



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:24 PM UTC

PDB ID : 4KNF / pdb_00004knf
Title : Crystal structure of a blue-light absorbing proteorhodopsin double-mutant D97N/Q105L from HOT75
Authors : Ran, T.; Ozorowski, G.; Gao, Y.; Wang, W.; Luecke, H.
Deposited on : 2013-05-09
Resolution : 2.60 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

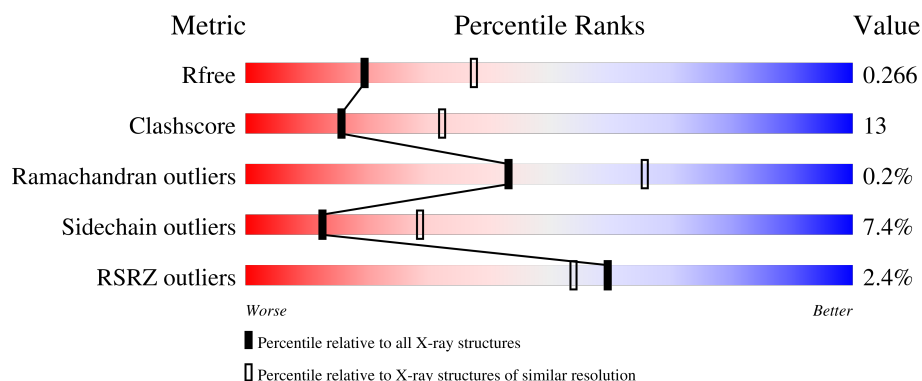
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	261	 2% 67% 20% 11%
1	B	261	 3% 61% 22% 14%
1	C	261	 % 74% 14% 11%
1	D	261	 2% 71% 14% 12%
1	E	261	 2% 69% 18% 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	RET	A	301	-	-	X	-
2	RET	C	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8858 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 Blue-light absorbing proteorhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1757	1172	268	306	11			
1	B	224	Total	C	N	O	S	0	0	0
			1727	1156	263	298	10			
1	C	232	Total	C	N	O	S	0	0	0
			1765	1178	269	307	11			
1	D	229	Total	C	N	O	S	0	0	0
			1745	1166	266	302	11			
1	E	230	Total	C	N	O	S	0	0	0
			1753	1170	267	305	11			

There are 60 discrepancies between the modelled and reference sequences:

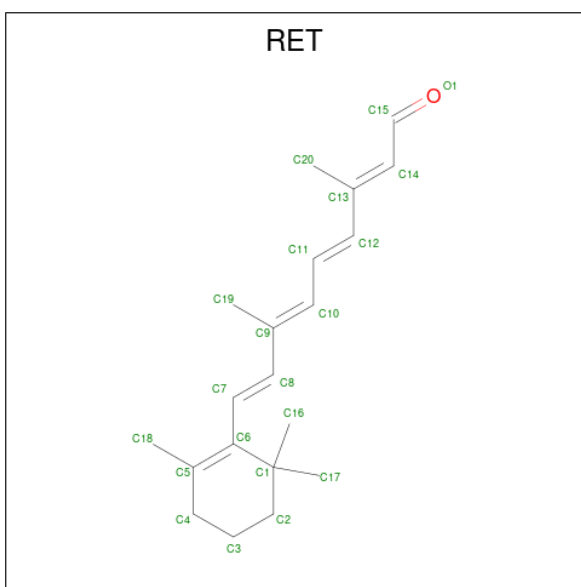
Chain	Residue	Modelled	Actual	Comment	Reference
A	97	ASN	ASP	engineered mutation	UNP Q9AFF7
A	105	LEU	GLN	engineered mutation	UNP Q9AFF7
A	251	LYS	-	expression tag	UNP Q9AFF7
A	252	LEU	-	expression tag	UNP Q9AFF7
A	253	LEU	-	expression tag	UNP Q9AFF7
A	254	GLU	-	expression tag	UNP Q9AFF7
A	255	HIS	-	expression tag	UNP Q9AFF7
A	256	HIS	-	expression tag	UNP Q9AFF7
A	257	HIS	-	expression tag	UNP Q9AFF7
A	258	HIS	-	expression tag	UNP Q9AFF7
A	259	HIS	-	expression tag	UNP Q9AFF7
A	260	HIS	-	expression tag	UNP Q9AFF7
B	97	ASN	ASP	engineered mutation	UNP Q9AFF7
B	105	LEU	GLN	engineered mutation	UNP Q9AFF7
B	251	LYS	-	expression tag	UNP Q9AFF7
B	252	LEU	-	expression tag	UNP Q9AFF7
B	253	LEU	-	expression tag	UNP Q9AFF7
B	254	GLU	-	expression tag	UNP Q9AFF7
B	255	HIS	-	expression tag	UNP Q9AFF7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	256	HIS	-	expression tag	UNP Q9AFF7
B	257	HIS	-	expression tag	UNP Q9AFF7
B	258	HIS	-	expression tag	UNP Q9AFF7
B	259	HIS	-	expression tag	UNP Q9AFF7
B	260	HIS	-	expression tag	UNP Q9AFF7
C	97	ASN	ASP	engineered mutation	UNP Q9AFF7
C	105	LEU	GLN	engineered mutation	UNP Q9AFF7
C	251	LYS	-	expression tag	UNP Q9AFF7
C	252	LEU	-	expression tag	UNP Q9AFF7
C	253	LEU	-	expression tag	UNP Q9AFF7
C	254	GLU	-	expression tag	UNP Q9AFF7
C	255	HIS	-	expression tag	UNP Q9AFF7
C	256	HIS	-	expression tag	UNP Q9AFF7
C	257	HIS	-	expression tag	UNP Q9AFF7
C	258	HIS	-	expression tag	UNP Q9AFF7
C	259	HIS	-	expression tag	UNP Q9AFF7
C	260	HIS	-	expression tag	UNP Q9AFF7
D	97	ASN	ASP	engineered mutation	UNP Q9AFF7
D	105	LEU	GLN	engineered mutation	UNP Q9AFF7
D	251	LYS	-	expression tag	UNP Q9AFF7
D	252	LEU	-	expression tag	UNP Q9AFF7
D	253	LEU	-	expression tag	UNP Q9AFF7
D	254	GLU	-	expression tag	UNP Q9AFF7
D	255	HIS	-	expression tag	UNP Q9AFF7
D	256	HIS	-	expression tag	UNP Q9AFF7
D	257	HIS	-	expression tag	UNP Q9AFF7
D	258	HIS	-	expression tag	UNP Q9AFF7
D	259	HIS	-	expression tag	UNP Q9AFF7
D	260	HIS	-	expression tag	UNP Q9AFF7
E	97	ASN	ASP	engineered mutation	UNP Q9AFF7
E	105	LEU	GLN	engineered mutation	UNP Q9AFF7
E	251	LYS	-	expression tag	UNP Q9AFF7
E	252	LEU	-	expression tag	UNP Q9AFF7
E	253	LEU	-	expression tag	UNP Q9AFF7
E	254	GLU	-	expression tag	UNP Q9AFF7
E	255	HIS	-	expression tag	UNP Q9AFF7
E	256	HIS	-	expression tag	UNP Q9AFF7
E	257	HIS	-	expression tag	UNP Q9AFF7
E	258	HIS	-	expression tag	UNP Q9AFF7
E	259	HIS	-	expression tag	UNP Q9AFF7
E	260	HIS	-	expression tag	UNP Q9AFF7

- Molecule 2 is RETINAL (CCD ID: RET) (formula: C₂₀H₂₈O).

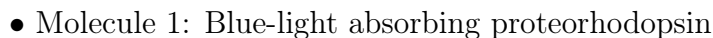
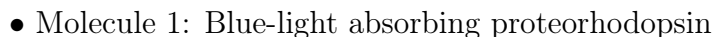


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C 20 20	0	0
2	B	1	Total C 20 20	0	0
2	C	1	Total C 20 20	0	0
2	D	1	Total C 20 20	0	0
2	E	1	Total C 20 20	0	0

- Molecule 3 is water.

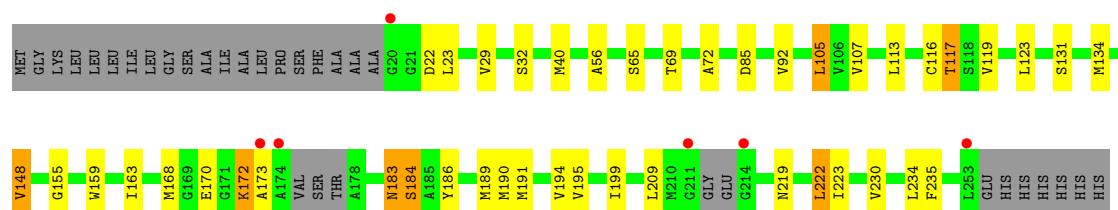
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total O 2 2	0	0
3	B	3	Total O 3 3	0	0
3	C	2	Total O 2 2	0	0
3	D	2	Total O 2 2	0	0
3	E	2	Total O 2 2	0	0

- Molecule 1: Blue-light absorbing proteorhodopsin



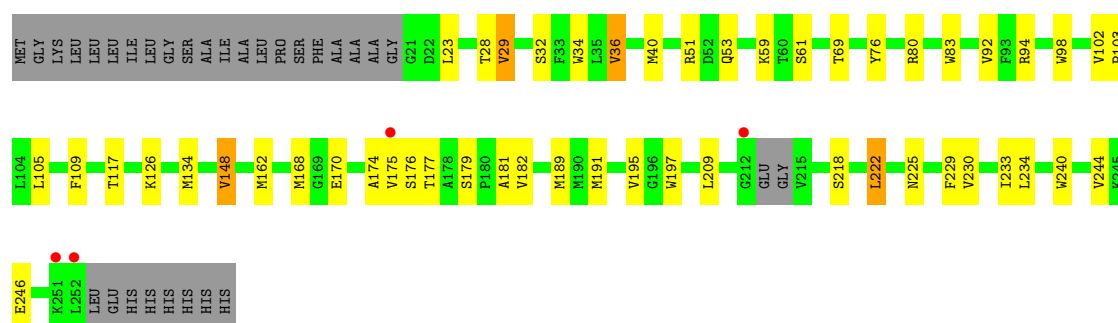
- Molecule 1: Blue-light absorbing proteorhodopsin

Chain D: 



- Molecule 1: Blue-light absorbing proteorhodopsin

Chain E: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	161.40Å 168.90Å 65.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.95 – 2.60 19.95 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.95-2.60) 98.7 (19.95-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.219 , 0.267 0.218 , 0.266	Depositor DCC
R_{free} test set	2842 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	64.9	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8858	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: RET

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	0.66	0/1803	0.99	1/2460 (0.0%)
1	B	0.62	0/1772	0.96	2/2414 (0.1%)
1	C	0.65	0/1811	1.02	1/2471 (0.0%)
1	D	0.65	0/1790	0.99	2/2440 (0.1%)
1	E	0.67	0/1799	1.02	3/2455 (0.1%)
All	All	0.65	0/8975	1.00	9/12240 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	148	VAL	N-CA-C	6.48	116.64	110.42
1	B	71	ILE	N-CA-C	-6.29	104.37	110.72
1	C	178	ALA	N-CA-C	6.22	118.58	111.11
1	E	148	VAL	N-CA-C	6.13	118.72	111.05
1	E	69	THR	CA-C-N	5.39	125.92	119.94
1	E	69	THR	C-N-CA	5.39	125.92	119.94
1	D	116	CYS	N-CA-C	5.32	118.56	111.75
1	A	120	ALA	N-CA-C	5.24	117.67	111.33
1	B	231	ASN	N-CA-C	5.23	117.06	111.36

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	1757	0	1778	48	0
1	B	1727	0	1744	61	0
1	C	1765	0	1789	38	0
1	D	1745	0	1767	34	0
1	E	1753	0	1775	43	0
2	A	20	0	27	22	0
2	B	20	0	27	8	0
2	C	20	0	27	9	0
2	D	20	0	27	5	0
2	E	20	0	27	6	0
3	A	2	0	0	0	0
3	B	3	0	0	2	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
All	All	8858	0	8988	237	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 13.

All (237) 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:40:MET:HE2	1:D:72:ALA:CB	1.62	1.29
1:C:40:MET:HE2	1:C:72:ALA:HB3	1.26	1.11
1:A:134:MET:CE	2:A:301:RET:H192	1.82	1.10
1:D:40:MET:HE2	1:D:72:ALA:HB3	1.21	1.09
1:C:79:MET:HE1	1:C:94:ARG:HD3	1.27	1.08
1:A:134:MET:HE2	2:A:301:RET:H7	1.10	1.06
1:D:191:MET:O	1:D:195:VAL:HG12	1.57	1.04
1:E:23:LEU:HD11	1:E:80:ARG:NH2	1.73	1.03
1:B:79:MET:HE1	1:B:94:ARG:HD3	1.41	1.02
1:D:40:MET:CE	1:D:72:ALA:HB3	1.91	1.01
1:A:134:MET:HE2	2:A:301:RET:C7	1.94	0.97
1:C:79:MET:HE1	1:C:94:ARG:CD	1.97	0.95
2:C:301:RET:H171	2:C:301:RET:H8	1.48	0.92
1:B:134:MET:SD	2:B:301:RET:H192	2.10	0.91
1:A:134:MET:CE	2:A:301:RET:H7	2.00	0.89
1:A:134:MET:CE	2:A:301:RET:C19	2.51	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:RET:H161	2:B:301:RET:H8	1.54	0.87
2:D:301:RET:H171	2:D:301:RET:H8	1.58	0.86
1:A:134:MET:HE1	2:A:301:RET:H192	1.59	0.85
1:D:219:ASN:O	1:D:223:ILE:HG12	1.77	0.84
1:B:94:ARG:NH2	3:B:406:HOH:O	1.98	0.84
1:C:40:MET:CE	1:C:69:THR:HA	2.07	0.84
1:A:149:LEU:O	1:A:153:ILE:HD12	1.78	0.84
1:C:40:MET:HE2	1:C:72:ALA:CB	2.09	0.82
1:D:40:MET:HE2	1:D:72:ALA:HB1	1.62	0.81
1:A:40:MET:HE1	1:A:69:THR:HG22	1.62	0.80
2:B:301:RET:H171	2:B:301:RET:C8	2.12	0.79
1:B:162:MET:O	1:B:166:LEU:HD12	1.82	0.79
1:A:134:MET:HE1	2:A:301:RET:C19	2.11	0.79
1:C:79:MET:HE3	1:C:94:ARG:CA	2.13	0.79
1:A:134:MET:HE3	2:A:301:RET:H181	1.66	0.78
1:B:79:MET:HE1	1:B:94:ARG:CD	2.14	0.76
1:E:23:LEU:HD11	1:E:80:ARG:HH22	1.51	0.76
1:B:79:MET:CE	1:B:94:ARG:HD3	2.17	0.75
1:E:29:VAL:HG22	1:E:83:TRP:CE2	2.20	0.75
1:A:134:MET:HE2	2:A:301:RET:H192	1.68	0.75
1:B:134:MET:HE2	1:B:159:TRP:HA	1.69	0.74
1:C:79:MET:HE3	1:C:94:ARG:HA	1.70	0.73
1:D:40:MET:CE	1:D:72:ALA:CB	2.53	0.73
1:E:29:VAL:HG22	1:E:83:TRP:NE1	2.03	0.73
1:E:181:ALA:CB	1:E:246:GLU:HG3	2.19	0.73
1:D:134:MET:SD	2:D:301:RET:H192	2.28	0.73
1:D:23:LEU:O	1:E:92:VAL:HG11	1.88	0.72
2:B:301:RET:H8	2:B:301:RET:H171	1.69	0.72
1:A:40:MET:CE	1:A:69:THR:HG22	2.20	0.72
2:B:301:RET:H8	2:B:301:RET:C16	2.20	0.71
2:A:301:RET:H8	2:A:301:RET:H161	1.72	0.71
1:E:175:VAL:HG12	1:E:176:SER:H	1.56	0.71
1:A:134:MET:CE	2:A:301:RET:H181	2.20	0.70
1:D:117:THR:HG22	1:D:173:ALA:HB3	1.72	0.70
1:A:134:MET:HE2	2:A:301:RET:C19	2.20	0.70
1:D:40:MET:HE3	1:D:69:THR:HA	1.73	0.70
1:B:40:MET:CE	1:B:69:THR:HG22	2.22	0.69
1:B:228:ASP:OD2	3:B:406:HOH:O	2.10	0.68
1:A:174:ALA:HB1	1:A:176:SER:OG	1.94	0.67
2:C:301:RET:H171	2:C:301:RET:C8	2.14	0.67
1:A:134:MET:HE3	1:A:159:TRP:HB2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:192:ILE:HG13	1:C:231:ASN:HD22	1.59	0.66
1:A:193:ILE:O	1:A:197:TRP:HB2	1.96	0.65
1:C:79:MET:CE	1:C:94:ARG:CD	2.75	0.65
1:B:191:MET:HE3	1:B:195:VAL:CG1	2.28	0.64
1:B:79:MET:HE3	1:B:94:ARG:HB2	1.81	0.63
1:D:105:LEU:HD11	1:D:235:PHE:CE2	2.33	0.63
1:B:79:MET:CE	1:B:94:ARG:NE	2.62	0.62
1:B:40:MET:HE3	1:B:40:MET:HA	1.80	0.62
1:D:22:ASP:CG	1:E:92:VAL:HG12	2.24	0.62
1:B:40:MET:HE1	1:B:69:THR:HG22	1.82	0.62
1:D:134:MET:HE2	1:D:159:TRP:HA	1.82	0.62
1:E:134:MET:SD	2:E:301:RET:H192	2.40	0.61
1:C:192:ILE:HG13	1:C:231:ASN:ND2	2.14	0.61
1:D:189:MET:HE3	1:D:190:MET:HE2	1.82	0.61
1:D:22:ASP:OD2	1:E:92:VAL:HG12	2.01	0.61
1:A:40:MET:HE2	1:A:69:THR:HA	1.82	0.61
1:C:79:MET:CE	1:C:94:ARG:HD3	2.17	0.60
1:C:147:PRO:HG2	1:C:150:PRO:HG2	1.83	0.60
1:B:79:MET:CE	1:B:94:ARG:CD	2.77	0.60
1:D:40:MET:CE	1:D:69:THR:HA	2.32	0.60
1:B:79:MET:CE	1:B:94:ARG:CZ	2.79	0.60
1:B:109:PHE:CE1	1:B:189:MET:HB2	2.37	0.60
1:B:98:TRP:CD1	2:B:301:RET:H14	2.37	0.60
1:C:79:MET:CE	1:C:94:ARG:HA	2.31	0.60
1:E:23:LEU:CD1	1:E:80:ARG:HH22	2.15	0.59
1:C:109:PHE:CD1	1:C:189:MET:HG3	2.38	0.59
1:C:40:MET:HE3	1:C:69:THR:HA	1.84	0.59
1:E:36:VAL:CG2	1:E:76:TYR:CE2	2.86	0.58
1:D:117:THR:CG2	1:D:173:ALA:HB3	2.33	0.58
2:D:301:RET:H171	2:D:301:RET:C8	2.22	0.58
1:E:51:ARG:NH1	1:E:59:LYS:HB3	2.19	0.58
1:A:30:GLY:O	1:A:80:ARG:NH1	2.37	0.57
1:D:186:TYR:CZ	1:D:190:MET:HE3	2.39	0.57
1:C:79:MET:HE3	1:C:94:ARG:CB	2.34	0.57
2:A:301:RET:C8	2:A:301:RET:H171	2.33	0.57
1:C:79:MET:HE3	1:C:94:ARG:HB2	1.87	0.57
1:A:219:ASN:O	1:A:223:ILE:HG13	2.05	0.56
1:B:126:LYS:NZ	1:B:165:GLU:OE2	2.36	0.56
1:B:152:PHE:CZ	1:B:156:MET:HE3	2.40	0.56
1:B:36:VAL:CG1	1:B:76:TYR:CE2	2.88	0.56
1:C:78:TYR:O	1:C:82:VAL:HG13	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ALA:HB1	1:B:238:ILE:HG21	1.88	0.56
1:D:170:GLU:C	1:D:172:LYS:H	2.13	0.56
1:A:134:MET:SD	2:A:301:RET:H192	2.46	0.55
1:B:79:MET:HE2	1:B:94:ARG:CZ	2.37	0.55
1:C:79:MET:HE1	1:C:94:ARG:NE	2.22	0.55
1:C:79:MET:CE	1:C:94:ARG:NE	2.70	0.55
1:C:40:MET:HE1	1:C:69:THR:HA	1.87	0.55
1:B:147:PRO:HG2	1:B:150:PRO:HG2	1.88	0.54
1:E:117:THR:HG22	1:E:174:ALA:HB3	1.90	0.54
1:B:149:LEU:HB3	1:B:150:PRO:HD3	1.90	0.54
1:E:179:SER:HB3	1:E:182:VAL:HG22	1.90	0.54
2:B:301:RET:H161	2:B:301:RET:C8	2.31	0.53
1:C:79:MET:CE	1:C:94:ARG:HB2	2.38	0.53
1:A:40:MET:HE3	1:A:40:MET:HA	1.90	0.53
1:B:32:SER:O	1:B:36:VAL:HG12	2.08	0.53
1:C:134:MET:SD	2:C:301:RET:H192	2.48	0.53
1:E:162:MET:HE1	1:E:197:TRP:HZ3	1.72	0.53
1:D:155:GLY:HA3	2:D:301:RET:H31	1.90	0.53
2:E:301:RET:H161	2:E:301:RET:H8	1.91	0.53
1:A:29:VAL:HG22	1:A:222:LEU:HD13	1.91	0.53
1:B:54:VAL:HG21	1:B:62:LEU:HD12	1.89	0.53
1:A:171:GLY:C	1:A:173:ALA:H	2.17	0.53
1:B:36:VAL:HG13	1:B:76:TYR:CE2	2.43	0.53
1:A:147:PRO:HB2	1:A:150:PRO:HD2	1.91	0.51
1:C:53:GLN:HE22	1:D:56:ALA:HA	1.73	0.51
1:E:105:LEU:HD23	1:E:197:TRP:HZ2	1.75	0.51
1:A:134:MET:HE1	2:A:301:RET:H193	1.92	0.51
1:B:40:MET:CE	1:B:40:MET:HA	2.41	0.51
1:D:22:ASP:CG	1:E:92:VAL:CG1	2.83	0.51
1:A:98:TRP:CD1	2:A:301:RET:H14	2.46	0.51
2:A:301:RET:H161	2:A:301:RET:C8	2.40	0.51
1:E:53:GLN:HA	1:E:53:GLN:NE2	2.25	0.51
1:A:35:LEU:HD21	1:B:96:ILE:HD12	1.92	0.50
2:A:301:RET:H8	2:A:301:RET:H171	1.92	0.50
1:E:29:VAL:CG2	1:E:83:TRP:CE2	2.90	0.50
1:C:171:GLY:C	1:C:173:ALA:H	2.20	0.50
1:C:40:MET:HE3	1:C:69:THR:CA	2.42	0.49
1:E:36:VAL:O	1:E:40:MET:HG2	2.13	0.49
2:C:301:RET:C8	2:C:301:RET:C17	2.89	0.49
1:A:186:TYR:O	1:A:190:MET:HG2	2.13	0.49
1:A:40:MET:CE	1:A:69:THR:HA	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:PRO:HB2	1:A:150:PRO:CG	2.42	0.49
2:E:301:RET:H8	2:E:301:RET:H171	1.94	0.49
1:A:147:PRO:HB2	1:A:150:PRO:CD	2.43	0.48
1:A:117:THR:HG22	1:A:174:ALA:HB3	1.96	0.48
1:B:134:MET:HE2	1:B:159:TRP:CA	2.41	0.48
1:C:79:MET:CE	1:C:94:ARG:CB	2.91	0.48
1:D:113:LEU:HD13	1:D:123:LEU:HD13	1.96	0.48
1:D:29:VAL:HG22	1:D:222:LEU:HD13	1.96	0.48
1:B:185:ALA:O	1:B:189:MET:CG	2.62	0.48
1:C:192:ILE:O	1:C:196:GLY:HA3	2.14	0.48
1:E:102:VAL:HB	1:E:103:PRO:HD3	1.96	0.48
1:A:113:LEU:HD13	1:A:123:LEU:HD13	1.95	0.47
1:B:103:PRO:O	1:B:107:VAL:HG23	2.14	0.47
1:B:36:VAL:HG11	1:B:76:TYR:CE2	2.48	0.47
2:A:301:RET:H8	2:A:301:RET:C16	2.43	0.47
1:E:181:ALA:HB2	1:E:246:GLU:HG3	1.94	0.47
1:B:188:ALA:CB	1:B:238:ILE:HG21	2.44	0.47
1:D:40:MET:HE3	1:D:69:THR:CA	2.43	0.46
1:B:36:VAL:HG13	1:B:76:TYR:CD2	2.50	0.46
1:D:199:ILE:HG22	1:D:223:ILE:HG22	1.97	0.46
1:B:109:PHE:CZ	1:B:189:MET:HB2	2.51	0.46
1:B:189:MET:O	1:B:193:ILE:HG12	2.16	0.46
1:B:191:MET:HE3	1:B:195:VAL:HG12	1.98	0.46
1:D:163:ILE:HD13	1:D:194:VAL:HG22	1.98	0.45
1:E:230:VAL:O	1:E:234:LEU:HB2	2.16	0.45
1:B:198:ALA:O	1:B:201:PRO:HG2	2.17	0.45
1:D:131:SER:HA	1:D:134:MET:HE3	1.98	0.45
1:B:156:MET:HE1	1:B:201:PRO:HB3	1.99	0.45
1:B:219:ASN:O	1:B:223:ILE:HG13	2.16	0.45
1:E:240:TRP:O	1:E:244:VAL:HG13	2.16	0.45
1:D:22:ASP:OD1	1:E:92:VAL:HG13	2.17	0.45
2:E:301:RET:H7	2:E:301:RET:H181	1.58	0.45
1:C:112:ILE:HG21	1:C:189:MET:HE2	1.99	0.44
1:C:200:TYR:N	1:C:201:PRO:HD2	2.32	0.44
1:D:183:ASN:HD22	1:D:184:SER:N	2.14	0.44
1:A:177:THR:HG23	1:A:183:ASN:HB2	2.00	0.44
1:E:23:LEU:CD1	1:E:80:ARG:NH2	2.61	0.44
1:A:52:ASP:C	1:A:52:ASP:OD1	2.59	0.44
1:B:33:PHE:HB3	1:B:80:ARG:HB2	2.00	0.44
1:B:35:LEU:HD21	1:C:96:ILE:HD12	2.00	0.44
2:E:301:RET:H171	2:E:301:RET:C8	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:MET:HE1	1:B:94:ARG:CZ	2.48	0.44
1:A:126:LYS:HG2	1:A:161:TYR:OH	2.17	0.44
1:A:166:LEU:HD12	1:A:193:ILE:HD12	1.99	0.44
1:B:185:ALA:O	1:B:189:MET:HG3	2.17	0.44
1:D:230:VAL:O	1:D:234:LEU:HB2	2.18	0.44
1:C:98:TRP:CD1	2:C:301:RET:H14	2.53	0.43
1:E:36:VAL:HG21	1:E:76:TYR:CE2	2.51	0.43
1:E:109:PHE:CD1	1:E:189:MET:HG3	2.53	0.43
1:B:191:MET:HE3	1:B:195:VAL:HG11	1.97	0.43
1:A:96:ILE:HG21	1:E:34:TRP:CD1	2.53	0.43
1:E:191:MET:SD	1:E:195:VAL:CG2	3.06	0.43
1:C:23:LEU:O	1:D:92:VAL:HG21	2.19	0.43
1:E:181:ALA:HB3	1:E:246:GLU:HG3	1.98	0.43
1:E:218:SER:O	1:E:222:LEU:HD22	2.17	0.42
2:D:301:RET:H7	2:D:301:RET:H181	1.61	0.42
2:A:301:RET:H7	2:A:301:RET:H181	1.81	0.42
1:B:34:TRP:CD1	1:C:96:ILE:HG21	2.53	0.42
1:E:23:LEU:HD11	1:E:80:ARG:HH21	1.76	0.42
1:A:40:MET:HE2	1:A:69:THR:CA	2.47	0.42
1:E:229:PHE:CD2	1:E:233:ILE:HD12	2.54	0.42
1:A:179:SER:HB3	1:A:182:VAL:HB	2.01	0.42
1:B:89:THR:HA	1:B:90:PRO:HD3	1.84	0.42
2:C:301:RET:H11	2:C:301:RET:H191	1.83	0.42
2:A:301:RET:H11	2:A:301:RET:H191	1.76	0.42
1:B:105:LEU:HD22	1:B:235:PHE:CE2	2.55	0.42
1:C:192:ILE:O	1:C:196:GLY:CA	2.68	0.42
1:E:36:VAL:HG11	1:E:225:ASN:HB3	2.02	0.42
1:A:175:VAL:O	1:A:176:SER:C	2.63	0.41
1:A:230:VAL:O	1:A:234:LEU:HB2	2.20	0.41
1:B:79:MET:HE3	1:B:94:ARG:CB	2.50	0.41
1:E:94:ARG:HD2	1:E:98:TRP:CZ2	2.55	0.41
1:E:175:VAL:HG12	1:E:176:SER:N	2.30	0.41
1:A:55:SER:OG	1:A:57:LYS:HB2	2.21	0.41
1:E:134:MET:CE	2:E:301:RET:H192	2.50	0.41
1:E:32:SER:O	1:E:36:VAL:HG13	2.21	0.41
1:A:200:TYR:HB3	2:A:301:RET:H161	2.02	0.41
1:B:175:VAL:O	1:B:176:SER:C	2.63	0.41
1:B:185:ALA:O	1:B:189:MET:HG2	2.21	0.41
1:D:183:ASN:HD22	1:D:184:SER:H	1.68	0.41
1:B:187:ASN:O	1:B:190:MET:HB2	2.20	0.41
1:C:134:MET:CE	2:C:301:RET:H192	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:VAL:O	1:A:37:THR:C	2.64	0.41
1:A:54:VAL:O	1:A:59:LYS:NZ	2.53	0.41
1:B:131:SER:HA	1:B:134:MET:HE3	2.02	0.41
1:C:89:THR:HA	1:C:90:PRO:HD3	1.78	0.41
1:B:80:ARG:NH1	1:B:84:ILE:HD11	2.35	0.41
2:C:301:RET:H181	2:C:301:RET:H7	1.63	0.41
1:B:119:VAL:HG21	1:B:173:ALA:CB	2.51	0.40
2:B:301:RET:H7	2:B:301:RET:H181	1.75	0.40
1:E:126:LYS:HZ2	1:E:170:GLU:CD	2.23	0.40
1:B:149:LEU:O	1:B:153:ILE:HG12	2.21	0.40
1:A:34:TRP:CD1	1:B:96:ILE:HG21	2.56	0.40
1:B:40:MET:HE2	1:B:69:THR:HA	2.01	0.40
1:B:109:PHE:CD1	1:B:189:MET:HE2	2.57	0.40
1:C:155:GLY:HA3	2:C:301:RET:H31	2.03	0.40
1:E:102:VAL:N	1:E:103:PRO:CD	2.85	0.40
1:E:105:LEU:HD23	1:E:197:TRP:CZ2	2.55	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	227/261 (87%)	215 (95%)	11 (5%)	1 (0%)	30	51
1	B	216/261 (83%)	207 (96%)	8 (4%)	1 (0%)	24	46
1	C	228/261 (87%)	216 (95%)	12 (5%)	0	100	100
1	D	223/261 (85%)	213 (96%)	10 (4%)	0	100	100
1	E	226/261 (87%)	219 (97%)	7 (3%)	0	100	100
All	All	1120/1305 (86%)	1070 (96%)	48 (4%)	2 (0%)	43	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	SER
1	B	174	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	178/200 (89%)	165 (93%)	13 (7%)	13	29
1	B	176/200 (88%)	159 (90%)	17 (10%)	8	17
1	C	179/200 (90%)	166 (93%)	13 (7%)	13	29
1	D	176/200 (88%)	162 (92%)	14 (8%)	11	25
1	E	178/200 (89%)	169 (95%)	9 (5%)	21	45
All	All	887/1000 (89%)	821 (93%)	66 (7%)	13	29

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	MET
1	A	57	LYS
1	A	61	SER
1	A	84	ILE
1	A	89	THR
1	A	193	ILE
1	A	195	VAL
1	A	209	LEU
1	A	218	SER
1	A	220	LEU
1	A	222	LEU
1	A	246	GLU
1	A	252	LEU
1	B	40	MET
1	B	61	SER
1	B	79	MET
1	B	84	ILE
1	B	108	GLU

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Mol	Chain	Res	Type
1	B	112	ILE
1	B	119	VAL
1	B	126	LYS
1	B	132	LEU
1	B	168	MET
1	B	175	VAL
1	B	183	ASN
1	B	189	MET
1	B	195	VAL
1	B	209	LEU
1	B	215	VAL
1	B	226	LEU
1	C	61	SER
1	C	82	VAL
1	C	86	THR
1	C	118	SER
1	C	125	LYS
1	C	142	GLU
1	C	166	LEU
1	C	191	MET
1	C	192	ILE
1	C	195	VAL
1	C	209	LEU
1	C	237	LEU
1	C	247	SER
1	D	32	SER
1	D	65	SER
1	D	85	ASP
1	D	105	LEU
1	D	107	VAL
1	D	117	THR
1	D	119	VAL
1	D	148	VAL
1	D	168	MET
1	D	172	LYS
1	D	183	ASN
1	D	184	SER
1	D	209	LEU
1	D	222	LEU
1	E	28	THR
1	E	29	VAL
1	E	36	VAL

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Mol	Chain	Res	Type
1	E	61	SER
1	E	148	VAL
1	E	168	MET
1	E	177	THR
1	E	209	LEU
1	E	222	LEU

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	241	ASN
1	B	53	GLN
1	B	97	ASN
1	B	255	HIS
1	C	53	GLN
1	C	183	ASN
1	C	231	ASN
1	C	249	ASN
1	D	221	ASN
1	E	53	GLN
1	E	219	ASN
1	E	241	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 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	RET	A	301	1	20,20,21	0.85	0	27,27,28	3.89	15 (55%)
2	RET	E	301	1	20,20,21	0.91	1 (5%)	27,27,28	3.53	12 (44%)
2	RET	D	301	1	20,20,21	0.74	0	27,27,28	3.81	16 (59%)
2	RET	C	301	1	20,20,21	0.96	1 (5%)	27,27,28	3.77	14 (51%)
2	RET	B	301	1	20,20,21	0.84	1 (5%)	27,27,28	4.10	12 (44%)

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	RET	A	301	1	-	0/13/30/31	0/1/1/1
2	RET	E	301	1	-	0/13/30/31	0/1/1/1
2	RET	D	301	1	-	0/13/30/31	0/1/1/1
2	RET	C	301	1	-	0/13/30/31	0/1/1/1
2	RET	B	301	1	-	0/13/30/31	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	RET	C14-C13	2.72	1.35	1.33
2	E	301	RET	C14-C13	2.31	1.35	1.33
2	B	301	RET	C14-C13	2.08	1.35	1.33

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	RET	C17-C1-C6	-13.94	88.38	110.24
2	A	301	RET	C17-C1-C6	-11.83	91.68	110.24
2	D	301	RET	C17-C1-C6	-8.99	96.14	110.24
2	C	301	RET	C17-C1-C6	-8.94	96.22	110.24
2	E	301	RET	C17-C1-C16	-8.75	83.57	108.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	RET	C16-C1-C6	8.02	122.82	110.24
2	E	301	RET	C17-C1-C6	-7.93	97.81	110.24
2	D	301	RET	C17-C1-C16	-7.74	86.46	108.63
2	D	301	RET	C16-C1-C6	7.56	122.10	110.24
2	C	301	RET	C17-C1-C16	-7.24	87.90	108.63
2	B	301	RET	C17-C1-C16	-7.21	87.98	108.63
2	A	301	RET	C17-C1-C16	-6.72	89.38	108.63
2	A	301	RET	C16-C1-C6	6.53	120.49	110.24
2	A	301	RET	C2-C1-C6	6.36	119.68	110.44
2	C	301	RET	C1-C6-C5	-6.24	114.10	122.64
2	D	301	RET	C17-C1-C2	-6.10	85.54	108.95
2	C	301	RET	C16-C1-C6	6.07	119.76	110.24
2	D	301	RET	C1-C6-C5	-5.92	114.55	122.64
2	C	301	RET	C17-C1-C2	-5.89	86.35	108.95
2	E	301	RET	C17-C1-C2	-5.87	86.43	108.95
2	B	301	RET	C2-C1-C6	5.83	118.90	110.44
2	E	301	RET	C16-C1-C6	5.77	119.29	110.24
2	A	301	RET	C17-C1-C2	-5.64	87.31	108.95
2	B	301	RET	C17-C1-C2	-5.41	88.19	108.95
2	D	301	RET	C1-C6-C7	4.95	129.09	115.65
2	E	301	RET	C2-C1-C6	4.90	117.56	110.44
2	C	301	RET	C1-C6-C7	4.75	128.55	115.65
2	C	301	RET	C2-C1-C6	4.58	117.09	110.44
2	C	301	RET	C7-C8-C9	-4.40	119.73	126.23
2	D	301	RET	C2-C1-C6	4.40	116.83	110.44
2	A	301	RET	C1-C6-C5	-4.34	116.71	122.64
2	E	301	RET	C1-C6-C7	4.32	127.36	115.65
2	E	301	RET	C7-C8-C9	-4.16	120.08	126.23
2	C	301	RET	C16-C1-C2	3.70	123.15	108.95
2	E	301	RET	C16-C1-C2	3.68	123.08	108.95
2	B	301	RET	C1-C6-C5	-3.66	117.63	122.64
2	D	301	RET	C7-C8-C9	-3.62	120.87	126.23
2	C	301	RET	C11-C10-C9	-3.60	122.23	127.28
2	A	301	RET	C20-C13-C12	3.55	123.52	118.09
2	E	301	RET	C1-C6-C5	-3.53	117.81	122.64
2	D	301	RET	C10-C11-C12	-3.42	113.28	123.20
2	A	301	RET	C11-C10-C9	-3.31	122.63	127.28
2	A	301	RET	C19-C9-C8	3.31	123.14	118.09
2	B	301	RET	C11-C10-C9	-3.27	122.69	127.28
2	E	301	RET	C10-C11-C12	-3.22	113.88	123.20
2	A	301	RET	C1-C6-C7	3.18	124.29	115.65
2	D	301	RET	C16-C1-C2	3.15	121.03	108.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	RET	C19-C9-C10	-3.13	117.75	122.82
2	B	301	RET	C20-C13-C12	3.09	122.80	118.09
2	A	301	RET	C7-C8-C9	-3.01	121.78	126.23
2	C	301	RET	C10-C11-C12	-2.96	114.61	123.20
2	C	301	RET	C20-C13-C12	2.95	122.59	118.09
2	E	301	RET	C7-C6-C5	-2.93	114.81	121.56
2	B	301	RET	C1-C6-C7	2.93	123.59	115.65
2	A	301	RET	C16-C1-C2	2.83	119.80	108.95
2	C	301	RET	C18-C5-C6	-2.70	121.54	124.48
2	B	301	RET	C18-C5-C6	-2.58	121.67	124.48
2	D	301	RET	C11-C10-C9	-2.44	123.86	127.28
2	D	301	RET	C18-C5-C6	-2.40	121.87	124.48
2	B	301	RET	C16-C1-C2	2.35	117.97	108.95
2	D	301	RET	C20-C13-C12	2.28	121.58	118.09
2	D	301	RET	C7-C6-C5	-2.26	116.34	121.56
2	E	301	RET	C3-C4-C5	-2.26	110.03	114.06
2	A	301	RET	C10-C11-C12	-2.20	116.81	123.20
2	C	301	RET	C19-C9-C8	2.20	121.44	118.09
2	D	301	RET	C19-C9-C8	2.13	121.35	118.09
2	D	301	RET	C2-C3-C4	2.11	115.92	111.28
2	B	301	RET	C19-C9-C8	2.08	121.27	118.09
2	A	301	RET	C12-C13-C14	-2.07	113.06	118.55

There are no chirality outliers.

There are no torsion outliers.

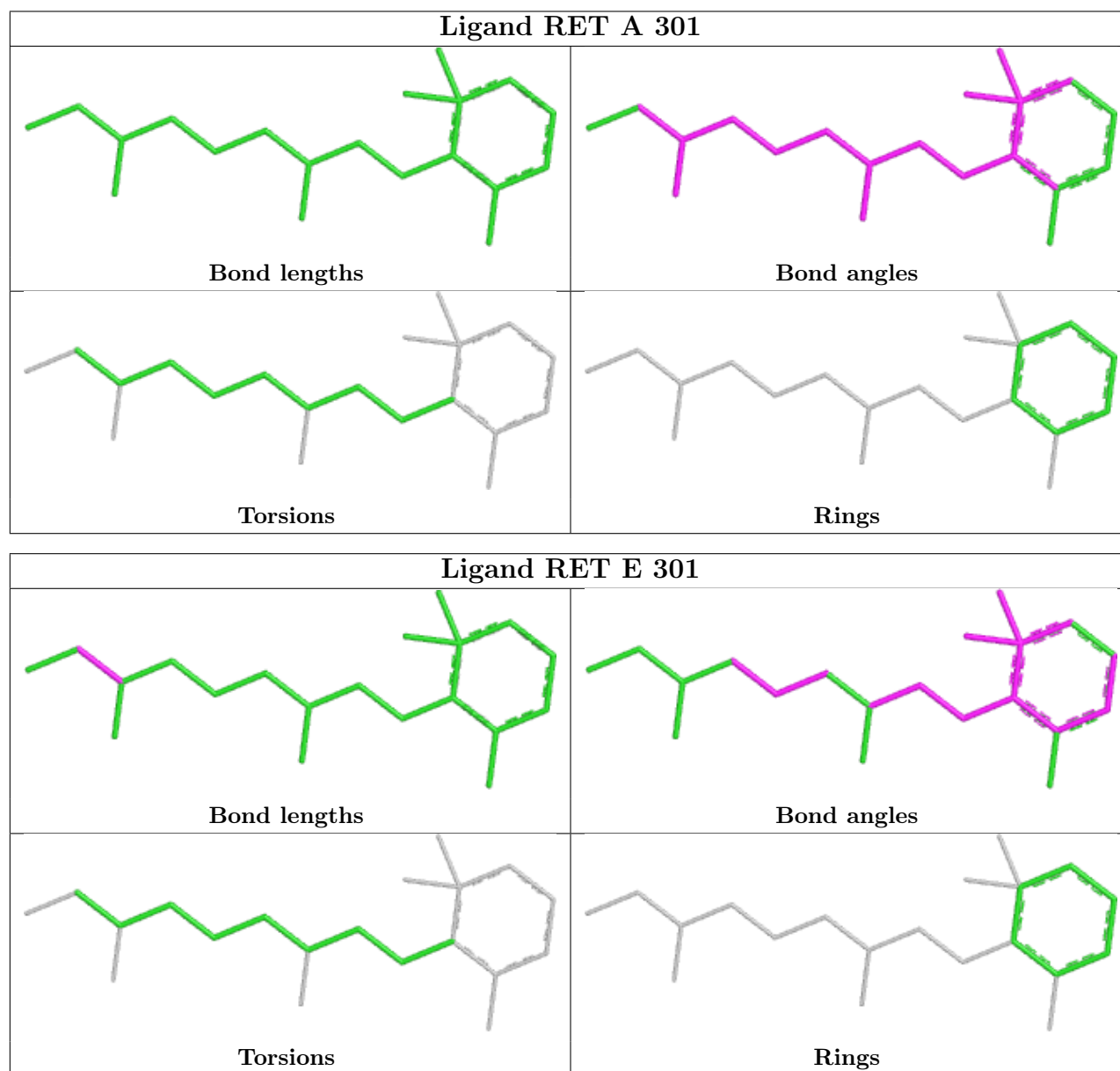
There are no ring outliers.

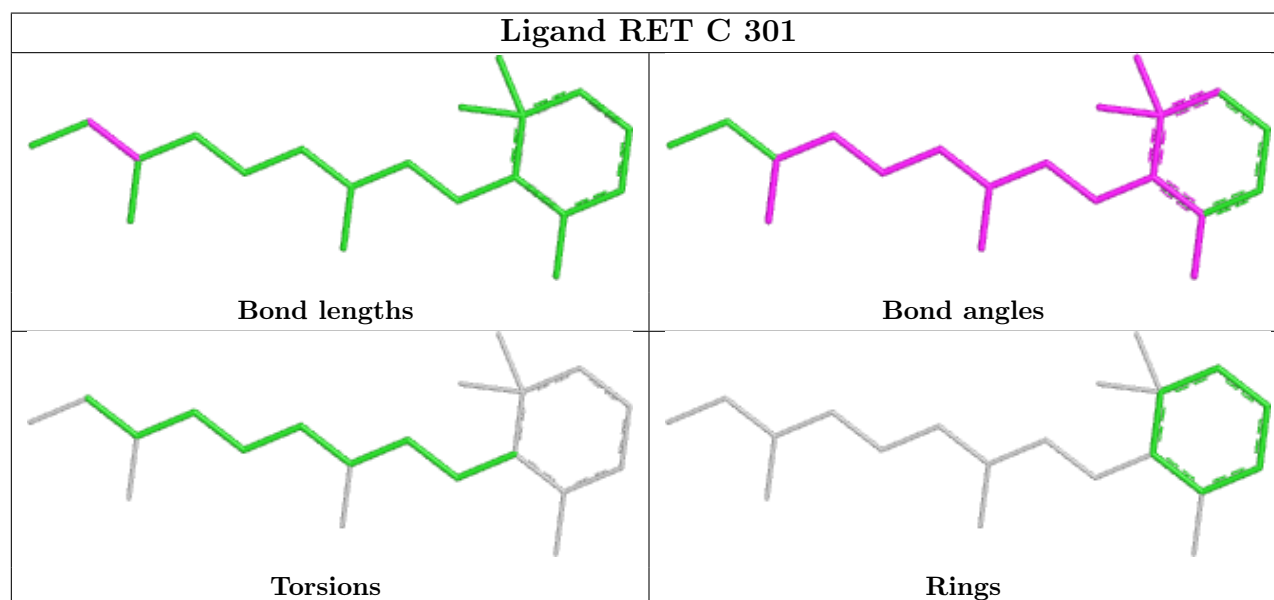
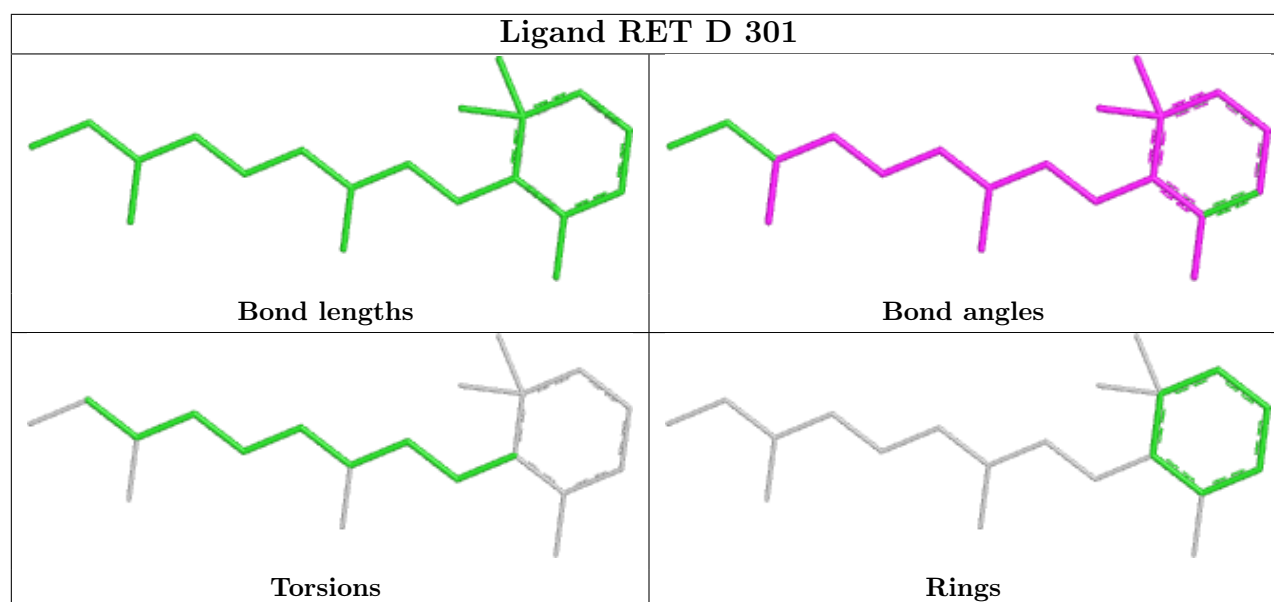
5 monomers are involved in 50 short contacts:

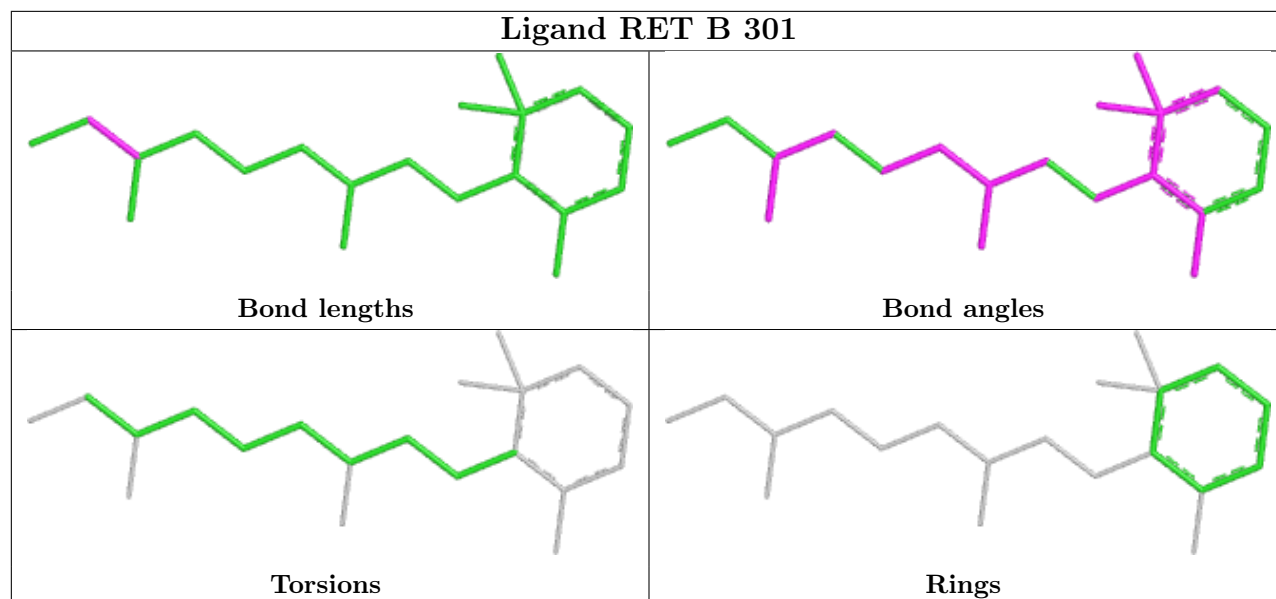
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	RET	22	0
2	E	301	RET	6	0
2	D	301	RET	5	0
2	C	301	RET	9	0
2	B	301	RET	8	0

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	231/261 (88%)	-0.14	6 (2%) 57 51	42, 67, 111, 143	0
1	B	224/261 (85%)	0.11	9 (4%) 42 37	43, 73, 126, 162	0
1	C	232/261 (88%)	-0.19	2 (0%) 81 78	40, 68, 105, 140	0
1	D	229/261 (87%)	-0.12	6 (2%) 57 51	46, 72, 111, 140	0
1	E	230/261 (88%)	-0.26	4 (1%) 69 64	46, 65, 102, 140	0
All	All	1146/1305 (87%)	-0.12	27 (2%) 59 54	40, 68, 112, 162	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	175	VAL	4.3
1	A	175	VAL	4.1
1	B	173	ALA	3.7
1	B	175	VAL	3.4
1	A	252	LEU	3.3
1	D	174	ALA	3.1
1	D	211	GLY	3.0
1	A	191	MET	2.9
1	E	212	GLY	2.8
1	B	185	ALA	2.8
1	B	168	MET	2.8
1	E	252	LEU	2.7
1	A	193	ILE	2.6
1	D	253	LEU	2.5
1	C	119	VAL	2.4
1	E	175	VAL	2.4
1	A	174	ALA	2.3
1	E	251	LYS	2.2
1	B	119	VAL	2.2
1	B	255	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	189	MET	2.2
1	D	20	GLY	2.2
1	D	214	GLY	2.2
1	D	173	ALA	2.2
1	B	186	TYR	2.2
1	B	253	LEU	2.2
1	A	212	GLY	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](#)

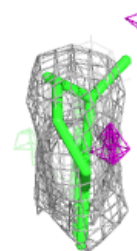
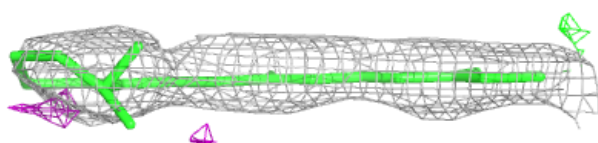
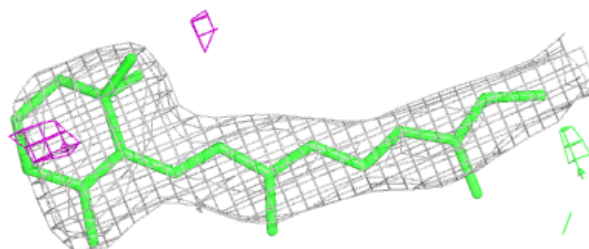
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	RET	D	301	20/21	0.92	0.17	71,78,86,86	0
2	RET	B	301	20/21	0.94	0.11	64,71,77,80	0
2	RET	A	301	20/21	0.94	0.10	54,63,68,69	0
2	RET	C	301	20/21	0.95	0.11	61,69,75,75	0
2	RET	E	301	20/21	0.95	0.10	47,57,69,77	0

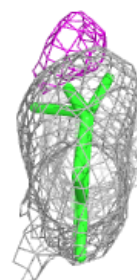
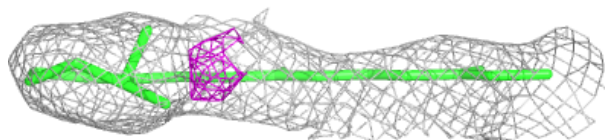
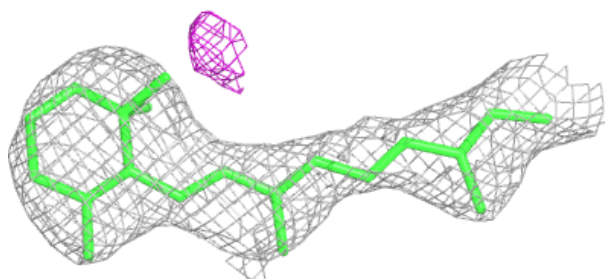
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around RET D 301:

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

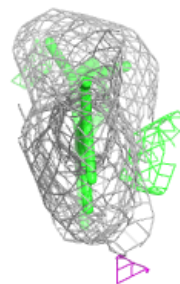
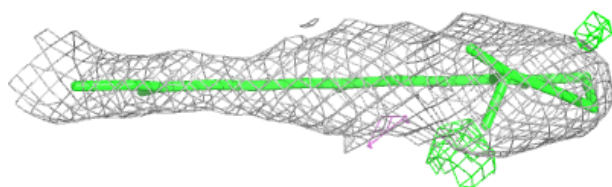
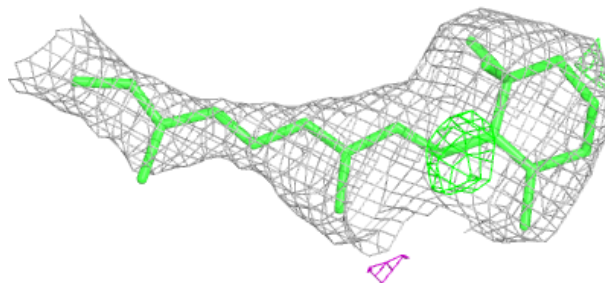
**Electron density around RET B 301:**

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

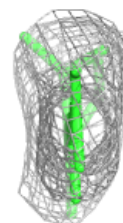
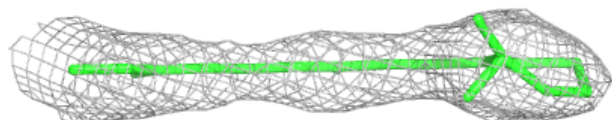
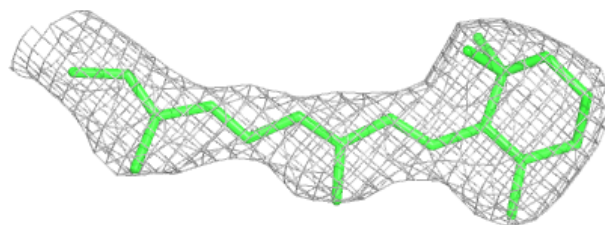


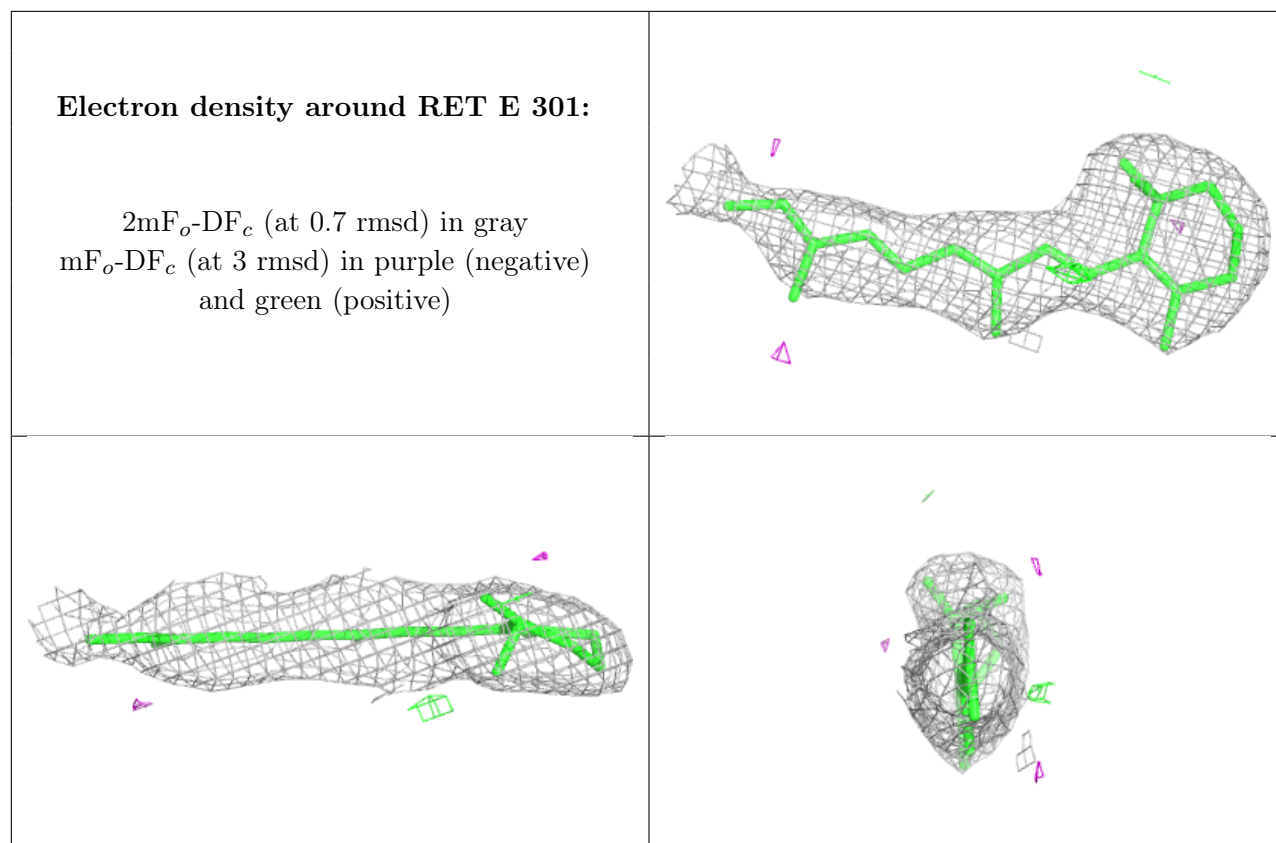
Electron density around RET A 301:

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

**Electron density around RET C 301:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.