



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:53 PM UTC

PDB ID : 6FLD / pdb_00006fld
Title : Carbamylated T. californica acetylcholinesterase bound to uncharged hybrid reactivator 1
Authors : De la Mora, E.; Santoni, G.; de Souza, J.; Sussman, J.; Silman, I.; Baati, R.; Weik, M.; Nachon, F.
Deposited on : 2018-01-25
Resolution : 2.40 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

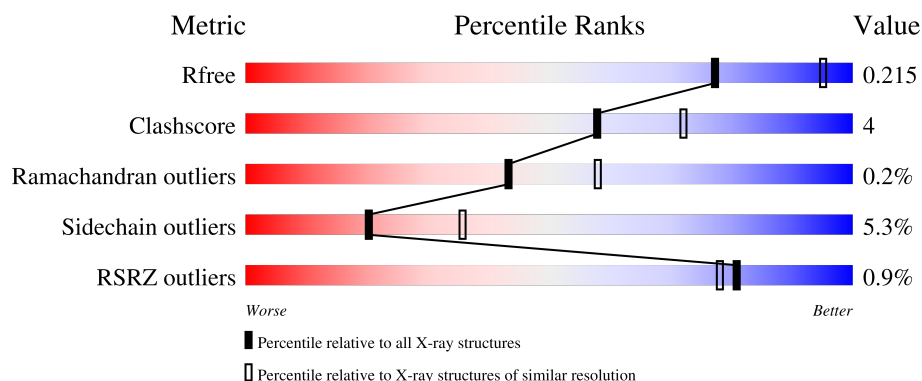
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	532	% 86% 11% .

2 Entry composition [i](#)

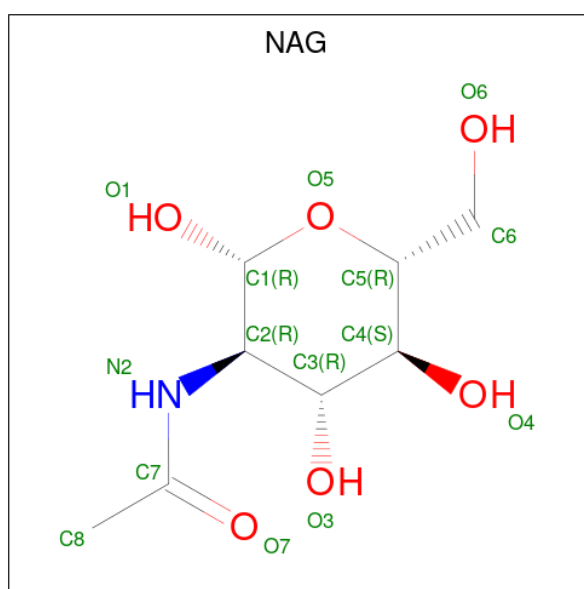
There are 7 unique types of molecules in this entry. The entry contains 4575 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Acetylcholinesterase.

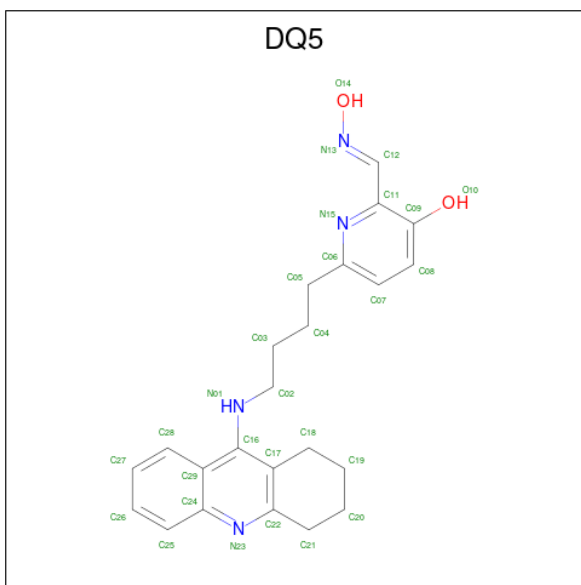
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	532	Total	C	N	O	S	0	3	0
			4259	2733	722	781	23			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



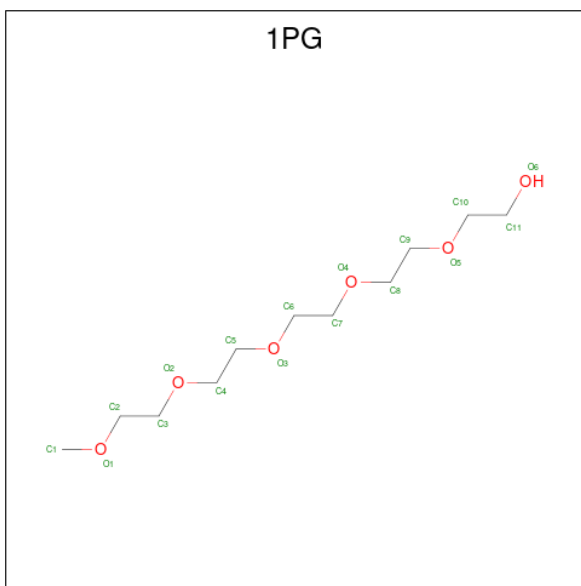
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is 2-[(E)-hydroxyiminomethyl]-6-[4-(1,2,3,4-tetrahydroacridin-9-ylamino)butyl]pyridin-3-ol (CCD ID: DQ5) (formula: $C_{23}H_{26}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	23	4	2		

- Molecule 4 is 2-(2-{2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXY}-ETHOXY)-ETHANOL (CCD ID: 1PG) (formula: $C_{11}H_{24}O_6$).

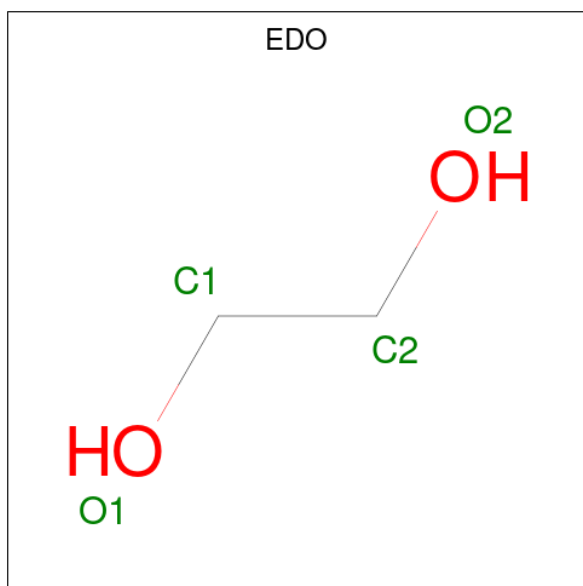


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	5	Total Cl 5 5	0	0

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0

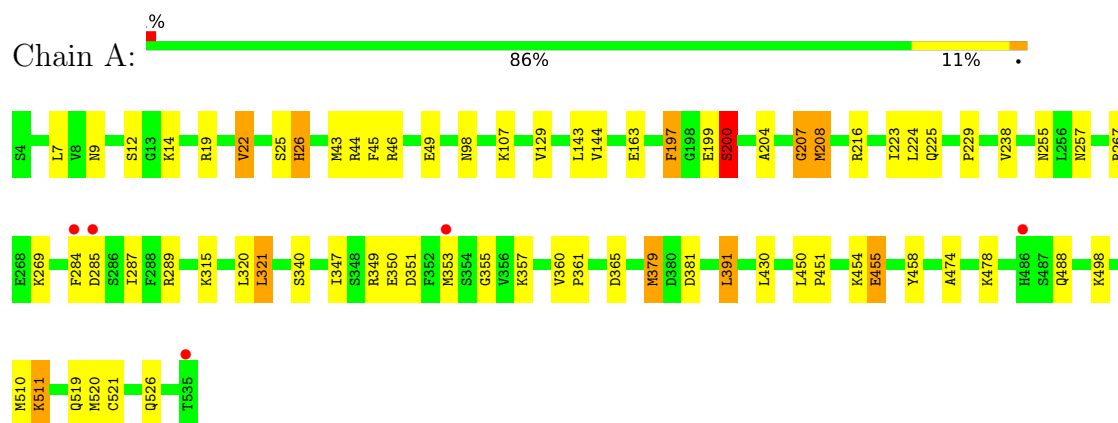
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	212	Total O 212 212	0	0

3 Residue-property plots

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: Acetylcholinesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.01Å 114.01Å 136.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.32 – 2.40 37.32 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (37.32-2.40) 90.2 (37.32-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.64 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.180 , 0.210 0.184 , 0.215	Depositor DCC
R_{free} test set	1992 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.705	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4575	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 1PG, BXT, DQ5, CL, EDO

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	0.59	3/4372 (0.1%)	0.59	3/5937 (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	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	450	LEU	C-O	-5.88	1.19	1.24
1	A	208	MET	C-O	-5.14	1.18	1.24
1	A	207	GLY	C-O	-5.01	1.17	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	521	CYS	CB-CA-C	7.17	123.04	110.85
1	A	269	LYS	N-CA-C	6.76	120.04	110.50
1	A	454	LYS	N-CA-C	5.31	117.07	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	200	BXT	Mainchain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4259	0	4080	36	0
2	A	42	0	39	0	0
3	A	29	0	0	0	0
4	A	16	0	21	2	0
5	A	5	0	0	0	0
6	A	12	0	18	0	0
7	A	212	0	0	6	0
All	All	4575	0	4158	38	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 4.

All (38) 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:455:GLU:N	1:A:455:GLU:OE1	1.88	1.06
1:A:520:MET:HG2	7:A:856:HOH:O	1.85	0.77
1:A:455:GLU:H	1:A:455:GLU:CD	1.97	0.70
1:A:44:ARG:O	1:A:45:PHE:HB2	1.92	0.70
1:A:355:GLY:HA3	1:A:391:LEU:HD11	1.75	0.69
1:A:360:VAL:O	1:A:360:VAL:HG23	2.00	0.61
1:A:204:ALA:O	1:A:208:MET:HG3	2.04	0.57
1:A:285:ASP:HB2	1:A:361:PRO:HG2	1.87	0.57
1:A:451:PRO:HA	1:A:458:TYR:CD1	2.40	0.56
1:A:350:GLU:HA	1:A:353[B]:MET:HE2	1.86	0.56
1:A:43:MET:HE1	1:A:49:GLU:HG2	1.89	0.54
1:A:44:ARG:O	1:A:45:PHE:CB	2.54	0.54
1:A:200:BXT:N	1:A:225:GLN:O	2.42	0.52
1:A:347:ILE:HG23	1:A:351:ASP:HB3	1.92	0.52
1:A:349:ARG:NH2	1:A:381:ASP:OD2	2.43	0.52
1:A:285:ASP:CB	1:A:361:PRO:HG2	2.40	0.51
1:A:207:GLY:HA3	1:A:229:PRO:HD3	1.95	0.49
1:A:255:ASN:HD21	1:A:257:ASN:CG	2.20	0.49
4:A:605:1PG:H101	7:A:805:HOH:O	2.14	0.46
1:A:9:ASN:OD1	1:A:14:LYS:HG2	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:605:1PG:H61	7:A:805:HOH:O	2.16	0.45
1:A:197:PHE:HB3	1:A:223:ILE:HB	1.99	0.45
1:A:107:LYS:N	7:A:706:HOH:O	2.43	0.44
1:A:46:ARG:HD2	7:A:701:HOH:O	2.17	0.44
1:A:474:ALA:O	1:A:478:LYS:HG3	2.19	0.43
1:A:197:PHE:CB	1:A:223:ILE:HB	2.48	0.43
1:A:223:ILE:HG12	1:A:320:LEU:HB3	2.01	0.43
1:A:224:LEU:HB2	1:A:321:LEU:HB3	2.00	0.43
1:A:285:ASP:OD1	1:A:285:ASP:N	2.41	0.42
1:A:25:SER:OG	1:A:26:HIS:N	2.51	0.42
1:A:216:ARG:HB3	1:A:315:LYS:HB2	2.01	0.42
1:A:498:LYS:HE3	1:A:498:LYS:HB2	1.81	0.42
1:A:19:ARG:HE	1:A:26:HIS:HB2	1.85	0.42
1:A:511:LYS:HE2	7:A:853:HOH:O	2.20	0.41
1:A:379:MET:HE2	1:A:379:MET:HB3	1.75	0.41
1:A:98:ASN:O	1:A:144:VAL:HA	2.21	0.41
1:A:163:GLU:HB3	1:A:267:ARG:HH22	1.87	0.40
1:A:199:GLU:O	1:A:200:BXT:C	2.69	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	532/532 (100%)	508 (96%)	23 (4%)	1 (0%)	43 58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	VAL

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	460/464 (99%)	436 (95%)	24 (5%)	21	36

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	12	SER
1	A	22	VAL
1	A	26	HIS
1	A	129	VAL
1	A	143	LEU
1	A	197	PHE
1	A	238	VAL
1	A	284	PHE
1	A	287	ILE
1	A	289	ARG
1	A	321	LEU
1	A	340	SER
1	A	357	LYS
1	A	365	ASP
1	A	379	MET
1	A	391	LEU
1	A	430	LEU
1	A	455	GLU
1	A	488	GLN
1	A	510	MET
1	A	511	LYS
1	A	519	GLN
1	A	526	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	255	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	BXT	A	200	1	9,11,12	2.43	4 (44%)	6,13,15	3.39	2 (33%)

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
1	BXT	A	200	1	-	5/11/13/15	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	200	BXT	O2-C1	-4.30	1.14	1.21
1	A	200	BXT	O-C	4.12	1.35	1.20
1	A	200	BXT	OG-CB	-2.38	1.39	1.45
1	A	200	BXT	C5-N3	-2.10	1.38	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	BXT	OG-C1-N3	7.25	118.29	111.12
1	A	200	BXT	CB-OG-C1	3.27	121.36	115.81

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	200	BXT	C-CA-CB-OG
1	A	200	BXT	OG-C1-N3-C5
1	A	200	BXT	O2-C1-N3-C5
1	A	200	BXT	C2-C4-N3-C1
1	A	200	BXT	N-CA-CB-OG

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	200	BXT	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 5 are monoatomic - leaving 8 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	1PG	A	605	-	15,15,16	1.03	1 (6%)	14,14,15	0.65	0
6	EDO	A	612	-	3,3,3	0.49	0	2,2,2	0.41	0
2	NAG	A	601	1	14,14,15	0.98	1 (7%)	17,19,21	0.68	1 (5%)
6	EDO	A	611	-	3,3,3	0.51	0	2,2,2	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	A	613	-	3,3,3	0.42	0	2,2,2	0.46	0
2	NAG	A	603	1	14,14,15	0.56	0	17,19,21	0.71	1 (5%)
2	NAG	A	602	1	14,14,15	0.36	0	17,19,21	0.91	1 (5%)
3	DQ5	A	604	-	31,32,32	2.05	6 (19%)	36,43,43	1.69	12 (33%)

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	1PG	A	605	-	-	6/13/13/14	-
6	EDO	A	612	-	-	1/1/1/1	-
2	NAG	A	601	1	-	0/6/23/26	0/1/1/1
6	EDO	A	611	-	-	1/1/1/1	-
6	EDO	A	613	-	-	1/1/1/1	-
2	NAG	A	603	1	-	2/6/23/26	0/1/1/1
2	NAG	A	602	1	-	0/6/23/26	0/1/1/1
3	DQ5	A	604	-	-	5/11/18/18	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	604	DQ5	C22-N23	6.42	1.40	1.32
3	A	604	DQ5	C16-C17	5.11	1.47	1.39
3	A	604	DQ5	C16-N01	3.70	1.48	1.37
2	A	601	NAG	O5-C1	3.22	1.49	1.43
3	A	604	DQ5	O10-C09	3.05	1.42	1.36
3	A	604	DQ5	C06-N15	2.43	1.38	1.34
3	A	604	DQ5	C29-C24	-2.14	1.39	1.42
4	A	605	1PG	C10-C11	2.05	1.60	1.49

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	604	DQ5	C05-C06-N15	3.34	121.12	116.06
3	A	604	DQ5	C21-C22-C17	3.31	124.78	121.40
2	A	602	NAG	C1-O5-C5	3.20	116.47	112.19
3	A	604	DQ5	C21-C22-N23	-2.83	112.37	116.75
3	A	604	DQ5	C29-C24-N23	-2.81	119.82	122.80
3	A	604	DQ5	C16-C17-C22	-2.69	117.03	118.68

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	604	DQ5	C22-N23-C24	2.63	120.78	117.67
3	A	604	DQ5	C16-C29-C24	2.58	121.10	117.27
2	A	603	NAG	C1-O5-C5	2.48	115.50	112.19
3	A	604	DQ5	O14-N13-C12	2.28	115.78	111.83
3	A	604	DQ5	C28-C29-C16	-2.28	120.60	124.70
2	A	601	NAG	C1-O5-C5	2.07	114.96	112.19
3	A	604	DQ5	O10-C09-C11	-2.05	115.51	120.44
3	A	604	DQ5	C18-C17-C22	-2.03	118.99	120.90
3	A	604	DQ5	C07-C06-N15	-2.01	119.86	122.40

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	605	1PG	O4-C8-C9-O5
4	A	605	1PG	C4-C5-O3-C6
3	A	604	DQ5	N01-C02-C03-C04
2	A	603	NAG	C4-C5-C6-O6
3	A	604	DQ5	C02-C03-C04-C05
4	A	605	1PG	O1-C2-C3-O2
6	A	612	EDO	O1-C1-C2-O2
4	A	605	1PG	O3-C6-C7-O4
2	A	603	NAG	O5-C5-C6-O6
6	A	611	EDO	O1-C1-C2-O2
4	A	605	1PG	C5-C4-O2-C3
3	A	604	DQ5	C03-C02-N01-C16
3	A	604	DQ5	C17-C16-N01-C02
6	A	613	EDO	O1-C1-C2-O2
3	A	604	DQ5	N15-C11-C12-N13
4	A	605	1PG	O5-C10-C11-O6

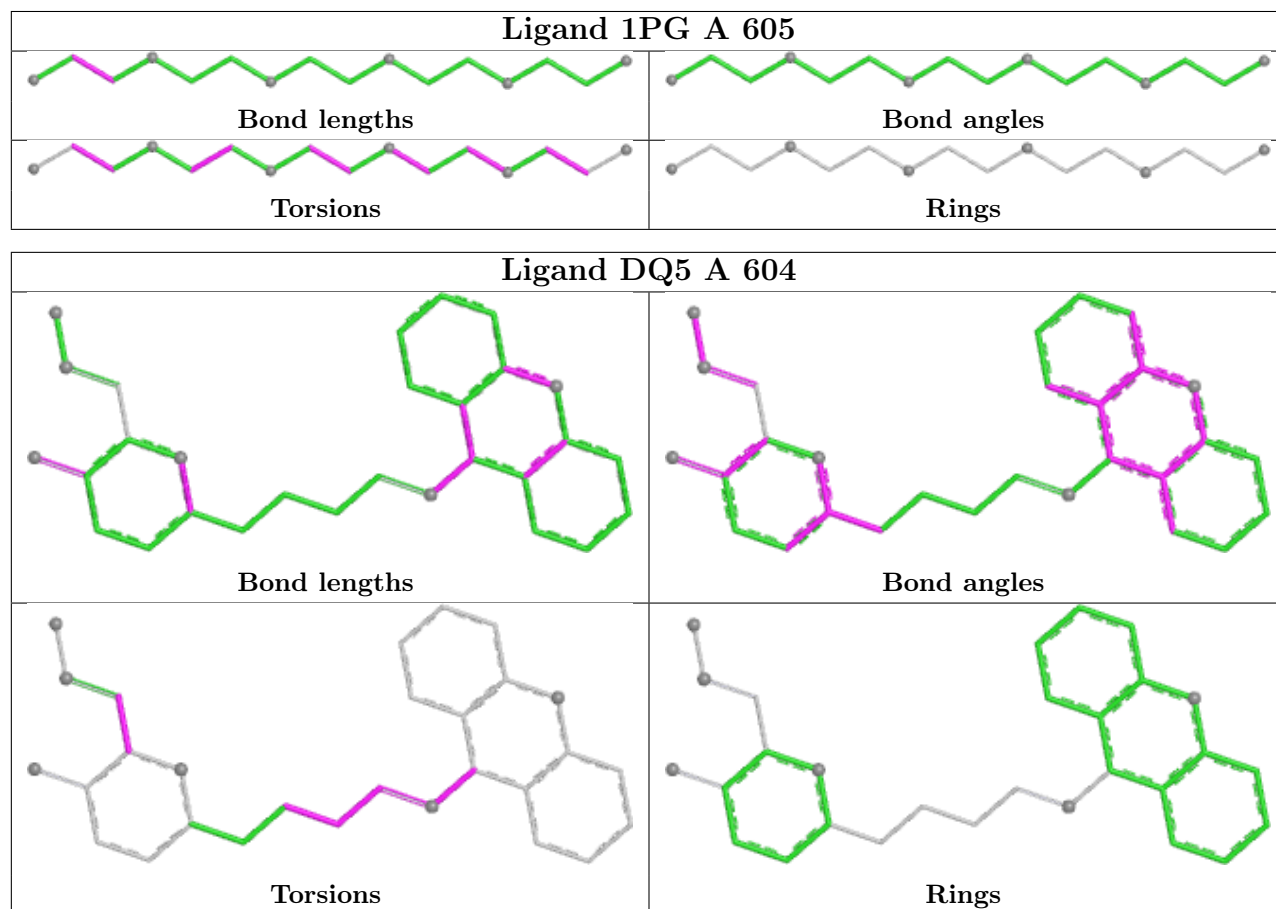
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	605	1PG	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 [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	531/532 (99%)	-0.18	5 (0%) 81 78	26, 49, 78, 114	5 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	353[A]	MET	3.9
1	A	284	PHE	3.8
1	A	486	HIS	3.6
1	A	285	ASP	2.5
1	A	535	THR	2.4

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
1	BXT	A	200	12/13	0.95	0.09	38,43,51,51	0

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,

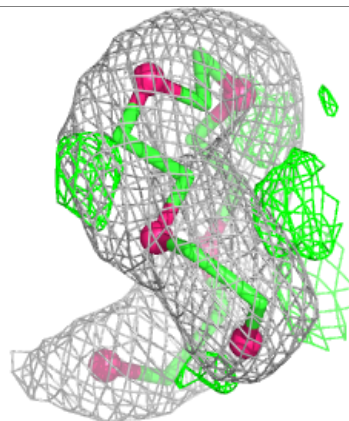
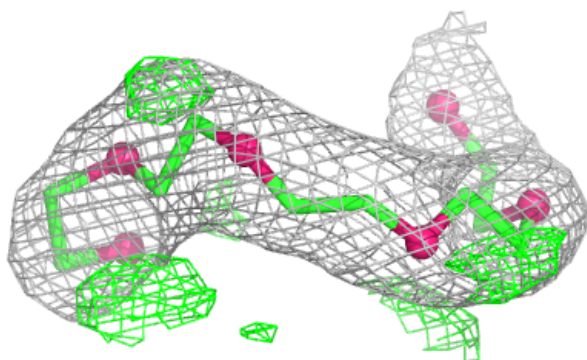
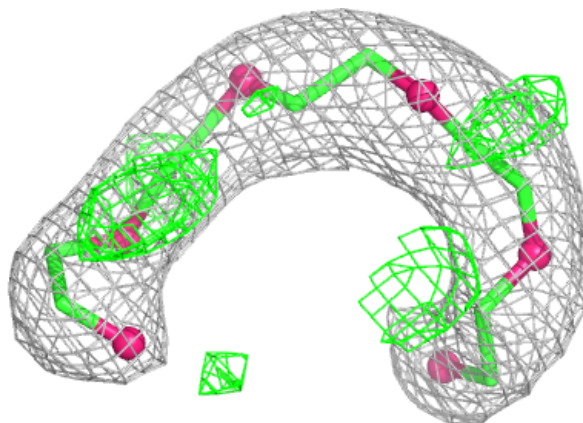
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	A	601	14/15	0.67	0.16	77,86,94,99	0
6	EDO	A	613	4/4	0.70	0.22	72,72,76,78	0
2	NAG	A	603	14/15	0.76	0.13	75,87,94,97	0
6	EDO	A	611	4/4	0.79	0.21	62,67,70,71	0
5	CL	A	607	1/1	0.83	0.10	96,96,96,96	0
6	EDO	A	612	4/4	0.84	0.36	51,58,61,62	0
2	NAG	A	602	14/15	0.86	0.10	50,55,64,69	0
5	CL	A	609	1/1	0.87	0.14	75,75,75,75	0
5	CL	A	608	1/1	0.87	0.12	73,73,73,73	0
4	1PG	A	605	16/17	0.93	0.13	38,48,59,60	0
3	DQ5	A	604	29/29	0.93	0.11	39,45,52,63	0
5	CL	A	610	1/1	0.93	0.17	82,82,82,82	0
5	CL	A	606	1/1	0.95	0.07	68,68,68,68	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.

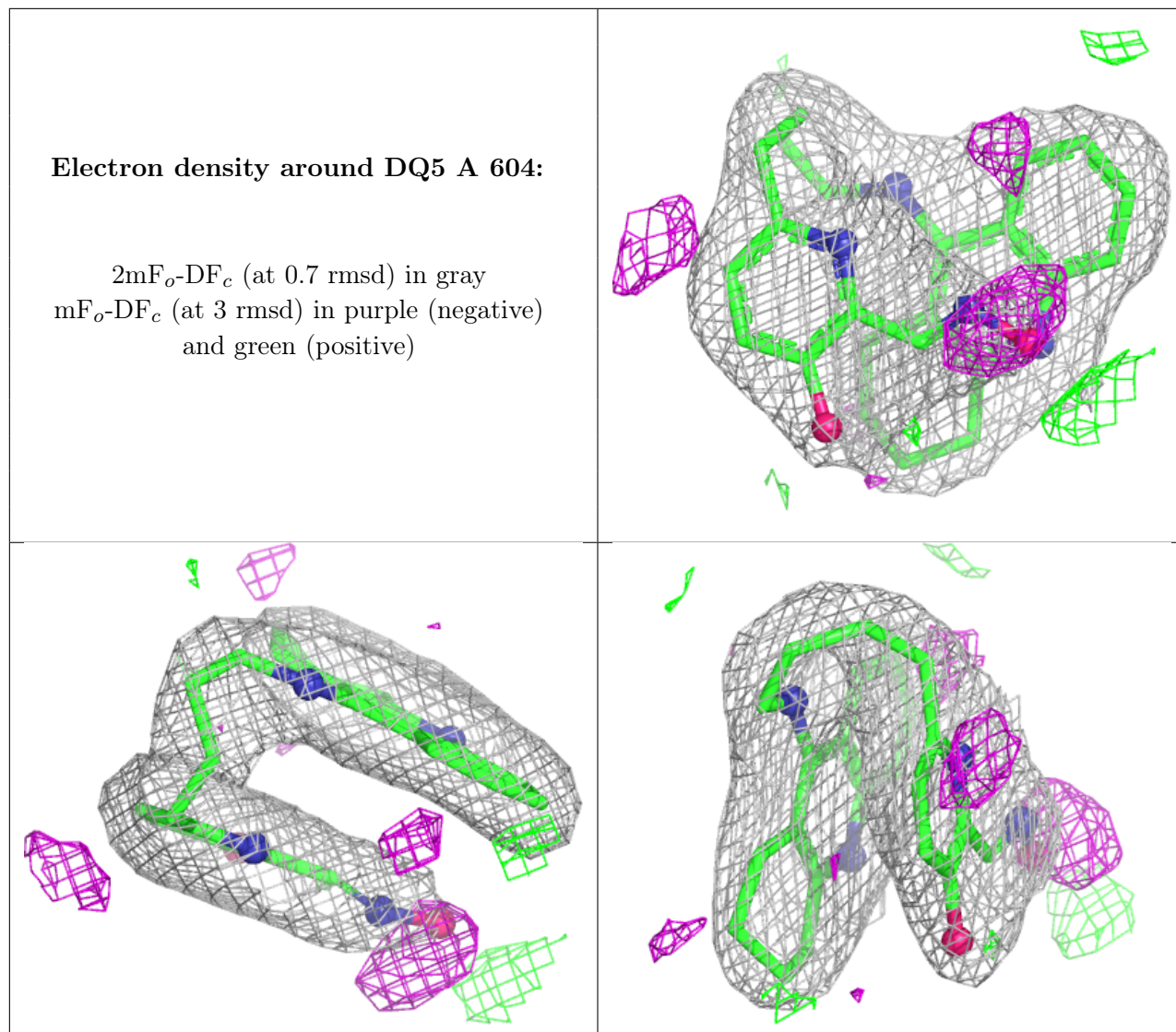
Electron density around 1PG A 605:

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 DQ5 A 604:

$2mF_o-DF_c$ (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.