



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:19 PM UTC

PDB ID : 1XF9 / pdb_00001xf9
Title : Structure of NBD1 from murine CFTR- F508S mutant
Authors : Thibodeau, P.H.; Brautigam, C.A.; Machius, M.; Thomas, P.J.
Deposited on : 2004-09-14
Resolution : 2.70 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

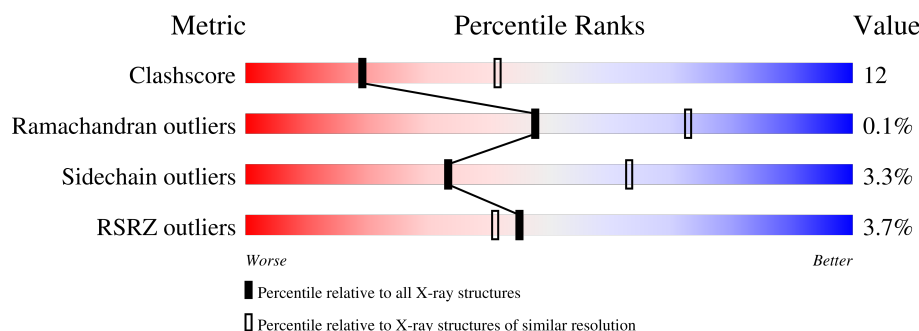
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	283	
1	B	283	
1	C	283	
1	D	283	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8601 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 Cystic fibrosis transmembrane conductance regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2078	1320	344	402	12			
1	B	265	Total	C	N	O	S	0	0	0
			2085	1324	345	404	12			
1	C	265	Total	C	N	O	S	0	0	0
			2085	1324	345	404	12			
1	D	260	Total	C	N	O	S	4	0	0
			2043	1297	339	395	12			

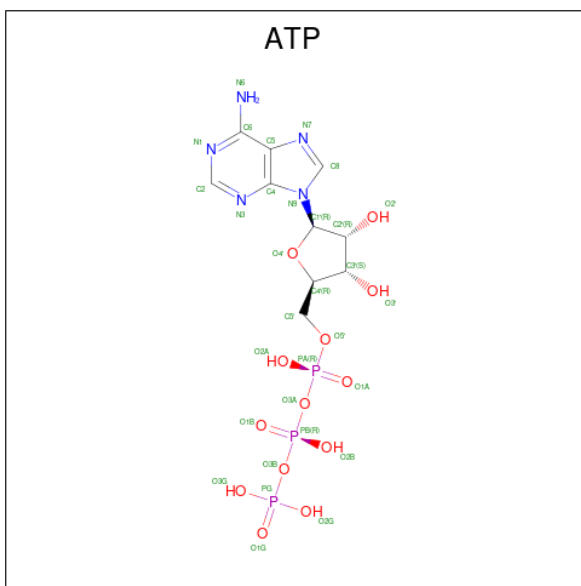
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	388	SER	MET	cloning artifact	UNP P26361
A	508	SER	PHE	engineered mutation	UNP P26361
B	388	SER	MET	cloning artifact	UNP P26361
B	508	SER	PHE	engineered mutation	UNP P26361
C	388	SER	MET	cloning artifact	UNP P26361
C	508	SER	PHE	engineered mutation	UNP P26361
D	388	SER	MET	cloning artifact	UNP P26361
D	508	SER	PHE	engineered mutation	UNP P26361

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

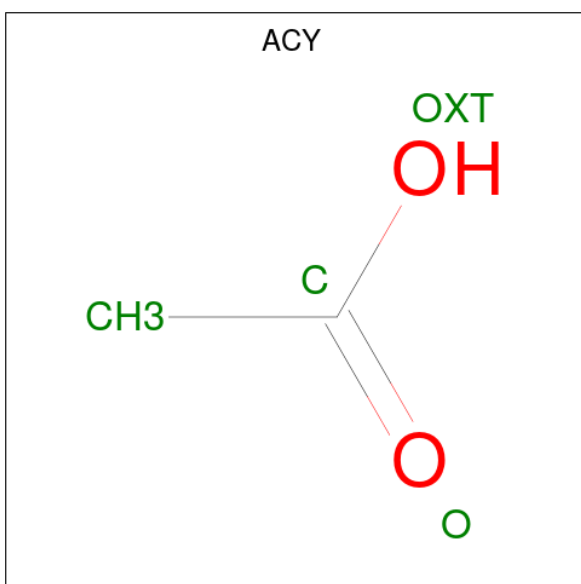
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	2	Total	Mg	0	0
			2	2		
2	C	2	Total	Mg	0	0
			2	2		
2	D	2	Total	Mg	0	0
			2	2		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $\text{C}_{10}\text{H}_{16}\text{N}_5\text{O}_{13}\text{P}_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 10	N 5	O 13	P 3	0	0
3	B	1	Total 31	C 10	N 5	O 13	P 3	0	0
3	C	1	Total 31	C 10	N 5	O 13	P 3	0	0
3	D	1	Total 31	C 10	N 5	O 13	P 3	0	0

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: $\text{C}_2\text{H}_4\text{O}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0

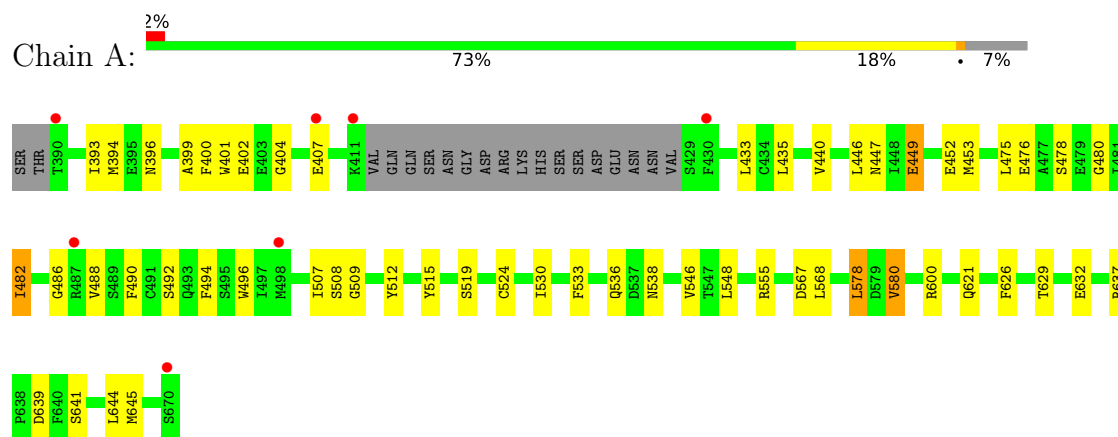
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	44	Total O 44 44	0	0
5	B	46	Total O 46 46	0	0
5	C	36	Total O 36 36	0	0
5	D	29	Total O 29 29	0	0

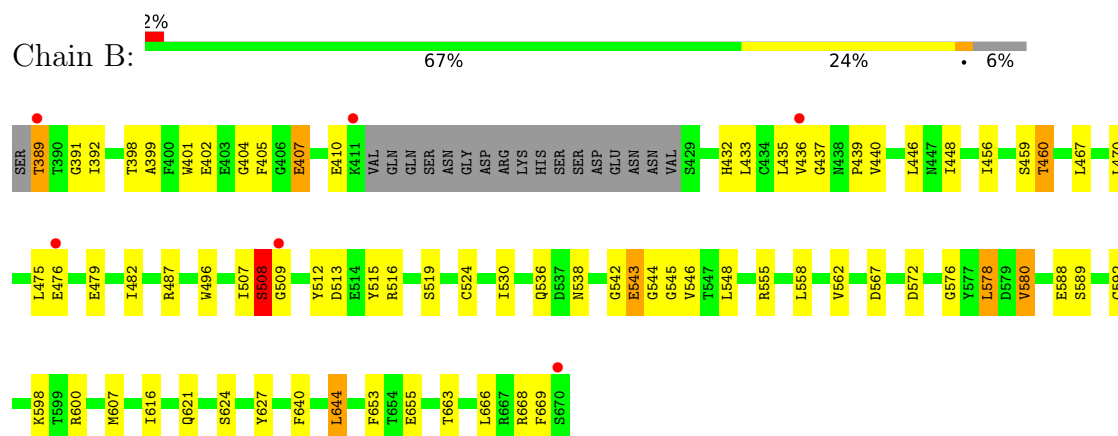
3 Residue-property plots [i](#)

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

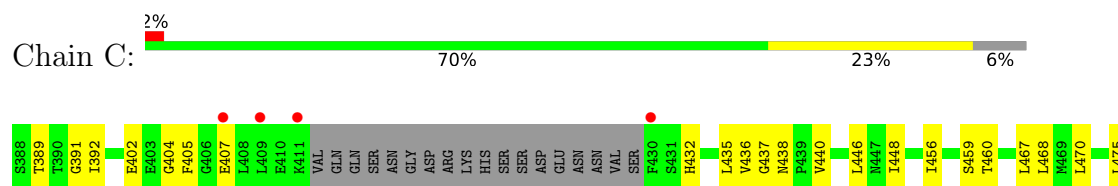
- Molecule 1: Cystic fibrosis transmembrane conductance regulator

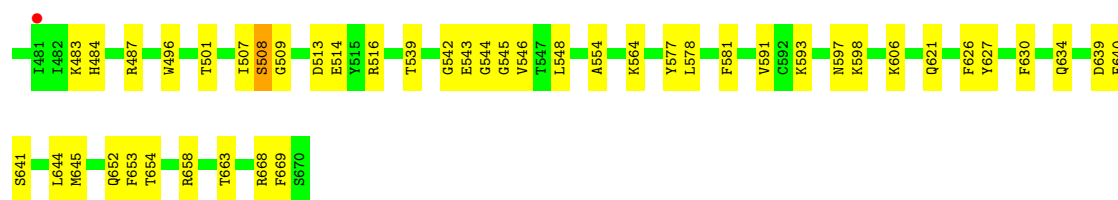


- Molecule 1: Cystic fibrosis transmembrane conductance regulator

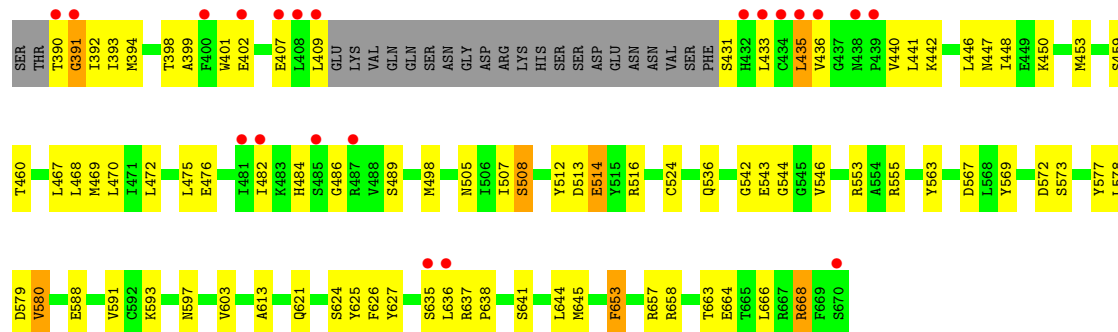


- Molecule 1: Cystic fibrosis transmembrane conductance regulator





● Molecule 1: Cystic fibrosis transmembrane conductance regulator



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	170.45Å 170.45Å 109.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.40 – 2.70 40.40 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.7 (40.40-2.70) 95.6 (40.40-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.253 0.197 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8601	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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: ATP, MG, ACY

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	0.62	0/2112	0.95	4/2838 (0.1%)
1	B	0.67	2/2119 (0.1%)	1.04	11/2848 (0.4%)
1	C	0.63	1/2119 (0.0%)	1.03	8/2848 (0.3%)
1	D	0.59	0/2076	1.01	13/2791 (0.5%)
All	All	0.63	3/8426 (0.0%)	1.01	36/11325 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	508	SER	C-O	-10.03	1.14	1.24
1	C	437	GLY	C-O	9.36	1.37	1.24
1	B	437	GLY	C-O	7.23	1.33	1.24

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	508	SER	CB-CA-C	-12.31	102.38	116.54
1	B	546	VAL	N-CA-C	10.07	120.90	110.62
1	C	546	VAL	N-CA-C	9.37	120.17	110.62
1	B	589	SER	N-CA-C	7.91	119.53	111.07
1	D	546	VAL	N-CA-C	7.88	119.61	110.62
1	B	653	PHE	N-CA-C	-7.16	101.60	110.41
1	D	508	SER	CB-CA-C	-7.00	107.72	117.23
1	D	514	GLU	N-CA-C	6.94	118.85	111.28
1	B	391	GLY	N-CA-C	-6.81	102.13	112.89
1	A	578	LEU	N-CA-C	6.52	118.39	111.28
1	B	578	LEU	N-CA-C	6.36	118.21	111.28
1	A	546	VAL	N-CA-C	6.33	119.90	111.17
1	D	591	VAL	N-CA-C	6.17	116.95	110.72
1	B	572	ASP	N-CA-C	6.12	119.49	107.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	391	GLY	N-CA-C	-6.00	102.32	112.77
1	C	578	LEU	N-CA-C	5.94	117.76	111.28
1	C	653	PHE	N-CA-C	-5.86	102.79	110.53
1	C	438	ASN	N-CA-CB	-5.67	99.30	109.68
1	D	536	GLN	CB-CA-C	-5.65	110.04	116.54
1	D	603	VAL	N-CA-C	-5.62	99.63	108.23
1	A	538	ASN	N-CA-C	-5.57	105.69	112.88
1	D	578	LEU	N-CA-C	5.57	118.28	111.82
1	C	438	ASN	N-CA-C	5.38	117.42	109.50
1	C	507	ILE	CB-CA-C	-5.38	105.02	112.24
1	B	580	VAL	N-CA-C	5.34	118.54	111.17
1	C	391	GLY	N-CA-C	-5.32	102.39	112.51
1	B	624	SER	N-CA-C	5.24	117.22	107.99
1	D	653	PHE	N-CA-C	-5.19	103.54	110.55
1	C	591	VAL	N-CA-C	5.19	115.96	110.72
1	D	624	SER	N-CA-C	5.16	117.72	108.17
1	D	513	ASP	N-CA-C	-5.11	98.95	107.80
1	B	538	ASN	N-CA-C	-5.06	106.52	113.30
1	B	508	SER	CA-C-O	5.06	127.25	119.21
1	A	580	VAL	N-CA-C	5.05	118.14	111.17
1	D	572	ASP	N-CA-C	5.03	116.84	107.99
1	D	577	TYR	N-CA-C	5.02	117.17	109.95

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	2078	0	2074	42	0
1	B	2085	0	2081	46	0
1	C	2085	0	2081	44	0
1	D	2043	0	2041	70	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2	0	0	0	0
3	A	31	0	12	0	0
3	B	31	0	12	1	0
3	C	31	0	12	0	0
3	D	31	0	12	5	0
4	A	8	0	6	0	0
4	B	4	0	3	0	0
4	C	4	0	3	0	0
4	D	8	0	6	0	0
5	A	44	0	0	3	0
5	B	46	0	0	0	0
5	C	36	0	0	2	0
5	D	29	0	0	5	0
All	All	8601	0	8343	196	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 12.

All (196) 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:D:402:GLU:HG3	1:D:476:GLU:OE2	1.48	1.13
1:B:402:GLU:HG3	1:B:476:GLU:OE2	1.47	1.13
1:C:487:ARG:HH21	1:C:564:LYS:HE3	1.13	1.11
1:A:402:GLU:HG3	1:A:476:GLU:OE2	1.54	1.07
1:C:389:THR:HG22	1:C:598:LYS:HD3	1.39	1.04
1:D:390:THR:HA	1:D:450:LYS:HE2	1.40	1.01
1:C:487:ARG:NH2	1:C:564:LYS:HE3	1.88	0.87
1:C:392:ILE:HG22	1:C:448:ILE:HB	1.57	0.86
1:D:390:THR:HG22	1:D:391:GLY:H	1.45	0.80
1:C:389:THR:CG2	1:C:598:LYS:HD3	2.13	0.78
1:D:402:GLU:CG	1:D:476:GLU:OE2	2.30	0.76
1:D:394:MET:HE1	1:D:470:LEU:HD23	1.74	0.69
1:D:453:MET:HE2	1:D:613:ALA:HA	1.75	0.69
1:C:389:THR:HG22	1:C:598:LYS:CD	2.19	0.68
1:D:467:LEU:HD12	1:D:467:LEU:O	1.93	0.68
1:C:392:ILE:CG2	1:C:448:ILE:HB	2.22	0.68
1:A:488:VAL:HG22	1:A:568:LEU:HB3	1.76	0.67
1:D:394:MET:HE2	1:D:482:ILE:HD12	1.76	0.67
1:A:639:ASP:HB2	5:A:148:HOH:O	1.93	0.67
1:D:409:LEU:HB3	3:D:4:ATP:H2	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:657:ARG:HH11	1:D:657:ARG:HG3	1.59	0.66
1:A:453:MET:HG2	1:A:600:ARG:HG3	1.77	0.66
1:A:496:TRP:HZ2	1:B:436:VAL:HG21	1.62	0.65
1:A:629:THR:OG1	1:A:632:GLU:HG3	1.97	0.65
1:D:409:LEU:HB3	3:D:4:ATP:C2	2.32	0.64
1:D:459:SER:HB3	1:D:666:LEU:HD12	1.79	0.64
1:C:459:SER:HB2	1:C:663:THR:HG23	1.80	0.63
1:A:402:GLU:CG	1:A:476:GLU:OE2	2.40	0.63
1:A:509:GLY:HA3	1:B:439:PRO:CG	2.29	0.63
1:D:392:ILE:HG13	1:D:484:HIS:HB3	1.80	0.63
1:B:405:PHE:CZ	1:B:475:LEU:HD22	2.33	0.62
1:D:401:TRP:CZ2	3:D:4:ATP:H4'	2.35	0.61
1:C:581:PHE:HE1	1:C:652:GLN:HE22	1.52	0.58
1:B:460:THR:HG22	1:B:663:THR:OG1	2.03	0.58
1:A:394:MET:HE2	1:A:482:ILE:HD12	1.85	0.58
1:C:654:THR:O	1:C:658:ARG:HG3	2.04	0.58
1:B:399:ALA:O	1:B:440:VAL:HG12	2.05	0.57
1:A:404:GLY:HA2	1:A:407:GLU:HG2	1.86	0.57
1:C:404:GLY:HA2	1:C:407:GLU:HG2	1.86	0.57
1:D:470:LEU:HD12	1:D:475:LEU:O	2.03	0.57
1:B:515:TYR:O	1:B:519:SER:HB2	2.05	0.56
1:D:393:ILE:HG12	1:D:447:ASN:OD1	2.05	0.56
1:B:470:LEU:HD12	1:B:475:LEU:O	2.06	0.56
1:C:668:ARG:NH1	1:C:669:PHE:CZ	2.74	0.55
1:C:487:ARG:HH21	1:C:564:LYS:CE	2.04	0.55
1:D:469:MET:HE2	1:D:472:LEU:HD12	1.87	0.55
1:D:635:SER:O	1:D:638:PRO:HD3	2.07	0.55
1:C:405:PHE:CZ	1:C:475:LEU:HD22	2.42	0.55
1:B:515:TYR:CE1	1:C:593:LYS:HE3	2.42	0.55
1:C:639:ASP:HB2	5:C:22:HOH:O	2.08	0.54
1:D:498:MET:HE1	1:D:508:SER:HA	1.90	0.54
1:D:516:ARG:HD3	1:D:563:TYR:CE1	2.43	0.54
1:D:626:PHE:C	1:D:626:PHE:CD1	2.86	0.54
1:B:398:THR:HB	1:B:479:GLU:HB2	1.89	0.54
1:B:410:GLU:HG2	3:B:2:ATP:N6	2.22	0.54
1:B:640:PHE:CE1	1:B:644:LEU:HD13	2.42	0.54
1:D:498:MET:HE2	1:D:505:ASN:HA	1.90	0.54
1:D:394:MET:HG3	1:D:446:LEU:HD11	1.90	0.53
1:D:446:LEU:HA	1:D:627:TYR:CZ	2.43	0.53
1:C:389:THR:HG21	1:C:597:ASN:O	2.08	0.53
1:B:402:GLU:CG	1:B:476:GLU:OE2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:483:LYS:O	1:C:484:HIS:HB3	2.07	0.53
1:D:446:LEU:HA	1:D:627:TYR:CE2	2.44	0.53
1:D:637:ARG:HG2	1:D:637:ARG:HH11	1.73	0.53
1:C:626:PHE:CD1	1:C:626:PHE:C	2.87	0.52
1:D:399:ALA:O	1:D:440:VAL:HG12	2.09	0.52
1:B:616:ILE:O	1:B:627:TYR:HA	2.11	0.51
1:A:446:LEU:HD12	1:A:446:LEU:O	2.11	0.51
1:B:524:CYS:O	1:B:555:ARG:HD2	2.11	0.51
1:C:640:PHE:CE1	1:C:644:LEU:HD13	2.46	0.50
1:C:516:ARG:HG3	1:C:516:ARG:HH11	1.76	0.50
1:C:630:PHE:O	1:C:634:GLN:HG3	2.12	0.50
1:D:390:THR:CA	1:D:450:LYS:HE2	2.27	0.50
1:A:440:VAL:HG11	1:A:475:LEU:HD21	1.93	0.50
1:D:588:GLU:HG2	1:D:593:LYS:HZ3	1.77	0.50
1:D:641:SER:O	1:D:645:MET:HG2	2.11	0.50
1:D:657:ARG:HG3	1:D:657:ARG:NH1	2.27	0.50
1:D:636:LEU:O	1:D:637:ARG:HG2	2.11	0.49
1:B:446:LEU:HD12	1:B:446:LEU:O	2.13	0.49
1:C:470:LEU:HD12	1:C:475:LEU:O	2.11	0.49
1:B:440:VAL:HG11	1:B:475:LEU:HD21	1.93	0.49
1:B:476:GLU:H	1:B:476:GLU:CD	2.20	0.49
1:A:637:ARG:HG2	1:A:637:ARG:HH11	1.78	0.48
1:A:486:GLY:HA3	1:A:567:ASP:OD2	2.12	0.48
1:B:516:ARG:HH11	1:B:516:ARG:HG3	1.79	0.48
1:D:469:MET:CE	1:D:472:LEU:HD12	2.43	0.48
1:D:440:VAL:HG11	1:D:475:LEU:HD21	1.96	0.48
1:A:515:TYR:O	1:A:519:SER:HB2	2.13	0.48
1:B:404:GLY:O	1:B:407:GLU:HG2	2.14	0.48
1:A:494:PHE:CD2	1:B:432:HIS:CE1	3.02	0.47
1:D:459:SER:HB2	1:D:663:THR:HG23	1.96	0.47
1:D:398:THR:HG23	1:D:442:LYS:HA	1.97	0.47
1:D:507:ILE:HD11	1:D:512:TYR:CD1	2.50	0.47
1:A:490:PHE:CE2	1:A:492:SER:HB3	2.50	0.47
1:C:508:SER:HB2	1:C:509:GLY:H	1.46	0.47
1:D:390:THR:HG23	1:D:450:LYS:HE2	1.96	0.47
1:D:401:TRP:CE2	1:D:433:LEU:HD13	2.49	0.47
1:C:501:THR:HA	1:C:539:THR:O	2.15	0.47
1:A:400:PHE:HD2	1:A:478:SER:HG	1.61	0.46
1:D:597:ASN:HA	5:D:170:HOH:O	2.13	0.46
1:A:524:CYS:O	1:A:555:ARG:HD2	2.15	0.46
1:B:588:GLU:O	1:B:592:CYS:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:456:ILE:HD13	1:C:467:LEU:HD23	1.97	0.46
1:D:653:PHE:O	1:D:658:ARG:NH1	2.49	0.46
1:A:626:PHE:C	1:A:626:PHE:CD1	2.94	0.46
1:C:389:THR:HG21	1:C:598:LYS:HA	1.97	0.46
1:D:392:ILE:HG22	1:D:448:ILE:HB	1.97	0.46
1:A:394:MET:CE	1:A:482:ILE:HD12	2.46	0.46
1:A:401:TRP:CE2	1:A:433:LEU:HD13	2.51	0.46
1:D:460:THR:HG22	1:D:663:THR:OG1	2.16	0.46
1:B:507:ILE:HD11	1:B:512:TYR:CD1	2.51	0.45
1:C:668:ARG:NH1	1:C:669:PHE:CE2	2.84	0.45
1:D:431:SER:O	1:D:435:LEU:HB2	2.16	0.45
1:B:459:SER:HB3	1:B:666:LEU:HD12	1.98	0.45
1:A:476:GLU:H	1:A:476:GLU:CD	2.24	0.45
1:B:542:GLY:O	1:B:543:GLU:C	2.59	0.45
1:C:404:GLY:O	1:C:407:GLU:HG2	2.16	0.45
1:A:404:GLY:O	1:A:407:GLU:HG2	2.16	0.45
1:D:486:GLY:HA3	1:D:567:ASP:OD2	2.16	0.45
1:D:588:GLU:HG2	1:D:593:LYS:NZ	2.31	0.45
1:B:607:MET:HA	1:B:607:MET:HE2	1.98	0.45
1:A:496:TRP:CZ2	1:B:436:VAL:HG21	2.47	0.45
1:D:512:TYR:OH	1:D:514:GLU:HG3	2.17	0.45
1:B:530:ILE:O	1:B:536:GLN:HA	2.16	0.45
1:C:641:SER:O	1:C:645:MET:HG2	2.17	0.45
1:C:402:GLU:OE1	1:C:402:GLU:HA	2.17	0.45
1:B:558:LEU:O	1:B:562:VAL:HG23	2.16	0.45
1:A:578:LEU:O	1:A:580:VAL:N	2.43	0.44
1:B:482:ILE:HG23	1:B:482:ILE:O	2.17	0.44
1:D:544:GLY:O	1:D:553:ARG:HD3	2.16	0.44
1:A:641:SER:O	1:A:645:MET:HG2	2.16	0.44
1:B:487:ARG:HH11	1:B:487:ARG:HG3	1.82	0.44
1:D:580:VAL:HG22	5:D:43:HOH:O	2.16	0.44
1:A:578:LEU:O	1:A:580:VAL:HG13	2.17	0.44
1:B:544:GLY:O	1:B:545:GLY:C	2.59	0.44
1:A:507:ILE:CD1	1:A:512:TYR:HD1	2.30	0.44
1:C:544:GLY:O	1:C:545:GLY:C	2.61	0.44
1:D:542:GLY:O	1:D:543:GLU:C	2.61	0.44
1:A:637:ARG:HG2	1:A:637:ARG:NH1	2.33	0.44
1:D:524:CYS:O	1:D:555:ARG:HD2	2.18	0.44
1:A:393:ILE:HG12	1:A:447:ASN:OD1	2.18	0.44
1:A:396:ASN:O	1:A:480:GLY:HA3	2.18	0.44
1:A:530:ILE:O	1:A:536:GLN:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:456:ILE:HD13	1:B:467:LEU:HD23	2.00	0.44
1:C:446:LEU:HA	1:C:627:TYR:CZ	2.53	0.44
1:D:626:PHE:CD1	1:D:627:TYR:N	2.86	0.44
1:C:440:VAL:HG11	1:C:475:LEU:HD21	2.00	0.43
1:C:432:HIS:O	1:C:436:VAL:HG23	2.17	0.43
1:A:507:ILE:HD11	1:A:512:TYR:HD1	1.82	0.43
1:C:640:PHE:CD1	1:C:640:PHE:C	2.96	0.43
1:D:436:VAL:O	1:D:436:VAL:HG12	2.19	0.43
1:D:498:MET:HG2	1:D:505:ASN:OD1	2.18	0.43
1:A:449:GLU:HB2	1:A:452:GLU:HG3	2.01	0.43
1:D:398:THR:HA	1:D:441:LEU:O	2.19	0.43
1:A:530:ILE:HA	1:A:533:PHE:CD2	2.54	0.43
1:A:399:ALA:O	1:A:440:VAL:HG12	2.19	0.42
1:B:392:ILE:HG22	1:B:448:ILE:HB	2.01	0.42
1:C:598:LYS:HD3	1:C:598:LYS:HA	1.88	0.42
1:D:567:ASP:O	5:D:86:HOH:O	2.21	0.42
1:A:555:ARG:CG	5:A:121:HOH:O	2.66	0.42
1:D:507:ILE:CD1	1:D:512:TYR:CD1	3.02	0.42
1:A:494:PHE:HD2	1:B:432:HIS:CE1	2.37	0.42
1:B:389:THR:HA	1:B:567:ASP:HA	2.01	0.42
1:B:401:TRP:CE2	1:B:433:LEU:HD13	2.55	0.42
1:B:668:ARG:NH1	1:B:669:PHE:CZ	2.88	0.42
1:D:390:THR:HG23	1:D:450:LYS:CE	2.49	0.42
1:A:476:GLU:OE1	1:A:476:GLU:N	2.48	0.42
1:B:496:TRP:H	1:B:496:TRP:CD1	2.37	0.42
1:B:513:ASP:OD1	1:B:513:ASP:C	2.62	0.42
1:B:576:GLY:HA3	1:B:655:GLU:HG3	2.01	0.42
1:D:390:THR:HG22	1:D:391:GLY:N	2.23	0.41
1:B:508:SER:HB2	1:B:509:GLY:H	1.07	0.41
1:B:578:LEU:O	1:B:580:VAL:HG13	2.20	0.41
1:D:625:TYR:OH	1:D:637:ARG:HD2	2.20	0.41
1:D:637:ARG:HG2	1:D:637:ARG:NH1	2.35	0.41
1:B:404:GLY:HA2	1:B:407:GLU:HG2	2.02	0.41
1:B:640:PHE:CZ	1:B:644:LEU:HD13	2.54	0.41
1:C:516:ARG:HG3	1:C:516:ARG:NH1	2.34	0.41
1:B:598:LYS:O	1:B:600:ARG:NH1	2.51	0.41
1:C:542:GLY:O	1:C:543:GLU:C	2.63	0.41
1:D:467:LEU:HD12	1:D:467:LEU:C	2.45	0.41
1:D:467:LEU:O	1:D:470:LEU:HB3	2.20	0.41
1:D:579:ASP:OD1	1:D:579:ASP:N	2.53	0.41
1:D:401:TRP:HZ2	3:D:4:ATP:O3'	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:573:SER:HA	5:D:164:HOH:O	2.20	0.41
1:C:513:ASP:O	1:C:514:GLU:C	2.64	0.41
1:C:513:ASP:OD1	1:C:513:ASP:C	2.65	0.40
1:D:409:LEU:CB	3:D:4:ATP:H2	2.32	0.40
1:D:489:SER:HB3	1:D:569:TYR:CD2	2.56	0.40
1:A:555:ARG:HG3	5:A:121:HOH:O	2.21	0.40
1:C:606:LYS:NZ	5:C:154:HOH:O	2.54	0.40
1:A:446:LEU:HD12	1:A:446:LEU:C	2.46	0.40
1:C:554:ALA:HB2	1:C:577:TYR:CE1	2.57	0.40
1:D:507:ILE:CD1	1:D:512:TYR:HD1	2.34	0.40
1:D:514:GLU:OE2	5:D:117:HOH:O	2.22	0.40
1:C:496:TRP:CD1	1:C:496:TRP:H	2.38	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	260/283 (92%)	248 (95%)	12 (5%)	0	100	100
1	B	261/283 (92%)	250 (96%)	11 (4%)	0	100	100
1	C	261/283 (92%)	244 (94%)	17 (6%)	0	100	100
1	D	256/283 (90%)	242 (94%)	13 (5%)	1 (0%)	30	54
All	All	1038/1132 (92%)	984 (95%)	53 (5%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	668	ARG

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	231/249 (93%)	224 (97%)	7 (3%)	36	66
1	B	232/249 (93%)	223 (96%)	9 (4%)	28	57
1	C	232/249 (93%)	226 (97%)	6 (3%)	40	70
1	D	227/249 (91%)	219 (96%)	8 (4%)	32	61
All	All	922/996 (93%)	892 (97%)	30 (3%)	33	63

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	435	LEU
1	A	449	GLU
1	A	482	ILE
1	A	508	SER
1	A	548	LEU
1	A	621	GLN
1	A	644	LEU
1	B	389	THR
1	B	407	GLU
1	B	435	LEU
1	B	460	THR
1	B	508	SER
1	B	543	GLU
1	B	548	LEU
1	B	621	GLN
1	B	644	LEU
1	C	435	LEU
1	C	460	THR
1	C	468	LEU
1	C	508	SER
1	C	548	LEU
1	C	621	GLN
1	D	407	GLU
1	D	435	LEU

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Mol	Chain	Res	Type
1	D	468	LEU
1	D	580	VAL
1	D	621	GLN
1	D	644	LEU
1	D	664	GLU
1	D	668	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	597	ASN
1	A	621	GLN
1	A	652	GLN
1	B	396	ASN
1	B	438	ASN
1	B	443	ASN
1	B	527	GLN
1	B	585	GLN
1	C	438	ASN
1	C	527	GLN
1	C	585	GLN
1	C	621	GLN
1	C	652	GLN
1	D	396	ASN
1	D	443	ASN
1	D	493	GLN
1	D	621	GLN
1	D	652	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 17 ligands modelled in this entry, 7 are monoatomic - leaving 10 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	ACY	B	6	-	3,3,3	1.18	0	3,3,3	0.96	0
4	ACY	A	5	-	3,3,3	1.30	0	3,3,3	0.98	0
4	ACY	C	7	-	3,3,3	1.14	0	3,3,3	0.89	0
3	ATP	B	2	2	32,33,33	0.68	1 (3%)	48,52,52	0.91	3 (6%)
4	ACY	A	9	-	3,3,3	1.12	0	3,3,3	0.95	0
4	ACY	D	8	-	3,3,3	1.15	0	3,3,3	0.93	0
4	ACY	D	10	-	3,3,3	0.83	0	3,3,3	1.96	2 (66%)
3	ATP	A	1	2	32,33,33	0.60	0	48,52,52	0.92	4 (8%)
3	ATP	D	4	2	32,33,33	0.66	0	48,52,52	0.90	3 (6%)
3	ATP	C	3	2	32,33,33	0.72	1 (3%)	48,52,52	0.94	3 (6%)

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	ATP	D	4	2	-	4/22/38/38	0/3/3/3
3	ATP	A	1	2	-	0/22/38/38	0/3/3/3
3	ATP	B	2	2	-	6/22/38/38	0/3/3/3
3	ATP	C	3	2	-	4/22/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2	ATP	PG-O2G	-2.29	1.46	1.54
3	C	3	ATP	PG-O2G	-2.06	1.47	1.54

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	3	ATP	O3G-PG-O2G	-3.13	96.07	107.80
3	B	2	ATP	O3G-PG-O2G	-2.92	96.85	107.80
3	A	1	ATP	O3G-PG-O2G	-2.55	98.23	107.80
3	D	4	ATP	O3G-PG-O2G	-2.46	98.58	107.80
4	D	10	ACY	OXT-C-CH3	2.40	125.10	115.05
4	D	10	ACY	OXT-C-O	-2.40	113.14	122.03
3	C	3	ATP	O2G-PG-O1G	2.39	120.13	110.83
3	B	2	ATP	O3G-PG-O1G	2.36	120.03	110.83
3	A	1	ATP	C2'-C1'-N9	2.23	118.86	113.30
3	A	1	ATP	O2G-PG-O1G	2.18	119.34	110.83
3	D	4	ATP	O2G-PG-O1G	2.17	119.29	110.83
3	A	1	ATP	O2B-PB-O3A	2.12	112.99	107.27
3	C	3	ATP	O3G-PG-O1G	2.08	118.93	110.83
3	B	2	ATP	O2B-PB-O3A	2.05	112.82	107.27
3	D	4	ATP	O3G-PG-O1G	2.03	118.75	110.83

There are no chirality outliers.

All (14) torsion outliers are listed below:

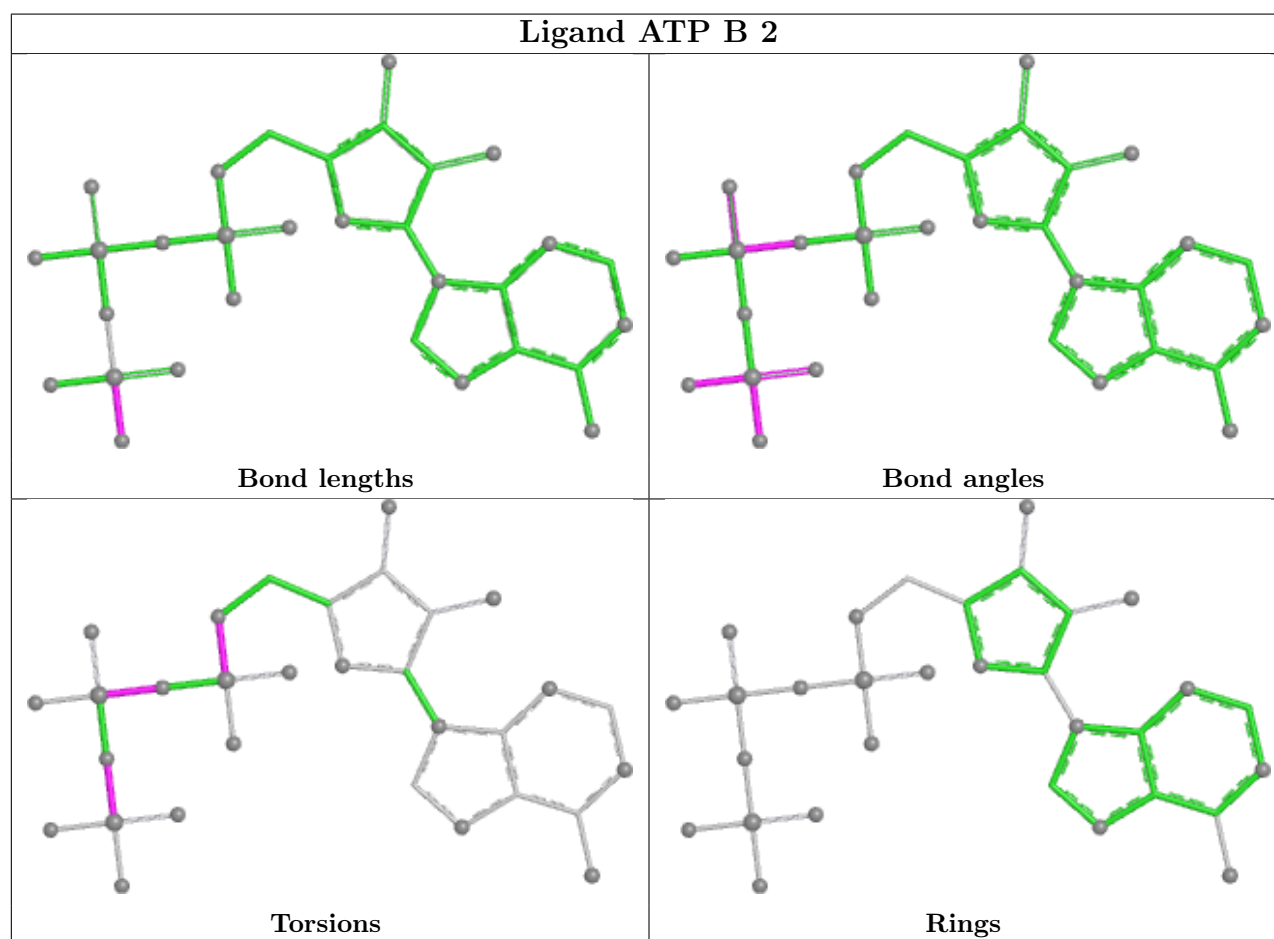
Mol	Chain	Res	Type	Atoms
3	B	2	ATP	C5'-O5'-PA-O1A
3	B	2	ATP	C5'-O5'-PA-O2A
3	B	2	ATP	C5'-O5'-PA-O3A
3	C	3	ATP	C5'-O5'-PA-O2A
3	C	3	ATP	C5'-O5'-PA-O3A
3	D	4	ATP	C5'-O5'-PA-O3A
3	D	4	ATP	O4'-C1'-N9-C4
3	D	4	ATP	O4'-C1'-N9-C8
3	B	2	ATP	PB-O3B-PG-O3G
3	B	2	ATP	PA-O3A-PB-O1B
3	C	3	ATP	C5'-O5'-PA-O1A
3	D	4	ATP	C5'-O5'-PA-O1A
3	B	2	ATP	PA-O3A-PB-O2B
3	C	3	ATP	PA-O3A-PB-O2B

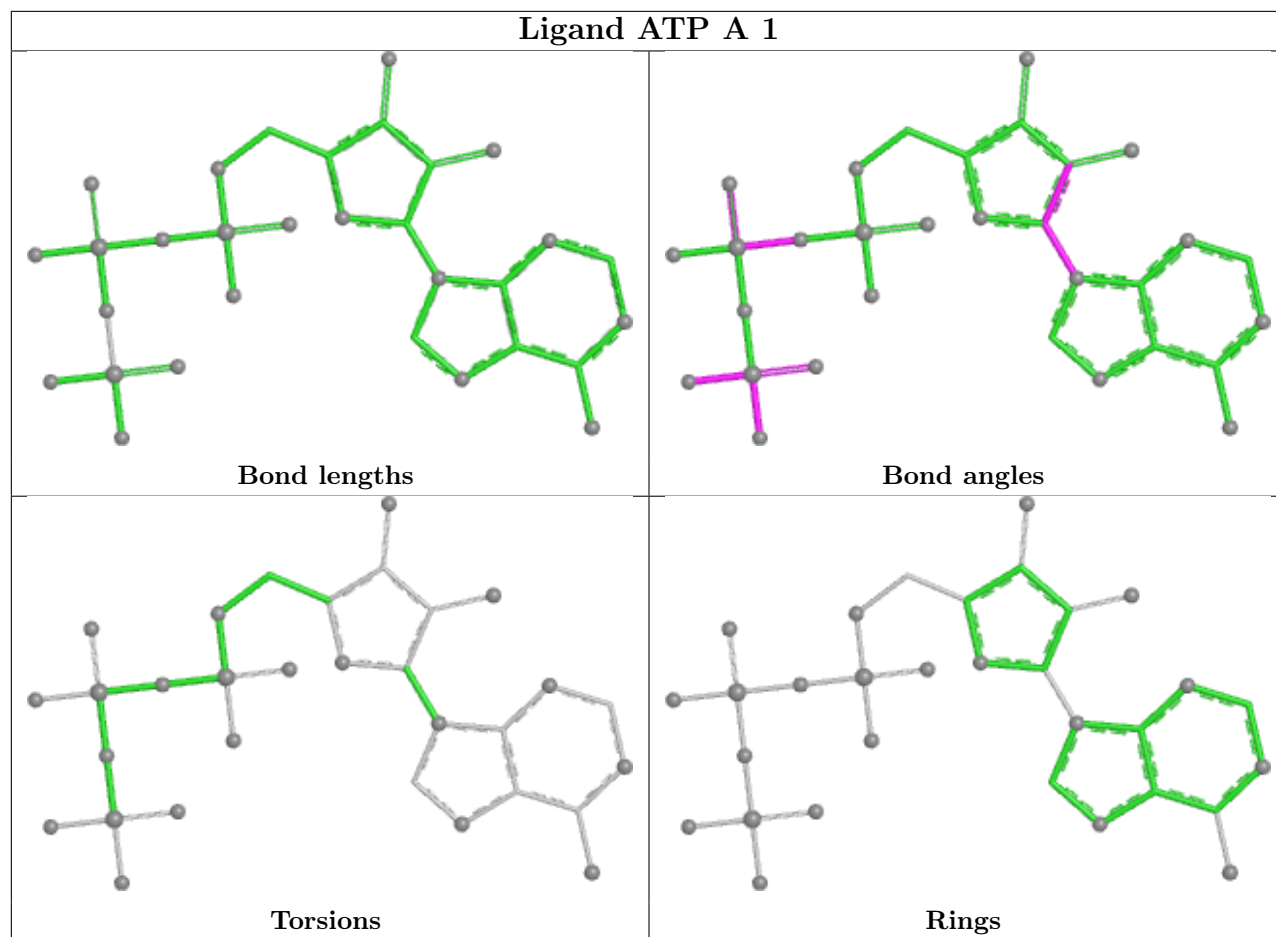
There are no ring outliers.

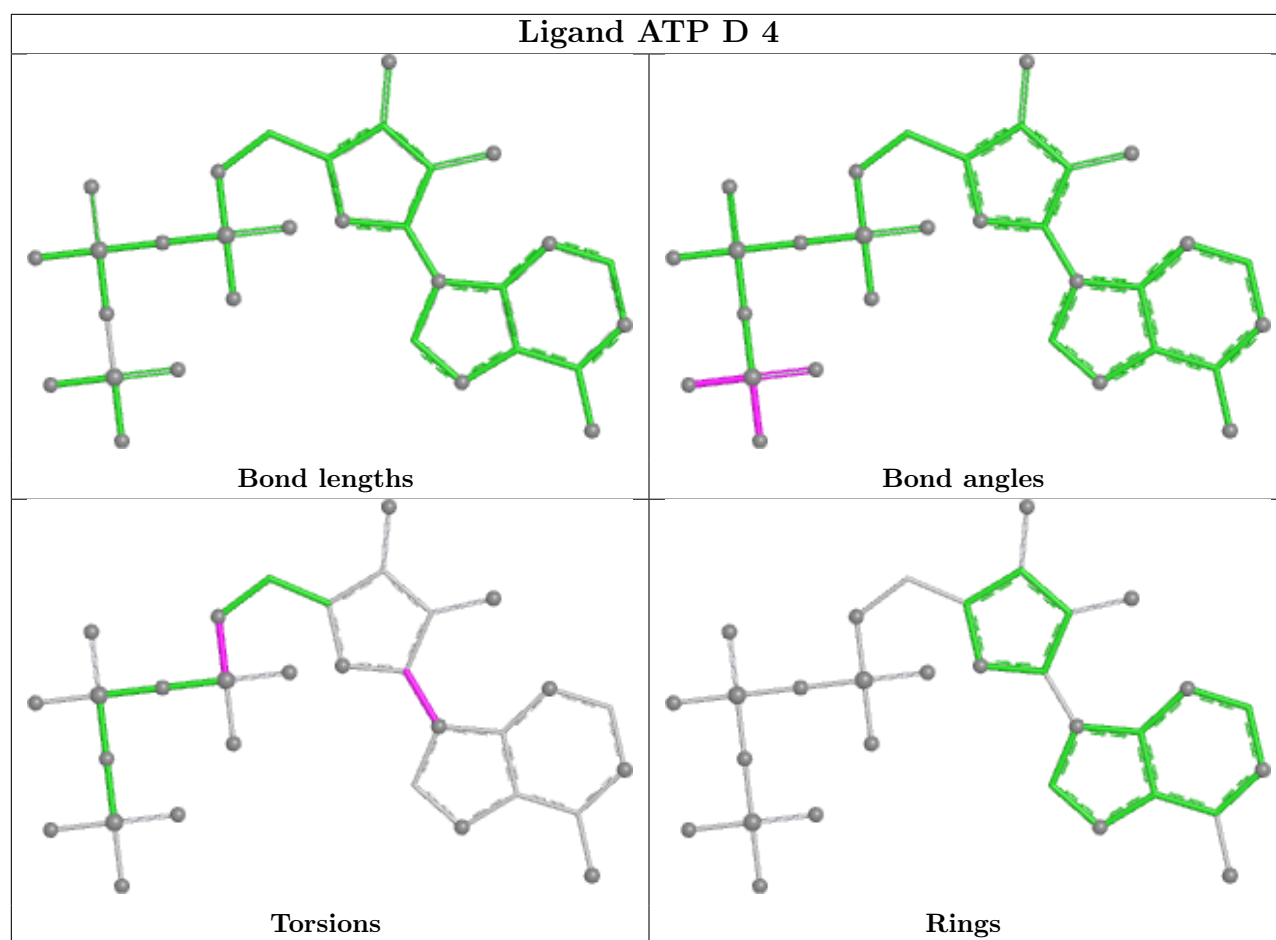
2 monomers are involved in 6 short contacts:

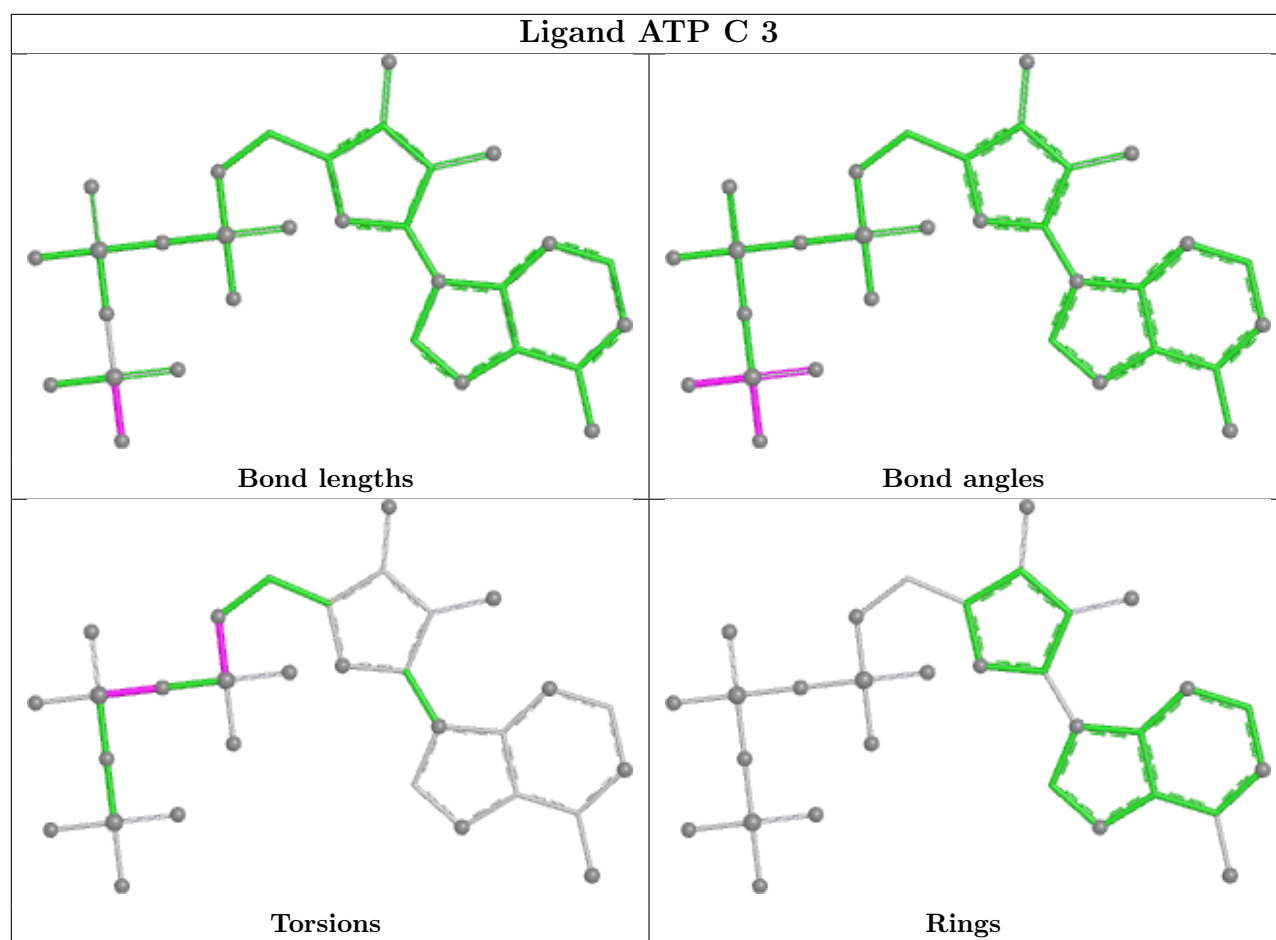
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2	ATP	1	0
3	D	4	ATP	5	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	264/283 (93%)	-0.08	7 (2%) 56 53	24, 46, 77, 99	0
1	B	265/283 (93%)	-0.09	6 (2%) 61 58	30, 46, 75, 99	0
1	C	265/283 (93%)	-0.10	5 (1%) 66 63	27, 46, 78, 96	0
1	D	260/283 (91%)	0.34	21 (8%) 18 15	29, 55, 103, 114	1 (0%)
All	All	1054/1132 (93%)	0.02	39 (3%) 45 41	24, 48, 86, 114	1 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	436	VAL	4.6
1	D	390	THR	4.2
1	D	433	LEU	4.1
1	D	435	LEU	3.6
1	D	432	HIS	3.6
1	D	481	ILE	3.4
1	D	402	GLU	3.4
1	D	482	ILE	3.2
1	D	408	LEU	3.2
1	D	487	ARG	3.1
1	A	498	MET	3.1
1	D	409	LEU	3.0
1	B	476	GLU	2.8
1	D	439	PRO	2.7
1	D	407	GLU	2.5
1	D	485	SER	2.5
1	D	635	SER	2.5
1	A	411	LYS	2.5
1	B	509	GLY	2.5
1	C	430	PHE	2.5
1	B	411	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	407	GLU	2.5
1	A	670	SER	2.4
1	A	430	PHE	2.4
1	D	670	SER	2.4
1	C	409	LEU	2.3
1	D	438	ASN	2.3
1	B	389	THR	2.3
1	C	481	ILE	2.3
1	C	411	LYS	2.2
1	B	670	SER	2.2
1	B	436	VAL	2.2
1	D	434	CYS	2.1
1	D	391	GLY	2.1
1	A	407	GLU	2.1
1	A	487	ARG	2.0
1	A	390	THR	2.0
1	D	636	LEU	2.0
1	D	400	PHE	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 [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
4	ACY	D	10	4/4	0.81	0.25	57,63,66,67	0
4	ACY	A	9	4/4	0.89	0.16	46,49,54,58	0
2	MG	D	17	1/1	0.89	0.10	106,106,106,106	0
3	ATP	D	4	31/31	0.90	0.12	59,105,125,126	0
2	MG	B	15	1/1	0.92	0.10	65,65,65,65	0

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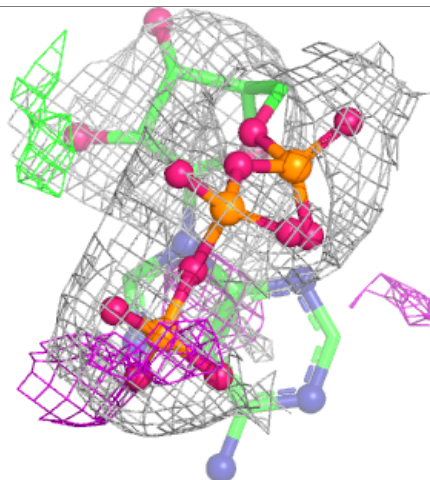
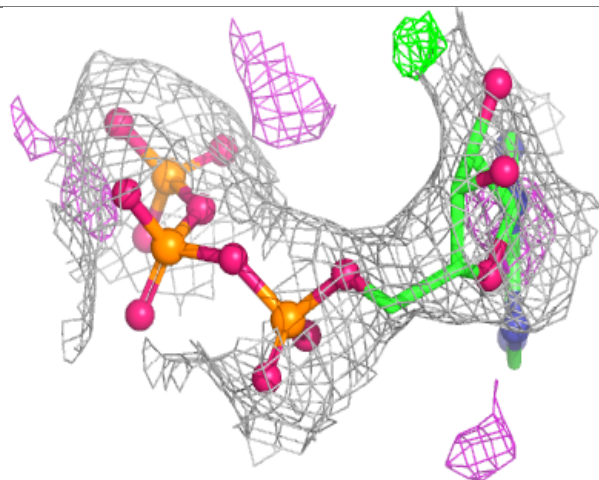
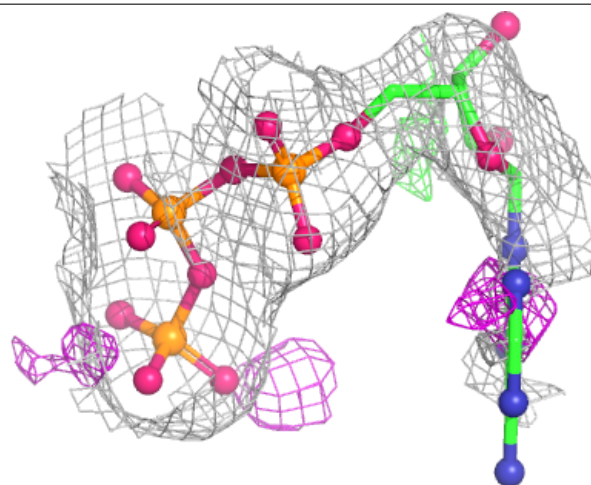
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	C	16	1/1	0.94	0.12	69,69,69,69	0
4	ACY	A	5	4/4	0.95	0.09	41,41,43,45	0
3	ATP	C	3	31/31	0.95	0.10	47,65,108,109	0
3	ATP	B	2	31/31	0.95	0.10	34,60,104,106	0
2	MG	D	14	1/1	0.96	0.06	57,57,57,57	0
4	ACY	B	6	4/4	0.96	0.09	35,40,42,43	0
4	ACY	C	7	4/4	0.96	0.09	38,41,43,44	0
3	ATP	A	1	31/31	0.96	0.10	38,77,106,107	0
2	MG	C	12	1/1	0.97	0.05	60,60,60,60	0
2	MG	A	13	1/1	0.98	0.05	47,47,47,47	0
4	ACY	D	8	4/4	0.98	0.07	43,45,46,49	0
2	MG	B	11	1/1	0.98	0.04	47,47,47,47	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.

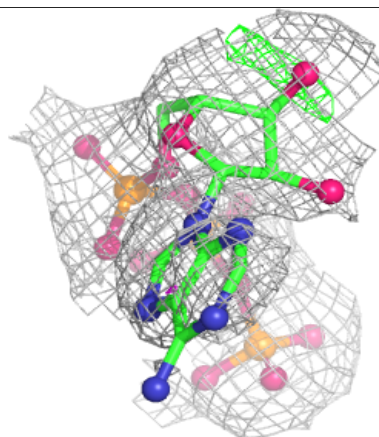
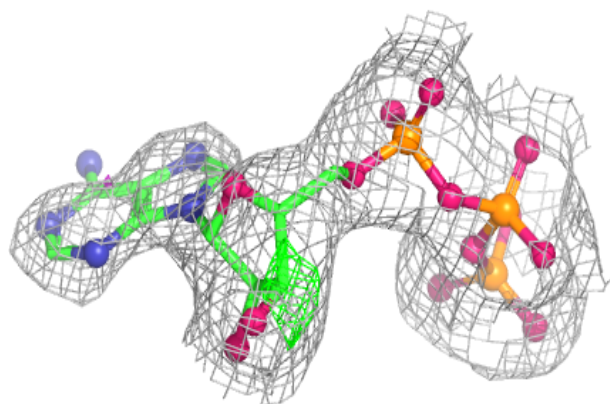
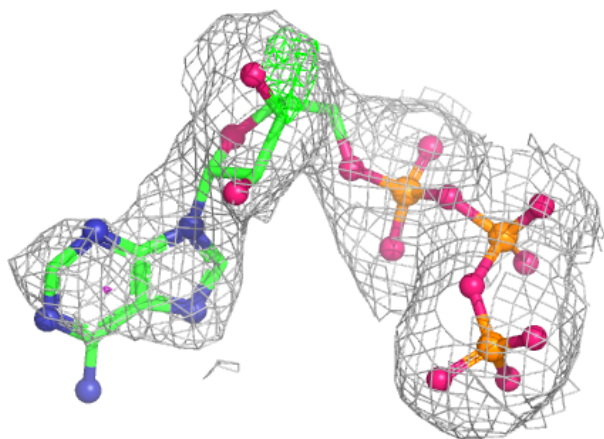
Electron density around ATP D 4:

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

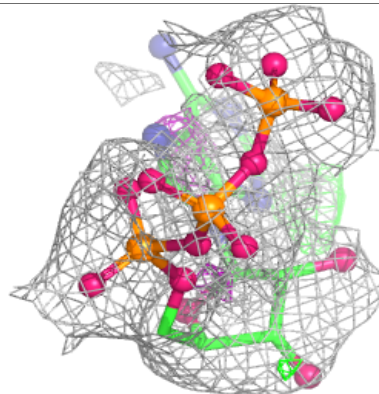
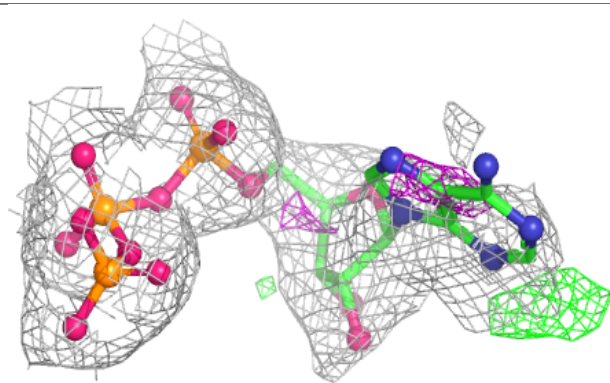
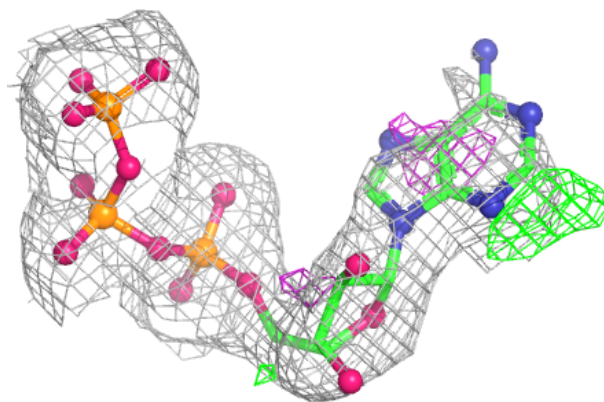


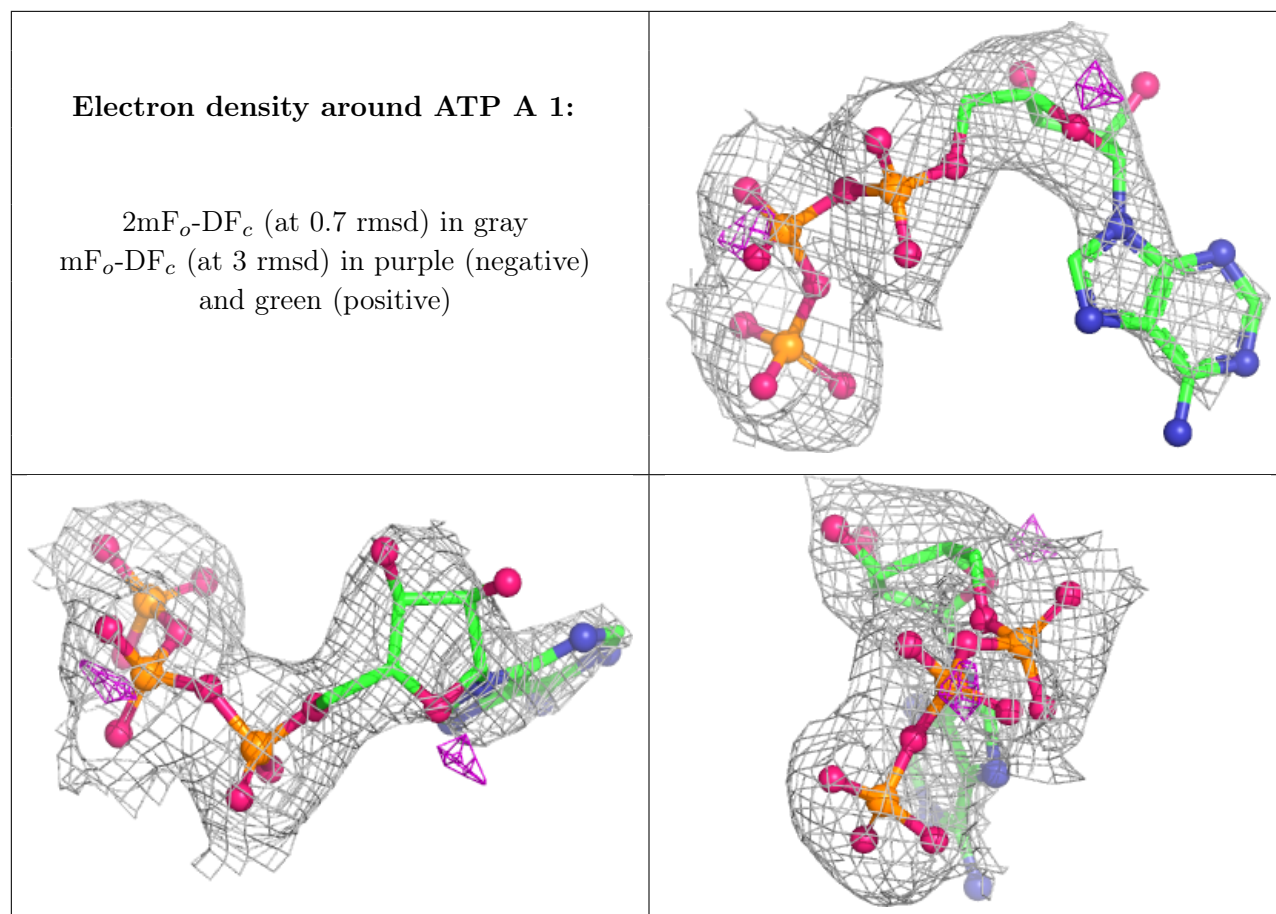
Electron density around ATP C 3:

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 ATP B 2:**

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.