



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:15 AM UTC

PDB ID : 6HYR / pdb_00006hyr
Title : THE GLIC PENTAMERIC LIGAND-GATED ION CHANNEL MUTANT Q193C+MMTS
Authors : Hu, H.D.; Delarue, M.
Deposited on : 2018-10-22
Resolution : 3.50 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

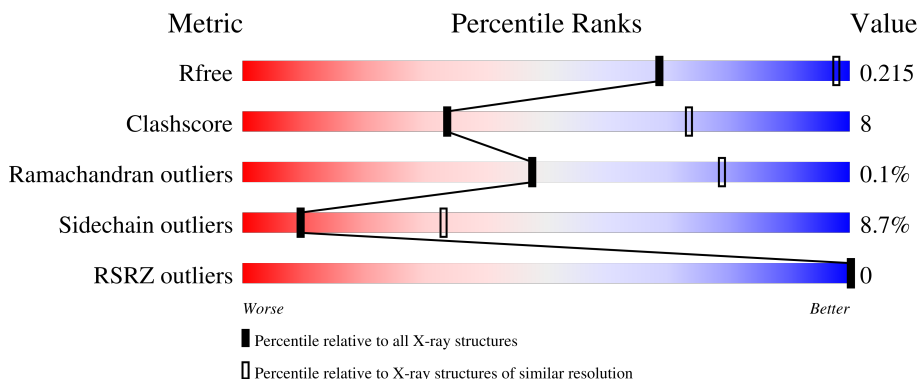
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	311	 77% 20% ..
1	B	311	 78% 18% ..
1	C	311	 79% 17% ..
1	D	311	 77% 21% ..
1	E	311	 78% 19% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12632 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.

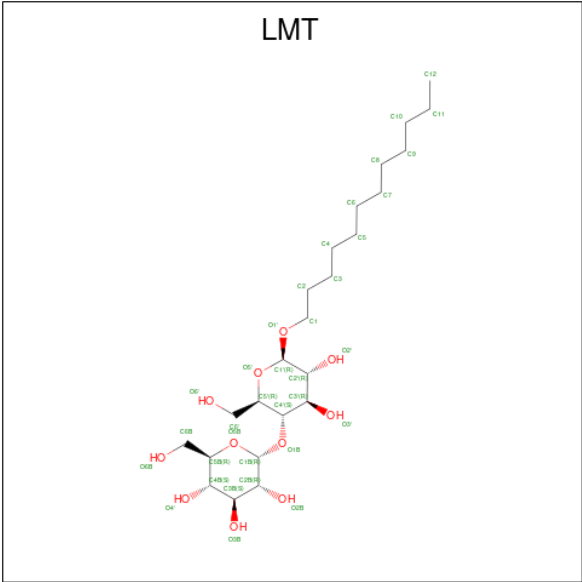
- Molecule 1 is a protein called Proton-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2524	1663	403	452	6			
1	B	311	Total	C	N	O	S	0	0	0
			2524	1663	403	452	6			
1	C	311	Total	C	N	O	S	0	0	0
			2524	1663	403	452	6			
1	D	311	Total	C	N	O	S	0	0	0
			2524	1663	403	452	6			
1	E	311	Total	C	N	O	S	0	0	0
			2524	1663	403	452	6			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	193	SCH	GLN	engineered mutation	UNP Q7NDN8
B	193	SCH	GLN	engineered mutation	UNP Q7NDN8
C	193	SCH	GLN	engineered mutation	UNP Q7NDN8
D	193	SCH	GLN	engineered mutation	UNP Q7NDN8
E	193	SCH	GLN	engineered mutation	UNP Q7NDN8

- Molecule 2 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: C₂₄H₄₆O₁₁).

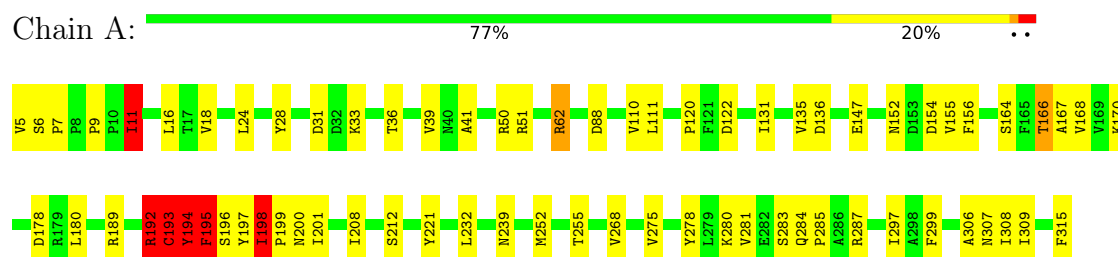


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C	0	0
			12	12		

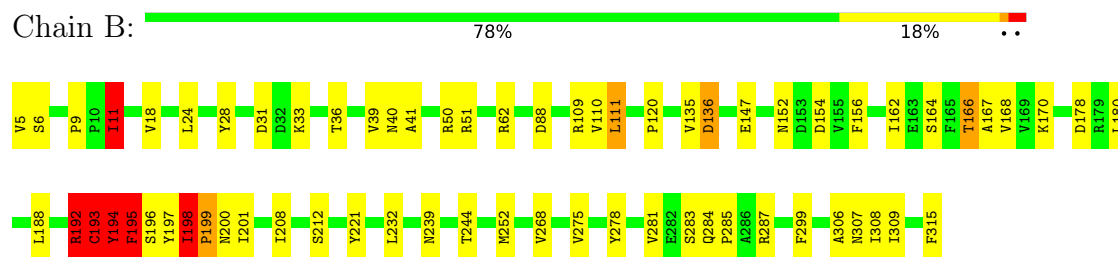
3 Residue-property plots

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.

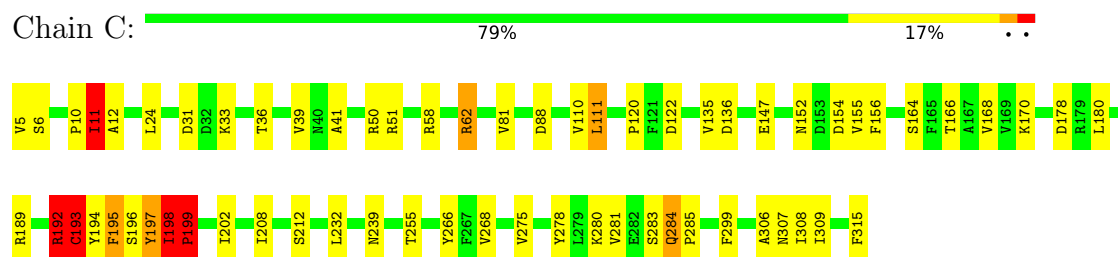
• Molecule 1: Proton-gated ion channel



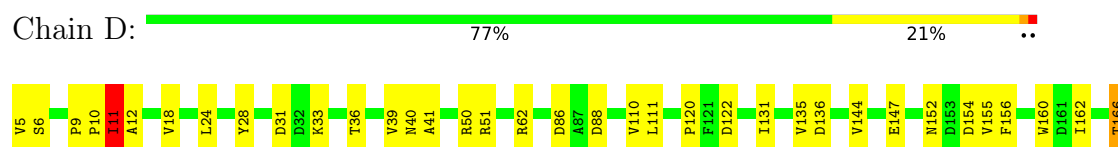
• Molecule 1: Proton-gated ion channel



• Molecule 1: Proton-gated ion channel

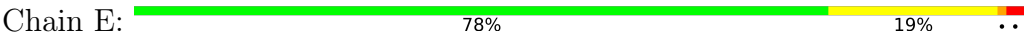


• Molecule 1: Proton-gated ion channel





● Molecule 1: Proton-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.98Å 128.28Å 160.43Å 90.00° 100.96° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 20.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	78.4 (20.00-3.50) 78.4 (20.00-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 3.51Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.231 , 0.241 0.155 , 0.215	Depositor DCC
R_{free} test set	1775 reflections (3.97%)	wwPDB-VP
Wilson B-factor (Å ²)	61.2	Xtriage
Anisotropy	0.563	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 71.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12632	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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: LMT, SCH

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	1.22	12/2583 (0.5%)	1.28	11/3530 (0.3%)
1	B	1.23	9/2583 (0.3%)	1.27	8/3530 (0.2%)
1	C	1.22	5/2583 (0.2%)	1.27	10/3530 (0.3%)
1	D	1.23	6/2583 (0.2%)	1.22	5/3530 (0.1%)
1	E	1.23	9/2583 (0.3%)	1.26	9/3530 (0.3%)
All	All	1.23	41/12915 (0.3%)	1.26	43/17650 (0.2%)

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	3
1	B	0	3
1	C	0	2
1	D	0	3
1	E	0	3
All	All	0	14

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	189	ARG	C-N	-10.86	1.18	1.33
1	B	195	PHE	C-O	6.92	1.32	1.24
1	B	278	TYR	C-O	-6.75	1.16	1.24
1	A	278	TYR	C-O	-6.60	1.16	1.24
1	E	278	TYR	C-O	-6.48	1.16	1.24
1	D	278	TYR	C-O	-6.39	1.16	1.24
1	C	278	TYR	C-O	-6.34	1.16	1.24
1	B	201	ILE	N-CA	6.27	1.51	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	201	ILE	N-CA	6.17	1.51	1.46
1	A	195	PHE	C-O	6.15	1.31	1.24
1	E	189	ARG	C-N	-6.14	1.25	1.33
1	A	189	ARG	C-N	-6.08	1.25	1.33
1	A	18	VAL	C-O	-5.99	1.17	1.24
1	B	308	ILE	C-O	-5.97	1.17	1.24
1	E	195	PHE	C-O	5.93	1.31	1.24
1	C	308	ILE	C-O	-5.86	1.17	1.24
1	E	164	SER	C-O	-5.80	1.16	1.23
1	B	164	SER	C-O	-5.76	1.16	1.23
1	A	164	SER	C-O	-5.68	1.16	1.23
1	C	41	ALA	C-O	-5.59	1.16	1.23
1	E	308	ILE	C-O	-5.57	1.17	1.24
1	E	41	ALA	C-O	-5.56	1.16	1.23
1	B	41	ALA	C-O	-5.55	1.16	1.23
1	E	201	ILE	N-CA	5.52	1.51	1.46
1	C	164	SER	C-O	-5.51	1.16	1.23
1	A	308	ILE	C-O	-5.48	1.17	1.24
1	A	201	ILE	N-CA	5.43	1.51	1.46
1	D	308	ILE	C-O	-5.43	1.17	1.24
1	E	7	PRO	CA-C	5.37	1.55	1.52
1	C	197	TYR	CA-C	-5.31	1.46	1.52
1	D	18	VAL	C-O	-5.31	1.18	1.24
1	A	41	ALA	C-O	-5.30	1.17	1.23
1	A	7	PRO	CA-C	5.24	1.55	1.52
1	A	221	TYR	C-O	-5.24	1.18	1.24
1	A	195	PHE	CA-C	5.23	1.59	1.52
1	B	221	TYR	C-O	-5.22	1.18	1.24
1	B	244	THR	C-O	-5.18	1.18	1.24
1	A	16	LEU	C-O	-5.17	1.17	1.24
1	B	18	VAL	C-O	-5.11	1.18	1.24
1	D	41	ALA	C-O	-5.01	1.17	1.23
1	E	11	ILE	CA-CB	5.00	1.59	1.54

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	198	ILE	CA-C-N	-9.95	109.40	120.45
1	C	198	ILE	C-N-CA	-9.95	109.40	120.45
1	A	194	TYR	O-C-N	-9.87	107.20	123.00
1	E	198	ILE	C-N-CD	-8.96	88.28	125.00
1	A	198	ILE	C-N-CD	-8.95	88.32	125.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	194	TYR	O-C-N	-8.73	109.03	123.00
1	B	198	ILE	C-N-CD	-8.54	89.98	125.00
1	C	189	ARG	O-C-N	8.47	133.10	123.19
1	B	194	TYR	O-C-N	-8.36	109.62	123.00
1	A	189	ARG	O-C-N	7.78	132.29	123.19
1	C	189	ARG	CA-C-N	-7.23	111.88	122.58
1	C	189	ARG	C-N-CA	-7.23	111.88	122.58
1	C	198	ILE	C-N-CD	-7.05	96.09	125.00
1	D	9	PRO	O-C-N	6.82	124.36	121.15
1	D	11	ILE	N-CA-CB	6.78	118.31	110.65
1	E	11	ILE	N-CA-CB	6.73	118.25	110.65
1	C	11	ILE	N-CA-CB	6.58	118.08	110.65
1	A	11	ILE	N-CA-CB	6.55	118.06	110.65
1	A	189	ARG	CA-C-N	-6.51	112.95	122.58
1	A	189	ARG	C-N-CA	-6.51	112.95	122.58
1	B	11	ILE	N-CA-CB	6.15	117.60	110.65
1	E	189	ARG	O-C-N	6.12	130.35	123.19
1	D	192	ARG	CB-CA-C	5.99	121.49	110.10
1	A	9	PRO	O-C-N	5.98	123.96	121.15
1	E	192	ARG	CB-CA-C	5.98	121.45	110.10
1	B	192	ARG	CB-CA-C	5.97	121.44	110.10
1	A	192	ARG	CB-CA-C	5.97	121.44	110.10
1	C	192	ARG	CB-CA-C	5.96	121.43	110.10
1	B	195	PHE	CA-C-O	5.91	126.82	120.20
1	A	195	PHE	CA-C-O	5.61	126.48	120.20
1	C	199	PRO	N-CA-C	5.57	121.32	114.35
1	C	62	ARG	CA-CB-CG	-5.56	102.98	114.10
1	E	197	TYR	CB-CA-C	-5.44	102.34	110.88
1	E	195	PHE	CA-C-O	5.34	126.18	120.20
1	B	9	PRO	O-C-N	5.25	123.62	121.15
1	A	287	ARG	CA-C-N	5.12	127.45	120.54
1	A	287	ARG	C-N-CA	5.12	127.45	120.54
1	D	144	VAL	N-CA-C	5.07	115.16	107.80
1	E	287	ARG	CA-C-N	5.05	127.35	120.54
1	E	287	ARG	C-N-CA	5.05	127.35	120.54
1	B	287	ARG	CA-C-N	5.02	127.32	120.54
1	B	287	ARG	C-N-CA	5.02	127.32	120.54
1	D	86	ASP	CA-C-O	-5.00	115.81	121.66

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	192	ARG	Mainchain
1	A	193	SCH	Mainchain
1	A	194	TYR	Mainchain
1	B	192	ARG	Mainchain
1	B	193	SCH	Mainchain
1	B	194	TYR	Mainchain
1	C	192	ARG	Mainchain
1	C	193	SCH	Mainchain
1	D	192	ARG	Mainchain
1	D	193	SCH	Mainchain
1	D	194	TYR	Mainchain
1	E	192	ARG	Mainchain
1	E	193	SCH	Mainchain
1	E	194	TYR	Mainchain

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	2524	0	2542	43	0
1	B	2524	0	2542	38	0
1	C	2524	0	2543	42	0
1	D	2524	0	2543	38	0
1	E	2524	0	2541	46	0
2	A	12	0	23	0	0
All	All	12632	0	12734	194	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 8.

All (194) 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:C:198:ILE:CG2	1:C:199:PRO:HD3	1.25	1.56
1:C:198:ILE:HG22	1:C:199:PRO:CD	1.31	1.55
1:B:198:ILE:HG22	1:B:199:PRO:CD	1.42	1.48
1:B:198:ILE:CG2	1:B:199:PRO:HD3	1.44	1.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:198:ILE:HG22	1:E:199:PRO:CD	1.44	1.46
1:A:198:ILE:CG2	1:A:199:PRO:HD3	1.45	1.45
1:E:198:ILE:CG2	1:E:199:PRO:HD3	1.47	1.42
1:A:198:ILE:HG22	1:A:199:PRO:CD	1.53	1.38
1:E:192:ARG:O	1:E:193:SCH:HE3	1.54	1.08
1:B:192:ARG:O	1:B:193:SCH:HE3	1.54	1.08
1:A:192:ARG:O	1:A:193:SCH:HE3	1.54	1.08
1:C:192:ARG:O	1:C:193:SCH:HE3	1.54	1.07
1:D:192:ARG:O	1:D:193:SCH:HE3	1.54	1.07
1:C:194:TYR:O	1:C:196:SER:N	1.99	0.95
1:A:198:ILE:CB	1:A:199:PRO:CD	2.48	0.89
1:C:197:TYR:HH	1:C:255:THR:HG1	0.90	0.87
1:E:198:ILE:CG2	1:E:199:PRO:CD	2.26	0.85
1:C:198:ILE:CG2	1:C:199:PRO:CD	2.14	0.84
1:A:194:TYR:CE1	1:A:196:SER:HB3	2.12	0.84
1:C:198:ILE:CB	1:C:199:PRO:HD3	2.08	0.83
1:E:196:SER:O	1:E:200:ASN:HB2	1.78	0.82
1:B:196:SER:O	1:B:200:ASN:HB2	1.80	0.81
1:C:198:ILE:CB	1:C:199:PRO:CD	2.55	0.81
1:A:196:SER:O	1:A:200:ASN:HB2	1.79	0.81
1:B:198:ILE:CG2	1:B:199:PRO:CD	2.25	0.79
1:D:196:SER:OG	1:D:200:ASN:ND2	2.17	0.78
1:A:198:ILE:CB	1:A:199:PRO:HD2	2.14	0.76
1:A:62:ARG:NH1	1:B:136:ASP:O	2.19	0.76
1:A:198:ILE:CG2	1:A:199:PRO:CD	2.28	0.76
1:B:194:TYR:O	1:B:195:PHE:C	2.29	0.74
1:D:120:PRO:HG3	1:D:197:TYR:CE2	2.22	0.74
1:E:194:TYR:CE1	1:E:196:SER:HB3	2.22	0.74
1:C:194:TYR:O	1:C:197:TYR:N	2.20	0.74
1:A:194:TYR:O	1:A:195:PHE:C	2.31	0.73
1:B:198:ILE:CB	1:B:199:PRO:CD	2.62	0.72
1:A:194:TYR:HE1	1:A:196:SER:HB3	1.54	0.71
1:A:198:ILE:HB	1:A:199:PRO:HD2	1.71	0.71
1:D:28:TYR:CG	1:E:111:LEU:HD11	2.23	0.71
1:D:122:ASP:OD1	1:D:192:ARG:NH1	2.23	0.71
1:D:208:ILE:HG22	1:E:232:LEU:HD21	1.73	0.71
1:E:194:TYR:O	1:E:196:SER:N	2.25	0.70
1:C:120:PRO:HG3	1:C:197:TYR:CD2	2.27	0.69
1:E:194:TYR:O	1:E:195:PHE:C	2.31	0.69
1:E:198:ILE:CB	1:E:199:PRO:CD	2.62	0.68
1:A:122:ASP:OD1	1:A:192:ARG:NH1	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:208:ILE:HG22	1:D:232:LEU:HD21	1.75	0.67
1:D:194:TYR:O	1:D:195:PHE:C	2.38	0.67
1:E:196:SER:O	1:E:200:ASN:N	2.28	0.66
1:E:198:ILE:HG22	1:E:199:PRO:HD2	1.69	0.65
1:D:194:TYR:O	1:D:196:SER:N	2.30	0.64
1:C:11:ILE:HG23	1:C:50:ARG:HH22	1.63	0.63
1:C:120:PRO:HG3	1:C:197:TYR:CE2	2.34	0.63
1:C:122:ASP:OD1	1:C:192:ARG:NH1	2.31	0.63
1:E:120:PRO:HG3	1:E:197:TYR:CD2	2.32	0.63
1:E:198:ILE:CB	1:E:199:PRO:HD2	2.28	0.63
1:E:11:ILE:HG23	1:E:50:ARG:HH22	1.63	0.63
1:D:11:ILE:HG23	1:D:50:ARG:HH22	1.64	0.63
1:A:208:ILE:HG22	1:B:232:LEU:HD21	1.80	0.62
1:C:120:PRO:CG	1:C:197:TYR:CD2	2.83	0.62
1:A:194:TYR:O	1:A:196:SER:N	2.33	0.62
1:A:196:SER:O	1:A:200:ASN:N	2.34	0.61
1:B:208:ILE:HG22	1:C:232:LEU:HD21	1.82	0.61
1:B:198:ILE:HG22	1:B:199:PRO:HD2	1.69	0.61
1:D:120:PRO:CG	1:D:197:TYR:CE2	2.84	0.61
1:B:194:TYR:O	1:B:196:SER:N	2.33	0.61
1:E:120:PRO:HG3	1:E:197:TYR:CE2	2.35	0.61
1:E:198:ILE:HG22	1:E:199:PRO:HD3	0.68	0.61
1:B:28:TYR:CG	1:C:111:LEU:HD11	2.36	0.60
1:E:198:ILE:CG2	1:E:199:PRO:HD2	2.30	0.60
1:B:198:ILE:CB	1:B:199:PRO:HD2	2.31	0.60
1:C:194:TYR:C	1:C:196:SER:N	2.60	0.60
1:E:197:TYR:OH	1:E:255:THR:OG1	1.66	0.59
1:B:196:SER:O	1:B:200:ASN:CB	2.51	0.59
1:B:198:ILE:HG22	1:B:199:PRO:HD3	0.65	0.58
1:E:196:SER:O	1:E:200:ASN:CB	2.52	0.58
1:B:284:GLN:N	1:B:285:PRO:HD3	2.18	0.58
1:C:194:TYR:O	1:C:195:PHE:C	2.47	0.58
1:A:284:GLN:N	1:A:285:PRO:HD3	2.18	0.58
1:D:120:PRO:HG3	1:D:197:TYR:CD2	2.39	0.57
1:A:197:TYR:HH	1:A:255:THR:HG1	0.59	0.57
1:C:284:GLN:N	1:C:285:PRO:HD3	2.18	0.57
1:D:284:GLN:N	1:D:285:PRO:HD3	2.20	0.57
1:E:284:GLN:N	1:E:285:PRO:HD3	2.18	0.57
1:A:232:LEU:HD21	1:E:208:ILE:HG22	1.85	0.57
1:B:196:SER:O	1:B:200:ASN:N	2.38	0.56
1:E:122:ASP:OD1	1:E:192:ARG:NH1	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:SER:OG	1:D:200:ASN:CG	2.47	0.56
1:A:199:PRO:HB2	1:B:252:MET:CE	2.37	0.55
1:E:24:LEU:HD22	1:E:39:VAL:HG21	1.88	0.55
1:A:196:SER:O	1:A:200:ASN:CB	2.52	0.55
1:C:120:PRO:CG	1:C:197:TYR:CE2	2.89	0.55
1:A:195:PHE:CD2	1:A:195:PHE:N	2.73	0.54
1:C:194:TYR:C	1:C:196:SER:H	2.15	0.54
1:A:120:PRO:HG3	1:A:197:TYR:CD2	2.43	0.54
1:C:24:LEU:HD22	1:C:39:VAL:HG21	1.89	0.54
1:E:120:PRO:CG	1:E:197:TYR:CE2	2.91	0.54
1:D:197:TYR:OH	1:D:255:THR:OG1	2.10	0.54
1:A:120:PRO:HG3	1:A:197:TYR:CE2	2.44	0.53
1:B:31:ASP:OD1	1:B:33:LYS:HG3	2.08	0.53
1:B:198:ILE:CG2	1:B:199:PRO:HD2	2.31	0.53
1:C:198:ILE:CB	1:C:199:PRO:HD2	2.36	0.53
1:D:31:ASP:OD1	1:D:33:LYS:HG3	2.09	0.52
1:A:252:MET:CE	1:E:199:PRO:HB2	2.39	0.52
1:D:268:VAL:HG12	1:D:299:PHE:CZ	2.45	0.52
1:C:31:ASP:OD1	1:C:33:LYS:HG3	2.09	0.52
1:B:268:VAL:HG12	1:B:299:PHE:CZ	2.44	0.52
1:A:268:VAL:HG12	1:A:299:PHE:CZ	2.44	0.52
1:E:268:VAL:HG12	1:E:299:PHE:CZ	2.45	0.52
1:A:31:ASP:OD1	1:A:33:LYS:HG3	2.09	0.52
1:E:31:ASP:OD1	1:E:33:LYS:HG3	2.09	0.52
1:A:198:ILE:HG22	1:A:199:PRO:HD3	0.64	0.52
1:C:268:VAL:HG12	1:C:299:PHE:CZ	2.45	0.51
1:A:197:TYR:OH	1:A:255:THR:OG1	1.73	0.50
1:E:120:PRO:CG	1:E:197:TYR:CD2	2.94	0.50
1:B:194:TYR:CE1	1:B:196:SER:HB3	2.46	0.50
1:B:120:PRO:HG3	1:B:197:TYR:CD2	2.47	0.49
1:B:40:ASN:ND2	1:C:81:VAL:O	2.46	0.49
1:A:24:LEU:HD22	1:A:39:VAL:HG21	1.95	0.48
1:D:120:PRO:HG3	1:D:197:TYR:CZ	2.47	0.48
1:A:36:THR:HA	1:A:110:VAL:O	2.13	0.48
1:B:36:THR:HA	1:B:110:VAL:O	2.13	0.48
1:C:36:THR:HA	1:C:110:VAL:O	2.14	0.48
1:E:36:THR:HA	1:E:110:VAL:O	2.14	0.48
1:C:120:PRO:CD	1:C:197:TYR:CE2	2.97	0.48
1:D:197:TYR:HB3	1:D:198:ILE:HD12	1.95	0.48
1:C:306:ALA:HA	1:C:309:ILE:HD12	1.96	0.47
1:A:28:TYR:CG	1:B:111:LEU:HD11	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:24:LEU:HD22	1:E:39:VAL:CG2	2.44	0.47
1:A:306:ALA:HA	1:A:309:ILE:HD12	1.96	0.47
1:B:152:ASN:ND2	1:B:154:ASP:HB2	2.30	0.47
1:B:306:ALA:HA	1:B:309:ILE:HD12	1.96	0.47
1:C:198:ILE:HG22	1:C:199:PRO:CG	2.28	0.47
1:D:306:ALA:HA	1:D:309:ILE:HD12	1.96	0.47
1:D:36:THR:HA	1:D:110:VAL:O	2.14	0.47
1:C:24:LEU:HD22	1:C:39:VAL:CG2	2.45	0.47
1:D:24:LEU:HD22	1:D:39:VAL:HG21	1.96	0.46
1:E:306:ALA:HA	1:E:309:ILE:HD12	1.96	0.46
1:B:24:LEU:HD22	1:B:39:VAL:HG21	1.98	0.46
1:D:152:ASN:ND2	1:D:154:ASP:HB2	2.30	0.46
1:D:120:PRO:CG	1:D:197:TYR:CD2	2.99	0.46
1:A:152:ASN:ND2	1:A:154:ASP:HB2	2.31	0.45
1:E:62:ARG:O	1:E:62:ARG:HD2	2.17	0.45
1:E:120:PRO:CD	1:E:197:TYR:CE2	3.00	0.45
1:C:120:PRO:HG3	1:C:197:TYR:CG	2.51	0.45
1:A:156:PHE:CD2	1:A:156:PHE:C	2.95	0.45
1:C:156:PHE:C	1:C:156:PHE:CD2	2.95	0.44
1:D:24:LEU:HD22	1:D:39:VAL:CG2	2.47	0.44
1:E:152:ASN:ND2	1:E:154:ASP:HB2	2.32	0.44
1:D:156:PHE:CD2	1:D:156:PHE:C	2.95	0.44
1:D:194:TYR:O	1:D:197:TYR:N	2.32	0.44
1:D:40:ASN:ND2	1:E:81:VAL:O	2.50	0.44
1:E:197:TYR:OH	1:E:255:THR:CB	2.62	0.44
1:C:198:ILE:HB	1:C:199:PRO:HD2	2.00	0.44
1:E:194:TYR:HE1	1:E:196:SER:HB3	1.74	0.44
1:A:11:ILE:HG23	1:A:50:ARG:NH2	2.33	0.44
1:C:152:ASN:ND2	1:C:154:ASP:HB2	2.32	0.44
1:A:24:LEU:HD22	1:A:39:VAL:CG2	2.48	0.44
1:B:166:THR:HG22	1:B:167:ALA:H	1.83	0.44
1:B:156:PHE:C	1:B:156:PHE:CD2	2.95	0.44
1:D:166:THR:HG22	1:D:167:ALA:H	1.83	0.43
1:E:156:PHE:CD2	1:E:156:PHE:C	2.95	0.43
1:C:283:SER:C	1:C:285:PRO:HD3	2.44	0.43
1:B:24:LEU:HD22	1:B:39:VAL:CG2	2.49	0.43
1:A:166:THR:HG22	1:A:167:ALA:H	1.84	0.43
1:B:283:SER:C	1:B:285:PRO:HD3	2.44	0.43
1:D:283:SER:C	1:D:285:PRO:HD3	2.44	0.43
1:E:283:SER:C	1:E:285:PRO:HD3	2.44	0.43
1:B:120:PRO:HG3	1:B:197:TYR:CE2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:ASN:HB3	1:D:155:VAL:HG23	2.01	0.42
1:E:152:ASN:HB3	1:E:155:VAL:HG23	2.01	0.42
1:A:283:SER:C	1:A:285:PRO:HD3	2.44	0.42
1:D:162:ILE:HG23	1:D:188:LEU:HD11	2.02	0.42
1:E:198:ILE:HB	1:E:199:PRO:HD2	2.00	0.42
1:D:160:TRP:CZ2	1:D:192:ARG:NH2	2.88	0.41
1:C:152:ASN:HB3	1:C:155:VAL:HG23	2.01	0.41
1:E:10:PRO:HB2	1:E:12:ALA:O	2.20	0.41
1:B:6:SER:O	1:B:51:ARG:NH2	2.53	0.41
1:C:10:PRO:HB2	1:C:12:ALA:O	2.20	0.41
1:D:10:PRO:HB2	1:D:12:ALA:O	2.20	0.41
1:E:6:SER:O	1:E:51:ARG:NH2	2.54	0.41
1:A:284:GLN:N	1:A:285:PRO:CD	2.84	0.41
1:D:196:SER:OG	1:D:196:SER:O	2.35	0.41
1:D:6:SER:O	1:D:51:ARG:NH2	2.53	0.41
1:B:11:ILE:HG12	1:B:50:ARG:NH2	2.36	0.41
1:A:120:PRO:CG	1:A:197:TYR:CE2	3.03	0.41
1:C:6:SER:O	1:C:51:ARG:NH2	2.54	0.41
1:C:120:PRO:HD3	1:C:197:TYR:CE2	2.55	0.41
1:A:6:SER:O	1:A:51:ARG:NH2	2.53	0.40
1:B:162:ILE:HG23	1:B:188:LEU:HD11	2.04	0.40
1:D:162:ILE:HA	1:D:190:ILE:HG22	2.03	0.40
1:E:268:VAL:HG12	1:E:299:PHE:HZ	1.86	0.40
1:A:152:ASN:HB3	1:A:155:VAL:HG23	2.03	0.40
1:C:232:LEU:HD13	1:C:266:TYR:HB3	2.04	0.40
1:D:268:VAL:HG12	1:D:299:PHE:HZ	1.85	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	308/311 (99%)	300 (97%)	8 (3%)	0	100	100
1	B	308/311 (99%)	300 (97%)	8 (3%)	0	100	100
1	C	308/311 (99%)	298 (97%)	9 (3%)	1 (0%)	36	67
1	D	308/311 (99%)	300 (97%)	8 (3%)	0	100	100
1	E	308/311 (99%)	299 (97%)	9 (3%)	0	100	100
All	All	1540/1555 (99%)	1497 (97%)	42 (3%)	1 (0%)	48	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	195	PHE

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	279/279 (100%)	254 (91%)	25 (9%)	9	32
1	B	279/279 (100%)	255 (91%)	24 (9%)	10	33
1	C	279/279 (100%)	253 (91%)	26 (9%)	8	31
1	D	279/279 (100%)	257 (92%)	22 (8%)	11	36
1	E	279/279 (100%)	255 (91%)	24 (9%)	10	33
All	All	1395/1395 (100%)	1274 (91%)	121 (9%)	9	33

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	11	ILE
1	A	62	ARG
1	A	88	ASP
1	A	111	LEU
1	A	131	ILE
1	A	135	VAL

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Mol	Chain	Res	Type
1	A	136	ASP
1	A	147	GLU
1	A	166	THR
1	A	168	VAL
1	A	170	LYS
1	A	178	ASP
1	A	180	LEU
1	A	192	ARG
1	A	195	PHE
1	A	198	ILE
1	A	212	SER
1	A	239	ASN
1	A	275	VAL
1	A	280	LYS
1	A	281	VAL
1	A	297	ILE
1	A	307	ASN
1	A	315	PHE
1	B	5	VAL
1	B	11	ILE
1	B	62	ARG
1	B	88	ASP
1	B	109	ARG
1	B	111	LEU
1	B	135	VAL
1	B	136	ASP
1	B	147	GLU
1	B	166	THR
1	B	168	VAL
1	B	170	LYS
1	B	178	ASP
1	B	180	LEU
1	B	192	ARG
1	B	195	PHE
1	B	198	ILE
1	B	199	PRO
1	B	212	SER
1	B	239	ASN
1	B	275	VAL
1	B	281	VAL
1	B	307	ASN
1	B	315	PHE

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Mol	Chain	Res	Type
1	C	5	VAL
1	C	11	ILE
1	C	58	ARG
1	C	62	ARG
1	C	88	ASP
1	C	111	LEU
1	C	135	VAL
1	C	136	ASP
1	C	147	GLU
1	C	166	THR
1	C	168	VAL
1	C	170	LYS
1	C	178	ASP
1	C	180	LEU
1	C	192	ARG
1	C	198	ILE
1	C	199	PRO
1	C	202	ILE
1	C	212	SER
1	C	239	ASN
1	C	275	VAL
1	C	280	LYS
1	C	281	VAL
1	C	284	GLN
1	C	307	ASN
1	C	315	PHE
1	D	5	VAL
1	D	11	ILE
1	D	62	ARG
1	D	88	ASP
1	D	111	LEU
1	D	131	ILE
1	D	135	VAL
1	D	136	ASP
1	D	147	GLU
1	D	166	THR
1	D	168	VAL
1	D	170	LYS
1	D	178	ASP
1	D	180	LEU
1	D	192	ARG
1	D	195	PHE

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Mol	Chain	Res	Type
1	D	212	SER
1	D	239	ASN
1	D	275	VAL
1	D	281	VAL
1	D	307	ASN
1	D	315	PHE
1	E	5	VAL
1	E	11	ILE
1	E	58	ARG
1	E	62	ARG
1	E	88	ASP
1	E	111	LEU
1	E	131	ILE
1	E	135	VAL
1	E	136	ASP
1	E	147	GLU
1	E	166	THR
1	E	168	VAL
1	E	170	LYS
1	E	178	ASP
1	E	180	LEU
1	E	192	ARG
1	E	198	ILE
1	E	212	SER
1	E	239	ASN
1	E	275	VAL
1	E	281	VAL
1	E	297	ILE
1	E	307	ASN
1	E	315	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	277	HIS
1	B	152	ASN
1	B	200	ASN
1	B	277	HIS
1	C	152	ASN
1	C	277	HIS
1	D	152	ASN

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Mol	Chain	Res	Type
1	D	200	ASN
1	D	277	HIS
1	E	152	ASN
1	E	277	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

5 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	SCH	B	193	1	6,7,8	1.40	2 (33%)	4,7,9	1.50	0
1	SCH	A	193	1	6,7,8	1.41	2 (33%)	4,7,9	1.50	0
1	SCH	C	193	1	6,7,8	1.40	2 (33%)	4,7,9	1.49	0
1	SCH	D	193	1	6,7,8	1.41	2 (33%)	4,7,9	1.49	0
1	SCH	E	193	1	6,7,8	1.40	2 (33%)	4,7,9	1.50	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
1	SCH	B	193	1	-	1/2/6/8	-
1	SCH	A	193	1	-	1/2/6/8	-
1	SCH	C	193	1	-	1/2/6/8	-
1	SCH	D	193	1	-	1/2/6/8	-
1	SCH	E	193	1	-	1/2/6/8	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	193	SCH	O-C	2.56	1.29	1.20
1	A	193	SCH	O-C	2.55	1.29	1.20
1	C	193	SCH	O-C	2.55	1.29	1.20
1	B	193	SCH	O-C	2.55	1.29	1.20
1	E	193	SCH	O-C	2.54	1.29	1.20
1	A	193	SCH	CB-SG	-2.17	1.74	1.81
1	B	193	SCH	CB-SG	-2.15	1.74	1.81
1	E	193	SCH	CB-SG	-2.15	1.74	1.81
1	D	193	SCH	CB-SG	-2.15	1.74	1.81
1	C	193	SCH	CB-SG	-2.14	1.74	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	193	SCH	N-CA-CB-SG
1	B	193	SCH	N-CA-CB-SG
1	C	193	SCH	N-CA-CB-SG
1	D	193	SCH	N-CA-CB-SG
1	E	193	SCH	N-CA-CB-SG

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	193	SCH	1	0
1	A	193	SCH	1	0
1	C	193	SCH	1	0
1	D	193	SCH	1	0
1	E	193	SCH	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	LMT	A	401	-	11,11,36	1.15	0	10,10,47	0.92	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
2	LMT	A	401	-	-	2/9/9/61	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

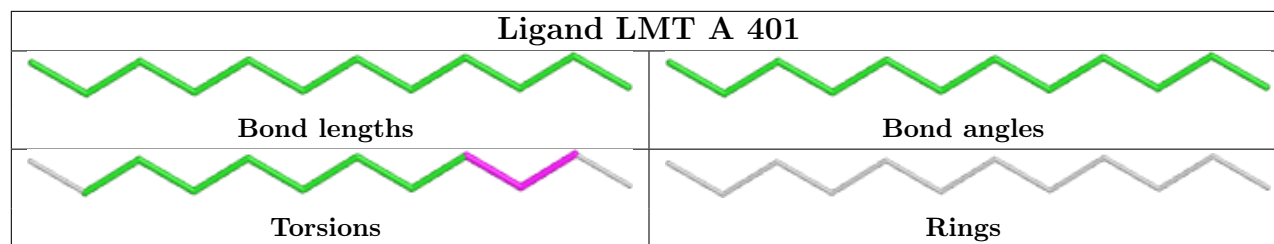
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	LMT	C2-C3-C4-C5
2	A	401	LMT	C1-C2-C3-C4

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	189:ARG	C	190:ILE	N	1.18

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	310/311 (99%)	-0.81	0 100 100	44, 89, 129, 174	0
1	B	310/311 (99%)	-0.88	0 100 100	45, 80, 113, 136	0
1	C	310/311 (99%)	-0.84	0 100 100	45, 81, 111, 133	0
1	D	310/311 (99%)	-0.87	0 100 100	47, 87, 119, 150	0
1	E	310/311 (99%)	-0.85	0 100 100	50, 85, 117, 144	0
All	All	1550/1555 (99%)	-0.85	0 100 100	44, 85, 118, 174	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SCH	D	193	8/9	0.77	0.20	84,91,126,150	0
1	SCH	A	193	8/9	0.80	0.21	84,91,126,150	0
1	SCH	E	193	8/9	0.89	0.14	84,91,126,150	0
1	SCH	C	193	8/9	0.92	0.12	84,91,126,150	0
1	SCH	B	193	8/9	0.96	0.09	84,91,126,150	0

6.3 Carbohydrates [i](#)

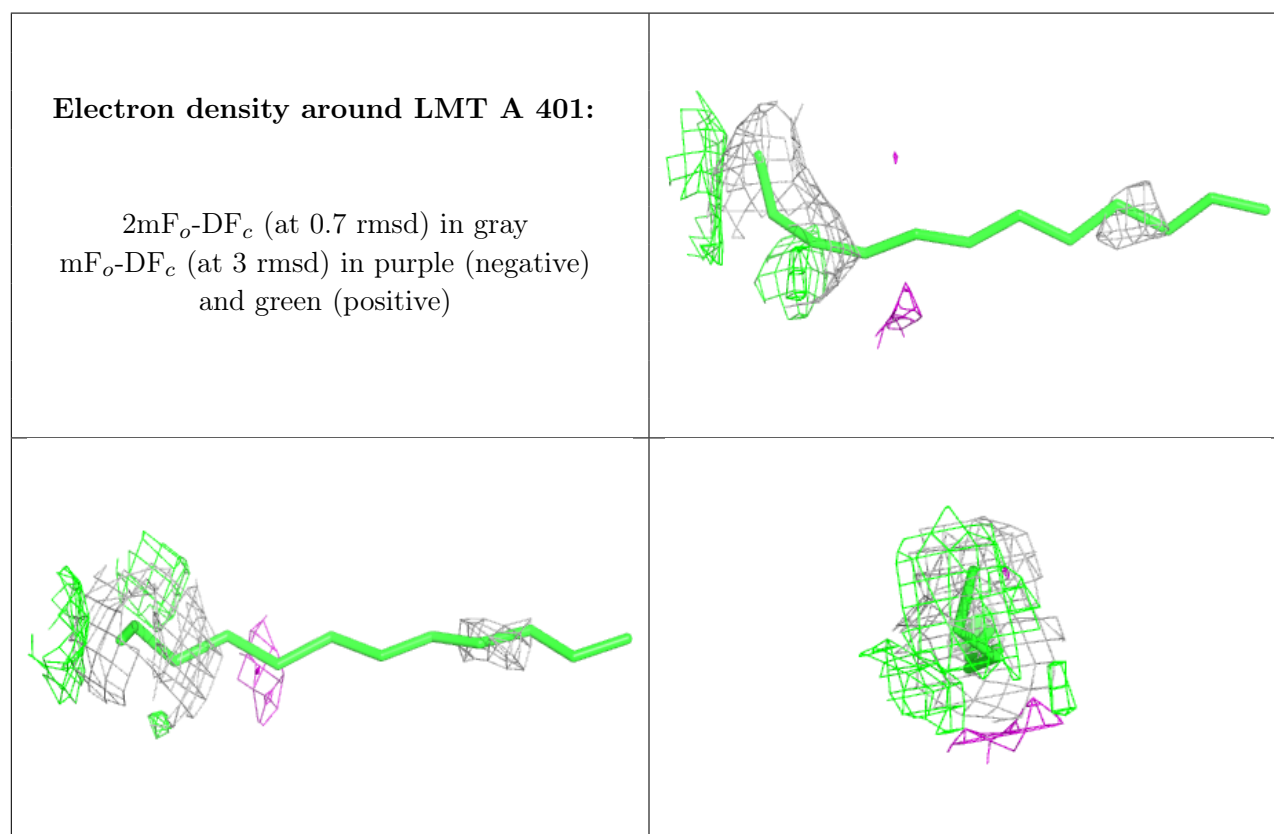
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	LMT	A	401	12/35	0.89	0.22	44,61,72,75	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.