



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:13 AM UTC

PDB ID : 3CD5 / pdb_00003cd5
Title : Thermodynamic and structure guided design of statin hmg-coa reductase inhibitors
Authors : Pavlovsky, A.; Sarver, R.W.; Harris, M.S.; Finzel, B.C.
Deposited on : 2008-02-26
Resolution : 2.39 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)
Xtrriage (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

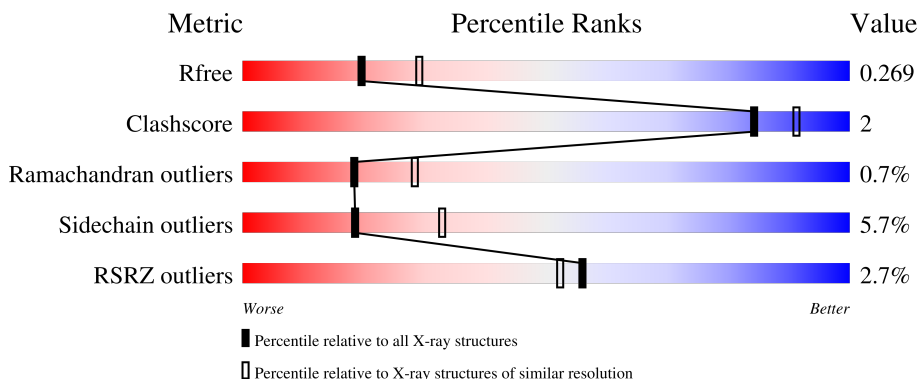
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	441	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 85%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 3%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">2% 85% 10% • •</p>
1	B	441	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 86%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 4%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">% 86% 10% •</p>
1	C	441	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 87%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 4%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">2% 87% 7% 5%</p>
1	D	441	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 86%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">6% 86% 7% • 6%</p>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13170 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 3-hydroxy-3-methylglutaryl-coenzyme A reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	425	3163	1971	556	606	30	0	0	0
1	B	425	3163	1971	556	606	30	0	0	0
1	C	418	3126	1947	552	597	30	0	4	0
1	D	416	3100	1932	546	592	30	0	1	0

There are 28 discrepancies between the modelled and reference sequences:

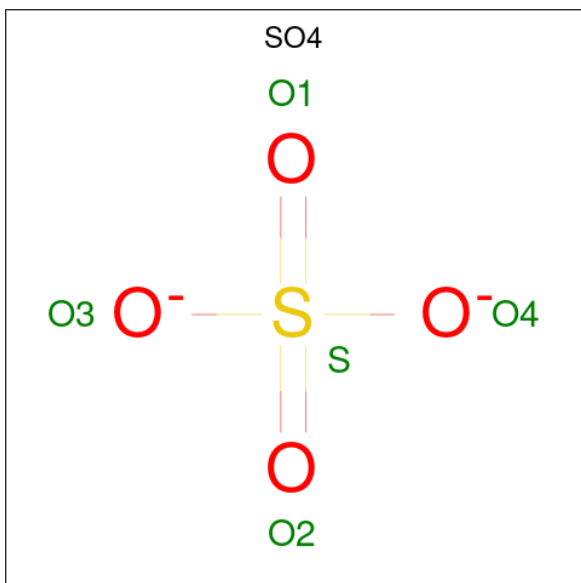
Chain	Residue	Modelled	Actual	Comment	Reference
A	435	HIS	-	expression tag	UNP P04035
A	436	HIS	-	expression tag	UNP P04035
A	437	HIS	-	expression tag	UNP P04035
A	438	HIS	-	expression tag	UNP P04035
A	439	HIS	-	expression tag	UNP P04035
A	440	HIS	-	expression tag	UNP P04035
A	485	ILE	MET	engineered mutation	UNP P04035
B	435	HIS	-	expression tag	UNP P04035
B	436	HIS	-	expression tag	UNP P04035
B	437	HIS	-	expression tag	UNP P04035
B	438	HIS	-	expression tag	UNP P04035
B	439	HIS	-	expression tag	UNP P04035
B	440	HIS	-	expression tag	UNP P04035
B	485	ILE	MET	engineered mutation	UNP P04035
C	435	HIS	-	expression tag	UNP P04035
C	436	HIS	-	expression tag	UNP P04035
C	437	HIS	-	expression tag	UNP P04035
C	438	HIS	-	expression tag	UNP P04035
C	439	HIS	-	expression tag	UNP P04035
C	440	HIS	-	expression tag	UNP P04035
C	485	ILE	MET	engineered mutation	UNP P04035

Continued on next page...

Continued from previous page...

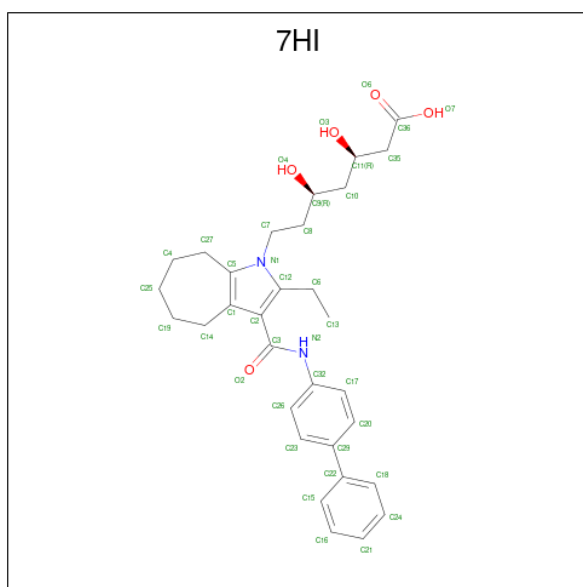
Chain	Residue	Modelled	Actual	Comment	Reference
D	435	HIS	-	expression tag	UNP P04035
D	436	HIS	-	expression tag	UNP P04035
D	437	HIS	-	expression tag	UNP P04035
D	438	HIS	-	expression tag	UNP P04035
D	439	HIS	-	expression tag	UNP P04035
D	440	HIS	-	expression tag	UNP P04035
D	485	ILE	MET	engineered mutation	UNP P04035

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is (3R,5R)-7-[3-(biphenyl-4-ylcarbonyl)-2-ethyl-5,6,7,8-tetrahydrocyclohepta[b]pyrrol-1(4H)-yl]-3,5-dihydroxyheptanoic acid (CCD ID: 7HI) (formula: C₃₁H₃₈N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	38	31	2	5	0	0
3	B	1	38	31	2	5	0	0
3	D	1	38	31	2	5	0	0
3	D	1	38	31	2	5	0	0

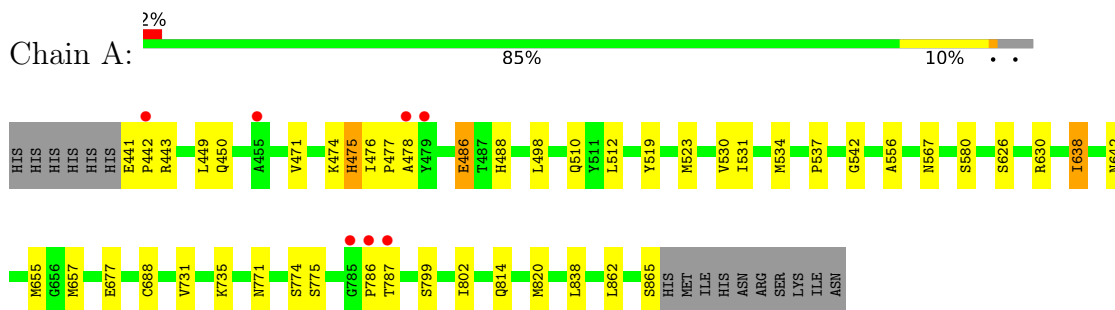
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	128	128	128	0	0
4	B	118	118	118	0	0
4	C	98	98	98	0	0
4	D	102	102	102	0	0

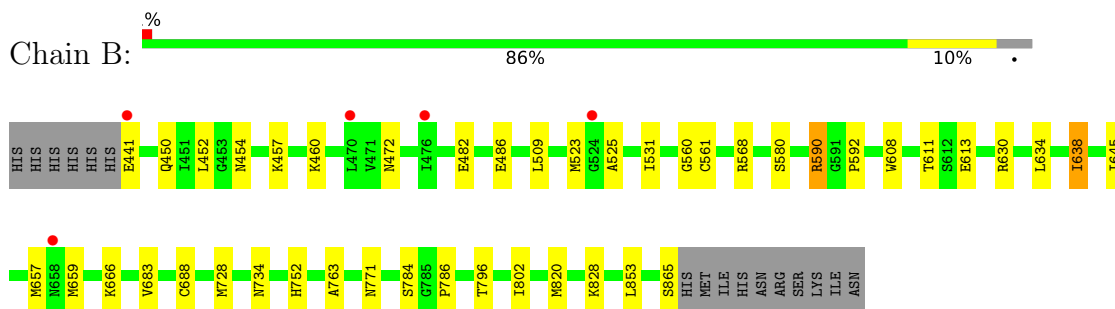
3 Residue-property plots [i](#)

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.

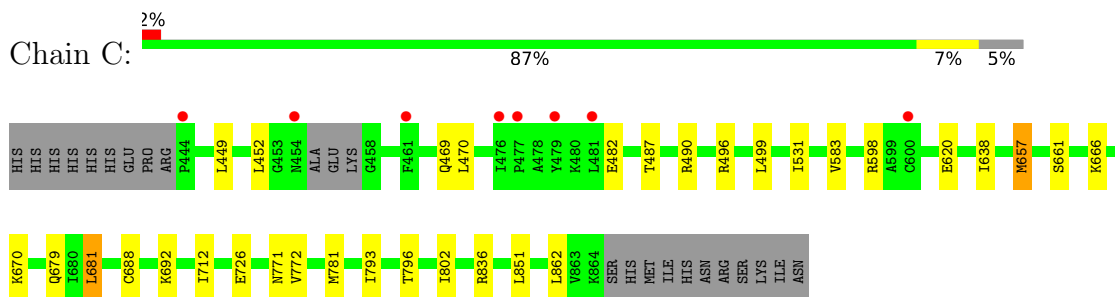
- Molecule 1: 3-hydroxy-3-methylglutaryl-coenzyme A reductase



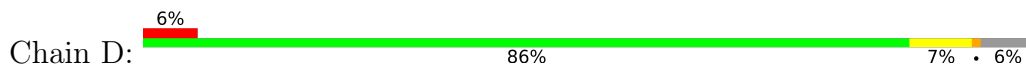
- Molecule 1: 3-hydroxy-3-methylglutaryl-coenzyme A reductase

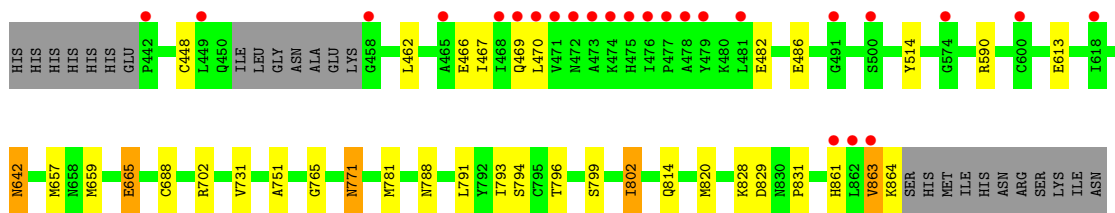


- Molecule 1: 3-hydroxy-3-methylglutaryl-coenzyme A reductase



- Molecule 1: 3-hydroxy-3-methylglutaryl-coenzyme A reductase





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.04Å 173.18Å 75.96Å 90.00° 118.76° 90.00°	Depositor
Resolution (Å)	50.00 – 2.39 50.00 – 2.39	Depositor EDS
% Data completeness (in resolution range)	60.8 (50.00-2.39) 60.8 (50.00-2.39)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.273 0.199 , 0.269	Depositor DCC
R_{free} test set	2012 reflections (3.07%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 26.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.004 for l,k,-h-l 0.004 for -h-l,k,h 0.044 for h,-k,-h-l 0.053 for l,-k,h 0.032 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13170	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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: 7HI, 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	0.48	0/3209	0.80	0/4338
1	B	0.49	0/3209	0.78	0/4338
1	C	0.48	0/3190	0.77	0/4309
1	D	0.48	0/3150	0.80	1/4257 (0.0%)
All	All	0.48	0/12758	0.79	1/17242 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	802	ILE	N-CA-C	5.61	115.71	107.75

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	3163	0	3205	22	0
1	B	3163	0	3205	16	0
1	C	3126	0	3172	14	0
1	D	3100	0	3141	23	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	38	0	37	0	0
3	B	38	0	37	1	0
3	D	76	0	74	4	0
4	A	128	0	0	0	0
4	B	118	0	0	0	0
4	C	98	0	0	0	0
4	D	102	0	0	0	0
All	All	13170	0	12871	60	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 2.

All (60) 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:D:863:VAL:HB	1:D:864:LYS:HA	1.39	1.04
1:D:863:VAL:CB	1:D:864:LYS:HA	1.94	0.98
1:C:771[B]:ASN:ND2	1:D:771[B]:ASN:OD1	2.10	0.83
1:D:863:VAL:CG2	1:D:864:LYS:HA	2.15	0.76
1:A:441:GLU:N	1:A:442:PRO:CD	2.62	0.63
1:D:863:VAL:HG23	1:D:864:LYS:HA	1.82	0.62
1:D:863:VAL:HB	1:D:864:LYS:CA	2.24	0.61
1:C:771[A]:ASN:ND2	1:D:771[A]:ASN:OD1	2.35	0.57
1:C:681:LEU:HD22	1:D:731:VAL:HG22	1.84	0.57
1:C:692:LYS:NZ	1:D:751:ALA:O	2.36	0.57
1:A:441:GLU:N	1:A:442:PRO:HD3	2.22	0.55
1:A:731:VAL:O	1:A:735:LYS:HB3	2.07	0.54
3:D:876:7HI:H13A	3:D:876:7HI:H8A	1.90	0.54
1:A:542:GLY:H	1:A:567:ASN:HD22	1.56	0.52
1:D:863:VAL:HG23	1:D:864:LYS:CA	2.39	0.52
1:A:531:ILE:HD11	1:B:820:MET:HE1	1.91	0.52
1:C:583:VAL:HG11	1:C:836[B]:ARG:HD2	1.93	0.50
1:D:642:ASN:HD22	1:D:642:ASN:N	2.09	0.50
1:A:523:MET:HE1	1:A:530:VAL:HG21	1.93	0.50
1:C:657:MET:HA	1:C:657:MET:HE2	1.94	0.49
1:B:638:ILE:O	1:C:796:THR:HG21	2.14	0.48
1:D:590:ARG:CZ	1:D:657:MET:HE1	2.44	0.47
1:C:499:LEU:HD21	1:D:820:MET:HE3	1.97	0.46
1:A:642:ASN:HD21	1:B:734:ASN:CG	2.23	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:462:LEU:HD12	1:D:467:ILE:HD11	1.99	0.45
1:C:496:ARG:HD2	1:C:531:ILE:O	2.17	0.45
1:A:774:SER:HA	1:A:799:SER:O	2.17	0.45
1:C:781:MET:HE2	1:C:793:ILE:HD12	1.99	0.44
1:D:861:HIS:HB3	3:D:3:7HI:H19A	1.99	0.44
1:A:771:ASN:OD1	1:A:775:SER:OG	2.34	0.44
1:B:560:GLY:O	1:B:561:CYS:HB2	2.18	0.43
1:C:712:ILE:HG13	1:C:851:LEU:HD11	1.99	0.43
1:A:531:ILE:CD1	1:B:820:MET:HE1	2.48	0.43
1:A:820:MET:HE2	1:B:509:LEU:CD2	2.48	0.43
1:B:608:TRP:O	1:B:611:THR:OG1	2.35	0.42
1:A:542:GLY:H	1:A:567:ASN:ND2	2.16	0.42
1:B:752:HIS:ND1	1:B:853:LEU:HD22	2.35	0.42
1:B:454:ASN:HB3	1:B:457:LYS:HB2	2.01	0.42
1:A:638:ILE:O	1:D:796:THR:HG21	2.20	0.42
1:B:592:PRO:HD2	1:B:645:ILE:O	2.19	0.42
3:D:3:7HI:H13A	3:D:3:7HI:H8A	2.02	0.42
1:B:796:THR:HG21	1:C:638:ILE:O	2.20	0.42
1:A:519:TYR:HB3	1:A:523:MET:HE3	2.02	0.42
1:A:471:VAL:HA	1:A:476:ILE:O	2.20	0.41
1:C:862:LEU:HB2	1:D:665:GLU:HG3	2.02	0.41
1:D:781:MET:HE2	1:D:793:ILE:HD12	2.02	0.41
1:B:590:ARG:CZ	1:B:657:MET:HE1	2.51	0.41
1:D:702:ARG:O	1:D:799:SER:HA	2.19	0.41
1:D:863:VAL:CG2	3:D:3:7HI:H14A	2.51	0.41
1:D:765:GLY:CA	1:D:814:GLN:HG2	2.51	0.41
1:A:820:MET:HE1	1:B:531:ILE:HD11	2.01	0.41
1:A:862:LEU:HD23	1:B:683:VAL:HB	2.02	0.41
1:D:863:VAL:HG23	1:D:864:LYS:HB3	2.03	0.41
1:B:568:ARG:HD3	3:B:1:7HI:C15	2.52	0.41
1:A:488:HIS:CD2	1:A:523:MET:HG3	2.56	0.40
1:C:772:VAL:HG23	1:D:771[B]:ASN:HD21	1.86	0.40
1:A:534:MET:HE3	1:B:763:ALA:O	2.21	0.40
1:A:537:PRO:O	1:A:556:ALA:HA	2.21	0.40
1:A:655:MET:SD	1:A:657:MET:HG2	2.61	0.40
1:A:802:ILE:HG12	1:A:838:LEU:HD23	2.02	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	423/441 (96%)	401 (95%)	16 (4%)	6 (1%)	9	13
1	B	423/441 (96%)	404 (96%)	17 (4%)	2 (0%)	24	37
1	C	418/441 (95%)	401 (96%)	17 (4%)	0	100	100
1	D	413/441 (94%)	391 (95%)	19 (5%)	3 (1%)	18	28
All	All	1677/1764 (95%)	1597 (95%)	69 (4%)	11 (1%)	18	28

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	474	LYS
1	A	478	ALA
1	A	486	GLU
1	B	525	ALA
1	A	786	PRO
1	B	786	PRO
1	A	475	HIS
1	A	477	PRO
1	D	514	TYR
1	D	831	PRO
1	D	863	VAL

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	339/355 (96%)	322 (95%)	17 (5%)	22	38
1	B	339/355 (96%)	316 (93%)	23 (7%)	14	25
1	C	337/355 (95%)	319 (95%)	18 (5%)	20	36
1	D	333/355 (94%)	314 (94%)	19 (6%)	18	33
All	All	1348/1420 (95%)	1271 (94%)	77 (6%)	18	33

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	443	ARG
1	A	449	LEU
1	A	450	GLN
1	A	475	HIS
1	A	486	GLU
1	A	498	LEU
1	A	510	GLN
1	A	512	LEU
1	A	580	SER
1	A	626	SER
1	A	630	ARG
1	A	638	ILE
1	A	677	GLU
1	A	688	CYS
1	A	787	THR
1	A	814	GLN
1	A	865	SER
1	B	441	GLU
1	B	450	GLN
1	B	452	LEU
1	B	460	LYS
1	B	472	ASN
1	B	482	GLU
1	B	486	GLU
1	B	523	MET
1	B	580	SER
1	B	590	ARG
1	B	613	GLU
1	B	630	ARG
1	B	634	LEU
1	B	638	ILE
1	B	659	MET
1	B	666	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	688	CYS
1	B	728	MET
1	B	771	ASN
1	B	784	SER
1	B	802	ILE
1	B	828	LYS
1	B	865	SER
1	C	449	LEU
1	C	452	LEU
1	C	469	GLN
1	C	470	LEU
1	C	482	GLU
1	C	487	THR
1	C	490	ARG
1	C	598	ARG
1	C	620	GLU
1	C	657	MET
1	C	661	SER
1	C	666	LYS
1	C	670	LYS
1	C	679	GLN
1	C	681	LEU
1	C	688	CYS
1	C	726	GLU
1	C	802	ILE
1	D	448	CYS
1	D	466	GLU
1	D	469	GLN
1	D	470	LEU
1	D	482	GLU
1	D	486	GLU
1	D	613	GLU
1	D	642	ASN
1	D	659	MET
1	D	665	GLU
1	D	688	CYS
1	D	771[A]	ASN
1	D	771[B]	ASN
1	D	788	ASN
1	D	791	LEU
1	D	794	SER
1	D	802	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	828	LYS
1	D	829	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	488	HIS
1	A	510	GLN
1	A	567	ASN
1	A	632	GLN
1	A	770	GLN
1	A	788	ASN
1	B	469	GLN
1	B	510	GLN
1	B	529	ASN
1	C	469	GLN
1	C	472	ASN
1	C	510	GLN
1	C	632	GLN
1	C	642	ASN
1	C	819	GLN
1	D	472	ASN
1	D	518	ASN
1	D	567	ASN
1	D	788	ASN
1	D	830	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

8 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	SO4	C	3	-	4,4,4	0.25	0	6,6,6	0.17	0
3	7HI	D	3	-	41,41,41	1.25	5 (12%)	48,56,56	1.35	7 (14%)
2	SO4	B	2	-	4,4,4	0.20	0	6,6,6	0.15	0
3	7HI	A	2	-	41,41,41	1.25	5 (12%)	48,56,56	1.18	7 (14%)
3	7HI	B	1	-	41,41,41	1.27	5 (12%)	48,56,56	1.23	7 (14%)
2	SO4	D	4	-	4,4,4	0.29	0	6,6,6	0.16	0
2	SO4	A	1	-	4,4,4	0.23	0	6,6,6	0.21	0
3	7HI	D	876	-	41,41,41	1.25	5 (12%)	48,56,56	1.27	7 (14%)

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	7HI	D	876	-	-	4/27/35/35	0/4/4/4
3	7HI	D	3	-	-	5/27/35/35	0/4/4/4
3	7HI	B	1	-	-	2/27/35/35	0/4/4/4
3	7HI	A	2	-	-	6/27/35/35	0/4/4/4

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3	7HI	C19-C25	-4.54	1.33	1.51
3	A	2	7HI	C19-C25	-4.52	1.33	1.51
3	B	1	7HI	C19-C25	-4.51	1.33	1.51
3	D	876	7HI	C19-C25	-4.48	1.33	1.51
3	A	2	7HI	O6-C36	3.63	1.33	1.22
3	D	876	7HI	O6-C36	3.44	1.33	1.22

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3	7HI	O6-C36	3.44	1.33	1.22
3	B	1	7HI	O6-C36	3.41	1.33	1.22
3	D	3	7HI	O7-C36	-2.78	1.21	1.30
3	D	876	7HI	C2-C1	-2.76	1.38	1.44
3	B	1	7HI	O7-C36	-2.67	1.22	1.30
3	B	1	7HI	C2-C1	-2.64	1.38	1.44
3	D	3	7HI	C2-C1	-2.63	1.38	1.44
3	B	1	7HI	C32-N2	-2.63	1.36	1.41
3	A	2	7HI	C2-C1	-2.63	1.38	1.44
3	D	876	7HI	O7-C36	-2.61	1.22	1.30
3	A	2	7HI	O7-C36	-2.47	1.22	1.30
3	D	876	7HI	C32-N2	-2.45	1.36	1.41
3	A	2	7HI	C32-N2	-2.44	1.36	1.41
3	D	3	7HI	C32-N2	-2.30	1.37	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	3	7HI	C25-C19-C14	3.63	124.04	114.74
3	D	876	7HI	O2-C3-C2	-3.50	116.28	121.53
3	D	3	7HI	C8-C7-N1	-3.05	106.83	112.16
3	D	3	7HI	O2-C3-C2	-3.03	116.99	121.53
3	B	1	7HI	O2-C3-C2	-3.00	117.03	121.53
3	D	3	7HI	C10-C11-C35	-2.74	107.37	112.96
3	D	876	7HI	C25-C19-C14	2.74	121.76	114.74
3	A	2	7HI	C8-C7-N1	-2.71	107.42	112.16
3	A	2	7HI	O2-C3-C2	-2.69	117.50	121.53
3	D	3	7HI	C6-C12-C2	-2.67	125.95	130.36
3	A	2	7HI	C25-C19-C14	2.58	121.35	114.74
3	B	1	7HI	C8-C7-N1	-2.56	107.67	112.16
3	B	1	7HI	C25-C19-C14	2.55	121.28	114.74
3	D	3	7HI	C2-C1-C5	-2.53	104.71	107.61
3	B	1	7HI	C10-C11-C35	-2.48	107.90	112.96
3	B	1	7HI	C6-C12-C2	-2.47	126.28	130.36
3	A	2	7HI	C6-C12-C2	-2.44	126.33	130.36
3	D	876	7HI	C2-C1-C5	-2.40	104.86	107.61
3	A	2	7HI	C2-C1-C5	-2.31	104.97	107.61
3	B	1	7HI	C2-C3-N2	2.26	119.67	115.81
3	A	2	7HI	C32-N2-C3	-2.23	123.50	127.45
3	D	876	7HI	C6-C12-C2	-2.20	126.72	130.36
3	B	1	7HI	C2-C1-C5	-2.20	105.09	107.61
3	D	876	7HI	C2-C3-N2	2.17	119.53	115.81

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	876	7HI	C32-N2-C3	-2.06	123.80	127.45
3	D	876	7HI	C8-C7-N1	-2.04	108.58	112.16
3	D	3	7HI	C25-C4-C27	-2.01	109.57	114.74
3	A	2	7HI	C2-C12-N1	-2.01	106.34	108.30

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	3	7HI	C18-C22-C29-C23
3	D	3	7HI	C15-C22-C29-C23
3	D	3	7HI	C15-C22-C29-C20
3	D	3	7HI	C18-C22-C29-C20
3	D	3	7HI	O3-C11-C35-C36
3	A	2	7HI	C15-C22-C29-C20
3	A	2	7HI	C15-C22-C29-C23
3	A	2	7HI	C18-C22-C29-C23
3	A	2	7HI	C26-C32-N2-C3
3	A	2	7HI	C17-C32-N2-C3
3	A	2	7HI	C18-C22-C29-C20
3	B	1	7HI	C17-C32-N2-C3
3	B	1	7HI	C26-C32-N2-C3
3	D	876	7HI	C18-C22-C29-C23
3	D	876	7HI	C15-C22-C29-C20
3	D	876	7HI	C18-C22-C29-C20
3	D	876	7HI	C15-C22-C29-C23

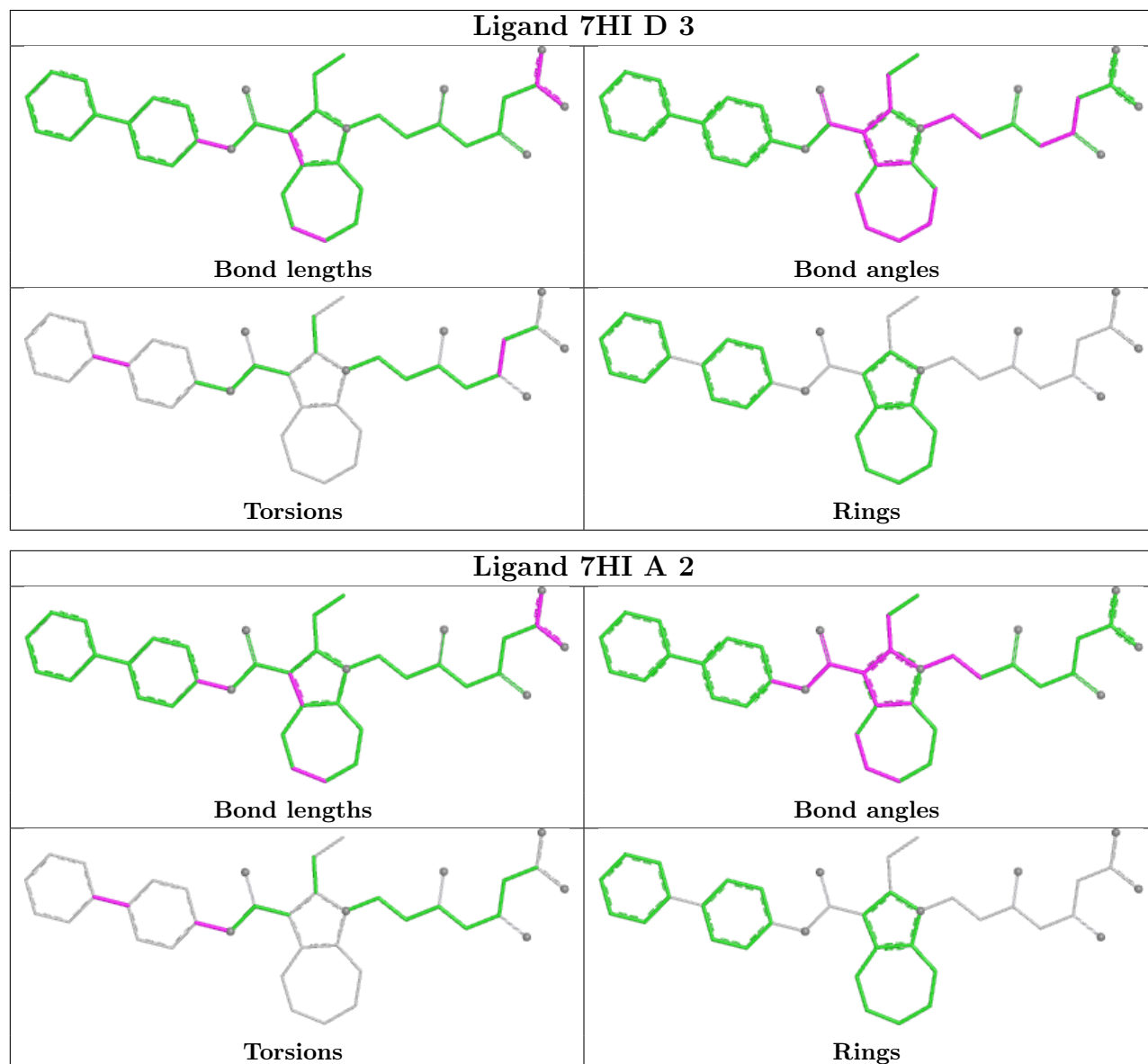
There are no ring outliers.

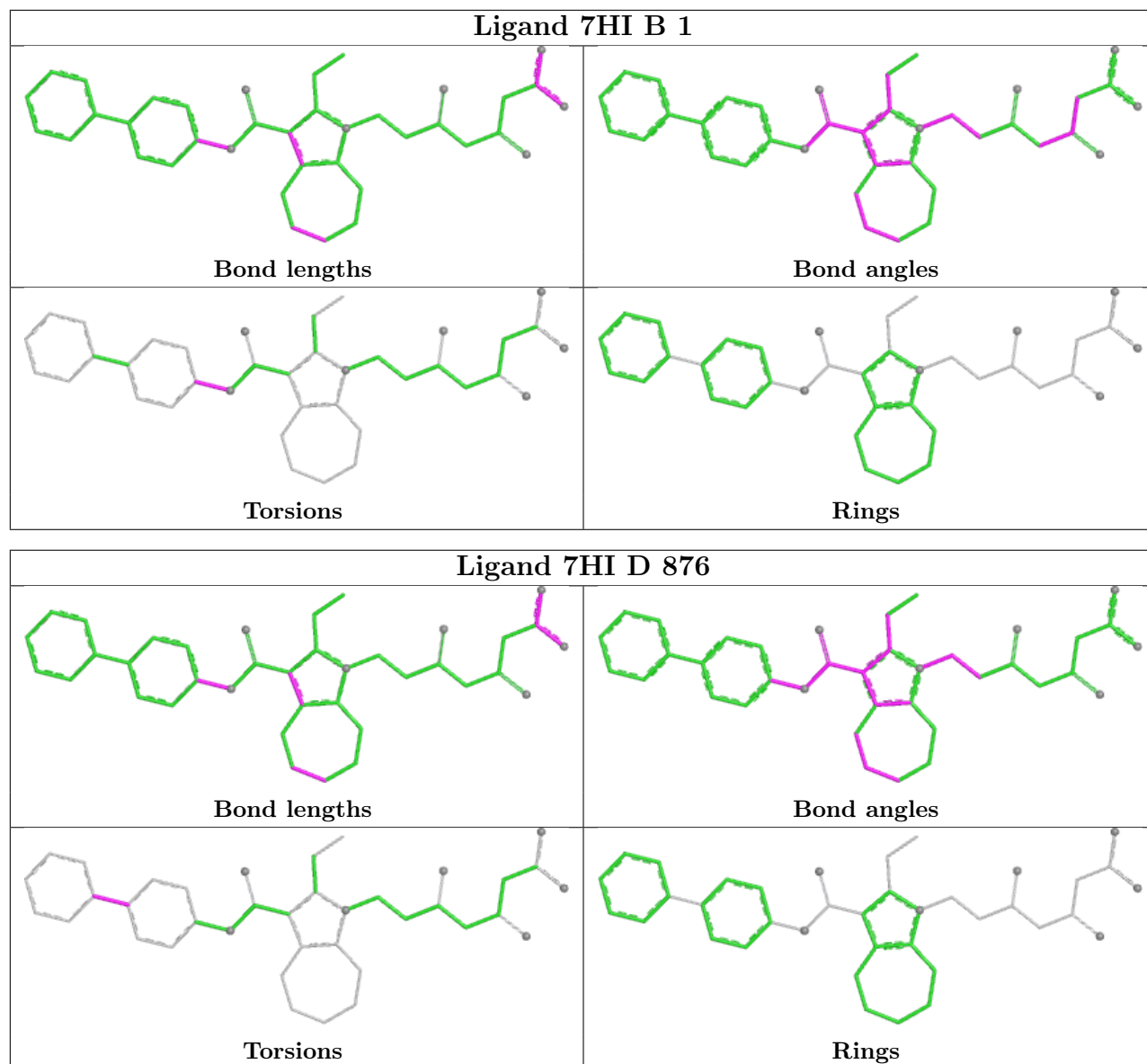
3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3	7HI	3	0
3	B	1	7HI	1	0
3	D	876	7HI	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	425/441 (96%)	0.07	7 (1%) 70 66	26, 39, 62, 70	0
1	B	425/441 (96%)	0.07	5 (1%) 76 73	26, 38, 66, 72	0
1	C	418/441 (94%)	0.10	8 (1%) 66 62	19, 39, 74, 80	4 (0%)
1	D	416/441 (94%)	0.29	25 (6%) 27 24	24, 40, 71, 76	1 (0%)
All	All	1684/1764 (95%)	0.13	45 (2%) 56 52	19, 39, 68, 80	5 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	473	ALA	6.2
1	D	862	LEU	4.3
1	D	469	GLN	4.3
1	D	475	HIS	3.7
1	D	479	TYR	3.6
1	D	481	LEU	3.6
1	D	476	ILE	3.3
1	D	442	PRO	3.2
1	D	478	ALA	3.0
1	C	477	PRO	2.8
1	D	471	VAL	2.8
1	D	491	GLY	2.7
1	D	449	LEU	2.7
1	A	479	TYR	2.7
1	D	477	PRO	2.7
1	B	658	ASN	2.7
1	D	863	VAL	2.6
1	D	468	ILE	2.6
1	D	861	HIS	2.6
1	C	479	TYR	2.6
1	D	618	ILE	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	478	ALA	2.6
1	D	472	ASN	2.6
1	D	600	CYS	2.4
1	C	461	PHE	2.3
1	A	785	GLY	2.3
1	B	470	LEU	2.3
1	A	786	PRO	2.3
1	C	481	LEU	2.2
1	D	458	GLY	2.2
1	D	474	LYS	2.2
1	B	476	ILE	2.2
1	B	524	GLY	2.2
1	A	787	THR	2.2
1	C	444	PRO	2.2
1	D	574	GLY	2.2
1	A	455	ALA	2.2
1	C	454	ASN	2.2
1	A	442	PRO	2.2
1	C	600	CYS	2.2
1	D	470	LEU	2.1
1	D	465	ALA	2.1
1	D	500	SER	2.1
1	B	441	GLU	2.1
1	C	476	ILE	2.1

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](#)

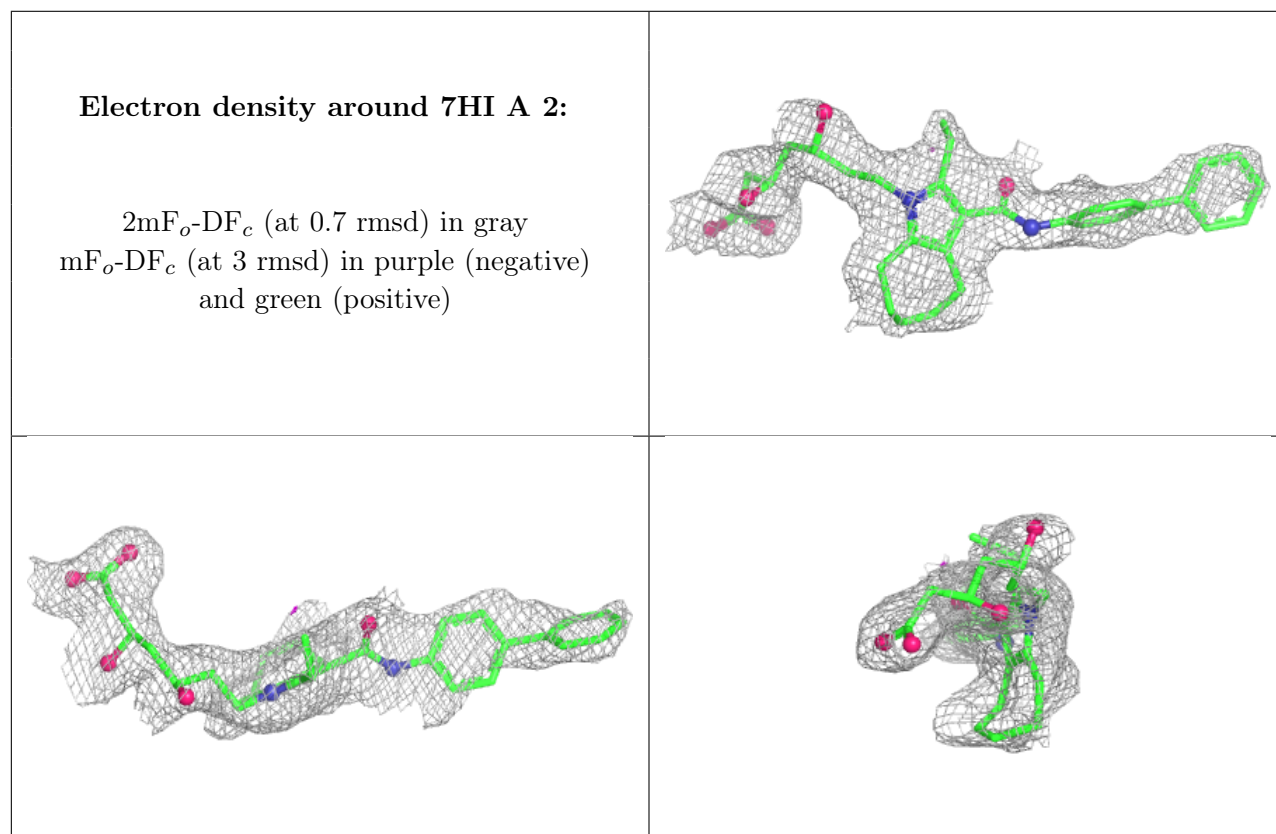
There are no oligosaccharides in this entry.

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.

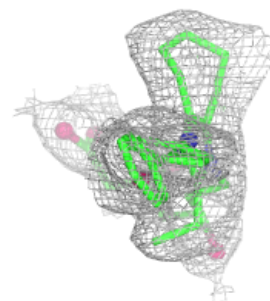
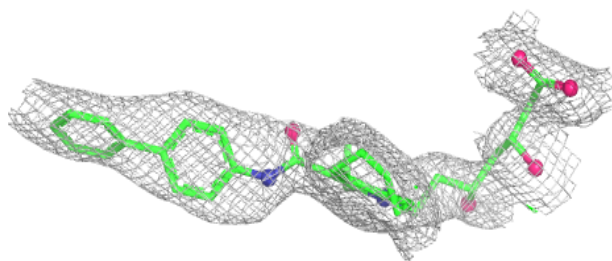
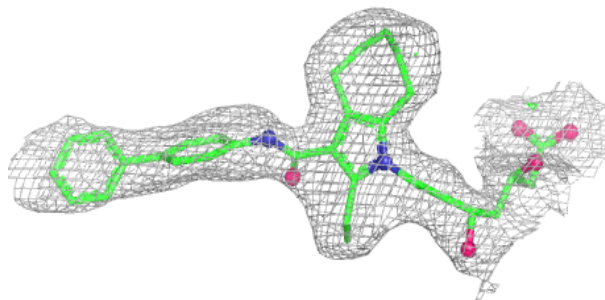
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	C	3	5/5	0.88	0.14	64,64,65,65	0
2	SO4	D	4	5/5	0.88	0.15	56,56,57,57	0
2	SO4	B	2	5/5	0.93	0.11	55,55,55,56	0
3	7HI	A	2	38/38	0.93	0.09	28,35,54,54	0
2	SO4	A	1	5/5	0.94	0.09	59,60,60,60	0
3	7HI	B	1	38/38	0.94	0.08	27,32,44,44	0
3	7HI	D	3	38/38	0.94	0.08	28,34,42,42	0
3	7HI	D	876	38/38	0.94	0.08	29,38,39,39	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.

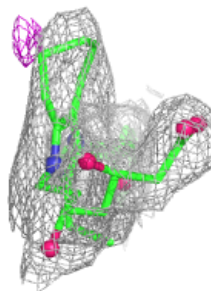
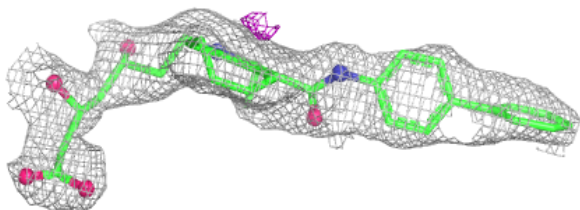
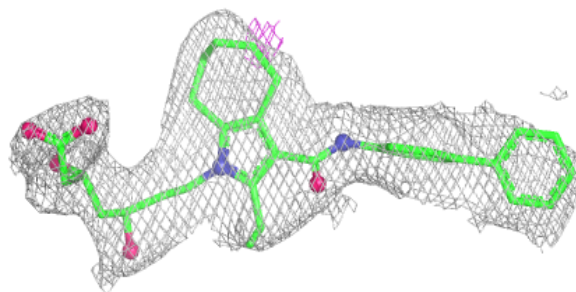


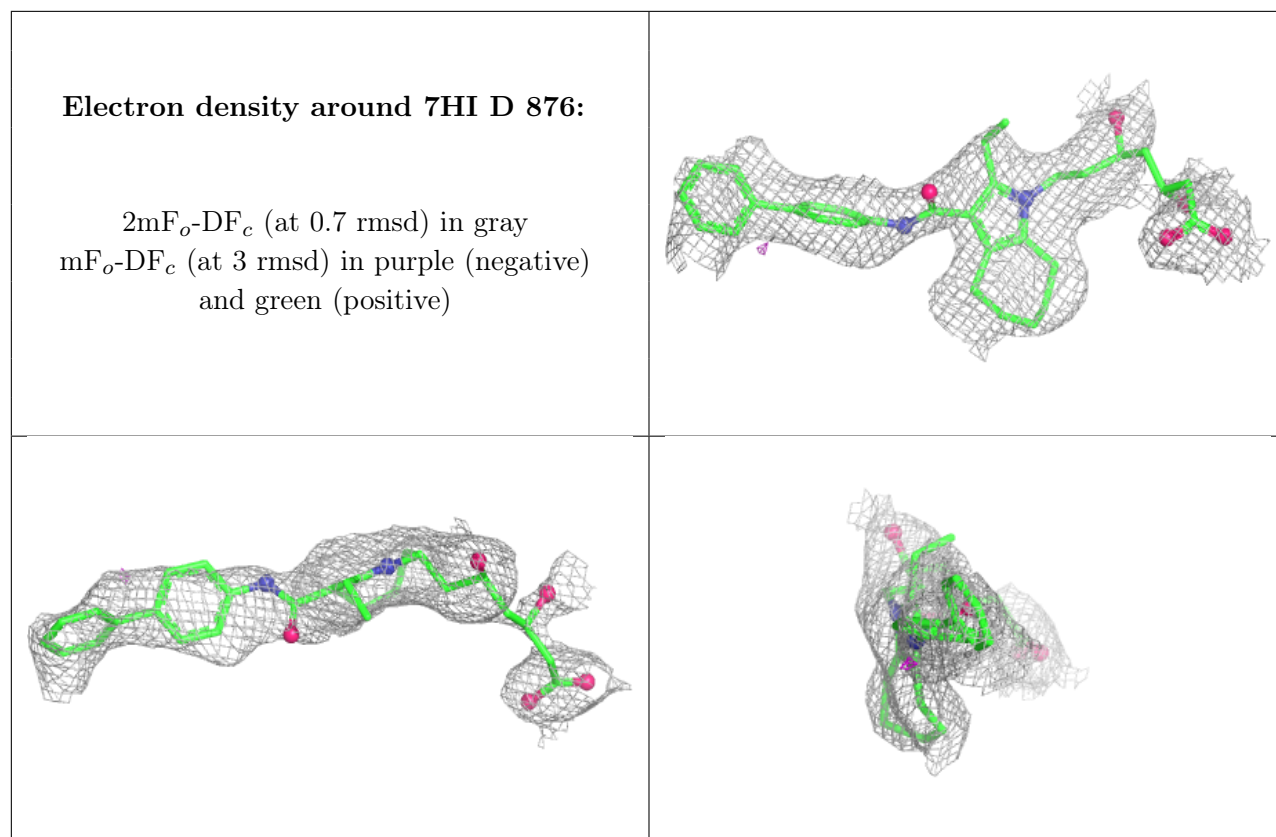
Electron density around 7HI B 1:

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

**Electron density around 7HI D 3:**

$2mF_o-DF_c$ (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.