



Full wwPDB EM Validation Report ⓘ

Mar 1, 2026 – 04:34 PM UTC

PDB ID : 1BRD / pdb_00001brd
Title : Model for the structure of Bacteriorhodopsin based on high-resolution Electron Cryo-microscopy
Authors : Henderson, R.; Baldwin, J.M.; Ceska, T.A.; Zemlin, F.; Beckmann, E.; Downing, K.H.
Deposited on : 1990-05-23
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at
<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
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)
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

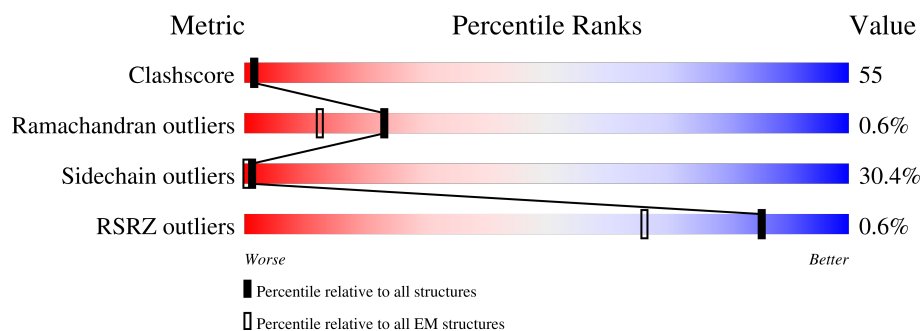
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON CRYSTALLOGRAPHY

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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RSRZ outliers	180361	209

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	248	

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	249	-	-	X	-

2 Entry composition [i](#)

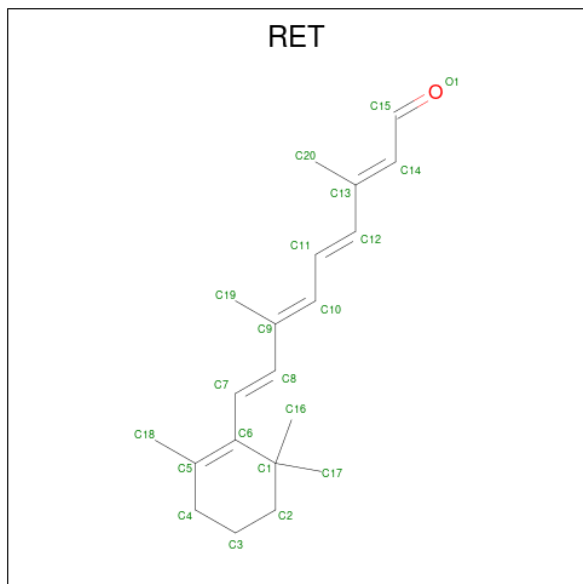
There are 2 unique types of molecules in this entry. The entry contains 1363 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 BACTERIORHODOPSIN PRECURSOR.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	177	Total	C	N	O	S	0	7
			1343	914	203	219	7		

- Molecule 2 is RETINAL (CCD ID: RET) (formula: $C_{20}H_{28}O$).

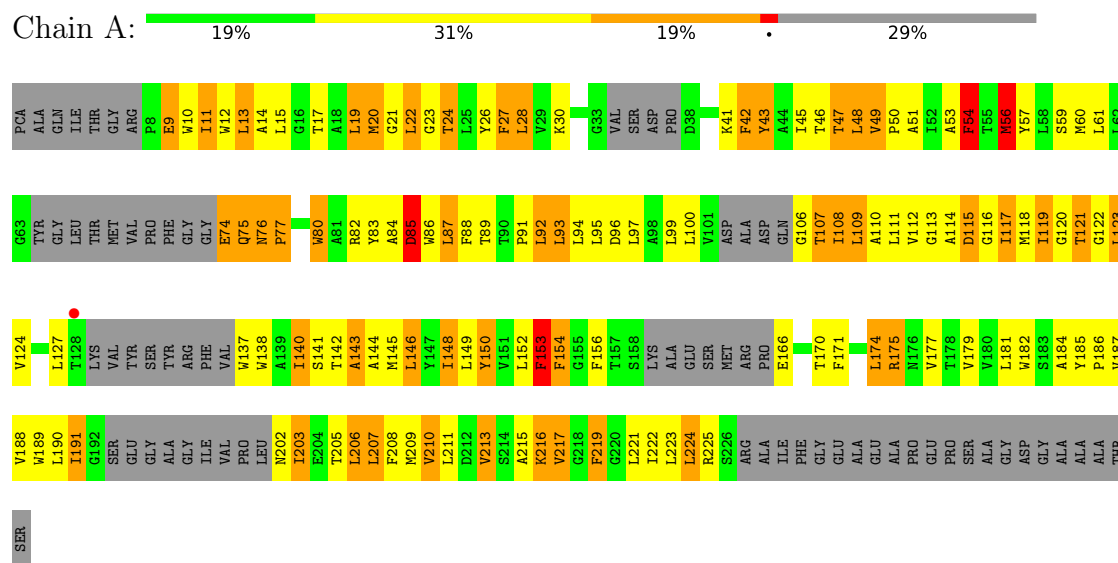


Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	C	0
			20	20	

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. 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. 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: BACTERIORHODOPSIN PRECURSOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	62.45Å 62.45Å 100.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 3.50 100.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-3.50) 50.4 (100.00-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	unknown	Depositor
R, R_{free}	(Not available) , (Not available) 0.387 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 336.7	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	1363	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.84	0/1375	2.11	56/1878 (3.0%)

There are no bond length outliers.

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	PHE	CA-CB-CG	12.99	126.79	113.80
1	A	216	LYS	N-CA-C	12.27	124.65	111.03
1	A	85	ASP	CA-CB-CG	12.18	124.78	112.60
1	A	219	PHE	CA-CB-CG	10.56	124.36	113.80
1	A	9	GLU	CB-CG-CD	8.98	127.87	112.60
1	A	153	PHE	CA-CB-CG	8.78	122.58	113.80
1	A	217	VAL	CB-CA-C	8.56	123.70	112.14
1	A	206	LEU	CA-C-N	8.46	131.96	120.54
1	A	206	LEU	C-N-CA	8.46	131.96	120.54
1	A	203	ILE	N-CA-C	7.92	117.97	110.53
1	A	27	PHE	CA-CB-CG	-7.63	106.17	113.80
1	A	154	PHE	CA-CB-CG	7.30	121.10	113.80
1	A	216	LYS	CB-CA-C	-7.17	99.84	110.95
1	A	42	PHE	N-CA-C	6.98	118.53	111.07
1	A	107	THR	CA-C-N	6.85	134.31	121.97
1	A	107	THR	C-N-CA	6.85	134.31	121.97
1	A	120	GLY	CA-C-N	6.64	130.01	120.79
1	A	120	GLY	C-N-CA	6.64	130.01	120.79
1	A	121	THR	N-CA-C	6.62	119.10	111.02
1	A	108	ILE	CA-C-N	6.59	129.12	120.28
1	A	108	ILE	C-N-CA	6.59	129.12	120.28
1	A	76	ASN	CA-C-N	6.16	125.38	118.97
1	A	76	ASN	C-N-CA	6.16	125.38	118.97
1	A	77	PRO	N-CA-C	6.04	122.11	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	75	GLN	N-CA-C	5.83	117.43	111.14
1	A	207	LEU	N-CA-CB	-5.78	101.32	109.94
1	A	42	PHE	CA-CB-CG	-5.74	108.06	113.80
1	A	210	VAL	CA-C-N	5.73	128.75	120.79
1	A	210	VAL	C-N-CA	5.73	128.75	120.79
1	A	110	ALA	CA-C-N	5.71	128.25	120.54
1	A	110	ALA	C-N-CA	5.71	128.25	120.54
1	A	121	THR	CA-C-N	5.64	126.21	119.94
1	A	121	THR	C-N-CA	5.64	126.21	119.94
1	A	115	ASP	CA-C-N	5.63	126.23	119.98
1	A	115	ASP	C-N-CA	5.63	126.23	119.98
1	A	56	MET	CB-CA-C	5.51	120.22	110.85
1	A	74	GLU	CA-CB-CG	5.48	125.07	114.10
1	A	11	ILE	CA-C-N	5.46	127.53	120.44
1	A	11	ILE	C-N-CA	5.46	127.53	120.44
1	A	9	GLU	CA-C-N	5.35	127.77	120.54
1	A	9	GLU	C-N-CA	5.35	127.77	120.54
1	A	116	GLY	N-CA-C	-5.29	106.38	112.73
1	A	123	LEU	N-CA-C	-5.24	105.48	111.14
1	A	19	LEU	CA-C-N	5.15	127.14	120.44
1	A	19	LEU	C-N-CA	5.15	127.14	120.44
1	A	148	ILE	CA-C-N	5.13	127.11	120.44
1	A	148	ILE	C-N-CA	5.13	127.11	120.44
1	A	92	LEU	CB-CA-C	5.11	119.04	110.92
1	A	117	ILE	CA-C-N	5.05	127.36	120.54
1	A	117	ILE	C-N-CA	5.05	127.36	120.54
1	A	80	TRP	N-CA-C	5.04	117.16	111.11
1	A	59	SER	CA-C-N	5.03	126.98	120.44
1	A	59	SER	C-N-CA	5.03	126.98	120.44
1	A	143	ALA	CA-C-N	5.02	127.01	120.28
1	A	143	ALA	C-N-CA	5.02	127.01	120.28
1	A	28	LEU	N-CA-C	-5.01	104.91	111.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1343	0	1390	137	0
2	A	20	0	27	28	0
All	All	1363	0	1417	154	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 55.

All (154) 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)
2:A:249:RET:C4	2:A:249:RET:C5	1.75	1.64
2:A:249:RET:C1	2:A:249:RET:C2	1.76	1.59
1:A:175:ARG:O	1:A:179:VAL:HG23	1.41	1.19
2:A:249:RET:C2	2:A:249:RET:C17	2.30	1.08
1:A:42:PHE:HB3	1:A:223:LEU:HD21	1.45	0.98
1:A:114:ALA:HA	1:A:117:ILE:HD12	1.55	0.89
1:A:215:ALA:O	1:A:219:PHE:HB2	1.75	0.86
1:A:118:MET:HG2	1:A:141:SER:HB3	1.57	0.86
1:A:91:PRO:HB3	1:A:112:VAL:HG22	1.63	0.79
1:A:121:THR:HB	1:A:137:TRP:HB3	1.65	0.79
2:A:249:RET:H173	2:A:249:RET:C3	2.13	0.78
1:A:149:LEU:HD21	1:A:179:VAL:HG13	1.64	0.78
2:A:249:RET:C5	2:A:249:RET:C2	2.54	0.78
2:A:249:RET:C17	2:A:249:RET:C3	2.62	0.77
1:A:76:ASN:HB2	1:A:77:PRO:HD3	1.66	0.77
1:A:145:MET:HA	1:A:148:ILE:HD12	1.70	0.73
1:A:80:TRP:HA	1:A:83:TYR:HD2	1.55	0.71
1:A:13:LEU:HD13	1:A:205:THR:HG23	1.73	0.71
1:A:49:VAL:HG13	1:A:50:PRO:HD3	1.73	0.70
1:A:117:ILE:O	1:A:121:THR:HG23	1.92	0.70
1:A:123:LEU:O	1:A:127:LEU:HB2	1.92	0.69
1:A:77:PRO:HA	1:A:80:TRP:CH2	2.28	0.69
2:A:249:RET:C2	2:A:249:RET:C16	2.65	0.69
1:A:189:TRP:CD1	2:A:249:RET:H21	2.28	0.68
1:A:80:TRP:HA	1:A:83:TYR:CD2	2.29	0.68
1:A:97:LEU:HD13	1:A:174:LEU:HD13	1.76	0.68
2:A:249:RET:C1	2:A:249:RET:C3	2.65	0.68
1:A:11:ILE:HB	1:A:12:TRP:CE3	2.28	0.67
2:A:249:RET:C4	2:A:249:RET:H173	2.25	0.67
1:A:20:MET:HE2	1:A:216:LYS:HD3	1.77	0.66
1:A:117:ILE:HD13	1:A:140:ILE:HD13	1.76	0.66
1:A:185:TYR:HB2	1:A:186:PRO:HD3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:THR:CB	1:A:137:TRP:HB3	2.26	0.66
1:A:42:PHE:HZ	1:A:100:LEU:HD13	1.60	0.66
1:A:123:LEU:HD12	1:A:127:LEU:HG	1.79	0.64
1:A:114:ALA:O	1:A:117:ILE:HB	1.99	0.63
2:A:249:RET:H173	2:A:249:RET:H31	1.82	0.61
1:A:13:LEU:HD21	1:A:60:MET:HE2	1.83	0.61
1:A:41:LYS:O	1:A:45:ILE:HG12	2.01	0.61
1:A:97:LEU:HD22	1:A:174:LEU:HD22	1.82	0.61
1:A:170:THR:O	1:A:174:LEU:HB2	2.01	0.61
2:A:249:RET:C4	2:A:249:RET:C6	2.66	0.61
1:A:76:ASN:HB3	1:A:127:LEU:HD13	1.82	0.60
1:A:12:TRP:CD1	1:A:206:LEU:HA	2.37	0.60
1:A:189:TRP:CG	2:A:249:RET:H21	2.37	0.59
1:A:85:ASP:O	1:A:89:THR:HG22	2.02	0.59
1:A:43:TYR:O	1:A:47:THR:HB	2.03	0.58
2:A:249:RET:C5	2:A:249:RET:C3	2.79	0.57
1:A:114:ALA:CB	1:A:144:ALA:HB1	2.33	0.57
1:A:11:ILE:HB	1:A:12:TRP:HE3	1.69	0.57
1:A:87:LEU:O	1:A:91:PRO:HD2	2.04	0.57
2:A:249:RET:C4	2:A:249:RET:C1	2.74	0.56
1:A:149:LEU:HD11	1:A:179:VAL:HA	1.87	0.56
1:A:184:ALA:HB3	1:A:211:LEU:HD21	1.88	0.56
1:A:46:THR:HG21	1:A:223:LEU:HD23	1.86	0.56
1:A:48:LEU:HD21	1:A:92:LEU:HD23	1.87	0.56
1:A:88:PHE:O	1:A:92:LEU:HD22	2.06	0.55
1:A:97:LEU:HD13	1:A:174:LEU:HD22	1.88	0.55
1:A:87:LEU:O	1:A:91:PRO:CD	2.54	0.55
1:A:42:PHE:CZ	1:A:100:LEU:HD13	2.41	0.55
1:A:175:ARG:O	1:A:179:VAL:CG2	2.35	0.55
1:A:185:TYR:HB2	1:A:186:PRO:CD	2.36	0.55
1:A:14:ALA:HA	1:A:61:LEU:HD22	1.89	0.55
1:A:12:TRP:O	1:A:15:LEU:HB3	2.06	0.55
1:A:121:THR:OG1	1:A:122:GLY:N	2.41	0.54
1:A:93:LEU:HD22	2:A:249:RET:C20	2.38	0.54
1:A:87:LEU:HG	1:A:88:PHE:N	2.24	0.53
1:A:77:PRO:HA	1:A:80:TRP:CZ3	2.44	0.53
1:A:89:THR:HG21	2:A:249:RET:C15	2.38	0.53
1:A:202:ASN:O	1:A:206:LEU:HB2	2.08	0.53
1:A:100:LEU:HG	1:A:171:PHE:HZ	1.73	0.53
1:A:106:GLY:HA2	1:A:109:LEU:CD1	2.39	0.53
1:A:117:ILE:HG23	1:A:137:TRP:HZ3	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:GLY:O	1:A:117:ILE:HG13	2.08	0.52
1:A:177:VAL:O	1:A:181:LEU:HD23	2.09	0.52
1:A:53:ALA:HB2	1:A:216:LYS:HE2	1.91	0.52
1:A:115:ASP:HA	1:A:118:MET:SD	2.50	0.51
1:A:117:ILE:HD13	1:A:140:ILE:CD1	2.39	0.51
1:A:203:ILE:HG12	1:A:207:LEU:HD22	1.91	0.51
1:A:21:GLY:HA2	1:A:54:PHE:CD2	2.46	0.51
1:A:156:PHE:CD2	1:A:171:PHE:HD2	2.28	0.51
1:A:19:LEU:O	1:A:22:LEU:HB3	2.11	0.50
1:A:47:THR:O	1:A:50:PRO:HD2	2.10	0.50
1:A:114:ALA:HB1	1:A:144:ALA:HB1	1.92	0.50
1:A:181:LEU:HD13	1:A:211:LEU:HD12	1.93	0.50
1:A:202:ASN:HD22	1:A:203:ILE:N	2.09	0.50
1:A:117:ILE:HG23	1:A:137:TRP:CZ3	2.46	0.50
1:A:9:GLU:O	1:A:205:THR:HG21	2.12	0.50
1:A:26:TYR:OH	1:A:224:LEU:HD22	2.12	0.50
1:A:181:LEU:O	1:A:211:LEU:HD11	2.12	0.49
1:A:100:LEU:HG	1:A:171:PHE:CZ	2.48	0.49
1:A:48:LEU:HD23	1:A:49:VAL:N	2.28	0.49
1:A:206:LEU:HD23	1:A:207:LEU:N	2.27	0.49
1:A:121:THR:HG22	1:A:137:TRP:CE3	2.48	0.49
1:A:48:LEU:HD21	1:A:92:LEU:CG	2.43	0.49
2:A:249:RET:C2	2:A:249:RET:H172	2.36	0.48
1:A:185:TYR:O	1:A:189:TRP:HB2	2.12	0.48
2:A:249:RET:C4	2:A:249:RET:C17	2.90	0.48
1:A:153:PHE:CE1	1:A:175:ARG:HD2	2.48	0.48
1:A:24:THR:O	1:A:28:LEU:HB2	2.13	0.48
1:A:97:LEU:HD22	1:A:174:LEU:CD2	2.43	0.48
1:A:106:GLY:O	1:A:107:THR:C	2.56	0.48
1:A:137:TRP:O	1:A:140:ILE:HG22	2.13	0.48
1:A:138:TRP:HE1	2:A:249:RET:H183	1.79	0.47
2:A:249:RET:C17	2:A:249:RET:H31	2.40	0.47
1:A:121:THR:CG2	1:A:137:TRP:HB3	2.44	0.47
1:A:20:MET:HB3	1:A:54:PHE:HB2	1.95	0.47
1:A:24:THR:OG1	1:A:51:ALA:HB2	2.14	0.47
2:A:249:RET:H173	2:A:249:RET:H41	1.96	0.47
1:A:53:ALA:O	1:A:57:TYR:HB2	2.14	0.47
1:A:114:ALA:O	1:A:118:MET:HG3	2.15	0.47
1:A:202:ASN:ND2	1:A:203:ILE:N	2.62	0.47
1:A:174:LEU:HD21	1:A:219:PHE:CZ	2.50	0.46
1:A:48:LEU:HD21	1:A:92:LEU:CD2	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:VAL:O	1:A:213:VAL:HG12	2.15	0.46
1:A:100:LEU:HB3	1:A:171:PHE:CE1	2.51	0.46
1:A:82:ARG:HH11	1:A:208:PHE:HD1	1.64	0.46
1:A:189:TRP:HD1	2:A:249:RET:H163	1.81	0.46
1:A:87:LEU:HA	1:A:115:ASP:OD2	2.15	0.45
1:A:86:TRP:HZ3	2:A:249:RET:H171	1.79	0.45
1:A:80:TRP:HB3	1:A:123:LEU:HD13	1.99	0.45
1:A:53:ALA:CB	1:A:216:LYS:HE2	2.46	0.45
1:A:56:MET:HB3	1:A:85:ASP:HB2	1.98	0.45
1:A:111:LEU:O	1:A:114:ALA:HB3	2.16	0.45
1:A:182:TRP:CE2	2:A:249:RET:H191	2.52	0.44
1:A:54:PHE:C	1:A:54:PHE:CD1	2.95	0.44
1:A:22:LEU:HD23	1:A:23:GLY:N	2.32	0.44
1:A:148:ILE:H	1:A:148:ILE:HG13	1.66	0.43
1:A:185:TYR:CB	1:A:186:PRO:CD	2.95	0.43
1:A:189:TRP:CD1	2:A:249:RET:H172	2.53	0.43
1:A:121:THR:HG22	1:A:137:TRP:HB3	2.00	0.43
1:A:188:VAL:HA	1:A:191:ILE:HB	2.00	0.43
1:A:10:TRP:O	1:A:11:ILE:C	2.61	0.43
1:A:45:ILE:O	1:A:48:LEU:HD22	2.18	0.43
1:A:84:ALA:O	1:A:87:LEU:HB3	2.19	0.43
1:A:188:VAL:HA	1:A:191:ILE:HD12	2.01	0.42
1:A:140:ILE:O	1:A:143:ALA:HB3	2.19	0.42
1:A:146:LEU:O	1:A:150:TYR:HB3	2.20	0.42
1:A:56:MET:HB3	1:A:85:ASP:CB	2.50	0.42
1:A:114:ALA:HB2	1:A:144:ALA:HB1	2.01	0.42
1:A:185:TYR:HB3	2:A:249:RET:C16	2.49	0.42
1:A:21:GLY:O	1:A:24:THR:HG22	2.20	0.42
1:A:115:ASP:O	1:A:119:ILE:HG22	2.20	0.42
1:A:123:LEU:CD1	1:A:127:LEU:HG	2.46	0.42
1:A:185:TYR:HB3	2:A:249:RET:H162	2.02	0.41
1:A:48:LEU:HD21	1:A:92:LEU:HG	2.03	0.41
1:A:149:LEU:HD11	1:A:179:VAL:HG22	2.02	0.41
1:A:182:TRP:CD1	1:A:185:TYR:HE1	2.39	0.41
1:A:153:PHE:CD1	1:A:175:ARG:HD2	2.56	0.41
1:A:87:LEU:O	1:A:91:PRO:HD3	2.21	0.41
1:A:12:TRP:NE1	1:A:206:LEU:HG	2.36	0.40
1:A:60:MET:HE3	1:A:82:ARG:HE	1.86	0.40
1:A:96:ASP:O	1:A:99:LEU:HB3	2.21	0.40
1:A:187:VAL:O	1:A:191:ILE:HG13	2.22	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 PDB entries followed by that with respect to all EM entries.

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	163/248 (66%)	141 (86%)	21 (13%)	1 (1%)	21	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	ILE

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 PDB entries followed by that with respect to all EM entries.

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	138/193 (72%)	96 (70%)	42 (30%)	0	2

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	17	THR
1	A	20	MET
1	A	22	LEU
1	A	24	THR
1	A	27	PHE
1	A	30	LYS
1	A	43	TYR
1	A	47	THR
1	A	48	LEU

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Mol	Chain	Res	Type
1	A	49	VAL
1	A	54	PHE
1	A	56	MET
1	A	74	GLU
1	A	75	GLN
1	A	85	ASP
1	A	87	LEU
1	A	93	LEU
1	A	94	LEU
1	A	95	LEU
1	A	108	ILE
1	A	109	LEU
1	A	119	ILE
1	A	124	VAL
1	A	140	ILE
1	A	142	THR
1	A	146	LEU
1	A	150	TYR
1	A	152	LEU
1	A	153	PHE
1	A	154	PHE
1	A	166	GLU
1	A	174	LEU
1	A	175	ARG
1	A	190	LEU
1	A	209	MET
1	A	213	VAL
1	A	217	VAL
1	A	221	LEU
1	A	222	ILE
1	A	224	LEU
1	A	225	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	RET	A	249	1	20,20,21	6.30	15 (75%)	27,27,28	5.42	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	249	1	-	0/13/30/31	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	249	RET	C4-C5	13.05	1.75	1.51
2	A	249	RET	C8-C9	11.15	1.69	1.46
2	A	249	RET	C2-C1	9.51	1.76	1.54
2	A	249	RET	C12-C13	9.28	1.65	1.46
2	A	249	RET	C7-C6	7.22	1.69	1.45
2	A	249	RET	C11-C10	6.68	1.63	1.43
2	A	249	RET	C16-C1	-6.52	1.41	1.53
2	A	249	RET	C1-C6	6.43	1.62	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	249	RET	C17-C1	-5.78	1.42	1.53
2	A	249	RET	C2-C3	-5.12	1.40	1.52
2	A	249	RET	C15-C14	4.62	1.66	1.49
2	A	249	RET	C14-C13	-4.49	1.30	1.33
2	A	249	RET	C18-C5	4.36	1.57	1.50
2	A	249	RET	C3-C4	3.25	1.62	1.52
2	A	249	RET	C20-C13	2.28	1.55	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	249	RET	C1-C6-C5	-20.31	94.88	122.64
2	A	249	RET	C11-C10-C9	11.40	143.26	127.28
2	A	249	RET	C1-C6-C7	9.63	141.78	115.65
2	A	249	RET	C20-C13-C12	5.51	126.51	118.09
2	A	249	RET	C2-C3-C4	-4.84	100.63	111.28
2	A	249	RET	C2-C1-C6	4.60	117.12	110.44
2	A	249	RET	C17-C1-C2	-4.44	91.90	108.95
2	A	249	RET	C18-C5-C6	-3.61	120.54	124.48
2	A	249	RET	C18-C5-C4	3.36	120.76	113.60
2	A	249	RET	C20-C13-C14	-3.30	114.54	123.63
2	A	249	RET	C4-C5-C6	-3.01	118.64	122.70
2	A	249	RET	C19-C9-C10	-2.10	119.42	122.82

There are no chirality outliers.

There are no torsion outliers.

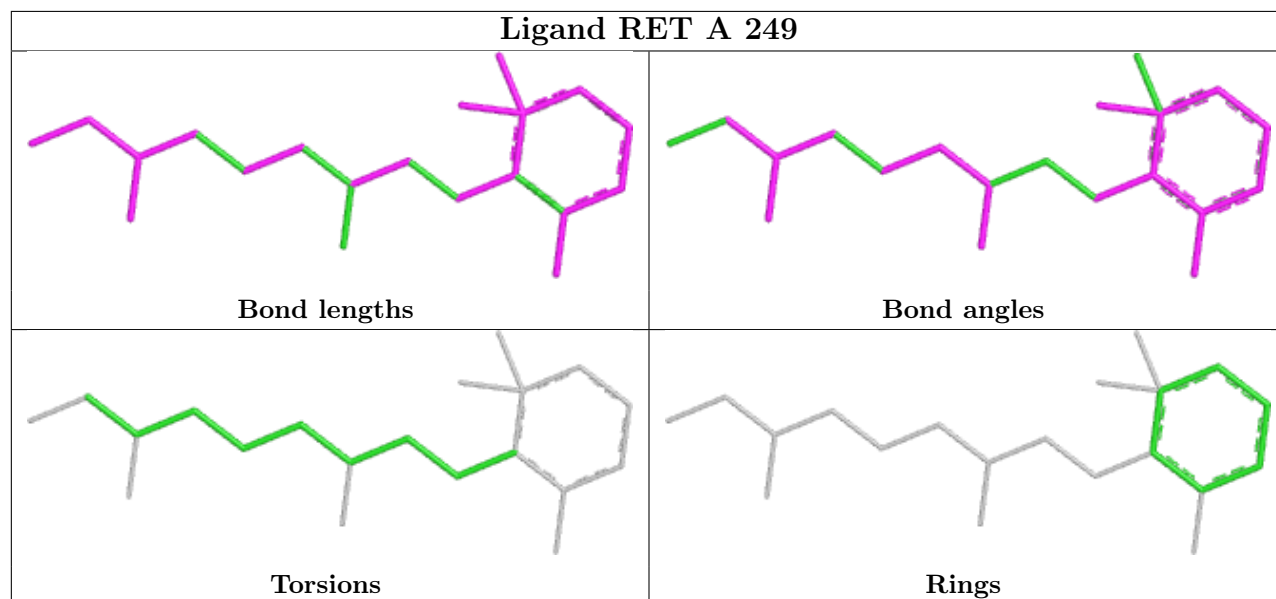
There are no ring outliers.

1 monomer is involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	249	RET	28	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.