



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 24, 2024 – 02:05 PM EDT

PDB ID : 5NZZ
Title : Crystal structure of phosphorylated p38aMAPK in complex with TAB1
Authors : Nichols, C.E.; De Nicola, G.F.
Deposited on : 2017-05-15
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.37.1
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) : 2.37.1

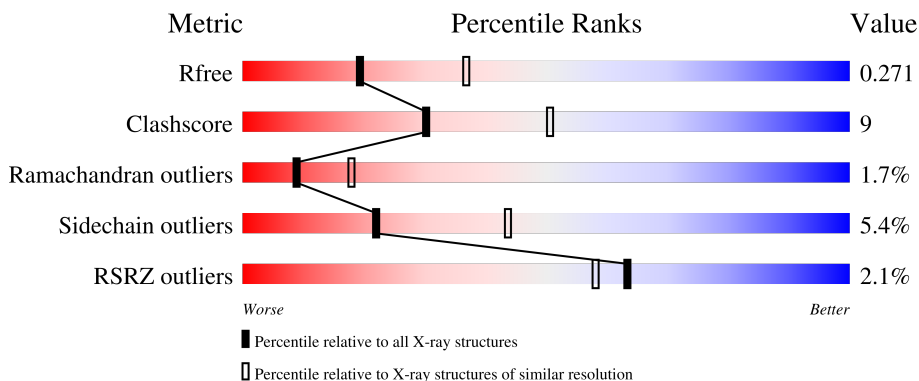
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	504	
1	B	504	
1	C	504	
1	D	504	
2	E	360	

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Mol	Chain	Length	Quality of chain
2	F	360	 76% 19% 5%
2	G	360	 75% 19% 6%
2	H	360	 11% 55% 29% 9% 6%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 22057 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 TGF-beta-activated kinase 1 and MAP3K7-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	378	Total	C	N	O	S	0	0	0
			2850	1790	492	558	10			
1	B	378	Total	C	N	O	S	0	0	0
			2856	1797	494	555	10			
1	C	367	Total	C	N	O	S	0	0	0
			2752	1730	477	535	10			
1	D	379	Total	C	N	O	S	0	0	0
			2843	1784	493	556	10			

- Molecule 2 is a protein called Mitogen-activated protein kinase 14.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
2	E	345	Total	C	N	O	P	S	0	1	0
			2743	1754	462	512	2	13			
2	F	343	Total	C	N	O	P	S	0	1	0
			2740	1756	463	506	2	13			
2	G	344	Total	C	N	O	P	S	0	2	0
			2723	1741	459	508	2	13			
2	H	337	Total	C	N	O	P	S	0	0	0
			2372	1501	411	448	2	10			

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

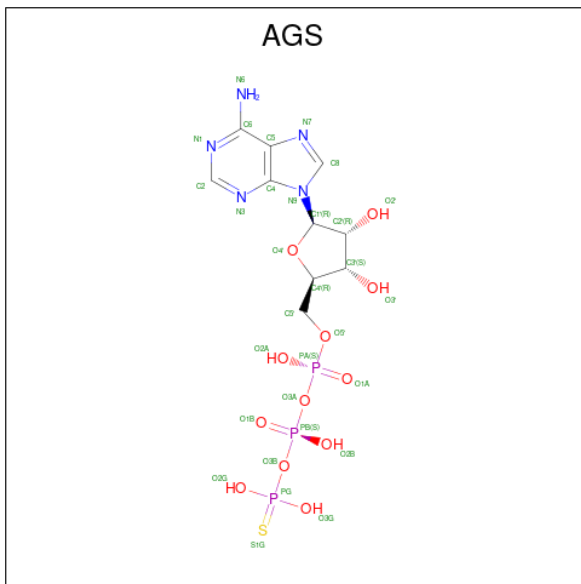
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total	0	0
			1		
3	B	1	Total	0	0
			1		
3	C	1	Total	0	0
			1		
3	D	1	Total	0	0
			1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total Ni 1 1	0	0
3	F	1	Total Ni 1 1	0	0

- Molecule 4 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (three-letter code: AGS) (formula: C₁₀H₁₆N₅O₁₂P₃S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	1	Total C N O P S 31 10 5 12 3 1	0	0
4	F	1	Total C N O P S 31 10 5 12 3 1	0	0
4	G	1	Total C N O P S 31 10 5 12 3 1	0	0
4	H	1	Total C N O P S 31 10 5 12 3 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	E	1	Total Mg 1 1	0	0
5	F	1	Total Mg 1 1	0	0
5	G	2	Total Mg 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	2	Total	Mg	0	0
			2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	O	0	0
			9	9		
6	B	5	Total	O	0	0
			5	5		
6	C	3	Total	O	0	0
			3	3		
6	D	3	Total	O	0	0
			3	3		
6	E	9	Total	O	0	0
			9	9		
6	F	10	Total	O	0	0
			10	10		
6	G	3	Total	O	0	0
			3	3		



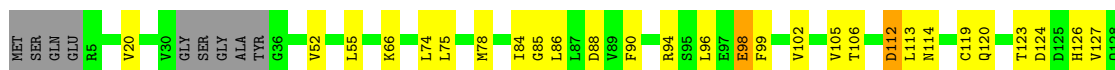
- Molecule 2: Mitogen-activated protein kinase 14

Chain F: 76% 19% 5%



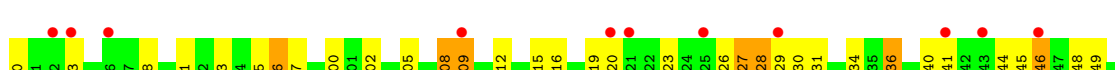
- Molecule 2: Mitogen-activated protein kinase 14

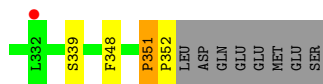
Chain G: 75% 19% . .



- Molecule 2: Mitogen-activated protein kinase 14

Chain H: 11% 55% 29% 9% 6%





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.88Å 107.63Å 144.20Å 89.26° 76.75° 84.28°	Depositor
Resolution (Å)	140.36 – 2.60 140.36 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.8 (140.36-2.60) 90.9 (140.36-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.62Å)	Xtrriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.219 , 0.271 0.219 , 0.271	Depositor DCC
R_{free} test set	5465 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	57.7	Xtrriage
Anisotropy	0.045	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.003 for -h,-k,-h+l	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	22057	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, AGS, PTR, MG, NI

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	0.51	0/2900	0.70	1/3938 (0.0%)
1	B	0.47	0/2906	0.69	0/3943
1	C	0.47	0/2799	0.67	0/3803
1	D	0.50	0/2894	0.65	0/3933
2	E	0.47	1/2778 (0.0%)	0.66	0/3783
2	F	0.50	0/2775	0.73	2/3774 (0.1%)
2	G	0.44	0/2757	0.66	2/3757 (0.1%)
2	H	0.43	0/2398	0.69	1/3292 (0.0%)
All	All	0.48	1/22207 (0.0%)	0.68	6/30223 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
1	D	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	223	PHE	C-N	6.63	1.46	1.34

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	173	ARG	NE-CZ-NH2	-8.54	116.03	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	205	ASP	CB-CG-OD1	6.00	123.69	118.30
2	F	149	ARG	NE-CZ-NH2	-5.48	117.56	120.30
2	F	67	ARG	NE-CZ-NH2	-5.39	117.61	120.30
2	G	229	ILE	CG1-CB-CG2	-5.30	99.75	111.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	59	GLU	Peptide
1	C	273	HIS	Peptide
1	C	274	GLY	Peptide
1	D	273	HIS	Peptide

5.2 Too-close contacts [\(i\)](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2850	0	2747	30	1
1	B	2856	0	2788	52	0
1	C	2752	0	2643	46	0
1	D	2843	0	2730	30	0
2	E	2743	0	2634	50	0
2	F	2740	0	2661	50	0
2	G	2723	0	2604	46	0
2	H	2372	0	1963	113	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	E	31	0	12	1	0
4	F	31	0	12	1	0
4	G	31	0	12	0	0
4	H	31	0	12	0	0
5	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	1	0	0	0	0
5	G	2	0	0	0	0
5	H	2	0	0	0	0
6	A	9	0	0	0	0
6	B	5	0	0	0	0
6	C	3	0	0	0	0
6	D	3	0	0	0	0
6	E	9	0	0	0	0
6	F	10	0	0	1	0
6	G	3	0	0	0	0
All	All	22057	0	20818	397	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 9.

The worst 5 of 397 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)
2:H:290:VAL:HG22	2:H:291:LEU:H	1.35	0.91
1:C:128:LYS:NZ	1:C:159:GLU:OE2	2.05	0.89
2:H:136:ARG:NH1	2:H:317:GLU:HA	1.89	0.88
2:H:220:ARG:HD2	2:H:221:THR:H	1.41	0.86
1:C:344:SER:O	1:C:346:GLY:N	2.14	0.81

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)
1:A:137:GLU:O	1:A:176:ASN:ND2[1_655]	2.18	0.02

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	372/504 (74%)	347 (93%)	22 (6%)	3 (1%)	19	39
1	B	372/504 (74%)	342 (92%)	27 (7%)	3 (1%)	19	39
1	C	359/504 (71%)	336 (94%)	16 (4%)	7 (2%)	8	15
1	D	373/504 (74%)	355 (95%)	12 (3%)	6 (2%)	9	19
2	E	340/360 (94%)	320 (94%)	16 (5%)	4 (1%)	13	27
2	F	338/360 (94%)	323 (96%)	14 (4%)	1 (0%)	41	64
2	G	340/360 (94%)	311 (92%)	26 (8%)	3 (1%)	17	35
2	H	327/360 (91%)	274 (84%)	32 (10%)	21 (6%)	1	1
All	All	2821/3456 (82%)	2608 (92%)	165 (6%)	48 (2%)	9	18

5 of 48 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	347	GLU
1	C	60	ASN
1	C	96	GLU
1	C	275	ALA
1	C	345	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	295/421 (70%)	278 (94%)	17 (6%)	20	40
1	B	300/421 (71%)	280 (93%)	20 (7%)	16	33
1	C	282/421 (67%)	267 (95%)	15 (5%)	22	45
1	D	294/421 (70%)	282 (96%)	12 (4%)	30	56
2	E	285/317 (90%)	273 (96%)	12 (4%)	30	55
2	F	287/317 (90%)	278 (97%)	9 (3%)	40	66
2	G	281/317 (89%)	275 (98%)	6 (2%)	53	77
2	H	189/317 (60%)	160 (85%)	29 (15%)	2	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2213/2952 (75%)	2093 (95%)	120 (5%)	22 44

5 of 120 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	190	LEU
2	H	228	HIS
2	E	293	SER
2	H	227	ASP
2	H	326	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	G	142	HIS
2	H	159	ASN
2	H	312	HIS
1	D	307	GLN
2	E	120	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

8 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TPO	E	180	2	8,10,11	1.66	2 (25%)	10,14,16	0.92	0
2	PTR	F	182	2	15,16,17	1.45	1 (6%)	19,22,24	0.86	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PTR	E	182	2	15,16,17	1.35	1 (6%)	19,22,24	0.65	0
2	PTR	G	182	2	15,16,17	1.32	1 (6%)	19,22,24	0.67	0
2	TPO	H	180	2	8,10,11	1.35	2 (25%)	10,14,16	1.64	2 (20%)
2	PTR	H	182	2	15,16,17	1.22	1 (6%)	19,22,24	1.14	2 (10%)
2	TPO	G	180	2	8,10,11	1.07	0	10,14,16	1.55	1 (10%)
2	TPO	F	180	2	8,10,11	1.10	0	10,14,16	1.57	3 (30%)

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	TPO	E	180	2	-	0/9/11/13	-
2	PTR	F	182	2	-	0/10/11/13	0/1/1/1
2	PTR	E	182	2	-	1/10/11/13	0/1/1/1
2	PTR	G	182	2	-	2/10/11/13	0/1/1/1
2	TPO	H	180	2	-	1/9/11/13	-
2	PTR	H	182	2	-	2/10/11/13	0/1/1/1
2	TPO	G	180	2	-	1/9/11/13	-
2	TPO	F	180	2	-	0/9/11/13	-

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	182	PTR	OH-CZ	-4.86	1.29	1.40
2	G	182	PTR	OH-CZ	-4.84	1.29	1.40
2	E	182	PTR	OH-CZ	-4.37	1.30	1.40
2	H	182	PTR	OH-CZ	-3.69	1.32	1.40
2	E	180	TPO	P-O1P	3.10	1.60	1.50

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	180	TPO	P-OG1-CB	-3.89	111.44	123.21
2	G	180	TPO	P-OG1-CB	-3.49	112.67	123.21
2	H	182	PTR	CB-CA-C	-3.22	105.43	111.47
2	F	180	TPO	O2P-P-O1P	-2.51	100.84	110.68
2	H	180	TPO	O3P-P-O1P	-2.34	101.50	110.68

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	182	PTR	O-C-CA-CB
2	G	182	PTR	CZ-OH-P-O1P
2	H	180	TPO	O-C-CA-CB
2	H	182	PTR	C-CA-CB-CG
2	H	182	PTR	N-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	182	PTR	1	0
2	H	180	TPO	2	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 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	AGS	H	500	5	26,33,33	0.76	1 (3%)	26,52,52	0.93	2 (7%)
4	AGS	F	401	5	26,33,33	1.00	2 (7%)	26,52,52	0.96	2 (7%)
4	AGS	E	401	5	26,33,33	0.85	1 (3%)	26,52,52	0.88	2 (7%)
4	AGS	G	500	5	26,33,33	0.88	1 (3%)	26,52,52	0.90	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
4	AGS	H	500	5	-	5/17/38/38	0/3/3/3
4	AGS	F	401	5	-	8/17/38/38	0/3/3/3
4	AGS	E	401	5	-	3/17/38/38	0/3/3/3
4	AGS	G	500	5	-	5/17/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	401	AGS	PA-O2A	-2.78	1.42	1.55
4	G	500	AGS	PG-S1G	2.61	1.96	1.90
4	H	500	AGS	PG-S1G	2.61	1.96	1.90
4	E	401	AGS	PG-S1G	2.59	1.96	1.90
4	F	401	AGS	PG-S1G	2.57	1.96	1.90

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	401	AGS	O3G-PG-O3B	3.08	114.94	104.64
4	H	500	AGS	PA-O3A-PB	-2.97	122.62	132.83
4	G	500	AGS	C5-C6-N6	2.42	124.03	120.35
4	F	401	AGS	C5-C6-N6	2.35	123.92	120.35
4	E	401	AGS	PA-O3A-PB	-2.22	125.21	132.83

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	401	AGS	PB-O3B-PG-O2G
4	F	401	AGS	PB-O3B-PG-O3G
4	F	401	AGS	C5'-O5'-PA-O1A
4	G	500	AGS	PB-O3B-PG-O2G
4	G	500	AGS	PB-O3B-PG-O3G

There are no ring outliers.

2 monomers are involved in 2 short contacts:

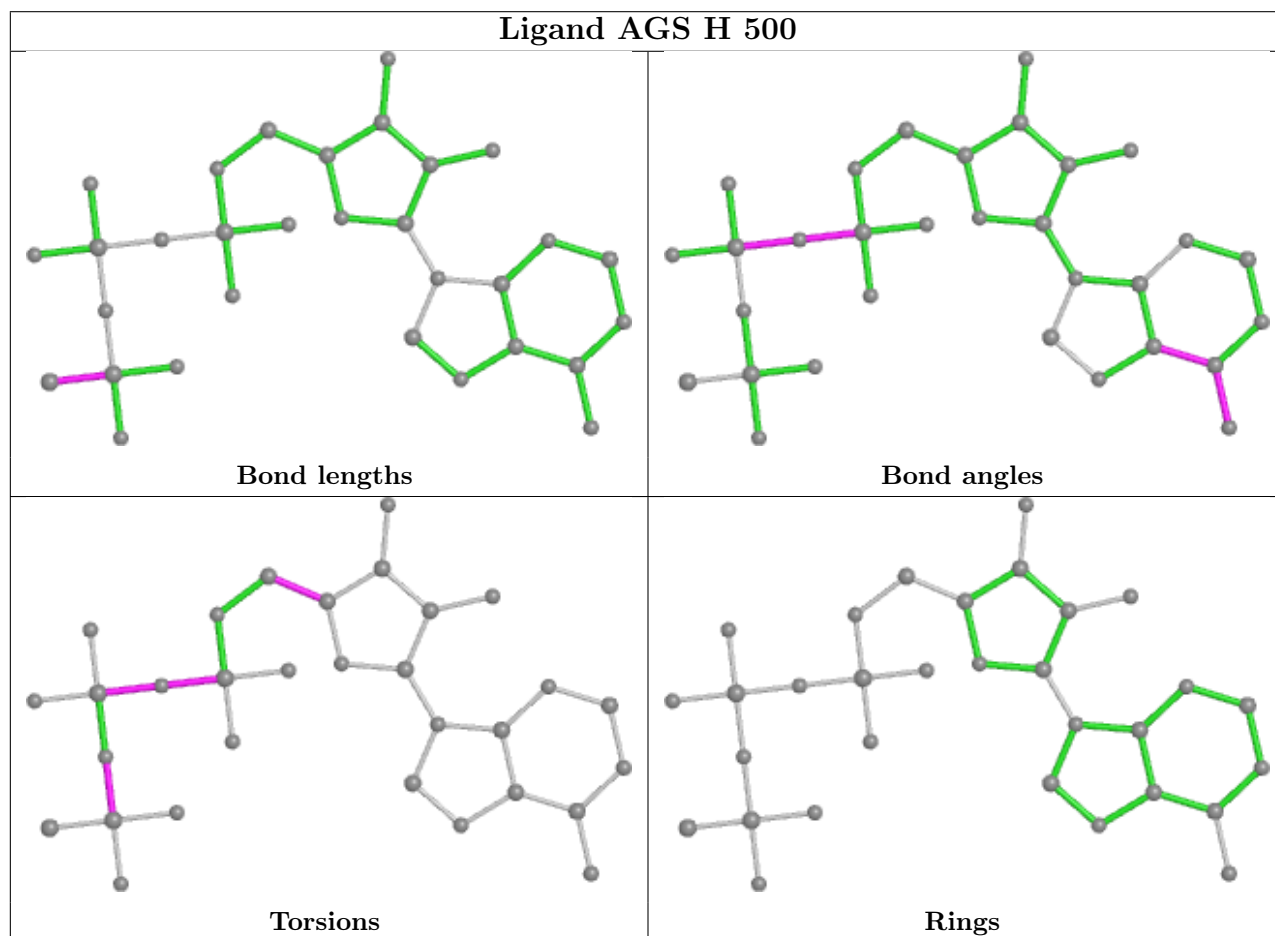
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	401	AGS	1	0

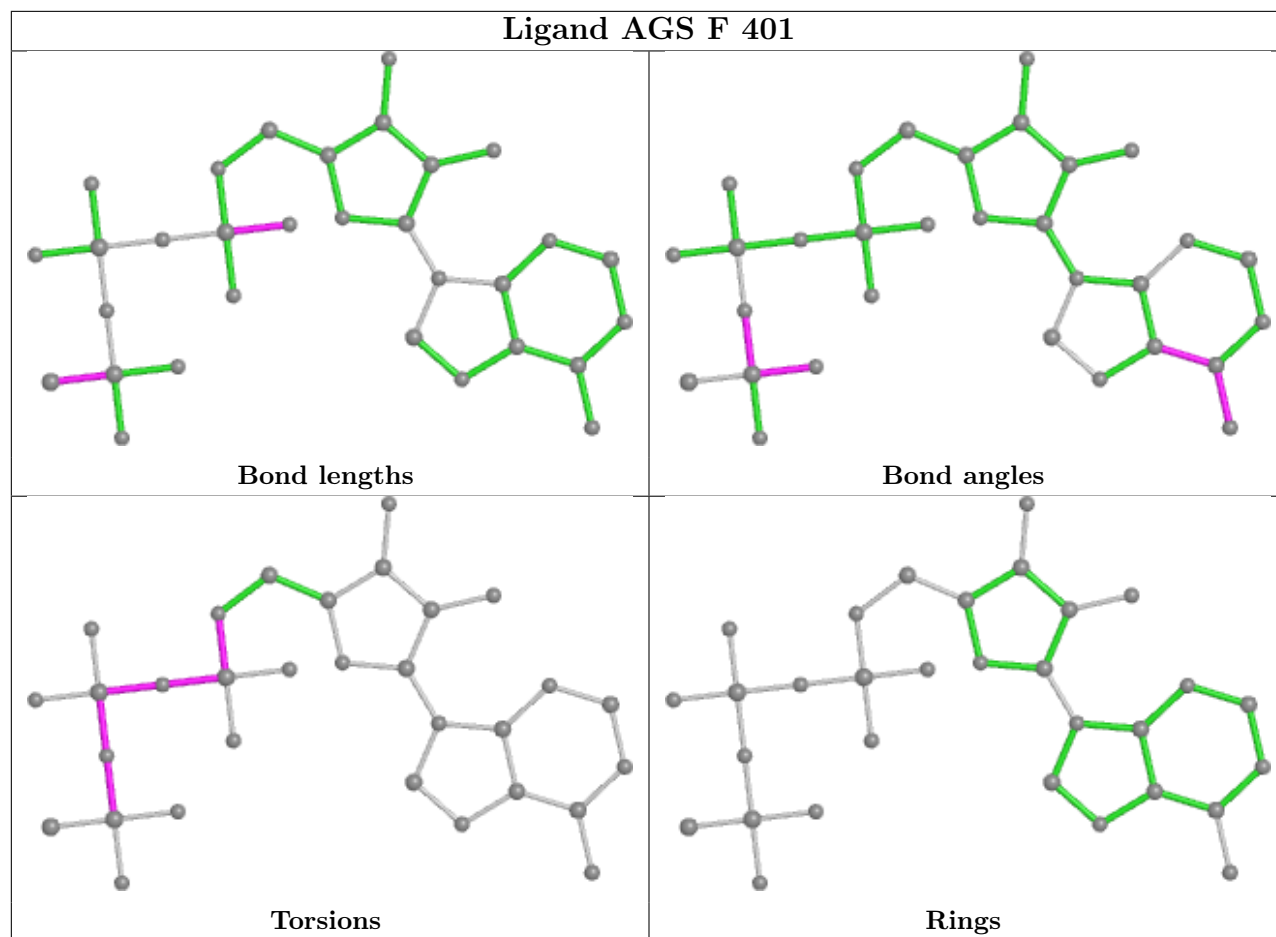
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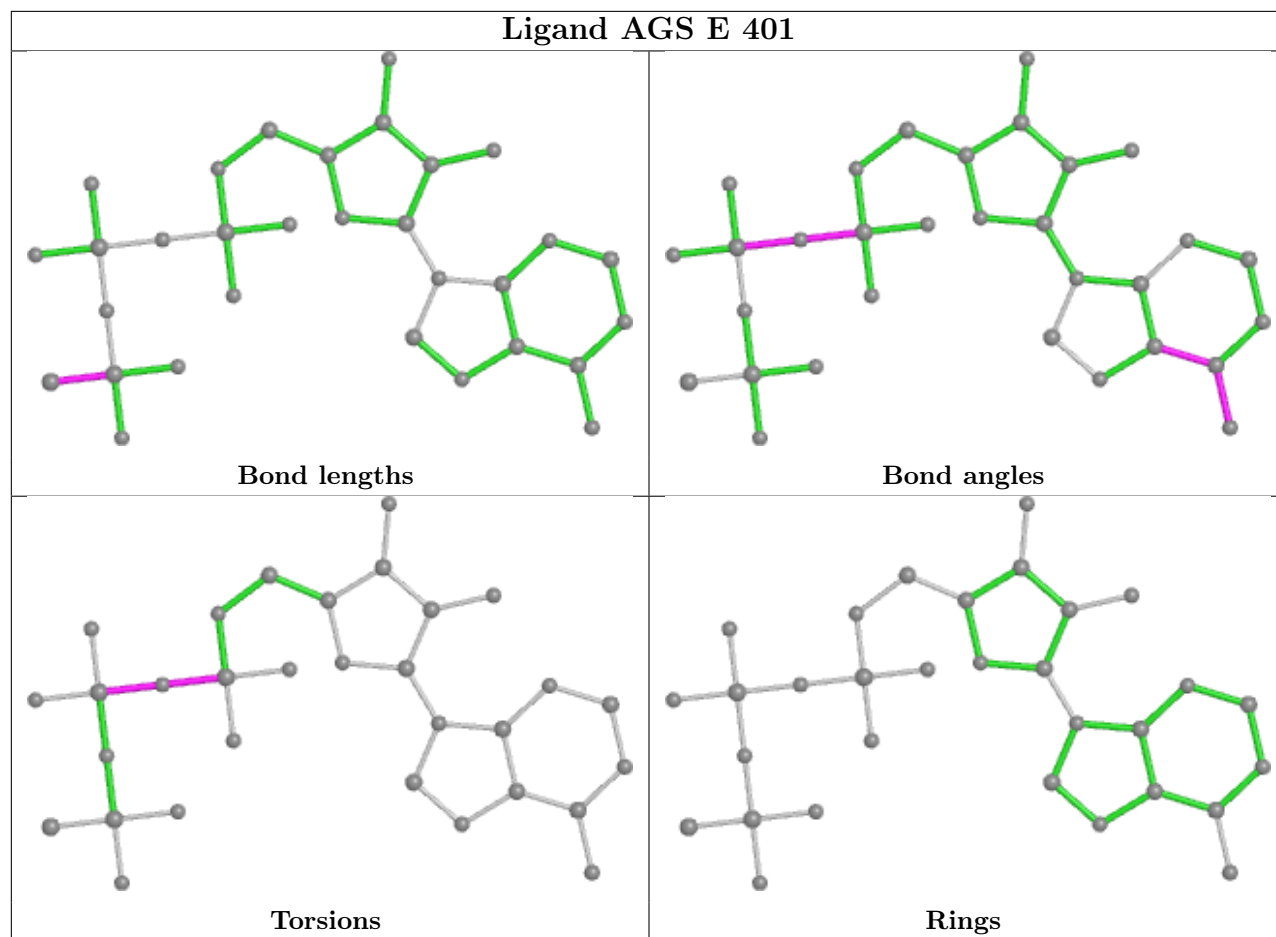
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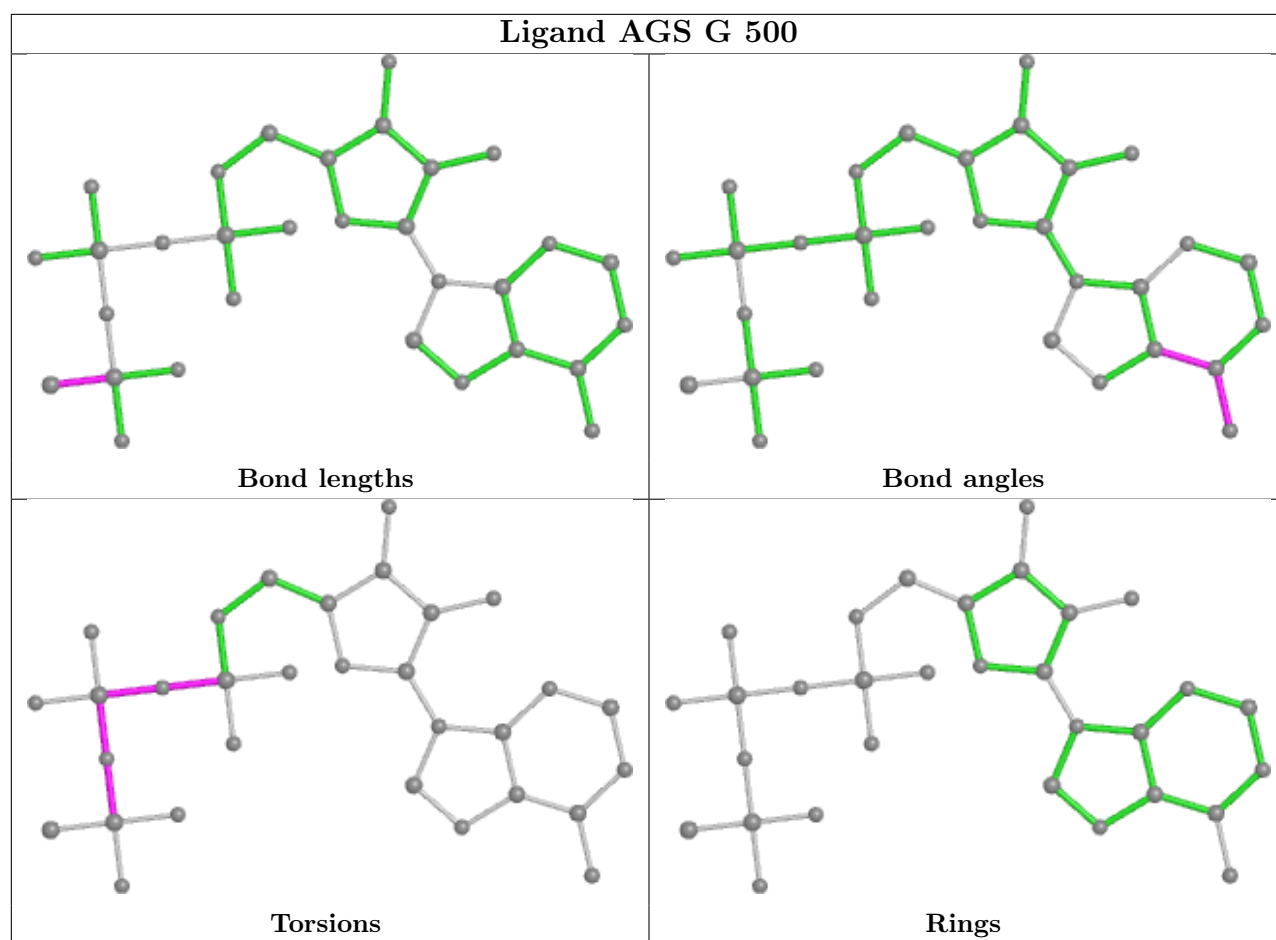
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	401	AGS	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	378/504 (75%)	-0.24	1 (0%) 94 93	32, 50, 82, 116	0
1	B	378/504 (75%)	-0.23	5 (1%) 77 73	33, 53, 91, 110	0
1	C	367/504 (72%)	-0.11	10 (2%) 54 48	32, 54, 98, 139	0
1	D	379/504 (75%)	-0.28	1 (0%) 94 93	33, 52, 83, 107	0
2	E	343/360 (95%)	-0.29	1 (0%) 94 93	35, 55, 78, 96	0
2	F	341/360 (94%)	-0.26	0 100 100	32, 51, 70, 98	0
2	G	342/360 (95%)	-0.18	2 (0%) 89 88	42, 62, 85, 101	0
2	H	335/360 (93%)	0.51	39 (11%) 4 3	69, 97, 119, 132	0
All	All	2863/3456 (82%)	-0.14	59 (2%) 63 58	32, 57, 102, 139	0

The worst 5 of 59 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	388	VAL	9.0
2	H	309	ALA	6.8
2	H	166	ILE	4.4
2	H	304	ALA	4.1
1	B	388	VAL	4.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	PTR	H	182	16/17	0.89	0.18	78,82,93,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PTR	G	182	16/17	0.90	0.14	52,65,80,84	0
2	TPO	H	180	11/12	0.93	0.14	69,85,93,95	0
2	PTR	F	182	16/17	0.94	0.13	42,51,66,69	0
2	PTR	E	182	16/17	0.95	0.14	51,59,71,73	0
2	TPO	G	180	11/12	0.98	0.13	37,46,49,51	0
2	TPO	F	180	11/12	0.98	0.13	23,40,42,46	0
2	TPO	E	180	11/12	0.99	0.14	24,39,44,51	0

6.3 Carbohydrates [i](#)

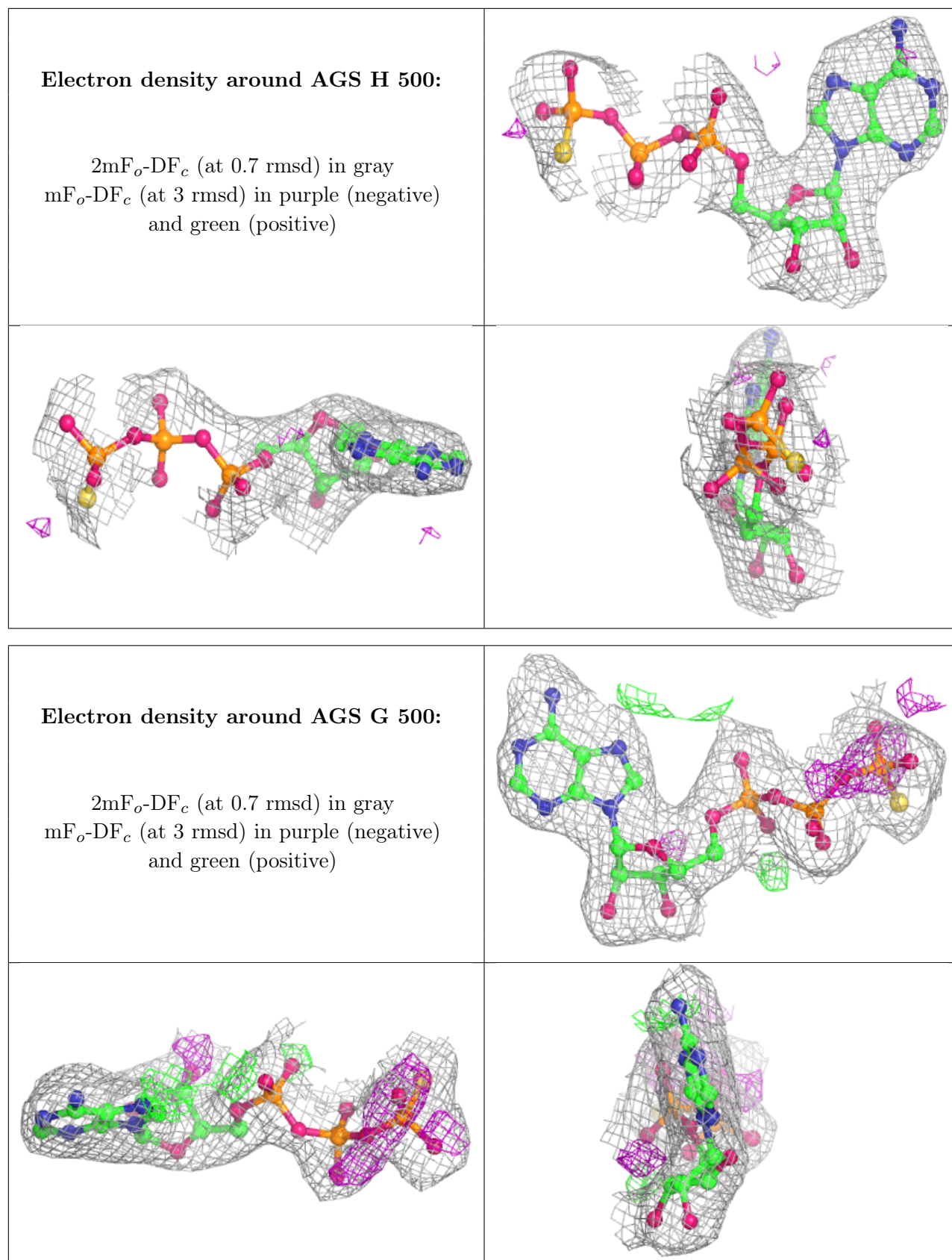
There are no monosaccharides in this entry.

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.

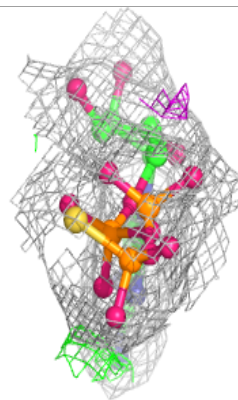
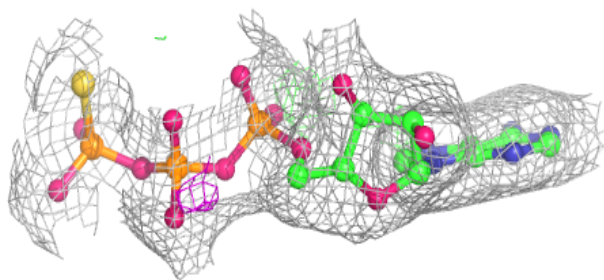
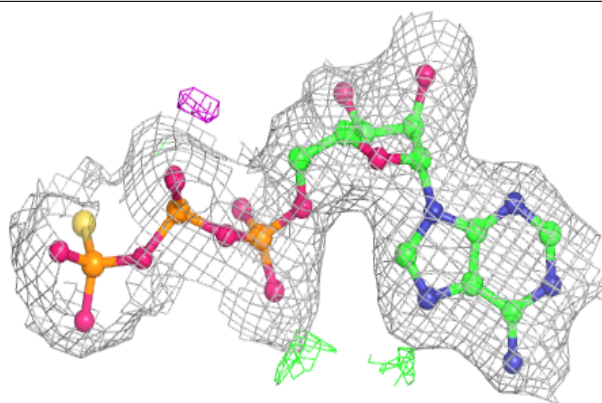
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	MG	G	501	1/1	0.78	0.11	49,49,49,49	0
3	NI	A	601	1/1	0.92	0.19	84,84,84,84	0
5	MG	H	502	1/1	0.92	0.12	82,82,82,82	0
4	AGS	H	500	31/31	0.93	0.16	75,90,102,114	0
5	MG	F	402	1/1	0.93	0.18	45,45,45,45	0
3	NI	E	403	1/1	0.94	0.11	87,87,87,87	0
4	AGS	G	500	31/31	0.94	0.15	38,55,67,79	0
3	NI	B	601	1/1	0.95	0.15	86,86,86,86	1
3	NI	D	601	1/1	0.95	0.14	98,98,98,98	0
5	MG	E	402	1/1	0.96	0.25	45,45,45,45	0
3	NI	C	601	1/1	0.96	0.14	97,97,97,97	0
3	NI	F	403	1/1	0.97	0.06	97,97,97,97	0
4	AGS	E	401	31/31	0.97	0.13	26,44,64,77	0
5	MG	G	502	1/1	0.97	0.16	68,68,68,68	0
4	AGS	F	401	31/31	0.97	0.15	32,45,66,87	0
5	MG	H	501	1/1	0.98	0.10	90,90,90,90	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.

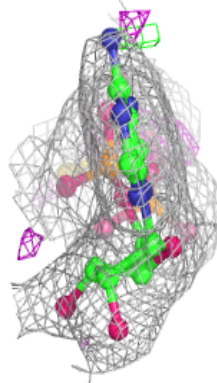
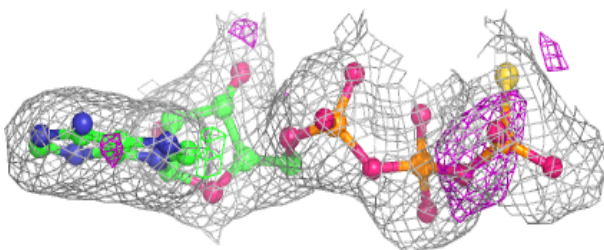
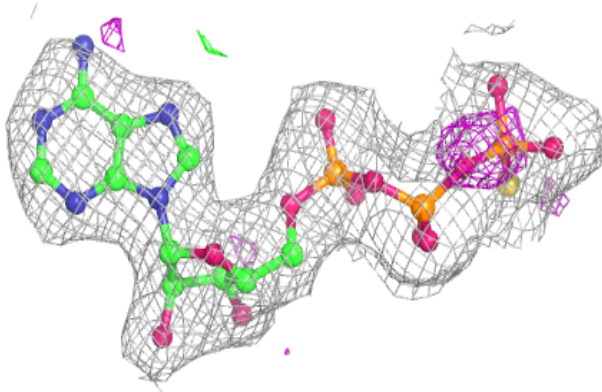


Electron density around AGS E 401:

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

**Electron density around AGS 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.