



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:55 AM GMT

PDB ID : 3QBK  
Title : Bromide-bound form of pharaonis halorhodopsin  
Authors : Kouyama, T.; Kanada, S.  
Deposited on : 2011-01-13  
Resolution : 2.20 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

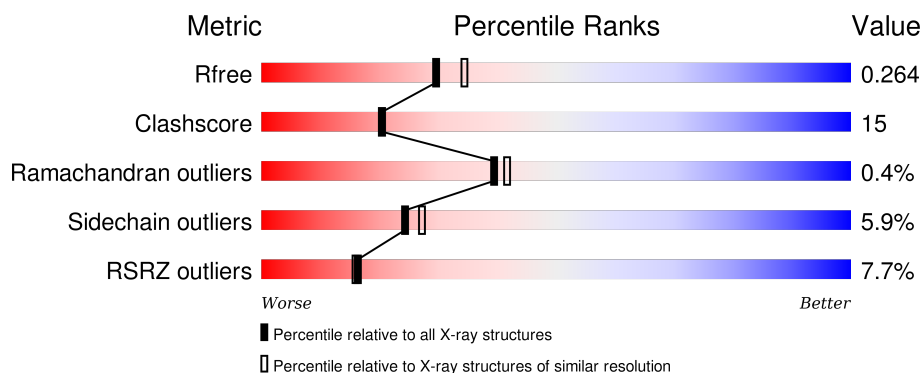
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	291	<div> <div>5%</div> <div>68%</div> <div>20%</div> <div>10%</div> </div>
1	B	291	<div> <div>3%</div> <div>68%</div> <div>18%</div> <div>11%</div> </div>
1	D	291	<div> <div>12%</div> <div>56%</div> <div>31%</div> <div>11%</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
3	22B	A	300	-	-	-	X
3	22B	B	300	-	-	-	X

## 2 Entry composition [i](#)

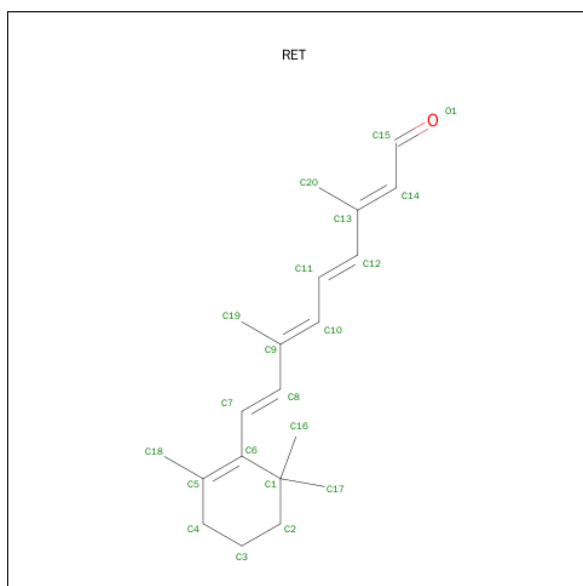
There are 5 unique types of molecules in this entry. The entry contains 6153 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Halorhodopsin.

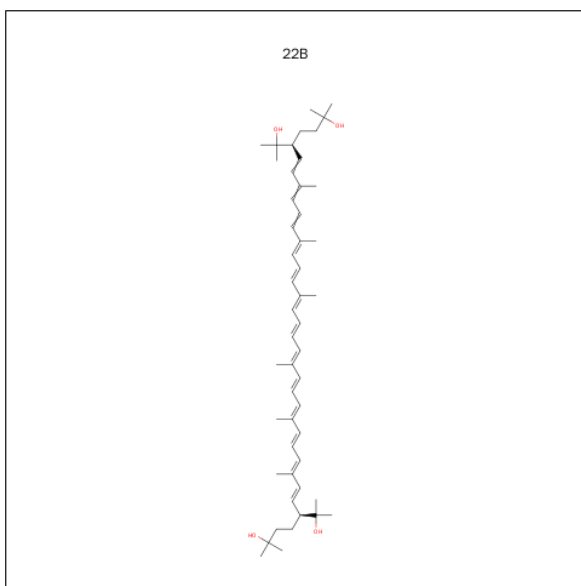
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			1968	1305	300	352	11			
1	B	259	Total	C	N	O	S	0	0	0
			1957	1299	298	349	11			
1	D	259	Total	C	N	O	S	0	0	0
			1957	1299	298	349	11			

- Molecule 2 is RETINAL (three-letter code: RET) (formula:  $C_{20}H_{28}O$ ).



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

- Molecule 3 is BACTERIORUBERIN (three-letter code: 22B) (formula:  $C_{50}H_{76}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			27	25	2		
3	B	1	Total	C	O	0	0
			27	25	2		

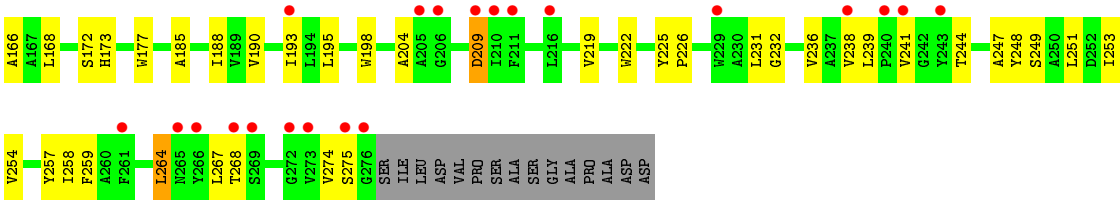
- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Br	0	0
			1	1		
4	A	3	Total	Br	0	0
			3	3		
4	D	2	Total	Br	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	52	Total	O	0	0
			52	52		
5	B	65	Total	O	0	0
			65	65		
5	D	34	Total	O	0	0
			34	34		





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	152.15Å 99.72Å 99.97Å 90.00° 127.80° 90.00°	Depositor
Resolution (Å)	15.00 – 2.20 49.91 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.3 (15.00-2.20) 97.2 (49.91-2.20)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.54 (at 2.20Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.248 , 0.267 0.245 , 0.264	Depositor DCC
$R_{free}$ test set	2936 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.2	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 80.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 58245 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6153	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: RET, BR, 22B

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	0/2014	0.59	0/2758
1	B	0.36	0/2003	0.56	0/2743
1	D	0.32	0/2003	0.54	0/2743
All	All	0.35	0/6020	0.56	0/8244

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 [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	1968	0	2019	46	0
1	B	1957	0	2009	50	0
1	D	1957	0	2009	79	0
2	A	20	0	27	1	0
2	B	20	0	27	1	0
2	D	20	0	27	2	0
3	A	27	0	37	2	0
3	B	27	0	37	1	0
4	A	3	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	2	0	0	0	0
5	A	52	0	0	2	1
5	B	65	0	0	2	0
5	D	34	0	0	2	0
All	All	6153	0	6192	178	1

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 (178) 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:A:238:VAL:HG23	1:A:239:LEU:HD13	1.47	0.96
1:B:236:VAL:HG13	1:B:238:VAL:HG13	1.56	0.88
1:D:145:ASN:HD21	1:D:148:LYS:HG3	1.49	0.78
1:D:254:VAL:HA	1:D:258:ILE:HG22	1.63	0.77
1:D:195:LEU:HD21	1:D:219:VAL:HG21	1.67	0.77
1:D:65:LYS:O	1:D:69:VAL:HG23	1.85	0.76
1:D:131:THR:OG1	1:D:132:PRO:HD3	1.87	0.74
1:B:20:THR:HG22	1:B:23:GLU:H	1.53	0.73
1:D:19:VAL:HG21	1:D:93:VAL:HG11	1.71	0.72
1:A:238:VAL:HG23	1:A:239:LEU:CD1	2.20	0.71
1:D:100:HIS:CD2	1:D:102:ALA:H	2.10	0.69
1:A:20:THR:HG22	1:A:22:ARG:H	1.59	0.68
1:D:185:ALA:O	1:D:188:ILE:HG12	1.94	0.68
1:D:45:ALA:HB3	1:D:79:ILE:HG23	1.77	0.67
1:D:108:MET:HB2	1:D:113:GLU:HG3	1.77	0.66
1:D:164:LEU:O	1:D:168:LEU:HD23	1.96	0.66
1:A:236:VAL:HG13	1:A:238:VAL:HG13	1.77	0.66
1:D:145:ASN:C	1:D:145:ASN:HD22	2.00	0.65
1:B:270:ASN:HB3	1:B:273:VAL:HG12	1.78	0.64
1:D:254:VAL:HA	1:D:258:ILE:CG2	2.27	0.64
1:D:28:VAL:HG13	1:D:34:LEU:HD12	1.79	0.63
1:A:107:VAL:CG2	1:A:109:LEU:HD13	2.29	0.63
1:D:36:SER:O	1:D:40:ILE:HG12	1.99	0.62
1:B:108:MET:HA	1:B:113:GLU:HA	1.81	0.62
1:D:108:MET:HA	1:D:113:GLU:HA	1.80	0.62
1:D:43:ALA:O	1:D:47:LEU:HD13	2.00	0.62
1:A:108:MET:HA	1:A:113:GLU:HA	1.81	0.61
1:D:238:VAL:HG23	1:D:239:LEU:CD1	2.31	0.60
1:D:238:VAL:HG23	1:D:239:LEU:HD13	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:THR:HG21	1:D:264:LEU:HD13	1.85	0.59
1:B:65:LYS:HE2	3:B:300:22B:H243	1.85	0.59
1:B:270:ASN:O	1:B:273:VAL:HG12	2.03	0.57
1:B:270:ASN:HB3	1:B:273:VAL:CG1	2.35	0.57
1:B:19:VAL:CG2	1:B:93:VAL:HG11	2.36	0.56
3:A:300:22B:H253	1:D:151:THR:HG21	1.86	0.56
1:D:135:LEU:HD11	1:D:222:TRP:CH2	2.40	0.56
1:D:54:PHE:HB3	1:D:57:ARG:NH1	2.20	0.56
1:B:93:VAL:HG13	1:B:93:VAL:O	2.05	0.56
1:A:234:GLU:HG3	5:A:511:HOH:O	2.04	0.56
1:A:253:ILE:O	1:A:258:ILE:HG23	2.06	0.56
1:D:93:VAL:HA	1:D:117:VAL:O	2.06	0.55
1:A:121:TRP:HZ3	1:A:125:LEU:HD22	1.70	0.55
1:B:19:VAL:HG21	1:B:93:VAL:HG11	1.89	0.55
1:B:29:LEU:O	1:B:30:ASN:HB2	2.05	0.55
1:D:254:VAL:CA	1:D:258:ILE:HG22	2.36	0.55
1:D:24:LEU:O	1:D:28:VAL:HG23	2.07	0.55
1:A:29:LEU:O	1:A:30:ASN:HB2	2.07	0.55
1:B:34:LEU:O	1:B:38:LEU:HD22	2.06	0.55
1:B:224:GLY:O	1:B:228:VAL:HG23	2.07	0.54
1:A:107:VAL:HG23	1:A:109:LEU:HD13	1.89	0.54
1:B:217:LEU:C	1:B:217:LEU:HD23	2.28	0.54
1:A:233:VAL:HG23	1:A:234:GLU:OE1	2.08	0.54
1:D:190:VAL:O	1:D:193:ILE:HG13	2.08	0.53
1:B:253:ILE:O	1:B:258:ILE:HG23	2.08	0.53
1:A:214:LEU:CD1	1:A:263:LEU:HD13	2.39	0.53
1:D:145:ASN:ND2	1:D:148:LYS:H	2.07	0.53
1:D:121:TRP:HA	1:D:124:TYR:CD2	2.44	0.53
1:D:76:VAL:HA	1:D:79:ILE:HD11	1.91	0.52
1:B:185:ALA:O	1:B:188:ILE:HG12	2.09	0.52
1:B:36:SER:O	1:B:40:ILE:HG12	2.09	0.52
1:D:28:VAL:HG13	1:D:34:LEU:CD1	2.39	0.52
1:A:26:GLU:O	1:A:29:LEU:O	2.27	0.52
1:B:20:THR:HG23	1:B:22:ARG:H	1.73	0.52
1:D:28:VAL:O	1:D:35:ALA:HB2	2.10	0.52
1:D:49:ILE:O	1:D:53:VAL:HG23	2.10	0.51
1:A:153:ILE:O	1:A:157:ILE:HG13	2.09	0.51
1:D:254:VAL:O	1:D:258:ILE:HG22	2.10	0.51
1:A:20:THR:HG22	1:A:21:GLN:N	2.26	0.51
1:A:264:LEU:O	1:A:268:THR:HG22	2.10	0.51
1:D:29:LEU:C	1:D:31:ASP:H	2.14	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:ALA:O	1:A:132:PRO:HG2	2.11	0.51
1:D:20:THR:O	1:D:24:LEU:HD22	2.11	0.51
1:B:193:ILE:C	1:B:193:ILE:HD12	2.31	0.51
1:B:71:THR:HG21	1:B:264:LEU:HD13	1.92	0.50
1:B:65:LYS:O	1:B:69:VAL:HG23	2.11	0.50
1:D:136:LEU:HD22	1:D:140:LEU:HD11	1.93	0.50
1:D:61:ASP:OD2	1:D:275:SER:HA	2.12	0.50
1:B:109:LEU:O	1:B:112:GLU:HG2	2.11	0.50
1:A:188:ILE:HG13	1:A:189:VAL:N	2.28	0.49
1:D:131:THR:O	1:D:135:LEU:HD13	2.13	0.49
1:B:207:THR:HG22	1:B:266:TYR:OH	2.13	0.49
1:A:232:GLY:HA2	1:A:239:LEU:HB2	1.95	0.49
1:D:100:HIS:CD2	1:D:102:ALA:HB3	2.48	0.49
1:B:38:LEU:HD12	1:B:82:TYR:CD1	2.48	0.49
1:D:81:SER:HB2	1:D:126:THR:HG21	1.95	0.49
1:D:267:LEU:HD11	1:D:274:VAL:HG21	1.94	0.49
1:B:157:ILE:O	1:B:161:VAL:HG23	2.12	0.48
1:B:101:PHE:CD2	1:B:236:VAL:HG23	2.48	0.48
1:A:38:LEU:HD23	1:A:86:ALA:HA	1.95	0.48
1:B:188:ILE:HG13	1:B:189:VAL:N	2.28	0.48
1:D:59:LEU:HD21	1:D:268:THR:HG22	1.96	0.48
1:B:108:MET:HA	1:B:112:GLU:O	2.14	0.48
1:B:234:GLU:HB2	5:B:509:HOH:O	2.14	0.47
1:D:106:SER:HB3	1:D:113:GLU:HG2	1.96	0.47
1:B:241:VAL:HA	1:B:244:THR:HG22	1.96	0.47
1:B:76:VAL:O	1:B:79:ILE:HG22	2.14	0.47
1:A:275:SER:O	1:A:276:GLY:C	2.53	0.47
1:A:33:LEU:HD22	1:A:246:TRP:HD1	1.80	0.47
1:D:23:GLU:HB3	1:D:109:LEU:HB3	1.96	0.47
1:D:29:LEU:C	1:D:31:ASP:N	2.69	0.47
1:D:63:ARG:HH11	1:D:63:ARG:HG2	1.80	0.47
1:D:145:ASN:C	1:D:145:ASN:ND2	2.67	0.47
1:D:248:TYR:HA	1:D:251:LEU:HD12	1.96	0.47
1:B:144:SER:OG	1:B:149:LEU:HD22	2.15	0.47
1:D:226:PRO:HA	2:D:292:RET:H172	1.97	0.47
1:D:231:LEU:HB3	1:D:239:LEU:HD13	1.96	0.46
1:A:20:THR:CG2	1:A:21:GLN:N	2.78	0.46
2:D:292:RET:H7	2:D:292:RET:H181	1.78	0.46
1:D:131:THR:HB	1:D:222:TRP:CH2	2.50	0.46
1:B:38:LEU:HD12	1:B:82:TYR:HD1	1.81	0.46
1:A:33:LEU:HD22	1:A:246:TRP:CD1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:ALA:HB2	5:D:518:HOH:O	2.16	0.46
1:D:247:ALA:O	1:D:251:LEU:HG	2.15	0.46
1:B:131:THR:OG1	1:B:132:PRO:HD3	2.16	0.46
1:D:145:ASN:ND2	1:D:148:LYS:HG3	2.26	0.46
1:D:163:GLY:O	1:D:166:ALA:HB3	2.15	0.46
1:D:93:VAL:O	1:D:93:VAL:HG13	2.16	0.45
1:B:109:LEU:HD22	1:B:109:LEU:N	2.30	0.45
1:D:241:VAL:HA	1:D:244:THR:OG1	2.15	0.45
1:D:128:ALA:O	1:D:132:PRO:HG2	2.17	0.45
1:A:29:LEU:O	1:A:30:ASN:CB	2.65	0.45
1:D:267:LEU:CD1	1:D:274:VAL:HG21	2.46	0.45
1:A:19:VAL:HG23	1:A:19:VAL:O	2.16	0.45
1:A:49:ILE:O	1:A:53:VAL:HG23	2.16	0.45
1:D:253:ILE:HA	1:D:257:TYR:CD2	2.52	0.45
1:D:109:LEU:N	1:D:109:LEU:HD22	2.32	0.44
1:D:107:VAL:O	1:D:108:MET:HB3	2.17	0.44
1:B:29:LEU:C	1:B:31:ASP:H	2.20	0.44
1:B:188:ILE:HA	1:B:191:LEU:HD12	1.99	0.44
1:A:131:THR:N	1:A:132:PRO:CD	2.80	0.44
1:A:65:LYS:HE2	3:A:300:22B:H243	2.00	0.44
1:D:136:LEU:HD22	1:D:140:LEU:CD1	2.48	0.44
1:D:249:SER:O	1:D:253:ILE:HG13	2.18	0.44
1:B:259:PHE:C	1:B:259:PHE:CD1	2.91	0.44
1:A:238:VAL:CG2	1:A:239:LEU:HD13	2.33	0.44
1:B:119:THR:HG23	1:B:234:GLU:OE1	2.17	0.44
1:B:28:VAL:C	1:B:29:LEU:O	2.56	0.43
2:B:292:RET:H181	2:B:292:RET:H7	1.84	0.43
1:A:44:LEU:HD23	1:A:44:LEU:HA	1.84	0.43
1:A:268:THR:HG23	5:A:543:HOH:O	2.19	0.43
1:A:258:ILE:HG13	1:A:259:PHE:N	2.33	0.43
1:A:252:ASP:O	1:A:256:LYS:HB2	2.18	0.43
1:D:72:ILE:O	1:D:75:PRO:HD2	2.18	0.43
2:A:292:RET:H7	2:A:292:RET:H181	1.84	0.43
1:B:241:VAL:HA	1:B:244:THR:CG2	2.49	0.43
1:D:100:HIS:CE1	5:D:512:HOH:O	2.72	0.42
1:A:106:SER:HB3	1:A:113:GLU:HG3	2.02	0.42
1:D:225:TYR:N	1:D:226:PRO:HD2	2.34	0.42
1:A:141:LEU:HD22	1:A:263:LEU:CD1	2.50	0.42
1:D:134:ILE:HG23	1:D:259:PHE:CZ	2.55	0.42
1:B:236:VAL:HG22	1:B:236:VAL:O	2.19	0.42
1:D:131:THR:HG1	1:D:132:PRO:HD3	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:THR:HG22	1:A:22:ARG:N	2.30	0.42
1:D:144:SER:OG	1:D:198:TRP:NE1	2.52	0.42
1:B:19:VAL:HG12	1:B:24:LEU:HD22	2.02	0.42
1:A:220:VAL:HG12	1:A:221:MET:CE	2.50	0.42
1:A:215:LYS:HB3	1:A:215:LYS:HE2	1.89	0.42
1:A:65:LYS:O	1:A:69:VAL:HG23	2.21	0.41
1:A:239:LEU:HB3	1:A:244:THR:HG22	2.02	0.41
1:B:268:THR:HG23	5:B:516:HOH:O	2.19	0.41
1:B:29:LEU:O	1:B:30:ASN:CB	2.69	0.41
1:A:131:THR:OG1	1:A:132:PRO:HD3	2.20	0.41
1:B:267:LEU:CD1	1:B:274:VAL:HG21	2.51	0.41
1:D:149:LEU:HD13	1:D:149:LEU:HA	1.94	0.41
1:D:173:HIS:CD2	1:D:177:TRP:HE1	2.39	0.41
1:A:231:LEU:O	1:A:238:VAL:HG22	2.20	0.41
1:D:145:ASN:HD21	1:D:148:LYS:H	1.68	0.41
1:D:81:SER:CB	1:D:126:THR:HG21	2.50	0.41
1:D:232:GLY:CA	1:D:244:THR:HG21	2.50	0.41
1:D:41:ASN:HD22	1:D:41:ASN:HA	1.61	0.41
1:A:240:PRO:O	1:A:244:THR:HG23	2.21	0.40
1:D:231:LEU:HA	1:D:236:VAL:HG12	2.04	0.40
1:B:20:THR:HG22	1:B:23:GLU:N	2.29	0.40
1:D:40:ILE:O	1:D:44:LEU:HG	2.22	0.40
1:A:210:ILE:O	1:A:214:LEU:HG	2.21	0.40
1:B:193:ILE:HD13	1:B:198:TRP:CE2	2.55	0.40
1:B:144:SER:HG	1:B:149:LEU:HD22	1.85	0.40
1:B:101:PHE:CE2	1:B:236:VAL:HG23	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:600:HOH:O	5:A:600:HOH:O[2_656]	1.55	0.65

## 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	259/291 (89%)	249 (96%)	9 (4%)	1 (0%)	39	42
1	B	257/291 (88%)	248 (96%)	8 (3%)	1 (0%)	39	42
1	D	257/291 (88%)	247 (96%)	9 (4%)	1 (0%)	39	42
All	All	773/873 (88%)	744 (96%)	26 (3%)	3 (0%)	39	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	GLY
1	D	209	ASP
1	B	108	MET

### 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	208/233 (89%)	194 (93%)	14 (7%)	20	21
1	B	207/233 (89%)	193 (93%)	14 (7%)	20	21
1	D	207/233 (89%)	198 (96%)	9 (4%)	35	43
All	All	622/699 (89%)	585 (94%)	37 (6%)	24	27

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	24	LEU
1	A	59	LEU
1	A	66	LEU
1	A	135	LEU
1	A	136	LEU
1	A	149	LEU

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Mol	Chain	Res	Type
1	A	168	LEU
1	A	172	SER
1	A	216	LEU
1	A	233	VAL
1	A	234	GLU
1	A	258	ILE
1	A	264	LEU
1	B	20	THR
1	B	24	LEU
1	B	38	LEU
1	B	66	LEU
1	B	108	MET
1	B	135	LEU
1	B	136	LEU
1	B	149	LEU
1	B	168	LEU
1	B	174	LEU
1	B	234	GLU
1	B	258	ILE
1	B	259	PHE
1	B	264	LEU
1	D	38	LEU
1	D	66	LEU
1	D	79	ILE
1	D	136	LEU
1	D	145	ASN
1	D	149	LEU
1	D	172	SER
1	D	209	ASP
1	D	264	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	B	41	ASN
1	B	265	ASN
1	D	41	ASN
1	D	100	HIS
1	D	145	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 for Mogul analysis.

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	292	1	19,20,21	2.16	6 (31%)	27,27,28	1.65	7 (25%)
3	22B	A	300	-	24,26,53	3.60	10 (41%)	28,35,72	1.98	9 (32%)
2	RET	B	292	1	19,20,21	2.05	4 (21%)	27,27,28	1.54	5 (18%)
3	22B	B	300	-	24,26,53	3.69	10 (41%)	28,35,72	1.99	8 (28%)
2	RET	D	292	1	19,20,21	1.93	4 (21%)	27,27,28	1.71	8 (29%)

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	292	1	-	0/13/30/31	0/1/1/1
3	22B	A	300	-	-	0/31/31/65	0/0/0/0
2	RET	B	292	1	-	0/13/30/31	0/1/1/1
3	22B	B	300	-	-	0/31/31/65	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	RET	D	292	1	-	0/13/30/31	0/1/1/1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	292	RET	C2-C3	-2.40	1.46	1.52
2	B	292	RET	C2-C3	-2.33	1.46	1.52
2	D	292	RET	C2-C3	-2.28	1.46	1.52
3	A	300	22B	C4-C3	2.10	1.37	1.32
3	B	300	22B	C4-C3	2.11	1.37	1.32
2	A	292	RET	C10-C9	2.23	1.38	1.35
3	A	300	22B	C18-C5	2.39	1.56	1.50
2	A	292	RET	C18-C5	2.46	1.55	1.51
2	B	292	RET	C7-C6	2.52	1.55	1.45
2	A	292	RET	C7-C6	2.54	1.55	1.45
3	B	300	22B	C18-C5	2.59	1.56	1.50
2	D	292	RET	C7-C6	2.69	1.55	1.45
3	B	300	22B	C17-C1	3.05	1.57	1.52
3	B	300	22B	C20-C13	3.07	1.57	1.50
3	A	300	22B	C20-C13	3.09	1.57	1.50
3	A	300	22B	C17-C1	3.44	1.58	1.52
3	A	300	22B	O27-C23	3.47	1.54	1.44
3	B	300	22B	O27-C23	3.73	1.54	1.44
3	B	300	22B	C16-C1	3.90	1.59	1.52
3	A	300	22B	C19-C9	3.91	1.59	1.50
3	B	300	22B	C19-C9	3.92	1.59	1.50
3	A	300	22B	C16-C1	3.98	1.59	1.52
2	D	292	RET	C5-C6	4.12	1.40	1.34
3	A	300	22B	C22-C21	4.27	1.62	1.53
2	B	292	RET	C5-C6	4.34	1.41	1.34
3	B	300	22B	C22-C21	4.64	1.63	1.53
2	A	292	RET	C5-C6	4.68	1.41	1.34
3	A	300	22B	C21-C2	5.38	1.61	1.54
2	D	292	RET	C1-C6	5.44	1.61	1.53
3	B	300	22B	C21-C2	5.50	1.61	1.54
2	A	292	RET	C1-C6	5.87	1.62	1.53
2	B	292	RET	C1-C6	6.03	1.62	1.53
3	A	300	22B	C2-C3	13.08	1.64	1.50
3	B	300	22B	C2-C3	13.41	1.64	1.50

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	300	22B	C17-C1-C16	-3.61	105.07	110.51
3	A	300	22B	C17-C1-C16	-3.43	105.34	110.51
3	B	300	22B	C4-C5-C6	-3.37	113.55	118.98
3	A	300	22B	C4-C5-C6	-3.34	113.60	118.98
2	D	292	RET	C1-C6-C5	-2.85	118.48	122.66
2	A	292	RET	C8-C9-C10	-2.76	114.54	118.98
2	A	292	RET	C1-C6-C5	-2.74	118.64	122.66
3	A	300	22B	C12-C13-C14	-2.61	110.42	118.92
2	B	292	RET	C1-C6-C5	-2.60	118.84	122.66
3	B	300	22B	C12-C13-C14	-2.33	111.34	118.92
2	B	292	RET	C8-C9-C10	-2.31	115.26	118.98
2	D	292	RET	C7-C6-C5	-2.31	116.09	121.37
3	A	300	22B	C8-C9-C10	-2.08	115.63	118.98
2	D	292	RET	C8-C9-C10	-2.03	115.72	118.98
2	B	292	RET	C7-C6-C5	-2.02	116.76	121.37
2	D	292	RET	C19-C9-C8	2.09	121.57	118.10
2	D	292	RET	C2-C1-C6	2.11	113.71	110.36
2	B	292	RET	C2-C1-C6	2.25	113.93	110.36
3	B	300	22B	C24-C23-C22	2.30	118.42	111.59
2	A	292	RET	C2-C1-C6	2.35	114.09	110.36
2	A	292	RET	C17-C1-C6	2.37	114.01	110.30
3	A	300	22B	C24-C23-C22	2.41	118.75	111.59
2	D	292	RET	C20-C13-C12	2.51	122.28	118.10
3	B	300	22B	C18-C5-C4	2.65	122.51	118.10
2	A	292	RET	C19-C9-C8	2.67	122.54	118.10
3	A	300	22B	C18-C5-C4	2.71	122.61	118.10
2	D	292	RET	C16-C1-C6	2.77	114.64	110.30
2	A	292	RET	C7-C8-C9	2.99	130.76	126.22
2	A	292	RET	C1-C6-C7	3.03	124.31	115.82
2	B	292	RET	C1-C6-C7	3.07	124.41	115.82
2	D	292	RET	C1-C6-C7	3.43	125.44	115.82
3	A	300	22B	C16-C1-C2	3.50	116.88	111.59
3	B	300	22B	C20-C13-C12	3.83	124.47	118.10
3	B	300	22B	C16-C1-C2	4.17	117.89	111.59
3	A	300	22B	C21-C22-C23	4.22	124.83	116.36
3	A	300	22B	C20-C13-C12	4.23	125.14	118.10
3	B	300	22B	C21-C22-C23	4.35	125.09	116.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	292	RET	1	0
3	A	300	22B	2	0
2	B	292	RET	1	0
3	B	300	22B	1	0
2	D	292	RET	2	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	261/291 (89%)	0.13	15 (5%) 27 27	20, 33, 59, 79	0
1	B	259/291 (89%)	0.10	9 (3%) 48 46	25, 38, 52, 66	0
1	D	259/291 (89%)	0.81	36 (13%) 4 3	32, 50, 71, 92	0
All	All	779/873 (89%)	0.35	60 (7%) 16 16	20, 40, 65, 92	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	GLY	7.1
1	D	276	GLY	7.1
1	A	17	ALA	6.7
1	D	111	GLY	6.6
1	D	109	LEU	6.1
1	A	110	GLY	5.6
1	A	111	GLY	4.8
1	D	112	GLU	4.7
1	D	243	TYR	4.7
1	D	110	GLY	4.4
1	D	241	VAL	4.1
1	D	29	LEU	4.1
1	D	269	SER	3.9
1	B	200	GLN	3.7
1	B	112	GLU	3.6
1	D	32	PRO	3.5
1	D	30	ASN	3.5
1	D	113	GLU	3.4
1	B	104	GLY	3.4
1	A	276	GLY	3.3
1	A	109	LEU	3.3
1	A	107	VAL	3.3
1	D	205	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	268	THR	3.1
1	D	275	SER	3.1
1	A	108	MET	3.1
1	A	30	ASN	3.1
1	D	273	VAL	3.0
1	D	193	ILE	3.0
1	D	59	LEU	3.0
1	D	216	LEU	3.0
1	B	98	ALA	2.9
1	D	240	PRO	2.7
1	A	112	GLU	2.7
1	A	57	ARG	2.7
1	D	206	GLY	2.7
1	D	265	ASN	2.6
1	B	243	TYR	2.6
1	D	266	TYR	2.6
1	D	107	VAL	2.5
1	D	272	GLY	2.5
1	A	236	VAL	2.5
1	A	54	PHE	2.5
1	B	203	LYS	2.4
1	D	114	VAL	2.4
1	D	261	PHE	2.4
1	B	99	GLY	2.4
1	D	229	TRP	2.4
1	D	210	ILE	2.3
1	A	277	SER	2.3
1	D	20	THR	2.2
1	A	18	GLU	2.2
1	A	104	GLY	2.2
1	D	211	PHE	2.1
1	D	238	VAL	2.1
1	D	34	LEU	2.1
1	B	209	ASP	2.0
1	D	108	MET	2.0
1	D	209	ASP	2.0
1	D	25	PHE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	22B	B	300	27/54	0.79	0.21	5.72	43,52,60,61	0
3	22B	A	300	27/54	0.61	0.32	2.88	72,77,85,86	0
2	RET	D	292	20/21	0.84	0.19	1.06	40,42,44,46	0
2	RET	A	292	20/21	0.96	0.14	0.60	19,23,27,28	0
4	BR	B	401	1/1	1.00	0.12	-0.29	32,32,32,32	0
2	RET	B	292	20/21	0.96	0.11	-0.47	25,28,30,30	0
4	BR	A	401	1/1	1.00	0.11	-1.41	28,28,28,28	0
4	BR	D	401	1/1	0.98	0.10	-2.36	50,50,50,50	0
4	BR	A	293	1/1	0.95	0.08	-2.72	58,58,58,58	0
4	BR	A	402	1/1	0.99	0.07	-	43,43,43,43	0
4	BR	D	402	1/1	0.98	0.04	-	58,58,58,58	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.