



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 08:20 PM UTC

PDB ID : 4UR9 / pdb_00004ur9
Title : Structure of ligand bound glycosylhydrolase
Authors : Darby, J.F.; Landstroem, J.; Roth, C.; He, Y.; Schultz, M.; Davies, G.J.; Hubbard, R.E.
Deposited on : 2014-06-27
Resolution : 2.20 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

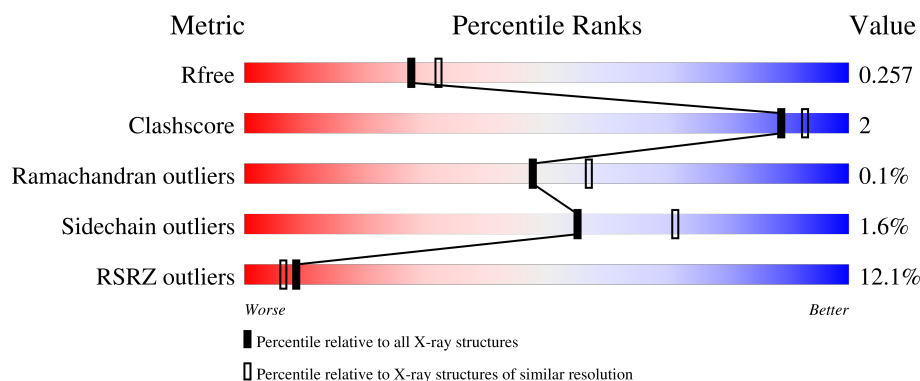
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	716	12% 89% 6% ••
1	B	716	11% 89% 5% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BK9	A	1717	-	X	-	-
2	BK9	B	1717	-	X	-	-

2 Entry composition [i](#)

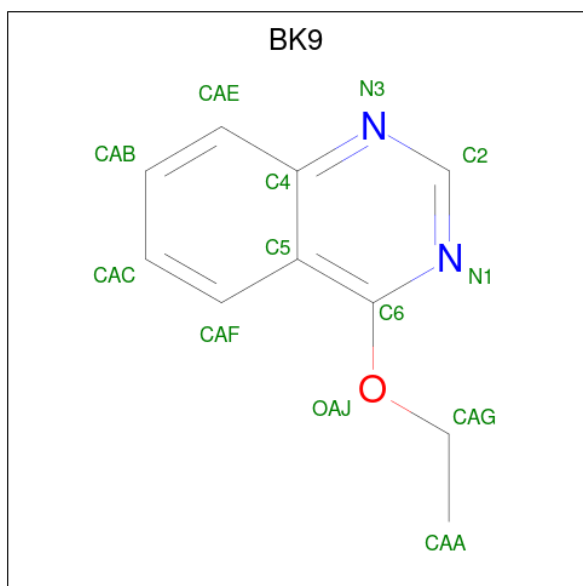
There are 5 unique types of molecules in this entry. The entry contains 22353 atoms, of which 10891 are hydrogens and 0 are deuteriums.

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 O-GLCNACASE BT_4395.

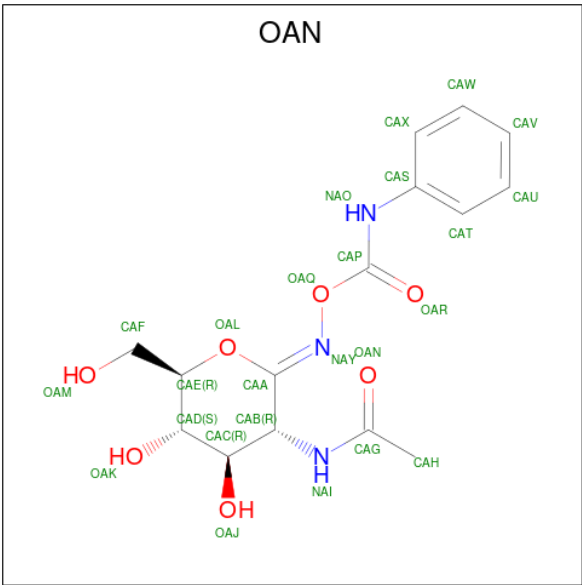
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	685	Total	C	H	N	O	S	0	0	0
			11049	3559	5490	941	1041	18			
1	B	667	Total	C	H	N	O	S	0	1	0
			10784	3471	5363	919	1013	18			

- Molecule 2 is 4-ethoxyquinazoline (CCD ID: BK9) (formula: $C_{10}H_{10}N_2O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	10	2	1		
2	B	1	Total	C	N	O	0	0
			13	10	2	1		

- Molecule 3 is O-(2-ACETAMIDO-2-DEOXY D-GLUCOPYRANOSYLIDENE) AMINO-N-PHENYLCARBAMATE (CCD ID: OAN) (formula: $C_{15}H_{19}N_3O_7$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			44	15	19	3	7		
3	B	1	Total	C	H	N	O	0	0
			44	15	19	3	7		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		

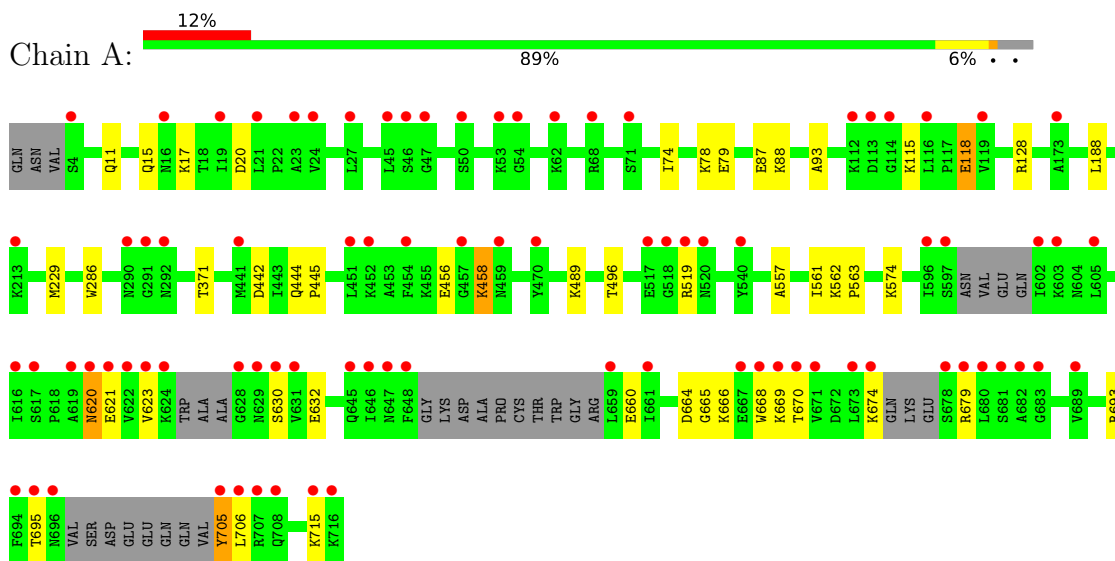
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	204	Total	O	0	0
			204	204		
5	B	201	Total	O	0	0
			201	201		

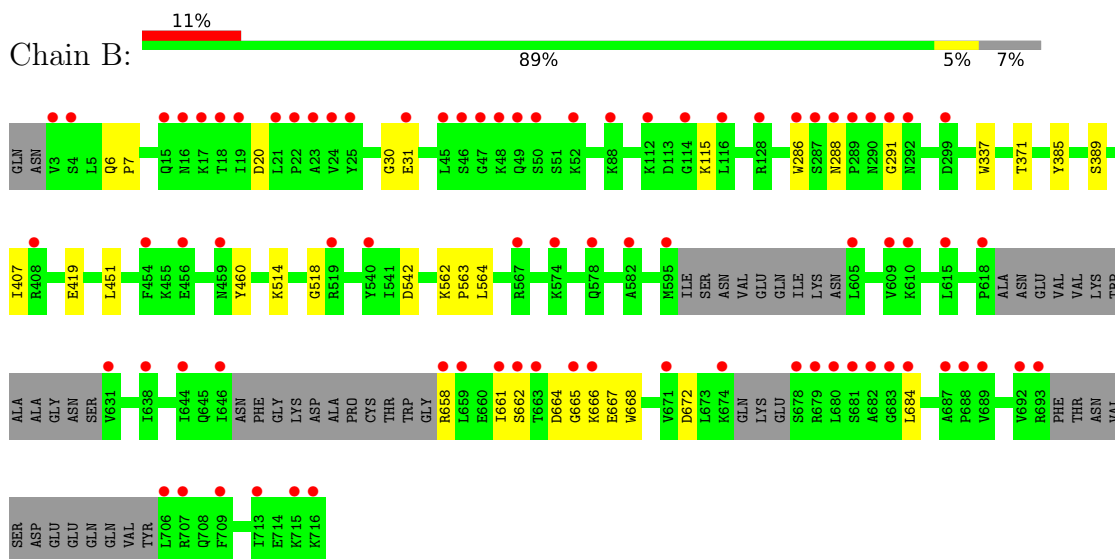
3 Residue-property plots [i](#)

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: O-GLCNACASE BT_4395



• Molecule 1: O-GLCNACASE BT_4395



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.49Å 162.74Å 223.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.79 – 2.20 48.79 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.79-2.20) 99.5 (48.79-2.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, R_{free}	0.222 , 0.252 0.230 , 0.257	Depositor DCC
R_{free} test set	4792 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.5	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	22353	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.85 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.4299e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, OAN, BK9

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	0.32	0/5693	0.63	0/7706
1	B	0.32	0/5555	0.63	1/7519 (0.0%)
All	All	0.32	0/11248	0.63	1/15225 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	30	GLY	N-CA-C	-5.43	106.22	112.73

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	5559	5490	5489	24	0
1	B	5421	5363	5362	17	0
2	A	13	0	8	0	0
2	B	13	0	8	0	0
3	A	25	19	19	2	0
3	B	25	19	19	1	0
4	B	1	0	0	0	0
5	A	204	0	0	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	201	0	0	2	0
All	All	11462	10891	10905	41	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 2.

All (41) 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:79:GLU:OE2	1:A:128:ARG:NH2	2.32	0.63
1:B:662:SER:OG	1:B:664:ASP:O	2.15	0.60
1:A:664:ASP:OD2	1:A:666:LYS:NZ	2.34	0.60
1:B:542:ASP:OD1	5:B:2133:HOH:O	2.16	0.58
1:A:621:GLU:OE1	5:A:2196:HOH:O	2.17	0.56
1:A:574:LYS:NZ	5:A:2188:HOH:O	2.38	0.53
1:A:632:GLU:OE1	1:A:693:ARG:NH2	2.42	0.52
1:B:288:ASN:OD1	1:B:291:GLY:N	2.44	0.51
1:B:7:PRO:HB3	1:B:389:SER:HA	1.91	0.51
1:B:6:GLN:HG2	1:B:385:TYR:CE1	2.47	0.50
1:A:489:LYS:NZ	5:A:2150:HOH:O	2.44	0.49
1:B:662:SER:OG	1:B:667:GLU:O	2.31	0.48
1:B:658:ARG:NH2	1:B:672:ASP:HA	2.29	0.47
1:A:442:ASP:N	1:A:442:ASP:OD1	2.47	0.47
1:B:286:TRP:CH2	3:B:1718:OAN:HAX	2.50	0.47
1:B:665:GLY:O	1:B:668:TRP:NE1	2.48	0.47
1:A:674:LYS:N	5:A:2202:HOH:O	2.46	0.46
1:B:20:ASP:OD1	1:B:115:LYS:NZ	2.47	0.46
1:B:562:LYS:HB3	1:B:563:PRO:HD3	1.97	0.45
1:B:661:ILE:HD11	1:B:684:LEU:HD22	1.99	0.45
1:A:557:ALA:HB1	1:A:561:ILE:HB	1.99	0.45
1:B:337:TRP:CD1	1:B:337:TRP:C	2.96	0.44
1:A:456:GLU:OE1	1:A:458:LYS:NZ	2.35	0.44
1:A:444:GLN:HB3	1:A:445:PRO:HD3	1.99	0.43
1:A:623:VAL:HG12	1:A:705:TYR:HA	1.99	0.43
1:B:451:LEU:HD21	1:B:564:LEU:HD12	2.00	0.43
1:A:20:ASP:OD1	1:A:115:LYS:NZ	2.49	0.43
1:A:562:LYS:HB3	1:A:563:PRO:HD3	2.01	0.43
1:B:460:TYR:OH	5:B:2170:HOH:O	2.22	0.43
1:A:665:GLY:O	1:A:668:TRP:NE1	2.52	0.42
1:B:407:ILE:HD13	1:B:419:GLU:HA	2.02	0.42
1:A:660:GLU:HG3	1:A:670:THR:HA	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:660:GLU:HG2	1:A:668:TRP:HB3	2.00	0.42
1:A:620:ASN:N	1:A:620:ASN:OD1	2.52	0.42
1:A:15:GLN:HB2	1:A:118:GLU:HG2	2.01	0.41
1:A:286:TRP:CH2	3:A:1718:OAN:HAX	2.55	0.41
1:A:188:LEU:HD13	1:A:229:MET:HE1	2.01	0.41
1:A:286:TRP:HH2	3:A:1718:OAN:HAX	1.85	0.41
1:A:660:GLU:HG3	1:A:669:LYS:C	2.45	0.40
1:A:74:ILE:HD11	1:A:93:ALA:HB1	2.03	0.40
1:B:514:LYS:O	1:B:518:GLY:N	2.48	0.40

There are no symmetry-related clashes.

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	673/716 (94%)	652 (97%)	21 (3%)	0	100	100
1	B	656/716 (92%)	631 (96%)	24 (4%)	1 (0%)	43	51
All	All	1329/1432 (93%)	1283 (96%)	45 (3%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	31	GLU

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	604/630 (96%)	587 (97%)	17 (3%)	38	52
1	B	589/630 (94%)	587 (100%)	2 (0%)	86	93
All	All	1193/1260 (95%)	1174 (98%)	19 (2%)	55	71

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	17	LYS
1	A	78	LYS
1	A	87	GLU
1	A	88	LYS
1	A	118	GLU
1	A	371	THR
1	A	458	LYS
1	A	496	THR
1	A	519	ARG
1	A	620	ASN
1	A	630	SER
1	A	679	ARG
1	A	695	THR
1	A	705	TYR
1	A	706	LEU
1	A	715	LYS
1	B	371	THR
1	B	666	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	469	GLN
1	B	73	GLN
1	B	77	HIS
1	B	372	ASN
1	B	500	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is 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$
3	OAN	B	1718	-	23,26,26	2.62	7 (30%)	24,35,35	1.24	4 (16%)
2	BK9	B	1717	-	14,14,14	3.31	8 (57%)	17,18,18	6.45	11 (64%)
3	OAN	A	1718	-	23,26,26	2.67	7 (30%)	24,35,35	1.15	3 (12%)
2	BK9	A	1717	-	14,14,14	3.33	8 (57%)	17,18,18	6.48	11 (64%)

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	OAN	B	1718	-	-	4/12/35/35	0/2/2/2
2	BK9	B	1717	-	-	2/3/3/3	0/2/2/2
3	OAN	A	1718	-	-	6/12/35/35	0/2/2/2
2	BK9	A	1717	-	-	2/3/3/3	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1718	OAN	CAP-NAO	8.81	1.53	1.36

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1718	OAN	CAP-NAO	8.70	1.53	1.36
2	A	1717	BK9	C5-C4	5.53	1.51	1.42
2	B	1717	BK9	C5-C4	5.17	1.50	1.42
2	B	1717	BK9	CAC-CAF	-5.16	1.26	1.36
2	A	1717	BK9	CAC-CAF	-5.09	1.26	1.36
2	B	1717	BK9	OAJ-CAG	-5.01	1.26	1.43
3	A	1718	OAN	CAG-NAI	5.00	1.50	1.34
2	A	1717	BK9	OAJ-CAG	-4.88	1.27	1.43
3	B	1718	OAN	CAG-NAI	4.86	1.50	1.34
2	A	1717	BK9	OAJ-C6	4.65	1.40	1.34
2	B	1717	BK9	OAJ-C6	4.56	1.40	1.34
2	B	1717	BK9	CAB-CAE	-4.54	1.27	1.36
2	A	1717	BK9	CAB-CAE	-4.51	1.27	1.36
3	A	1718	OAN	OAQ-NAY	-4.03	1.35	1.44
3	B	1718	OAN	OAQ-NAY	-3.78	1.36	1.44
2	B	1717	BK9	C6-N1	3.19	1.36	1.31
2	A	1717	BK9	C6-N1	3.11	1.36	1.31
2	A	1717	BK9	CAE-C4	-3.04	1.36	1.41
3	B	1718	OAN	OAR-CAP	-3.00	1.16	1.21
3	A	1718	OAN	CAC-CAB	-3.00	1.49	1.52
2	B	1717	BK9	CAE-C4	-2.97	1.36	1.41
3	B	1718	OAN	CAC-CAB	-2.73	1.49	1.52
3	A	1718	OAN	OAR-CAP	-2.70	1.16	1.21
3	B	1718	OAN	CAS-NAO	2.62	1.46	1.41
3	A	1718	OAN	CAS-NAO	2.54	1.46	1.41
3	A	1718	OAN	OAN-CAG	2.51	1.28	1.23
3	B	1718	OAN	OAN-CAG	2.45	1.28	1.23
2	B	1717	BK9	CAA-CAG	-2.22	1.31	1.47
2	A	1717	BK9	CAA-CAG	-2.17	1.32	1.47

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1717	BK9	CAG-OAJ-C6	14.83	136.41	117.32
2	B	1717	BK9	CAG-OAJ-C6	14.79	136.35	117.32
2	A	1717	BK9	OAJ-CAG-CAA	13.06	148.80	108.04
2	B	1717	BK9	OAJ-CAG-CAA	12.99	148.58	108.04
2	A	1717	BK9	CAF-C5-C6	8.42	134.22	123.45
2	B	1717	BK9	CAF-C5-C6	8.36	134.15	123.45
2	A	1717	BK9	CAB-CAE-C4	7.68	130.59	120.09
2	B	1717	BK9	CAB-CAE-C4	7.65	130.55	120.09
2	A	1717	BK9	CAF-C5-C4	-7.62	109.86	118.36

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1717	BK9	CAF-C5-C4	-7.57	109.92	118.36
2	B	1717	BK9	CAC-CAF-C5	6.24	129.34	120.91
2	A	1717	BK9	CAC-CAF-C5	6.16	129.23	120.91
2	A	1717	BK9	CAE-C4-N3	5.24	126.99	118.55
2	B	1717	BK9	CAE-C4-N3	5.23	126.97	118.55
2	A	1717	BK9	CAE-C4-C5	-4.29	114.51	119.13
2	B	1717	BK9	CAE-C4-C5	-4.28	114.52	119.13
2	A	1717	BK9	C5-C4-N3	-4.07	118.50	122.82
2	B	1717	BK9	CAC-CAB-CAE	-4.07	114.95	120.40
2	A	1717	BK9	CAC-CAB-CAE	-4.07	114.95	120.40
2	B	1717	BK9	C5-C4-N3	-4.06	118.52	122.82
2	A	1717	BK9	OAJ-C6-C5	3.07	118.25	115.01
3	A	1718	OAN	OAL-CAE-CAF	2.98	109.04	106.05
3	B	1718	OAN	OAL-CAE-CAF	2.98	109.04	106.05
2	B	1717	BK9	OAJ-C6-C5	2.90	118.08	115.01
3	B	1718	OAN	CAS-NAO-CAP	-2.80	121.96	126.39
3	A	1718	OAN	CAS-NAO-CAP	-2.37	122.64	126.39
3	B	1718	OAN	OAR-CAP-NAO	-2.30	121.22	126.12
3	B	1718	OAN	CAH-CAG-NAI	2.22	119.79	116.12
3	A	1718	OAN	CAH-CAG-NAI	2.05	119.53	116.12

There are no chirality outliers.

All (14) torsion outliers are listed below:

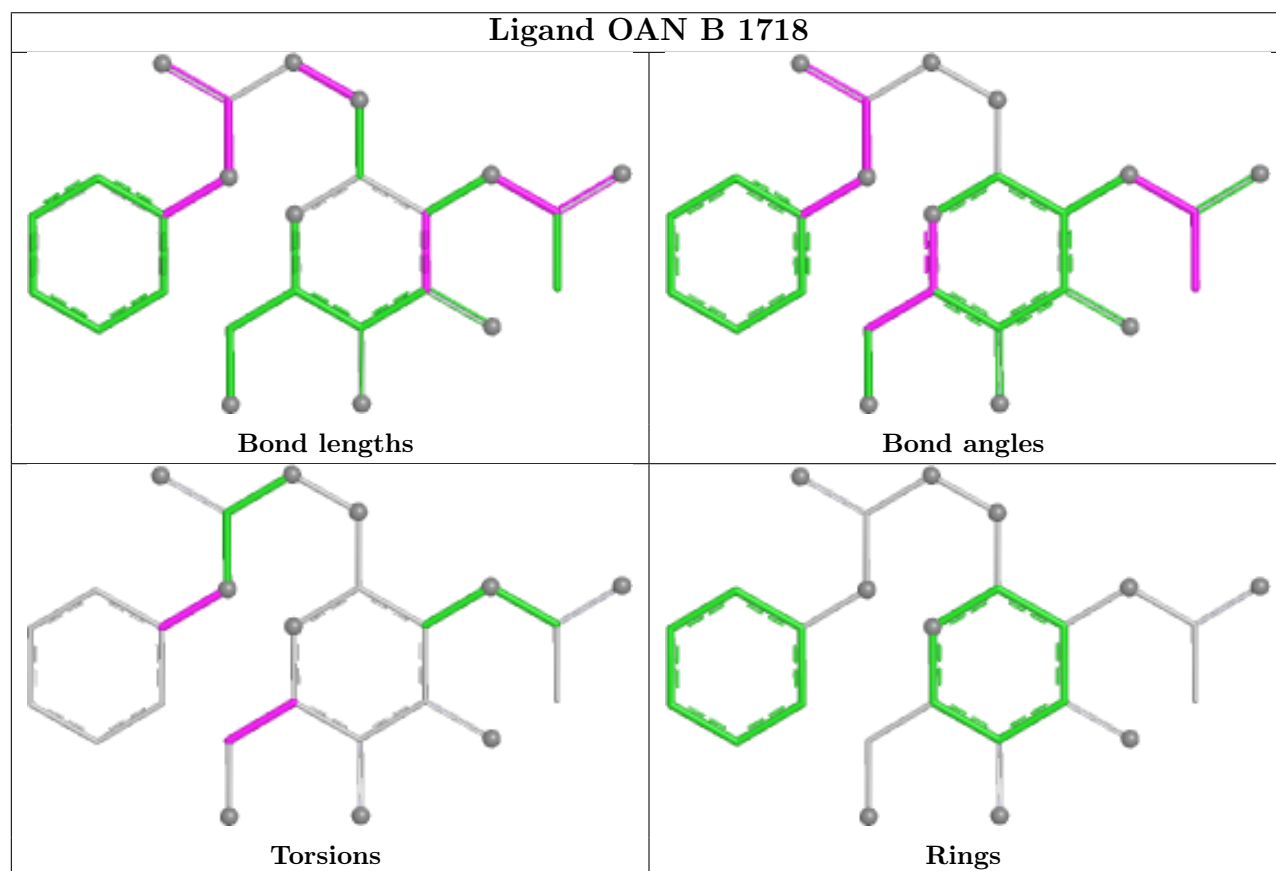
Mol	Chain	Res	Type	Atoms
2	A	1717	BK9	N1-C6-OAJ-CAG
2	A	1717	BK9	C5-C6-OAJ-CAG
2	B	1717	BK9	N1-C6-OAJ-CAG
2	B	1717	BK9	C5-C6-OAJ-CAG
3	A	1718	OAN	OAR-CAP-OAQ-NAY
3	A	1718	OAN	NAO-CAP-OAQ-NAY
3	B	1718	OAN	CAD-CAE-CAF-OAM
3	B	1718	OAN	OAL-CAE-CAF-OAM
3	A	1718	OAN	OAR-CAP-NAO-CAS
3	B	1718	OAN	CAT-CAS-NAO-CAP
3	B	1718	OAN	CAX-CAS-NAO-CAP
3	A	1718	OAN	CAD-CAE-CAF-OAM
3	A	1718	OAN	OAQ-CAP-NAO-CAS
3	A	1718	OAN	OAL-CAE-CAF-OAM

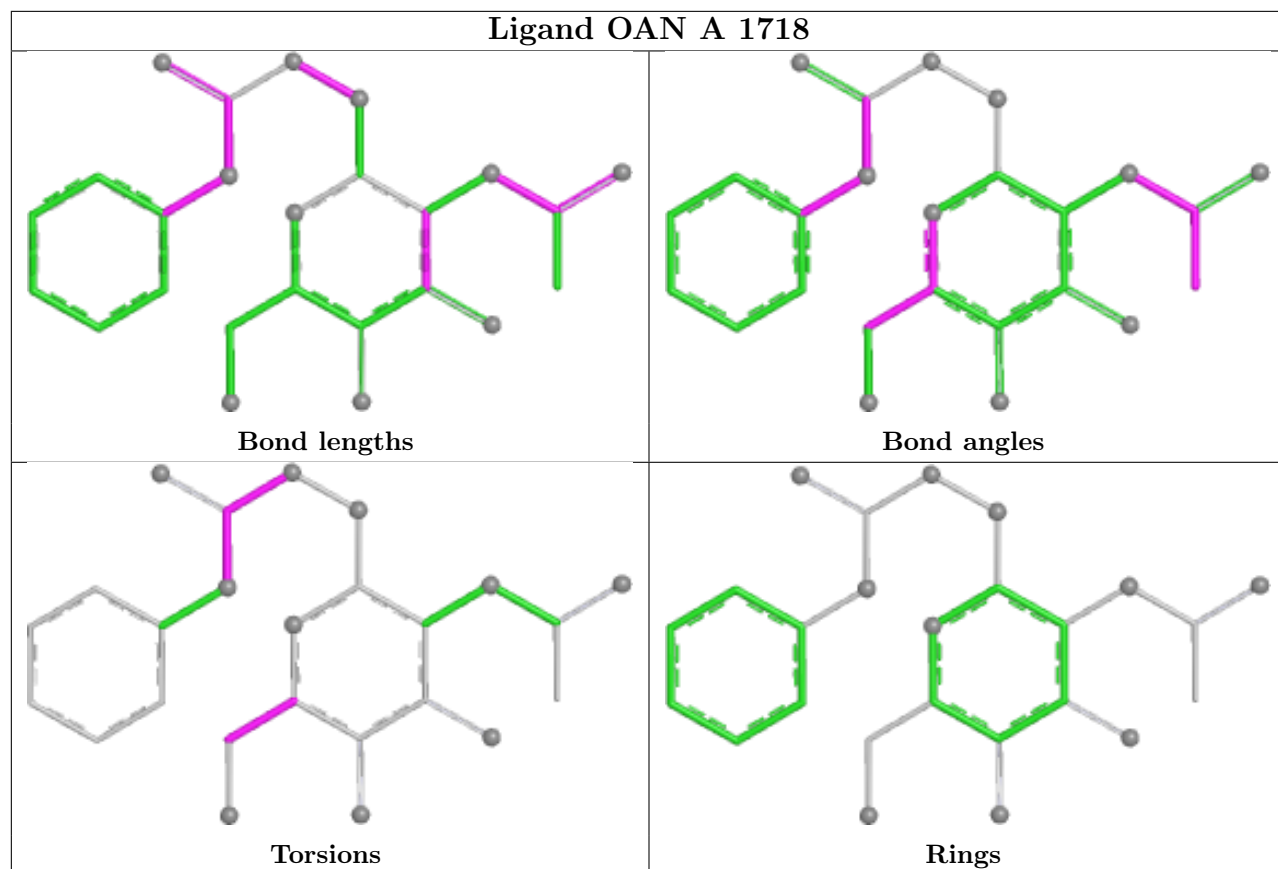
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1718	OAN	1	0
3	A	1718	OAN	2	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	685/716 (95%)	0.57	84 (12%) 8 6	37, 58, 137, 195	0
1	B	667/716 (93%)	0.54	80 (11%) 9 6	30, 56, 139, 196	1 (0%)
All	All	1352/1432 (94%)	0.55	164 (12%) 8 6	30, 57, 139, 196	1 (0%)

All (164) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	648	PHE	8.0
1	A	705	TYR	6.9
1	A	622	VAL	6.6
1	A	457	GLY	6.4
1	A	706	LEU	6.3
1	B	289	PRO	6.3
1	B	678	SER	6.1
1	A	678	SER	5.9
1	A	628	GLY	5.8
1	B	3	VAL	5.7
1	A	624	LYS	5.5
1	A	518	GLY	5.5
1	A	623	VAL	5.5
1	A	696	ASN	4.9
1	B	18	THR	4.7
1	B	631	VAL	4.7
1	B	290	ASN	4.7
1	B	286	TRP	4.5
1	B	4[A]	SER	4.4
1	B	595	MET	4.4
1	A	661	ILE	4.4
1	B	674	LYS	4.3
1	A	674	LYS	4.2
1	B	605	LEU	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	682	ALA	4.0
1	A	291	GLY	4.0
1	B	688	PRO	4.0
1	A	597	SER	4.0
1	B	689	VAL	3.9
1	A	621	GLU	3.9
1	A	646	ILE	3.9
1	B	31	GLU	3.8
1	B	706	LEU	3.8
1	A	659	LEU	3.7
1	B	680	LEU	3.7
1	B	663	THR	3.7
1	B	291	GLY	3.7
1	A	23	ALA	3.7
1	B	49	GLN	3.6
1	A	680	LEU	3.6
1	B	618	PRO	3.6
1	B	665	GLY	3.5
1	B	46	SER	3.5
1	A	679	ARG	3.5
1	B	658	ARG	3.5
1	B	21	LEU	3.5
1	B	48	LYS	3.4
1	B	15	GLN	3.4
1	A	459	ASN	3.4
1	B	661	ILE	3.4
1	A	605	LEU	3.3
1	A	45	LEU	3.3
1	B	540	TYR	3.3
1	A	681	SER	3.3
1	A	707	ARG	3.3
1	B	646	ILE	3.2
1	A	630	SER	3.2
1	A	441	MET	3.2
1	B	292	ASN	3.2
1	A	673	LEU	3.2
1	A	647	ASN	3.2
1	A	519	ARG	3.1
1	A	694	PHE	3.1
1	B	582	ALA	3.1
1	A	682	ALA	3.1
1	B	128	ARG	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	602	ILE	3.0
1	B	671	VAL	3.0
1	A	19	ILE	3.0
1	A	695	THR	3.0
1	B	114	GLY	3.0
1	A	715	LYS	3.0
1	A	47	GLY	3.0
1	A	629	ASN	2.9
1	B	519	ARG	2.9
1	B	713	ILE	2.9
1	A	603	LYS	2.9
1	A	596	ILE	2.9
1	B	716	LYS	2.9
1	B	687	ALA	2.8
1	B	681	SER	2.8
1	B	683	GLY	2.8
1	A	470	TYR	2.8
1	A	292	ASN	2.8
1	A	619	ALA	2.8
1	B	659	LEU	2.8
1	A	4	SER	2.7
1	A	113	ASP	2.7
1	A	517	GLU	2.7
1	A	290	ASN	2.7
1	B	709	PHE	2.6
1	A	689	VAL	2.6
1	B	24	VAL	2.6
1	B	666	LYS	2.6
1	B	644	ILE	2.6
1	B	19	ILE	2.6
1	B	408	ARG	2.6
1	A	668	TRP	2.6
1	A	631	VAL	2.6
1	B	692	VAL	2.6
1	A	54	GLY	2.5
1	B	454	PHE	2.5
1	A	617	SER	2.5
1	B	715	LYS	2.5
1	B	288	ASN	2.5
1	A	24	VAL	2.5
1	A	540	TYR	2.5
1	A	46	SER	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	609	VAL	2.5
1	A	53	LYS	2.5
1	A	670	THR	2.4
1	A	116	LEU	2.4
1	B	47	GLY	2.4
1	A	452	LYS	2.4
1	B	679	ARG	2.4
1	B	17	LYS	2.4
1	A	50	SER	2.4
1	B	567	ARG	2.3
1	B	693	ARG	2.3
1	B	638	ILE	2.3
1	B	45	LEU	2.3
1	A	27	LEU	2.3
1	A	71	SER	2.3
1	B	287	SER	2.3
1	A	645	GLN	2.3
1	A	213	LYS	2.3
1	B	112	LYS	2.3
1	A	21	LEU	2.2
1	A	68	ARG	2.2
1	A	683	GLY	2.2
1	A	451	LEU	2.2
1	A	16	ASN	2.2
1	A	173	ALA	2.2
1	B	299	ASP	2.2
1	B	88	LYS	2.2
1	A	708	GLN	2.2
1	B	615	LEU	2.1
1	A	669	LYS	2.1
1	B	456	GLU	2.1
1	B	574	LYS	2.1
1	B	707	ARG	2.1
1	A	620	ASN	2.1
1	A	671	VAL	2.1
1	B	22	PRO	2.1
1	B	662	SER	2.1
1	B	52	LYS	2.1
1	A	667	GLU	2.1
1	A	616	ILE	2.1
1	B	684	LEU	2.1
1	A	119	VAL	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	716	LYS	2.1
1	B	16	ASN	2.1
1	B	459	ASN	2.1
1	B	578	GLN	2.1
1	A	114	GLY	2.1
1	A	454	PHE	2.1
1	B	25	TYR	2.0
1	A	112	LYS	2.0
1	B	610	LYS	2.0
1	B	50	SER	2.0
1	B	23	ALA	2.0
1	A	520	ASN	2.0
1	B	116	LEU	2.0
1	A	62	LYS	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](#)

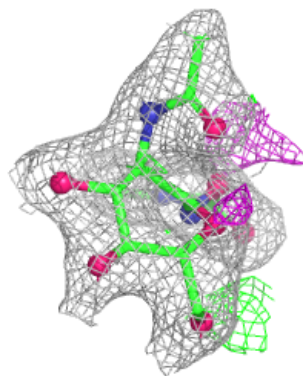
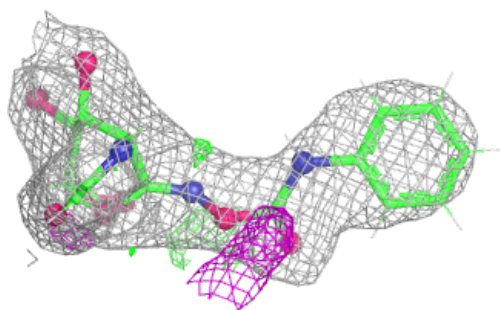
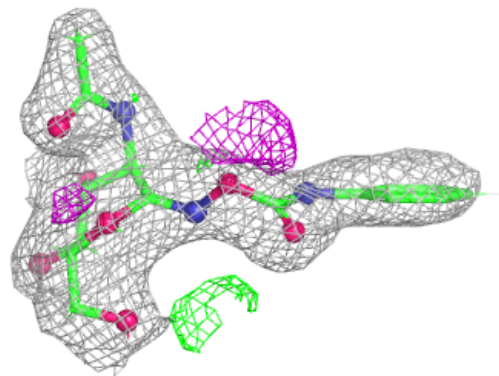
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BK9	A	1717	13/13	0.91	0.14	48,54,58,59	0
2	BK9	B	1717	13/13	0.93	0.13	56,64,67,68	0
3	OAN	A	1718	25/25	0.93	0.10	34,48,60,66	0
3	OAN	B	1718	25/25	0.94	0.09	36,51,73,76	0
4	CA	B	1719	1/1	0.99	0.03	43,43,43,43	0

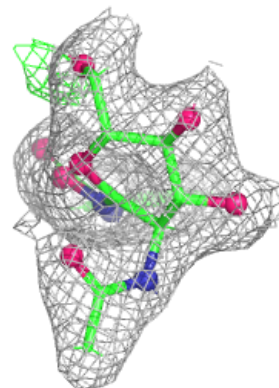
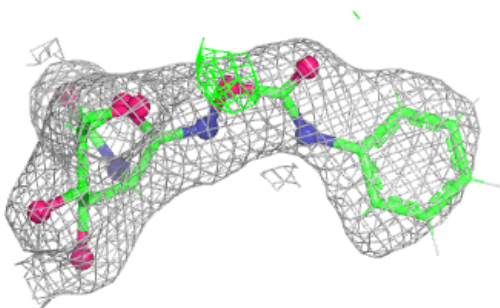
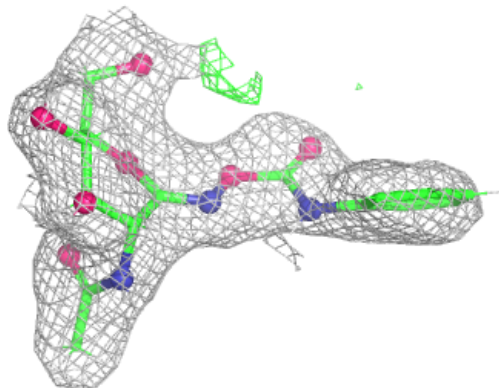
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OAN A 1718:

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

**Electron density around OAN B 1718:**

$2mF_o-DF_c$ (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.