-PB-O1B
2	H	401	ANP	PA-O3A-PB-O1B
2	A	401	ANP	PG-N3B-PB-O3A
2	B	401	ANP	PB-O3A-PA-O1A

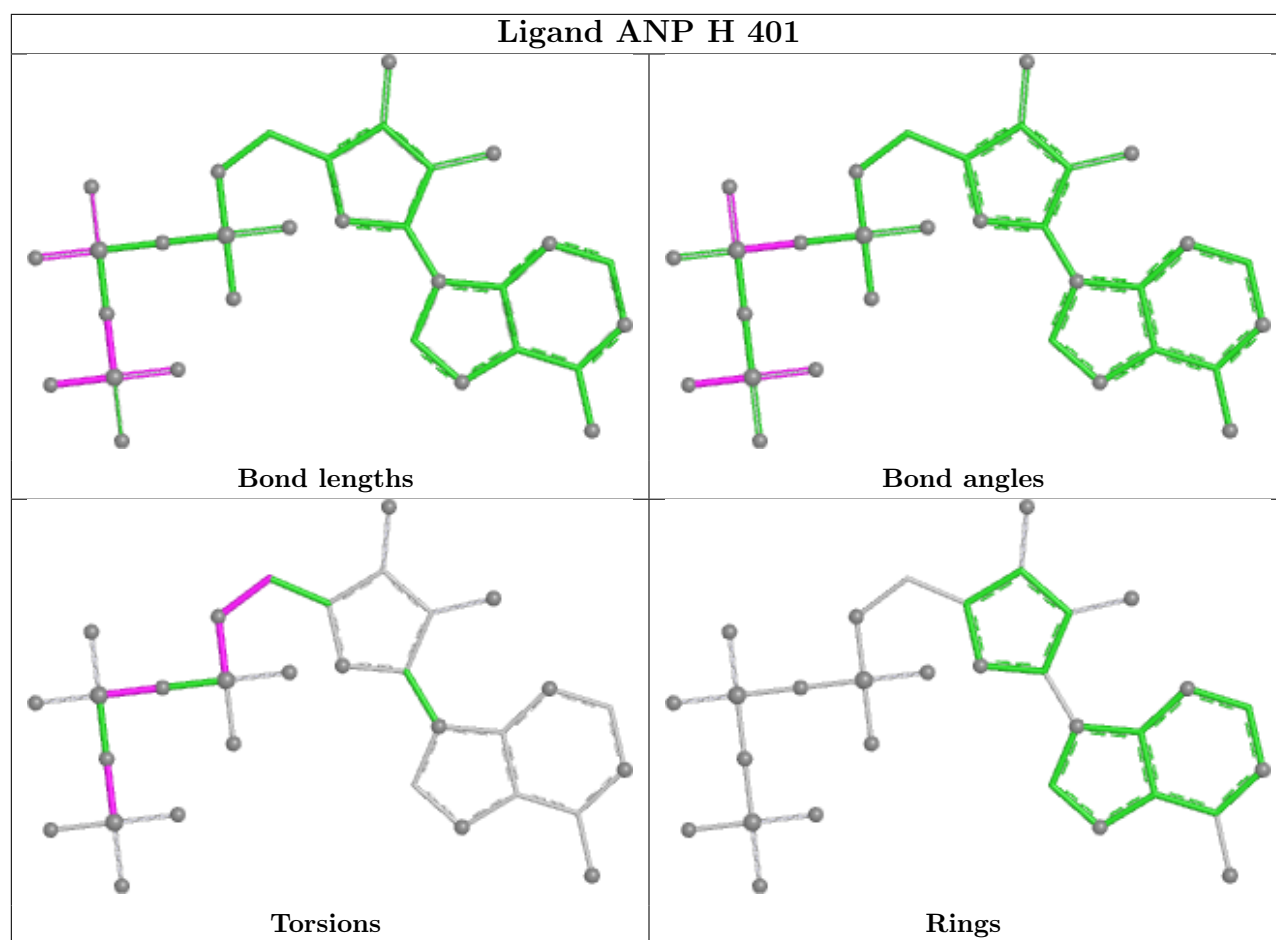
There are no ring outliers.

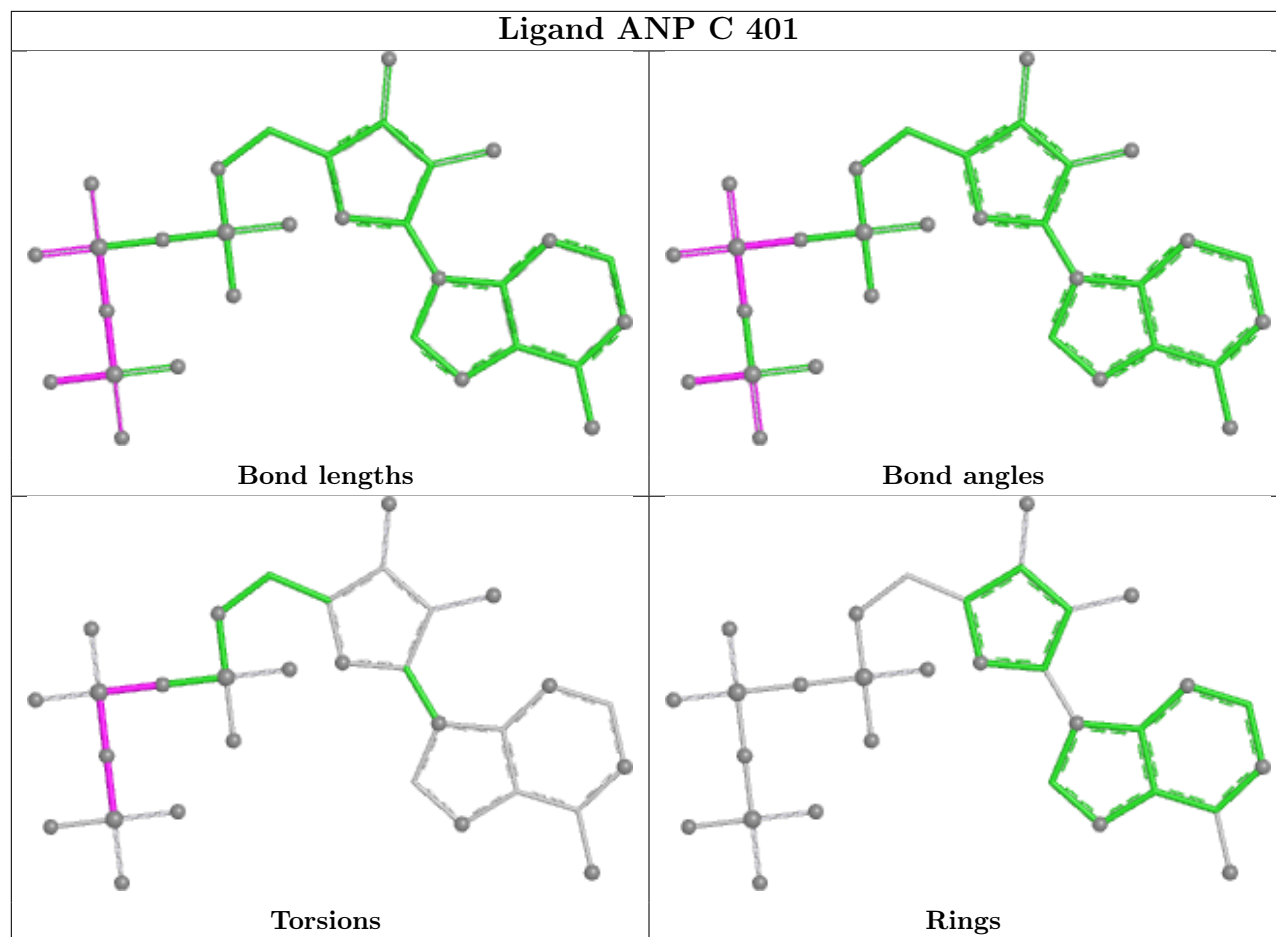
10 monomers are involved in 12 short contacts:

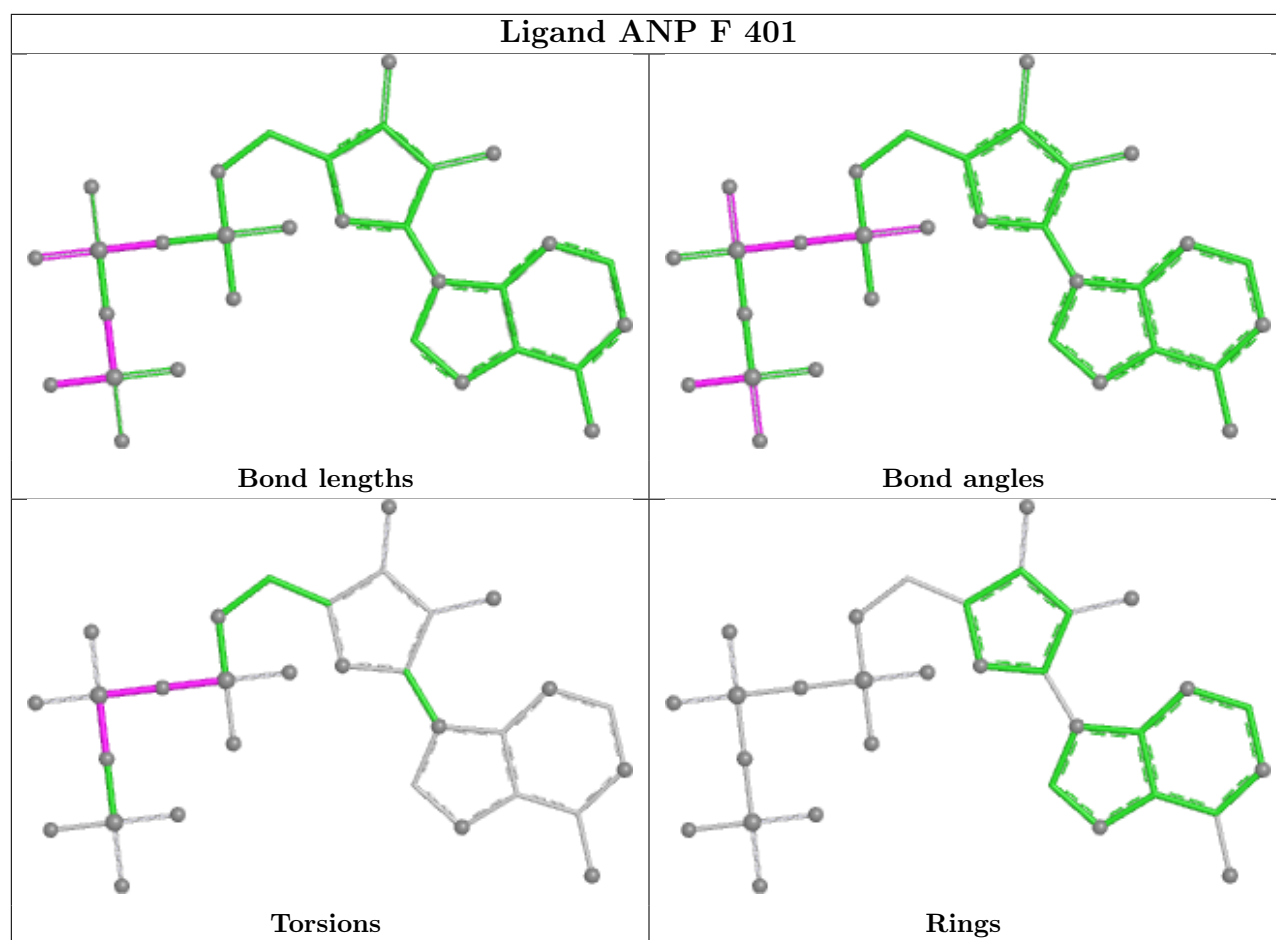
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ANP	1	0
2	H	401	ANP	2	0
4	C	404	SO4	1	0
4	H	403	SO4	1	0
4	G	404	SO4	1	0
2	B	401	ANP	1	0
4	A	405	SO4	1	0
4	D	403	SO4	1	0
4	F	403	SO4	2	0
4	C	403	SO4	1	0

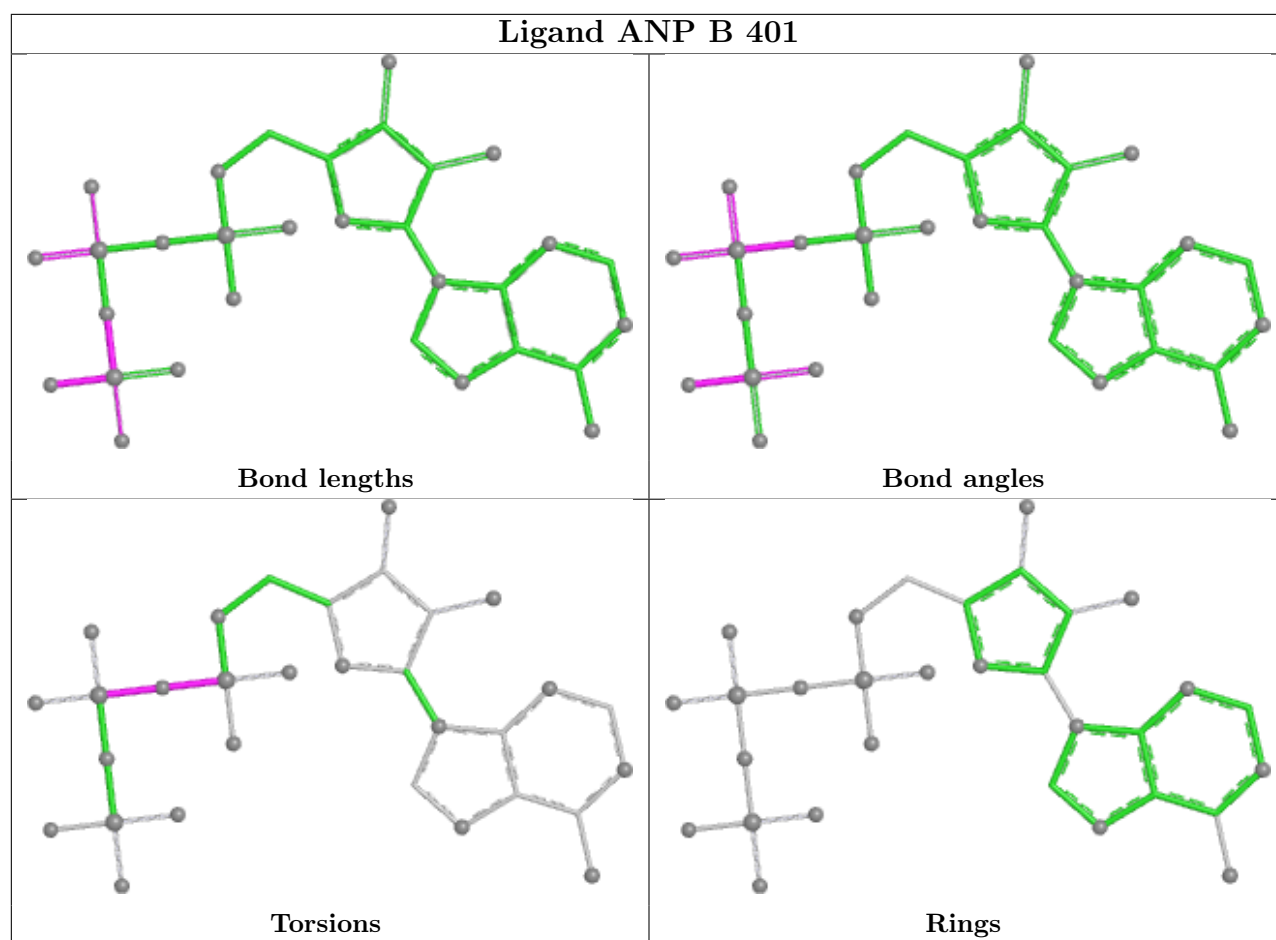
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

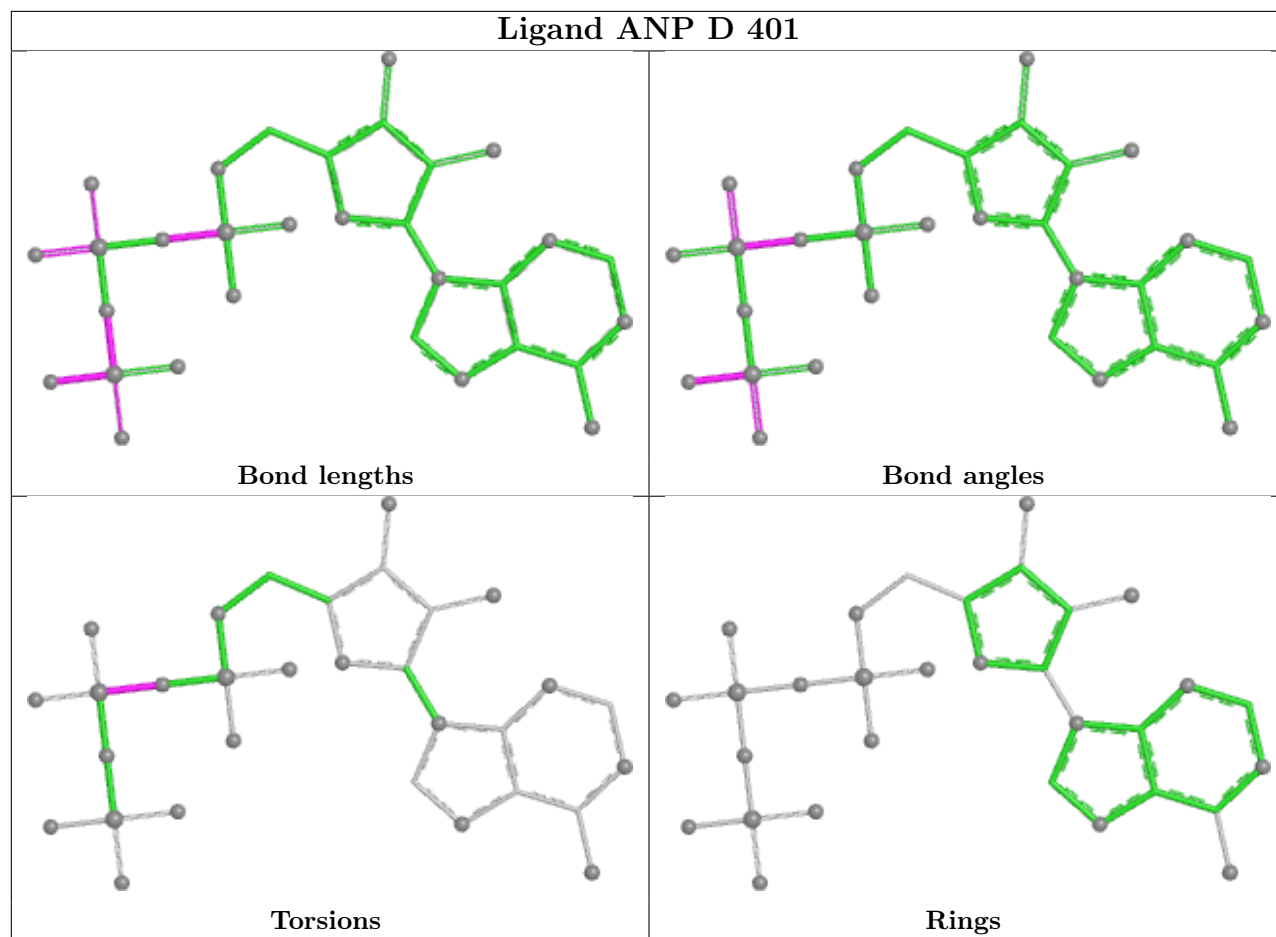


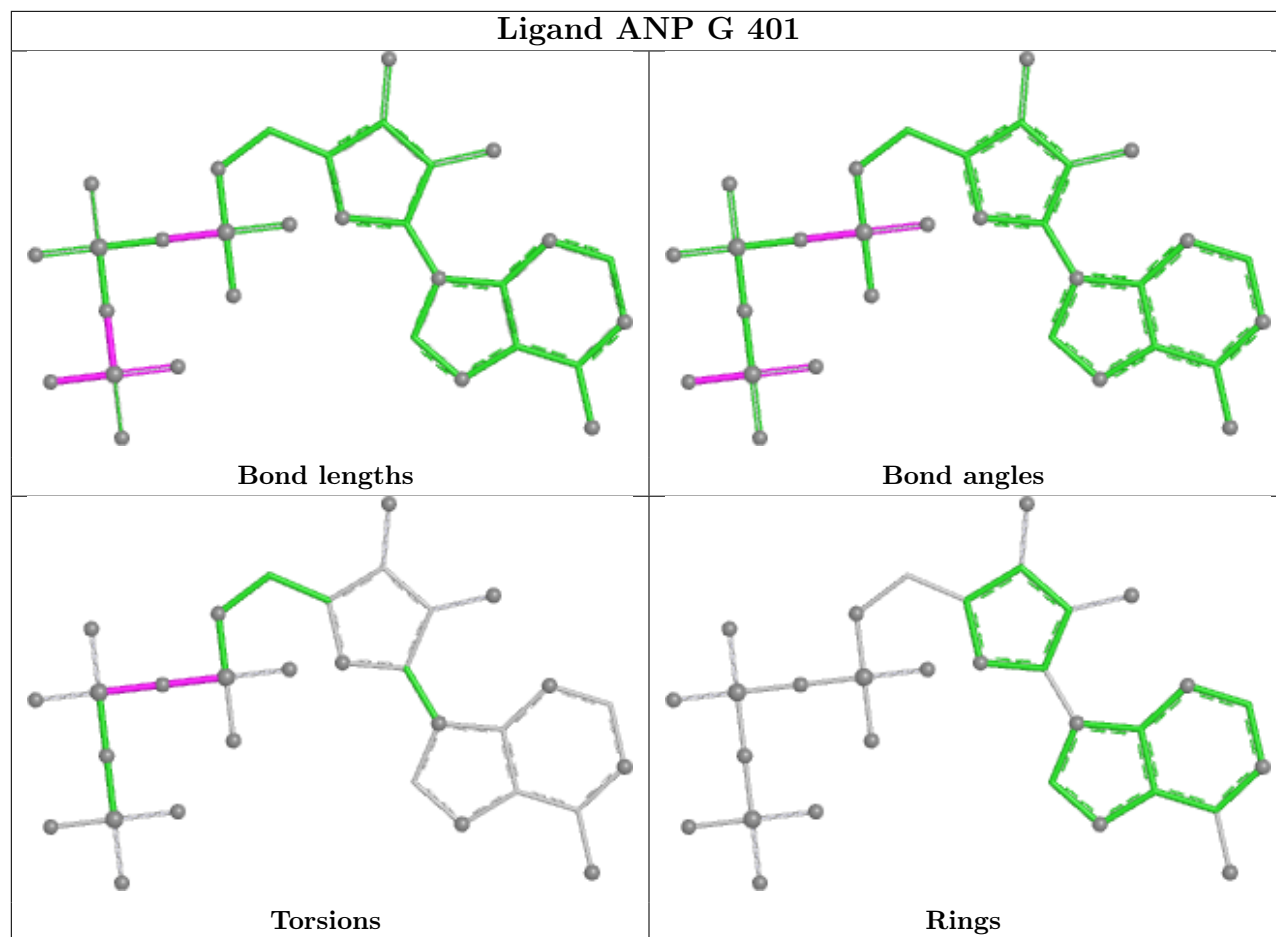


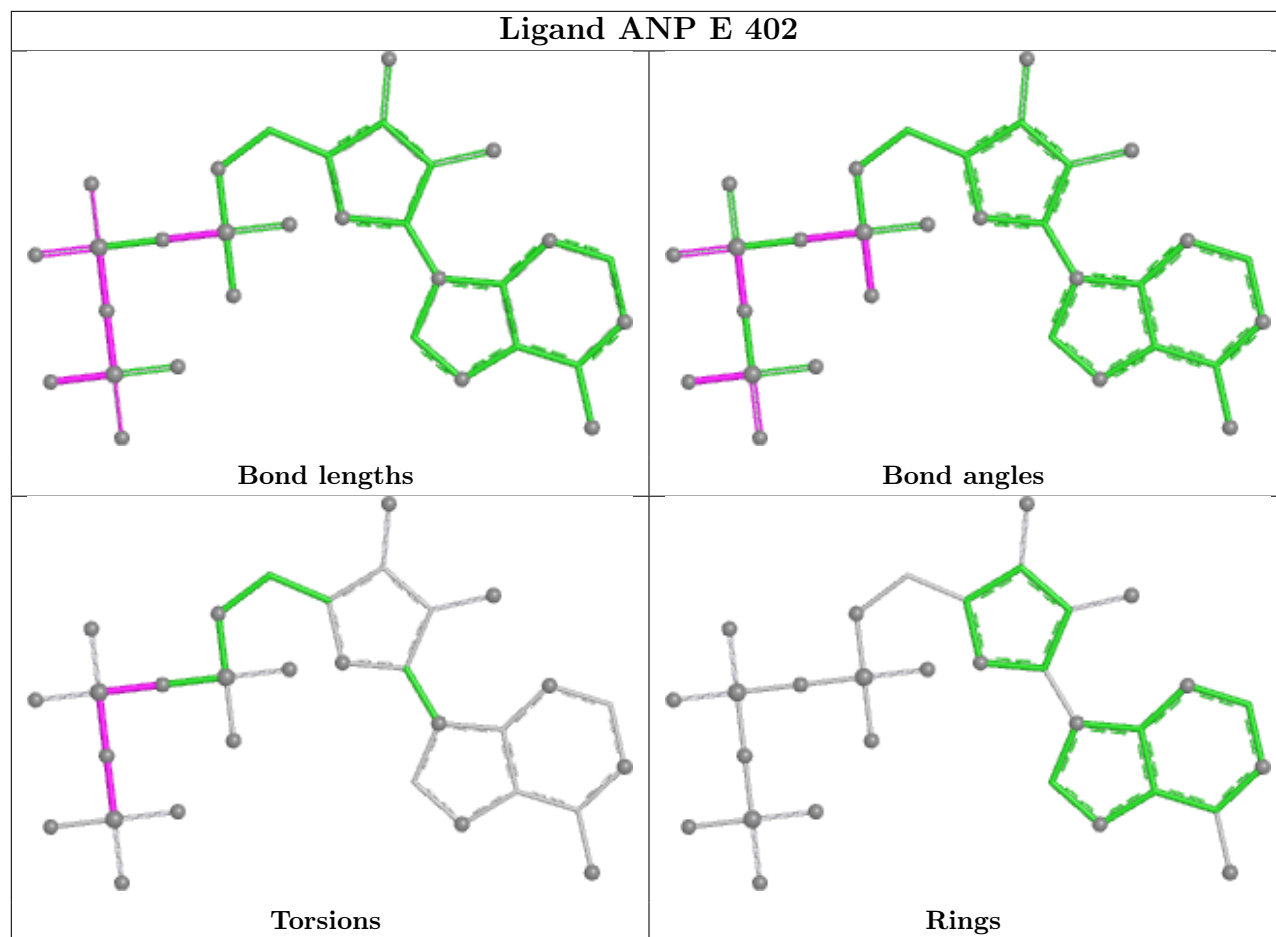












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	292/306 (95%)	-0.28	2 (0%) 84 85	22, 33, 48, 61	0
1	B	292/306 (95%)	0.74	37 (12%) 8 7	27, 52, 81, 103	0
1	C	290/306 (94%)	0.01	5 (1%) 69 70	28, 39, 52, 69	0
1	D	292/306 (95%)	0.20	8 (2%) 56 58	31, 44, 68, 97	0
1	E	292/306 (95%)	-0.11	7 (2%) 59 61	25, 35, 54, 64	0
1	F	292/306 (95%)	0.22	10 (3%) 48 48	27, 42, 74, 102	0
1	G	292/306 (95%)	-0.37	3 (1%) 79 81	21, 32, 46, 73	0
1	H	292/306 (95%)	0.33	15 (5%) 33 33	28, 43, 73, 117	0
All	All	2334/2448 (95%)	0.09	87 (3%) 45 45	21, 39, 65, 117	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	45	ALA	5.0
1	F	45	ALA	4.8
1	F	43	PRO	4.2
1	F	44	ALA	4.0
1	B	214	LEU	3.9
1	C	45	ALA	3.8
1	B	201	ILE	3.7
1	C	42	ASN	3.4
1	B	269	GLY	3.3
1	B	273	THR	3.3
1	F	68	ALA	3.3
1	H	214	LEU	3.2
1	B	202	GLU	3.2
1	B	292	THR	3.2
1	B	200	ALA	3.1
1	B	222	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	H	136	ALA	3.1
1	H	217	GLY	3.0
1	F	42	ASN	3.0
1	B	220	ALA	3.0
1	D	44	ALA	3.0
1	D	45	ALA	3.0
1	H	208	GLY	2.9
1	B	219	ALA	2.9
1	B	66	GLY	2.9
1	B	215	HIS	2.9
1	H	215	HIS	2.8
1	F	41	ILE	2.8
1	C	43	PRO	2.8
1	B	232	ILE	2.8
1	B	199	ALA	2.8
1	H	219	ALA	2.8
1	H	211	ARG	2.7
1	H	210	PHE	2.7
1	B	43	PRO	2.7
1	E	169	GLY	2.7
1	G	214	LEU	2.6
1	B	221	SER	2.6
1	B	205	ALA	2.6
1	B	188	ILE	2.6
1	H	291	THR	2.6
1	B	185	GLY	2.6
1	E	45	ALA	2.6
1	B	133	VAL	2.5
1	B	225	ILE	2.5
1	D	291	THR	2.5
1	D	292	THR	2.5
1	H	220	ALA	2.5
1	A	214	LEU	2.5
1	F	208	GLY	2.5
1	H	154	ALA	2.5
1	E	214	LEU	2.5
1	D	42	ASN	2.5
1	H	206	LYS	2.5
1	B	217	GLY	2.4
1	G	292	THR	2.4
1	B	44	ALA	2.4
1	F	292	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	215	HIS	2.4
1	B	276	ILE	2.4
1	E	292	THR	2.3
1	B	249	ILE	2.3
1	F	291	THR	2.3
1	C	44	ALA	2.3
1	B	208	GLY	2.3
1	H	167	PHE	2.3
1	E	123	PRO	2.2
1	C	214	LEU	2.2
1	D	43	PRO	2.2
1	H	137	PRO	2.2
1	E	44	ALA	2.2
1	E	142	LEU	2.2
1	B	11	THR	2.2
1	B	227	PRO	2.1
1	B	274	THR	2.2
1	B	216	ARG	2.1
1	B	186	CYS	2.1
1	D	66	GLY	2.1
1	F	218	GLY	2.1
1	B	13	TYR	2.1
1	B	204	ARG	2.1
1	B	218	GLY	2.1
1	B	197	VAL	2.1
1	A	292	THR	2.0
1	B	267	LEU	2.0
1	H	216	ARG	2.0
1	G	291	THR	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 ⓘ

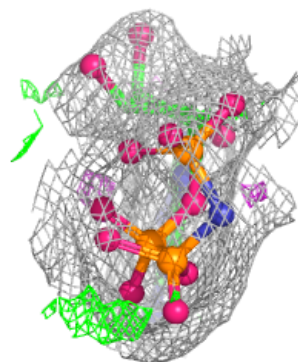
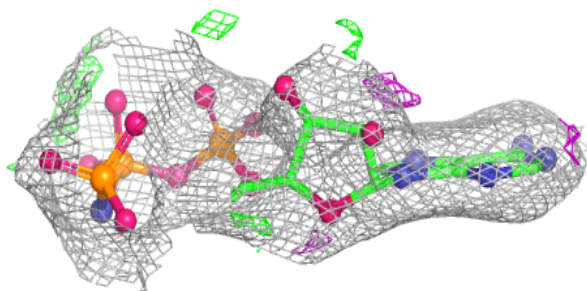
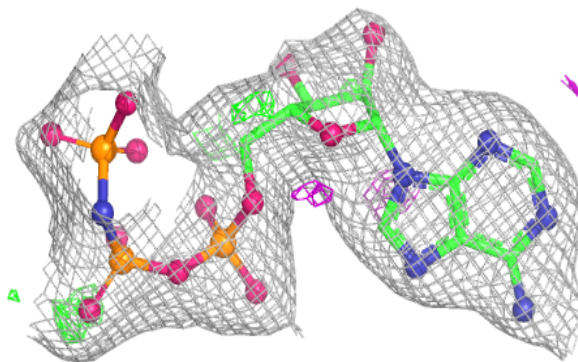
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	E	401	5/5	0.74	0.11	57,61,62,64	0
4	SO4	G	403	5/5	0.80	0.12	50,52,57,57	0
3	MG	B	402	1/1	0.83	0.09	73,73,73,73	0
4	SO4	A	405	5/5	0.83	0.10	53,61,69,92	0
3	MG	F	402	1/1	0.86	0.08	62,62,62,62	0
2	ANP	B	401	31/31	0.87	0.10	48,64,83,90	0
2	ANP	H	401	31/31	0.88	0.09	42,65,92,97	0
4	SO4	H	403	5/5	0.88	0.10	50,58,63,64	0
4	SO4	B	403	5/5	0.89	0.08	63,65,72,75	0
2	ANP	F	401	31/31	0.89	0.09	32,57,76,83	0
4	SO4	D	403	5/5	0.90	0.09	59,59,67,76	0
2	ANP	D	401	31/31	0.90	0.09	33,54,78,85	0
3	MG	H	402	1/1	0.91	0.10	77,77,77,77	0
4	SO4	E	404	5/5	0.92	0.09	52,53,66,66	0
3	MG	D	402	1/1	0.92	0.05	76,76,76,76	0
4	SO4	C	403	5/5	0.92	0.10	48,50,53,59	0
4	SO4	F	403	5/5	0.93	0.08	51,57,61,64	0
4	SO4	C	404	5/5	0.93	0.07	50,56,63,64	0
4	SO4	G	404	5/5	0.93	0.08	50,54,59,64	0
4	SO4	A	404	5/5	0.93	0.08	43,45,51,59	0
4	SO4	C	405	5/5	0.94	0.09	43,44,46,52	0
2	ANP	E	402	31/31	0.96	0.06	27,36,50,55	0
2	ANP	C	401	31/31	0.97	0.06	33,40,47,51	0
2	ANP	A	401	31/31	0.97	0.06	23,32,42,46	0
3	MG	E	403	1/1	0.98	0.07	43,43,43,43	0
4	SO4	A	403	5/5	0.98	0.07	32,34,36,39	0
2	ANP	G	401	31/31	0.98	0.05	28,36,42,47	0
3	MG	C	402	1/1	0.99	0.03	41,41,41,41	0
3	MG	A	402	1/1	0.99	0.04	35,35,35,35	0
3	MG	G	402	1/1	0.99	0.03	36,36,36,36	0

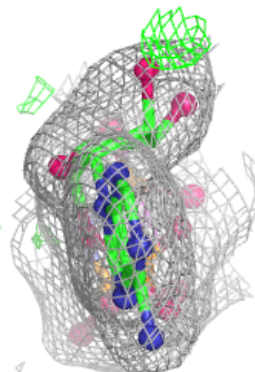
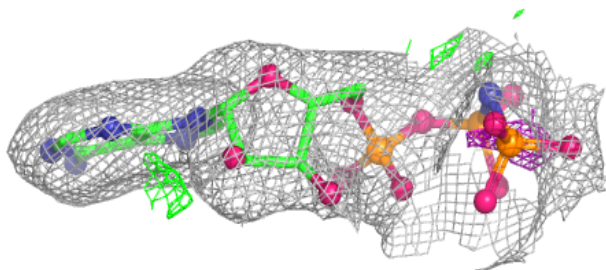
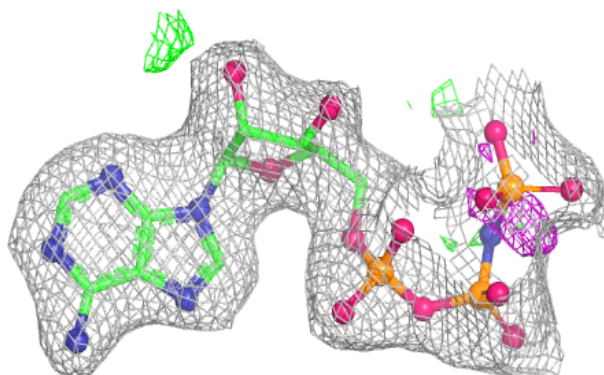
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP B 401:

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

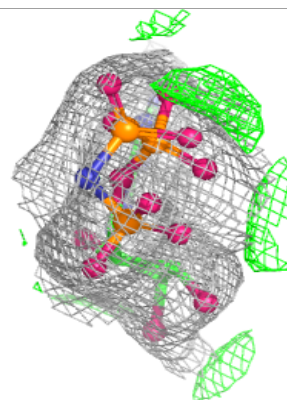
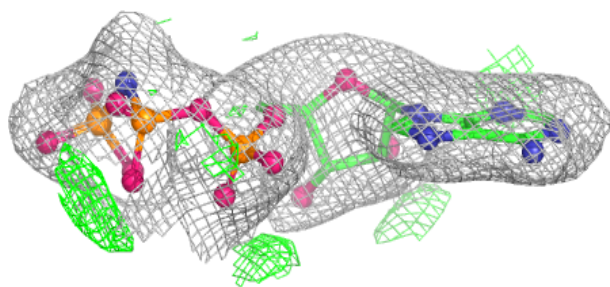
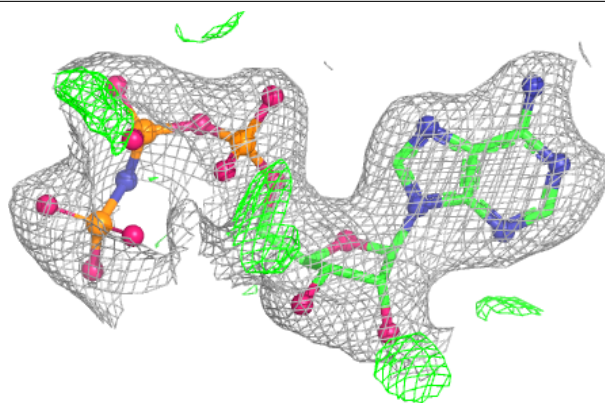
**Electron density around ANP H 401:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

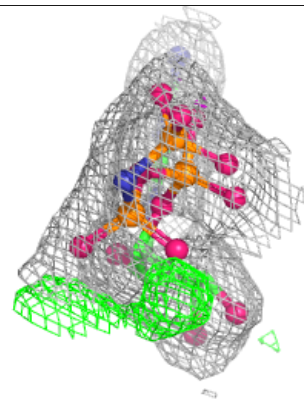
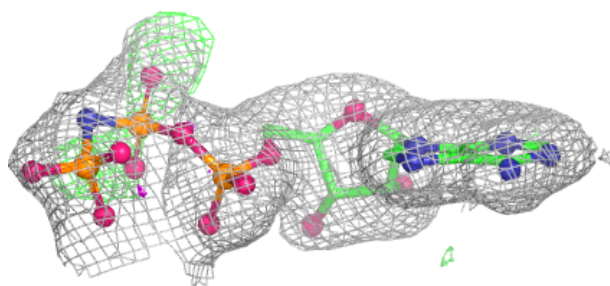
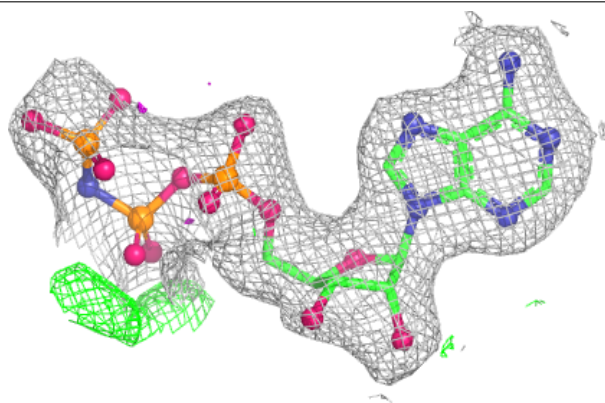


Electron density around ANP F 401:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

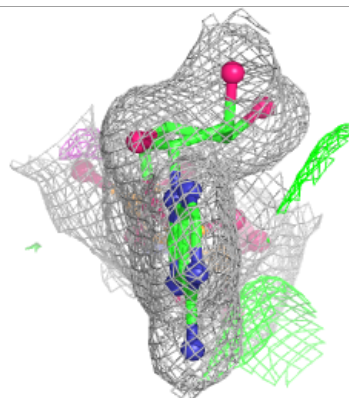
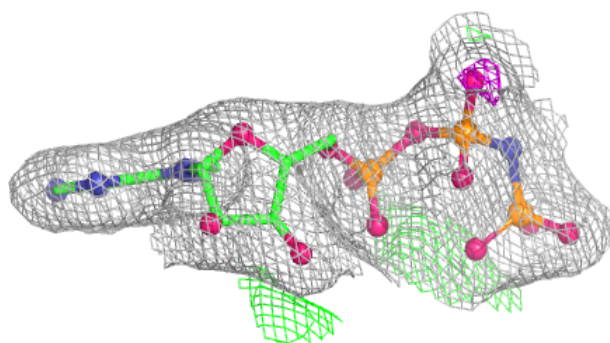
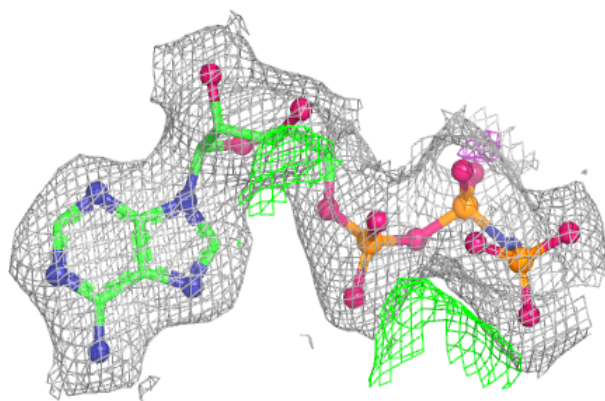
**Electron density around ANP D 401:**

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

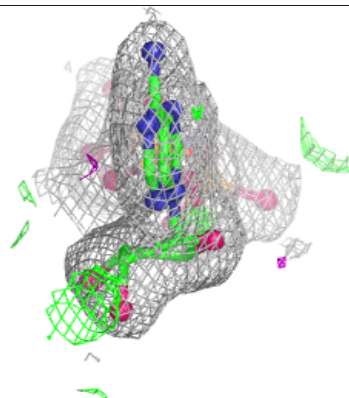
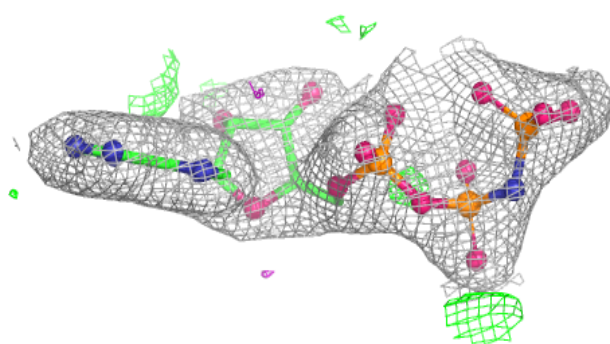
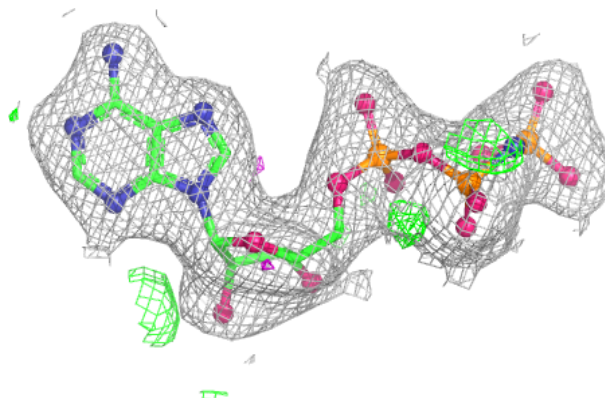


Electron density around ANP E 402:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

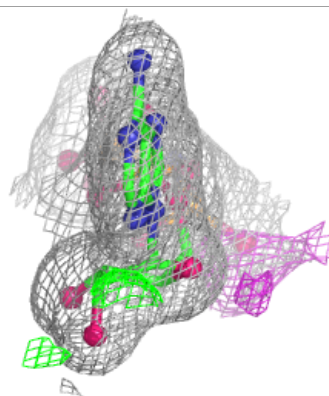
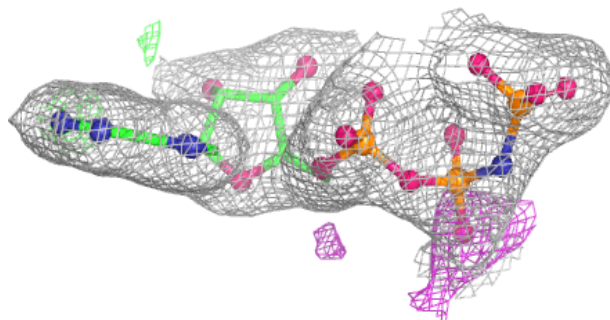
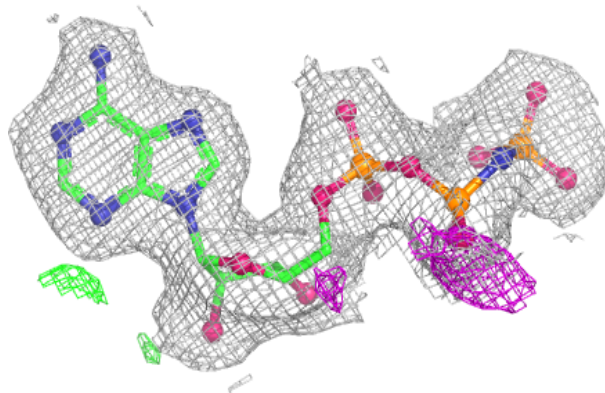
**Electron density around ANP C 401:**

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

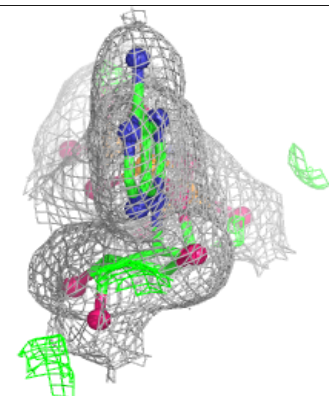
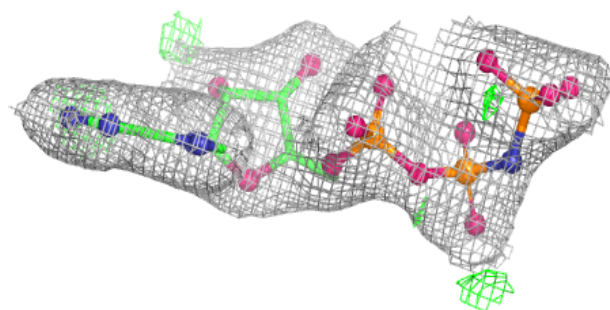
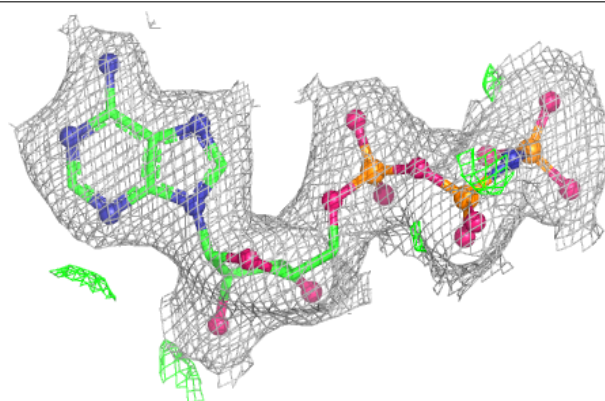


Electron density around ANP A 401:

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

**Electron density around ANP G 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.