Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 03:50 AM BST

PDB ID	:	6SC7
Title	:	dAb3/HOIP-RBR-Ligand3
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Deposited on	:	2019-07-23
Resolution	:	2.56 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:

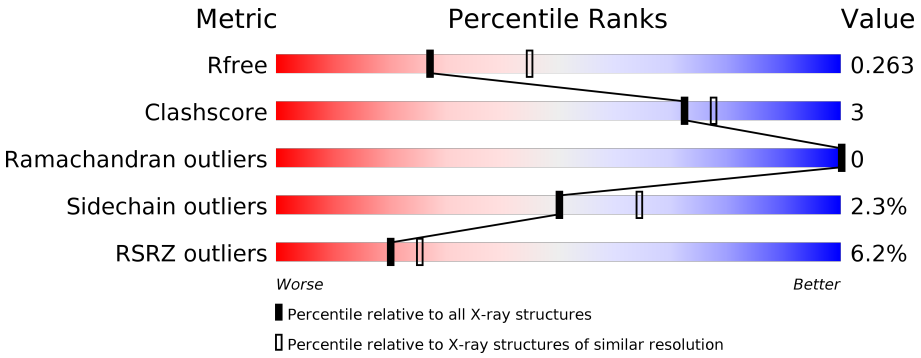
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 on 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 <=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	376	8% 85% 10% 5%
2	B	120	5% 86% 13%
2	C	120	92% 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4785 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.

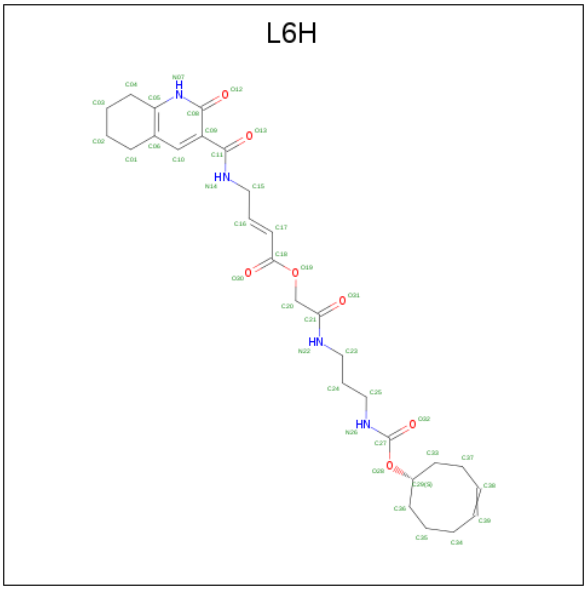
- Molecule 1 is a protein called E3 ubiquitin-protein ligase RNF31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total	C	N	O	S	0	0	0
			2800	1768	486	512	34			

- Molecule 2 is a protein called Single domain antibody.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	119	Total	C	N	O	S	0	0	0
			903	573	152	174	4			
2	C	120	Total	C	N	O	S	0	0	0
			917	581	155	177	4			

- Molecule 3 is [2-[3-(cyclooct-4-en-1-yloxycarbonylamino)propylamino]-2-oxidanylidene-ethyl] ({E})-4-[(2-oxidanylidene-5,6,7,8-tetrahydro-1 {H}-quinolin-3-yl)carbonylamino]but-2-enoate (three-letter code: L6H) (formula: $C_{28}H_{38}N_4O_7$) (labeled as "Ligand of Interest" by author).

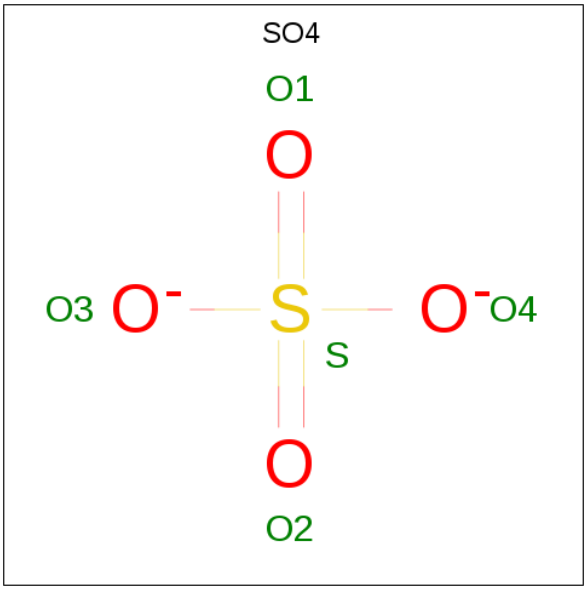


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			62	44	6	12		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	Zn	0	0
			8	8		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		

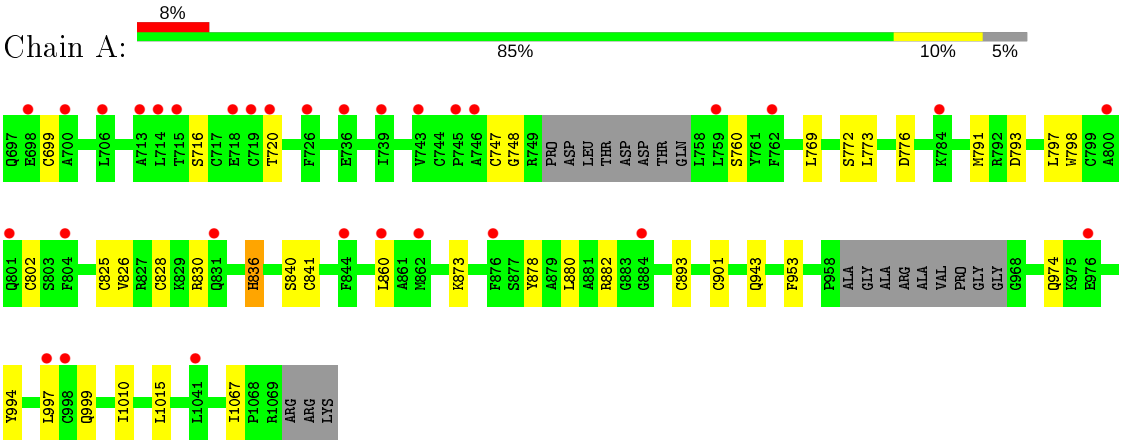
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	35	Total O 35 35	0	0
7	B	11	Total O 11 11	0	0
7	C	28	Total O 28 28	0	0

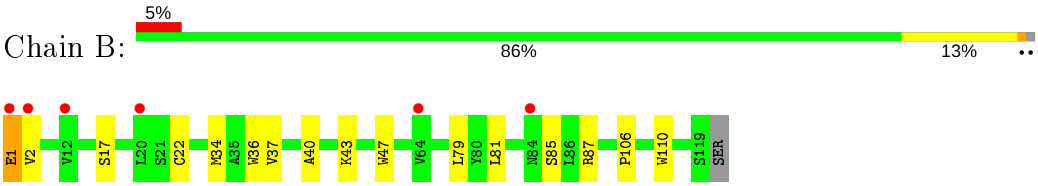
3 Residue-property plots (i)

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.

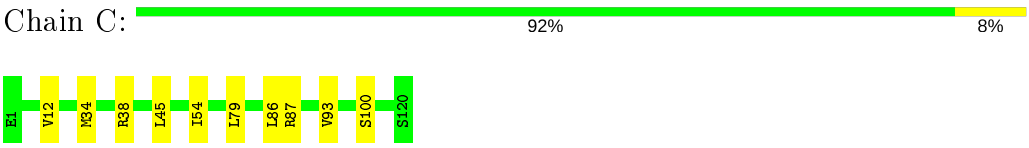
- Molecule 1: E3 ubiquitin-protein ligase RNF31



- Molecule 2: Single domain antibody



- Molecule 2: Single domain antibody



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	65.71Å 86.70Å 240.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.39 – 2.56 63.39 – 2.56	Depositor EDS
% Data completeness (in resolution range)	99.9 (63.39-2.56) 99.9 (63.39-2.56)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.49 (at 2.55Å)	Xtriage
Refinement program	REFMAC 1.15.2_3472, PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.224 , 0.267 0.234 , 0.263	Depositor DCC
R_{free} test set	1132 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	63.6	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 47.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4785	wwPDB-VP
Average B, all atoms (Å^2)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, CL, L6H, SO4

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	0.24	0/2874	0.40	0/3897
2	B	0.25	0/925	0.46	0/1255
2	C	0.26	0/939	0.45	0/1273
All	All	0.25	0/4738	0.42	0/6425

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

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	2800	0	2602	20	0
2	B	903	0	860	9	0
2	C	917	0	885	5	0
3	A	62	0	0	0	0
4	A	8	0	0	0	0
5	A	10	0	0	0	0
5	B	10	0	0	0	0
6	A	1	0	0	0	0
7	A	35	0	0	1	0
7	B	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	28	0	0	0	0
All	All	4785	0	4347	31	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 3.

All (31) 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:828:CYS:HB3	1:A:830:ARG:HG2	1.64	0.79
2:B:1:GLU:OE1	2:B:2:VAL:HG23	1.90	0.71
1:A:747:CYS:SG	1:A:748:GLY:N	2.73	0.61
1:A:1010:ILE:HG23	1:A:1015:LEU:HB2	1.84	0.59
1:A:825:CYS:HB3	1:A:828:CYS:HB2	1.86	0.57
2:B:85:SER:O	2:B:87:ARG:NH2	2.40	0.55
1:A:828:CYS:SG	1:A:841:CYS:HB3	2.48	0.54
1:A:836:HIS:O	1:A:836:HIS:ND1	2.43	0.51
1:A:791:MET:SD	1:A:791:MET:N	2.83	0.51
1:A:699:CYS:HA	1:A:720:THR:O	2.12	0.49
1:A:860:LEU:HB2	1:A:997:LEU:HD11	1.96	0.48
1:A:769:LEU:O	1:A:773:LEU:HB2	2.14	0.47
1:A:943:GLN:HB2	1:A:953:PHE:CZ	2.50	0.47
1:A:882:ARG:NH2	7:A:2104:HOH:O	2.49	0.46
1:A:901:CYS:HA	1:A:1067:ILE:HG12	1.98	0.46
2:B:40:ALA:HB3	2:B:43:LYS:HE2	1.98	0.46
2:C:34:MET:HB3	2:C:79:LEU:HD22	1.99	0.44
2:B:34:MET:HB3	2:B:79:LEU:HD22	1.99	0.43
2:B:110:TRP:HB2	2:C:45:LEU:HB2	2.00	0.43
1:A:994:TYR:CE2	1:A:999:GLN:HG3	2.54	0.43
1:A:873:LYS:HD3	1:A:893:CYS:SG	2.58	0.43
2:B:22:CYS:HB3	2:B:79:LEU:HB3	2.00	0.42
2:C:38:ARG:HA	2:C:93:VAL:O	2.20	0.41
2:B:37:VAL:HG12	2:B:47:TRP:HA	2.02	0.41
1:A:793:ASP:OD2	2:C:54:ILE:HG12	2.20	0.41
1:A:880:LEU:HD21	1:A:974:GLN:HG3	2.02	0.41
1:A:878:TYR:HB3	1:A:880:LEU:HD12	2.02	0.41
1:A:797:LEU:HD21	1:A:826:VAL:HG12	2.03	0.41
2:B:36:TRP:NE1	2:B:81:LEU:HB2	2.35	0.40
2:C:12:VAL:HG11	2:C:86:LEU:HD12	2.03	0.40
1:A:798:TRP:CD2	2:B:106:PRO:HD3	2.57	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	350/376 (93%)	338 (97%)	12 (3%)	0	100	100
2	B	117/120 (98%)	117 (100%)	0	0	100	100
2	C	118/120 (98%)	117 (99%)	1 (1%)	0	100	100
All	All	585/616 (95%)	572 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	293/326 (90%)	286 (98%)	7 (2%)	49	63
2	B	92/97 (95%)	90 (98%)	2 (2%)	52	66
2	C	96/97 (99%)	94 (98%)	2 (2%)	53	67
All	All	481/520 (92%)	470 (98%)	11 (2%)	50	64

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	716	SER
1	A	760	SER

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Mol	Chain	Res	Type
1	A	772	SER
1	A	776	ASP
1	A	802	CYS
1	A	836	HIS
1	A	840	SER
2	B	1	GLU
2	B	17	SER
2	C	87	ARG
2	C	100	SER

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

Mol	Chain	Res	Type
1	A	851	ASN
1	A	865	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 15 ligands modelled in this entry, 9 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	L6H	A	2001[A]	-	41,41,41	2.75	14 (34%)	42,52,52	2.21	12 (28%)
5	SO4	B	201	-	4,4,4	0.14	0	6,6,6	0.06	0
5	SO4	B	202	-	4,4,4	0.14	0	6,6,6	0.05	0
5	SO4	A	2011	-	4,4,4	0.14	0	6,6,6	0.05	0
5	SO4	A	2010	-	4,4,4	0.15	0	6,6,6	0.05	0
3	L6H	A	2001[B]	-	41,41,41	2.74	14 (34%)	42,52,52	2.19	14 (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
3	L6H	A	2001[A]	-	-	14/28/45/45	1/3/3/3
3	L6H	A	2001[B]	-	-	10/28/45/45	1/3/3/3

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2001[A]	L6H	C17-C16	8.25	1.53	1.32
3	A	2001[B]	L6H	C17-C16	8.21	1.53	1.32
3	A	2001[A]	L6H	O12-C08	6.28	1.40	1.24
3	A	2001[B]	L6H	O12-C08	6.28	1.40	1.24
3	A	2001[A]	L6H	C21-N22	5.59	1.46	1.33
3	A	2001[A]	L6H	C11-N14	5.59	1.46	1.33
3	A	2001[B]	L6H	C11-N14	5.59	1.46	1.33
3	A	2001[B]	L6H	C21-N22	5.55	1.46	1.33
3	A	2001[A]	L6H	C27-N26	5.23	1.45	1.34
3	A	2001[A]	L6H	C39-C38	5.21	1.53	1.32
3	A	2001[B]	L6H	C39-C38	5.21	1.53	1.32
3	A	2001[B]	L6H	C27-N26	5.19	1.45	1.34
3	A	2001[A]	L6H	C04-C05	3.55	1.56	1.50
3	A	2001[B]	L6H	C04-C05	3.55	1.56	1.50
3	A	2001[B]	L6H	O28-C27	3.04	1.40	1.35
3	A	2001[A]	L6H	O28-C27	2.99	1.40	1.35
3	A	2001[A]	L6H	O19-C18	2.74	1.41	1.34
3	A	2001[B]	L6H	O19-C18	2.69	1.40	1.34
3	A	2001[A]	L6H	C06-C05	-2.60	1.37	1.41
3	A	2001[B]	L6H	C06-C05	-2.60	1.37	1.41
3	A	2001[A]	L6H	C17-C18	2.46	1.54	1.48
3	A	2001[B]	L6H	C17-C18	2.35	1.53	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2001[A]	L6H	O13-C11	-2.15	1.18	1.23
3	A	2001[B]	L6H	O13-C11	-2.15	1.18	1.23
3	A	2001[A]	L6H	O31-C21	-2.13	1.18	1.23
3	A	2001[B]	L6H	O31-C21	-2.12	1.19	1.23
3	A	2001[A]	L6H	C10-C06	2.05	1.43	1.39
3	A	2001[B]	L6H	C10-C06	2.05	1.43	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	2001[A]	L6H	C09-C08-N07	-6.85	119.63	124.40
3	A	2001[B]	L6H	C09-C08-N07	-6.85	119.63	124.40
3	A	2001[B]	L6H	O28-C27-N26	5.09	120.05	111.11
3	A	2001[A]	L6H	O28-C27-N26	4.89	119.69	111.11
3	A	2001[A]	L6H	C16-C17-C18	-4.37	111.79	122.92
3	A	2001[A]	L6H	C10-C09-C08	3.66	119.88	117.58
3	A	2001[B]	L6H	C10-C09-C08	3.66	119.88	117.58
3	A	2001[B]	L6H	C16-C17-C18	-3.50	113.98	122.92
3	A	2001[B]	L6H	O32-C27-N26	-3.31	119.88	124.96
3	A	2001[A]	L6H	O19-C18-C17	3.29	120.30	111.55
3	A	2001[A]	L6H	O32-C27-N26	-3.14	120.14	124.96
3	A	2001[B]	L6H	O19-C18-C17	2.96	119.43	111.55
3	A	2001[B]	L6H	O28-C27-O32	-2.96	120.07	124.53
3	A	2001[A]	L6H	C08-N07-C05	2.96	122.57	116.16
3	A	2001[B]	L6H	C08-N07-C05	2.96	122.57	116.16
3	A	2001[A]	L6H	O28-C27-O32	-2.90	120.16	124.53
3	A	2001[A]	L6H	C09-C11-N14	2.84	120.13	116.30
3	A	2001[B]	L6H	C09-C11-N14	2.84	120.13	116.30
3	A	2001[B]	L6H	C34-C39-C38	-2.75	113.65	126.85
3	A	2001[A]	L6H	C34-C39-C38	-2.72	113.81	126.85
3	A	2001[A]	L6H	C06-C05-N07	-2.72	120.81	122.81
3	A	2001[B]	L6H	C06-C05-N07	-2.72	120.81	122.81
3	A	2001[A]	L6H	C37-C38-C39	-2.57	114.52	126.85
3	A	2001[B]	L6H	C37-C38-C39	-2.54	114.68	126.85
3	A	2001[B]	L6H	C29-O28-C27	-2.20	113.33	116.48
3	A	2001[B]	L6H	C20-C21-N22	2.03	119.94	116.58

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2001[A]	L6H	C15-C16-C17-C18

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Mol	Chain	Res	Type	Atoms
3	A	2001[A]	L6H	C20-C21-N22-C23
3	A	2001[A]	L6H	O28-C27-N26-C25
3	A	2001[A]	L6H	O32-C27-N26-C25
3	A	2001[B]	L6H	C17-C18-O19-C20
3	A	2001[B]	L6H	O30-C18-O19-C20
3	A	2001[B]	L6H	O28-C27-N26-C25
3	A	2001[B]	L6H	N26-C27-O28-C29
3	A	2001[B]	L6H	O32-C27-O28-C29
3	A	2001[A]	L6H	C17-C18-O19-C20
3	A	2001[B]	L6H	O32-C27-N26-C25
3	A	2001[A]	L6H	O31-C21-N22-C23
3	A	2001[A]	L6H	C16-C17-C18-O19
3	A	2001[A]	L6H	O30-C18-O19-C20
3	A	2001[A]	L6H	N22-C23-C24-C25
3	A	2001[A]	L6H	C16-C17-C18-O30
3	A	2001[B]	L6H	O19-C20-C21-N22
3	A	2001[A]	L6H	O19-C20-C21-O31
3	A	2001[B]	L6H	O19-C20-C21-O31
3	A	2001[B]	L6H	C16-C17-C18-O30
3	A	2001[B]	L6H	C16-C17-C18-O19
3	A	2001[A]	L6H	O19-C20-C21-N22
3	A	2001[A]	L6H	C21-C20-O19-C18
3	A	2001[A]	L6H	N14-C15-C16-C17

All (2) ring outliers are listed below:

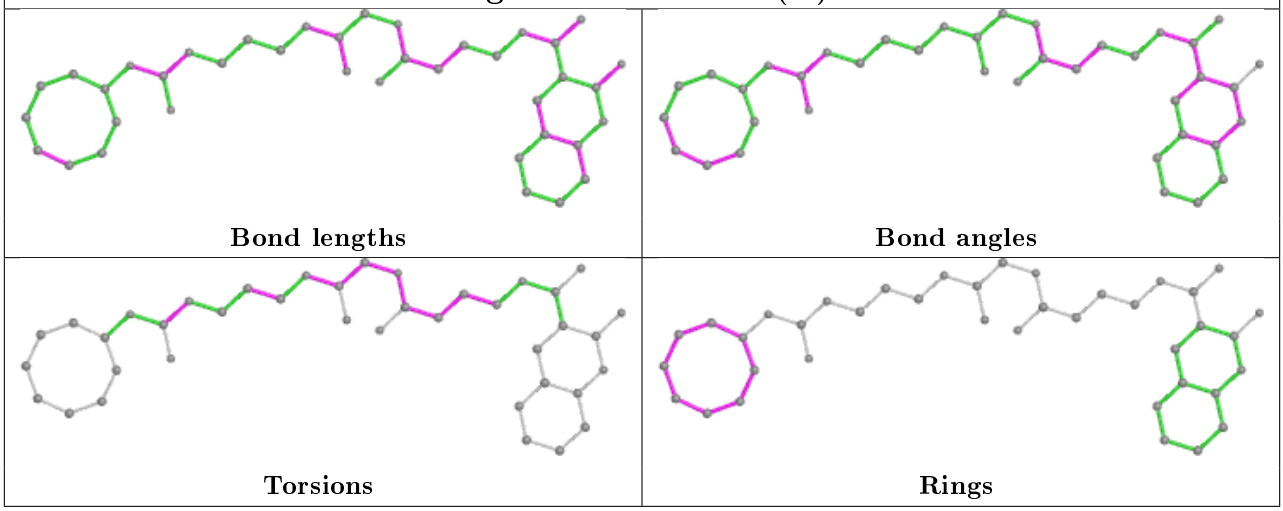
Mol	Chain	Res	Type	Atoms
3	A	2001[B]	L6H	C29-C33-C34-C35-C36-C37-C38-C39
3	A	2001[A]	L6H	C29-C33-C34-C35-C36-C37-C38-C39

No monomer is involved in short contacts.

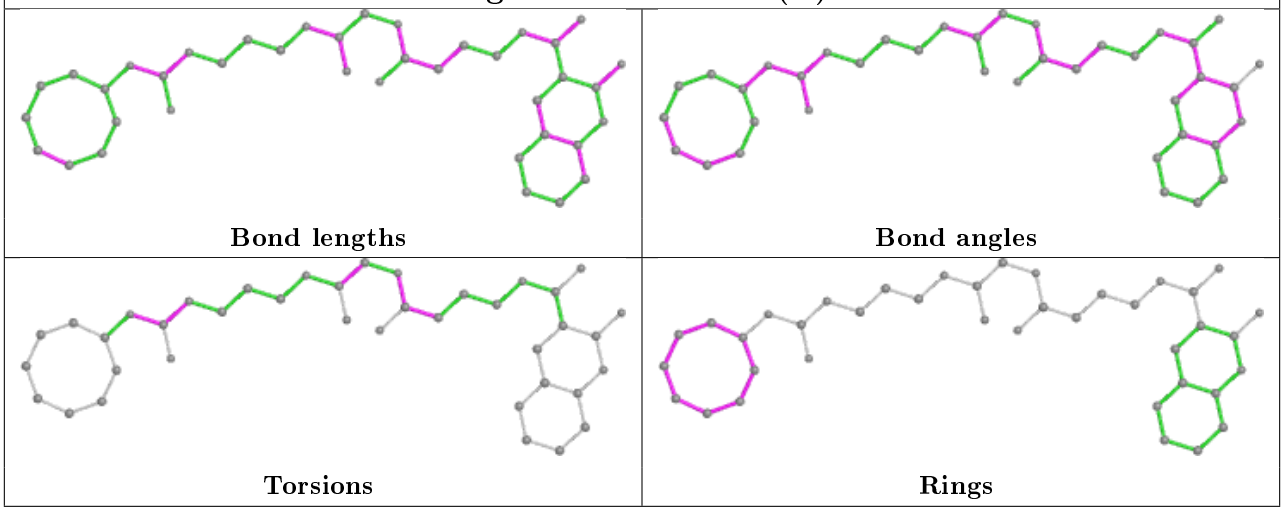
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

Ligand L6H A 2001 (A)



Ligand L6H A 2001 (B)



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, 95^{th} 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	356/376 (94%)	0.68	31 (8%) 10 14	46, 70, 99, 109	0
2	B	119/120 (99%)	0.47	6 (5%) 28 36	48, 62, 79, 91	0
2	C	120/120 (100%)	0.27	0 100 100	44, 52, 67, 77	0
All	All	595/616 (96%)	0.55	37 (6%) 20 25	44, 63, 93, 109	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	715	THR	5.2
1	A	739	ILE	4.9
1	A	759	LEU	4.7
1	A	801	GLN	4.1
1	A	762	PHE	4.0
1	A	746	ALA	4.0
1	A	1041	LEU	3.8
1	A	706	LEU	3.5
1	A	714	LEU	3.4
2	B	20	LEU	3.1
1	A	713	ALA	3.1
1	A	745	PRO	2.9
1	A	726	PHE	2.8
2	B	84	ASN	2.7
1	A	844	PHE	2.6
1	A	800	ALA	2.5
2	B	2	VAL	2.5
1	A	743	VAL	2.5
1	A	736	GLU	2.5
1	A	784	LYS	2.5
2	B	1	GLU	2.4
1	A	718	GLU	2.4
1	A	860	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	12	VAL	2.4
1	A	698	GLU	2.4
1	A	884	GLY	2.2
1	A	720	THR	2.2
1	A	976	GLU	2.2
1	A	998	CYS	2.2
1	A	719	CYS	2.2
1	A	804	PHE	2.1
1	A	862	MET	2.1
1	A	831	GLN	2.1
1	A	876	PHE	2.1
2	B	64	VAL	2.1
1	A	700	ALA	2.0
1	A	997	LEU	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, 95^{th} 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	L6H	A	2001[A]	39/39	0.80	0.33	66,75,80,80	23
3	L6H	A	2001[B]	39/39	0.80	0.33	66,72,79,81	23
5	SO4	B	202	5/5	0.92	0.18	76,87,89,104	0
5	SO4	A	2011	5/5	0.93	0.13	70,73,81,91	0
4	ZN	A	2003	1/1	0.94	0.13	105,105,105,105	0
5	SO4	B	201	5/5	0.95	0.15	63,67,71,72	0
4	ZN	A	2002	1/1	0.98	0.12	70,70,70,70	0
4	ZN	A	2009	1/1	0.98	0.24	76,76,76,76	0
5	SO4	A	2010	5/5	0.98	0.14	48,50,53,53	0

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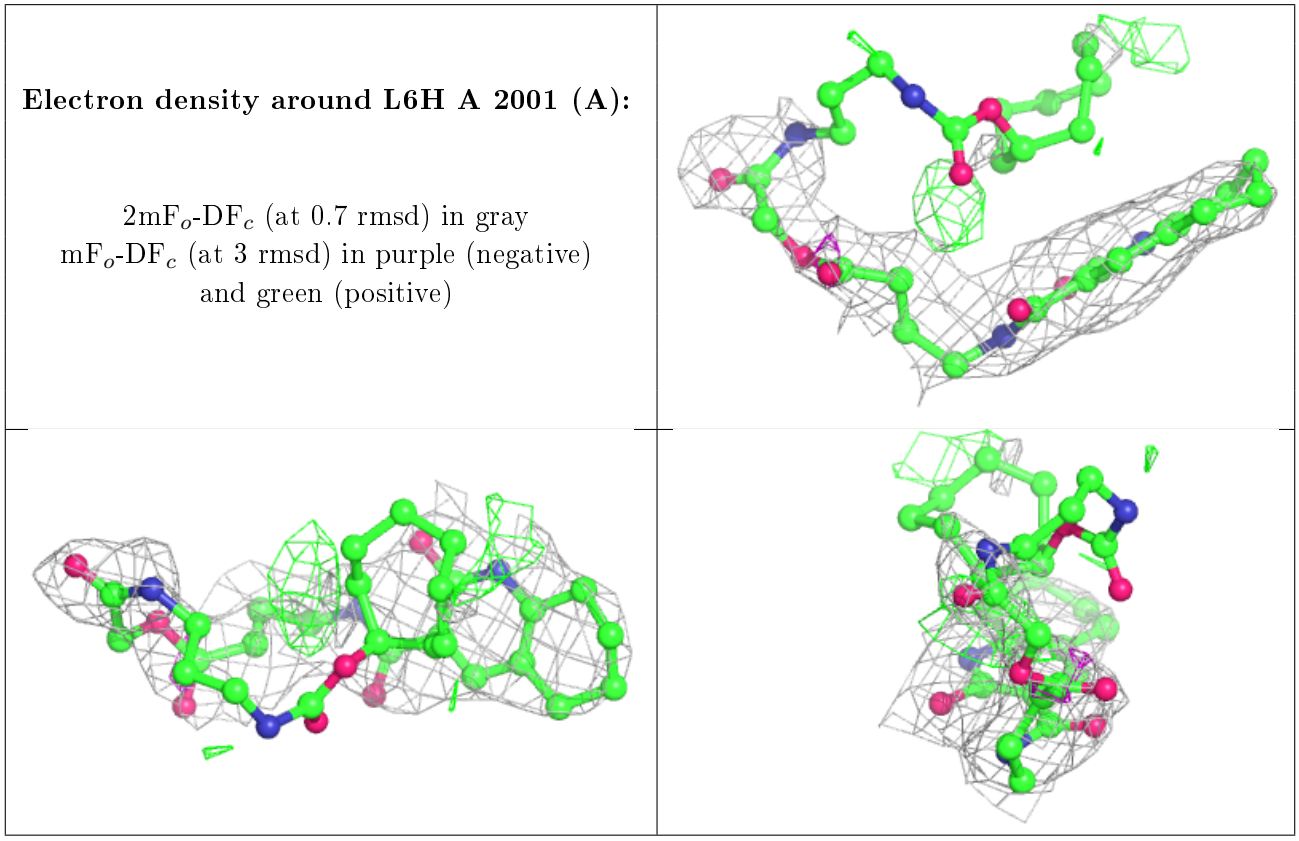
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	CL	A	2012	1/1	0.98	0.06	57,57,57,57	0
4	ZN	A	2006	1/1	0.98	0.19	61,61,61,61	0
4	ZN	A	2005	1/1	0.99	0.16	61,61,61,61	0
4	ZN	A	2007	1/1	0.99	0.19	46,46,46,46	0
4	ZN	A	2004	1/1	0.99	0.14	61,61,61,61	0
4	ZN	A	2008	1/1	0.99	0.15	55,55,55,55	0

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.

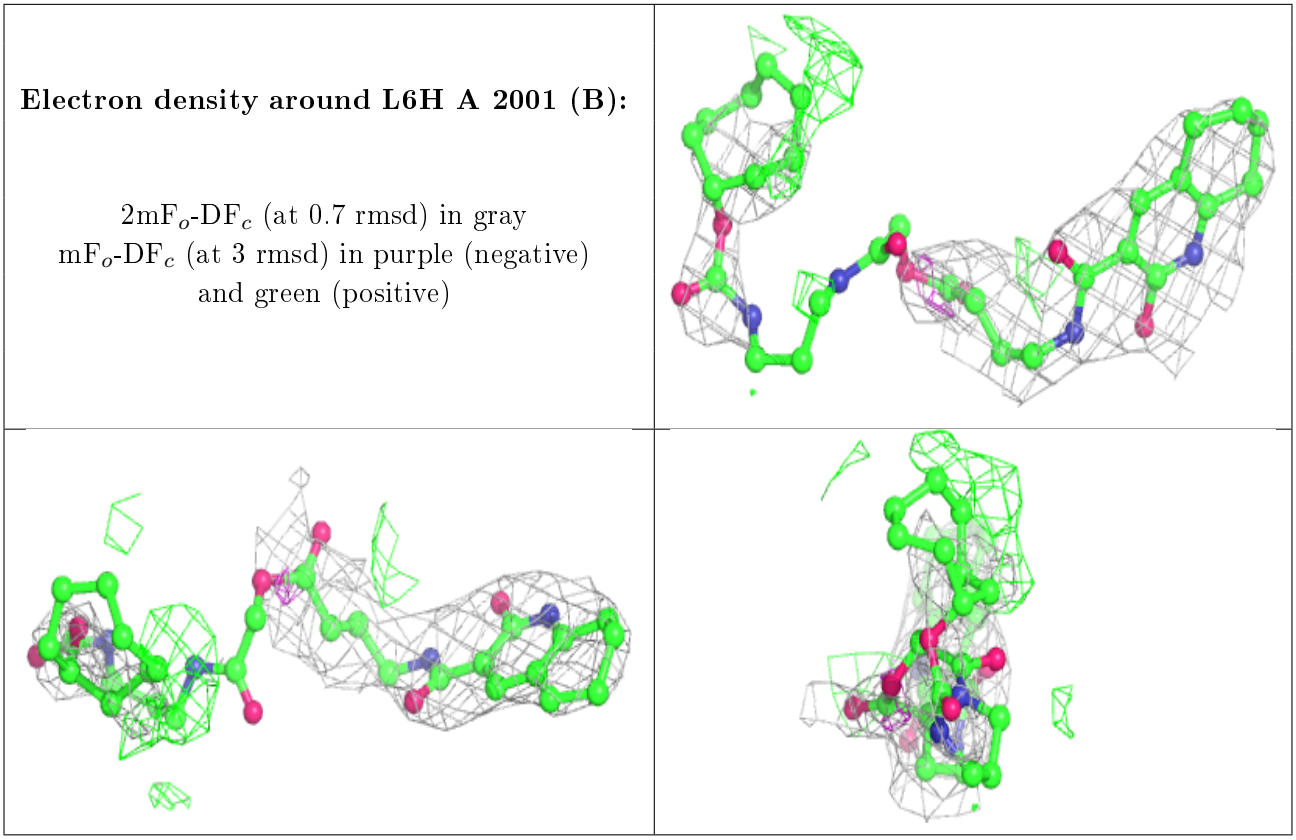
Electron density around L6H A 2001 (A):

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



Electron density around L6H A 2001 (B):

$2mF_o$-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.