



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 02:46 AM BST

PDB ID : 2A2C
Title : x-ray structure of human N-acetyl galactosamine kinase complexed with Mg-ADP and N-acetyl galactosamine 1-phosphate
Authors : Thoden, J.B.; Holden, H.M.
Deposited on : 2005-06-22
Resolution : 1.65 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

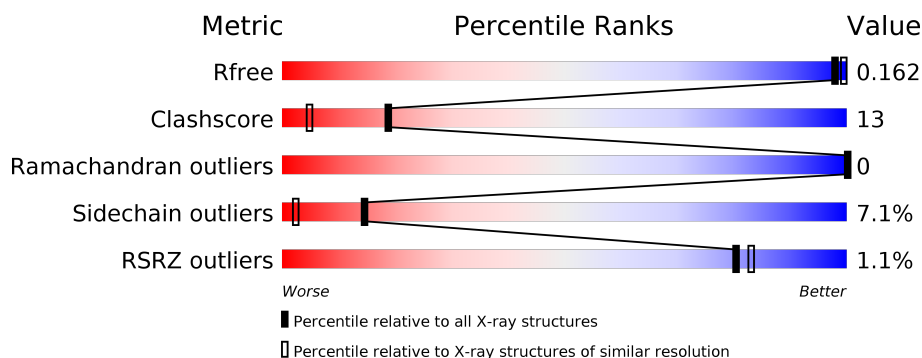
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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 on 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	478	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3976 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.

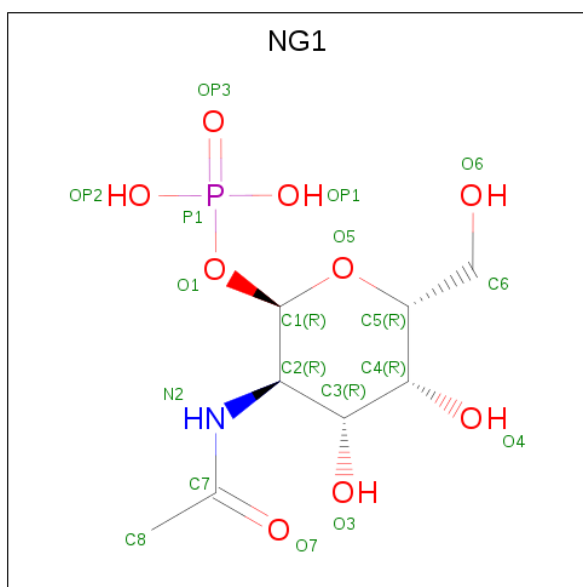
- Molecule 1 is a protein called N-acetylgalactosamine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	446	3470	2206	589	645	30	0	9	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP Q01415
A	-18	GLY	-	cloning artifact	UNP Q01415
A	-17	SER	-	cloning artifact	UNP Q01415
A	-16	SER	-	cloning artifact	UNP Q01415
A	-15	HIS	-	expression tag	UNP Q01415
A	-14	HIS	-	expression tag	UNP Q01415
A	-13	HIS	-	expression tag	UNP Q01415
A	-12	HIS	-	expression tag	UNP Q01415
A	-11	HIS	-	expression tag	UNP Q01415
A	-10	HIS	-	expression tag	UNP Q01415
A	-9	SER	-	cloning artifact	UNP Q01415
A	-8	SER	-	cloning artifact	UNP Q01415
A	-7	GLU	-	cloning artifact	UNP Q01415
A	-6	ASN	-	cloning artifact	UNP Q01415
A	-5	LEU	-	cloning artifact	UNP Q01415
A	-4	TYR	-	cloning artifact	UNP Q01415
A	-3	PHE	-	cloning artifact	UNP Q01415
A	-2	GLN	-	cloning artifact	UNP Q01415
A	-1	GLY	-	cloning artifact	UNP Q01415
A	0	HIS	-	cloning artifact	UNP Q01415

- Molecule 2 is 2-acetamido-2-deoxy-1-O-phosphono-alpha-D-galactopyranose (three-letter code: NG1) (formula: C₈H₁₆NO₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			19	8	1	9	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

- Molecule 6 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

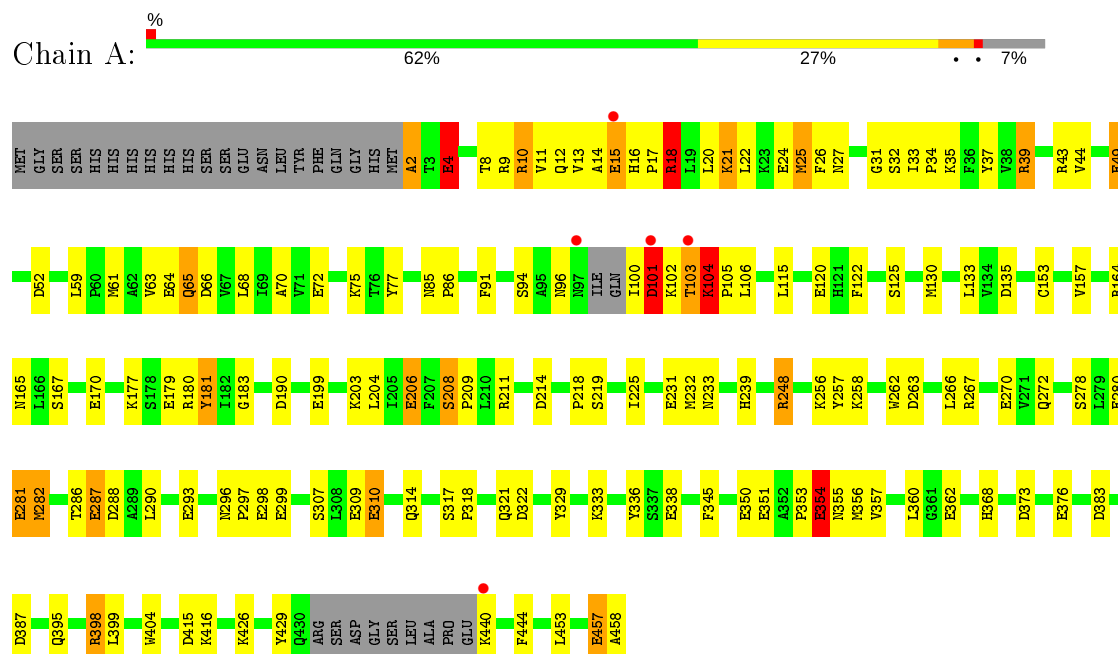
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	457	Total	O	0	0
			457	457		

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: N-acetylgalactosamine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	123.80 Å 123.80 Å 60.10 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 1.65 29.74 – 1.65	Depositor EDS
% Data completeness (in resolution range)	94.3 (50.00-1.65) 94.4 (29.74-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.86 (at 1.65 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.167 , 0.201 0.165 , 0.162	Depositor DCC
R_{free} test set	5980 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 86.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.054 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3976	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: NA, MG, NG1, ADP, CL

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	0.99	25/3572 (0.7%)	1.40	48/4828 (1.0%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	49	GLU	CD-OE2	8.03	1.34	1.25
1	A	354	GLU	CD-OE2	7.10	1.33	1.25
1	A	376	GLU	CD-OE2	6.81	1.33	1.25
1	A	338	GLU	CD-OE2	6.62	1.32	1.25
1	A	350	GLU	CD-OE2	6.51	1.32	1.25
1	A	4	GLU	CD-OE2	6.42	1.32	1.25
1	A	299	GLU	CD-OE2	6.39	1.32	1.25
1	A	309	GLU	CD-OE2	6.39	1.32	1.25
1	A	179	GLU	CD-OE2	6.37	1.32	1.25
1	A	206	GLU	CD-OE1	-6.22	1.18	1.25
1	A	120	GLU	CD-OE2	6.08	1.32	1.25
1	A	64	GLU	CD-OE2	6.04	1.32	1.25
1	A	280	GLU	CD-OE2	5.93	1.32	1.25
1	A	457	GLU	CD-OE2	5.83	1.32	1.25
1	A	351	GLU	CD-OE2	5.81	1.32	1.25
1	A	293	GLU	CD-OE2	5.73	1.31	1.25
1	A	310	GLU	CD-OE2	5.67	1.31	1.25
1	A	298	GLU	CD-OE2	5.59	1.31	1.25
1	A	199	GLU	CD-OE2	5.59	1.31	1.25
1	A	24	GLU	CD-OE2	5.45	1.31	1.25
1	A	281	GLU	CD-OE2	5.34	1.31	1.25
1	A	362	GLU	CD-OE2	5.25	1.31	1.25
1	A	15	GLU	CD-OE2	5.24	1.31	1.25
1	A	287	GLU	CD-OE2	5.22	1.31	1.25
1	A	231	GLU	CD-OE2	5.14	1.31	1.25

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	398	ARG	NE-CZ-NH2	-10.22	115.19	120.30
1	A	10	ARG	NE-CZ-NH2	-9.91	115.35	120.30
1	A	398	ARG	NE-CZ-NH1	8.83	124.71	120.30
1	A	52	ASP	CB-CG-OD2	-8.52	110.63	118.30
1	A	248[A]	ARG	NE-CZ-NH1	8.45	124.52	120.30
1	A	248[B]	ARG	NE-CZ-NH1	8.45	124.52	120.30
1	A	101	ASP	CB-CG-OD2	-8.36	110.78	118.30
1	A	288	ASP	CB-CG-OD2	-8.30	110.83	118.30
1	A	383	ASP	CB-CG-OD2	-8.23	110.89	118.30
1	A	248[A]	ARG	NE-CZ-NH2	-8.20	116.20	120.30
1	A	248[B]	ARG	NE-CZ-NH2	-8.20	116.20	120.30
1	A	373	ASP	CB-CG-OD2	-7.86	111.23	118.30
1	A	181	TYR	CB-CG-CD2	-7.59	116.45	121.00
1	A	373	ASP	CB-CG-OD1	7.39	124.95	118.30
1	A	214	ASP	CB-CG-OD2	-7.24	111.79	118.30
1	A	453	LEU	CB-CA-C	-7.19	96.53	110.20
1	A	18	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	A	180	ARG	NE-CZ-NH2	-7.13	116.73	120.30
1	A	322	ASP	CB-CG-OD2	-7.00	112.00	118.30
1	A	8	THR	CA-CB-CG2	-6.90	102.74	112.40
1	A	288	ASP	CB-CG-OD1	6.79	124.41	118.30
1	A	383	ASP	CB-CG-OD1	6.76	124.39	118.30
1	A	257	TYR	CB-CG-CD2	6.55	124.93	121.00
1	A	190	ASP	CB-CG-OD1	6.50	124.15	118.30
1	A	190	ASP	CB-CG-OD2	-6.46	112.49	118.30
1	A	257	TYR	CB-CG-CD1	-6.46	117.13	121.00
1	A	52	ASP	CB-CG-OD1	6.38	124.04	118.30
1	A	135	ASP	CB-CG-OD1	6.36	124.02	118.30
1	A	387	ASP	CB-CG-OD2	-6.30	112.63	118.30
1	A	345	PHE	CB-CG-CD2	-6.19	116.47	120.80
1	A	2	ALA	CB-CA-C	-6.07	101.00	110.10
1	A	429	TYR	CB-CG-CD2	-5.91	117.45	121.00
1	A	263	ASP	CB-CG-OD2	-5.73	113.15	118.30
1	A	135	ASP	CB-CG-OD2	-5.68	113.18	118.30
1	A	429	TYR	CB-CG-CD1	5.67	124.41	121.00
1	A	415	ASP	CB-CG-OD1	5.62	123.36	118.30
1	A	387	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	91	PHE	CB-CG-CD1	-5.56	116.91	120.80
1	A	180	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	164	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	A	39	ARG	NE-CZ-NH2	5.43	123.02	120.30
1	A	68	LEU	CB-CG-CD2	-5.32	101.97	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	104	LYS	N-CA-CB	-5.25	101.15	110.60
1	A	122	PHE	CB-CG-CD1	5.23	124.46	120.80
1	A	66	ASP	CB-CG-OD1	5.22	123.00	118.30
1	A	164	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	122	PHE	CB-CG-CD2	-5.18	117.18	120.80
1	A	101	ASP	CB-CG-OD1	5.12	122.91	118.30

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	3470	0	3519	93	0
2	A	19	0	14	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
6	A	27	0	12	0	0
7	A	457	0	0	13	1
All	All	3976	0	3545	93	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (93) 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:13:VAL:HG12	1:A:20:LEU:HD12	1.15	1.10
1:A:103:THR:HB	1:A:104:LYS:HE2	1.39	1.04
1:A:13:VAL:HG12	1:A:20:LEU:CD1	2.04	0.86
1:A:12:GLN:HG3	1:A:458:ALA:CB	2.07	0.84
1:A:103:THR:CB	1:A:104:LYS:HE2	2.07	0.84
1:A:102:LYS:HE2	1:A:183:GLY:HA2	1.60	0.82
1:A:12:GLN:HG3	1:A:458:ALA:HB3	1.62	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:GLN:HB2	1:A:15:GLU:CD	2.07	0.75
1:A:100:ILE:HD13	1:A:181:TYR:HB3	1.69	0.74
1:A:307:SER:OG	1:A:310:GLU:HG3	1.87	0.74
1:A:65:GLN:H	1:A:65:GLN:HE21	1.36	0.74
1:A:21:LYS:HE2	1:A:25[B]:MET:HE2	1.71	0.73
1:A:317:SER:HB2	1:A:318:PRO:HD2	1.72	0.72
1:A:2:ALA:HB1	7:A:914:HOH:O	1.89	0.71
1:A:13:VAL:CG1	1:A:20:LEU:HD12	2.09	0.70
1:A:44:VAL:HG21	1:A:225[A]:ILE:HD11	1.74	0.69
1:A:100:ILE:HD13	1:A:181:TYR:CB	2.23	0.69
1:A:21:LYS:HE2	1:A:25[B]:MET:CE	2.23	0.69
1:A:248[A]:ARG:NH2	7:A:484:HOH:O	2.26	0.68
1:A:256:LYS:HD2	1:A:262:TRP:CD1	2.29	0.68
1:A:102:LYS:HE2	1:A:183:GLY:CA	2.25	0.67
1:A:59[A]:LEU:HD12	1:A:203:LYS:O	1.96	0.66
1:A:21:LYS:HG3	1:A:25[B]:MET:HE3	1.77	0.66
1:A:103:THR:HB	1:A:104:LYS:CE	2.21	0.65
1:A:206:GLU:OE2	1:A:211:ARG:NH1	2.30	0.65
1:A:12:GLN:HB2	1:A:15:GLU:OE1	1.96	0.64
1:A:278:SER:OG	1:A:281:GLU:HG3	1.97	0.64
1:A:13:VAL:HG23	1:A:457:GLU:HG2	1.82	0.60
1:A:65:GLN:H	1:A:65:GLN:NE2	1.99	0.60
1:A:25[A]:MET:HG2	7:A:796:HOH:O	2.03	0.58
1:A:165:ASN:HB2	7:A:811:HOH:O	2.03	0.58
1:A:18:ARG:HB2	7:A:884:HOH:O	2.04	0.58
1:A:272:GLN:HB2	1:A:282:MET:HG2	1.85	0.57
1:A:256:LYS:HD2	1:A:262:TRP:NE1	2.19	0.57
1:A:77:TYR:HB3	7:A:802:HOH:O	2.06	0.54
1:A:167:SER:OG	1:A:170:GLU:HG3	2.09	0.53
1:A:225[B]:ILE:HD12	1:A:444:PHE:CE1	2.44	0.53
1:A:14:ALA:HA	1:A:20:LEU:HD13	1.90	0.51
1:A:11:VAL:O	1:A:458:ALA:N	2.32	0.51
1:A:101:ASP:O	1:A:102:LYS:C	2.50	0.49
1:A:232[B]:MET:HG3	1:A:404:TRP:HB2	1.94	0.49
1:A:368:HIS:HB2	1:A:398:ARG:CB	2.43	0.49
1:A:208:SER:OG	1:A:208:SER:O	2.30	0.48
1:A:27:ASN:HD22	1:A:32:SER:C	2.16	0.48
1:A:34:PRO:HG2	1:A:37:TYR:OH	2.14	0.48
1:A:63:VAL:HG12	7:A:709:HOH:O	2.13	0.47
1:A:218:PRO:HG3	1:A:360:LEU:HD12	1.95	0.47
1:A:287:GLU:OE1	1:A:333:LYS:NZ	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:TYR:CE2	1:A:333:LYS:HE3	2.50	0.47
1:A:115:LEU:HD12	1:A:130:MET:CE	2.45	0.46
1:A:13:VAL:CG2	1:A:457:GLU:HG2	2.45	0.46
1:A:232[B]:MET:HE1	1:A:239:HIS:HD2	1.80	0.46
1:A:296:ASN:HB2	1:A:297:PRO:HD2	1.98	0.46
1:A:354:GLU:HG2	1:A:355:ASN:N	2.31	0.46
1:A:16:HIS:O	1:A:17:PRO:C	2.54	0.46
1:A:12:GLN:OE1	1:A:15:GLU:OE1	2.34	0.46
1:A:353:PRO:HB2	1:A:355:ASN:OD1	2.16	0.46
1:A:310:GLU:O	1:A:314:GLN:N	2.44	0.45
1:A:317:SER:CB	1:A:318:PRO:HD2	2.37	0.45
1:A:356:MET:HE1	7:A:591:HOH:O	2.16	0.45
1:A:26:PHE:HZ	1:A:72:GLU:HB3	1.80	0.45
1:A:310:GLU:O	1:A:314:GLN:HB2	2.17	0.45
1:A:218:PRO:HG3	1:A:360:LEU:CD1	2.47	0.44
1:A:354:GLU:CG	1:A:355:ASN:N	2.80	0.44
1:A:232[B]:MET:CE	1:A:239:HIS:CD2	3.00	0.44
1:A:457:GLU:HG2	7:A:883:HOH:O	2.16	0.44
1:A:282:MET:HG3	1:A:336:TYR:CE1	2.53	0.44
1:A:286:THR:HG23	1:A:290:LEU:HD12	1.99	0.44
1:A:106:LEU:HD22	7:A:647:HOH:O	2.19	0.43
1:A:266:LEU:HA	1:A:266:LEU:HD23	1.81	0.43
1:A:22:LEU:HA	1:A:22:LEU:HD12	1.85	0.43
1:A:262:TRP:HZ3	7:A:726:HOH:O	2.01	0.43
1:A:357:VAL:HG12	1:A:395:GLN:HG3	2.01	0.43
1:A:59[A]:LEU:HD13	1:A:204:LEU:HA	2.00	0.43
1:A:153:CYS:O	1:A:157:VAL:HG23	2.18	0.43
1:A:399:LEU:C	1:A:399:LEU:HD12	2.39	0.42
1:A:232[B]:MET:HE2	1:A:239:HIS:CD2	2.54	0.42
1:A:33:ILE:HB	1:A:34:PRO:HD2	2.01	0.42
1:A:25[A]:MET:HB3	1:A:133:LEU:HD21	2.00	0.42
1:A:248[A]:ARG:HD3	1:A:266:LEU:HD22	2.00	0.42
1:A:177:LYS:NZ	7:A:815:HOH:O	2.51	0.42
1:A:248[A]:ARG:HD3	1:A:266:LEU:CD2	2.50	0.42
1:A:368:HIS:HB2	1:A:398:ARG:HB3	2.01	0.42
1:A:22:LEU:HG	1:A:70:ALA:HB2	2.00	0.42
1:A:49:GLU:HB2	1:A:398:ARG:NH2	2.34	0.41
1:A:85:ASN:HA	1:A:86:PRO:HD2	1.93	0.41
1:A:2:ALA:HB3	1:A:4:GLU:H	1.85	0.41
1:A:27:ASN:O	1:A:31:GLY:N	2.42	0.41
1:A:282:MET:HG3	1:A:336:TYR:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ARG:O	1:A:270:GLU:HB2	2.21	0.41
1:A:208:SER:HA	1:A:209:PRO:HA	1.80	0.41
1:A:61:MET:HE3	7:A:709:HOH:O	2.20	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:662:HOH:O	7:A:713:HOH:O[2_655]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	449/478 (94%)	434 (97%)	15 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	387/407 (95%)	358 (92%)	29 (8%)	13	2

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	9	ARG
1	A	10	ARG
1	A	18	ARG
1	A	21	LYS
1	A	25[A]	MET
1	A	25[B]	MET
1	A	35	LYS
1	A	39	ARG
1	A	43	ARG
1	A	65	GLN
1	A	75	LYS
1	A	94	SER
1	A	101	ASP
1	A	103	THR
1	A	104	LYS
1	A	105	PRO
1	A	125[A]	SER
1	A	125[B]	SER
1	A	208	SER
1	A	219	SER
1	A	233	ASN
1	A	258	LYS
1	A	282	MET
1	A	321	GLN
1	A	354	GLU
1	A	416	LYS
1	A	426	LYS
1	A	440	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	27	ASN
1	A	65	GLN
1	A	109	ASN
1	A	165	ASN
1	A	233	ASN
1	A	239	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NG1	A	459	3	18,19,19	1.41	2 (11%)	27,28,28	1.52	4 (14%)
6	ADP	A	463	3	24,29,29	0.99	1 (4%)	29,45,45	1.57	3 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NG1	A	459	3	-	4/10/31/31	0/1/1/1
6	ADP	A	463	3	-	3/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	459	NG1	C1-C2	3.33	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	459	NG1	P1-O1	2.56	1.64	1.59
6	A	463	ADP	C6-N6	-2.07	1.26	1.34

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	459	NG1	O5-C1-O1	-5.01	104.82	111.36
6	A	463	ADP	C5-C6-N1	-4.54	110.07	120.35
6	A	463	ADP	C5-C6-N6	4.53	127.24	120.35
6	A	463	ADP	C2-N1-C6	3.53	124.80	118.75
2	A	459	NG1	O1-C1-C2	2.60	113.10	108.40
2	A	459	NG1	C8-C7-N2	2.05	119.57	116.10
2	A	459	NG1	OP2-P1-OP3	-2.03	102.72	110.68

There are no chirality outliers.

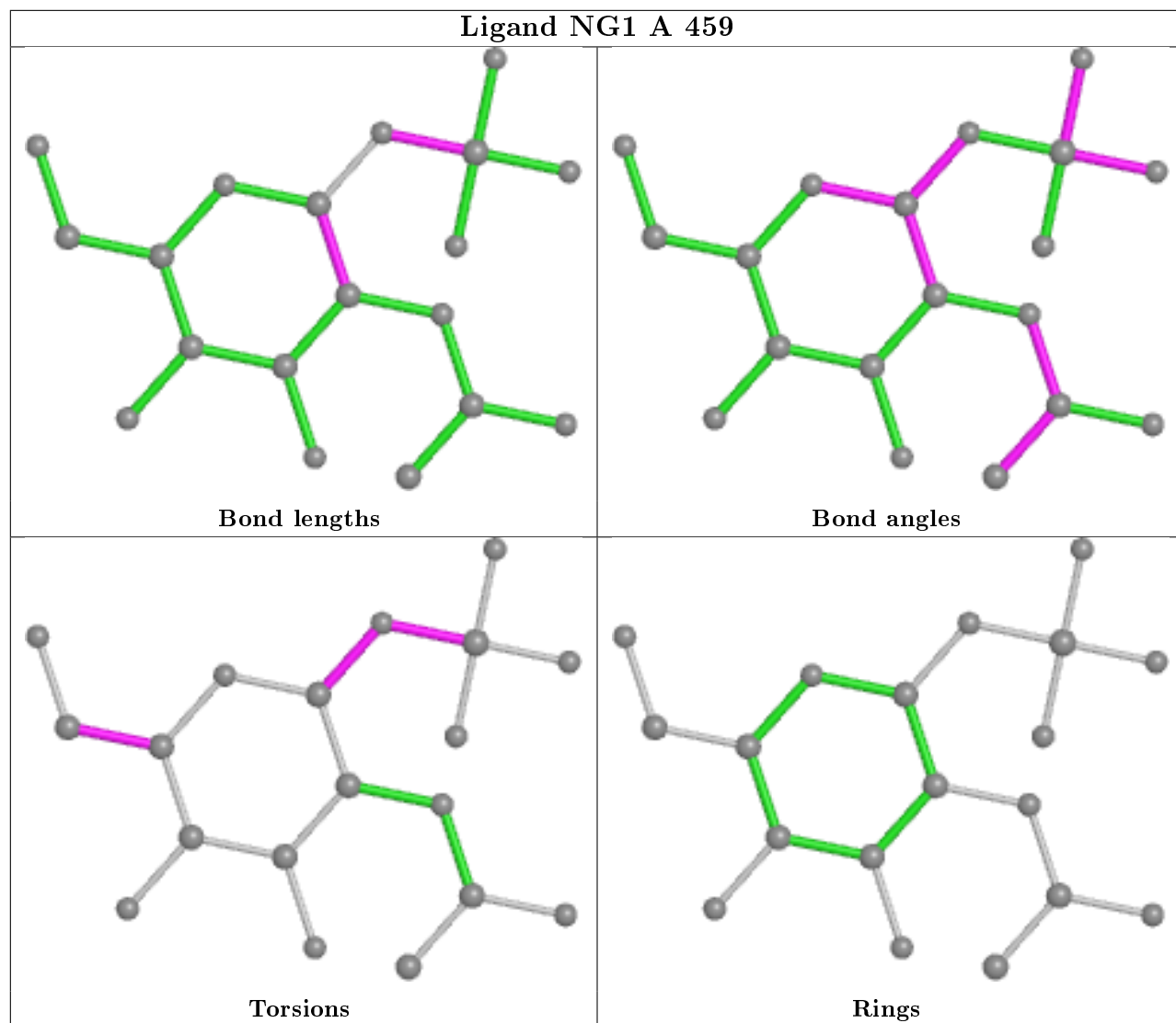
All (7) torsion outliers are listed below:

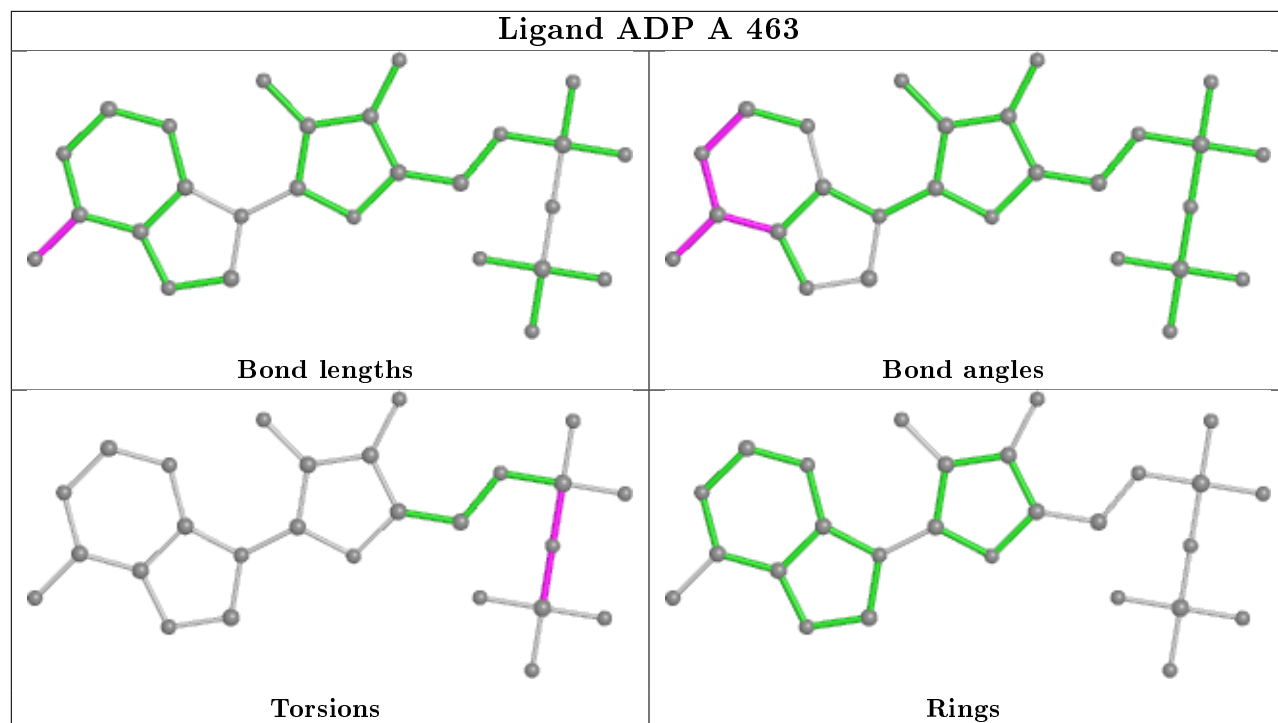
Mol	Chain	Res	Type	Atoms
2	A	459	NG1	O5-C5-C6-O6
6	A	463	ADP	PB-O3A-PA-O1A
2	A	459	NG1	O5-C1-O1-P1
6	A	463	ADP	PB-O3A-PA-O2A
6	A	463	ADP	PA-O3A-PB-O2B
2	A	459	NG1	C1-O1-P1-OP2
2	A	459	NG1	C1-O1-P1-OP3

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	446/478 (93%)	-0.32	5 (1%) 80 83	13, 23, 57, 91	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	103	THR	7.3
1	A	101	ASP	4.0
1	A	440	LYS	3.5
1	A	97	ASN	2.7
1	A	15	GLU	2.6

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
5	NA	A	462	1/1	0.94	0.17	56,56,56,56	0
2	NG1	A	459	19/19	0.97	0.10	10,13,27,36	0

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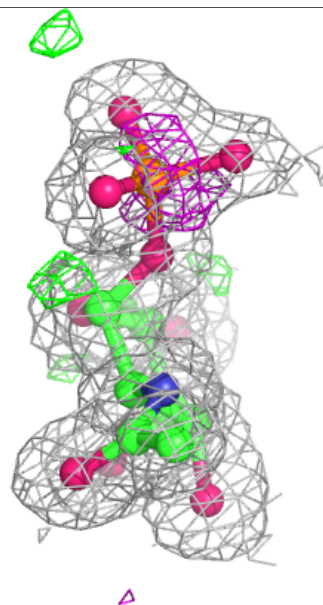
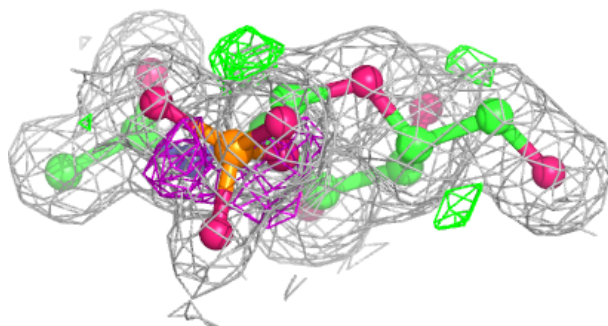
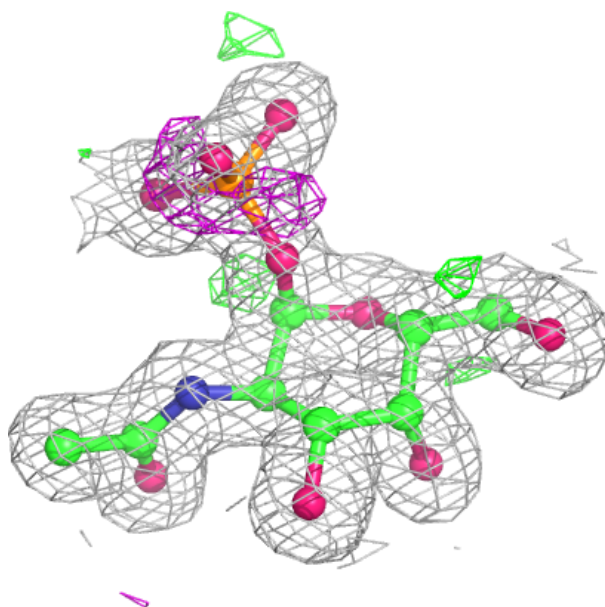
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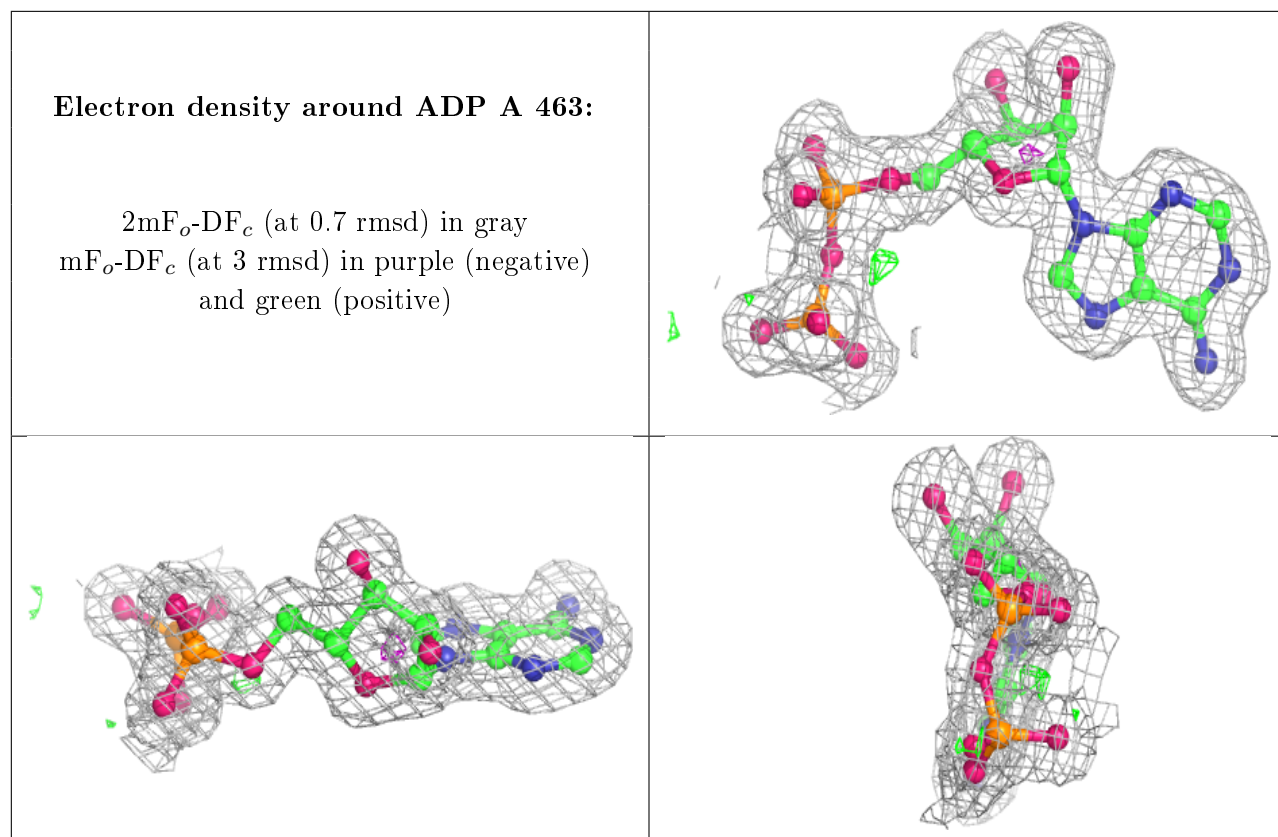
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	A	461	1/1	0.99	0.04	26,26,26,26	0
6	ADP	A	463	27/27	0.99	0.05	12,16,19,23	0
3	MG	A	460	1/1	1.00	0.10	9,9,9,9	0

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 NG1 A 459:

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.