



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 04:42 pm BST

PDB ID : 3O82
Title : Structure of BasE N-terminal domain from *Acinetobacter baumannii* bound to 5'-O-[N-(2,3-dihydroxybenzoyl)sulfamoyl] adenosine
Authors : Drake, E.J.; Duckworth, B.P.; Neres, J.; Aldrich, C.C.; Gulick, A.M.
Deposited on : 2010-08-02
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 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 : 2.11
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) : 2.11

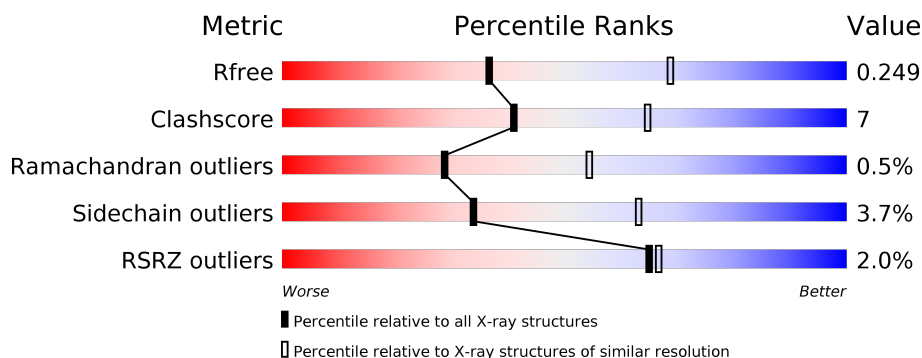
1 Overall quality at a glance

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))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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 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	544	<div> <div>2%</div> <div> <div></div> <div>68%</div> <div>11%</div> <div>•</div> <div>20%</div> </div> </div>
1	B	544	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>16%</div> <div>•</div> <div>20%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6881 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.

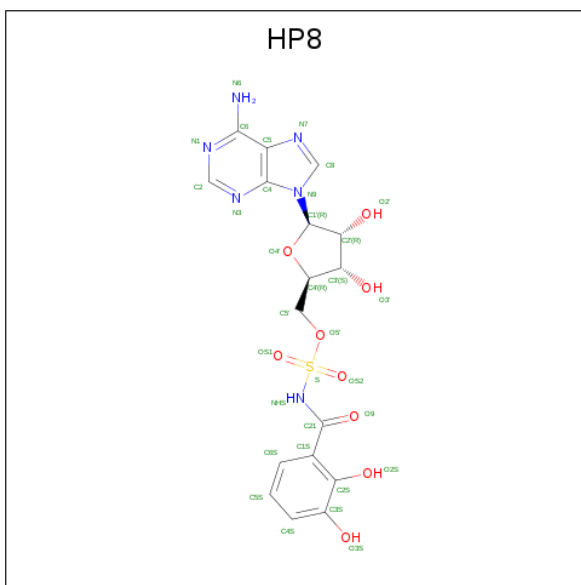
- Molecule 1 is a protein called Peptide arylation enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3378	2153	578	633	14			
1	B	435	Total	C	N	O	S	0	0	0
			3384	2155	579	636	14			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP B2HVG8
A	0	HIS	-	EXPRESSION TAG	UNP B2HVG8
A	45	LEU	PRO	ENGINEERED MUTATION	UNP B2HVG8
A	68	THR	SER	SEE REMARK 999	UNP B2HVG8
A	149	ASP	GLU	SEE REMARK 999	UNP B2HVG8
A	180	PHE	LEU	SEE REMARK 999	UNP B2HVG8
A	226	GLY	ASP	SEE REMARK 999	UNP B2HVG8
A	329	LEU	ILE	SEE REMARK 999	UNP B2HVG8
A	378	ARG	LYS	SEE REMARK 999	UNP B2HVG8
B	-1	GLY	-	EXPRESSION TAG	UNP B2HVG8
B	0	HIS	-	EXPRESSION TAG	UNP B2HVG8
B	45	LEU	PRO	ENGINEERED MUTATION	UNP B2HVG8
B	68	THR	SER	SEE REMARK 999	UNP B2HVG8
B	149	ASP	GLU	SEE REMARK 999	UNP B2HVG8
B	180	PHE	LEU	SEE REMARK 999	UNP B2HVG8
B	226	GLY	ASP	SEE REMARK 999	UNP B2HVG8
B	329	LEU	ILE	SEE REMARK 999	UNP B2HVG8
B	378	ARG	LYS	SEE REMARK 999	UNP B2HVG8

- Molecule 2 is 5'-O-{[(2,3-dihydroxyphenyl)carbonyl]sulfamoyl}adenosine (three-letter code: HP8) (formula: C₁₇H₁₈N₆O₉S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			33	17	6	9	1		
2	B	1	Total	C	N	O	S	0	0
			33	17	6	9	1		

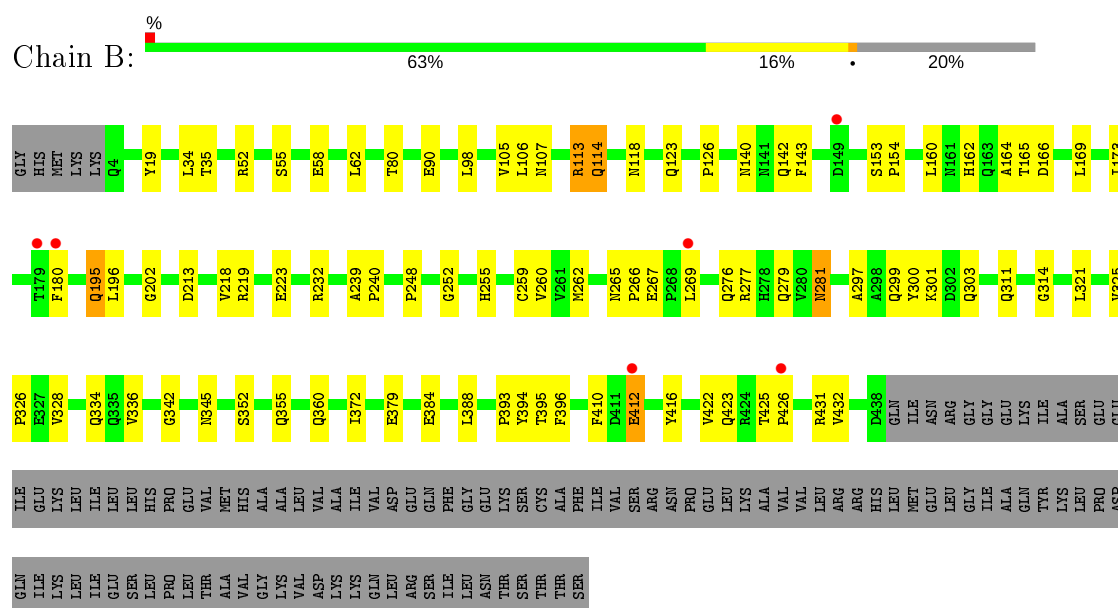
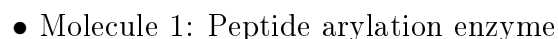
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		
4	B	24	Total	O	0	0
			24	24		

- Molecule 1: Peptide arylation enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.45 Å 145.26 Å 149.06 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.70 39.57 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.5 (40.00-2.70) 97.5 (39.57-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.69 Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.200 , 0.251 0.202 , 0.249	Depositor DCC
R_{free} test set	1969 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.547	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6881	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: CA, HP8

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	0.85	2/3457 (0.1%)	0.84	2/4710 (0.0%)
1	B	0.85	0/3462	0.85	0/4715
All	All	0.85	2/6919 (0.0%)	0.84	2/9425 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	14	GLU	CG-CD	5.67	1.60	1.51
1	A	375	GLU	CG-CD	5.11	1.59	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	302	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	232	ARG	NE-CZ-NH1	5.20	122.90	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3378	0	3284	35	0
1	B	3384	0	3297	60	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	33	0	17	0	0
2	B	33	0	17	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	25	0	0	0	1
4	B	24	0	0	2	0
All	All	6881	0	6615	95	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 7.

All (95) 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:B:360:GLN:HE21	1:B:432:VAL:H	1.11	0.93
1:B:232:ARG:H	1:B:281:ASN:HD21	1.20	0.88
1:B:106:LEU:HD11	1:B:196:LEU:HD21	1.63	0.81
1:B:360:GLN:NE2	1:B:432:VAL:H	1.81	0.79
1:A:352:SER:H	1:A:355:GLN:HE21	1.29	0.78
1:A:423:GLN:HE21	1:A:431:ARG:HH11	1.35	0.74
1:A:106:LEU:HD11	1:A:196:LEU:HD21	1.69	0.74
1:B:52:ARG:NH2	1:B:90:GLU:OE2	2.22	0.73
1:B:352:SER:H	1:B:355:GLN:HE21	1.38	0.72
1:B:213:ASP:HB2	1:B:394:TYR:HA	1.71	0.71
1:A:381:PRO:HG2	1:A:384:GLU:HG3	1.74	0.69
1:B:232:ARG:H	1:B:281:ASN:ND2	1.89	0.69
1:A:114:GLN:H	1:A:114:GLN:HE21	1.41	0.67
1:A:213:ASP:HB2	1:A:394:TYR:HA	1.78	0.66
1:A:118:ASN:HD22	1:A:143:PHE:HE1	1.42	0.65
1:A:195:GLN:HE21	1:A:196:LEU:H	1.46	0.63
1:B:260:VAL:HG12	1:B:262:MET:HE3	1.81	0.63
1:A:77:LYS:NZ	1:A:184:SER:OG	2.31	0.62
1:B:352:SER:H	1:B:355:GLN:NE2	2.01	0.57
1:B:423:GLN:HE21	1:B:431:ARG:HH11	1.52	0.57
1:A:75:LEU:HA	1:A:79:ASP:OD2	2.04	0.57
1:B:55:SER:OG	1:B:58:GLU:HG2	2.04	0.57
1:B:52:ARG:HH22	1:B:90:GLU:CD	2.09	0.56
1:A:113:ARG:NH2	1:A:136:GLU:CB	2.69	0.55
1:B:202:GLY:HA2	4:B:561:HOH:O	2.06	0.54
1:A:114:GLN:H	1:A:114:GLN:NE2	2.05	0.54
1:B:412:GLU:CD	1:B:412:GLU:H	2.11	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:VAL:HG12	1:B:262:MET:CE	2.39	0.52
1:B:372:ILE:HD13	1:B:422:VAL:HG23	1.92	0.52
1:A:195:GLN:NE2	1:A:196:LEU:H	2.08	0.52
1:A:180:PHE:O	1:A:180:PHE:CD2	2.64	0.51
1:A:352:SER:H	1:A:355:GLN:NE2	2.03	0.51
1:B:195:GLN:HE21	1:B:196:LEU:H	1.59	0.51
1:A:150:VAL:O	1:A:150:VAL:HG12	2.11	0.51
1:B:248:PRO:HD3	1:B:311:GLN:HE22	1.74	0.51
1:B:118:ASN:ND2	1:B:143:PHE:HE1	2.09	0.51
1:B:113:ARG:HB3	1:B:114:GLN:HE21	1.76	0.50
1:A:237:LEU:HB2	1:A:244:MET:SD	2.51	0.50
1:B:342:GLY:HA3	1:B:395:THR:HA	1.94	0.50
1:B:239:ALA:N	1:B:240:PRO:HD2	2.27	0.50
1:B:260:VAL:CG1	1:B:262:MET:CE	2.90	0.49
1:B:219:ARG:HG2	1:B:255:HIS:CE1	2.47	0.49
1:B:160:LEU:HD13	1:B:169:LEU:HD23	1.95	0.49
1:B:297:ALA:HB3	1:B:328:VAL:HG12	1.95	0.48
1:A:195:GLN:HE21	1:A:196:LEU:N	2.11	0.48
1:B:410:PHE:CE2	1:B:416:TYR:HB2	2.49	0.47
1:B:166:ASP:OD1	4:B:554:HOH:O	2.20	0.47
1:B:314:GLY:HA2	2:B:602:HP8:H6S	1.96	0.47
1:B:98:LEU:HD13	1:B:105:VAL:HB	1.97	0.46
1:B:321:LEU:HD23	1:B:321:LEU:HA	1.74	0.46
1:A:311:GLN:HA	1:A:334:GLN:O	2.17	0.45
1:B:232:ARG:N	1:B:281:ASN:HD21	2.01	0.45
1:B:336:VAL:HA	1:B:345:ASN:O	2.17	0.45
1:A:181:VAL:CG1	1:A:182:ASP:N	2.81	0.44
1:A:335:GLN:HB2	1:A:347:THR:HG23	1.99	0.44
1:B:80:THR:OG1	1:B:126:PRO:HA	2.17	0.44
1:B:162:HIS:HB2	1:B:166:ASP:HB2	1.99	0.44
1:B:62:LEU:HD22	1:B:173:ILE:HG22	2.00	0.44
1:A:285:LEU:HD11	1:A:293:TRP:CE2	2.53	0.44
1:B:34:LEU:HA	1:B:34:LEU:HD12	1.50	0.44
1:B:140:ASN:CG	1:B:142:GLN:HE21	2.21	0.43
1:B:55:SER:H	1:B:58:GLU:HG3	1.83	0.43
1:B:218:VAL:HG13	1:B:252:GLY:HA2	2.00	0.43
1:B:388:LEU:HD13	1:B:422:VAL:HG21	2.00	0.43
1:B:352:SER:N	1:B:355:GLN:HE21	2.11	0.43
1:B:113:ARG:NH1	1:B:113:ARG:HG3	2.32	0.43
1:A:114:GLN:N	1:A:114:GLN:HE21	2.12	0.43
1:B:360:GLN:NE2	1:B:432:VAL:N	2.58	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLU:CD	1:A:277:ARG:HH12	2.22	0.42
1:A:113:ARG:HH22	1:A:136:GLU:CB	2.32	0.42
1:A:219:ARG:HG2	1:A:255:HIS:CE1	2.54	0.42
1:B:265:ASN:HB2	1:B:266:PRO:CD	2.50	0.42
1:B:34:LEU:O	1:B:35:THR:C	2.58	0.42
1:B:232:ARG:HA	1:B:259:CYS:O	2.20	0.42
1:A:239:ALA:N	1:A:240:PRO:HD2	2.34	0.42
1:A:243:PHE:HA	1:A:247:SER:HB2	2.01	0.42
1:A:300:TYR:O	1:A:301:LYS:C	2.58	0.42
1:B:266:PRO:O	1:B:267:GLU:C	2.56	0.42
1:A:181:VAL:HG12	1:A:182:ASP:N	2.35	0.42
1:A:19:TYR:CE1	1:A:393:PRO:HA	2.55	0.41
1:B:311:GLN:HA	1:B:334:GLN:O	2.20	0.41
1:B:153:SER:HA	1:B:154:PRO:HD2	1.92	0.41
1:B:260:VAL:CG1	1:B:262:MET:HE1	2.50	0.41
1:B:425:THR:HB	1:B:426:PRO:HD2	2.02	0.41
1:A:203:THR:HA	1:A:204:PRO:HD3	1.98	0.41
1:A:180:PHE:HD2	1:A:180:PHE:O	2.02	0.41
1:B:262:MET:HA	1:B:262:MET:HE2	2.03	0.41
1:B:19:TYR:CE1	1:B:393:PRO:HA	2.55	0.41
1:A:323:ARG:NH2	1:A:354:GLU:OE2	2.53	0.41
1:B:106:LEU:HD12	1:B:107:ASN:N	2.36	0.41
1:B:300:TYR:O	1:B:301:LYS:C	2.60	0.41
1:A:267:GLU:HA	1:A:268:PRO:HD3	1.92	0.40
1:B:276:GLN:HB2	1:B:303:GLN:O	2.21	0.40
1:B:325:VAL:HB	1:B:326:PRO:HD3	2.02	0.40
1:B:423:GLN:NE2	1:B:431:ARG:HH11	2.16	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)
1:B:279:GLN:OE1	4:A:554:HOH:O[1_655]	1.97	0.23

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	433/544 (80%)	406 (94%)	25 (6%)	2 (0%)	29	54
1	B	433/544 (80%)	411 (95%)	20 (5%)	2 (0%)	29	54
All	All	866/1088 (80%)	817 (94%)	45 (5%)	4 (0%)	29	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	164	ALA
1	B	165	THR
1	A	343	LEU
1	A	202	GLY

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	363/472 (77%)	350 (96%)	13 (4%)	35	64
1	B	365/472 (77%)	351 (96%)	14 (4%)	33	62
All	All	728/944 (77%)	701 (96%)	27 (4%)	34	63

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	ARG
1	A	113	ARG
1	A	114	GLN
1	A	134	GLN
1	A	180	PHE
1	A	195	GLN
1	A	200	SER
1	A	201	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	299	GLN
1	A	320	SER
1	A	370	ILE
1	A	396	PHE
1	A	436	ILE
1	B	113	ARG
1	B	114	GLN
1	B	123	GLN
1	B	180	PHE
1	B	195	GLN
1	B	223	GLU
1	B	269	LEU
1	B	277	ARG
1	B	281	ASN
1	B	299	GLN
1	B	379	GLU
1	B	384	GLU
1	B	396	PHE
1	B	412	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	GLN
1	A	195	GLN
1	A	270	ASN
1	A	303	GLN
1	A	355	GLN
1	A	423	GLN
1	B	41	HIS
1	B	53	GLN
1	B	114	GLN
1	B	118	ASN
1	B	134	GLN
1	B	162	HIS
1	B	195	GLN
1	B	270	ASN
1	B	281	ASN
1	B	303	GLN
1	B	335	GLN
1	B	355	GLN
1	B	360	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	423	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 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	HP8	B	602	-	33,36,36	1.49	4 (12%)	39,54,54	1.89	8 (20%)
2	HP8	A	601	-	33,36,36	1.42	3 (9%)	39,54,54	1.82	12 (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	HP8	B	602	-	-	2/14/35/35	0/4/4/4
2	HP8	A	601	-	-	2/14/35/35	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HP8	C21-NHS	-5.46	1.32	1.39
2	B	602	HP8	C21-NHS	-4.90	1.33	1.39
2	B	602	HP8	OS1-S	4.44	1.46	1.42
2	A	601	HP8	C3S-C2S	-2.49	1.37	1.40
2	B	602	HP8	O5'-S	-2.29	1.54	1.59
2	A	601	HP8	O5'-S	-2.19	1.55	1.59
2	B	602	HP8	C3S-C2S	-2.00	1.37	1.40

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	HP8	C1S-C2S-C3S	-5.34	116.61	119.99
2	B	602	HP8	C21-NHS-S	-4.92	117.19	123.38
2	A	601	HP8	C4S-C3S-C2S	4.45	124.62	120.06
2	B	602	HP8	N3-C2-N1	-4.15	122.19	128.68
2	A	601	HP8	N3-C2-N1	-4.04	122.36	128.68
2	B	602	HP8	C4S-C3S-C2S	3.97	124.13	120.06
2	A	601	HP8	C4-C5-N7	-3.81	105.43	109.40
2	A	601	HP8	C3'-C2'-C1'	3.44	106.16	100.98
2	A	601	HP8	C1S-C21-NHS	2.98	118.52	114.34
2	A	601	HP8	C21-NHS-S	-2.94	119.67	123.38
2	A	601	HP8	C1S-C2S-C3S	-2.74	118.25	119.99
2	A	601	HP8	C5S-C4S-C3S	-2.69	116.64	120.05
2	B	602	HP8	C1S-C21-NHS	2.66	118.08	114.34
2	B	602	HP8	C6S-C1S-C2S	2.65	121.34	118.74
2	B	602	HP8	C3'-C2'-C1'	2.42	104.62	100.98
2	A	601	HP8	O5'-S-NHS	2.33	112.10	105.60
2	A	601	HP8	O5'-S-OS2	2.25	112.47	105.59
2	B	602	HP8	O9-C21-NHS	2.22	123.81	121.08
2	A	601	HP8	O4'-C1'-C2'	-2.15	103.79	106.93
2	A	601	HP8	OS2-S-OS1	-2.10	117.48	120.76

There are no chirality outliers.

All (4) torsion outliers are listed below:

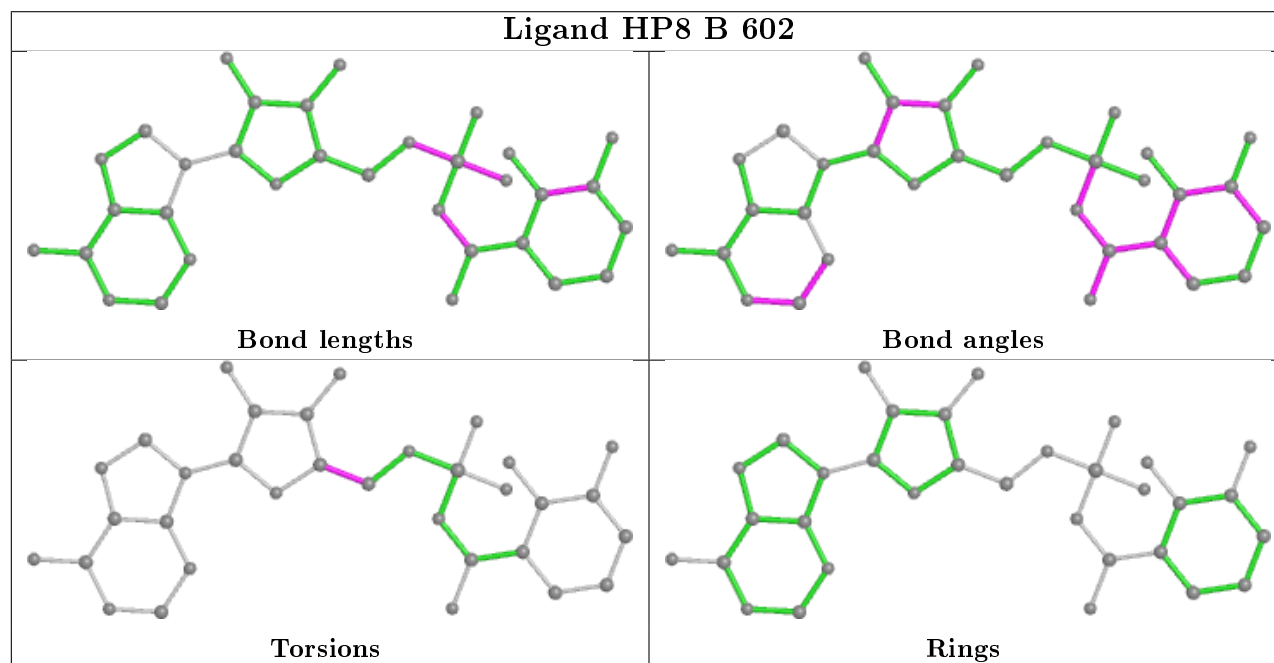
Mol	Chain	Res	Type	Atoms
2	A	601	HP8	C5'-O5'-S-NHS
2	A	601	HP8	C21-NHS-S-OS1
2	B	602	HP8	O4'-C4'-C5'-O5'
2	B	602	HP8	C3'-C4'-C5'-O5'

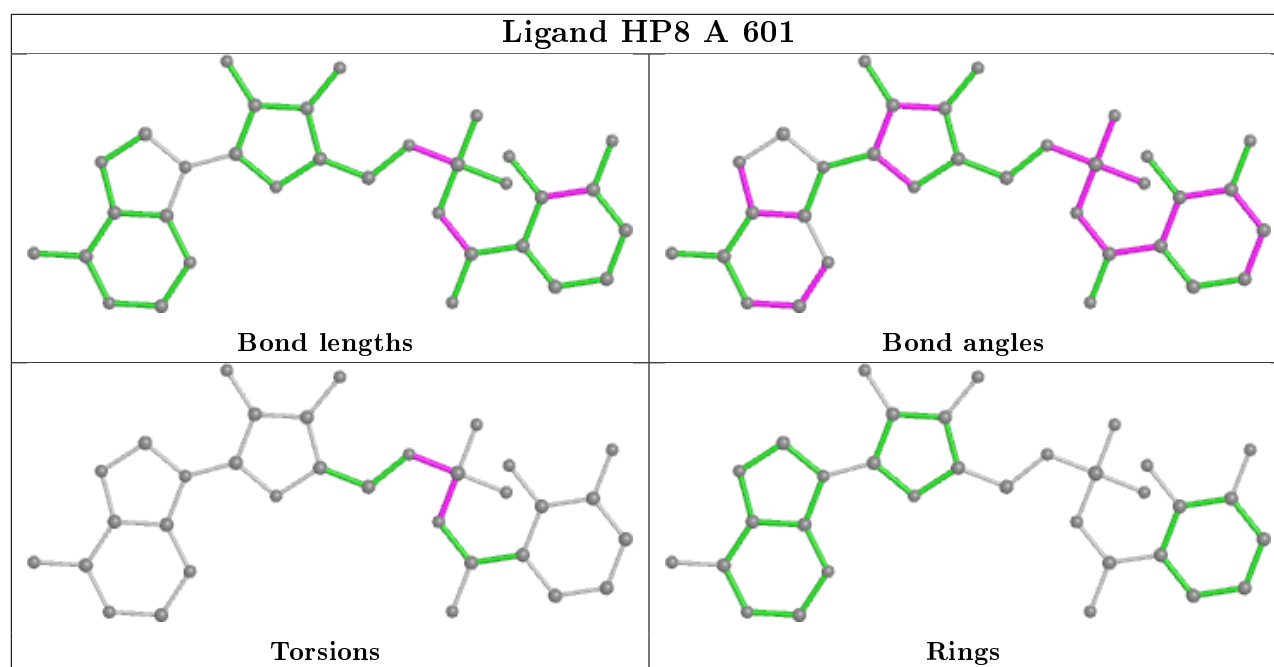
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	602	HP8	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	435/544 (79%)	-0.12	11 (2%) 57 59	41, 55, 81, 96	0
1	B	435/544 (79%)	-0.16	6 (1%) 75 77	41, 56, 80, 93	0
All	All	870/1088 (79%)	-0.14	17 (1%) 65 67	41, 55, 80, 96	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	202	GLY	4.9
1	B	180	PHE	4.1
1	A	426	PRO	3.4
1	A	412	GLU	3.4
1	A	376	GLN	3.4
1	B	426	PRO	3.4
1	A	142	GLN	2.8
1	A	180	PHE	2.6
1	B	269	LEU	2.4
1	A	200	SER	2.4
1	B	412	GLU	2.4
1	A	203	THR	2.3
1	B	179	THR	2.2
1	A	167	PHE	2.2
1	A	165	THR	2.2
1	A	404	GLU	2.2
1	B	149	ASP	2.1

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

There are no carbohydrates in this entry.

6.4 Ligands

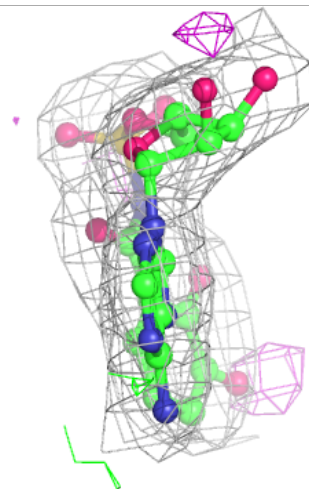
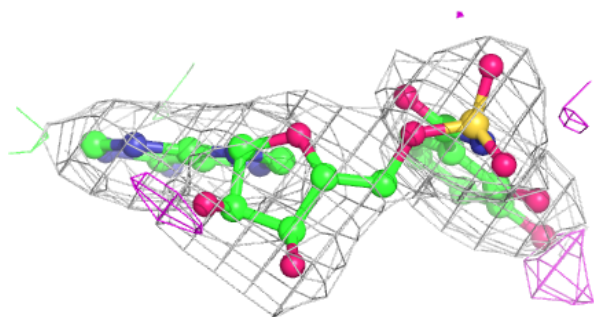
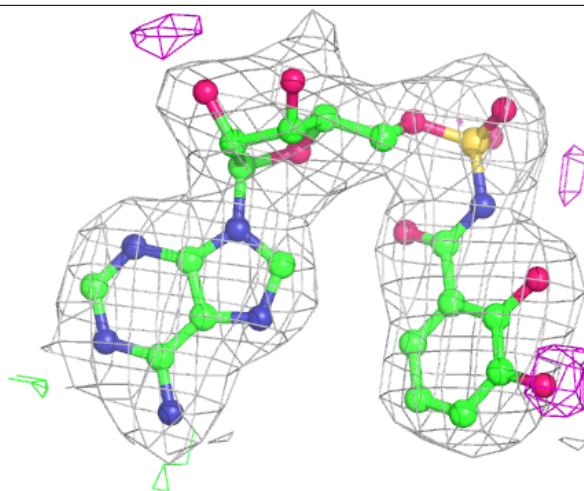
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CA	B	543	1/1	0.95	0.12	89,89,89,89	0
3	CA	B	544	1/1	0.95	0.15	62,62,62,62	0
3	CA	A	544	1/1	0.96	0.14	56,56,56,56	0
3	CA	A	543	1/1	0.97	0.12	53,53,53,53	0
2	HP8	A	601	33/33	0.98	0.15	36,40,49,55	0
2	HP8	B	602	33/33	0.98	0.15	44,48,51,55	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.

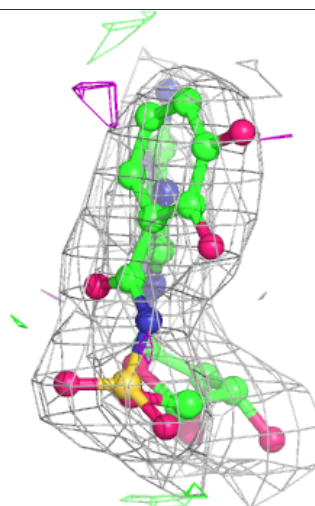
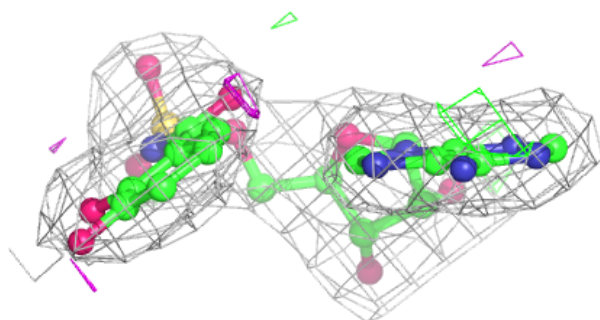
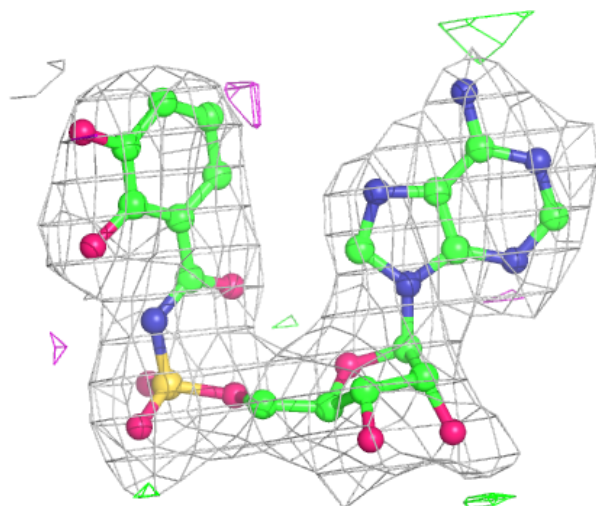
Electron density around HP8 A 601:

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



Electron density around HP8 B 602:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.