



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:51 PM GMT

PDB ID : 1VGO  
Title : Crystal Structure of Archaelhodopsin-2  
Authors : Yoshimura, K.; Enami, N.; Murakami, M.; Okumura, H.; Ihara, K.; Kouyama, T.  
Deposited on : 2004-04-28  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

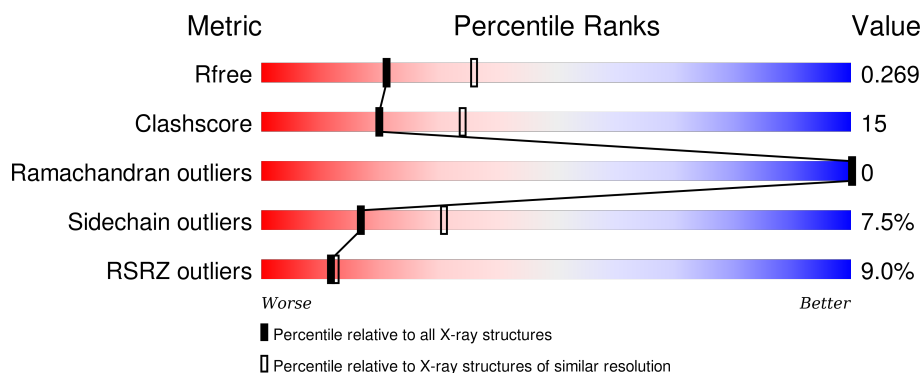
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	253	<div> <div>8%</div> <div>62%</div> <div>26%</div> <div>8%</div> </div>
1	B	253	<div> <div>9%</div> <div>60%</div> <div>32%</div> <div>7%</div> </div>

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	BNG	A	302	-	-	-	X
2	BNG	A	305	-	-	-	X
2	BNG	A	306	-	-	-	X
2	BNG	A	307	-	-	-	X
2	BNG	A	309	-	-	-	X
2	BNG	B	310	-	-	-	X
2	BNG	B	311	-	-	-	X
3	SO4	B	402	-	-	-	X

## 2 Entry composition [i](#)

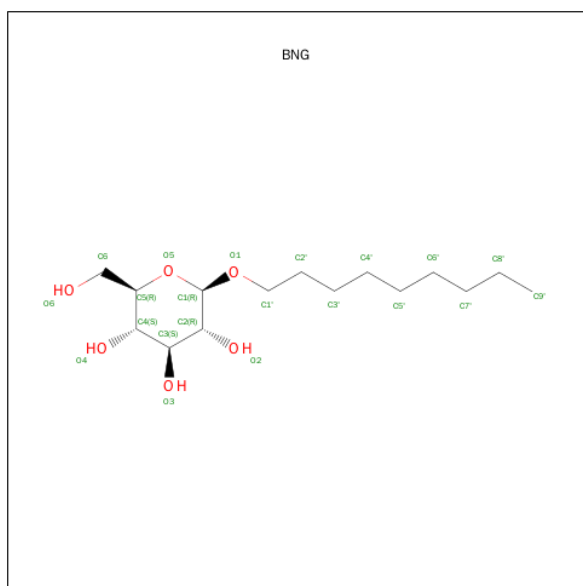
There are 5 unique types of molecules in this entry. The entry contains 3956 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Archaelhodopsin 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1802	1202	280	316	4			
1	B	235	Total	C	N	O	S	0	0	0
			1818	1211	282	321	4			

- Molecule 2 is B-NONYLGLUCOSIDE (three-letter code: BNG) (formula: C<sub>15</sub>H<sub>30</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		

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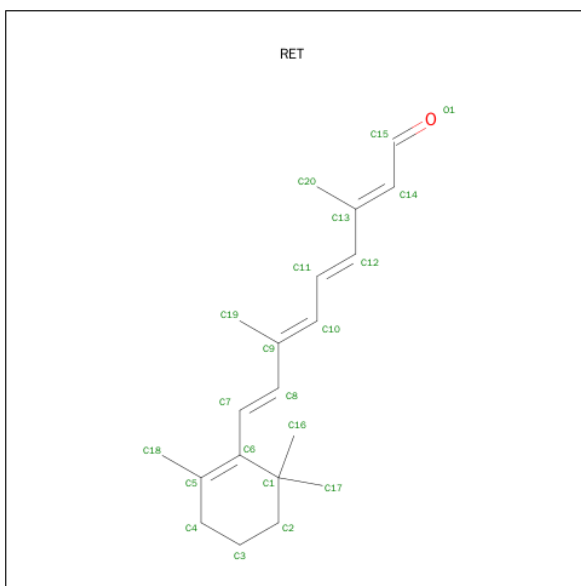
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	A	1	Total	C	O	0	0
			21	15	6		
2	B	1	Total	C	O	0	0
			21	15	6		
2	B	1	Total	C	O	0	0
			21	15	6		
2	B	1	Total	C	O	0	0
			21	15	6		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 4 is RETINAL (three-letter code: RET) (formula:  $C_{20}H_{28}O$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C 20 20	0	0
4	B	1	Total C 20 20	0	0

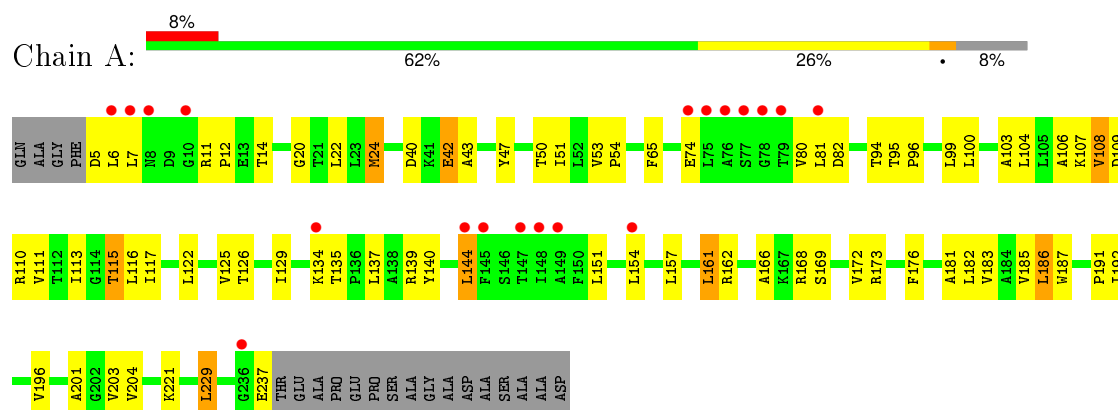
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	16	Total O 16 16	0	0
5	B	13	Total O 13 13	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: Archaelhodopsin 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.88Å 139.48Å 108.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.94 – 2.50 40.58 – 2.50	Depositor EDS
% Data completeness (in resolution range)	83.5 (14.94-2.50) 83.3 (40.58-2.50)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.241 , 0.268 0.241 , 0.269	Depositor DCC
$R_{free}$ test set	1313 reflections (4.87%)	DCC
Wilson B-factor (Å <sup>2</sup> )	57.4	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 62.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 27030 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3956	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: BNG, SO4, 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	0.39	0/1843	0.58	0/2519
1	B	0.38	0/1859	0.59	0/2541
All	All	0.39	0/3702	0.58	0/5060

There are no bond length outliers.

There are no bond angle outliers.

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	1802	0	1870	55	0
1	B	1818	0	1883	62	0
2	A	189	0	269	4	0
2	B	63	0	90	1	0
3	A	5	0	0	1	0
3	B	10	0	0	0	0
4	A	20	0	27	3	0
4	B	20	0	27	4	0
5	A	16	0	0	0	0
5	B	13	0	0	0	0
All	All	3956	0	4166	120	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 15.

All (120) 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:B:94:THR:HG21	1:B:221:LYS:HE3	1.39	1.03
1:A:94:THR:HG21	1:A:221:LYS:HE3	1.48	0.95
1:A:108:VAL:HG11	1:A:161:LEU:HD22	1.67	0.77
1:A:53:VAL:HG13	1:A:94:THR:HG23	1.68	0.76
1:B:94:THR:HG21	1:B:221:LYS:CE	2.18	0.73
1:A:203:VAL:HG23	1:A:204:VAL:HG23	1.72	0.72
1:A:166:ALA:HA	1:A:173:ARG:HD3	1.72	0.72
1:B:75:LEU:HD22	1:B:139:ARG:NH2	2.05	0.72
1:B:147:THR:O	1:B:151:LEU:HD13	1.90	0.71
1:B:232:ARG:O	1:B:235:LEU:HB2	1.91	0.70
1:A:135:THR:HG22	1:A:137:LEU:H	1.57	0.69
1:B:74:GLU:HA	1:B:80:VAL:HA	1.76	0.67
1:B:50:THR:O	1:B:54:PRO:HD2	1.94	0.67
1:B:108:VAL:HB	1:B:112:THR:HG21	1.77	0.67
1:B:214:MET:O	1:B:218:VAL:HG12	1.95	0.67
1:B:53:VAL:HG13	1:B:94:THR:HG23	1.79	0.64
1:A:94:THR:HG21	1:A:221:LYS:CE	2.25	0.64
1:B:94:THR:CG2	1:B:221:LYS:HE3	2.24	0.62
1:A:196:VAL:O	1:A:203:VAL:HG22	1.98	0.62
1:A:50:THR:O	1:A:54:PRO:HD2	1.99	0.61
1:B:42:GLU:OE2	1:B:238:THR:HG23	2.01	0.61
1:B:75:LEU:HD22	1:B:139:ARG:HH21	1.64	0.61
1:B:63:ALA:HA	1:B:68:ILE:HD12	1.83	0.60
1:B:93:PHE:C	1:B:96:PRO:HD2	2.21	0.60
1:A:95:THR:OG1	1:A:96:PRO:HD3	2.04	0.58
1:B:139:ARG:HH22	2:B:312:BNG:H61	1.68	0.57
1:B:179:LEU:O	1:B:183:VAL:HG23	2.05	0.57
1:A:80:VAL:C	1:A:81:LEU:HD12	2.25	0.57
1:A:135:THR:HG22	1:A:137:LEU:N	2.19	0.56
1:A:157:LEU:O	1:A:162:ARG:HB2	2.04	0.56
1:B:103:ALA:HA	1:B:161:LEU:HD11	1.87	0.56
1:B:24:MET:HB3	1:B:58:SER:HB3	1.88	0.56
1:B:103:ALA:HA	1:B:161:LEU:CD1	2.34	0.56
1:B:107:LYS:HG3	1:B:168:ARG:HD3	1.87	0.55
1:B:16:TRP:CE3	1:B:211:LEU:HD23	2.42	0.55
1:A:181:ALA:O	1:A:185:VAL:HG12	2.06	0.54
1:B:185:VAL:HG13	1:B:186:LEU:N	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ILE:HD13	2:A:301:BNG:H5'1	1.90	0.53
1:B:113:ILE:O	1:B:117:ILE:HD13	2.09	0.53
1:B:46:TYR:CD2	1:B:234:ILE:HB	2.44	0.53
1:B:109:ASP:O	1:B:112:THR:HG22	2.08	0.53
1:A:137:LEU:HA	2:A:302:BNG:H2'2	1.90	0.52
1:A:20:GLY:O	1:A:24:MET:HB2	2.09	0.52
1:A:201:ALA:HA	2:A:304:BNG:H1	1.90	0.52
1:B:40:ASP:HB3	1:B:43:ALA:HB3	1.91	0.52
1:A:42:GLU:HG3	1:A:237:GLU:HB3	1.91	0.52
1:A:182:LEU:O	1:A:186:LEU:HB2	2.10	0.52
1:B:191:PRO:HB3	4:B:261:RET:H183	1.91	0.52
1:B:230:ARG:O	1:B:230:ARG:HG2	2.09	0.52
1:A:144:LEU:HD11	2:A:303:BNG:H4'2	1.90	0.52
1:B:144:LEU:HD13	1:B:148:ILE:HD11	1.92	0.52
1:A:122:LEU:O	1:A:125:VAL:HG12	2.10	0.51
1:A:47:TYR:O	1:A:51:ILE:HG13	2.10	0.51
1:B:185:VAL:HG13	1:B:186:LEU:H	1.75	0.51
1:A:111:VAL:O	1:A:115:THR:HG22	2.11	0.51
1:B:91:TRP:CD1	4:B:261:RET:H14	2.46	0.51
1:B:109:ASP:O	1:B:112:THR:CG2	2.59	0.51
1:B:121:ALA:O	1:B:125:VAL:HG12	2.10	0.51
1:A:125:VAL:O	1:A:129:ILE:HG13	2.12	0.50
1:A:40:ASP:HB3	1:A:43:ALA:HB3	1.92	0.50
1:A:192:ILE:O	1:A:196:VAL:HG22	2.11	0.50
1:A:5:ASP:HA	1:A:12:PRO:HA	1.93	0.49
1:A:7:LEU:N	1:A:7:LEU:HD12	2.28	0.49
1:B:203:VAL:HG23	1:B:204:VAL:HG23	1.93	0.49
1:B:144:LEU:HD13	1:B:148:ILE:CD1	2.43	0.48
1:B:92:LEU:O	1:B:96:PRO:HG2	2.12	0.48
1:B:66:PHE:HB2	1:B:68:ILE:HG13	1.95	0.48
1:A:11:ARG:O	1:A:14:THR:HG23	2.13	0.48
1:B:195:ILE:O	1:B:200:GLY:HA3	2.14	0.48
1:B:134:LYS:NZ	1:B:134:LYS:HB3	2.29	0.48
1:A:110:ARG:HG3	1:A:110:ARG:HH11	1.78	0.48
1:A:125:VAL:HG13	1:A:126:THR:N	2.29	0.47
1:A:110:ARG:NH1	1:A:110:ARG:HG3	2.30	0.47
1:A:106:ALA:O	1:A:107:LYS:HB2	2.15	0.47
1:A:80:VAL:O	1:A:81:LEU:HD12	2.16	0.46
1:B:95:THR:N	1:B:96:PRO:CD	2.78	0.46
1:B:53:VAL:HB	1:B:54:PRO:CD	2.45	0.46
1:B:128:LEU:O	1:B:132:LEU:HG	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:VAL:HA	1:B:83:ILE:O	2.15	0.46
1:A:95:THR:N	1:A:96:PRO:CD	2.78	0.46
1:A:191:PRO:HB3	4:A:260:RET:H183	1.97	0.46
1:A:122:LEU:HA	1:A:125:VAL:HG12	1.99	0.45
1:B:169:SER:O	1:B:173:ARG:HG2	2.16	0.45
4:B:261:RET:H161	4:B:261:RET:H8	1.99	0.45
4:A:260:RET:H8	4:A:260:RET:H161	1.98	0.45
1:B:93:PHE:O	1:B:97:LEU:HG	2.16	0.45
1:B:144:LEU:O	1:B:148:ILE:HG13	2.17	0.45
1:A:169:SER:OG	1:A:172:VAL:HG23	2.17	0.45
1:B:72:GLU:HA	1:B:81:LEU:O	2.16	0.45
1:B:181:ALA:O	1:B:185:VAL:HG12	2.18	0.44
1:B:93:PHE:O	1:B:96:PRO:HD2	2.18	0.44
1:B:107:LYS:CG	1:B:168:ARG:HD3	2.47	0.44
4:A:260:RET:H7	4:A:260:RET:H181	1.89	0.44
1:B:109:ASP:CG	1:B:112:THR:HG22	2.38	0.43
1:B:162:ARG:HA	1:B:176:PHE:CE2	2.52	0.43
1:A:135:THR:O	1:A:139:ARG:HG3	2.18	0.43
1:B:211:LEU:C	1:B:211:LEU:HD13	2.38	0.43
1:A:134:LYS:O	1:A:134:LYS:HD2	2.19	0.43
1:B:196:VAL:O	1:B:203:VAL:HG22	2.19	0.42
1:A:53:VAL:HB	1:A:54:PRO:CD	2.50	0.42
1:B:95:THR:OG1	1:B:96:PRO:HD3	2.19	0.42
1:B:90:ASP:OD1	1:B:90:ASP:C	2.56	0.42
1:B:74:GLU:O	1:B:134:LYS:HD2	2.20	0.42
1:A:94:THR:CG2	1:A:221:LYS:HE3	2.35	0.42
1:A:183:VAL:HG13	1:A:187:TRP:CD2	2.55	0.42
1:A:40:ASP:HB2	1:B:237:GLU:OE1	2.21	0.41
1:B:144:LEU:CD1	1:B:148:ILE:HD11	2.50	0.41
1:A:7:LEU:H	1:A:7:LEU:HD12	1.85	0.41
4:B:261:RET:H7	4:B:261:RET:H181	1.90	0.41
1:A:74:GLU:O	1:A:134:LYS:HG2	2.20	0.41
1:A:6:LEU:HD12	1:A:65:PHE:HE1	1.85	0.41
1:A:168:ARG:HB3	3:A:400:SO4:O1	2.20	0.41
1:A:103:ALA:HB1	1:A:113:ILE:HG12	2.03	0.41
1:B:192:ILE:O	1:B:195:ILE:HG22	2.20	0.41
1:A:47:TYR:CE2	1:A:229:LEU:HG	2.56	0.40
1:A:109:ASP:OD2	1:A:111:VAL:HB	2.21	0.40
1:B:87:ARG:HD3	1:B:91:TRP:CH2	2.56	0.40
1:A:113:ILE:O	1:A:117:ILE:HD13	2.20	0.40
1:A:161:LEU:HB3	1:A:176:PHE:CZ	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:TYR:N	1:A:140:TYR:CD2	2.87	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	231/253 (91%)	224 (97%)	7 (3%)	0	100	100
1	B	233/253 (92%)	223 (96%)	10 (4%)	0	100	100
All	All	464/506 (92%)	447 (96%)	17 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	185/196 (94%)	169 (91%)	16 (9%)	13	24
1	B	187/196 (95%)	175 (94%)	12 (6%)	22	39
All	All	372/392 (95%)	344 (92%)	28 (8%)	17	31

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	24	MET
1	A	42	GLU
1	A	82	ASP
1	A	99	LEU
1	A	100	LEU
1	A	104	LEU
1	A	108	VAL
1	A	115	THR
1	A	116	LEU
1	A	144	LEU
1	A	151	LEU
1	A	154	LEU
1	A	161	LEU
1	A	186	LEU
1	A	229	LEU
1	B	9	ASP
1	B	11	ARG
1	B	82	ASP
1	B	99	LEU
1	B	100	LEU
1	B	104	LEU
1	B	116	LEU
1	B	134	LYS
1	B	144	LEU
1	B	154	LEU
1	B	157	LEU
1	B	229	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	ASN
1	B	177	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

17 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$
4	RET	A	260	1	19,20,21	2.98	5 (26%)	27,27,28	1.70	8 (29%)
2	BNG	A	301	-	21,21,21	1.58	5 (23%)	26,26,26	0.80	0
2	BNG	A	302	-	21,21,21	1.90	7 (33%)	26,26,26	0.86	0
2	BNG	A	303	-	21,21,21	2.39	6 (28%)	26,26,26	0.88	1 (3%)
2	BNG	A	304	-	21,21,21	1.83	6 (28%)	26,26,26	0.86	0
2	BNG	A	305	-	21,21,21	1.96	7 (33%)	26,26,26	1.08	2 (7%)
2	BNG	A	306	-	21,21,21	2.87	7 (33%)	26,26,26	1.30	4 (15%)
2	BNG	A	307	-	21,21,21	2.61	6 (28%)	26,26,26	1.13	2 (7%)
2	BNG	A	308	-	21,21,21	2.79	6 (28%)	26,26,26	1.30	4 (15%)
2	BNG	A	309	-	21,21,21	1.54	5 (23%)	26,26,26	0.83	1 (3%)
3	SO4	A	400	-	4,4,4	0.17	0	6,6,6	0.14	0
4	RET	B	261	1	19,20,21	3.15	4 (21%)	27,27,28	1.64	8 (29%)
2	BNG	B	310	-	21,21,21	2.71	6 (28%)	26,26,26	0.80	0
2	BNG	B	311	-	21,21,21	2.55	6 (28%)	26,26,26	0.88	0
2	BNG	B	312	-	21,21,21	2.74	5 (23%)	26,26,26	1.75	3 (11%)
3	SO4	B	401	-	4,4,4	0.27	0	6,6,6	0.09	0
3	SO4	B	402	-	4,4,4	0.33	0	6,6,6	0.12	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
4	RET	A	260	1	-	0/13/30/31	0/1/1/1
2	BNG	A	301	-	-	0/12/32/32	0/1/1/1
2	BNG	A	302	-	-	0/12/32/32	0/1/1/1
2	BNG	A	303	-	-	0/12/32/32	0/1/1/1
2	BNG	A	304	-	-	0/12/32/32	0/1/1/1
2	BNG	A	305	-	-	0/12/32/32	0/1/1/1
2	BNG	A	306	-	-	0/12/32/32	0/1/1/1
2	BNG	A	307	-	-	0/12/32/32	0/1/1/1
2	BNG	A	308	-	-	0/12/32/32	0/1/1/1
2	BNG	A	309	-	-	0/12/32/32	0/1/1/1
3	SO4	A	400	-	-	0/0/0/0	0/0/0/0
4	RET	B	261	1	-	0/13/30/31	0/1/1/1
2	BNG	B	310	-	-	0/12/32/32	0/1/1/1
2	BNG	B	311	-	-	0/12/32/32	0/1/1/1
2	BNG	B	312	-	-	0/12/32/32	0/1/1/1
3	SO4	B	401	-	-	0/0/0/0	0/0/0/0
3	SO4	B	402	-	-	0/0/0/0	0/0/0/0

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	260	RET	C2-C3	-2.08	1.46	1.52
2	A	301	BNG	O5-C5	2.03	1.49	1.44
2	A	307	BNG	C2'-C1'	2.03	1.60	1.50
2	A	302	BNG	O5-C1	2.04	1.47	1.41
2	A	309	BNG	O4-C4	2.04	1.47	1.43
2	B	311	BNG	C4-C5	2.08	1.57	1.53
2	A	308	BNG	C2'-C1'	2.18	1.60	1.50
2	A	306	BNG	C2'-C1'	2.24	1.60	1.50
2	A	306	BNG	C1-C2	2.28	1.59	1.52
2	A	305	BNG	C1-C2	2.30	1.59	1.52
2	B	311	BNG	C2'-C1'	2.34	1.61	1.50
2	A	305	BNG	O5-C1	2.34	1.47	1.41
2	A	301	BNG	C3-C2	2.37	1.58	1.52
2	B	310	BNG	C2'-C1'	2.44	1.61	1.50
2	A	304	BNG	O5-C5	2.45	1.50	1.44
2	A	307	BNG	O5-C1	2.49	1.48	1.41
4	B	261	RET	C2-C1	2.51	1.60	1.54
2	A	309	BNG	C4-C5	2.52	1.58	1.53
2	A	301	BNG	C1-C2	2.52	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	304	BNG	O4-C4	2.60	1.49	1.43
2	A	309	BNG	O5-C5	2.60	1.50	1.44
2	A	309	BNG	C3-C2	2.67	1.59	1.52
2	A	308	BNG	O5-C1	2.68	1.48	1.41
2	B	312	BNG	C1-C2	2.69	1.60	1.52
4	A	260	RET	C2-C1	2.69	1.60	1.54
2	A	304	BNG	C1-C2	2.71	1.60	1.52
2	A	306	BNG	O5-C1	2.76	1.48	1.41
2	A	305	BNG	O4-C4	2.78	1.49	1.43
2	A	304	BNG	C3-C2	2.82	1.59	1.52
2	B	310	BNG	O5-C1	2.86	1.49	1.41
4	A	260	RET	C7-C6	2.90	1.56	1.45
2	A	301	BNG	C4-C5	2.98	1.59	1.53
2	A	305	BNG	O1-C1	2.99	1.45	1.40
2	A	302	BNG	C3-C2	3.00	1.60	1.52
2	A	302	BNG	O4-C4	3.01	1.50	1.43
4	B	261	RET	C7-C6	3.06	1.57	1.45
2	A	302	BNG	O5-C5	3.18	1.52	1.44
2	A	302	BNG	C1-C2	3.25	1.62	1.52
2	A	303	BNG	C3-C2	3.31	1.61	1.52
2	A	305	BNG	O5-C5	3.37	1.52	1.44
2	A	304	BNG	O1-C1	3.39	1.46	1.40
2	A	303	BNG	O4-C4	3.40	1.51	1.43
2	A	302	BNG	O1-C1	3.61	1.46	1.40
2	A	303	BNG	C1-C2	3.79	1.63	1.52
2	A	305	BNG	C3-C2	3.81	1.62	1.52
2	A	309	BNG	O1-C1	3.91	1.47	1.40
2	A	302	BNG	C4-C5	4.00	1.61	1.53
2	A	301	BNG	O1-C1	4.10	1.47	1.40
2	A	303	BNG	O1-C1	4.12	1.47	1.40
2	A	308	BNG	O4-C4	4.24	1.53	1.43
2	A	305	BNG	C4-C5	4.25	1.62	1.53
2	A	306	BNG	O4-C4	4.33	1.53	1.43
2	A	307	BNG	O4-C4	4.41	1.53	1.43
2	A	307	BNG	C3-C2	4.47	1.64	1.52
2	A	304	BNG	C4-C5	4.54	1.62	1.53
2	A	308	BNG	C3-C2	4.56	1.64	1.52
2	B	310	BNG	O5-C5	4.62	1.55	1.44
2	B	310	BNG	O4-C4	4.68	1.54	1.43
2	B	311	BNG	O1-C1	4.71	1.48	1.40
2	A	306	BNG	C3-C2	4.71	1.64	1.52
2	A	306	BNG	O5-C5	4.75	1.56	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	312	BNG	C3-C2	4.78	1.64	1.52
2	A	307	BNG	O5-C5	4.86	1.56	1.44
2	A	308	BNG	O5-C5	4.96	1.56	1.44
2	B	311	BNG	O4-C4	5.06	1.55	1.43
2	B	312	BNG	O4-C4	5.10	1.55	1.43
2	A	303	BNG	O5-C5	5.16	1.57	1.44
2	B	310	BNG	C3-C2	5.19	1.66	1.52
2	B	311	BNG	O5-C5	5.49	1.58	1.44
2	B	311	BNG	C3-C2	5.50	1.66	1.52
2	A	303	BNG	C4-C5	5.51	1.64	1.53
2	B	312	BNG	O5-C5	5.97	1.59	1.44
4	A	260	RET	C5-C6	6.66	1.45	1.34
2	B	312	BNG	O1-C1	6.70	1.52	1.40
4	B	261	RET	C5-C6	6.98	1.45	1.34
2	B	310	BNG	O1-C1	7.01	1.52	1.40
2	A	307	BNG	O1-C1	7.22	1.53	1.40
2	A	308	BNG	O1-C1	8.34	1.55	1.40
2	A	306	BNG	O1-C1	8.79	1.56	1.40
4	A	260	RET	C1-C6	9.63	1.67	1.53
4	B	261	RET	C1-C6	10.54	1.68	1.53

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	260	RET	C1-C6-C5	-3.17	118.00	122.66
4	A	260	RET	C18-C5-C6	-3.04	121.62	124.61
4	B	261	RET	C18-C5-C6	-3.02	121.64	124.61
4	B	261	RET	C1-C6-C5	-2.98	118.28	122.66
2	A	305	BNG	C1'-O1-C1	-2.91	108.86	113.94
4	A	260	RET	C8-C9-C10	-2.79	114.49	118.98
4	B	261	RET	C7-C8-C9	-2.60	122.25	126.22
4	B	261	RET	C8-C9-C10	-2.46	115.02	118.98
4	A	260	RET	C10-C11-C12	-2.45	115.65	123.13
4	B	261	RET	C10-C11-C12	-2.39	115.83	123.13
2	A	307	BNG	C3-C4-C5	-2.33	106.14	110.20
2	A	307	BNG	C6-C5-C4	-2.33	107.28	113.02
4	A	260	RET	C7-C8-C9	-2.31	122.69	126.22
2	A	308	BNG	C3-C4-C5	-2.23	106.32	110.20
2	A	309	BNG	C1'-O1-C1	-2.21	110.08	113.94
2	B	312	BNG	C4-C3-C2	-2.08	106.91	110.79
2	A	306	BNG	C3-C4-C5	-2.04	106.64	110.20
2	A	308	BNG	C6-C5-C4	-2.03	108.00	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	306	BNG	C6-C5-C4	-2.02	108.02	113.02
2	A	305	BNG	O5-C5-C6	2.14	111.77	106.36
4	B	261	RET	C20-C13-C12	2.15	121.67	118.10
2	A	303	BNG	O5-C5-C6	2.32	112.22	106.36
4	B	261	RET	C1-C6-C7	2.33	122.35	115.82
4	A	260	RET	C1-C6-C7	2.45	122.68	115.82
4	A	260	RET	C20-C13-C12	2.46	122.19	118.10
2	A	306	BNG	C1'-O1-C1	2.61	118.50	113.94
4	B	261	RET	C19-C9-C8	2.98	123.06	118.10
2	A	308	BNG	O1-C1-C2	3.07	111.92	108.04
2	A	308	BNG	C1'-O1-C1	3.12	119.39	113.94
4	A	260	RET	C19-C9-C8	3.14	123.31	118.10
2	A	306	BNG	O1-C1-C2	3.53	112.50	108.04
2	B	312	BNG	C1'-O1-C1	3.60	120.24	113.94
2	B	312	BNG	O1-C1-C2	7.03	116.92	108.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	260	RET	3	0
2	A	301	BNG	1	0
2	A	302	BNG	1	0
2	A	303	BNG	1	0
2	A	304	BNG	1	0
3	A	400	SO4	1	0
4	B	261	RET	4	0
2	B	312	BNG	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	233/253 (92%)	0.26	19 (8%) 14 15	40, 51, 82, 94	0
1	B	235/253 (92%)	0.32	23 (9%) 10 10	41, 58, 96, 109	0
All	All	468/506 (92%)	0.29	42 (8%) 12 12	40, 54, 88, 109	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	78	GLY	6.2
1	A	76	ALA	5.4
1	B	79	THR	5.2
1	B	239	GLU	4.5
1	B	7	LEU	4.2
1	B	238	THR	3.9
1	A	148	ILE	3.7
1	B	81	LEU	3.3
1	B	71	THR	3.3
1	B	80	VAL	3.3
1	B	9	ASP	3.2
1	A	6	LEU	3.2
1	A	77	SER	2.9
1	B	10	GLY	2.9
1	B	68	ILE	2.8
1	B	74	GLU	2.8
1	A	134	LYS	2.8
1	A	145	PHE	2.8
1	B	70	VAL	2.7
1	B	73	VAL	2.7
1	A	75	LEU	2.7
1	B	8	ASN	2.7
1	A	236	GLY	2.5
1	A	144	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	78	GLY	2.4
1	A	147	THR	2.4
1	B	75	LEU	2.4
1	B	111	VAL	2.4
1	B	85	TYR	2.3
1	A	149	ALA	2.3
1	A	74	GLU	2.3
1	A	154	LEU	2.3
1	B	108	VAL	2.3
1	A	7	LEU	2.3
1	B	12	PRO	2.2
1	B	76	ALA	2.2
1	A	81	LEU	2.1
1	B	83	ILE	2.1
1	A	10	GLY	2.0
1	A	8	ASN	2.0
1	B	110	ARG	2.0
1	A	79	THR	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	SO4	B	402	5/5	0.61	0.64	28.64	157,157,158,158	0
2	BNG	A	309	21/21	0.50	0.58	9.21	127,135,137,138	0
2	BNG	A	307	21/21	0.57	0.35	5.25	112,121,123,124	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	BNG	A	306	21/21	0.59	0.32	4.82	105,110,112,112	0
2	BNG	B	311	21/21	0.77	0.26	4.80	77,91,98,100	0
2	BNG	A	305	21/21	0.54	0.42	4.05	92,103,108,108	0
2	BNG	A	302	21/21	0.50	0.37	2.99	86,113,118,120	0
2	BNG	B	310	21/21	0.64	0.33	2.92	67,95,101,105	0
2	BNG	B	312	21/21	0.73	0.36	1.94	107,111,113,114	0
4	RET	B	261	20/21	0.97	0.17	1.67	47,50,52,54	0
2	BNG	A	304	21/21	0.71	0.29	1.61	89,100,106,106	0
2	BNG	A	301	21/21	0.71	0.29	1.22	74,77,79,84	0
4	RET	A	260	20/21	0.96	0.22	0.73	39,43,46,46	0
3	SO4	B	401	5/5	0.97	0.17	-	79,79,82,83	0
2	BNG	A	308	21/21	0.57	0.52	-	128,138,140,140	0
2	BNG	A	303	21/21	0.11	0.80	-	126,137,141,141	0
3	SO4	A	400	5/5	0.97	0.12	-	74,75,75,77	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.