



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 07:53 PM EDT

PDB ID : 3FO2
Title : Crystal structure of hapten complex of catalytic elimination antibody 13G5 (Glu(L39)Gln mutant)
Authors : Debler, E.W.; Wilson, I.A.
Deposited on : 2008-12-27
Resolution : 2.18 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

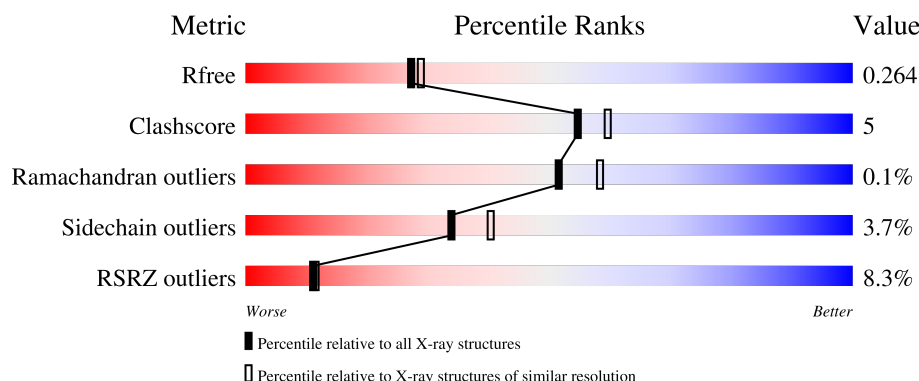
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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	219	<div> <div>9%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
1	L	219	<div> <div>10%</div> <div>89%</div> <div>9%</div> <div>.</div> </div>
2	B	229	<div> <div>7%</div> <div>77%</div> <div>15%</div> <div>7%</div> </div>
2	H	229	<div> <div>6%</div> <div>83%</div> <div>10%</div> <div>7%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6738 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 Catalytic antibody Fab 13G5 kappa light chain chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1668	1049	283	331	5			
1	A	216	Total	C	N	O	S	0	0	0
			1668	1049	283	331	5			

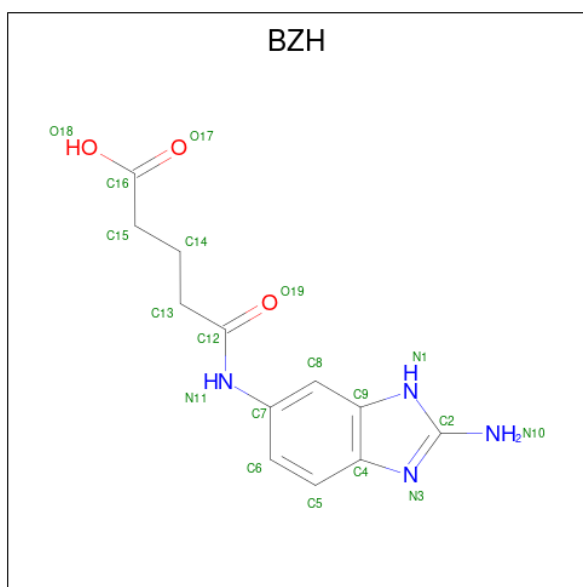
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	39	GLN	GLU	engineered mutation	PDB 3FO2
A	39	GLN	GLU	engineered mutation	PDB 3FO2

- Molecule 2 is a protein called Catalytic antibody Fab 13G5 IgG2b heavy chain chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	214	Total	C	N	O	S	0	0	0
			1601	1018	262	314	7			
2	B	214	Total	C	N	O	S	0	0	0
			1601	1018	262	314	7			

- Molecule 3 is 5-[(2-AMINO-1H-BENZIMIDAZOL-6-YL)AMINO]-5-OXOPENTANOIC ACID (three-letter code: BZH) (formula: C₁₂H₁₄N₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	H	1	Total	C	N	O	0	0
			19	12	4	3		
3	A	1	Total	C	N	O	0	0
			19	12	4	3		

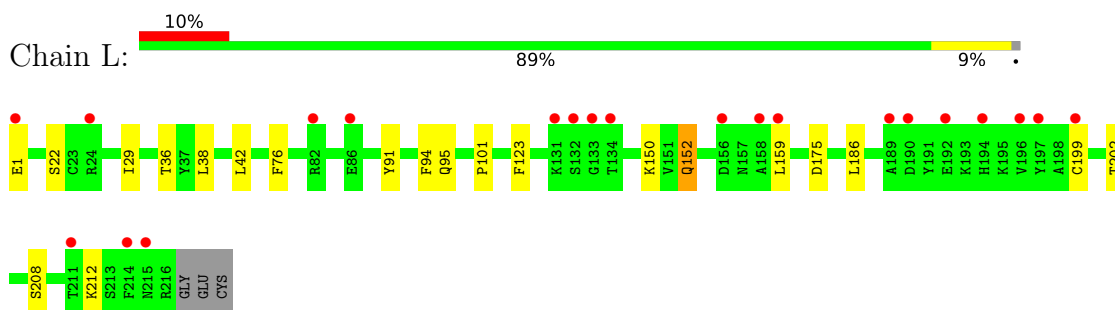
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	49	Total	O	0	0
			49	49		
4	H	40	Total	O	0	0
			40	40		
4	A	40	Total	O	0	0
			40	40		
4	B	33	Total	O	0	0
			33	33		

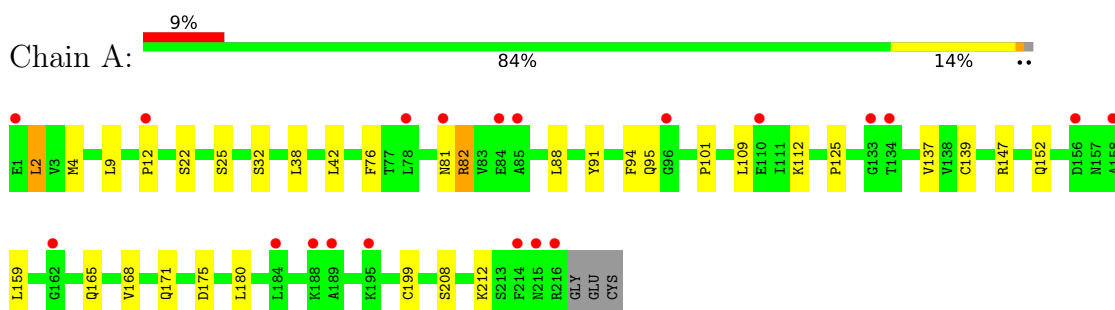
3 Residue-property plots [i](#)

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.

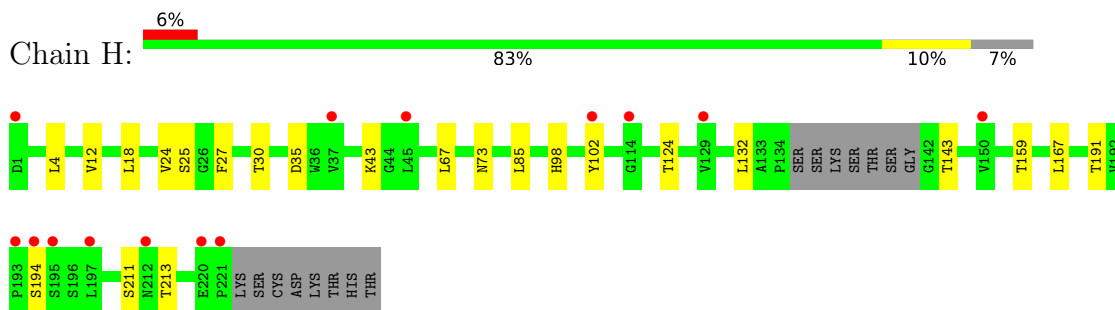
- Molecule 1: Catalytic antibody Fab 13G5 kappa light chain chimera



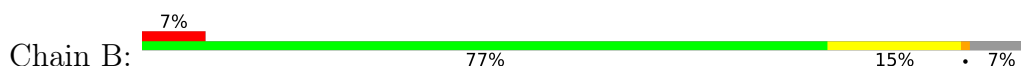
- Molecule 1: Catalytic antibody Fab 13G5 kappa light chain chimera

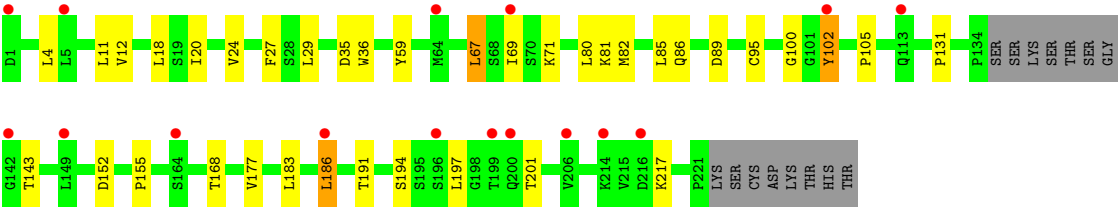


- Molecule 2: Catalytic antibody Fab 13G5 IgG2b heavy chain chimera



- Molecule 2: Catalytic antibody Fab 13G5 IgG2b heavy chain chimera





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.12Å 86.22Å 113.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.27 – 2.18 31.26 – 2.18	Depositor EDS
% Data completeness (in resolution range)	91.9 (31.27-2.18) 92.0 (31.26-2.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.256 0.212 , 0.264	Depositor DCC
R_{free} test set	1994 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6738	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 50.59 % 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 6.3334e-05. 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BZH

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	0.71	2/1705 (0.1%)	0.75	0/2313
1	L	0.74	1/1705 (0.1%)	0.75	0/2313
2	B	0.64	0/1644	0.74	0/2248
2	H	0.66	0/1644	0.75	0/2248
All	All	0.69	3/6698 (0.0%)	0.75	0/9122

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	199	CYS	CB-SG	-6.53	1.71	1.82
1	A	139	CYS	CB-SG	-5.45	1.73	1.81
1	L	199	CYS	CB-SG	-5.20	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1668	0	1637	17	0
1	L	1668	0	1637	12	0
2	B	1601	0	1562	27	0
2	H	1601	0	1562	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	19	0	13	0	0
3	H	19	0	13	0	0
4	A	40	0	0	1	0
4	B	33	0	0	0	0
4	H	40	0	0	1	0
4	L	49	0	0	1	0
All	All	6738	0	6424	67	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 5.

All (67) 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)
2:B:12:VAL:HG11	2:B:85:LEU:HD12	1.49	0.95
1:L:152:GLN:NE2	1:L:159:LEU:HD11	1.81	0.94
2:B:12:VAL:HG11	2:B:85:LEU:CD1	1.99	0.93
2:H:211:SER:OG	2:H:213:THR:HG23	1.77	0.84
2:H:18:LEU:HB2	2:H:85:LEU:HD11	1.59	0.83
2:H:12:VAL:HG11	2:H:85:LEU:HD12	1.72	0.72
1:A:165:GLN:OE1	2:B:177:VAL:HG11	1.93	0.68
1:A:2:LEU:HD21	1:A:25:SER:HB2	1.77	0.66
1:A:165:GLN:OE1	2:B:177:VAL:CG1	2.44	0.66
2:H:24:VAL:HG13	2:H:27:PHE:CE1	2.31	0.65
2:B:186:LEU:C	2:B:186:LEU:HD12	2.17	0.65
2:H:211:SER:OG	2:H:213:THR:CG2	2.46	0.63
2:B:18:LEU:HB2	2:B:85:LEU:HD11	1.80	0.63
2:B:12:VAL:CG1	2:B:85:LEU:HD12	2.25	0.63
2:H:159:THR:HG23	4:H:265:HOH:O	2.00	0.61
2:B:143:THR:HG23	2:B:191:THR:HG23	1.82	0.60
1:L:186:LEU:HD23	4:L:260:HOH:O	2.03	0.59
2:B:59:TYR:HD1	2:B:67:LEU:HD13	1.68	0.58
2:H:12:VAL:HG11	2:H:85:LEU:CD1	2.33	0.58
2:B:59:TYR:CD1	2:B:67:LEU:HD13	2.39	0.57
2:B:80:LEU:HD23	2:B:81:LYS:N	2.20	0.56
1:A:32:SER:HB2	4:A:242:HOH:O	2.06	0.55
1:A:88:LEU:HD21	1:A:171:GLN:HB3	1.88	0.55
2:H:18:LEU:CB	2:H:85:LEU:HD11	2.35	0.55
2:B:80:LEU:HD23	2:B:80:LEU:C	2.26	0.55
1:L:42:LEU:HD13	1:L:91:TYR:CZ	2.43	0.54
1:L:29:ILE:HG13	1:L:36:THR:HG23	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:38:LEU:HD22	1:L:76:PHE:CG	2.44	0.53
1:L:1:GLU:HB3	1:A:9:LEU:HD12	1.91	0.52
2:B:12:VAL:CG1	2:B:85:LEU:CD1	2.83	0.52
1:A:152:GLN:OE1	1:A:159:LEU:HD11	2.10	0.52
2:H:4:LEU:HG	2:H:24:VAL:HG22	1.93	0.50
1:L:152:GLN:HE22	1:L:159:LEU:HD11	1.71	0.50
1:A:180:LEU:C	1:A:180:LEU:HD23	2.31	0.50
2:B:36:TRP:HD1	2:B:69:ILE:HD13	1.77	0.49
1:A:42:LEU:HD13	1:A:91:TYR:CZ	2.48	0.48
2:B:100:GLY:O	2:B:105:PRO:HA	2.12	0.48
2:H:124:THR:HG21	2:H:211:SER:HA	1.95	0.48
1:L:29:ILE:HD11	1:L:76:PHE:CE1	2.49	0.48
1:A:125:PRO:HD3	1:A:137:VAL:HG22	1.96	0.47
1:L:150:LYS:HB3	1:L:202:THR:HB	1.97	0.47
2:B:82:MET:HE3	2:B:89:ASP:CG	2.36	0.47
1:A:81:ASN:O	1:A:82:ARG:HB2	2.15	0.46
1:A:38:LEU:HD22	1:A:76:PHE:CG	2.51	0.46
1:L:94:PHE:CZ	1:L:101:PRO:HB2	2.52	0.44
2:B:24:VAL:HG13	2:B:27:PHE:CE1	2.52	0.44
2:B:18:LEU:HD21	2:B:20:ILE:HD11	2.00	0.44
2:B:102:TYR:C	2:B:102:TYR:CD1	2.91	0.43
2:B:11:LEU:HB2	2:B:155:PRO:HG3	1.99	0.43
2:B:35:ASP:O	2:B:95:CYS:HA	2.18	0.43
1:A:12:PRO:HB2	1:A:112:LYS:HD2	2.00	0.43
2:B:82:MET:HE2	2:B:85:LEU:HD23	1.99	0.43
2:B:4:LEU:HG	2:B:24:VAL:HG22	2.01	0.42
2:B:131:PRO:HD3	2:B:217:LYS:HE3	2.02	0.42
1:L:29:ILE:HD11	1:L:76:PHE:CZ	2.55	0.42
2:H:30:THR:HG22	2:H:73:ASN:ND2	2.34	0.42
2:H:167:LEU:HD12	2:H:167:LEU:HA	1.92	0.41
2:B:152:ASP:HB3	2:B:183:LEU:HD13	2.03	0.41
2:B:186:LEU:C	2:B:186:LEU:CD1	2.86	0.41
1:A:147:ARG:HH11	1:A:168:VAL:HG11	1.85	0.41
2:H:24:VAL:CG1	2:H:27:PHE:CZ	3.04	0.41
1:L:123:PHE:CB	2:H:132:LEU:HD22	2.51	0.41
1:A:42:LEU:HD13	1:A:91:TYR:CE1	2.57	0.40
2:B:29:LEU:O	2:B:71:LYS:NZ	2.55	0.40
1:A:94:PHE:CZ	1:A:101:PRO:HB2	2.56	0.40
2:H:35:ASP:OD2	2:H:98:HIS:NE2	2.54	0.40
1:A:4:MET:HE1	1:A:38:LEU:CD1	2.51	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	214/219 (98%)	209 (98%)	4 (2%)	1 (0%)	29	28
1	L	214/219 (98%)	210 (98%)	4 (2%)	0	100	100
2	B	210/229 (92%)	205 (98%)	5 (2%)	0	100	100
2	H	210/229 (92%)	203 (97%)	7 (3%)	0	100	100
All	All	848/896 (95%)	827 (98%)	20 (2%)	1 (0%)	51	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	ARG

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	192/194 (99%)	185 (96%)	7 (4%)	35	42
1	L	192/194 (99%)	186 (97%)	6 (3%)	40	48
2	B	183/197 (93%)	175 (96%)	8 (4%)	28	33
2	H	183/197 (93%)	176 (96%)	7 (4%)	33	39
All	All	750/782 (96%)	722 (96%)	28 (4%)	34	40

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	22	SER
1	L	95	GLN
1	L	152	GLN
1	L	175	ASP
1	L	208	SER
1	L	212	LYS
2	H	25	SER
2	H	43	LYS
2	H	67	LEU
2	H	102	TYR
2	H	143	THR
2	H	191	THR
2	H	194	SER
1	A	2	LEU
1	A	22	SER
1	A	95	GLN
1	A	109	LEU
1	A	175	ASP
1	A	208	SER
1	A	212	LYS
2	B	67	LEU
2	B	86	GLN
2	B	102	TYR
2	B	168	THR
2	B	186	LEU
2	B	194	SER
2	B	197	LEU
2	B	201	THR

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

Mol	Chain	Res	Type
2	H	172	HIS
2	B	172	HIS

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are 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$
3	BZH	A	401	-	17,20,20	1.49	3 (17%)	16,27,27	1.03	1 (6%)
3	BZH	H	401	-	17,20,20	1.46	2 (11%)	16,27,27	0.83	0

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	BZH	A	401	-	-	0/8/10/10	0/2/2/2
3	BZH	H	401	-	-	0/8/10/10	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	BZH	C7-N11	-3.45	1.34	1.41
3	H	401	BZH	C7-N11	-3.32	1.34	1.41
3	A	401	BZH	C8-C9	-2.41	1.38	1.41
3	A	401	BZH	C4-N3	-2.12	1.32	1.38
3	H	401	BZH	C13-C12	2.08	1.55	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	BZH	C8-C7-N11	-2.43	115.04	123.13

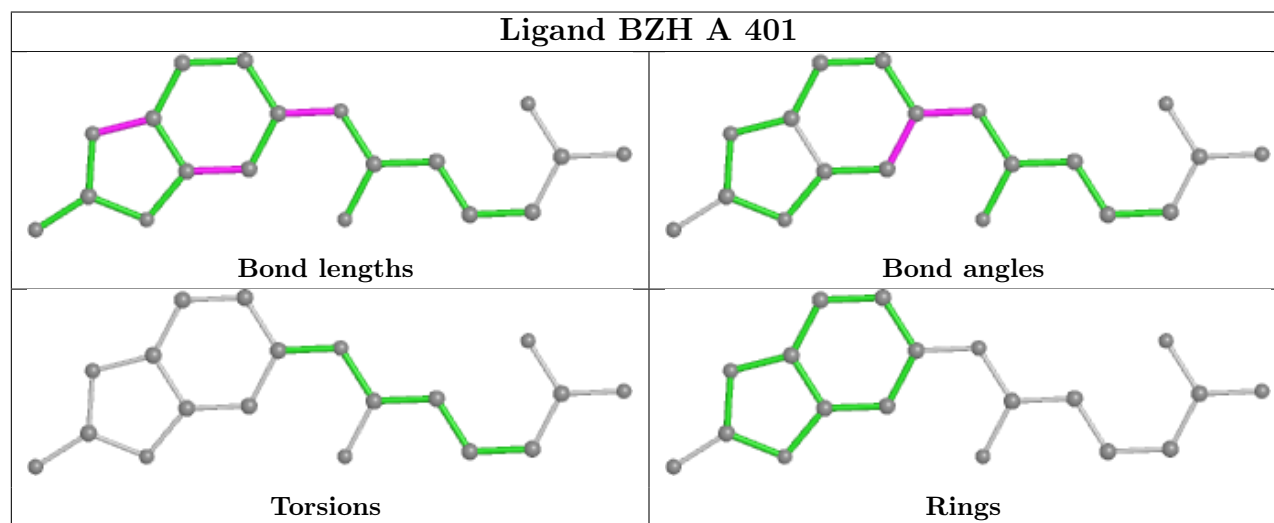
There are no chirality outliers.

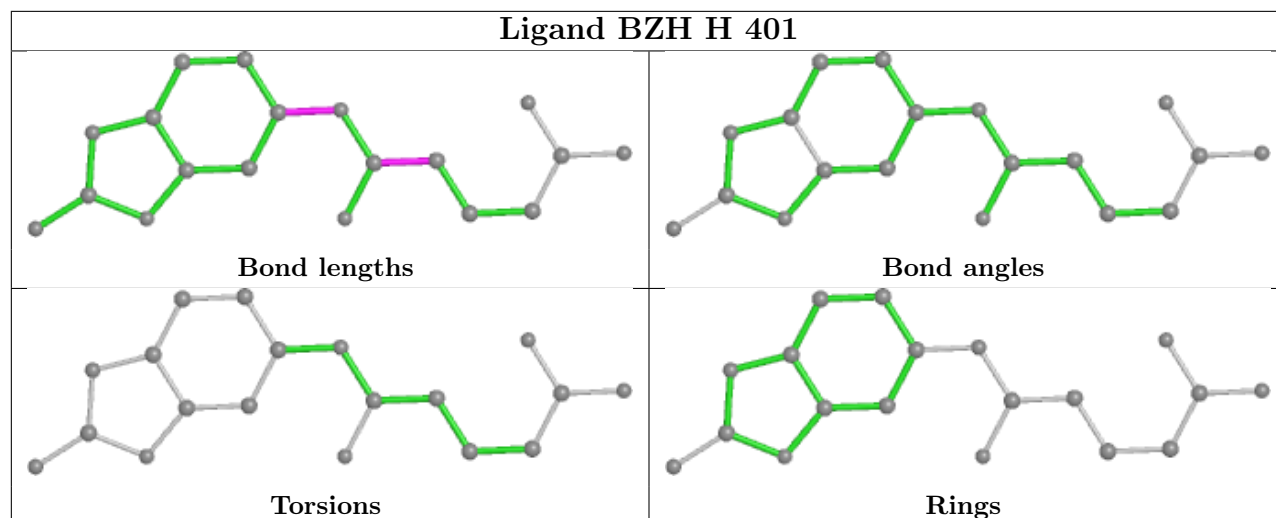
There are no torsion outliers.

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	216/219 (98%)	0.96	20 (9%) 8 9	44, 52, 59, 67	0
1	L	216/219 (98%)	0.88	21 (9%) 7 8	44, 52, 59, 70	0
2	B	214/229 (93%)	0.89	16 (7%) 14 15	45, 52, 59, 69	0
2	H	214/229 (93%)	0.89	14 (6%) 18 19	44, 52, 59, 69	0
All	All	860/896 (95%)	0.91	71 (8%) 11 11	44, 52, 59, 70	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	1	ASP	5.5
1	L	1	GLU	4.8
2	B	142	GLY	4.7
2	H	221	PRO	4.3
2	H	102	TYR	4.3
2	B	1	ASP	4.2
1	L	215	ASN	4.2
1	A	215	ASN	4.2
1	A	12	PRO	4.1
2	H	194	SER	4.0
2	B	199	THR	4.0
2	B	102	TYR	4.0
1	A	156	ASP	3.8
1	A	189	ALA	3.7
1	A	158	ALA	3.7
1	L	199	CYS	3.7
1	A	1	GLU	3.7
1	A	162	GLY	3.6
1	L	214	PHE	3.4
2	B	5	LEU	3.4
1	L	189	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
2	H	114	GLY	3.3
1	L	197	TYR	3.2
1	A	134	THR	3.2
1	L	132	SER	3.1
1	L	82	ARG	3.1
2	H	129	VAL	3.0
2	H	220	GLU	3.0
2	H	197	LEU	3.0
1	L	196	VAL	2.9
1	L	24	ARG	2.8
1	A	81	ASN	2.8
2	B	186	LEU	2.7
1	L	190	ASP	2.7
1	L	131	LYS	2.7
1	L	133	GLY	2.7
1	A	110	GLU	2.7
2	B	69	ILE	2.6
1	L	158	ALA	2.6
2	B	164	SER	2.5
1	L	134	THR	2.5
1	A	133	GLY	2.5
2	B	149	LEU	2.4
1	L	156	ASP	2.4
2	B	216	ASP	2.4
1	A	78	LEU	2.4
1	A	96	GLY	2.4
2	H	212	ASN	2.3
2	H	150	VAL	2.3
1	L	194	HIS	2.3
2	H	193	PRO	2.3
1	A	184	LEU	2.3
1	A	84	GLU	2.3
1	A	85	ALA	2.2
2	H	45	LEU	2.2
2	H	37	VAL	2.2
1	A	188	LYS	2.2
1	A	214	PHE	2.2
1	A	195	LYS	2.2
2	B	200	GLN	2.2
2	B	64	MET	2.2
2	B	206	VAL	2.2
1	L	192	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
2	H	195	SER	2.1
1	L	159	LEU	2.1
1	A	216	ARG	2.1
2	B	113	GLN	2.1
2	B	196	SER	2.1
1	L	211	THR	2.1
1	L	86	GLU	2.1
2	B	214	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 monosaccharides in this entry.

6.4 Ligands [i](#)

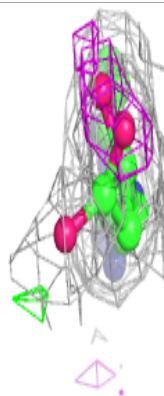
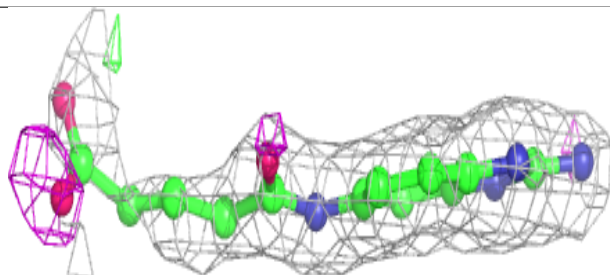
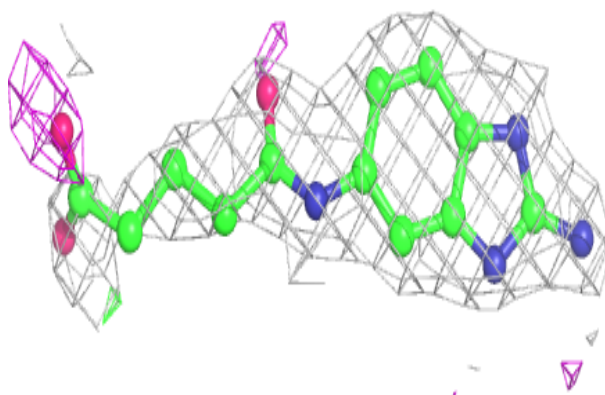
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	BZH	A	401	19/19	0.86	0.22	43,49,80,82	0
3	BZH	H	401	19/19	0.89	0.20	41,46,68,69	0

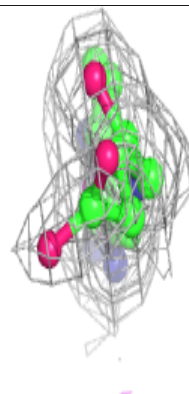
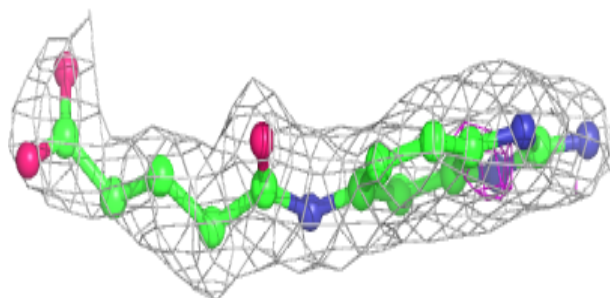
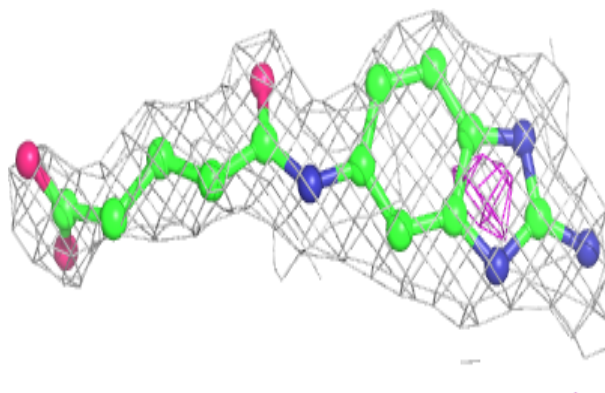
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BZH A 401:

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

**Electron density around BZH H 401:**

$2mF_o-DF_c$ (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.