



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 01:03 PM EDT

PDB ID : 6BO6 / pdb_00006bo6
Title : Eubacterium eligens beta-glucuronidase bound to UNC4917 glucuronic acid conjugate
Authors : Pellock, S.J.; Walton, W.G.; Redinbo, M.R.
Deposited on : 2017-11-18
Resolution : 2.80 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

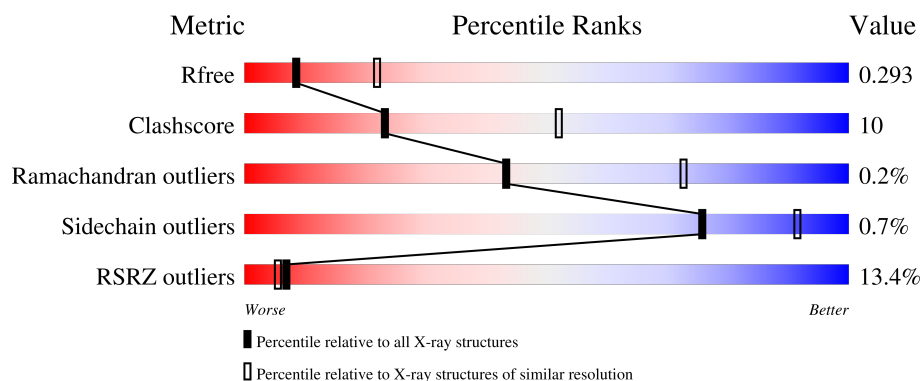
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	611	11% 61% 19% 19%

2 Entry composition [i](#)

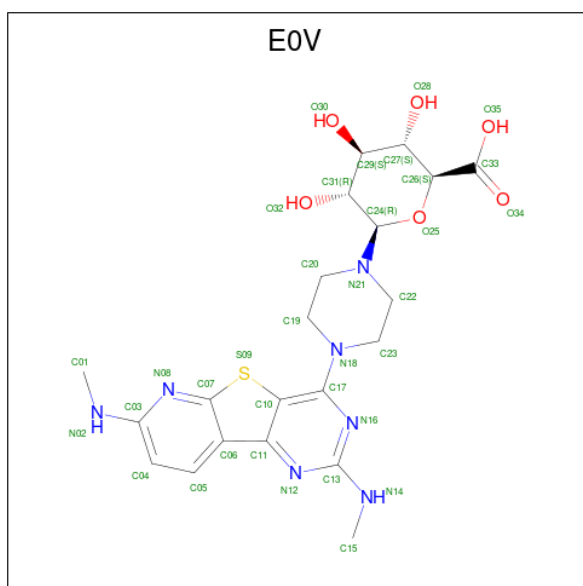
There are 3 unique types of molecules in this entry. The entry contains 4036 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycoside Hydrolase Family 2 candidate b-glucuronidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			3982	2545	663	753	21			

- Molecule 2 is 4-(4-beta-D-glucopyranuronosylpiperazin-1-yl)-2,7-bis(methylamino)pyrido[3',2':4,5]thieno[3,2-d]pyrimidine (CCD ID: E0V) (formula: C₂₁H₂₇N₇O₆S) (labeled as "Ligand of Interest" by depositor).



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	179.44Å 179.44Å 133.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.37 – 2.80 29.37 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.37-2.80) 97.6 (29.37-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.80Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.258 , 0.292 0.260 , 0.293	Depositor DCC
R_{free} test set	2000 reflections (6.30%)	wwPDB-VP
Wilson B-factor (Å ²)	65.1	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4036	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: E0V

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	0.23	0/4089	0.43	0/5538

There are no bond length outliers.

There are no bond angle outliers.

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	3982	0	3771	80	0
2	A	35	0	0	0	0
3	A	19	0	0	1	0
All	All	4036	0	3771	80	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 10.

All (80) 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:288:ARG:NH2	1:A:298:GLU:OE1	2.01	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ARG:HH22	1:A:298:GLU:CD	1.77	0.92
1:A:448:ASP:OD1	1:A:452:ARG:NH1	2.11	0.83
1:A:25:ASN:N	1:A:26:GLY:HA2	2.01	0.76
1:A:377:GLY:HA3	1:A:384:ARG:NH1	2.02	0.74
1:A:93:TYR:OH	1:A:105:GLU:OE2	2.06	0.74
1:A:408:ILE:O	1:A:412:LYS:HG2	1.87	0.73
1:A:266:LEU:HD11	1:A:355:ASP:HB3	1.72	0.72
1:A:462:THR:HG21	1:A:467:ASP:OD2	1.90	0.70
1:A:104:CYS:HB2	1:A:115:VAL:HG11	1.73	0.70
1:A:119:ASP:OD2	1:A:120:ASP:N	2.25	0.69
1:A:522:VAL:HB	1:A:525:LEU:HD23	1.73	0.69
1:A:258:LEU:HD23	1:A:284:VAL:HG23	1.75	0.68
1:A:424:ASN:HA	1:A:457:VAL:HG23	1.77	0.66
1:A:110:PHE:HE1	1:A:317:ARG:HD2	1.61	0.65
1:A:303:PHE:HB2	1:A:557:GLY:HA3	1.79	0.63
1:A:441:TYR:OH	1:A:473:SER:OG	2.15	0.59
1:A:260:GLN:OE1	1:A:297:ASN:ND2	2.37	0.58
1:A:25:ASN:H	1:A:26:GLY:HA2	1.68	0.58
1:A:499:ARG:NH1	1:A:551:GLU:OE2	2.36	0.58
1:A:334:GLN:HA	1:A:597:ARG:HG3	1.87	0.56
1:A:425:GLU:HB3	1:A:459:VAL:HG22	1.86	0.56
1:A:288:ARG:NH1	1:A:298:GLU:OE2	2.38	0.55
1:A:169:PRO:HB3	1:A:171:PHE:CZ	2.41	0.55
1:A:37:PRO:HA	1:A:69:GLN:HE22	1.72	0.55
1:A:562:ASN:OD1	1:A:563:PHE:N	2.40	0.55
1:A:368:VAL:HG23	1:A:424:ASN:HD22	1.72	0.54
1:A:420:TRP:HZ2	1:A:452:ARG:NH1	2.06	0.53
1:A:121:LEU:HD22	1:A:126:ASN:HD22	1.73	0.53
1:A:72:ILE:HG22	1:A:126:ASN:HB2	1.90	0.53
1:A:462:THR:HG23	1:A:467:ASP:HB2	1.89	0.53
1:A:424:ASN:HA	1:A:457:VAL:CG2	2.40	0.51
1:A:457:VAL:HG12	1:A:477:CYS:HB3	1.94	0.50
1:A:305:GLY:HA3	1:A:338:SER:O	2.12	0.50
1:A:69:GLN:OE1	1:A:129:THR:OG1	2.28	0.50
1:A:512:VAL:O	1:A:555:VAL:HA	2.12	0.50
1:A:370:VAL:HG22	3:A:805:HOH:O	2.12	0.49
1:A:101:LYS:HD3	1:A:120:ASP:OD2	2.12	0.49
1:A:484:TRP:CZ2	1:A:520:ASP:HB2	2.46	0.49
1:A:434:TYR:CE1	1:A:472:LEU:HD11	2.47	0.49
1:A:98:LEU:HB2	1:A:103:ILE:HD11	1.93	0.48
1:A:288:ARG:HH21	1:A:288:ARG:HG2	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASP:OD2	1:A:26:GLY:HA2	2.14	0.48
1:A:328:ILE:HG21	1:A:357:GLU:HG3	1.94	0.48
1:A:420:TRP:CZ2	1:A:452:ARG:NH1	2.81	0.48
1:A:133:ASN:ND2	1:A:135:VAL:H	2.11	0.48
1:A:102:LEU:HD12	1:A:102:LEU:H	1.78	0.48
1:A:407:LEU:HD22	1:A:420:TRP:HZ3	1.80	0.47
1:A:348:GLU:O	1:A:352:ARG:HG3	2.14	0.47
1:A:370:VAL:O	1:A:370:VAL:HG23	2.15	0.47
1:A:362:ILE:HD11	1:A:513:MET:HE1	1.97	0.47
1:A:97:TYR:CE1	1:A:102:LEU:HG	2.51	0.46
1:A:401:LYS:H	1:A:401:LYS:HG2	1.59	0.46
1:A:260:GLN:HB2	1:A:263:ASN:HB3	1.97	0.46
1:A:295:LEU:HA	1:A:299:LYS:O	2.16	0.46
1:A:304:LYS:NZ	1:A:603:ILE:O	2.41	0.46
1:A:482:TYR:OH	1:A:501:GLU:OE1	2.27	0.46
1:A:561:TRP:HA	1:A:562:ASN:HA	1.62	0.45
1:A:25:ASN:N	1:A:25:ASN:OD1	2.49	0.45
1:A:91:THR:HA	1:A:92:HIS:HA	1.62	0.45
1:A:315:ASN:OD1	1:A:315:ASN:N	2.50	0.45
1:A:341:THR:HA	1:A:346:TYR:CZ	2.52	0.44
1:A:342:SER:HA	1:A:343:HIS:HA	1.62	0.44
1:A:90:VAL:O	1:A:109:GLY:HA2	2.17	0.44
1:A:18:VAL:HA	1:A:45:PRO:HA	2.00	0.43
1:A:295:LEU:HD23	1:A:300:PRO:HA	2.01	0.43
1:A:370:VAL:HG23	1:A:426:PRO:HB3	2.00	0.43
1:A:404:ILE:HD11	1:A:422:ILE:HG23	2.00	0.43
1:A:147:ALA:O	1:A:168:ASN:HB3	2.19	0.42
1:A:260:GLN:HE22	1:A:299:LYS:NZ	2.18	0.42
1:A:463:THR:HG23	1:A:466:THR:H	1.84	0.42
1:A:480:ARG:HD3	1:A:480:ARG:HA	1.84	0.41
1:A:174:PHE:O	1:A:317:ARG:NH2	2.52	0.41
1:A:24:ASP:CG	1:A:26:GLY:HA2	2.45	0.41
1:A:23:LEU:HD11	1:A:59:PHE:CD2	2.56	0.41
1:A:38:LEU:CD1	1:A:38:LEU:H	2.34	0.41
1:A:285:ARG:NH1	1:A:450:GLN:OE1	2.32	0.41
1:A:123:ASP:HB3	1:A:124:GLY:H	1.59	0.41
1:A:327:ASP:O	1:A:331:MET:HG3	2.21	0.41
1:A:353:LEU:HD12	1:A:353:LEU:HA	1.95	0.40

There are no symmetry-related clashes.

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	483/611 (79%)	454 (94%)	28 (6%)	1 (0%)	43 72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	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	426/532 (80%)	423 (99%)	3 (1%)	76 91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LEU
1	A	143	VAL
1	A	306	TYR

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	133	ASN
1	A	148	ASN

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Mol	Chain	Res	Type
1	A	392	HIS
1	A	397	GLN
1	A	424	ASN

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](#)

1 ligand 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	$\# Z > 2$
2	E0V	A	701	-	37,39,39	2.88	10 (27%)	43,58,58	2.79	15 (34%)

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	E0V	A	701	-	-	4/16/46/46	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	E0V	C13-N14	12.51	1.44	1.34
2	A	701	E0V	C03-N02	6.32	1.49	1.36
2	A	701	E0V	C17-N18	5.85	1.51	1.37
2	A	701	E0V	C23-N18	2.86	1.51	1.46
2	A	701	E0V	C22-N21	-2.81	1.41	1.47
2	A	701	E0V	O34-C33	2.75	1.30	1.22
2	A	701	E0V	O25-C24	2.64	1.48	1.42
2	A	701	E0V	O25-C26	2.33	1.47	1.43
2	A	701	E0V	C19-N18	2.19	1.50	1.46
2	A	701	E0V	C01-N02	-2.18	1.41	1.45

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	E0V	C06-C07-S09	-9.44	105.49	112.58
2	A	701	E0V	C10-S09-C07	9.08	100.41	90.40
2	A	701	E0V	C11-C10-S09	-6.34	105.31	112.41
2	A	701	E0V	C01-N02-C03	-5.14	118.21	122.96
2	A	701	E0V	C29-C31-C24	4.04	115.38	108.94
2	A	701	E0V	C15-N14-C13	-3.14	119.80	123.59
2	A	701	E0V	N16-C13-N12	-2.92	121.47	126.26
2	A	701	E0V	N14-C13-N12	2.67	119.53	116.96
2	A	701	E0V	C05-C06-C11	-2.45	126.23	131.36
2	A	701	E0V	C05-C06-C07	2.31	119.33	117.11
2	A	701	E0V	C10-C17-N16	-2.23	119.68	122.33
2	A	701	E0V	C13-N16-C17	2.16	120.91	113.99
2	A	701	E0V	O25-C24-C31	2.14	117.64	110.79
2	A	701	E0V	N14-C13-N16	2.12	119.00	116.96
2	A	701	E0V	C22-N21-C20	2.02	112.98	109.13

There are no chirality outliers.

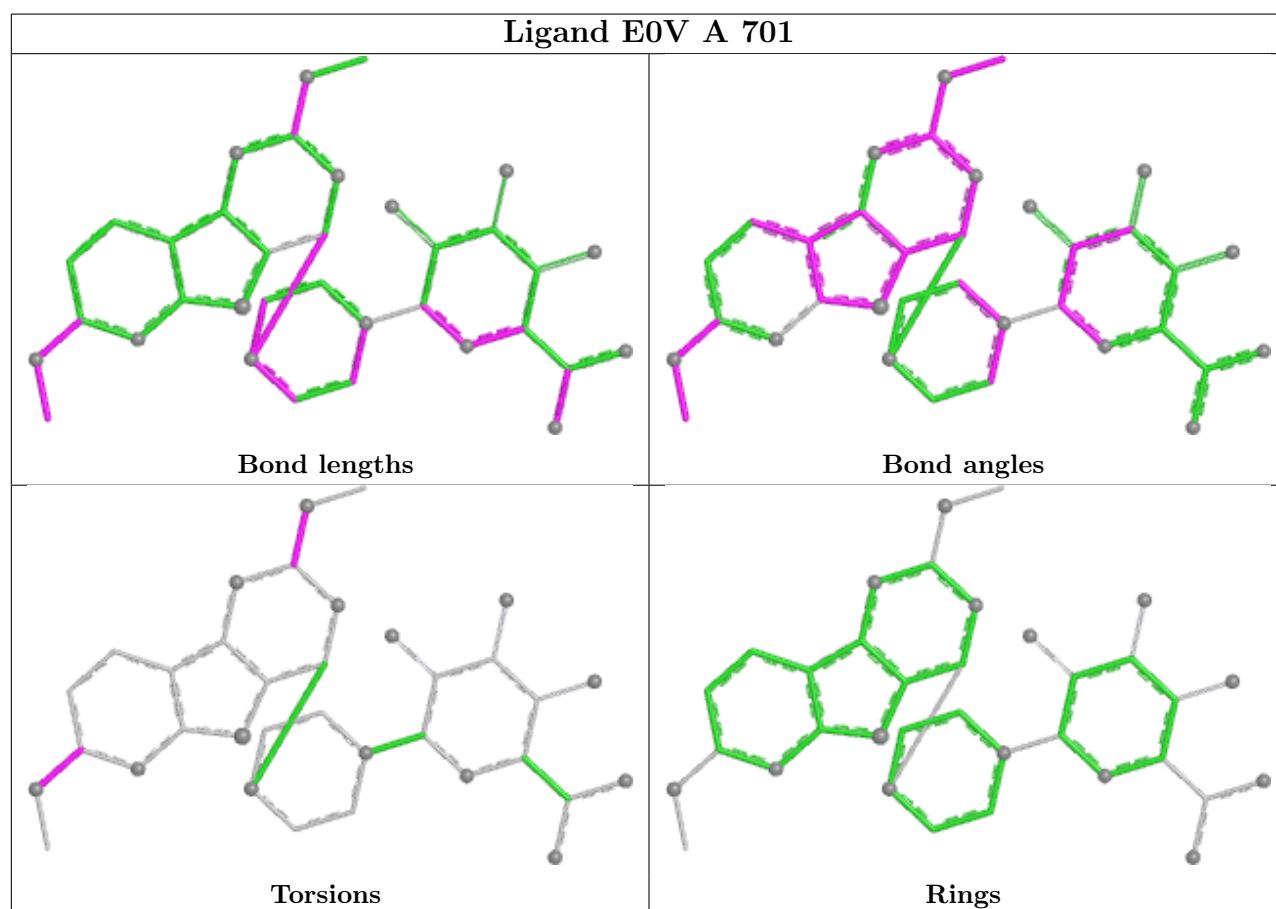
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	E0V	C04-C03-N02-C01
2	A	701	E0V	N08-C03-N02-C01
2	A	701	E0V	N12-C13-N14-C15
2	A	701	E0V	N16-C13-N14-C15

There are no ring outliers.

No monomer is involved in short contacts.

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	493/611 (80%)	0.82	66 (13%) 7 5	35, 67, 103, 123	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	186	TYR	5.1
1	A	73	SER	4.8
1	A	74	VAL	4.5
1	A	32	LYS	4.4
1	A	375	GLY	4.3
1	A	85	LEU	4.3
1	A	148	ASN	4.2
1	A	102	LEU	3.9
1	A	13	SER	3.9
1	A	121	LEU	3.8
1	A	268	LYS	3.8
1	A	98	LEU	3.8
1	A	104	CYS	3.5
1	A	103	ILE	3.5
1	A	99	ASN	3.4
1	A	115	VAL	3.3
1	A	88	ALA	3.3
1	A	258	LEU	3.3
1	A	374	PHE	3.2
1	A	265	TYR	3.1
1	A	39	LYS	3.1
1	A	101	LYS	3.1
1	A	380	PHE	3.1
1	A	36	LYS	3.0
1	A	279	THR	3.0
1	A	449	PRO	2.9
1	A	38	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	117	LEU	2.8
1	A	165	PRO	2.8
1	A	25	ASN	2.8
1	A	24	ASP	2.8
1	A	278	TYR	2.7
1	A	378	ALA	2.7
1	A	259	TRP	2.6
1	A	391	GLU	2.6
1	A	126	ASN	2.6
1	A	86	ARG	2.6
1	A	123	ASP	2.6
1	A	146	LYS	2.5
1	A	69	GLN	2.5
1	A	97	TYR	2.4
1	A	390	LYS	2.4
1	A	127	LEU	2.4
1	A	108	GLY	2.4
1	A	72	ILE	2.4
1	A	412	LYS	2.3
1	A	27	LYS	2.3
1	A	410	ARG	2.3
1	A	409	SER	2.3
1	A	446	GLU	2.2
1	A	96	ILE	2.2
1	A	105	GLU	2.2
1	A	100	GLY	2.2
1	A	116	GLU	2.2
1	A	26	GLY	2.1
1	A	437	PHE	2.1
1	A	34	TYR	2.1
1	A	286	SER	2.1
1	A	288	ARG	2.1
1	A	184	LYS	2.1
1	A	122	GLN	2.1
1	A	295	LEU	2.0
1	A	381	GLY	2.0
1	A	608	TYR	2.0
1	A	185	ILE	2.0
1	A	336	ALA	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](#)

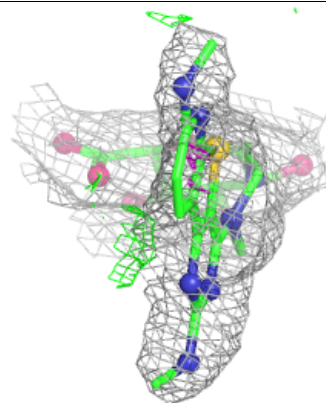
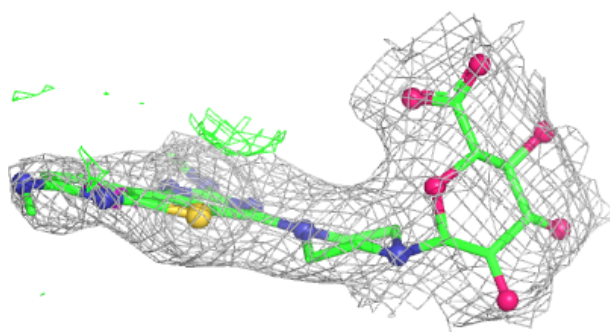
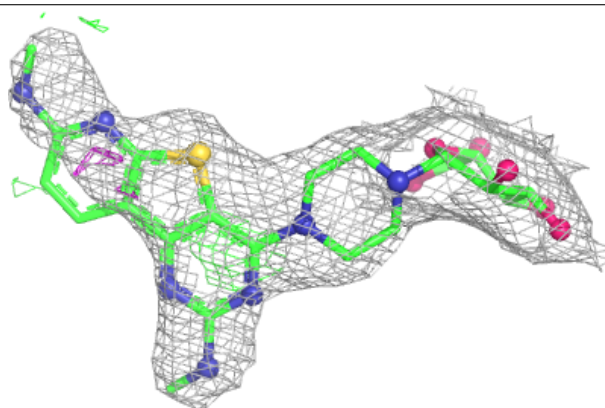
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	E0V	A	701	35/35	0.89	0.15	53,70,101,103	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 E0V A 701:

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.