



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:48 PM UTC

PDB ID : 2CBJ / pdb_00002cbj
Title : Structure of the Clostridium perfringens NagJ family 84 glycoside hydrolase, a homologue of human O-GlcNAcase in complex with PUGNAc
Authors : Rao, F.V.; Dorfmueller, H.C.; Villa, F.; Allwood, M.; Eggleston, I.M.; van Aalten, D.M.F.
Deposited on : 2006-01-05
Resolution : 2.35 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

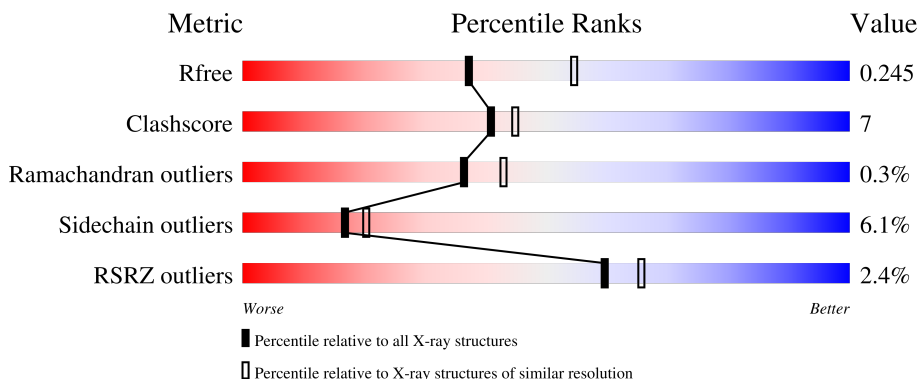
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	594	 3% 81% 14% ..
1	B	594	 2% 82% 15% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9589 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.

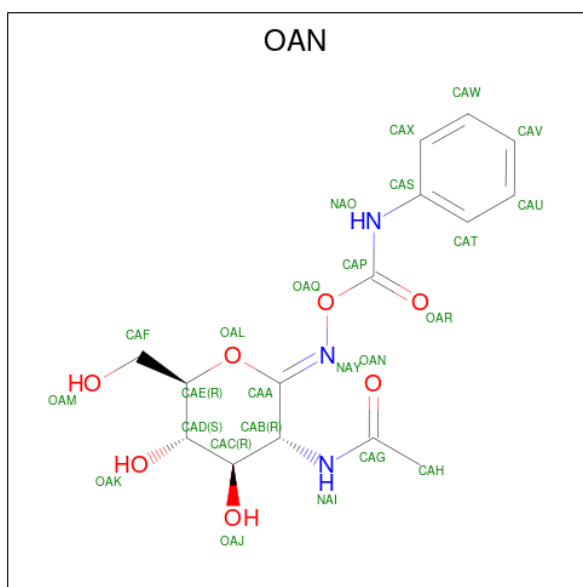
- Molecule 1 is a protein called HYALURONIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	585	4628	2913	757	941	17	0	0	0
1	B	585	4629	2913	758	941	17	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	196	GLN	LYS	conflict	UNP Q8XL08
A	234	SER	ASN	conflict	UNP Q8XL08
A	244	ASN	ASP	conflict	UNP Q8XL08
A	268	ASP	GLU	conflict	UNP Q8XL08
A	279	THR	ALA	conflict	UNP Q8XL08
A	348	ALA	THR	conflict	UNP Q8XL08
B	196	GLN	LYS	conflict	UNP Q8XL08
B	234	SER	ASN	conflict	UNP Q8XL08
B	244	ASN	ASP	conflict	UNP Q8XL08
B	268	ASP	GLU	conflict	UNP Q8XL08
B	279	THR	ALA	conflict	UNP Q8XL08
B	348	ALA	THR	conflict	UNP Q8XL08

- Molecule 2 is O-(2-ACETAMIDO-2-DEOXY D-GLUCOPYRANOSYLIDENE) AMINO-N-PHENYLCARBAMATE (CCD ID: OAN) (formula: C₁₅H₁₉N₃O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
2	A	1	Total	C	N	O	0	0
			25	15	3	7		
2	B	1	Total	C	N	O	0	0
			25	15	3	7		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		

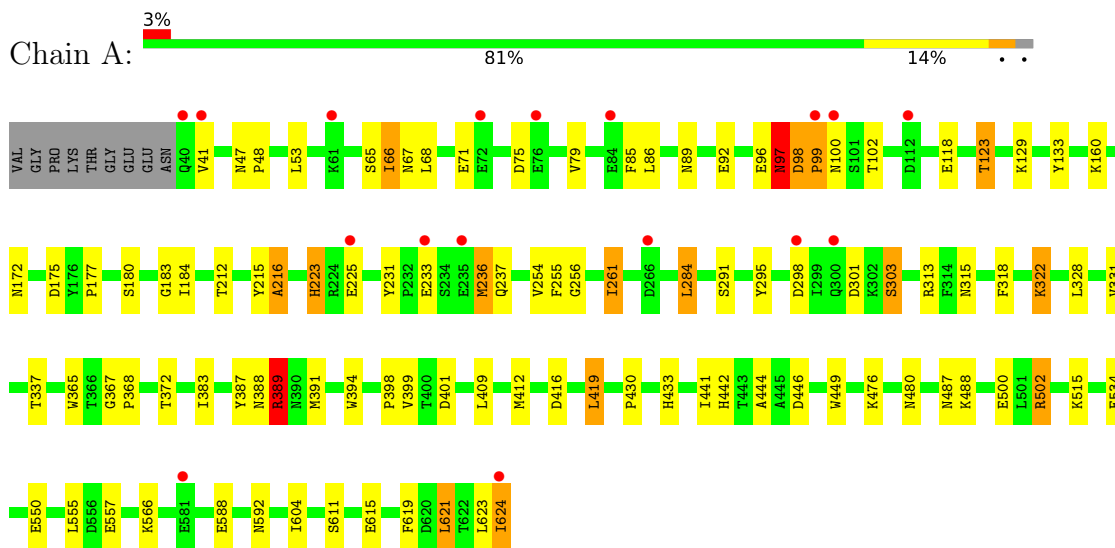
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		
4	B	137	Total	O	0	0
			137	137		

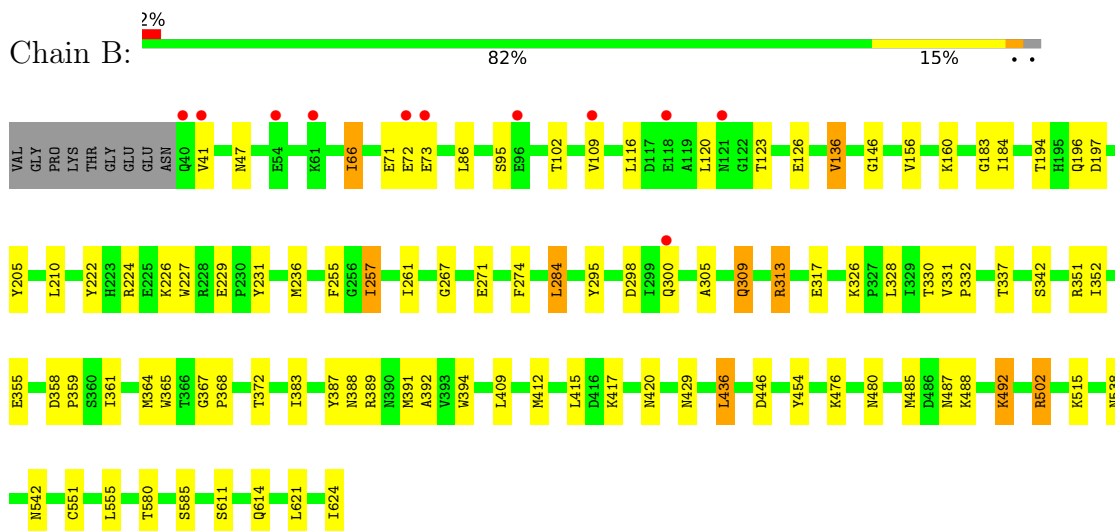
3 Residue-property plots [i](#)

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: HYALURONIDASE



• Molecule 1: HYALURONIDASE



4 Data and refinement statistics i

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	129.61Å 145.75Å 152.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.35 20.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	93.5 (20.00-2.35) 93.3 (20.00-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 2.35Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.193 , 0.248 0.192 , 0.245	Depositor DCC
R_{free} test set	562 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtrriage
Anisotropy	0.291	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 36.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.011 for -h,-l,-k	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9589	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: CL, OAN

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	0.77	0/4724	0.97	6/6413 (0.1%)
1	B	0.76	1/4725 (0.0%)	0.95	4/6415 (0.1%)
All	All	0.77	1/9449 (0.0%)	0.96	10/12828 (0.1%)

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	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	267	GLY	C-O	6.31	1.32	1.23

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	367	GLY	CA-C-N	6.72	126.42	119.56
1	A	367	GLY	C-N-CA	6.72	126.42	119.56
1	B	367	GLY	CA-C-N	6.05	125.73	119.56
1	B	367	GLY	C-N-CA	6.05	125.73	119.56
1	A	216	ALA	CA-C-N	-6.00	113.35	120.13
1	A	216	ALA	C-N-CA	-6.00	113.35	120.13
1	A	98	ASP	N-CA-C	5.27	121.45	109.81
1	A	389	ARG	N-CA-C	5.13	115.31	108.38
1	B	429	ASN	CA-C-N	5.08	126.18	119.84
1	B	429	ASN	C-N-CA	5.08	126.18	119.84

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	97	ASN	Peptide
1	A	98	ASP	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	4628	0	4418	66	0
1	B	4629	0	4420	56	0
2	A	25	0	19	1	0
2	B	25	0	19	1	0
3	B	2	0	0	0	0
4	A	143	0	0	2	0
4	B	137	0	0	0	0
All	All	9589	0	8876	117	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 7.

All (117) 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:123:THR:HG21	1:A:172:ASN:HD21	1.05	1.09
1:A:123:THR:HG21	1:A:172:ASN:ND2	1.72	1.03
1:B:305:ALA:O	1:B:309:GLN:HG2	1.61	1.00
1:A:619:PHE:HE2	1:A:621:LEU:HD13	1.29	0.93
1:A:619:PHE:CE2	1:A:621:LEU:HD13	2.10	0.87
1:B:417:LYS:HE2	1:B:454:TYR:O	1.75	0.87
1:A:215:TYR:O	1:A:223:HIS:HE1	1.62	0.83
1:A:223:HIS:CE1	1:A:256:GLY:O	2.39	0.75
1:B:387:TYR:HD2	1:B:389:ARG:NH1	1.86	0.74
1:A:261:ILE:H	1:A:261:ILE:HD13	1.53	0.73
1:B:502:ARG:O	1:B:502:ARG:HD3	1.87	0.73
1:A:387:TYR:HB3	1:A:389:ARG:HD2	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:TYR:HB2	1:A:236:MET:SD	2.31	0.69
1:B:359:PRO:HA	1:B:389:ARG:HH21	1.58	0.69
1:A:295:TYR:CD2	1:A:331:VAL:HG12	2.27	0.69
1:A:215:TYR:O	1:A:223:HIS:CE1	2.47	0.67
1:A:66:ILE:HG22	1:A:102:THR:HB	1.77	0.66
1:B:66:ILE:HG22	1:B:102:THR:HB	1.80	0.63
1:B:156:VAL:O	1:B:160:LYS:HG3	1.98	0.63
1:B:47:ASN:HD21	1:B:446:ASP:HB2	1.64	0.63
1:A:254:VAL:HG22	1:A:291:SER:HB3	1.80	0.62
2:A:1625:OAN:HAV	1:B:621:LEU:HD23	1.81	0.62
1:A:216:ALA:HA	1:A:223:HIS:CE1	2.35	0.61
1:A:65:SER:HB3	1:A:92:GLU:HB3	1.83	0.60
1:A:75:ASP:O	1:A:79:VAL:HG23	2.01	0.60
1:A:416:ASP:HB3	1:A:419:LEU:HD13	1.84	0.60
1:A:96:GLU:O	1:A:97:ASN:HB2	2.01	0.59
1:A:502:ARG:O	1:A:502:ARG:HD3	2.04	0.58
1:A:183:GLY:C	1:A:184:ILE:HD12	2.29	0.57
1:B:502:ARG:HD3	1:B:502:ARG:C	2.25	0.57
1:B:352:ILE:HA	1:B:355:GLU:HG2	1.88	0.56
1:A:255:PHE:CE2	1:A:284:LEU:HD13	2.41	0.56
1:B:295:TYR:CD2	1:B:331:VAL:HG12	2.41	0.55
1:A:502:ARG:NH2	1:A:615:GLU:OE2	2.38	0.53
1:B:183:GLY:C	1:B:184:ILE:HD12	2.33	0.53
1:A:398:PRO:HD2	1:A:430:PRO:HA	1.89	0.53
1:A:85:PHE:O	1:A:89:ASN:ND2	2.34	0.53
1:B:194:THR:O	1:B:197:ASP:HB2	2.09	0.53
1:B:387:TYR:HD2	1:B:389:ARG:HH11	1.57	0.53
1:A:500:GLU:OE1	4:A:2098:HOH:O	2.19	0.52
1:A:67:ASN:OD1	1:A:97:ASN:O	2.28	0.52
1:A:515:LYS:NZ	1:B:298:ASP:HA	2.24	0.52
1:B:227:TRP:HB2	1:B:257:ILE:HD11	1.92	0.51
1:B:47:ASN:ND2	1:B:446:ASP:HB2	2.25	0.51
1:B:255:PHE:CE2	1:B:284:LEU:HD13	2.46	0.50
1:B:487:ASN:O	1:B:488:LYS:HB2	2.11	0.50
1:A:133:TYR:CZ	1:A:175:ASP:HB3	2.47	0.50
1:A:99:PRO:O	1:A:100:ASN:C	2.55	0.49
1:A:502:ARG:HD3	1:A:502:ARG:C	2.34	0.49
1:A:476:LYS:O	1:A:480:ASN:HB2	2.13	0.49
1:A:624:ILE:OXT	1:A:624:ILE:HG22	2.11	0.49
1:B:351:ARG:O	1:B:355:GLU:HG2	2.13	0.49
1:A:624:ILE:HG21	1:B:224:ARG:HH12	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:538:ASN:OD1	1:B:542:ASN:ND2	2.46	0.48
1:A:254:VAL:HG22	1:A:291:SER:CB	2.44	0.48
1:B:417:LYS:CE	1:B:454:TYR:O	2.57	0.48
1:B:109:VAL:HG11	1:B:126:GLU:HB2	1.96	0.48
1:A:68:LEU:HG	1:A:71:GLU:HG3	1.96	0.48
1:B:476:LYS:O	1:B:480:ASN:HB2	2.14	0.48
1:B:611:SER:HA	1:B:614:GLN:HE21	1.77	0.48
1:A:133:TYR:CE2	1:A:175:ASP:HB3	2.48	0.47
1:A:624:ILE:CG2	1:B:224:ARG:HH12	2.27	0.47
1:B:313:ARG:HD3	1:B:317:GLU:OE2	2.15	0.47
1:B:205:TYR:HB3	1:B:210:LEU:HB2	1.97	0.47
1:A:216:ALA:HA	1:A:223:HIS:ND1	2.29	0.47
1:B:387:TYR:HB3	1:B:389:ARG:HG2	1.96	0.46
1:A:96:GLU:O	1:A:97:ASN:CB	2.62	0.46
1:A:261:ILE:H	1:A:261:ILE:CD1	2.26	0.46
1:A:487:ASN:O	1:A:488:LYS:HB2	2.15	0.46
1:A:123:THR:CG2	1:A:172:ASN:HD21	1.99	0.46
1:B:337:THR:HB	1:B:368:PRO:HA	1.97	0.46
1:B:231:TYR:HB2	1:B:236:MET:SD	2.56	0.46
1:A:53:LEU:C	1:A:53:LEU:HD23	2.40	0.46
1:A:388:ASN:C	1:A:389:ARG:HG3	2.41	0.46
1:A:557:GLU:OE1	1:A:604:ILE:HG22	2.15	0.46
1:A:177:PRO:HB3	1:A:449:TRP:CE3	2.51	0.45
1:A:515:LYS:HZ3	1:B:298:ASP:HA	1.80	0.45
1:B:222:TYR:HB2	1:B:231:TYR:CE1	2.51	0.45
1:B:364:MET:HA	1:B:392:ALA:O	2.16	0.45
1:A:383:ILE:CG2	1:A:391:MET:HE2	2.47	0.45
1:A:433:HIS:HB3	1:A:550:GLU:OE1	2.16	0.45
1:A:48:PRO:HB3	1:A:446:ASP:HA	1.98	0.45
1:A:184:ILE:HD11	1:A:441:ILE:HG23	1.98	0.44
1:A:298:ASP:HA	1:B:515:LYS:HE2	1.98	0.44
1:B:120:LEU:HD23	1:B:136:VAL:HG23	1.98	0.44
1:B:226:LYS:HB3	1:B:229:GLU:HG3	1.99	0.44
1:B:387:TYR:O	1:B:389:ARG:HG2	2.18	0.44
1:B:368:PRO:HD2	1:B:372:THR:HG21	1.99	0.44
1:A:337:THR:HB	1:A:368:PRO:HA	1.99	0.44
1:B:383:ILE:HG22	1:B:391:MET:HE2	2.00	0.44
1:A:223:HIS:NE2	1:A:256:GLY:O	2.51	0.43
1:A:301:ASP:OD2	1:A:303:SER:OG	2.35	0.43
1:A:85:PHE:CD1	1:A:160:LYS:HG2	2.53	0.43
1:B:387:TYR:CD2	1:B:389:ARG:NH1	2.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:GLU:CD	1:B:313:ARG:HH22	2.27	0.43
1:B:394:TRP:C	1:B:394:TRP:CD1	2.97	0.43
1:A:412:MET:HG3	1:A:444:ALA:HA	1.99	0.43
1:A:383:ILE:HG22	1:A:391:MET:HE2	2.01	0.43
1:B:436:LEU:HD21	1:B:551:CYS:HB3	2.01	0.42
1:B:358:ASP:HB2	1:B:361:ILE:HD12	2.01	0.42
1:B:412:MET:HB3	1:B:415:LEU:HD12	2.01	0.42
1:A:315:ASN:ND2	4:A:2024:HOH:O	2.53	0.42
1:A:215:TYR:HD2	1:A:223:HIS:NE2	2.18	0.42
1:B:274:PHE:CZ	1:B:313:ARG:HD2	2.54	0.42
1:A:318:PHE:O	1:A:322:LYS:HG3	2.19	0.42
1:B:109:VAL:HG23	1:B:146:GLY:O	2.20	0.41
1:A:183:GLY:HA3	1:A:212:THR:O	2.20	0.41
1:B:330:THR:O	1:B:364:MET:HG3	2.20	0.41
1:A:399:VAL:HG22	1:A:401:ASP:H	1.86	0.41
1:B:274:PHE:CE2	1:B:313:ARG:HD2	2.56	0.41
1:A:47:ASN:CG	1:A:442:HIS:CE1	2.99	0.40
1:B:383:ILE:CG2	1:B:391:MET:HE2	2.51	0.40
1:B:331:VAL:HA	1:B:332:PRO:HD3	1.97	0.40
1:B:485:MET:O	1:B:492:LYS:HA	2.21	0.40
1:B:298:ASP:OD1	2:B:1625:OAN:HAB	2.22	0.40
1:A:394:TRP:CD1	1:A:394:TRP:C	3.00	0.40
1:A:566:LYS:C	1:A:566:LYS:HD3	2.47	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	583/594 (98%)	555 (95%)	24 (4%)	4 (1%)	18 20
1	B	583/594 (98%)	558 (96%)	25 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1166/1188 (98%)	1113 (96%)	49 (4%)	4 (0%)	36	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	VAL
1	A	97	ASN
1	A	99	PRO
1	A	233	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	500/507 (99%)	470 (94%)	30 (6%)	17	21
1	B	500/507 (99%)	469 (94%)	31 (6%)	16	19
All	All	1000/1014 (99%)	939 (94%)	61 (6%)	17	20

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ILE
1	A	86	LEU
1	A	118	GLU
1	A	123	THR
1	A	129	LYS
1	A	180	SER
1	A	223	HIS
1	A	225	GLU
1	A	236	MET
1	A	237	GLN
1	A	261	ILE
1	A	284	LEU
1	A	303	SER
1	A	313	ARG

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Mol	Chain	Res	Type
1	A	322	LYS
1	A	328	LEU
1	A	365	TRP
1	A	372	THR
1	A	389	ARG
1	A	409	LEU
1	A	419	LEU
1	A	502	ARG
1	A	534	GLU
1	A	555	LEU
1	A	588	GLU
1	A	592	ASN
1	A	611	SER
1	A	621	LEU
1	A	623	LEU
1	A	624	ILE
1	B	41	VAL
1	B	66	ILE
1	B	71	GLU
1	B	72	GLU
1	B	73	GLU
1	B	86	LEU
1	B	95	SER
1	B	116	LEU
1	B	123	THR
1	B	136	VAL
1	B	196	GLN
1	B	257	ILE
1	B	261	ILE
1	B	284	LEU
1	B	300	GLN
1	B	309	GLN
1	B	313	ARG
1	B	326	LYS
1	B	328	LEU
1	B	342	SER
1	B	365	TRP
1	B	388	ASN
1	B	409	LEU
1	B	420	ASN
1	B	436	LEU
1	B	492	LYS

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Mol	Chain	Res	Type
1	B	502	ARG
1	B	555	LEU
1	B	580	THR
1	B	585	SER
1	B	624	ILE

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

Mol	Chain	Res	Type
1	A	52	ASN
1	A	167	ASN
1	A	172	ASN
1	A	201	GLN
1	A	223	HIS
1	A	249	ASN
1	A	442	HIS
1	A	537	ASN
1	A	592	ASN
1	B	47	ASN
1	B	52	ASN
1	B	249	ASN
1	B	309	GLN
1	B	388	ASN
1	B	420	ASN
1	B	442	HIS
1	B	461	ASN
1	B	614	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](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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	OAN	B	1625	-	23,26,26	1.42	3 (13%)	24,35,35	1.77	6 (25%)
2	OAN	A	1625	-	23,26,26	1.40	3 (13%)	24,35,35	1.23	3 (12%)

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	OAN	B	1625	-	-	3/12/35/35	0/2/2/2
2	OAN	A	1625	-	-	3/12/35/35	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1625	OAN	CAS-NAO	-3.49	1.34	1.41
2	B	1625	OAN	OAQ-NAY	-3.07	1.37	1.44
2	A	1625	OAN	CAA-NAY	3.04	1.34	1.27
2	B	1625	OAN	CAH-CAG	2.95	1.56	1.50
2	A	1625	OAN	CAC-CAB	2.67	1.56	1.52
2	A	1625	OAN	OAQ-NAY	-2.40	1.39	1.44

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1625	OAN	CAS-NAO-CAP	-5.99	116.91	126.39
2	B	1625	OAN	CAC-CAB-NAI	-3.78	106.18	112.26
2	A	1625	OAN	CAF-CAE-CAD	2.53	119.23	113.02
2	B	1625	OAN	OAR-CAP-NAO	-2.45	120.92	126.12
2	B	1625	OAN	OAL-CAA-CAB	2.27	124.54	118.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1625	OAN	CAC-CAB-NAI	-2.18	108.75	112.26
2	B	1625	OAN	CAB-NAI-CAG	2.07	125.37	121.82
2	B	1625	OAN	OAL-CAE-CAF	2.06	108.12	106.05
2	A	1625	OAN	OAL-CAE-CAF	2.04	108.10	106.05

There are no chirality outliers.

All (6) torsion outliers are listed below:

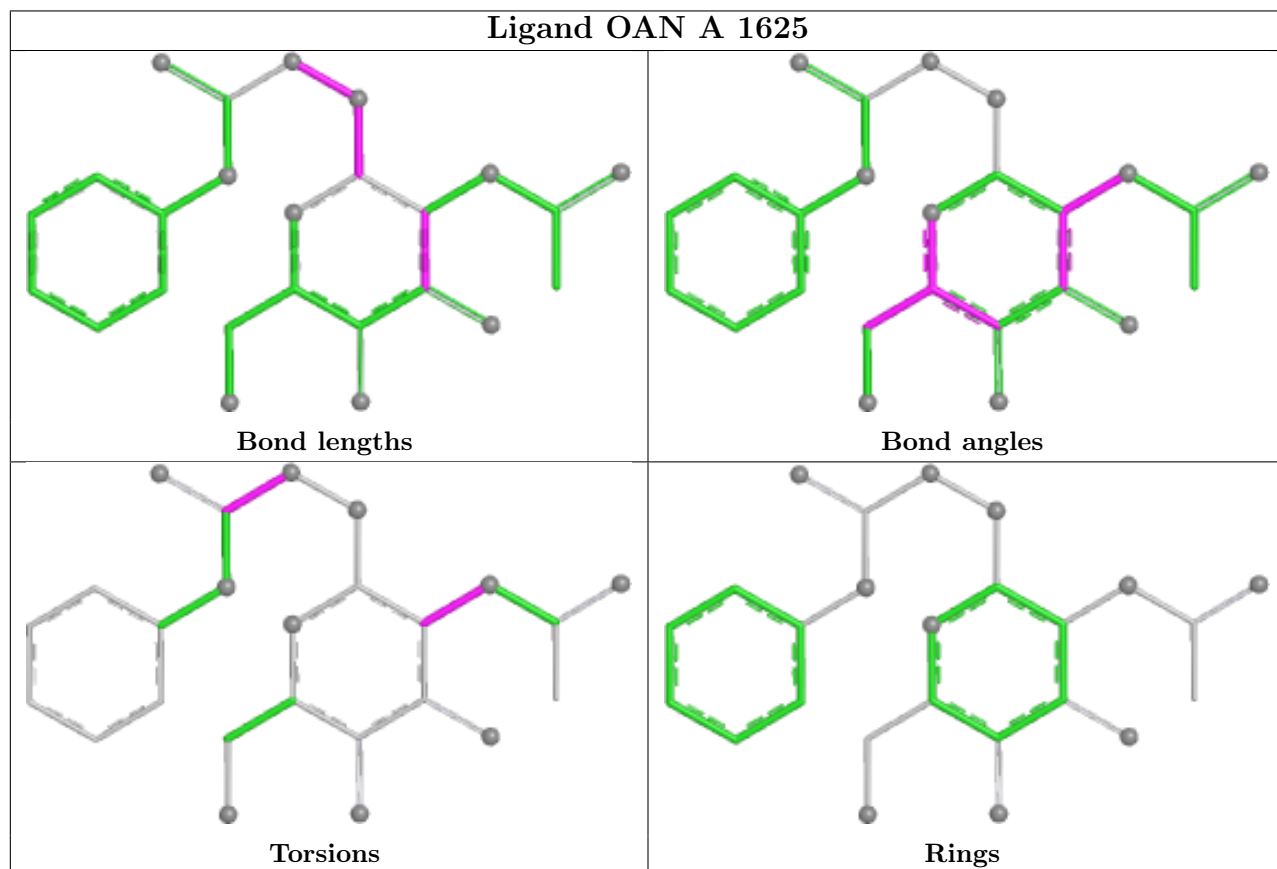
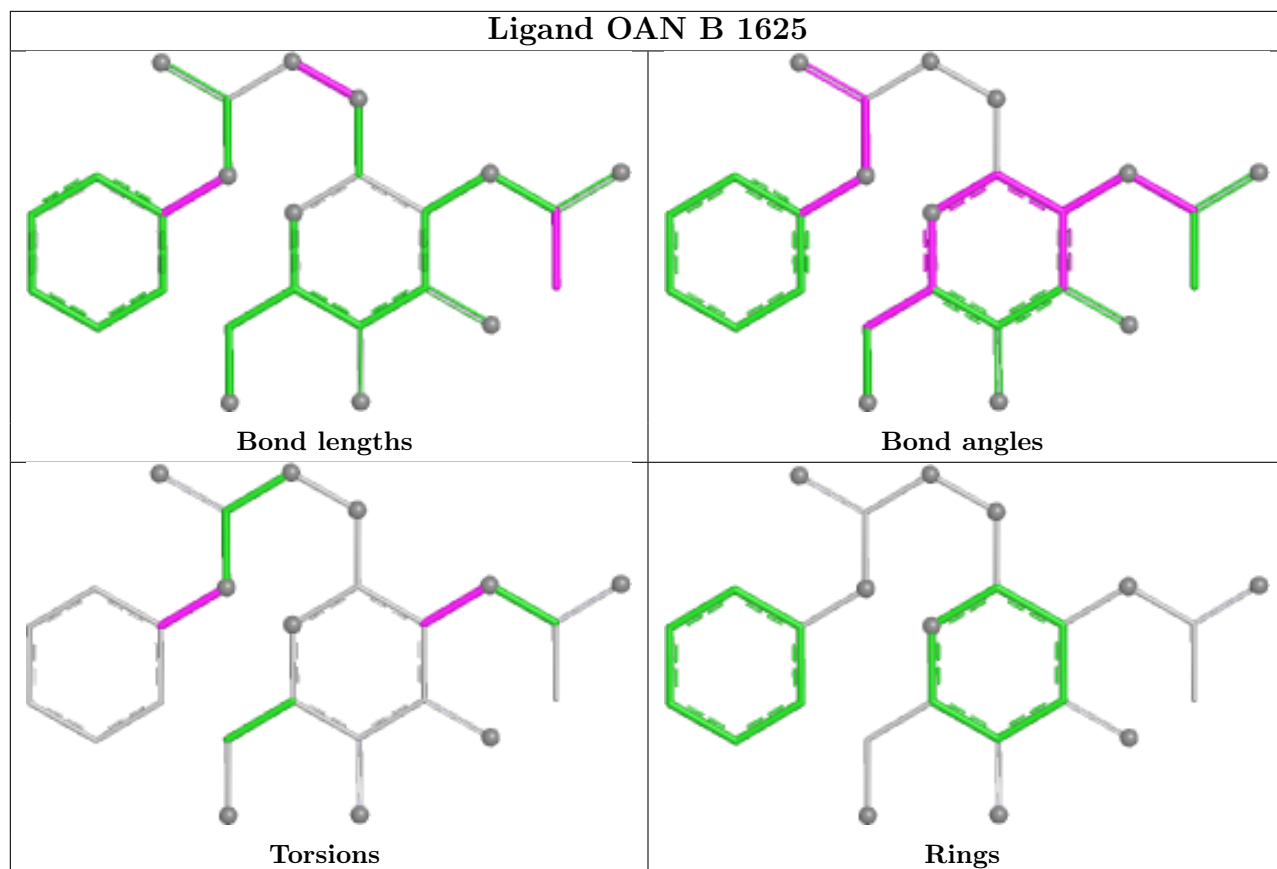
Mol	Chain	Res	Type	Atoms
2	A	1625	OAN	OAR-CAP-OAQ-NAY
2	B	1625	OAN	CAT-CAS-NAO-CAP
2	B	1625	OAN	CAX-CAS-NAO-CAP
2	A	1625	OAN	CAA-CAB-NAI-CAG
2	B	1625	OAN	CAA-CAB-NAI-CAG
2	A	1625	OAN	NAO-CAP-OAQ-NAY

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1625	OAN	1	0
2	A	1625	OAN	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	585/594 (98%)	-0.12	17 (2%) 53 60	1, 44, 71, 85	14 (2%)
1	B	585/594 (98%)	-0.17	11 (1%) 66 71	1, 45, 78, 96	10 (1%)
All	All	1170/1188 (98%)	-0.15	28 (2%) 59 66	1, 45, 76, 96	24 (2%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	GLN	24.9
1	B	40	GLN	16.5
1	A	300	GLN	9.5
1	A	581	GLU	8.0
1	A	72	GLU	7.0
1	A	235	GLU	6.5
1	A	233	GLU	6.4
1	A	266	ASP	6.2
1	B	61	LYS	5.8
1	A	76	GLU	5.4
1	B	73	GLU	5.4
1	B	121	ASN	5.3
1	A	112	ASP	5.1
1	A	61	LYS	4.5
1	B	72	GLU	4.5
1	A	84	GLU	4.3
1	A	99	PRO	4.2
1	B	300	GLN	4.1
1	B	96	GLU	4.1
1	A	225	GLU	3.9
1	B	54	GLU	3.3
1	B	118	GLU	3.0
1	A	624	ILE	3.0
1	A	41	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	109	VAL	2.9
1	A	298	ASP	2.7
1	A	100	ASN	2.7
1	B	41	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

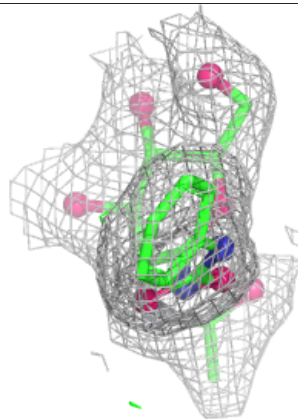
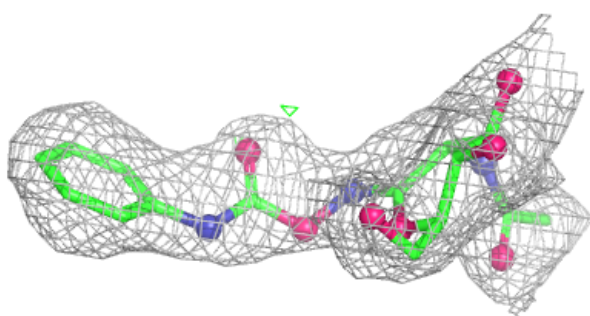
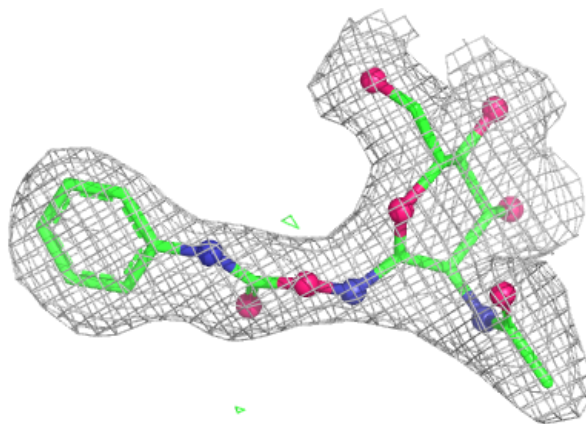
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
3	CL	B	1627	1/1	0.78	0.25	59,59,59,59	1
3	CL	B	1626	1/1	0.95	0.19	63,63,63,63	0
2	OAN	A	1625	25/25	0.95	0.07	25,37,49,54	0
2	OAN	B	1625	25/25	0.96	0.08	22,35,68,70	0

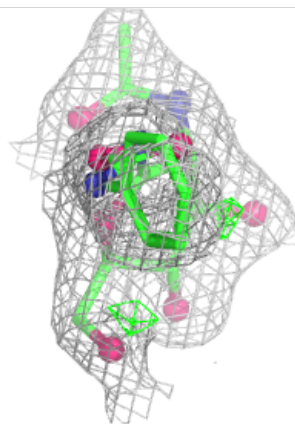
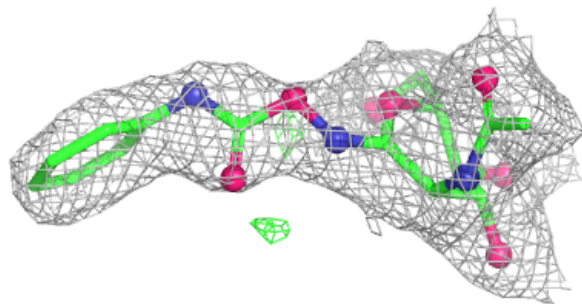
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OAN A 1625:

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

**Electron density around OAN B 1625:**

$2mF_o-DF_c$ (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.