



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:02 PM UTC

PDB ID : 9EO0 / pdb_00009eo0
Title : Small-Molecule Inhibitors of Programmed Cell Death-1/Programmed Death-Ligand 1
Authors : Plewka, J.; Hec, A.; Sitar, T.; Holak, T.
Deposited on : 2024-03-14
Resolution : 2.50 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

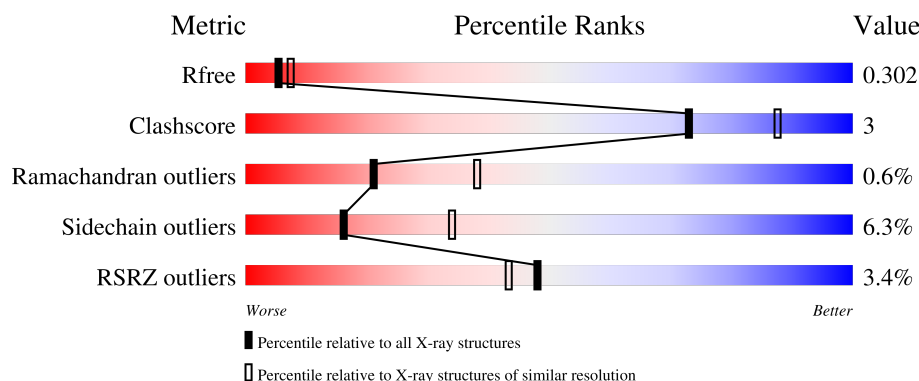
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	131	4% 76% 16% 8%
1	B	131	3% 76% 13% 10%
1	C	131	5% 80% 9% 9%
1	D	131	4% 76% 12% 9%
1	E	131	0% 77% 9% 11%

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Mol	Chain	Length	Quality of chain
1	F	131	2% 73% 13% 5% 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11430 atoms, of which 5597 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 Programmed cell death 1 ligand 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	120	Total	C	H	N	O	S	0	0	0
			1864	605	916	158	180	5			
1	B	118	Total	C	H	N	O	S	0	0	0
			1859	602	918	157	177	5			
1	C	119	Total	C	H	N	O	S	0	0	0
			1869	605	922	158	179	5			
1	D	119	Total	C	H	N	O	S	0	0	0
			1859	603	917	157	177	5			
1	E	117	Total	C	H	N	O	S	0	0	0
			1834	596	902	155	176	5			
1	F	118	Total	C	H	N	O	S	0	0	0
			1874	605	929	158	177	5			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q9NZQ7
A	2	GLY	-	expression tag	UNP Q9NZQ7
A	3	SER	-	expression tag	UNP Q9NZQ7
A	121	ALA	-	expression tag	UNP Q9NZQ7
A	122	ALA	-	expression tag	UNP Q9NZQ7
A	123	ALA	-	expression tag	UNP Q9NZQ7
A	124	LEU	-	expression tag	UNP Q9NZQ7
A	125	GLU	-	expression tag	UNP Q9NZQ7
A	126	HIS	-	expression tag	UNP Q9NZQ7
A	127	HIS	-	expression tag	UNP Q9NZQ7
A	128	HIS	-	expression tag	UNP Q9NZQ7
A	129	HIS	-	expression tag	UNP Q9NZQ7
A	130	HIS	-	expression tag	UNP Q9NZQ7
A	131	HIS	-	expression tag	UNP Q9NZQ7
B	1	MET	-	initiating methionine	UNP Q9NZQ7
B	2	GLY	-	expression tag	UNP Q9NZQ7
B	3	SER	-	expression tag	UNP Q9NZQ7

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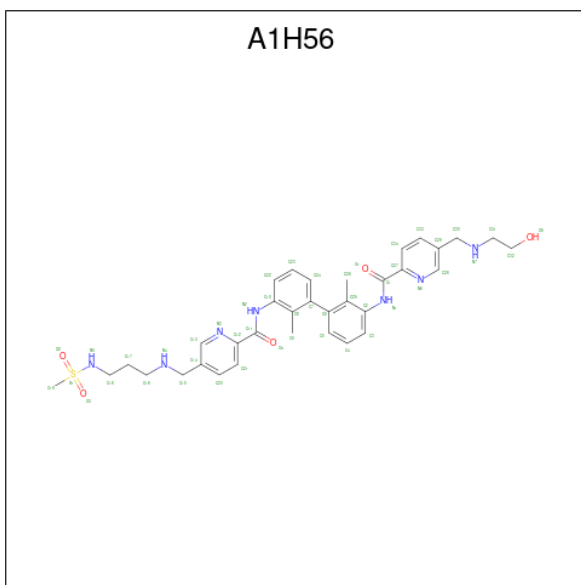
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B	121	ALA	-	expression tag	UNP Q9NZQ7
B	122	ALA	-	expression tag	UNP Q9NZQ7
B	123	ALA	-	expression tag	UNP Q9NZQ7
B	124	LEU	-	expression tag	UNP Q9NZQ7
B	125	GLU	-	expression tag	UNP Q9NZQ7
B	126	HIS	-	expression tag	UNP Q9NZQ7
B	127	HIS	-	expression tag	UNP Q9NZQ7
B	128	HIS	-	expression tag	UNP Q9NZQ7
B	129	HIS	-	expression tag	UNP Q9NZQ7
B	130	HIS	-	expression tag	UNP Q9NZQ7
B	131	HIS	-	expression tag	UNP Q9NZQ7
C	1	MET	-	initiating methionine	UNP Q9NZQ7
C	2	GLY	-	expression tag	UNP Q9NZQ7
C	3	SER	-	expression tag	UNP Q9NZQ7
C	121	ALA	-	expression tag	UNP Q9NZQ7
C	122	ALA	-	expression tag	UNP Q9NZQ7
C	123	ALA	-	expression tag	UNP Q9NZQ7
C	124	LEU	-	expression tag	UNP Q9NZQ7
C	125	GLU	-	expression tag	UNP Q9NZQ7
C	126	HIS	-	expression tag	UNP Q9NZQ7
C	127	HIS	-	expression tag	UNP Q9NZQ7
C	128	HIS	-	expression tag	UNP Q9NZQ7
C	129	HIS	-	expression tag	UNP Q9NZQ7
C	130	HIS	-	expression tag	UNP Q9NZQ7
C	131	HIS	-	expression tag	UNP Q9NZQ7
D	1	MET	-	initiating methionine	UNP Q9NZQ7
D	2	GLY	-	expression tag	UNP Q9NZQ7
D	3	SER	-	expression tag	UNP Q9NZQ7
D	121	ALA	-	expression tag	UNP Q9NZQ7
D	122	ALA	-	expression tag	UNP Q9NZQ7
D	123	ALA	-	expression tag	UNP Q9NZQ7
D	124	LEU	-	expression tag	UNP Q9NZQ7
D	125	GLU	-	expression tag	UNP Q9NZQ7
D	126	HIS	-	expression tag	UNP Q9NZQ7
D	127	HIS	-	expression tag	UNP Q9NZQ7
D	128	HIS	-	expression tag	UNP Q9NZQ7
D	129	HIS	-	expression tag	UNP Q9NZQ7
D	130	HIS	-	expression tag	UNP Q9NZQ7
D	131	HIS	-	expression tag	UNP Q9NZQ7
E	1	MET	-	initiating methionine	UNP Q9NZQ7
E	2	GLY	-	expression tag	UNP Q9NZQ7
E	3	SER	-	expression tag	UNP Q9NZQ7

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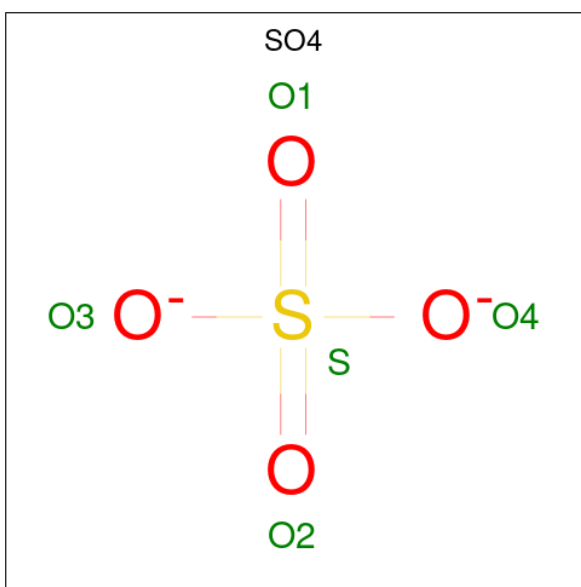
Chain	Residue	Modelled	Actual	Comment	Reference
E	121	ALA	-	expression tag	UNP Q9NZQ7
E	122	ALA	-	expression tag	UNP Q9NZQ7
E	123	ALA	-	expression tag	UNP Q9NZQ7
E	124	LEU	-	expression tag	UNP Q9NZQ7
E	125	GLU	-	expression tag	UNP Q9NZQ7
E	126	HIS	-	expression tag	UNP Q9NZQ7
E	127	HIS	-	expression tag	UNP Q9NZQ7
E	128	HIS	-	expression tag	UNP Q9NZQ7
E	129	HIS	-	expression tag	UNP Q9NZQ7
E	130	HIS	-	expression tag	UNP Q9NZQ7
E	131	HIS	-	expression tag	UNP Q9NZQ7
F	1	MET	-	initiating methionine	UNP Q9NZQ7
F	2	GLY	-	expression tag	UNP Q9NZQ7
F	3	SER	-	expression tag	UNP Q9NZQ7
F	121	ALA	-	expression tag	UNP Q9NZQ7
F	122	ALA	-	expression tag	UNP Q9NZQ7
F	123	ALA	-	expression tag	UNP Q9NZQ7
F	124	LEU	-	expression tag	UNP Q9NZQ7
F	125	GLU	-	expression tag	UNP Q9NZQ7
F	126	HIS	-	expression tag	UNP Q9NZQ7
F	127	HIS	-	expression tag	UNP Q9NZQ7
F	128	HIS	-	expression tag	UNP Q9NZQ7
F	129	HIS	-	expression tag	UNP Q9NZQ7
F	130	HIS	-	expression tag	UNP Q9NZQ7
F	131	HIS	-	expression tag	UNP Q9NZQ7

- Molecule 2 is {N}-[3-[3-[[5-[(2-hydroxyethylamino)methyl]pyridin-2-yl]carbonylamino]-2-methyl-phenyl]-2-methyl-phenyl]-5-[[3-(methylsulfonylamino)propylamino]methyl]pyridine-2-carboxamide (CCD ID: A1H56) (formula: C₃₄H₄₁N₇O₅S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	H	N	O	S	0	0
			78	34	31	7	5	1		
2	C	1	Total	C	H	N	O	S	0	0
			78	34	31	7	5	1		
2	D	1	Total	C	H	N	O	S	0	0
			78	34	31	7	5	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	O S	0	0
			5	4 1		

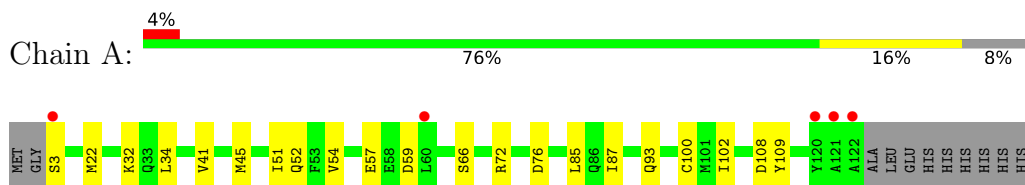
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total 11	O 11	0	0
4	B	3	Total 3	O 3	0	0
4	C	3	Total 3	O 3	0	0
4	D	6	Total 6	O 6	0	0
4	E	3	Total 3	O 3	0	0
4	F	6	Total 6	O 6	0	0

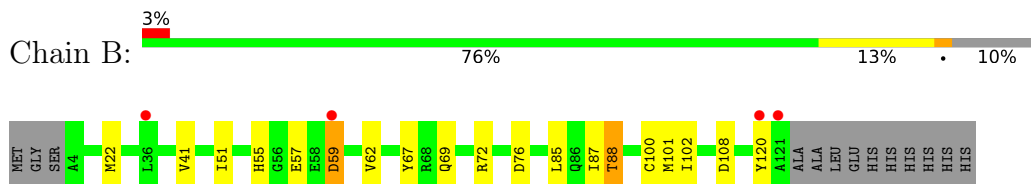
3 Residue-property plots [i](#)

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.

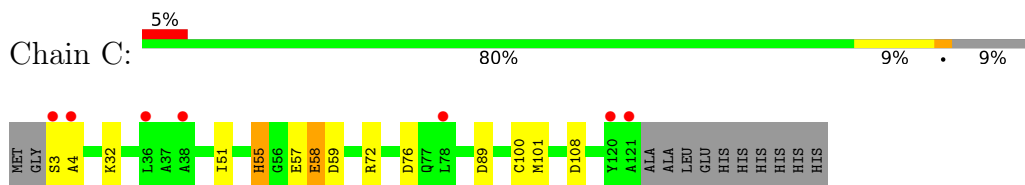
- Molecule 1: Programmed cell death 1 ligand 1



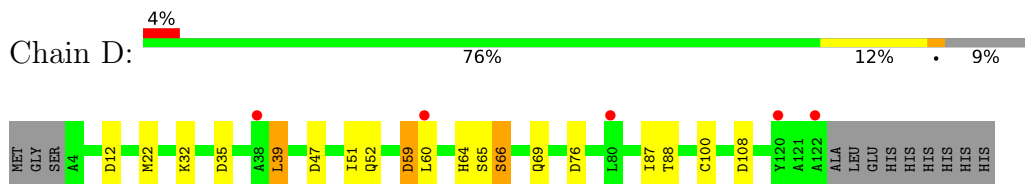
- Molecule 1: Programmed cell death 1 ligand 1



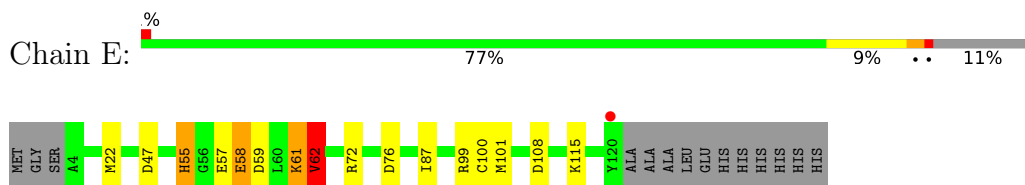
- Molecule 1: Programmed cell death 1 ligand 1



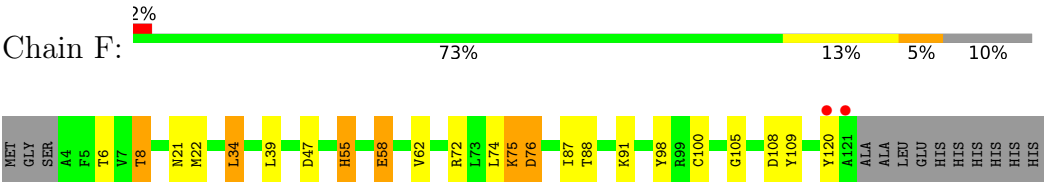
- Molecule 1: Programmed cell death 1 ligand 1



- Molecule 1: Programmed cell death 1 ligand 1



- Molecule 1: Programmed cell death 1 ligand 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.75Å 75.41Å 75.92Å 90.00° 97.25° 90.00°	Depositor
Resolution (Å)	44.39 – 2.50 44.39 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (44.39-2.50) 99.1 (44.39-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0425, REFMAC 5.8.0425	Depositor
R, R_{free}	0.253 , 0.302 0.253 , 0.302	Depositor DCC
R_{free} test set	1380 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11430	wwPDB-VP
Average B, all atoms (Å ²)	65.0	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 49.41 % 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 7.4937e-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: SO4, A1H56

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	0/965	1.14	3/1308 (0.2%)
1	B	0.60	0/958	1.18	4/1297 (0.3%)
1	C	0.61	0/964	1.13	5/1305 (0.4%)
1	D	0.61	0/959	1.15	7/1299 (0.5%)
1	E	0.61	0/949	1.21	6/1286 (0.5%)
1	F	0.61	0/962	1.28	6/1301 (0.5%)
All	All	0.61	0/5757	1.18	31/7796 (0.4%)

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
1	B	0	1
1	D	0	1
All	All	0	3

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	55	HIS	CA-CB-CG	13.63	127.44	113.80
1	E	55	HIS	CA-CB-CG	13.32	127.12	113.80
1	F	8	THR	CA-CB-OG1	-12.65	90.63	109.60
1	F	76	ASP	CA-CB-CG	8.47	121.07	112.60
1	B	55	HIS	CB-CA-C	8.30	124.22	112.11
1	C	55	HIS	CB-CA-C	7.22	122.66	112.11
1	D	59	ASP	CA-CB-CG	7.06	119.66	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	62	VAL	N-CA-CB	6.72	122.32	111.23
1	D	108	ASP	CA-CB-CG	6.19	118.79	112.60
1	F	58	GLU	N-CA-CB	6.03	119.46	110.29
1	C	108	ASP	CA-CB-CG	5.99	118.59	112.60
1	D	76	ASP	CA-CB-CG	5.84	118.44	112.60
1	C	58	GLU	N-CA-CB	5.79	119.09	110.29
1	C	76	ASP	CA-CB-CG	5.74	118.34	112.60
1	E	76	ASP	CA-CB-CG	5.74	118.34	112.60
1	E	108	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	108	ASP	CA-CB-CG	5.64	118.24	112.60
1	F	108	ASP	CA-CB-CG	5.63	118.23	112.60
1	B	76	ASP	CA-CB-CG	5.63	118.23	112.60
1	E	58	GLU	N-CA-CB	5.58	118.77	110.29
1	D	47	ASP	CA-CB-CG	5.58	118.18	112.60
1	B	108	ASP	CA-CB-CG	5.51	118.11	112.60
1	E	61	LYS	N-CA-C	-5.37	102.45	110.23
1	A	76	ASP	CA-CB-CG	5.26	117.86	112.60
1	F	91	LYS	CB-CG-CD	5.23	123.33	111.30
1	C	89	ASP	CA-CB-CG	5.20	117.80	112.60
1	D	35	ASP	CA-CB-CG	5.18	117.78	112.60
1	D	12	ASP	CA-CB-CG	5.17	117.77	112.60
1	B	22	MET	CG-SD-CE	5.15	112.23	100.90
1	D	32	LYS	CB-CA-C	5.12	120.61	110.42
1	A	32	LYS	CB-CA-C	5.08	120.54	110.42

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	ARG	Sidechain
1	B	72	ARG	Sidechain
1	D	66	SER	Peptide

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	948	916	936	6	0
1	B	941	918	937	9	0
1	C	947	922	942	6	0
1	D	942	917	936	6	0
1	E	932	902	921	6	0
1	F	945	929	948	10	0
2	B	47	31	0	0	0
2	C	47	31	0	1	0
2	D	47	31	0	0	0
3	C	5	0	0	0	0
4	A	11	0	0	0	0
4	B	3	0	0	1	0
4	C	3	0	0	0	0
4	D	6	0	0	2	0
4	E	3	0	0	0	0
4	F	6	0	0	1	0
All	All	5833	5597	5620	37	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 (37) 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:F:98:TYR:OH	4:F:201:HOH:O	2.04	0.75
1:B:67:TYR:OH	4:B:301:HOH:O	2.03	0.73
1:B:59:ASP:OD2	1:B:62:VAL:HG22	1.92	0.70
1:B:88:THR:HG21	1:C:72:ARG:HH21	1.66	0.60
1:D:64:HIS:CD2	1:D:65:SER:O	2.57	0.58
1:F:21:ASN:HD22	1:F:88:THR:HA	1.68	0.57
1:D:60:LEU:O	1:D:60:LEU:HD23	2.04	0.56
1:B:41:VAL:HG22	1:B:102:ILE:HD12	1.87	0.56
1:E:61:LYS:O	1:E:62:VAL:HG12	2.05	0.56
1:A:41:VAL:HG22	1:A:102:ILE:HD12	1.89	0.55
1:B:69:GLN:O	1:C:72:ARG:CZ	2.55	0.55
1:A:22:MET:HE2	1:A:87:ILE:HB	1.91	0.53
2:C:201:A1H56:N5	1:F:6:THR:HG22	2.24	0.53
1:C:3:SER:HA	1:F:62:VAL:HG11	1.92	0.51
1:E:99:ARG:CD	1:E:101:MET:HE3	2.40	0.51
1:A:109:TYR:CE1	1:B:101:MET:HE1	2.47	0.50
1:D:52:GLN:N	4:D:301:HOH:O	2.45	0.50
1:C:101:MET:HE1	1:F:109:TYR:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:99:ARG:HD2	1:E:101:MET:HE3	1.95	0.47
1:A:52:GLN:HG2	1:A:54:VAL:HG23	1.97	0.46
1:F:22:MET:HE2	1:F:87:ILE:HB	1.96	0.46
1:B:100:CYS:SG	1:B:100:CYS:O	2.73	0.46
1:E:22:MET:HE2	1:E:87:ILE:HB	1.97	0.46
1:B:85:LEU:HG	1:B:87:ILE:HD12	1.99	0.45
1:A:100:CYS:O	1:A:100:CYS:SG	2.75	0.44
1:D:100:CYS:SG	1:D:100:CYS:O	2.74	0.44
1:C:100:CYS:O	1:C:100:CYS:SG	2.76	0.44
1:E:100:CYS:O	1:E:100:CYS:SG	2.76	0.44
1:A:85:LEU:HG	1:A:87:ILE:HD12	1.99	0.43
1:D:22:MET:HE2	1:D:87:ILE:HB	2.00	0.43
1:F:100:CYS:O	1:F:100:CYS:SG	2.76	0.43
1:D:39:LEU:N	4:D:302:HOH:O	2.53	0.42
1:F:34:LEU:HD11	1:F:39:LEU:HD22	2.01	0.41
1:B:85:LEU:HD21	1:B:87:ILE:HD11	2.02	0.41
1:C:55:HIS:NE2	1:F:105:GLY:O	2.53	0.41
1:E:99:ARG:HD3	1:E:101:MET:HE3	2.02	0.41
1:F:74:LEU:O	1:F:76:ASP:N	2.54	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	118/131 (90%)	112 (95%)	6 (5%)	0	100	100
1	B	116/131 (88%)	110 (95%)	6 (5%)	0	100	100
1	C	117/131 (89%)	110 (94%)	6 (5%)	1 (1%)	14	27
1	D	117/131 (89%)	112 (96%)	5 (4%)	0	100	100
1	E	115/131 (88%)	111 (96%)	3 (3%)	1 (1%)	14	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	116/131 (88%)	108 (93%)	6 (5%)	2 (2%)	7	13
All	All	699/786 (89%)	663 (95%)	32 (5%)	4 (1%)	21	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	62	VAL
1	F	75	LYS
1	F	120	TYR
1	C	4	ALA

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	101/112 (90%)	93 (92%)	8 (8%)	11	24
1	B	101/112 (90%)	96 (95%)	5 (5%)	22	44
1	C	102/112 (91%)	97 (95%)	5 (5%)	22	45
1	D	100/112 (89%)	94 (94%)	6 (6%)	17	36
1	E	100/112 (89%)	93 (93%)	7 (7%)	14	29
1	F	102/112 (91%)	95 (93%)	7 (7%)	14	30
All	All	606/672 (90%)	568 (94%)	38 (6%)	16	34

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	34	LEU
1	A	45	MET
1	A	51	ILE
1	A	57	GLU
1	A	59	ASP
1	A	66	SER

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Mol	Chain	Res	Type
1	A	93	GLN
1	B	51	ILE
1	B	57	GLU
1	B	59	ASP
1	B	88	THR
1	B	120	TYR
1	C	32	LYS
1	C	51	ILE
1	C	57	GLU
1	C	58	GLU
1	C	59	ASP
1	D	39	LEU
1	D	51	ILE
1	D	59	ASP
1	D	66	SER
1	D	69	GLN
1	D	88	THR
1	E	47	ASP
1	E	55	HIS
1	E	57	GLU
1	E	58	GLU
1	E	59	ASP
1	E	72	ARG
1	E	115	LYS
1	F	8	THR
1	F	34	LEU
1	F	47	ASP
1	F	55	HIS
1	F	58	GLU
1	F	72	ARG
1	F	75	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	86	GLN
1	A	117	ASN
1	B	64	HIS
1	B	69	GLN
1	C	52	GLN
1	C	63	GLN

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Mol	Chain	Res	Type
1	C	64	HIS
1	D	52	GLN
1	D	63	GLN
1	D	64	HIS
1	D	69	GLN
1	E	64	HIS
1	E	86	GLN
1	F	21	ASN
1	F	63	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
2	A1H56	C	201	-	50,50,50	0.78	2 (4%)	67,68,68	0.92	2 (2%)
2	A1H56	B	201	-	50,50,50	0.77	2 (4%)	67,68,68	0.92	5 (7%)
3	SO4	C	202	-	4,4,4	0.41	0	6,6,6	0.08	0
2	A1H56	D	201	-	50,50,50	0.79	2 (4%)	67,68,68	1.03	3 (4%)

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	A1H56	C	201	-	-	10/35/35/35	0/4/4/4
2	A1H56	B	201	-	-	15/35/35/35	0/4/4/4
2	A1H56	D	201	-	-	13/35/35/35	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	201	A1H56	C7-C8	3.12	1.45	1.40
2	B	201	A1H56	C7-C8	3.11	1.45	1.40
2	B	201	A1H56	C19-S1	2.97	1.82	1.75
2	C	201	A1H56	C6-C25	2.95	1.45	1.40
2	C	201	A1H56	C19-S1	2.95	1.82	1.75
2	D	201	A1H56	C19-S1	2.53	1.81	1.75

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	201	A1H56	C6-C7-C8	3.51	125.04	121.44
2	C	201	A1H56	C7-C6-C25	3.22	124.74	121.44
2	D	201	A1H56	C10-N2-C11	3.11	135.54	126.90
2	C	201	A1H56	C6-C7-C8	3.03	124.55	121.44
2	D	201	A1H56	C7-C6-C25	2.91	124.43	121.44
2	B	201	A1H56	C7-C6-C25	2.83	124.35	121.44
2	B	201	A1H56	C6-C7-C8	2.55	124.06	121.44
2	B	201	A1H56	C10-N2-C11	2.48	133.79	126.90
2	B	201	A1H56	C2-N1-C1	2.14	132.86	126.90
2	B	201	A1H56	O5-C32-C31	2.04	119.01	111.50

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	201	A1H56	C16-C17-C18-N5
2	B	201	A1H56	C18-N5-S1-C19
2	B	201	A1H56	C18-N5-S1-O2
2	B	201	A1H56	C18-N5-S1-O3
2	C	201	A1H56	C16-C17-C18-N5

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Mol	Chain	Res	Type	Atoms
2	C	201	A1H56	C18-N5-S1-C19
2	C	201	A1H56	N7-C31-C32-O5
2	D	201	A1H56	C16-C17-C18-N5
2	D	201	A1H56	C18-N5-S1-C19
2	D	201	A1H56	C18-N5-S1-O2
2	D	201	A1H56	N2-C11-C12-C21
2	D	201	A1H56	O4-C11-C12-C21
2	C	201	A1H56	N1-C1-C27-C34
2	D	201	A1H56	O4-C11-C12-N3
2	C	201	A1H56	O1-C1-C27-C34
2	C	201	A1H56	N1-C1-C27-N6
2	D	201	A1H56	N2-C11-C12-N3
2	C	201	A1H56	O1-C1-C27-N6
2	B	201	A1H56	N1-C1-C27-C34
2	B	201	A1H56	O1-C1-C27-C34
2	B	201	A1H56	N1-C1-C27-N6
2	B	201	A1H56	N2-C11-C12-N3
2	B	201	A1H56	N2-C11-C12-C21
2	B	201	A1H56	O4-C11-C12-C21
2	B	201	A1H56	O1-C1-C27-N6
2	B	201	A1H56	O4-C11-C12-N3
2	B	201	A1H56	C14-C15-N4-C16
2	D	201	A1H56	C32-C31-N7-C30
2	C	201	A1H56	C14-C15-N4-C16
2	C	201	A1H56	C29-C30-N7-C31
2	D	201	A1H56	N7-C31-C32-O5
2	B	201	A1H56	C29-C30-N7-C31
2	D	201	A1H56	C18-N5-S1-O3
2	D	201	A1H56	O1-C1-C27-N6
2	B	201	A1H56	C32-C31-N7-C30
2	D	201	A1H56	C17-C16-N4-C15
2	D	201	A1H56	O1-C1-C27-C34
2	C	201	A1H56	C32-C31-N7-C30

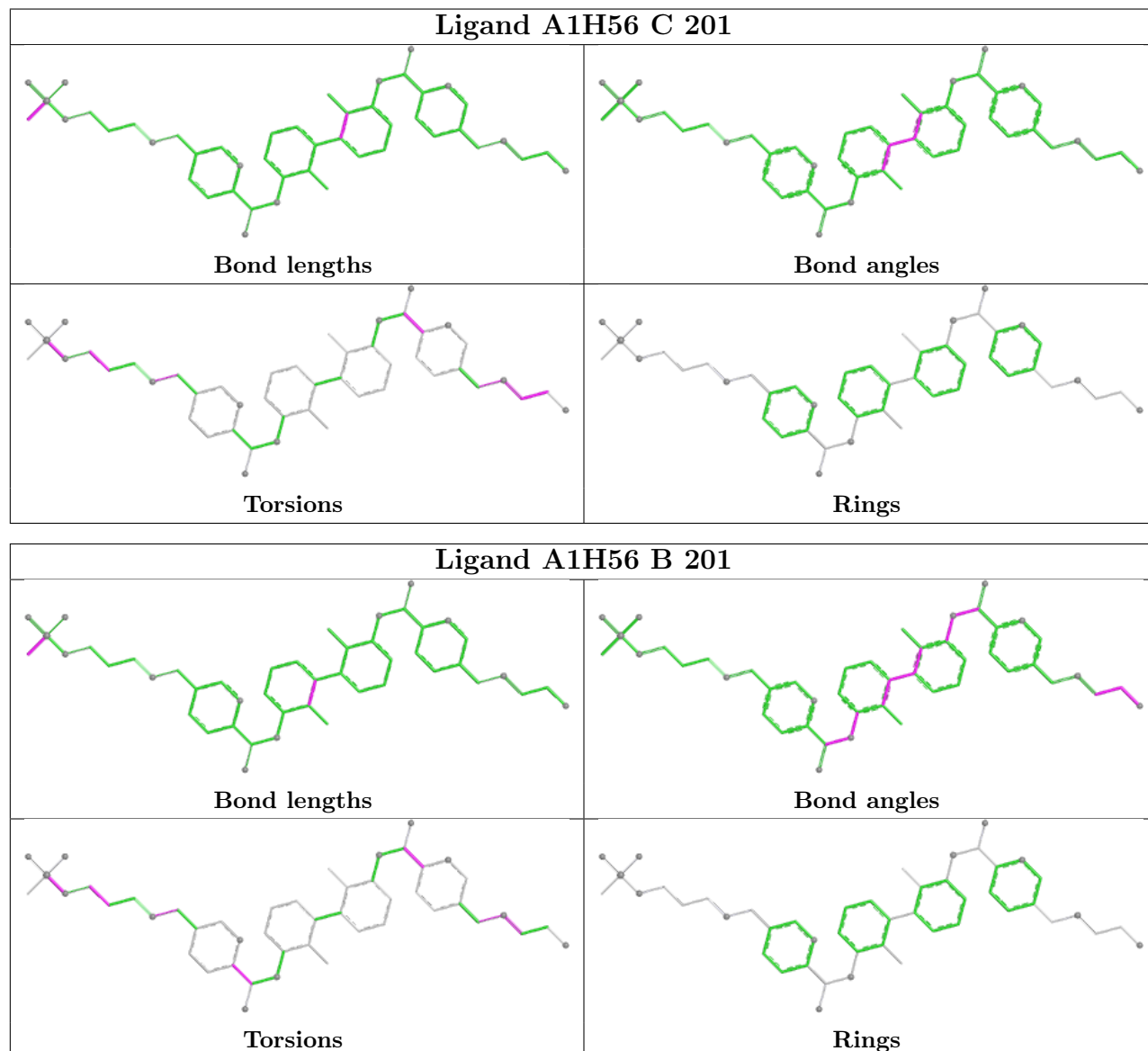
There are no ring outliers.

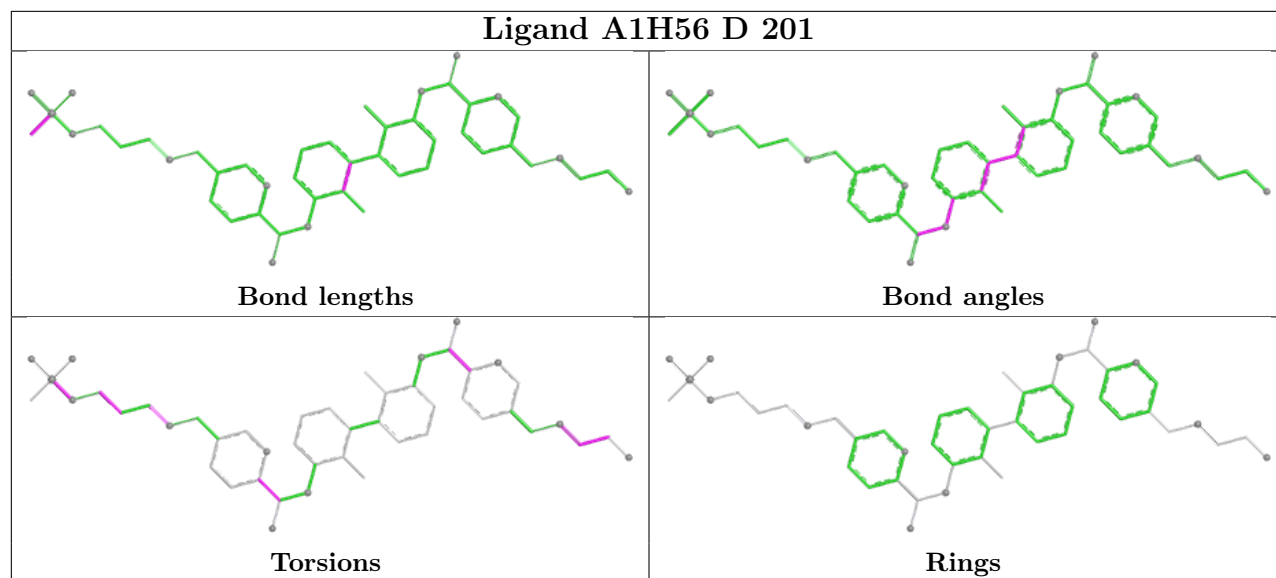
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	201	A1H56	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	120/131 (91%)	0.16	5 (4%)	40	36	32, 51, 97, 122	0
1	B	118/131 (90%)	0.33	4 (3%)	48	43	30, 56, 101, 144	0
1	C	119/131 (90%)	0.54	7 (5%)	28	24	36, 67, 117, 132	0
1	D	119/131 (90%)	0.45	5 (4%)	40	36	41, 65, 110, 152	0
1	E	117/131 (89%)	0.23	1 (0%)	81	78	43, 65, 92, 114	0
1	F	118/131 (90%)	0.19	2 (1%)	69	65	36, 57, 99, 144	0
All	All	711/786 (90%)	0.32	24 (3%)	48	43	30, 60, 105, 152	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	121	ALA	5.0
1	E	120	TYR	3.9
1	A	121	ALA	3.7
1	B	121	ALA	3.4
1	C	121	ALA	3.4
1	F	120	TYR	3.3
1	D	122	ALA	3.1
1	C	120	TYR	2.8
1	A	122	ALA	2.7
1	B	59	ASP	2.7
1	B	120	TYR	2.7
1	D	38	ALA	2.6
1	D	60	LEU	2.5
1	C	4	ALA	2.5
1	C	78	LEU	2.5
1	A	60	LEU	2.3
1	C	3	SER	2.3
1	A	120	TYR	2.2
1	B	36	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	36	LEU	2.1
1	C	38	ALA	2.1
1	A	3	SER	2.0
1	D	80	LEU	2.0
1	D	120	TYR	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](#)

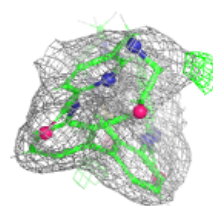
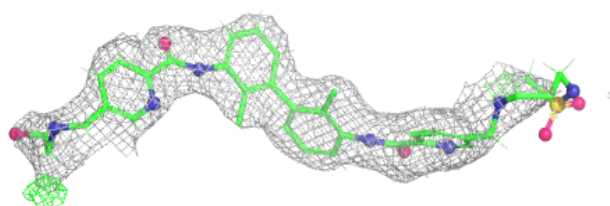
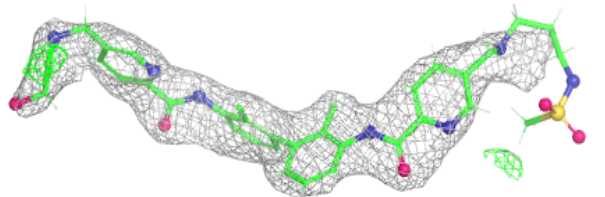
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	C	202	5/5	0.80	0.19	286,302,323,325	0
2	A1H56	D	201	47/47	0.89	0.12	41,59,203,217	0
2	A1H56	B	201	47/47	0.90	0.12	31,39,151,187	0
2	A1H56	C	201	47/47	0.91	0.11	39,54,155,202	0

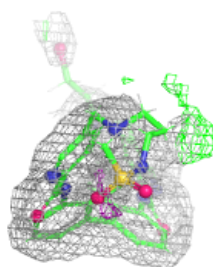
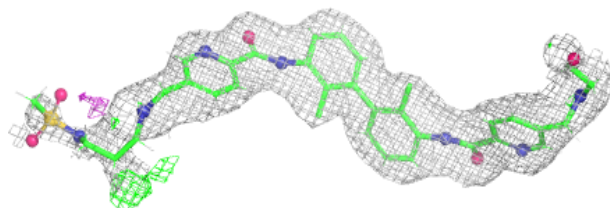
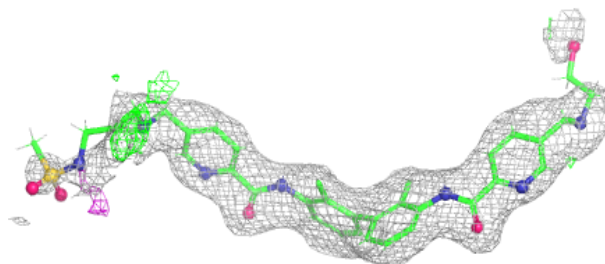
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

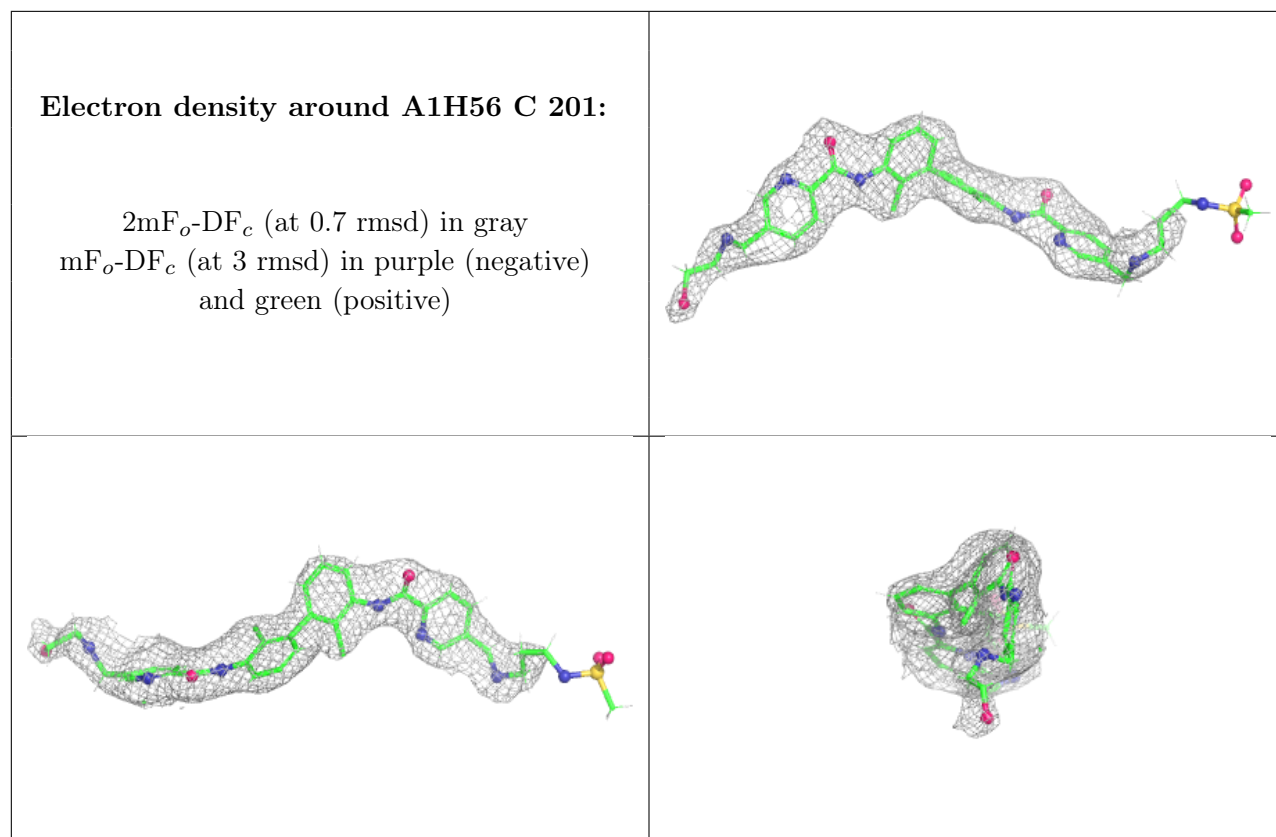
Electron density around A1H56 D 201:

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

**Electron density around A1H56 B 201:**

$2mF_o-DF_c$ (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.