



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 29, 2024 – 08:54 am BST

PDB ID : 4BLO
Title : P4 PROTEIN FROM BACTERIOPHAGE PHI6 IN COMPLEX WITH ADP
Authors : El Omari, K.; Meier, C.; Kainov, D.; Sutton, G.; Grimes, J.M.; Poranen, M.M.; Bamford, D.H.; Tuma, R.; Stuart, D.I.; Mancini, E.J.
Deposited on : 2013-05-04
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
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

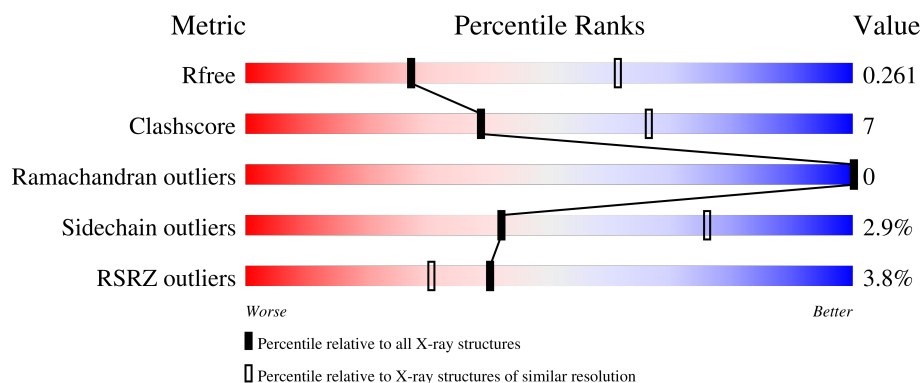
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	309	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	309	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>12%</div> <div>•</div> <div>13%</div> </div> </div>
1	C	309	<div> <div>4%</div> <div> <div></div> <div>72%</div> <div>13%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	309	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>15%</div> <div>•</div> <div>13%</div> </div> </div>
1	E	309	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>15%</div> <div>•</div> <div>13%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	309	<div><div>4%</div><div><div></div><div>72%</div><div>14%</div><div>•</div><div>13%</div></div></div>
1	G	309	<div><div>5%</div><div><div></div><div>71%</div><div>14%</div><div>•</div><div>13%</div></div></div>
1	H	309	<div><div>3%</div><div><div></div><div>71%</div><div>15%</div><div>•</div><div>13%</div></div></div>
1	I	309	<div><div>2%</div><div><div></div><div>72%</div><div>13%</div><div>•</div><div>13%</div></div></div>
1	J	309	<div><div>3%</div><div><div></div><div>72%</div><div>14%</div><div>•</div><div>13%</div></div></div>
1	K	309	<div><div>2%</div><div><div></div><div>72%</div><div>13%</div><div>•</div><div>13%</div></div></div>
1	L	309	<div><div>2%</div><div><div></div><div>70%</div><div>17%</div><div>•</div><div>11%</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24437 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 PACKAGING ENZYME P4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2003	1276	335	382	10			
1	B	270	Total	C	N	O	S	0	0	0
			2011	1281	336	383	11			
1	C	270	Total	C	N	O	S	0	0	0
			2011	1281	336	383	11			
1	D	270	Total	C	N	O	S	0	0	0
			2011	1281	336	383	11			
1	E	269	Total	C	N	O	S	0	0	0
			2006	1278	335	382	11			
1	F	269	Total	C	N	O	S	0	0	0
			2006	1278	335	382	11			
1	G	268	Total	C	N	O	S	0	0	0
			1998	1274	333	380	11			
1	H	269	Total	C	N	O	S	0	0	0
			2003	1276	335	382	10			
1	I	269	Total	C	N	O	S	0	0	0
			2003	1276	335	382	10			
1	J	269	Total	C	N	O	S	0	0	0
			2006	1278	335	382	11			
1	K	269	Total	C	N	O	S	0	0	0
			2003	1276	335	382	10			
1	L	275	Total	C	N	O	S	0	0	0
			2040	1300	341	388	11			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

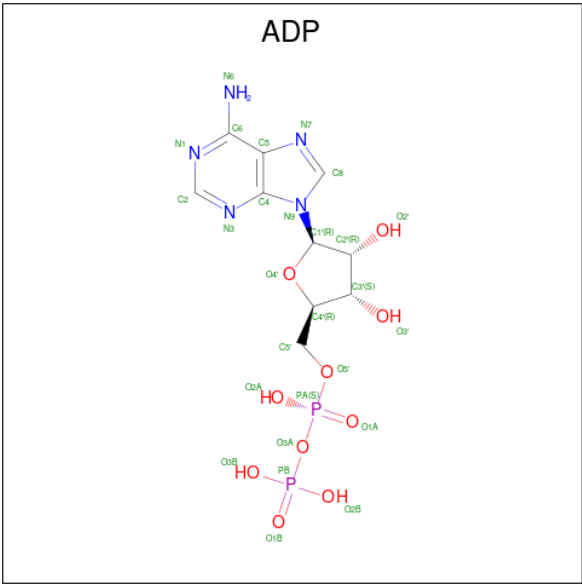
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	E	1	Total	Ca	0	0
			1	1		
2	F	1	Total	Ca	0	0
			1	1		
2	G	1	Total	Ca	0	0
			1	1		
2	H	1	Total	Ca	0	0
			1	1		
2	I	1	Total	Ca	0	0
			1	1		
2	J	1	Total	Ca	0	0
			1	1		
2	K	1	Total	Ca	0	0
			1	1		
2	L	1	Total	Ca	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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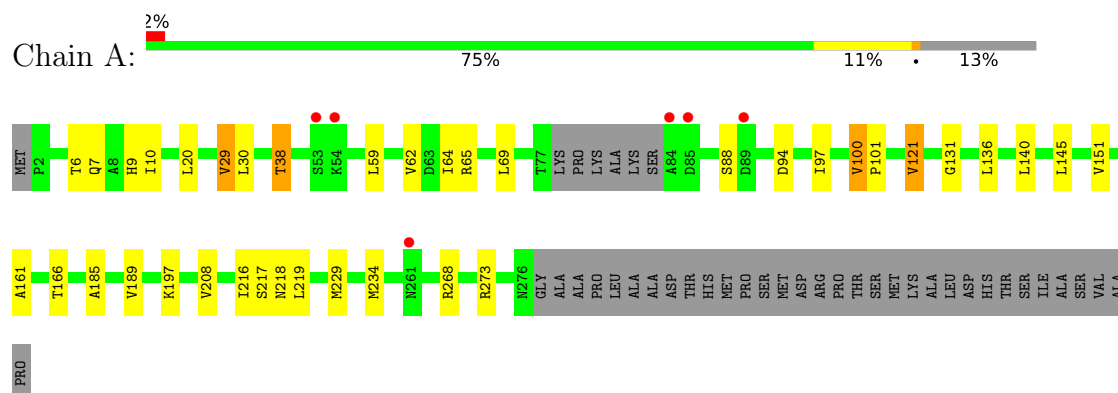
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	I	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	J	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	K	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	L	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

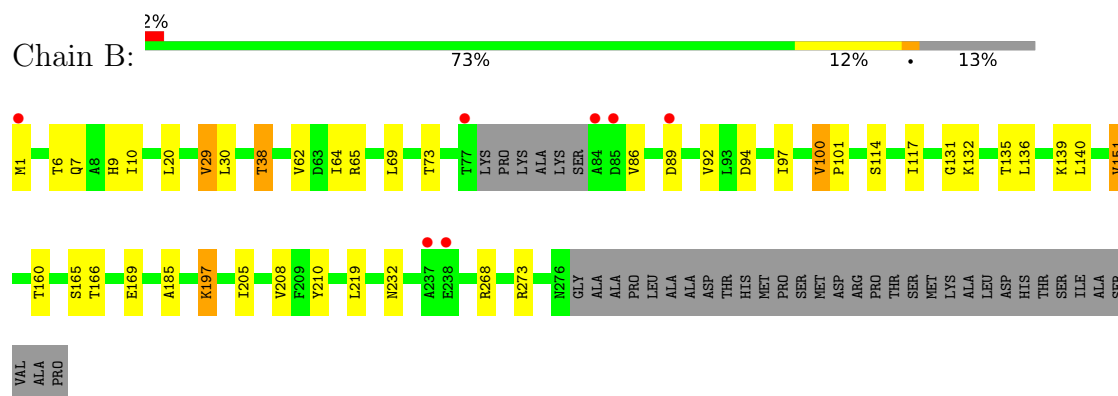
3 Residue-property plots [i](#)

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.

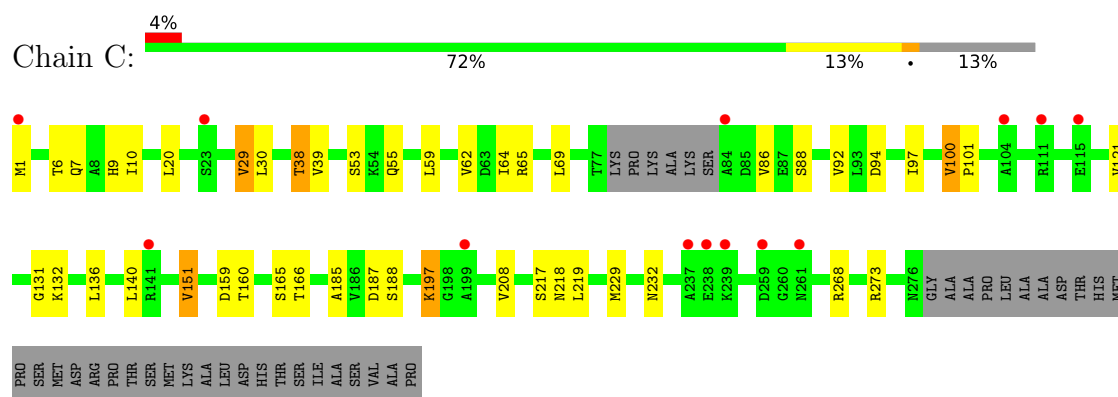
• Molecule 1: PACKAGING ENZYME P4



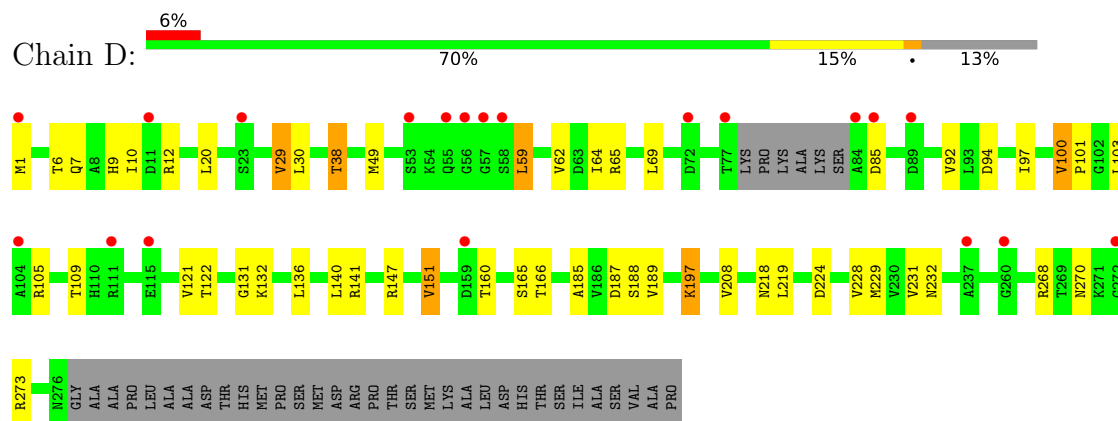
• Molecule 1: PACKAGING ENZYME P4



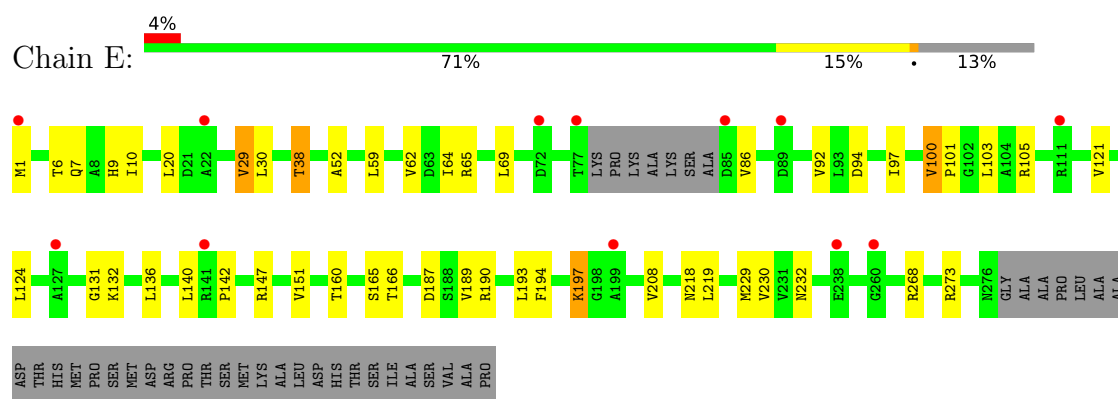
• Molecule 1: PACKAGING ENZYME P4



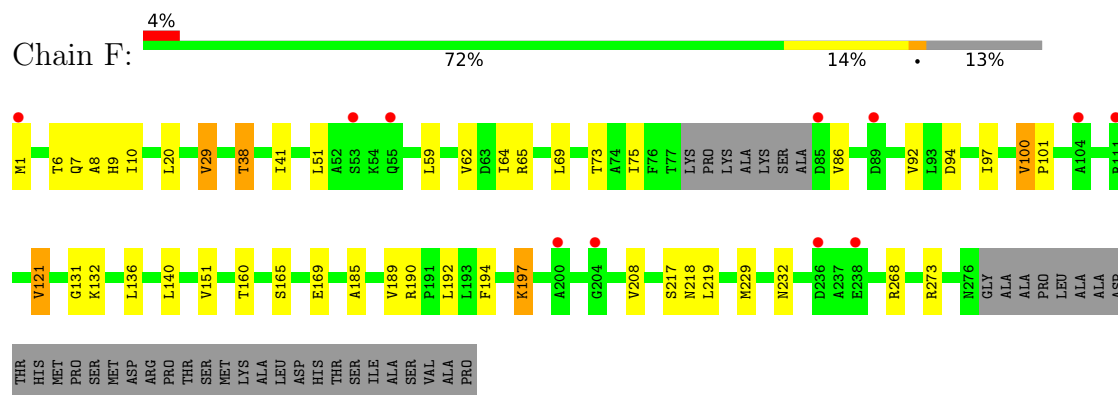
- Molecule 1: PACKAGING ENZYME P4



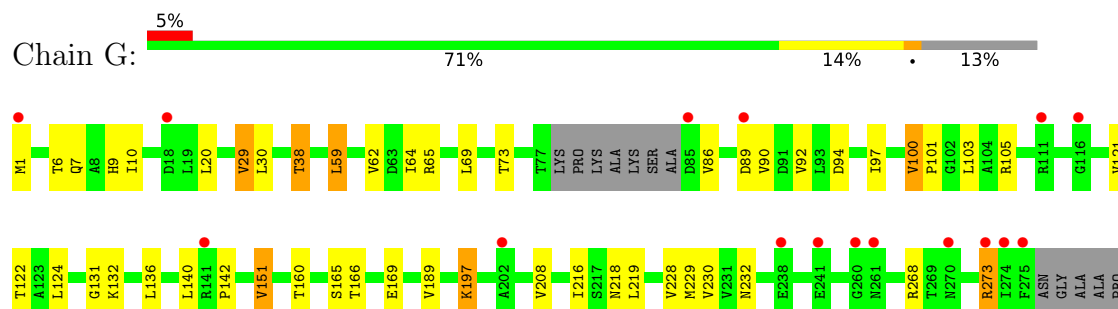
- Molecule 1: PACKAGING ENZYME P4

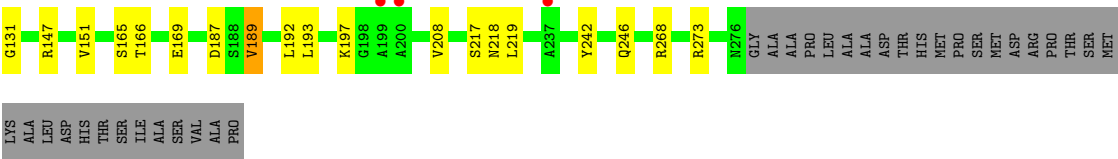


- Molecule 1: PACKAGING ENZYME P4

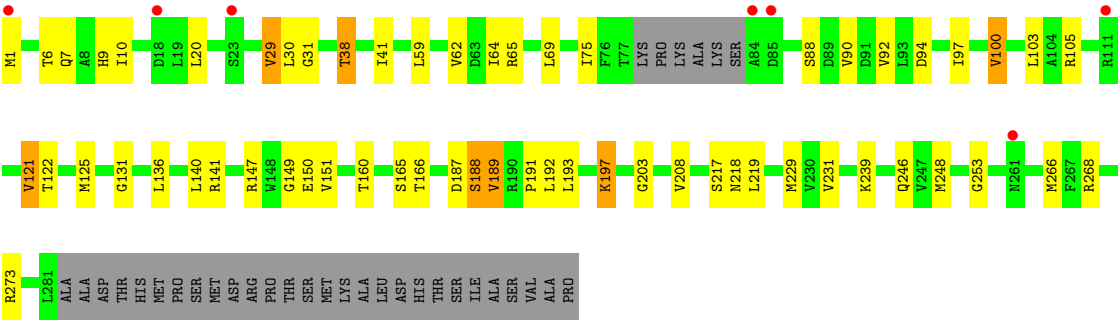


- Molecule 1: PACKAGING ENZYME P4





● Molecule 1: PACKAGING ENZYME P4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	93.84Å 95.26Å 98.33Å 82.46° 86.17° 64.85°	Depositor
Resolution (Å)	19.99 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (19.99-2.80) 96.9 (20.00-2.80)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.79Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.217 , 0.244 0.235 , 0.261	Depositor DCC
R_{free} test set	3664 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	39.7	Xtriage
Anisotropy	0.657	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	24437	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.71% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CA, ADP

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	0.47	0/2033	0.66	2/2767 (0.1%)
1	B	0.48	0/2041	0.67	2/2778 (0.1%)
1	C	0.44	0/2041	0.67	2/2778 (0.1%)
1	D	0.45	0/2041	0.66	2/2778 (0.1%)
1	E	0.44	0/2036	0.66	2/2771 (0.1%)
1	F	0.46	0/2036	0.66	2/2771 (0.1%)
1	G	0.46	0/2028	0.68	2/2760 (0.1%)
1	H	0.46	0/2033	0.66	2/2767 (0.1%)
1	I	0.48	0/2033	0.70	2/2767 (0.1%)
1	J	0.47	0/2036	0.68	2/2771 (0.1%)
1	K	0.46	0/2033	0.68	2/2767 (0.1%)
1	L	0.54	1/2071 (0.0%)	0.70	2/2820 (0.1%)
All	All	0.47	1/24462 (0.0%)	0.67	24/33295 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	31	GLY	C-N	10.48	1.58	1.34

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	29	VAL	CB-CA-C	-7.84	96.51	111.40
1	I	29	VAL	CB-CA-C	-7.83	96.53	111.40
1	J	29	VAL	CB-CA-C	-7.83	96.53	111.40
1	G	29	VAL	CB-CA-C	-7.82	96.53	111.40
1	K	29	VAL	CB-CA-C	-7.82	96.53	111.40
1	D	29	VAL	CB-CA-C	-7.82	96.54	111.40
1	L	29	VAL	CB-CA-C	-7.82	96.55	111.40
1	A	29	VAL	CB-CA-C	-7.82	96.55	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	29	VAL	CB-CA-C	-7.82	96.55	111.40
1	F	29	VAL	CB-CA-C	-7.81	96.55	111.40
1	E	29	VAL	CB-CA-C	-7.81	96.56	111.40
1	C	29	VAL	CB-CA-C	-7.80	96.58	111.40
1	H	29	VAL	N-CA-C	7.79	132.05	111.00
1	G	29	VAL	N-CA-C	7.79	132.03	111.00
1	E	29	VAL	N-CA-C	7.79	132.03	111.00
1	I	29	VAL	N-CA-C	7.79	132.02	111.00
1	K	29	VAL	N-CA-C	7.78	132.02	111.00
1	F	29	VAL	N-CA-C	7.78	132.01	111.00
1	D	29	VAL	N-CA-C	7.78	132.00	111.00
1	A	29	VAL	N-CA-C	7.78	132.00	111.00
1	C	29	VAL	N-CA-C	7.78	132.00	111.00
1	L	29	VAL	N-CA-C	7.78	132.00	111.00
1	J	29	VAL	N-CA-C	7.77	131.99	111.00
1	B	29	VAL	N-CA-C	7.77	131.98	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2003	0	2052	28	0
1	B	2011	0	2063	37	0
1	C	2011	0	2063	35	0
1	D	2011	0	2063	36	0
1	E	2006	0	2058	30	0
1	F	2006	0	2058	37	0
1	G	1998	0	2052	39	0
1	H	2003	0	2052	40	0
1	I	2003	0	2052	32	0
1	J	2006	0	2058	40	0
1	K	2003	0	2052	40	0
1	L	2040	0	2094	39	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	27	0	12	1	0
3	B	27	0	12	1	0
3	C	27	0	12	1	0
3	D	27	0	12	1	0
3	E	27	0	12	1	0
3	F	27	0	12	1	0
3	G	27	0	12	1	0
3	H	27	0	12	1	0
3	I	27	0	12	0	0
3	J	27	0	12	1	0
3	K	27	0	12	1	0
3	L	27	0	12	1	0
All	All	24437	0	24861	368	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 7.

All (368) 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:29:VAL:O	1:A:208:VAL:HG21	1.42	1.18
1:F:29:VAL:O	1:F:208:VAL:HG21	1.45	1.17
1:L:29:VAL:O	1:L:208:VAL:HG21	1.43	1.17
1:K:29:VAL:O	1:K:208:VAL:HG21	1.42	1.16
1:H:29:VAL:O	1:H:208:VAL:HG21	1.47	1.13
1:G:29:VAL:O	1:G:208:VAL:HG21	1.46	1.13
1:I:29:VAL:O	1:I:208:VAL:HG21	1.49	1.12
1:E:29:VAL:O	1:E:208:VAL:HG21	1.45	1.12
1:C:29:VAL:O	1:C:208:VAL:HG21	1.49	1.12
1:D:29:VAL:O	1:D:208:VAL:HG21	1.47	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:29:VAL:O	1:J:208:VAL:HG21	1.46	1.10
1:B:29:VAL:O	1:B:208:VAL:HG21	1.49	1.10
1:C:30:LEU:O	1:D:197:LYS:NZ	2.03	0.92
1:D:30:LEU:O	1:E:197:LYS:NZ	2.01	0.91
1:K:52:ALA:HA	1:K:59:LEU:HD11	1.51	0.91
1:J:88:SER:O	1:K:197:LYS:HD2	1.75	0.85
1:J:73:THR:HG21	1:J:169:GLU:CG	2.07	0.85
1:H:73:THR:HG21	1:H:169:GLU:CG	2.07	0.84
1:B:73:THR:HG21	1:B:169:GLU:CG	2.08	0.83
1:C:197:LYS:NZ	1:H:30:LEU:O	2.12	0.82
1:D:94:ASP:OD2	1:E:86:VAL:HG12	1.79	0.82
1:G:73:THR:HG21	1:G:169:GLU:CG	2.10	0.82
1:K:73:THR:HG21	1:K:169:GLU:CG	2.10	0.82
1:F:73:THR:HG21	1:F:169:GLU:CG	2.10	0.82
1:B:30:LEU:O	1:I:197:LYS:NZ	2.14	0.81
1:C:6:THR:H	1:C:9:HIS:HD2	1.31	0.78
1:K:6:THR:H	1:K:9:HIS:HD2	1.30	0.78
1:B:6:THR:H	1:B:9:HIS:HD2	1.31	0.77
1:G:6:THR:H	1:G:9:HIS:HD2	1.31	0.77
1:H:6:THR:H	1:H:9:HIS:HD2	1.33	0.76
1:L:6:THR:H	1:L:9:HIS:HD2	1.32	0.75
1:J:6:THR:H	1:J:9:HIS:HD2	1.34	0.75
1:J:90:VAL:O	1:K:197:LYS:HE3	1.87	0.75
1:G:89:ASP:OD1	1:H:88:SER:HB2	1.87	0.75
1:I:6:THR:H	1:I:9:HIS:HD2	1.33	0.74
1:D:6:THR:H	1:D:9:HIS:HD2	1.32	0.74
1:A:6:THR:H	1:A:9:HIS:HD2	1.34	0.74
1:C:86:VAL:HG12	1:H:94:ASP:OD2	1.89	0.73
1:E:52:ALA:HA	1:E:59:LEU:HD11	1.71	0.73
1:A:30:LEU:O	1:B:197:LYS:NZ	2.21	0.73
1:E:30:LEU:O	1:F:197:LYS:NZ	2.21	0.73
1:G:30:LEU:O	1:H:197:LYS:NZ	2.22	0.72
1:F:136:LEU:HA	1:F:140:LEU:HD12	1.74	0.70
1:H:73:THR:HG21	1:H:169:GLU:HG2	1.72	0.70
1:B:73:THR:HG21	1:B:169:GLU:HG3	1.74	0.70
1:C:136:LEU:HA	1:C:140:LEU:HD12	1.74	0.70
1:E:94:ASP:OD2	1:F:86:VAL:HG12	1.92	0.70
1:B:94:ASP:OD2	1:I:86:VAL:HG12	1.92	0.69
1:J:73:THR:HG21	1:J:169:GLU:HG3	1.73	0.69
1:F:94:ASP:OD2	1:G:86:VAL:HG12	1.93	0.69
1:A:136:LEU:HA	1:A:140:LEU:HD12	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:136:LEU:HA	1:I:140:LEU:HD12	1.75	0.69
1:E:136:LEU:HA	1:E:140:LEU:HD12	1.75	0.69
1:H:73:THR:HG21	1:H:169:GLU:HG3	1.75	0.68
1:K:73:THR:HG21	1:K:169:GLU:HG3	1.75	0.68
1:B:136:LEU:HA	1:B:140:LEU:HD12	1.76	0.68
1:G:136:LEU:HA	1:G:140:LEU:HD12	1.77	0.67
1:F:73:THR:HG21	1:F:169:GLU:HG2	1.75	0.67
1:G:73:THR:HG21	1:G:169:GLU:HG3	1.77	0.66
1:J:73:THR:HG21	1:J:169:GLU:HG2	1.78	0.66
1:E:6:THR:H	1:E:9:HIS:HD2	1.44	0.65
1:K:73:THR:HG21	1:K:169:GLU:HG2	1.79	0.65
1:D:101:PRO:HG3	1:E:165:SER:HB3	1.78	0.65
1:J:94:ASP:OD2	1:K:86:VAL:HG12	1.97	0.65
1:G:73:THR:HG21	1:G:169:GLU:HG2	1.79	0.64
1:C:101:PRO:HG3	1:D:165:SER:HB3	1.79	0.64
1:J:73:THR:CG2	1:J:169:GLU:HG3	2.28	0.63
1:A:94:ASP:OD2	1:B:86:VAL:HG12	1.98	0.63
1:J:90:VAL:O	1:K:197:LYS:CE	2.45	0.63
1:F:73:THR:HG21	1:F:169:GLU:HG3	1.79	0.63
1:H:136:LEU:HA	1:H:140:LEU:HD12	1.80	0.63
1:C:7:GLN:HA	1:C:10:ILE:HD12	1.80	0.63
1:B:73:THR:HG21	1:B:169:GLU:HG2	1.78	0.63
1:I:132:LYS:HE3	1:I:232:ASN:HB3	1.80	0.62
1:I:94:ASP:OD2	1:J:86:VAL:HG12	1.99	0.62
1:I:88:SER:O	1:J:197:LYS:HD2	2.00	0.62
1:B:73:THR:CG2	1:B:169:GLU:HG3	2.30	0.61
1:L:189:VAL:HG13	1:L:193:LEU:HD13	1.81	0.61
1:J:73:THR:CG2	1:J:169:GLU:CG	2.79	0.61
1:B:101:PRO:HG3	1:I:165:SER:HB3	1.83	0.60
1:B:89:ASP:OD1	1:I:88:SER:HB2	2.02	0.60
1:K:73:THR:CG2	1:K:169:GLU:HG3	2.32	0.59
1:C:88:SER:HB2	1:H:89:ASP:OD1	2.02	0.59
1:K:52:ALA:CA	1:K:59:LEU:HD11	2.28	0.59
1:G:94:ASP:OD2	1:H:86:VAL:HG12	2.03	0.59
1:C:94:ASP:OD2	1:D:85:ASP:HA	2.02	0.59
1:F:6:THR:H	1:F:9:HIS:HD2	1.51	0.58
1:J:7:GLN:HA	1:J:10:ILE:HD12	1.85	0.58
1:G:73:THR:CG2	1:G:169:GLU:HG3	2.32	0.58
1:L:203:GLY:HA2	1:L:239:LYS:HD2	1.85	0.58
1:B:205:ILE:HD11	1:B:210:TYR:OH	2.02	0.58
1:B:73:THR:CG2	1:B:169:GLU:CG	2.81	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:73:THR:CG2	1:G:169:GLU:CG	2.82	0.57
1:L:136:LEU:HA	1:L:140:LEU:HD12	1.86	0.57
1:B:7:GLN:HA	1:B:10:ILE:HD12	1.86	0.57
1:D:62:VAL:O	1:D:65:ARG:HD2	2.05	0.57
1:F:73:THR:CG2	1:F:169:GLU:CG	2.82	0.57
1:K:62:VAL:O	1:K:65:ARG:HD2	2.04	0.57
1:G:62:VAL:O	1:G:65:ARG:HD2	2.04	0.56
1:L:7:GLN:HA	1:L:10:ILE:HD12	1.88	0.56
1:A:7:GLN:HA	1:A:10:ILE:HD12	1.88	0.56
1:F:73:THR:CG2	1:F:169:GLU:HG3	2.34	0.56
1:G:7:GLN:HA	1:G:10:ILE:HD12	1.88	0.56
1:H:62:VAL:O	1:H:65:ARG:HD2	2.06	0.56
1:H:73:THR:CG2	1:H:169:GLU:CG	2.83	0.56
1:H:73:THR:CG2	1:H:169:GLU:HG3	2.35	0.56
1:I:7:GLN:HA	1:I:10:ILE:HD12	1.87	0.56
1:K:189:VAL:HG13	1:K:193:LEU:HD13	1.88	0.56
1:C:121:VAL:HG12	1:C:217:SER:OG	2.06	0.56
1:F:6:THR:HG22	1:F:8:ALA:H	1.71	0.56
1:A:62:VAL:O	1:A:65:ARG:HD2	2.06	0.55
1:D:12:ARG:HH22	1:D:224:ASP:HB2	1.70	0.55
1:F:62:VAL:O	1:F:65:ARG:HD2	2.06	0.55
1:J:101:PRO:HG3	1:K:165:SER:HB3	1.87	0.55
1:K:7:GLN:HA	1:K:10:ILE:HD12	1.88	0.55
1:B:62:VAL:O	1:B:65:ARG:HD2	2.06	0.55
1:I:50:GLU:O	1:I:54:LYS:HG2	2.06	0.55
1:J:62:VAL:O	1:J:65:ARG:HD2	2.07	0.55
1:K:88:SER:O	1:L:197:LYS:HD2	2.06	0.55
1:F:7:GLN:HA	1:F:10:ILE:HD12	1.88	0.55
1:D:7:GLN:HA	1:D:10:ILE:HD12	1.89	0.55
1:C:165:SER:HB3	1:H:101:PRO:HG3	1.90	0.54
1:C:62:VAL:O	1:C:65:ARG:HD2	2.07	0.54
1:E:62:VAL:O	1:E:65:ARG:HD2	2.06	0.54
1:L:62:VAL:O	1:L:65:ARG:HD2	2.07	0.54
1:C:53:SER:O	1:C:55:GLN:O	2.24	0.54
1:F:6:THR:H	1:F:9:HIS:CD2	2.26	0.54
1:I:62:VAL:O	1:I:65:ARG:HD2	2.08	0.54
1:K:59:LEU:HD23	1:K:64:ILE:CG2	2.38	0.54
1:C:39:VAL:HG11	1:J:262:VAL:O	2.08	0.54
1:K:131:GLY:HA2	3:K:401:ADP:H5'1	1.90	0.54
1:F:20:LEU:HB3	1:F:38:THR:HG23	1.90	0.53
1:A:30:LEU:HD12	1:A:197:LYS:O	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:7:GLN:HA	1:E:10:ILE:HD12	1.89	0.53
1:H:7:GLN:HA	1:H:10:ILE:HD12	1.89	0.53
1:K:73:THR:CG2	1:K:169:GLU:CG	2.82	0.53
1:J:41:ILE:HD11	1:J:75:ILE:HD11	1.90	0.53
1:E:190:ARG:HD2	1:E:194:PHE:CE2	2.44	0.53
1:A:166:THR:HA	1:L:97:ILE:O	2.09	0.53
1:G:218:ASN:HD21	1:H:151:VAL:HA	1.73	0.52
1:E:189:VAL:HG13	1:E:193:LEU:HD13	1.92	0.52
1:B:92:VAL:CG1	1:B:97:ILE:HD11	2.40	0.52
1:C:20:LEU:HB3	1:C:38:THR:HG23	1.91	0.52
1:A:197:LYS:HE3	1:L:90:VAL:O	2.10	0.51
1:D:121:VAL:HG21	1:D:229:MET:HE2	1.92	0.51
1:G:101:PRO:HG3	1:H:165:SER:HB3	1.92	0.51
1:A:121:VAL:HG21	1:A:229:MET:HE2	1.92	0.51
1:I:20:LEU:HB3	1:I:38:THR:HG23	1.92	0.51
1:K:20:LEU:HB3	1:K:38:THR:HG23	1.91	0.51
1:K:89:ASP:OD1	1:L:88:SER:HB2	2.11	0.51
1:I:268:ARG:HG2	1:I:273:ARG:HG2	1.92	0.51
1:E:20:LEU:HB3	1:E:38:THR:HG23	1.92	0.51
1:H:92:VAL:HG12	1:H:97:ILE:HD11	1.91	0.51
1:J:20:LEU:HB3	1:J:38:THR:HG23	1.93	0.51
1:C:92:VAL:HG12	1:C:97:ILE:HD11	1.91	0.51
1:B:20:LEU:HB3	1:B:38:THR:HG23	1.91	0.51
1:C:218:ASN:HD21	1:D:151:VAL:HA	1.75	0.51
1:B:268:ARG:HG2	1:B:273:ARG:HG2	1.92	0.51
1:L:121:VAL:HG21	1:L:229:MET:HE2	1.92	0.51
1:D:20:LEU:HB3	1:D:38:THR:HG23	1.91	0.51
1:K:59:LEU:HD23	1:K:64:ILE:HG21	1.93	0.50
1:E:101:PRO:HG3	1:F:165:SER:HB3	1.92	0.50
1:K:268:ARG:HG2	1:K:273:ARG:HG2	1.94	0.50
1:I:218:ASN:HD21	1:J:151:VAL:HA	1.76	0.50
1:A:197:LYS:HD2	1:L:88:SER:O	2.12	0.50
1:C:151:VAL:HA	1:H:218:ASN:HD21	1.77	0.50
1:L:20:LEU:HB3	1:L:38:THR:HG23	1.94	0.50
1:L:189:VAL:HG11	1:L:231:VAL:HG12	1.94	0.50
1:A:101:PRO:HG3	1:B:165:SER:HB3	1.94	0.49
1:D:136:LEU:HD11	1:D:185:ALA:HB1	1.93	0.49
1:J:268:ARG:HG2	1:J:273:ARG:HG2	1.94	0.49
1:G:20:LEU:HB3	1:G:38:THR:HG23	1.93	0.49
1:I:221:THR:HG21	1:J:152:ALA:HB2	1.95	0.49
1:A:20:LEU:HB3	1:A:38:THR:HG23	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:248:MET:HE2	1:L:266:MET:HB2	1.95	0.49
1:H:265:THR:O	1:H:275:PHE:O	2.30	0.49
1:H:268:ARG:HG2	1:H:273:ARG:HG2	1.94	0.49
1:B:1:MET:O	1:B:160:THR:HG23	2.13	0.49
1:C:268:ARG:HG2	1:C:273:ARG:HG2	1.94	0.49
1:H:92:VAL:CG1	1:H:97:ILE:HD11	2.42	0.49
1:G:97:ILE:O	1:H:166:THR:HA	2.12	0.49
1:E:268:ARG:HG2	1:E:273:ARG:HG2	1.94	0.49
1:G:1:MET:O	1:G:160:THR:HG23	2.13	0.49
1:J:136:LEU:O	1:J:140:LEU:HB2	2.12	0.49
1:G:122:THR:HB	1:G:228:VAL:HG22	1.95	0.49
1:G:124:LEU:HB3	1:G:230:VAL:HA	1.95	0.49
1:L:92:VAL:HG12	1:L:97:ILE:HD11	1.95	0.49
1:L:147:ARG:HA	1:L:187:ASP:O	2.12	0.48
1:D:136:LEU:O	1:D:140:LEU:HB2	2.13	0.48
1:G:132:LYS:HE3	1:G:232:ASN:HB2	1.95	0.48
1:D:268:ARG:HG2	1:D:273:ARG:HG2	1.96	0.48
1:A:97:ILE:O	1:B:166:THR:HA	2.14	0.48
1:A:268:ARG:HG2	1:A:273:ARG:HG2	1.96	0.48
1:H:132:LYS:HE3	1:H:232:ASN:HB2	1.96	0.48
1:K:97:ILE:O	1:L:166:THR:HA	2.14	0.48
1:J:1:MET:O	1:J:160:THR:HG23	2.14	0.48
1:K:29:VAL:O	1:K:208:VAL:CG2	2.36	0.48
1:L:122:THR:HA	1:L:253:GLY:O	2.13	0.48
1:J:73:THR:CG2	1:J:169:GLU:HG2	2.43	0.47
1:F:1:MET:O	1:F:160:THR:HG23	2.14	0.47
1:C:132:LYS:HE3	1:C:232:ASN:HB2	1.97	0.47
1:H:41:ILE:HD11	1:H:75:ILE:HD11	1.96	0.47
1:A:145:LEU:O	1:A:161:ALA:HA	2.14	0.47
1:D:92:VAL:HG12	1:D:97:ILE:HD11	1.96	0.47
1:D:109:THR:HG21	1:D:270:ASN:HD22	1.79	0.47
1:H:216:ILE:HG21	1:H:229:MET:HE1	1.95	0.47
1:J:132:LYS:HE3	1:J:232:ASN:HB2	1.96	0.47
1:D:1:MET:O	1:D:160:THR:HG23	2.14	0.47
1:I:101:PRO:HG3	1:J:165:SER:HB3	1.97	0.47
1:G:92:VAL:HG12	1:G:97:ILE:HD11	1.97	0.47
1:D:6:THR:H	1:D:9:HIS:CD2	2.23	0.46
1:F:92:VAL:HG12	1:F:97:ILE:HD11	1.96	0.46
1:L:1:MET:O	1:L:160:THR:HG23	2.15	0.46
1:K:147:ARG:HA	1:K:187:ASP:O	2.16	0.46
1:L:100:VAL:HG11	1:L:219:LEU:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:52:ALA:CA	1:E:59:LEU:HD11	2.44	0.46
1:E:121:VAL:HG21	1:E:229:MET:HE2	1.98	0.46
1:C:121:VAL:HG21	1:C:229:MET:HE2	1.97	0.46
1:C:100:VAL:HG11	1:C:219:LEU:HD13	1.98	0.46
1:F:268:ARG:HG2	1:F:273:ARG:HG2	1.96	0.46
1:H:100:VAL:HG11	1:H:219:LEU:HD13	1.98	0.46
1:B:132:LYS:HE3	1:B:232:ASN:HB2	1.96	0.46
1:E:1:MET:O	1:E:160:THR:HG23	2.16	0.45
1:F:97:ILE:O	1:G:166:THR:HA	2.16	0.45
1:E:132:LYS:HE3	1:E:232:ASN:HB2	1.99	0.45
1:B:6:THR:H	1:B:9:HIS:CD2	2.22	0.45
1:I:51:LEU:HD23	1:I:59:LEU:HD22	1.99	0.45
1:D:132:LYS:HE3	1:D:232:ASN:HB2	1.98	0.45
1:C:1:MET:O	1:C:160:THR:HG23	2.17	0.45
1:A:100:VAL:HG11	1:A:219:LEU:HD13	1.99	0.45
1:F:131:GLY:HA2	3:F:401:ADP:H5'1	1.99	0.45
1:F:190:ARG:HD2	1:F:194:PHE:CE2	2.52	0.45
1:G:216:ILE:HG21	1:G:229:MET:HE1	1.99	0.45
1:G:131:GLY:HA2	3:G:401:ADP:H5'1	1.99	0.45
1:J:136:LEU:HD11	1:J:185:ALA:HB1	1.98	0.45
1:A:234:MET:HE1	1:L:246:GLN:OE1	2.16	0.45
1:G:268:ARG:HG2	1:G:273:ARG:HG3	1.99	0.45
1:A:218:ASN:HD21	1:B:151:VAL:HA	1.82	0.45
1:G:100:VAL:HG11	1:G:219:LEU:HD13	1.99	0.45
1:F:6:THR:HB	1:F:9:HIS:CD2	2.53	0.44
1:I:64:ILE:HA	1:I:69:LEU:HD11	1.99	0.44
1:J:90:VAL:H	1:K:197:LYS:HE3	1.82	0.44
1:C:136:LEU:HD11	1:C:185:ALA:HB1	1.99	0.44
1:A:88:SER:O	1:B:197:LYS:HD2	2.18	0.44
1:E:147:ARG:HA	1:E:187:ASP:O	2.17	0.44
1:B:100:VAL:HG11	1:B:219:LEU:HD13	2.00	0.44
1:I:92:VAL:CG1	1:I:97:ILE:HD11	2.48	0.44
1:L:131:GLY:HA2	3:L:401:ADP:H5'1	1.99	0.44
1:B:92:VAL:HG12	1:B:97:ILE:HD11	2.00	0.44
1:F:121:VAL:HG21	1:F:229:MET:HE2	1.98	0.44
1:L:30:LEU:HD12	1:L:197:LYS:O	2.17	0.44
1:F:64:ILE:HA	1:F:69:LEU:HD11	2.00	0.44
1:L:41:ILE:HD11	1:L:75:ILE:HD11	1.99	0.44
1:D:131:GLY:HA2	3:D:401:ADP:H5'1	2.01	0.43
1:E:92:VAL:HG12	1:E:97:ILE:HD11	1.98	0.43
1:E:100:VAL:HG11	1:E:219:LEU:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:20:LEU:HB3	1:H:38:THR:HB	2.00	0.43
1:I:55:GLN:OE1	1:I:59:LEU:HD23	2.18	0.43
1:D:64:ILE:HA	1:D:69:LEU:HD11	2.00	0.43
1:D:147:ARG:HA	1:D:187:ASP:O	2.18	0.43
1:H:73:THR:CG2	1:H:169:GLU:HG2	2.43	0.43
1:L:100:VAL:CG2	1:L:218:ASN:HB2	2.48	0.43
1:B:30:LEU:HD12	1:B:197:LYS:O	2.18	0.43
1:B:136:LEU:HD11	1:B:185:ALA:HB1	2.00	0.43
1:D:100:VAL:HG11	1:D:219:LEU:HD13	2.00	0.43
1:E:103:LEU:HD22	1:E:105:ARG:HG2	2.00	0.43
1:E:131:GLY:HA2	3:E:401:ADP:H5'1	2.00	0.43
1:F:73:THR:CG2	1:F:169:GLU:HG2	2.43	0.43
1:F:92:VAL:CG1	1:F:97:ILE:HD11	2.48	0.43
1:G:64:ILE:HA	1:G:69:LEU:HD11	2.01	0.43
1:I:90:VAL:O	1:J:197:LYS:HE3	2.18	0.43
1:B:64:ILE:HA	1:B:69:LEU:HD11	2.00	0.43
1:A:131:GLY:HA2	3:A:401:ADP:H5'1	2.00	0.43
1:C:6:THR:H	1:C:9:HIS:CD2	2.22	0.43
1:H:64:ILE:HA	1:H:69:LEU:HD11	2.01	0.43
1:I:100:VAL:HG11	1:I:219:LEU:HD13	2.01	0.43
1:K:64:ILE:HA	1:K:69:LEU:HD11	1.99	0.43
1:I:149:GLY:HA3	1:I:191:PRO:HG3	2.01	0.43
1:J:64:ILE:HA	1:J:69:LEU:HD11	2.00	0.43
1:K:98:TYR:HA	1:L:165:SER:O	2.19	0.43
1:C:64:ILE:HA	1:C:69:LEU:HD11	2.01	0.43
1:K:121:VAL:HG12	1:K:217:SER:HB2	2.00	0.43
1:L:149:GLY:HA3	1:L:191:PRO:HG3	2.00	0.43
1:F:101:PRO:HG3	1:G:165:SER:HB3	2.01	0.43
1:I:197:LYS:H	1:I:197:LYS:HG2	1.63	0.43
1:J:30:LEU:HD12	1:J:197:LYS:O	2.19	0.43
1:J:100:VAL:HG11	1:J:219:LEU:HD13	2.01	0.43
1:K:242:TYR:O	1:K:246:GLN:HG2	2.19	0.43
1:L:141:ARG:NH1	1:L:141:ARG:HB3	2.34	0.43
1:A:136:LEU:HD11	1:A:185:ALA:HB1	2.00	0.42
1:D:92:VAL:CG1	1:D:97:ILE:HD11	2.49	0.42
1:I:92:VAL:HG12	1:I:97:ILE:HD11	2.01	0.42
1:C:131:GLY:HA2	3:C:401:ADP:H5'1	2.01	0.42
1:H:100:VAL:CG2	1:H:218:ASN:HB2	2.49	0.42
1:K:103:LEU:HD22	1:K:105:ARG:HG2	2.01	0.42
1:B:73:THR:CG2	1:B:169:GLU:HG2	2.46	0.42
1:G:30:LEU:HD12	1:G:197:LYS:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:101:PRO:HG3	1:L:165:SER:HB3	1.99	0.42
1:K:100:VAL:HG11	1:K:219:LEU:HD13	2.00	0.42
1:B:114:SER:O	1:B:117:ILE:HG22	2.19	0.42
1:B:131:GLY:HA2	3:B:401:ADP:H5'1	2.01	0.42
1:F:41:ILE:HD11	1:F:75:ILE:HD11	2.01	0.42
1:L:268:ARG:HG2	1:L:273:ARG:HG2	2.02	0.42
1:A:197:LYS:CD	1:L:88:SER:O	2.68	0.42
1:E:100:VAL:CG2	1:E:218:ASN:HB2	2.50	0.42
1:A:64:ILE:HA	1:A:69:LEU:HD11	2.01	0.42
1:D:103:LEU:HD22	1:D:105:ARG:HG2	2.01	0.42
1:D:189:VAL:HG11	1:D:231:VAL:HG12	2.01	0.42
1:E:64:ILE:HA	1:E:69:LEU:HD11	2.01	0.42
1:F:100:VAL:CG2	1:F:218:ASN:HB2	2.50	0.42
1:F:100:VAL:HG11	1:F:219:LEU:HD13	2.01	0.42
1:F:136:LEU:HD11	1:F:185:ALA:HB1	2.01	0.42
1:H:103:LEU:HD22	1:H:105:ARG:HG2	2.02	0.42
1:A:100:VAL:CG2	1:A:218:ASN:HB2	2.50	0.42
1:F:132:LYS:HE3	1:F:232:ASN:HB2	2.01	0.42
1:H:131:GLY:HA2	3:H:401:ADP:H5'1	2.01	0.42
1:L:103:LEU:HD22	1:L:105:ARG:HG2	2.02	0.42
1:A:121:VAL:HG12	1:A:217:SER:HB2	2.02	0.42
1:C:166:THR:HA	1:H:97:ILE:O	2.19	0.42
1:J:131:GLY:HA2	3:J:401:ADP:H5'1	2.02	0.42
1:K:100:VAL:CG2	1:K:218:ASN:HB2	2.50	0.41
1:C:97:ILE:O	1:D:166:THR:HA	2.20	0.41
1:C:100:VAL:CG2	1:C:218:ASN:HB2	2.50	0.41
1:D:10:ILE:HG12	1:D:49:MET:HE1	2.02	0.41
1:G:73:THR:CG2	1:G:169:GLU:HG2	2.45	0.41
1:I:136:LEU:HD11	1:I:185:ALA:HB1	2.02	0.41
1:L:64:ILE:HA	1:L:69:LEU:HD11	2.01	0.41
1:F:218:ASN:HD21	1:G:151:VAL:HA	1.85	0.41
1:J:97:ILE:O	1:K:166:THR:HA	2.19	0.41
1:K:73:THR:CG2	1:K:169:GLU:HG2	2.45	0.41
1:G:103:LEU:HD22	1:G:105:ARG:HG2	2.01	0.41
1:H:149:GLY:HA3	1:H:191:PRO:HG3	2.02	0.41
1:J:103:LEU:HD22	1:J:105:ARG:HG2	2.02	0.41
1:G:92:VAL:CG1	1:G:97:ILE:HD11	2.50	0.41
1:L:121:VAL:HG12	1:L:217:SER:HB2	2.02	0.41
1:D:122:THR:HB	1:D:228:VAL:HG22	2.02	0.41
1:F:121:VAL:HG12	1:F:217:SER:HB2	2.01	0.41
1:G:6:THR:H	1:G:9:HIS:CD2	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:121:VAL:HG21	1:J:229:MET:HE2	2.02	0.41
1:D:97:ILE:O	1:E:166:THR:HA	2.21	0.41
1:G:92:VAL:HG22	1:H:197:LYS:HE3	2.01	0.41
1:I:140:LEU:HD13	1:I:228:VAL:HG11	2.03	0.41
1:C:30:LEU:HD12	1:C:197:LYS:O	2.21	0.41
1:C:187:ASP:HA	1:C:188:SER:HA	1.84	0.41
1:D:59:LEU:HD12	1:D:59:LEU:HA	1.94	0.41
1:I:97:ILE:O	1:J:166:THR:HA	2.20	0.41
1:I:121:VAL:HG21	1:I:229:MET:HE2	2.03	0.41
1:K:6:THR:H	1:K:9:HIS:CD2	2.21	0.41
1:C:92:VAL:CG1	1:C:97:ILE:HD11	2.51	0.41
1:D:100:VAL:HA	1:D:101:PRO:HD3	1.98	0.41
1:D:100:VAL:CG2	1:D:218:ASN:HB2	2.51	0.41
1:G:59:LEU:HD12	1:G:59:LEU:HA	1.92	0.41
1:I:112:TRP:CE3	1:I:183:ASN:HB3	2.56	0.41
1:J:100:VAL:CG2	1:J:218:ASN:HB2	2.50	0.41
1:L:248:MET:CE	1:L:266:MET:HB2	2.51	0.41
1:G:90:VAL:HA	1:H:87:GLU:HA	2.03	0.40
1:K:100:VAL:HA	1:K:101:PRO:HD3	1.98	0.40
1:L:150:GLU:OE2	1:L:188:SER:HB2	2.21	0.40
1:A:216:ILE:HG21	1:A:229:MET:HE2	2.03	0.40
1:F:51:LEU:HD23	1:F:59:LEU:HD22	2.02	0.40
1:H:124:LEU:HB3	1:H:230:VAL:HA	2.03	0.40
1:B:197:LYS:H	1:B:197:LYS:HG2	1.61	0.40
1:C:1:MET:O	1:C:159:ASP:HB2	2.21	0.40
1:L:92:VAL:CG1	1:L:97:ILE:HD11	2.51	0.40
1:B:135:THR:HA	1:B:139:LYS:HE3	2.04	0.40
1:E:100:VAL:HG22	1:E:218:ASN:HB2	2.04	0.40
1:E:124:LEU:HB3	1:E:230:VAL:HA	2.04	0.40
1:J:113:PRO:HB3	1:J:140:LEU:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	265/309 (86%)	260 (98%)	5 (2%)	0	100	100
1	B	266/309 (86%)	261 (98%)	5 (2%)	0	100	100
1	C	266/309 (86%)	258 (97%)	8 (3%)	0	100	100
1	D	266/309 (86%)	257 (97%)	9 (3%)	0	100	100
1	E	265/309 (86%)	260 (98%)	5 (2%)	0	100	100
1	F	265/309 (86%)	259 (98%)	6 (2%)	0	100	100
1	G	264/309 (85%)	257 (97%)	7 (3%)	0	100	100
1	H	265/309 (86%)	259 (98%)	6 (2%)	0	100	100
1	I	265/309 (86%)	261 (98%)	4 (2%)	0	100	100
1	J	265/309 (86%)	260 (98%)	5 (2%)	0	100	100
1	K	265/309 (86%)	256 (97%)	9 (3%)	0	100	100
1	L	271/309 (88%)	266 (98%)	5 (2%)	0	100	100
All	All	3188/3708 (86%)	3114 (98%)	74 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	219/250 (88%)	213 (97%)	6 (3%)	44	78
1	B	220/250 (88%)	216 (98%)	4 (2%)	59	86
1	C	220/250 (88%)	215 (98%)	5 (2%)	50	82
1	D	220/250 (88%)	213 (97%)	7 (3%)	39	73
1	E	220/250 (88%)	215 (98%)	5 (2%)	50	82
1	F	220/250 (88%)	213 (97%)	7 (3%)	39	73
1	G	219/250 (88%)	210 (96%)	9 (4%)	30	64

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	219/250 (88%)	214 (98%)	5 (2%)	50	82
1	I	219/250 (88%)	213 (97%)	6 (3%)	44	78
1	J	220/250 (88%)	216 (98%)	4 (2%)	59	86
1	K	219/250 (88%)	212 (97%)	7 (3%)	39	73
1	L	222/250 (89%)	211 (95%)	11 (5%)	24	56
All	All	2637/3000 (88%)	2561 (97%)	76 (3%)	42	76

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	59	LEU
1	A	100	VAL
1	A	121	VAL
1	A	151	VAL
1	A	189	VAL
1	B	38	THR
1	B	100	VAL
1	B	151	VAL
1	B	197	LYS
1	C	38	THR
1	C	59	LEU
1	C	100	VAL
1	C	151	VAL
1	C	197	LYS
1	D	38	THR
1	D	59	LEU
1	D	100	VAL
1	D	141	ARG
1	D	151	VAL
1	D	188	SER
1	D	197	LYS
1	E	38	THR
1	E	100	VAL
1	E	142	PRO
1	E	151	VAL
1	E	197	LYS
1	F	38	THR
1	F	100	VAL
1	F	121	VAL

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Mol	Chain	Res	Type
1	F	151	VAL
1	F	189	VAL
1	F	192	LEU
1	F	197	LYS
1	G	38	THR
1	G	59	LEU
1	G	100	VAL
1	G	121	VAL
1	G	142	PRO
1	G	151	VAL
1	G	189	VAL
1	G	197	LYS
1	G	273	ARG
1	H	100	VAL
1	H	121	VAL
1	H	136	LEU
1	H	151	VAL
1	H	197	LYS
1	I	38	THR
1	I	54	LYS
1	I	65	ARG
1	I	100	VAL
1	I	151	VAL
1	I	197	LYS
1	J	38	THR
1	J	100	VAL
1	J	121	VAL
1	J	151	VAL
1	K	38	THR
1	K	92	VAL
1	K	94	ASP
1	K	100	VAL
1	K	151	VAL
1	K	189	VAL
1	K	192	LEU
1	L	38	THR
1	L	59	LEU
1	L	94	ASP
1	L	100	VAL
1	L	121	VAL
1	L	125	MET
1	L	151	VAL

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Mol	Chain	Res	Type
1	L	188	SER
1	L	189	VAL
1	L	192	LEU
1	L	197	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	37	ASN
1	A	183	ASN
1	A	218	ASN
1	B	9	HIS
1	B	37	ASN
1	B	183	ASN
1	B	218	ASN
1	C	9	HIS
1	C	37	ASN
1	C	183	ASN
1	C	218	ASN
1	D	9	HIS
1	D	37	ASN
1	D	183	ASN
1	D	218	ASN
1	E	9	HIS
1	E	37	ASN
1	E	183	ASN
1	E	218	ASN
1	F	9	HIS
1	F	37	ASN
1	F	183	ASN
1	F	218	ASN
1	G	9	HIS
1	G	37	ASN
1	G	183	ASN
1	G	218	ASN
1	H	9	HIS
1	H	37	ASN
1	H	183	ASN
1	H	218	ASN
1	I	9	HIS
1	I	37	ASN

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Mol	Chain	Res	Type
1	I	218	ASN
1	J	9	HIS
1	J	37	ASN
1	J	183	ASN
1	J	218	ASN
1	K	9	HIS
1	K	37	ASN
1	K	183	ASN
1	K	218	ASN
1	L	9	HIS
1	L	37	ASN
1	L	183	ASN
1	L	218	ASN
1	L	232	ASN

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](#)

Of 24 ligands modelled in this entry, 12 are monoatomic - leaving 12 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	ADP	C	401	2	24,29,29	0.69	0	29,45,45	0.73	1 (3%)
3	ADP	B	401	2	24,29,29	0.72	0	29,45,45	0.70	1 (3%)
3	ADP	H	401	2	24,29,29	0.70	0	29,45,45	0.76	1 (3%)
3	ADP	L	401	2	24,29,29	0.73	0	29,45,45	0.70	1 (3%)
3	ADP	E	401	2	24,29,29	0.73	0	29,45,45	0.71	1 (3%)
3	ADP	I	401	2	24,29,29	0.65	0	29,45,45	0.72	1 (3%)
3	ADP	D	401	2	24,29,29	0.68	0	29,45,45	0.70	1 (3%)
3	ADP	G	401	2	24,29,29	0.71	0	29,45,45	0.71	1 (3%)
3	ADP	A	401	2	24,29,29	0.77	0	29,45,45	0.72	1 (3%)
3	ADP	J	401	2	24,29,29	0.70	0	29,45,45	0.72	1 (3%)
3	ADP	F	401	2	24,29,29	0.69	0	29,45,45	0.70	1 (3%)
3	ADP	K	401	2	24,29,29	0.72	1 (4%)	29,45,45	0.68	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	ADP	C	401	2	-	4/12/32/32	0/3/3/3
3	ADP	B	401	2	-	4/12/32/32	0/3/3/3
3	ADP	H	401	2	-	4/12/32/32	0/3/3/3
3	ADP	L	401	2	-	4/12/32/32	0/3/3/3
3	ADP	E	401	2	-	4/12/32/32	0/3/3/3
3	ADP	I	401	2	-	3/12/32/32	0/3/3/3
3	ADP	D	401	2	-	4/12/32/32	0/3/3/3
3	ADP	G	401	2	-	4/12/32/32	0/3/3/3
3	ADP	A	401	2	-	4/12/32/32	0/3/3/3
3	ADP	J	401	2	-	4/12/32/32	0/3/3/3
3	ADP	F	401	2	-	4/12/32/32	0/3/3/3
3	ADP	K	401	2	-	4/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	401	ADP	C8-N7	-2.01	1.31	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	401	ADP	C5-C6-N6	2.47	124.10	120.35
3	I	401	ADP	C5-C6-N6	2.43	124.05	120.35
3	A	401	ADP	C5-C6-N6	2.43	124.05	120.35
3	C	401	ADP	C5-C6-N6	2.43	124.04	120.35
3	H	401	ADP	C5-C6-N6	2.38	123.97	120.35
3	D	401	ADP	C5-C6-N6	2.37	123.95	120.35
3	E	401	ADP	C5-C6-N6	2.35	123.93	120.35
3	B	401	ADP	C5-C6-N6	2.29	123.83	120.35
3	L	401	ADP	C5-C6-N6	2.28	123.81	120.35
3	K	401	ADP	C5-C6-N6	2.28	123.81	120.35
3	G	401	ADP	C5-C6-N6	2.27	123.80	120.35
3	F	401	ADP	C5-C6-N6	2.26	123.78	120.35

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ADP	PA-O3A-PB-O2B
3	A	401	ADP	C5'-O5'-PA-O3A
3	B	401	ADP	PA-O3A-PB-O2B
3	B	401	ADP	C5'-O5'-PA-O3A
3	C	401	ADP	PA-O3A-PB-O2B
3	D	401	ADP	PA-O3A-PB-O2B
3	D	401	ADP	C5'-O5'-PA-O3A
3	E	401	ADP	PA-O3A-PB-O2B
3	E	401	ADP	C5'-O5'-PA-O3A
3	F	401	ADP	PA-O3A-PB-O2B
3	F	401	ADP	C5'-O5'-PA-O3A
3	G	401	ADP	PA-O3A-PB-O2B
3	G	401	ADP	C5'-O5'-PA-O3A
3	H	401	ADP	PA-O3A-PB-O2B
3	I	401	ADP	PA-O3A-PB-O2B
3	I	401	ADP	PA-O3A-PB-O3B
3	J	401	ADP	PA-O3A-PB-O2B
3	J	401	ADP	C5'-O5'-PA-O3A
3	K	401	ADP	PA-O3A-PB-O2B
3	K	401	ADP	C5'-O5'-PA-O3A
3	L	401	ADP	PA-O3A-PB-O2B
3	L	401	ADP	C5'-O5'-PA-O3A
3	C	401	ADP	C5'-O5'-PA-O3A
3	H	401	ADP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
3	A	401	ADP	C5'-O5'-PA-O2A
3	B	401	ADP	C5'-O5'-PA-O2A
3	E	401	ADP	C5'-O5'-PA-O2A
3	F	401	ADP	C5'-O5'-PA-O2A
3	G	401	ADP	C5'-O5'-PA-O2A
3	H	401	ADP	C5'-O5'-PA-O2A
3	J	401	ADP	C5'-O5'-PA-O2A
3	K	401	ADP	C5'-O5'-PA-O2A
3	L	401	ADP	C5'-O5'-PA-O2A
3	A	401	ADP	PA-O3A-PB-O1B
3	B	401	ADP	PA-O3A-PB-O1B
3	C	401	ADP	PA-O3A-PB-O1B
3	F	401	ADP	PA-O3A-PB-O1B
3	G	401	ADP	PA-O3A-PB-O1B
3	H	401	ADP	PA-O3A-PB-O1B
3	C	401	ADP	C5'-O5'-PA-O2A
3	D	401	ADP	C5'-O5'-PA-O2A
3	D	401	ADP	PA-O3A-PB-O1B
3	E	401	ADP	PA-O3A-PB-O1B
3	I	401	ADP	PA-O3A-PB-O1B
3	J	401	ADP	PA-O3A-PB-O1B
3	K	401	ADP	PA-O3A-PB-O1B
3	L	401	ADP	PA-O3A-PB-O1B

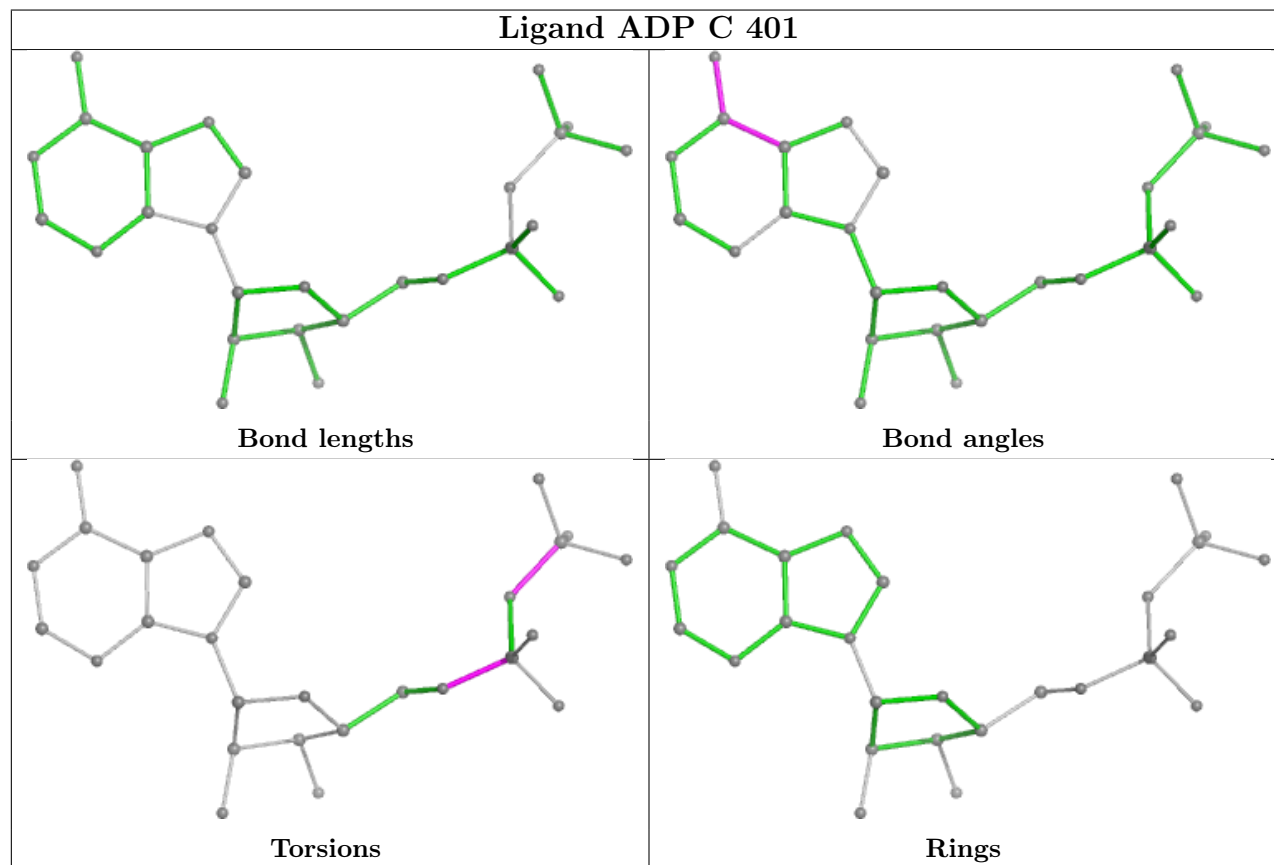
There are no ring outliers.

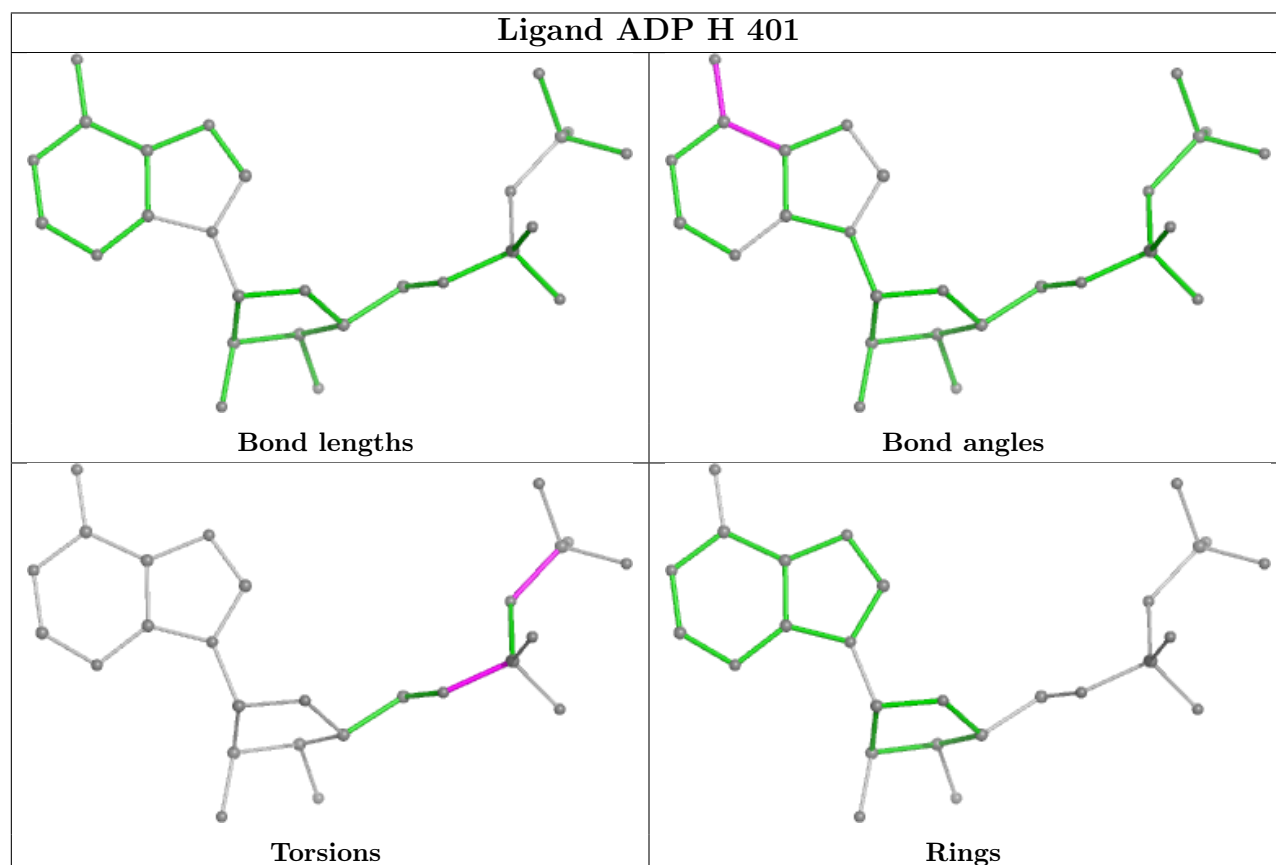
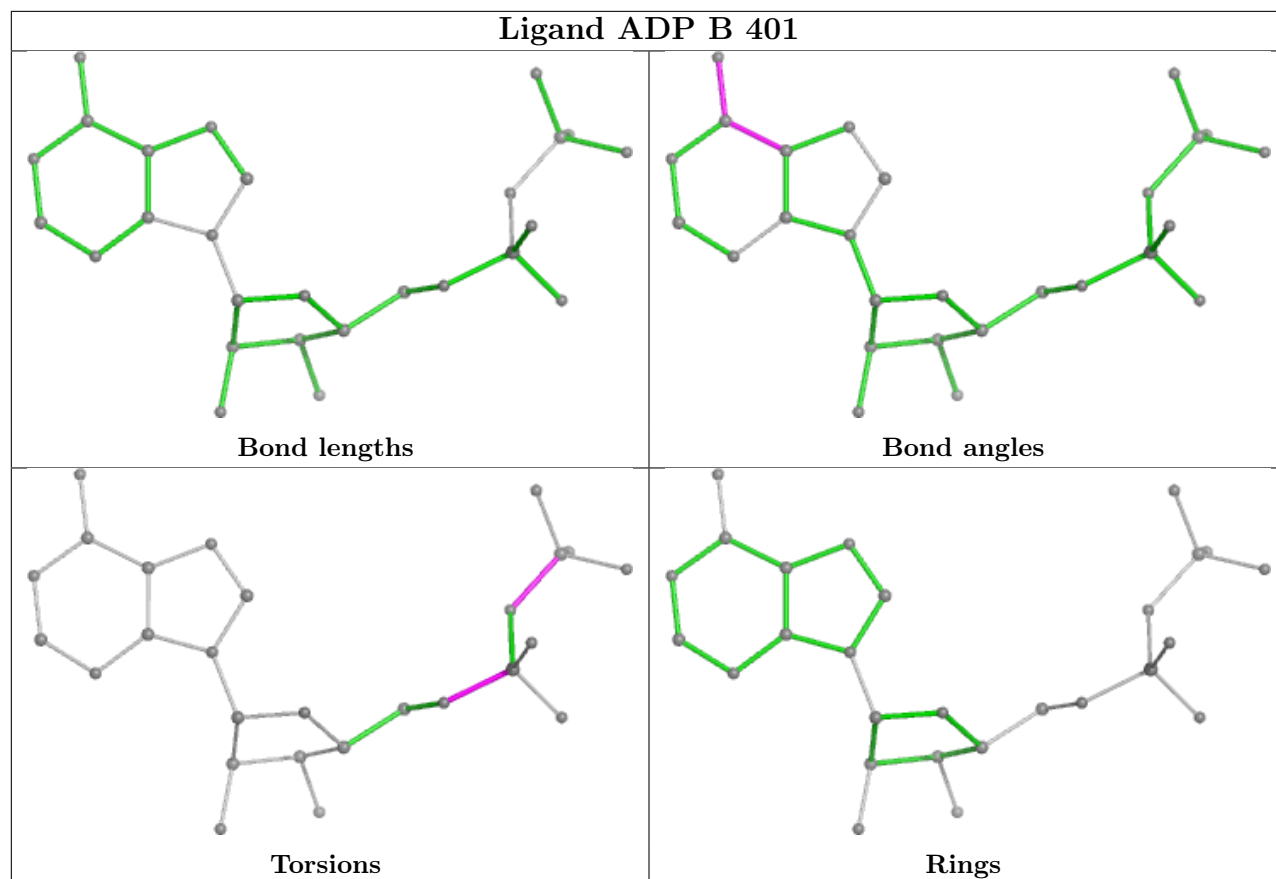
11 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	401	ADP	1	0
3	B	401	ADP	1	0
3	H	401	ADP	1	0
3	L	401	ADP	1	0
3	E	401	ADP	1	0
3	D	401	ADP	1	0
3	G	401	ADP	1	0
3	A	401	ADP	1	0
3	J	401	ADP	1	0
3	F	401	ADP	1	0
3	K	401	ADP	1	0

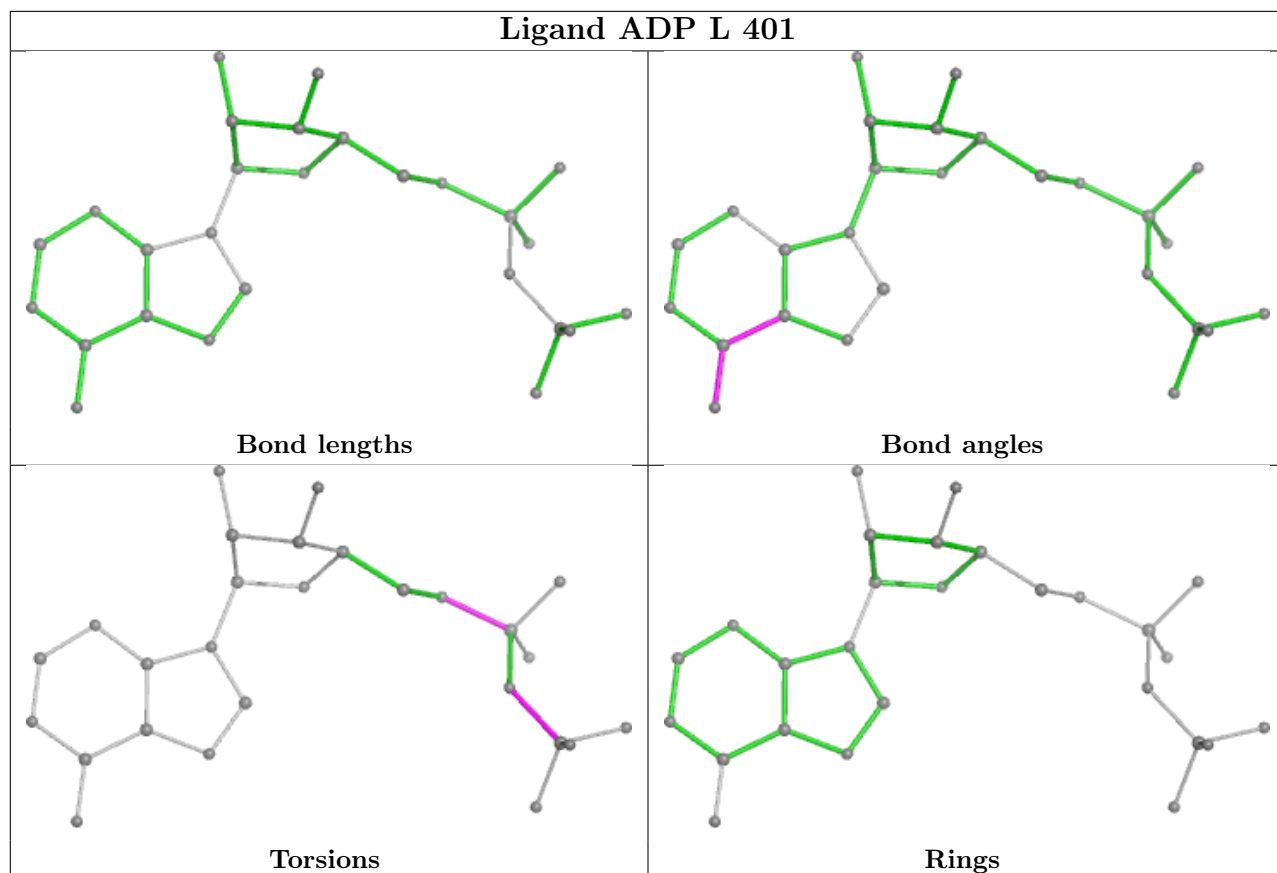
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.

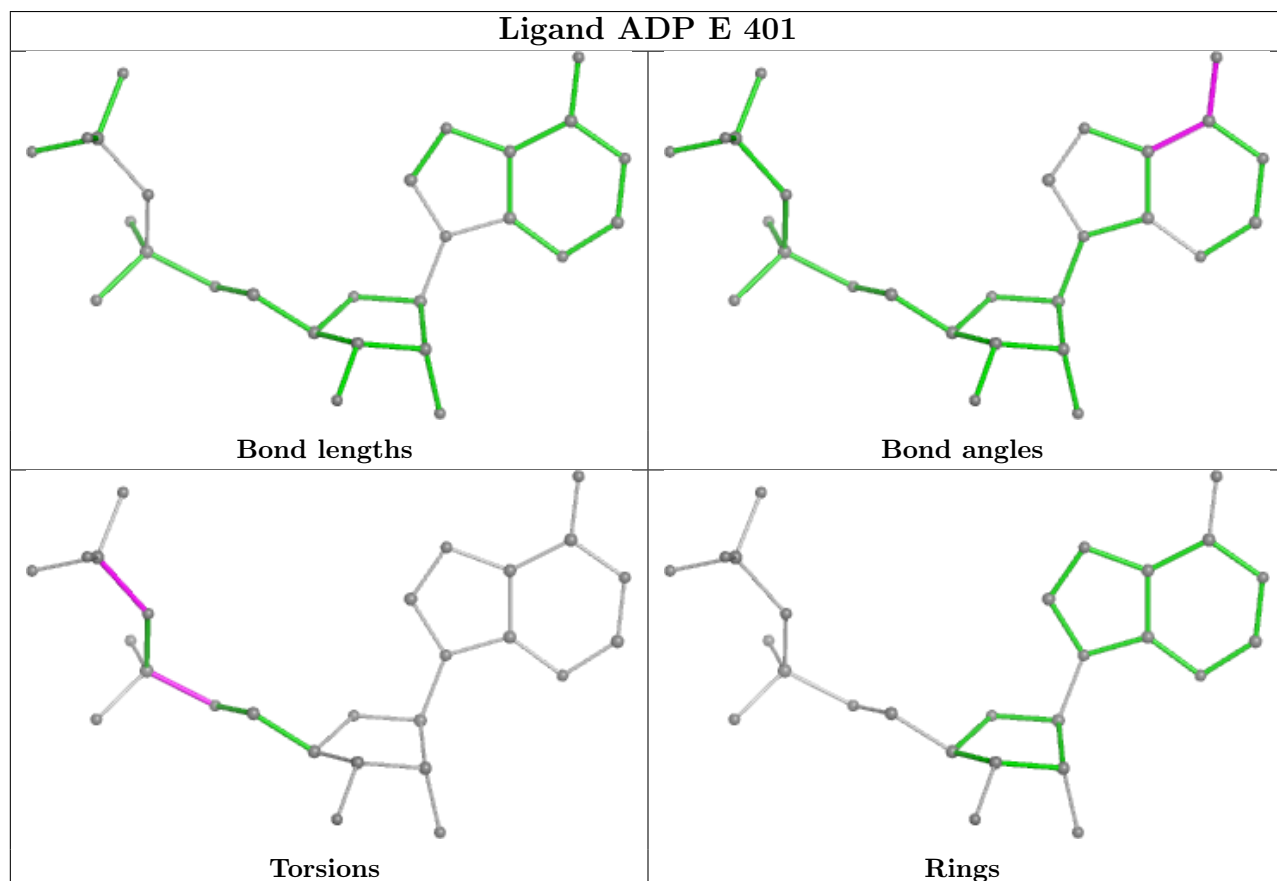


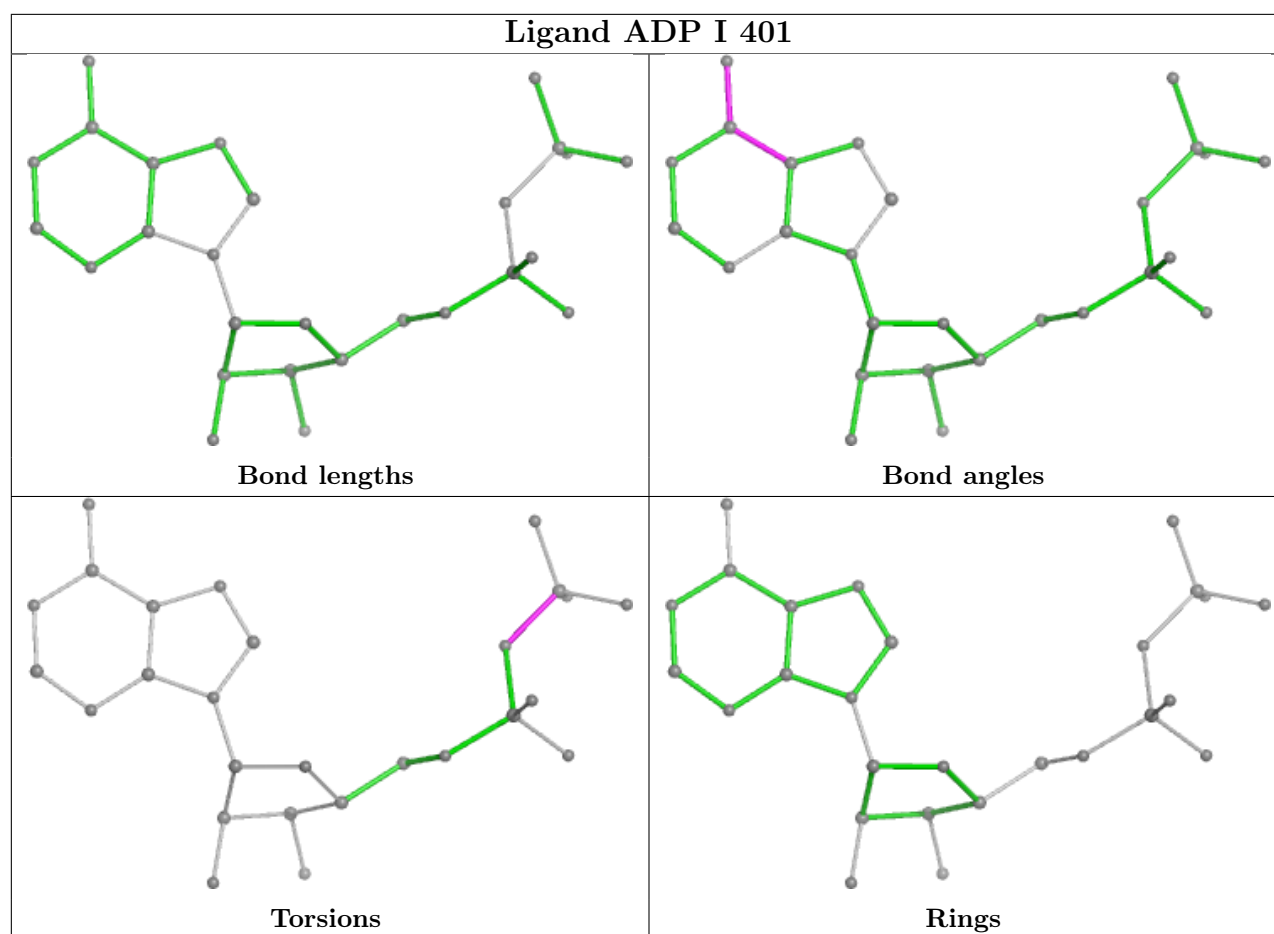


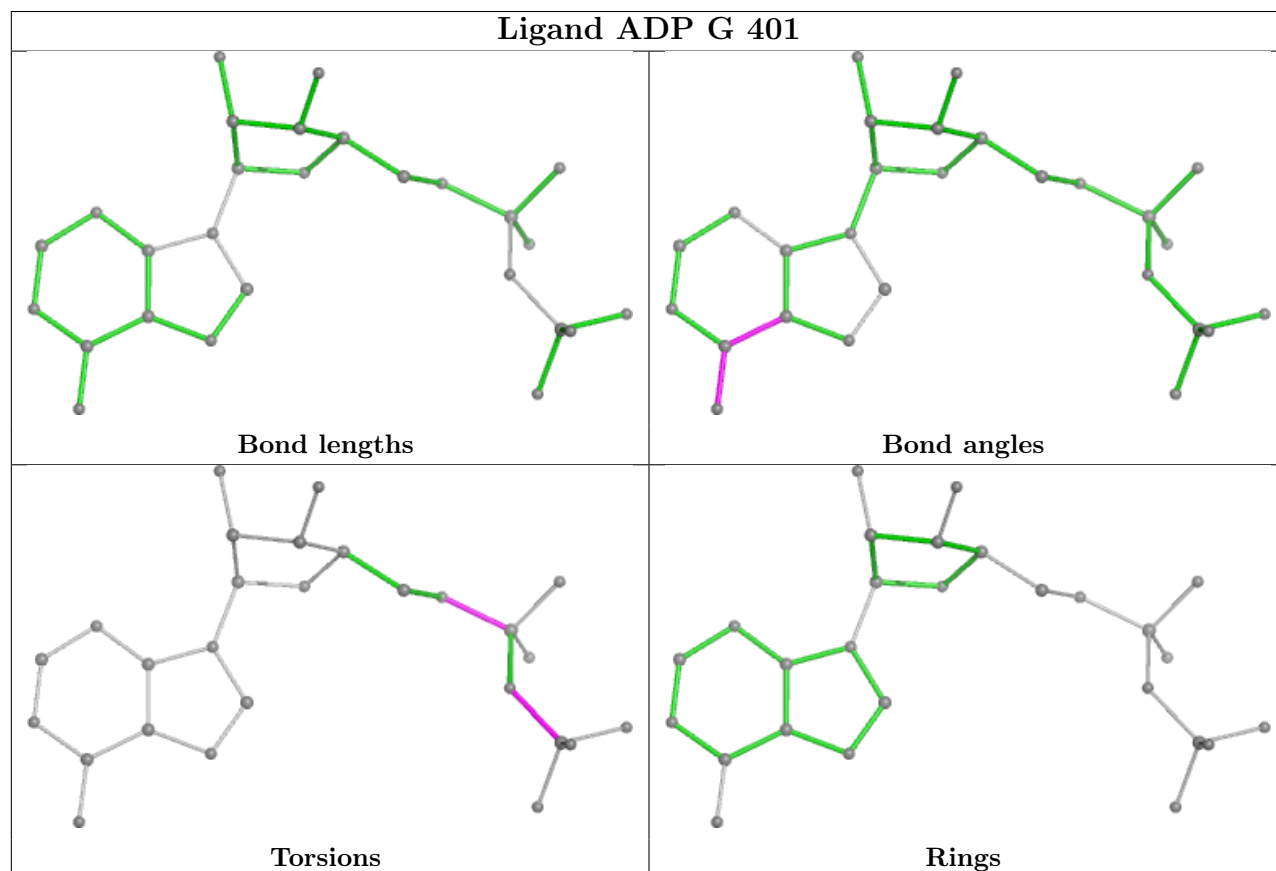
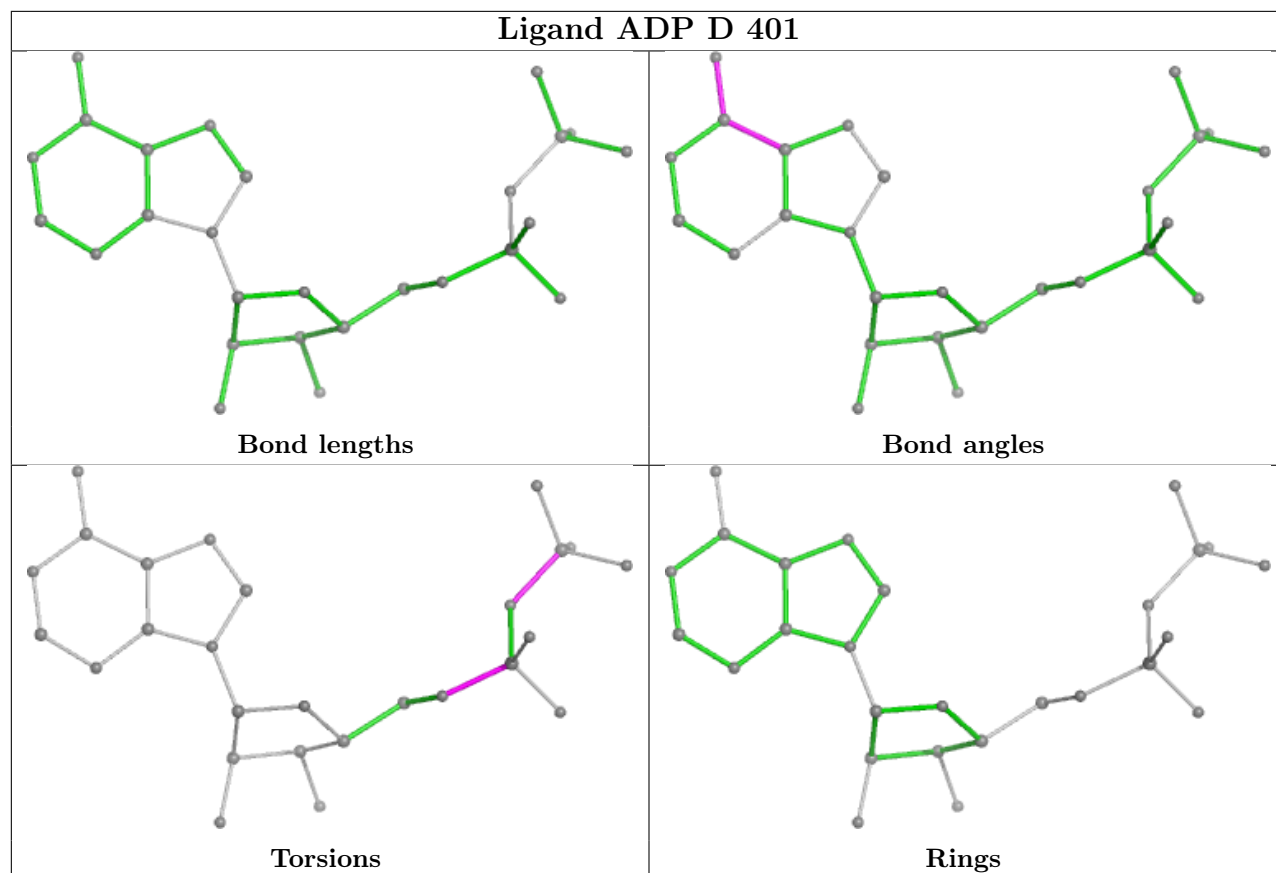
Ligand ADP L 401

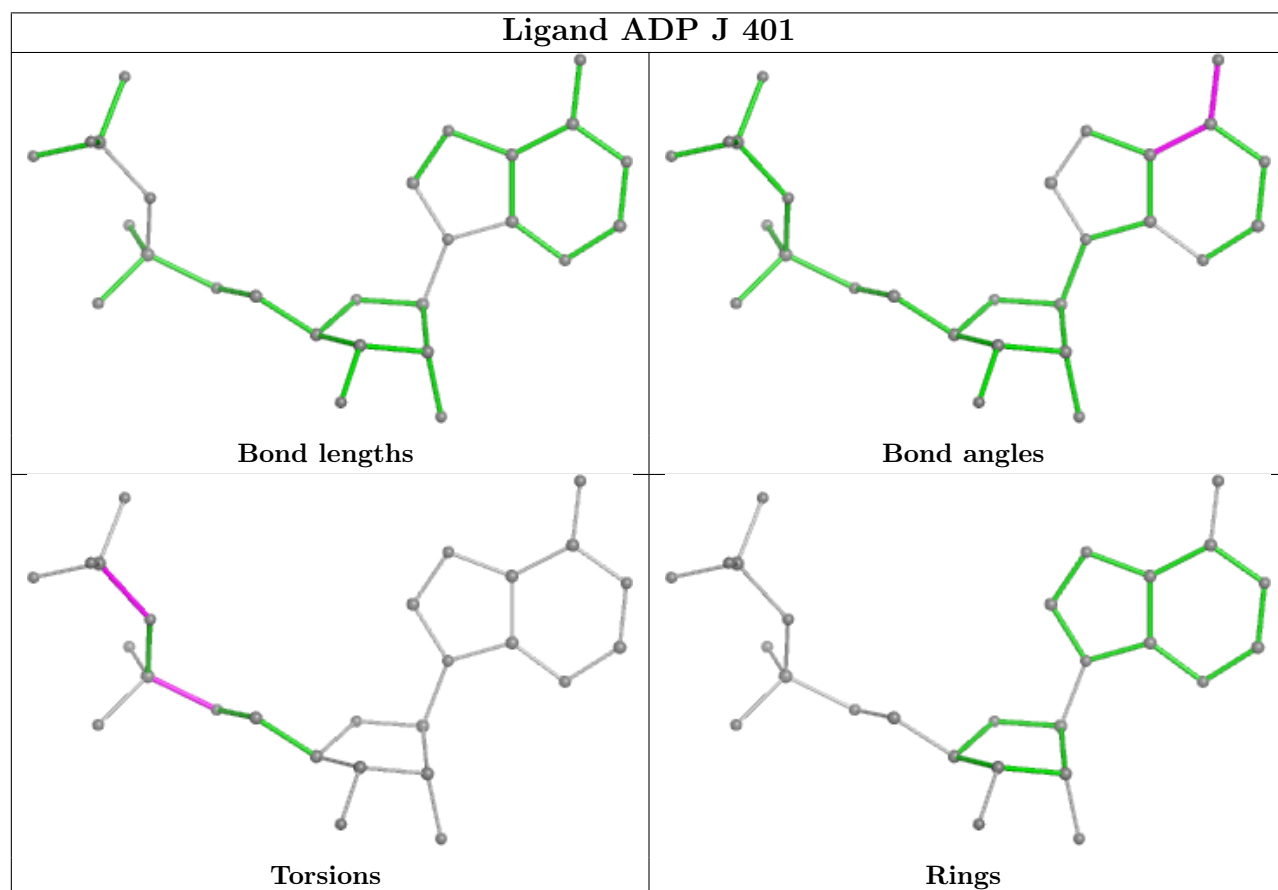
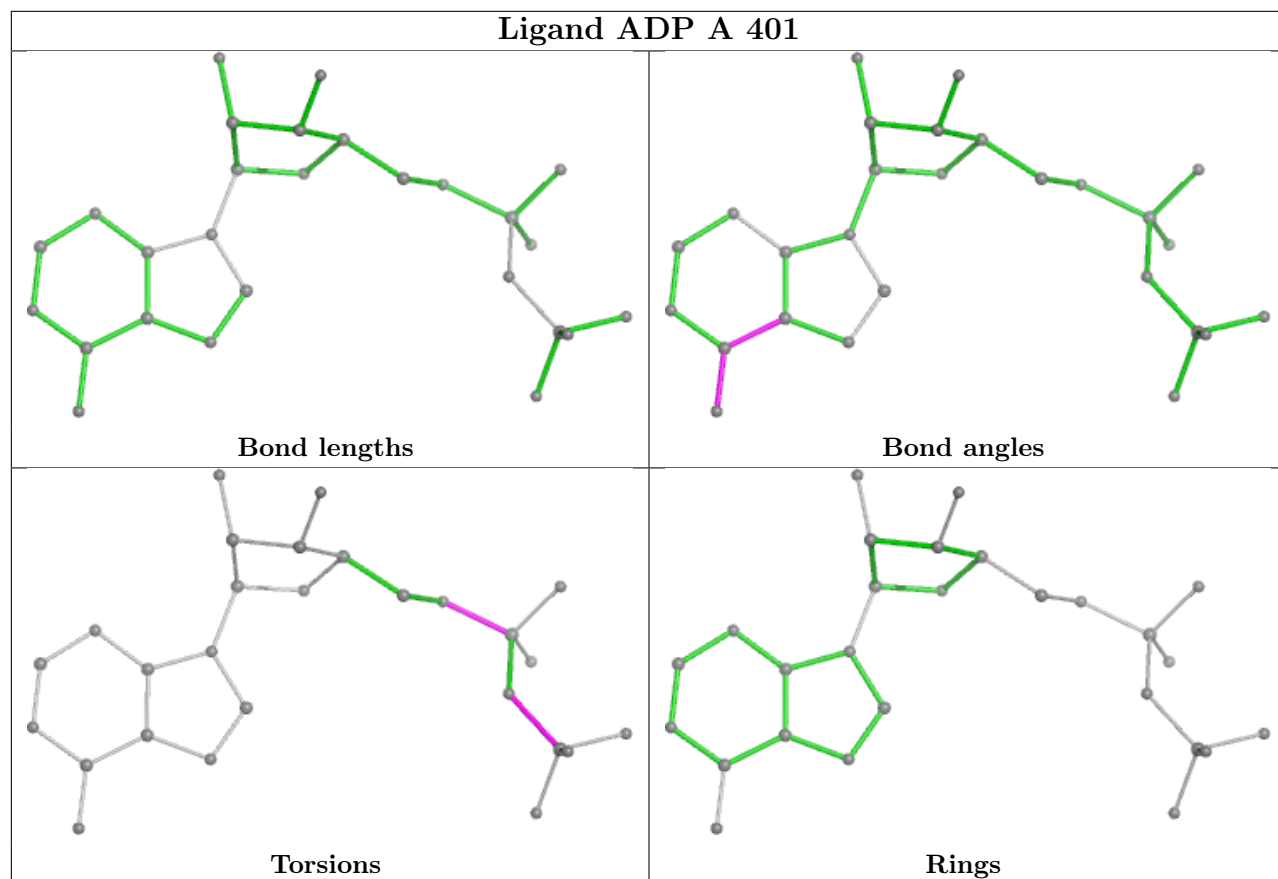


Ligand ADP E 401

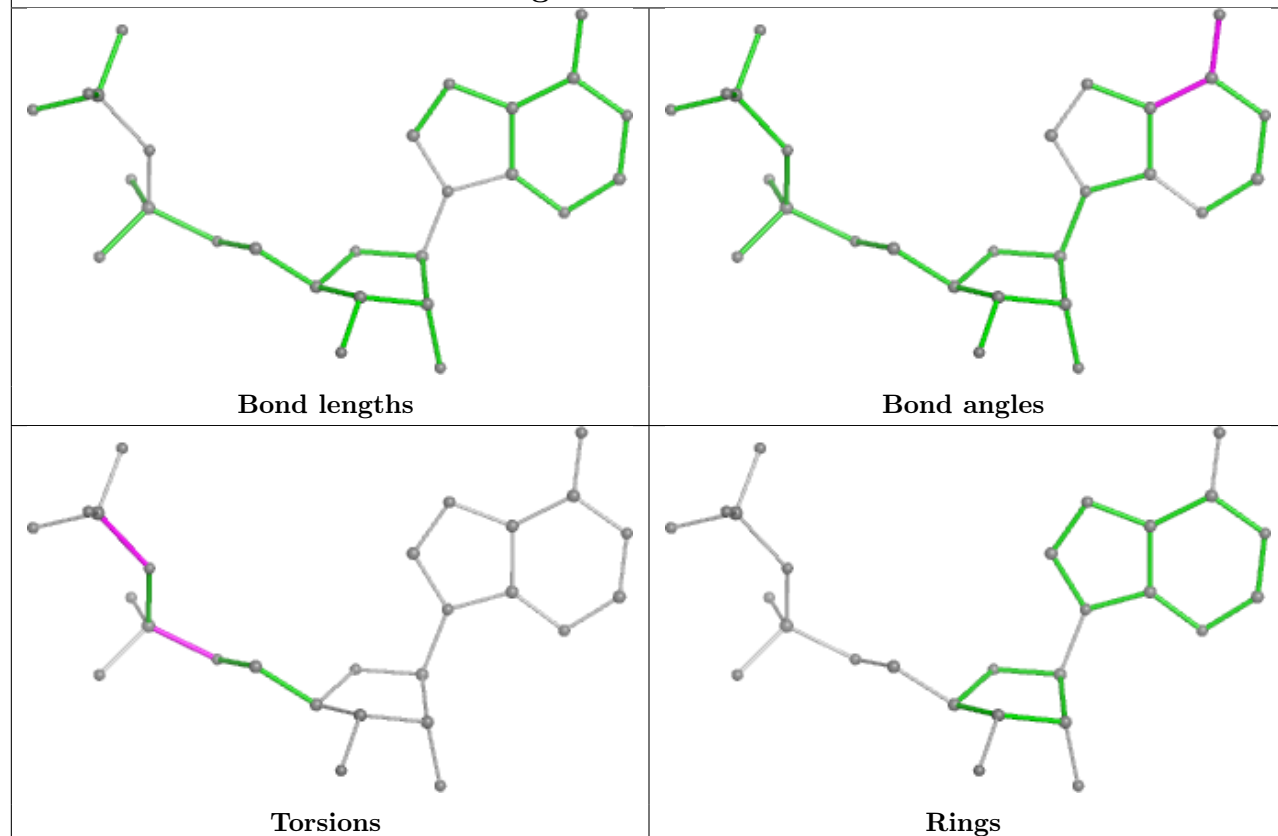




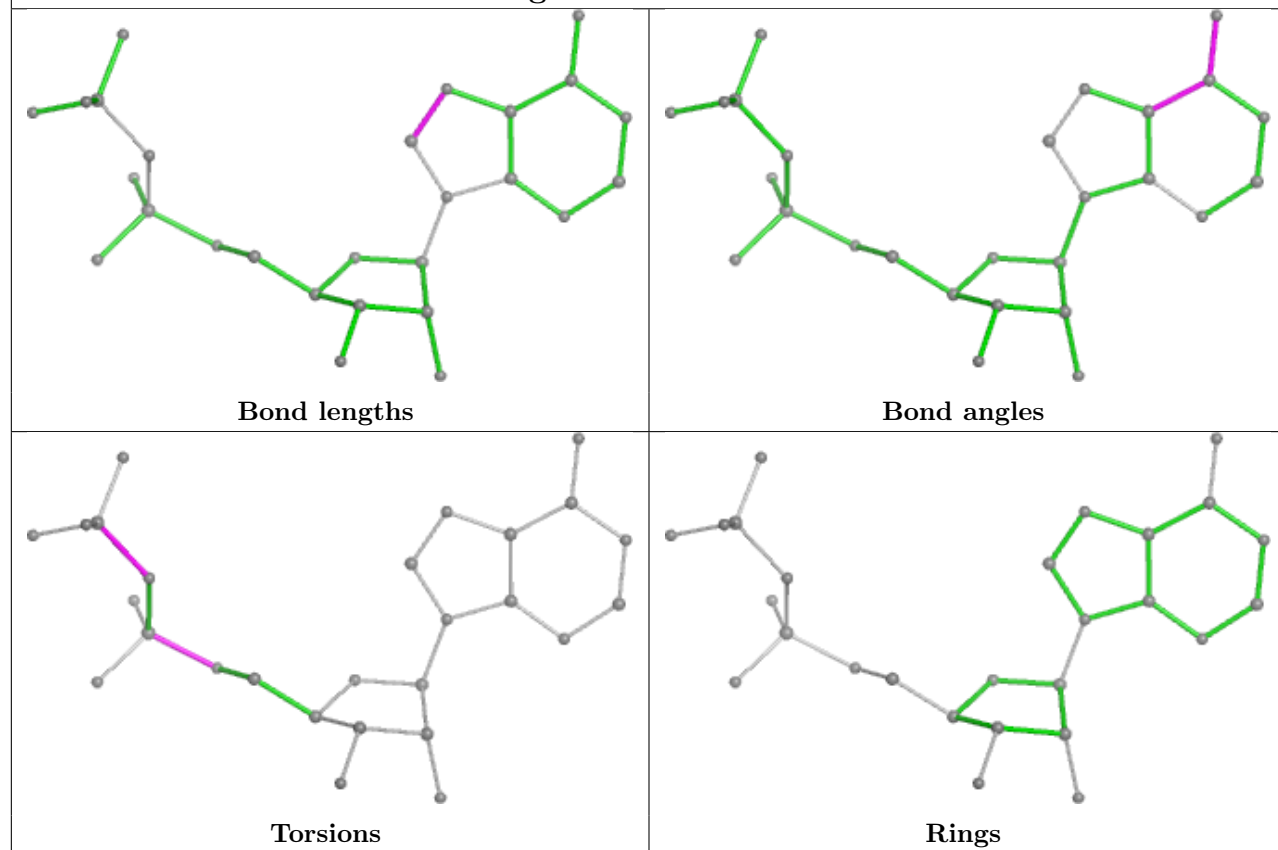




Ligand ADP F 401



Ligand ADP K 401



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	269/309 (87%)	-0.19	6 (2%) 62 52	22, 44, 77, 113	0
1	B	270/309 (87%)	-0.18	7 (2%) 56 46	23, 41, 73, 100	0
1	C	270/309 (87%)	0.25	13 (4%) 30 21	26, 64, 107, 130	0
1	D	270/309 (87%)	0.32	20 (7%) 14 8	26, 73, 112, 158	0
1	E	269/309 (87%)	0.17	12 (4%) 33 23	26, 63, 98, 115	0
1	F	269/309 (87%)	0.06	11 (4%) 37 27	26, 54, 87, 120	0
1	G	268/309 (86%)	0.08	16 (5%) 21 14	26, 52, 93, 137	0
1	H	269/309 (87%)	0.06	9 (3%) 46 36	26, 55, 92, 123	0
1	I	269/309 (87%)	-0.20	7 (2%) 56 46	25, 41, 73, 95	0
1	J	269/309 (87%)	-0.14	8 (2%) 50 40	26, 44, 80, 107	0
1	K	269/309 (87%)	-0.15	7 (2%) 56 46	26, 45, 81, 106	0
1	L	275/309 (88%)	-0.14	7 (2%) 57 47	26, 48, 82, 101	0
All	All	3236/3708 (87%)	-0.01	123 (3%) 40 30	22, 52, 93, 158	0

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	56	GLY	7.8
1	C	1	MET	6.3
1	J	1	MET	6.1
1	F	1	MET	6.1
1	I	84	ALA	5.9
1	D	57	GLY	5.8
1	E	1	MET	5.4
1	F	238	GLU	4.9
1	K	85	ASP	4.8
1	E	85	ASP	4.8
1	G	1	MET	4.7

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Mol	Chain	Res	Type	RSRZ
1	C	104	ALA	4.6
1	B	1	MET	4.5
1	K	84	ALA	4.1
1	C	261	ASN	4.1
1	C	84	ALA	4.1
1	G	275	PHE	4.0
1	H	85	ASP	3.8
1	B	238	GLU	3.6
1	L	23	SER	3.6
1	H	84	ALA	3.5
1	G	85	ASP	3.5
1	E	22	ALA	3.5
1	G	261	ASN	3.5
1	F	85	ASP	3.5
1	D	237	ALA	3.4
1	C	238	GLU	3.4
1	B	89	ASP	3.4
1	I	85	ASP	3.4
1	L	85	ASP	3.3
1	B	84	ALA	3.3
1	J	238	GLU	3.3
1	E	199	ALA	3.1
1	B	77	THR	3.0
1	C	237	ALA	3.0
1	D	89	ASP	3.0
1	D	260	GLY	3.0
1	L	1	MET	3.0
1	E	77	THR	2.9
1	D	1	MET	2.9
1	C	199	ALA	2.9
1	D	272	GLY	2.9
1	L	261	ASN	2.9
1	H	271	LYS	2.9
1	I	261	ASN	2.9
1	D	77	THR	2.8
1	I	89	ASP	2.8
1	L	111	ARG	2.8
1	A	84	ALA	2.8
1	D	84	ALA	2.8
1	C	111	ARG	2.8
1	A	54	LYS	2.8
1	D	23	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	239	LYS	2.8
1	H	237	ALA	2.8
1	G	260	GLY	2.8
1	D	85	ASP	2.7
1	K	199	ALA	2.7
1	H	71	LYS	2.7
1	C	259	ASP	2.7
1	J	85	ASP	2.7
1	G	202	ALA	2.7
1	G	111	ARG	2.7
1	A	85	ASP	2.7
1	D	53	SER	2.7
1	K	89	ASP	2.6
1	L	18	ASP	2.6
1	G	238	GLU	2.6
1	E	89	ASP	2.6
1	I	22	ALA	2.6
1	F	104	ALA	2.6
1	I	77	THR	2.6
1	E	260	GLY	2.6
1	J	86	VAL	2.6
1	A	261	ASN	2.6
1	G	274	ILE	2.5
1	D	72	ASP	2.5
1	F	89	ASP	2.5
1	G	270	ASN	2.5
1	B	85	ASP	2.5
1	E	238	GLU	2.5
1	C	141	ARG	2.5
1	I	238	GLU	2.5
1	G	89	ASP	2.5
1	D	104	ALA	2.5
1	J	271	LYS	2.4
1	F	236	ASP	2.4
1	F	53	SER	2.4
1	J	22	ALA	2.4
1	D	55	GLN	2.4
1	F	111	ARG	2.4
1	D	58	SER	2.4
1	L	84	ALA	2.4
1	D	115	GLU	2.3
1	G	273	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	89	ASP	2.3
1	E	141	ARG	2.3
1	H	22	ALA	2.3
1	D	11	ASP	2.3
1	H	72	ASP	2.3
1	B	237	ALA	2.2
1	H	276	ASN	2.2
1	C	23	SER	2.2
1	E	127	ALA	2.2
1	C	115	GLU	2.2
1	K	237	ALA	2.2
1	J	111	ARG	2.2
1	J	77	THR	2.1
1	G	241	GLU	2.1
1	E	72	ASP	2.1
1	K	71	LYS	2.1
1	K	200	ALA	2.1
1	F	204	GLY	2.1
1	D	111	ARG	2.1
1	E	111	ARG	2.1
1	G	141	ARG	2.1
1	A	53	SER	2.1
1	G	116	GLY	2.1
1	F	55	GLN	2.0
1	G	18	ASP	2.0
1	D	159	ASP	2.0
1	F	200	ALA	2.0
1	H	111	ARG	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,

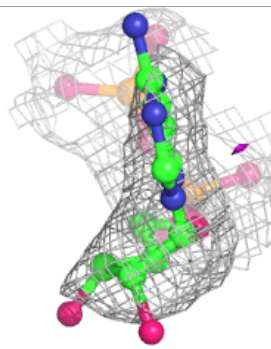
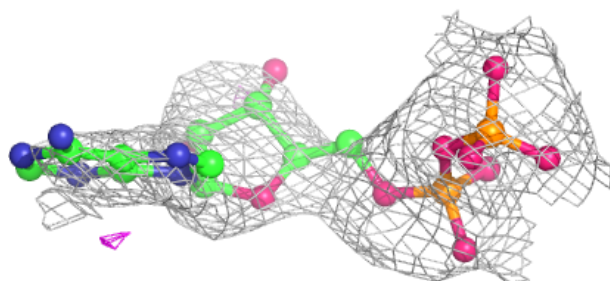
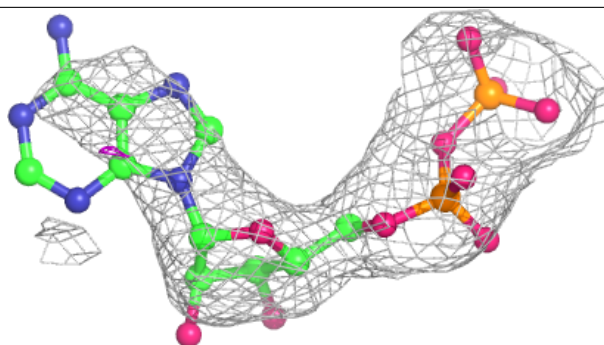
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ADP	C	401	27/27	0.82	0.28	70,100,119,121	0
3	ADP	H	401	27/27	0.82	0.41	77,127,140,143	0
3	ADP	D	401	27/27	0.83	0.27	82,120,135,138	0
3	ADP	F	401	27/27	0.85	0.25	62,91,106,108	0
2	CA	F	400	1/1	0.88	0.25	70,70,70,70	0
3	ADP	E	401	27/27	0.88	0.27	62,87,96,98	0
3	ADP	G	401	27/27	0.89	0.25	63,88,99,102	0
3	ADP	J	401	27/27	0.89	0.23	48,59,67,70	0
3	ADP	K	401	27/27	0.90	0.27	49,73,79,80	0
3	ADP	B	401	27/27	0.91	0.21	24,57,64,65	0
3	ADP	L	401	27/27	0.91	0.23	39,71,76,77	0
3	ADP	I	401	27/27	0.92	0.24	33,63,77,79	0
2	CA	I	400	1/1	0.92	0.13	49,49,49,49	0
2	CA	J	400	1/1	0.92	0.18	48,48,48,48	0
2	CA	B	400	1/1	0.92	0.19	55,55,55,55	0
3	ADP	A	401	27/27	0.93	0.19	33,57,62,63	0
2	CA	C	400	1/1	0.93	0.12	68,68,68,68	0
2	CA	A	400	1/1	0.93	0.25	65,65,65,65	0
2	CA	E	400	1/1	0.94	0.16	76,76,76,76	0
2	CA	H	400	1/1	0.96	0.15	68,68,68,68	0
2	CA	G	400	1/1	0.97	0.16	80,80,80,80	0
2	CA	D	400	1/1	0.97	0.17	70,70,70,70	0
2	CA	K	400	1/1	0.97	0.24	61,61,61,61	0
2	CA	L	400	1/1	0.98	0.15	50,50,50,50	0

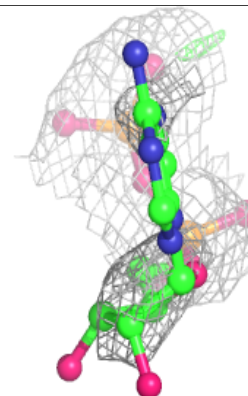
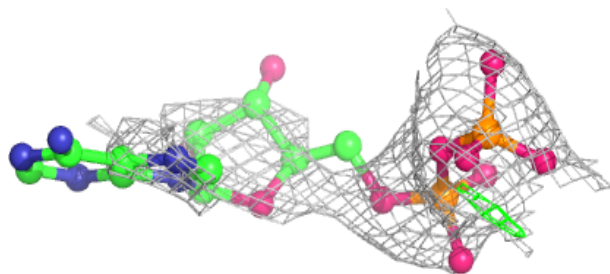
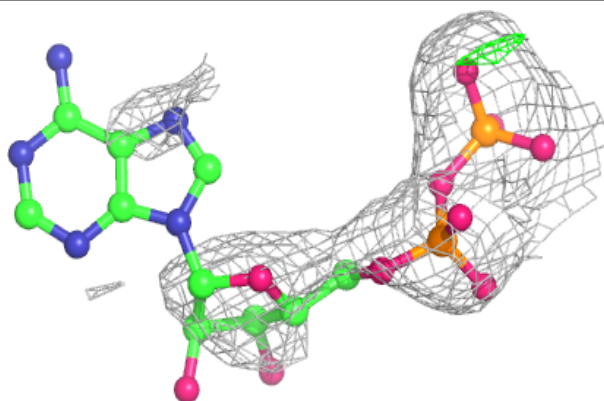
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP 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)

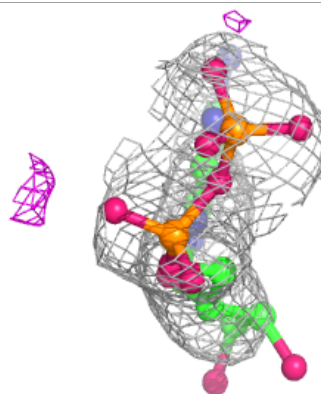
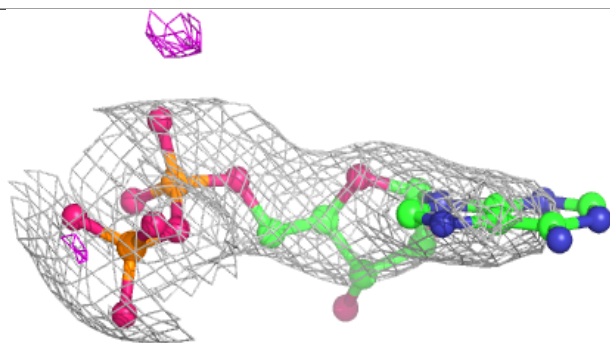
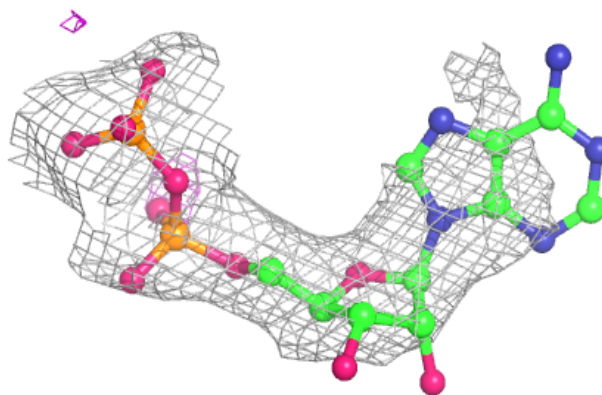
**Electron density around ADP H 401:**

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

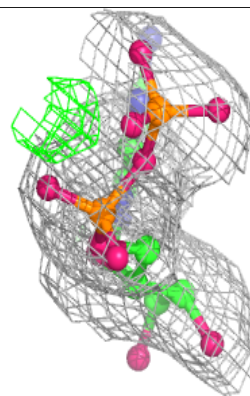
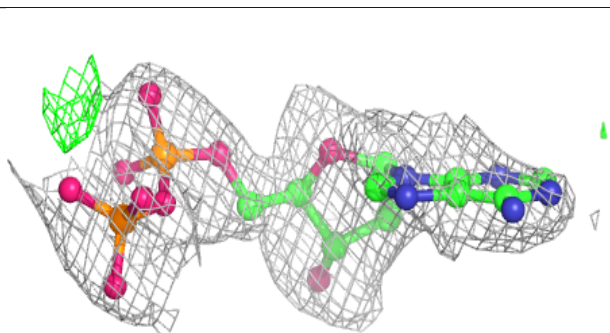
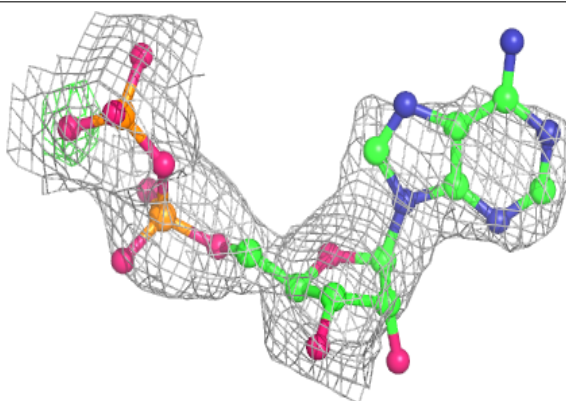


Electron density around ADP D 401:

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

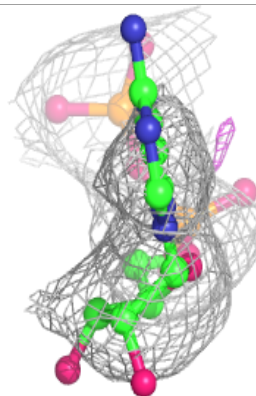
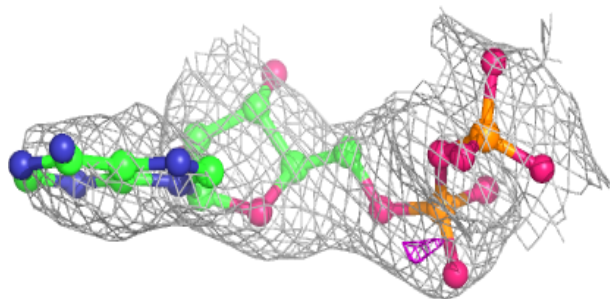
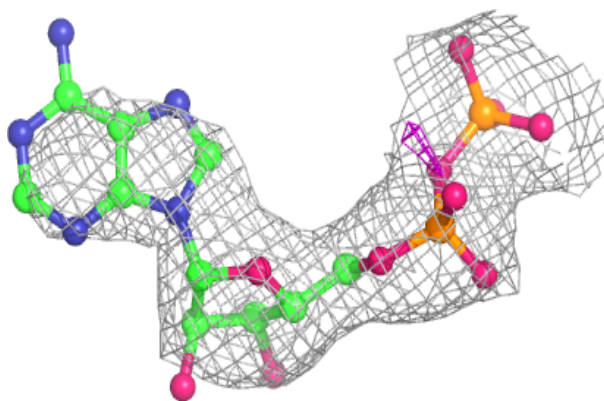
**Electron density around ADP F 401:**

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

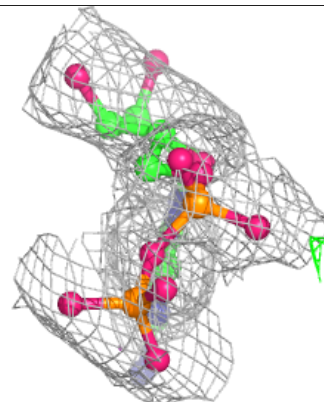
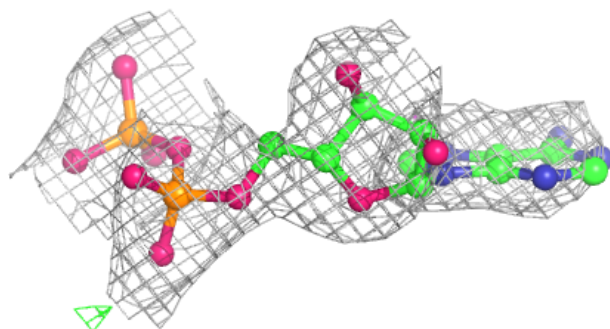
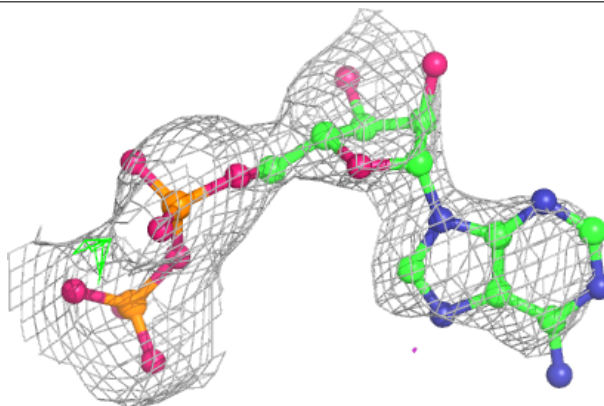


Electron density around ADP E 401:

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

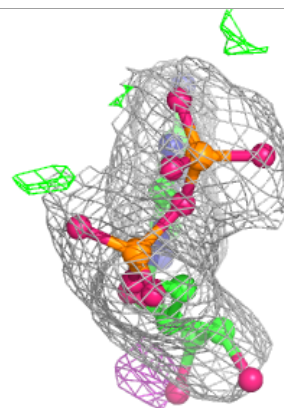
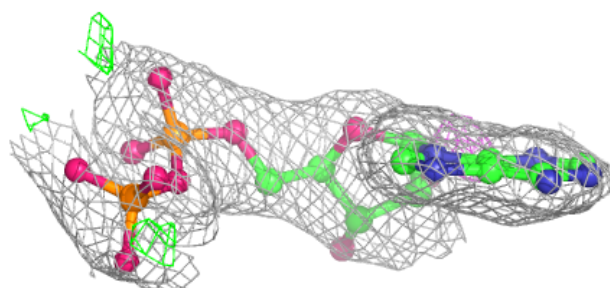
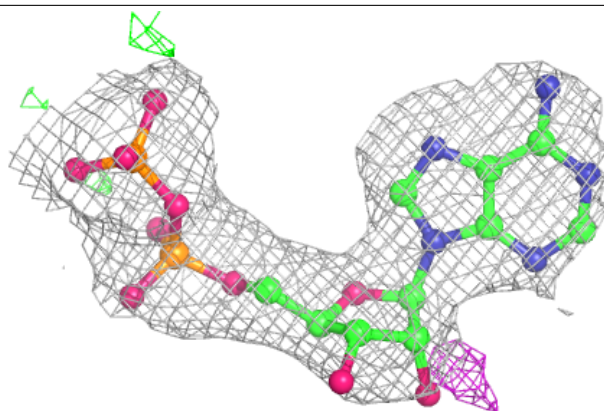
**Electron density around ADP G 401:**

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

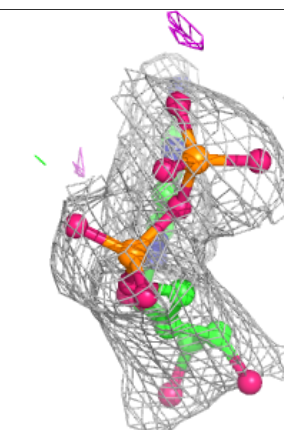
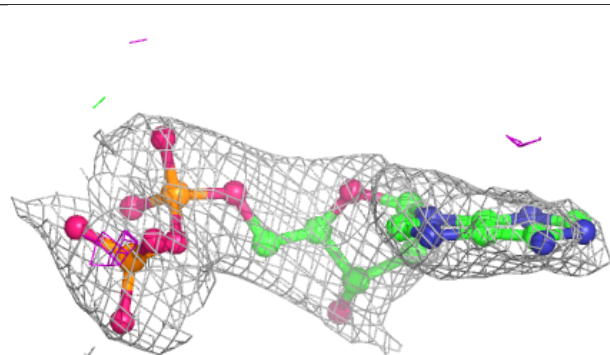
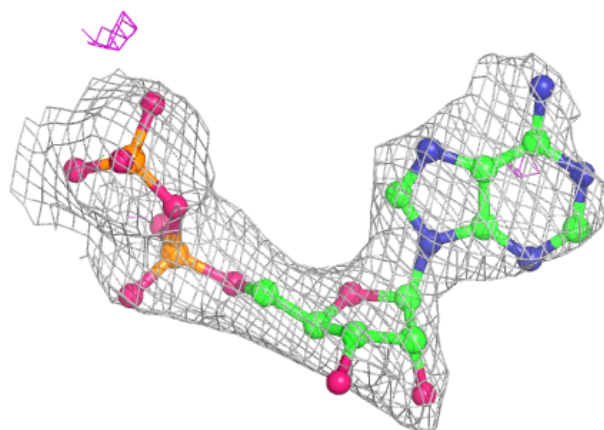


Electron density around ADP J 401:

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

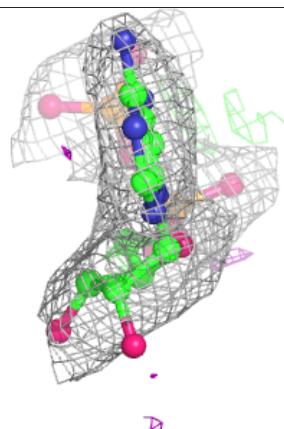
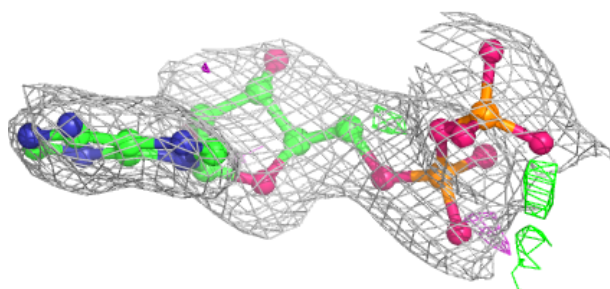
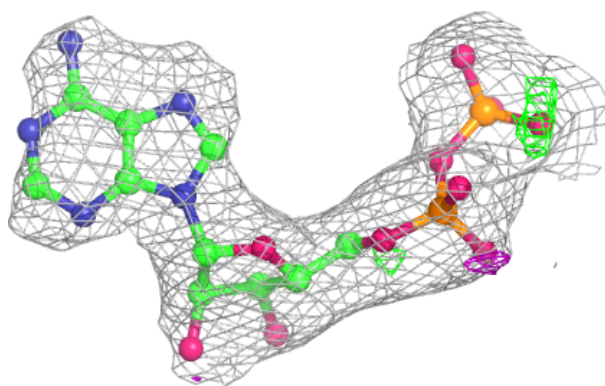
**Electron density around ADP K 401:**

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

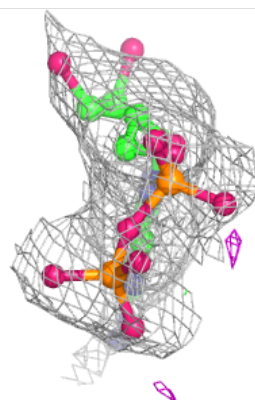
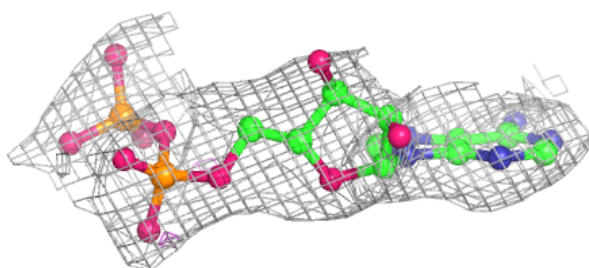
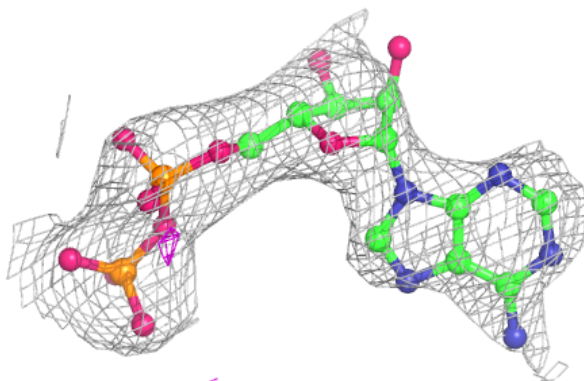


Electron density around ADP B 401:

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

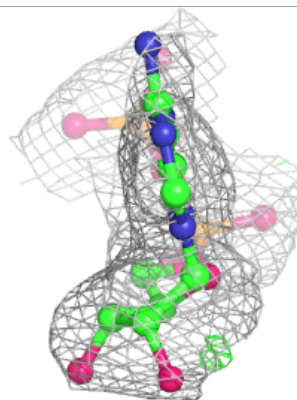
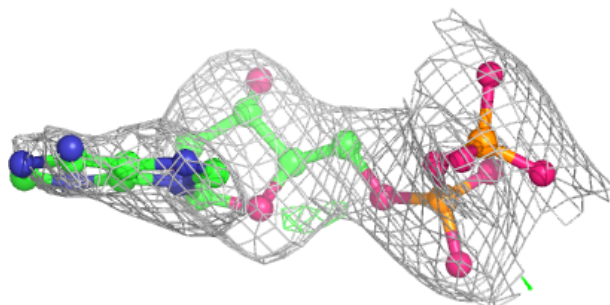
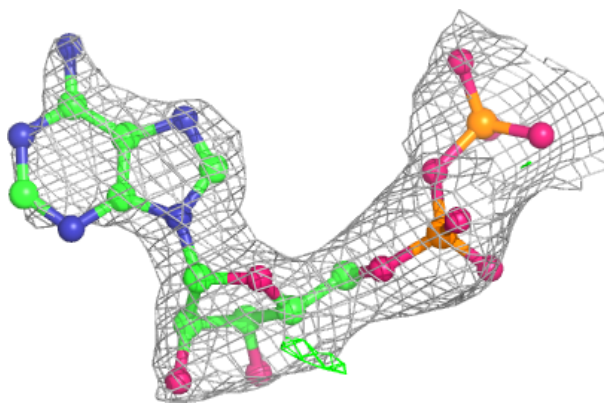
**Electron density around ADP L 401:**

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

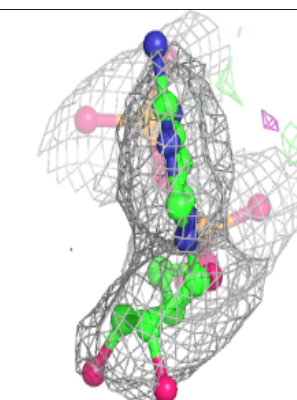
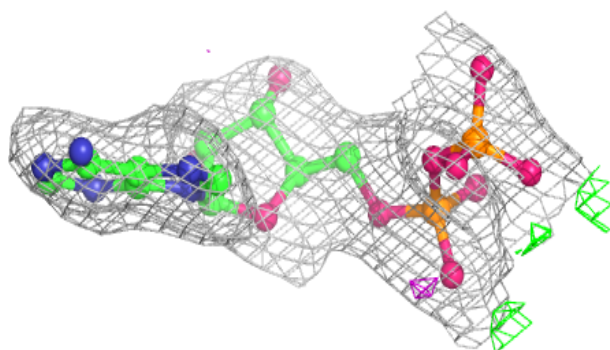
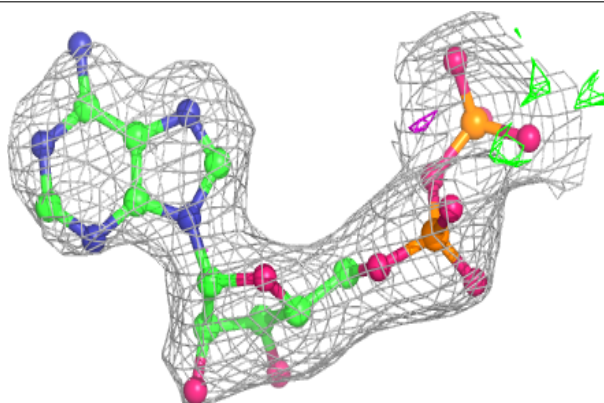


Electron density around ADP I 401:

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

**Electron density around ADP 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.