



Full wwPDB EM Validation Report ⓘ

Sep 21, 2025 – 12:28 AM JST

PDB ID : 9UT9 / pdb_00009ut9
EMDB ID : EMD-64485
Title : The VFT domains of human sweet taste receptor TAS1R2 and TAS1R3 in the apo state
Authors : Shi, Z.J.; Xu, W.X.; Yue, X.L.; Wu, L.J.; Hua, T.; Liu, Z.J.
Deposited on : 2025-05-03
Resolution : 3.18 Å(reported)

This is a Full wwPDB EM 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/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:

EMDB validation analysis : **FAILED**
MolProbity : 4-5-2 with Phenix2.0rc1
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

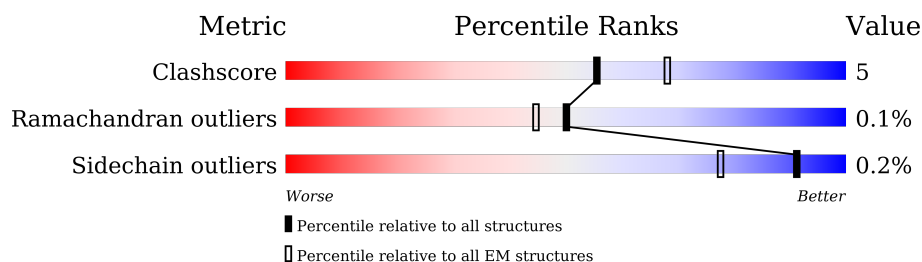
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.18 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1078	
2	B	1130	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7958 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 Taste receptor type 1 member 2, Engineered red fluorescent protein mScarlet3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	490	Total	C	N	O	S	0	0
			3962	2541	665	733	23		

There are 29 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MET	-	initiating methionine	UNP Q8TE23
A	-2	LYS	-	expression tag	UNP Q8TE23
A	-1	THR	-	expression tag	UNP Q8TE23
A	0	ILE	-	expression tag	UNP Q8TE23
A	1	ILE	-	expression tag	UNP Q8TE23
A	2	ALA	-	expression tag	UNP Q8TE23
A	3	LEU	-	expression tag	UNP Q8TE23
A	4	SER	-	expression tag	UNP Q8TE23
A	5	TYR	-	expression tag	UNP Q8TE23
A	6	ILE	-	expression tag	UNP Q8TE23
A	7	PHE	-	expression tag	UNP Q8TE23
A	8	CYS	-	expression tag	UNP Q8TE23
A	9	LEU	-	expression tag	UNP Q8TE23
A	10	VAL	-	expression tag	UNP Q8TE23
A	11	PHE	-	expression tag	UNP Q8TE23
A	12	ALA	-	expression tag	UNP Q8TE23
A	13	GLY	-	expression tag	UNP Q8TE23
A	14	SER	-	expression tag	UNP Q8TE23
A	15	ASP	-	expression tag	UNP Q8TE23
A	16	TYR	-	expression tag	UNP Q8TE23
A	17	LYS	-	expression tag	UNP Q8TE23
A	18	ASP	-	expression tag	UNP Q8TE23
A	19	ASP	-	expression tag	UNP Q8TE23
A	20	ASP	-	expression tag	UNP Q8TE23
A	21	ASP	-	expression tag	UNP Q8TE23
A	22	LYS	-	expression tag	UNP Q8TE23
A	23	GLY	-	expression tag	UNP Q8TE23

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Chain	Residue	Modelled	Actual	Comment	Reference
A	24	SER	-	expression tag	UNP Q8TE23
A	25	ALA	-	expression tag	UNP Q8TE23

- Molecule 2 is a protein called Taste receptor type 1 member 3,mNeonGreen.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	514	Total	C	N	O	S	0	0
			3996	2526	704	733	33		

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	initiating methionine	UNP Q7RTX0
B	-6	LYS	-	expression tag	UNP Q7RTX0
B	-5	THR	-	expression tag	UNP Q7RTX0
B	-4	ILE	-	expression tag	UNP Q7RTX0
B	-3	ILE	-	expression tag	UNP Q7RTX0
B	-2	ALA	-	expression tag	UNP Q7RTX0
B	-1	LEU	-	expression tag	UNP Q7RTX0
B	0	SER	-	expression tag	UNP Q7RTX0
B	1	TYR	-	expression tag	UNP Q7RTX0
B	2	ILE	-	expression tag	UNP Q7RTX0
B	3	PHE	-	expression tag	UNP Q7RTX0
B	4	CYS	-	expression tag	UNP Q7RTX0
B	5	LEU	-	expression tag	UNP Q7RTX0
B	6	VAL	-	expression tag	UNP Q7RTX0
B	7	PHE	-	expression tag	UNP Q7RTX0
B	8	ALA	-	expression tag	UNP Q7RTX0
B	9	GLY	-	expression tag	UNP Q7RTX0
B	10	SER	-	expression tag	UNP Q7RTX0
B	11	ASP	-	expression tag	UNP Q7RTX0
B	12	TYR	-	expression tag	UNP Q7RTX0
B	13	LYS	-	expression tag	UNP Q7RTX0
B	14	ASP	-	expression tag	UNP Q7RTX0
B	15	ASP	-	expression tag	UNP Q7RTX0
B	16	ASP	-	expression tag	UNP Q7RTX0
B	17	ASP	-	expression tag	UNP Q7RTX0
B	18	LYS	-	expression tag	UNP Q7RTX0
B	19	GLY	-	expression tag	UNP Q7RTX0
B	20	SER	-	expression tag	UNP Q7RTX0
B	853	GLY	-	linker	UNP Q7RTX0
B	854	SER	-	linker	UNP Q7RTX0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	855	SER	-	linker	UNP Q7RTX0
B	856	GLY	-	linker	UNP Q7RTX0
B	857	LEU	-	linker	UNP Q7RTX0
B	858	GLU	-	linker	UNP Q7RTX0
B	859	VAL	-	linker	UNP Q7RTX0
B	860	LEU	-	linker	UNP Q7RTX0
B	861	PHE	-	linker	UNP Q7RTX0
B	862	GLN	-	linker	UNP Q7RTX0
B	863	GLY	-	linker	UNP Q7RTX0
B	864	PRO	-	linker	UNP Q7RTX0
B	865	SER	-	linker	UNP Q7RTX0
B	866	GLY	-	linker	UNP Q7RTX0
B	867	GLY	-	linker	UNP Q7RTX0
B	1103	GLY	-	expression tag	UNP A0A1S4NYF2
B	1104	SER	-	expression tag	UNP A0A1S4NYF2
B	1105	GLU	-	expression tag	UNP A0A1S4NYF2
B	1106	ASN	-	expression tag	UNP A0A1S4NYF2
B	1107	LEU	-	expression tag	UNP A0A1S4NYF2
B	1108	TYR	-	expression tag	UNP A0A1S4NYF2
B	1109	PHE	-	expression tag	UNP A0A1S4NYF2
B	1110	GLN	-	expression tag	UNP A0A1S4NYF2
B	1111	SER	-	expression tag	UNP A0A1S4NYF2
B	1112	SER	-	expression tag	UNP A0A1S4NYF2
B	1113	GLY	-	expression tag	UNP A0A1S4NYF2
B	1114	HIS	-	expression tag	UNP A0A1S4NYF2
B	1115	HIS	-	expression tag	UNP A0A1S4NYF2
B	1116	HIS	-	expression tag	UNP A0A1S4NYF2
B	1117	HIS	-	expression tag	UNP A0A1S4NYF2
B	1118	HIS	-	expression tag	UNP A0A1S4NYF2
B	1119	HIS	-	expression tag	UNP A0A1S4NYF2
B	1120	HIS	-	expression tag	UNP A0A1S4NYF2
B	1121	HIS	-	expression tag	UNP A0A1S4NYF2
B	1122	HIS	-	expression tag	UNP A0A1S4NYF2

PHE	ARG	THR	PHE	SER	LEU	ALA	ASP	THR	PHE	M429	L139	MET
LYS	LYS	ALA	ALA	ASN	GLY	THR	THR	PHE	HIS	F437	F157	THR
THR	THR	TRP	VAL	GLY	PRO	THR	MET	GLN	HIS			ILE
LEU	LEU	CYS	VAL	ASP	GLY	TRP	PRO	ALA	ASP	D455	F160	ILE
LYS	HIS	ARG	ASP	PHE	ASP	VAL	THR	ALA	SER	L486	A248	ALA
SER	SER	LYS	SER	MET	ALA	SER	GLU	GLU	PRO	K457	ASP	LEU
LYS	LYS	TRP	GLY	VAL	GLN	PHE	ALA	ILE	LEU		ASP	THR
THR	THR	THR	GLN	VAL	GLN	VAL	VAL	VAL	VAL	H469	SER	ILE
GLU	GLU	PRO	GLN	GLN	ASN	LEU	HIS	GLU	ALA	D470	ARG	CYS
LEU	LEU	ASP	ASP	GLY	ASP	LEU	CYS	SER	SER	V471	L253	LEU
ASN	ASN	ASN	GLY	THR	GLY	ALA	ARG	GLU	GLY	D491	L260	VAL
PHE	PHE	ASP	ARG	THR	ASN	VAL	THR	LEU	GLY			PHE
LYS	LYS	LYS	THR	ASN	THR	VAL	ARG	PRO	PRO	H515	S267	ALA
THR	THR	THR	MET	PRO	GLY	GLN	SER	LEU	LEU			GLY
ILE	ILE	ILE	GLN	ASN	ASN	VAL	TRP	SER	ALA	Y529	L283	ASP
GLN	GLN	ILE	PHE	ASP	GLN	VAL	VAL	TRP	CYS			SER
LYS	LYS	SER	GLU	GLY	GLY	LEU	VAL	ALA	ALA	I536	I288	THR
ALA	ALA	THR	ASP	TYR	LYS	ARG	PHE	ASP	PHE			LYS
PHE	PHE	PHE	GLY	GLU	HIS	PRO	GLY	ARG	GLY	T539	L312	ASP
THR	THR	LYS	ALA	GLU	GLY	ALA	LEU	LEU	VAL			ASP
TRP	ASP	TRP	SER	LEU	GLY	VAL	ALA	SER	CYS	W546	T320	ASP
VAL	VAL	SER	ASN	ASN	SER	GLN	HIS	GLY	LEU		V321	LYS
MET	MET	TYR	THR	LEU	SER	MET	ALA	CYS	GLY	F555	L322	GLY
GLY	GLY	THR	VAL	LYS	GLY	GLY	THR	LEU	LEU	R556		SER
MET	MET	THR	ASN	SER	LEU	ALA	ALA	ARG	CYS	ARG	R327	ALA
GLY	GLY	GLY	THR	THR	GLU	LEU	ALA	GLY	VAL	ARG	Q330	PRO
ASN	ASN	ASN	ARG	LYS	VAL	LEU	THR	PRO	THR	SER		LEU
GLY	GLY	GLY	THR	GLY	LEU	LEU	LEU	TRP	SER	ARG		LEU
LYS	LYS	LYS	THR	ASP	PHE	CYS	ALA	ALA	VAL	PHE	P335	LEU
ARG	ARG	THR	THR	LEU	GLN	VAL	PHE	TRP	LEU	LEU	V338	ALA
THR	THR	THR	GLU	GLN	GLY	LEU	LEU	VAL	LEU	TRP	K339	GLY
ARG	SER	ARG	GLY	PHE	PRO	GLY	CYS	VAL	PHE	GLY		GLU
SER	GLU	SER	SER	SER	SER	ILE	PHE	VAL	PRO	GLY	L344	ALA
ASN	ASN	THR	HIS	PRO	GLY	LEU	LEU	LEU	GLY	GLY		LEU
THR	THR	ALA	ILE	TRP	GLY	ALA	GLY	ALA	GLN	PRO	E356	LEU
ARG	ARG	THR	GLY	ILE	VAL	ALA	THR	ALA	PRO	ALA		LEU
THR	THR	THR	GLY	LEU	SER	PHE	PHE	MET	SER	VAL	ARG	GLY
GLN	GLN	THR	GLU	VAL	LYS	HIS	LEU	VAL	PRO	LEU	S53	S53
SER	SER	TYR	ALA	PRO	GLY	LEU	VAL	VAL	ALA	LEU	R54	GLY
THR	THR	THR	GLN	ILE	GLU	PRO	ARG	GLU	ARG	LEU		GLY
PHE	GLY	PHE	VAL	ILE	GLU	ARG	SER	VAL	CYS	LEU	S58	LEU
HIS	HIS	ALA	LYS	GLY	ASP	GLN	GLN	VAL	LEU	LEU		GLU
HIS	HIS	LYS	GLY	TYR	ASN	TYR	PRO	LEU	ALA	LEU	K36	GLU
HIS	HIS	PRO	THR	GLY	MET	LEU	GLY	CYS	GLN	LEU	ASP	ASP
HIS	HIS	MET	GLY	PHE	ALA	LEU	CYS	THR	GLN	SER	VAL	VAL
HIS	HIS	ALA	PHE	HIS	SER	MET	TYR	TRP	PRO	LEU	VAL	VAL
HIS	HIS	ALA	PRO	GLN	LEU	ARG	ASN	TYR	LEU	ALA	GLY	T102

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	189062	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.16	0/4062	0.31	0/5534
2	B	0.18	0/4092	0.35	0/5551
All	All	0.17	0/8154	0.33	0/11085

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	3962	0	3854	41	0
2	B	3996	0	3883	36	0
All	All	7958	0	7737	76	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 5.

All (76) 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:284:PHE:O	1:A:288:VAL:HG23	1.86	0.76
1:A:88:LEU:O	1:A:400:HIS:ND1	2.19	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:TYR:HD1	1:A:165:SER:HG	1.38	0.72
1:A:243:LEU:HD21	1:A:253:GLU:OE1	1.90	0.71
1:A:243:LEU:HD21	1:A:253:GLU:CD	2.20	0.67
2:B:546:TRP:N	2:B:555:PHE:O	2.29	0.66
1:A:100:ASP:OD1	1:A:101:VAL:N	2.29	0.65
1:A:89:LEU:HD22	1:A:92:VAL:HB	1.83	0.61
2:B:260:LEU:HD21	2:B:283:LEU:HD11	1.84	0.60
2:B:98:ASP:OD1	2:B:369:ARG:NH2	2.36	0.57
1:A:183:THR:HA	1:A:444:VAL:HG22	1.85	0.57
2:B:139:LEU:HD21	2:B:426:LEU:CD1	2.36	0.55
1:A:33:TYR:C	1:A:34:LEU:HD23	2.31	0.55
1:A:192:GLU:OE1	1:A:226:ARG:NH1	2.38	0.55
1:A:278:ASP:N	1:A:278:ASP:OD1	2.40	0.54
1:A:307:ASP:O	1:A:307:ASP:OD1	2.25	0.54
2:B:131:TYR:O	2:B:132:THR:HB	2.07	0.54
2:B:330:GLN:N	2:B:330:GLN:OE1	2.41	0.54
1:A:85:ASP:OD1	1:A:86:SER:N	2.40	0.54
2:B:536:ILE:HG13	2:B:536:ILE:O	2.08	0.53
1:A:263:LYS:NZ	2:B:515:HIS:O	2.38	0.53
2:B:267:SER:O	2:B:267:SER:OG	2.28	0.51
2:B:407:THR:OG1	2:B:429:ASN:OD1	2.20	0.51
2:B:54:ARG:NH2	2:B:58:SER:O	2.44	0.51
1:A:265:GLN:HG2	1:A:265:GLN:O	2.11	0.50
1:A:439:ASP:OD1	1:A:443:ASP:N	2.44	0.50
2:B:139:LEU:HD21	2:B:426:LEU:HD11	1.94	0.50
1:A:315:GLU:HA	1:A:315:GLU:OE2	2.12	0.49
2:B:344:LEU:HD12	2:B:344:LEU:O	2.12	0.49
2:B:132:THR:HG22	2:B:132:THR:O	2.13	0.49
1:A:67:ILE:O	1:A:70:ASN:OD1	2.30	0.49
2:B:411:ASN:OD1	2:B:411:ASN:C	2.55	0.49
2:B:123:ARG:CG	2:B:123:ARG:O	2.61	0.49
1:A:167:ILE:HD11	1:A:302:GLU:OE2	2.13	0.48
2:B:117:LEU:HD13	2:B:160:PHE:CD1	2.47	0.48
2:B:529:TYR:CD1	2:B:539:THR:HG23	2.48	0.48
1:A:552:PHE:O	1:A:553:LYS:C	2.55	0.48
2:B:335:PRO:O	2:B:339:LYS:HG2	2.13	0.48
1:A:227:VAL:HG13	1:A:232:ILE:HB	1.95	0.48
1:A:311:HIS:ND1	1:A:311:HIS:O	2.47	0.47
2:B:86:LYS:HB2	2:B:86:LYS:NZ	2.30	0.46
1:A:35:LEU:HD21	1:A:395:VAL:HG11	1.97	0.46
1:A:140:GLY:N	1:A:162:ILE:O	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:139:LEU:HD11	2:B:404:LEU:HD13	1.98	0.46
1:A:474:GLN:OE1	1:A:476:LYS:NZ	2.41	0.46
1:A:411:THR:O	1:A:412:LYS:HB2	2.15	0.46
1:A:242:THR:O	1:A:242:THR:HG23	2.16	0.46
1:A:521:LEU:O	1:A:524:THR:OG1	2.31	0.45
2:B:529:TYR:CD1	2:B:529:TYR:N	2.83	0.45
2:B:456:LEU:HD23	2:B:457:LYS:N	2.31	0.45
2:B:322:LEU:HD22	2:B:456:LEU:HD21	1.97	0.45
1:A:307:ASP:OD1	1:A:307:ASP:C	2.60	0.44
1:A:262:ASP:OD1	1:A:266:GLN:NE2	2.50	0.44
2:B:117:LEU:HD11	2:B:157:PHE:HD1	1.83	0.44
2:B:437:VAL:CG1	2:B:437:VAL:O	2.66	0.44
2:B:288:ILE:CD1	2:B:312:LEU:HD23	2.47	0.44
1:A:207:ILE:HD13	1:A:264:LEU:HD23	1.99	0.43
1:A:287:GLU:O	1:A:287:GLU:OE1	2.37	0.43
1:A:426:LYS:O	1:A:426:LYS:HG3	2.18	0.43
2:B:491:ASP:OD1	2:B:491:ASP:N	2.36	0.43
2:B:102:THR:O	2:B:108:VAL:HG13	2.19	0.43
2:B:456:LEU:HD22	2:B:471:VAL:CG2	2.49	0.42
1:A:439:ASP:HB2	1:A:440:PRO:CD	2.50	0.42
2:B:320:THR:HG22	2:B:321:VAL:N	2.35	0.42
1:A:60:LYS:H	1:A:60:LYS:HD2	1.84	0.42
1:A:318:HIS:O	1:A:318:HIS:ND1	2.53	0.42
1:A:243:LEU:HD11	1:A:249:MET:HE2	2.01	0.41
1:A:386:TYR:OH	1:A:431:LEU:O	2.39	0.41
2:B:123:ARG:O	2:B:123:ARG:HG3	2.21	0.41
1:A:515:PHE:CD1	1:A:515:PHE:C	2.98	0.41
2:B:30:ARG:NH1	2:B:30:ARG:HB3	2.36	0.41
2:B:456:LEU:HD23	2:B:456:LEU:C	2.46	0.41
1:A:243:LEU:CD1	1:A:249:MET:HE2	2.51	0.41
1:A:552:PHE:O	1:A:552:PHE:CD2	2.73	0.40
2:B:335:PRO:O	2:B:338:VAL:HG12	2.21	0.40
2:B:327:ARG:HD2	2:B:455:ASP:OD1	2.21	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	484/1078 (45%)	440 (91%)	43 (9%)	1 (0%)	44	73
2	B	506/1130 (45%)	476 (94%)	30 (6%)	0	100	100
All	All	990/2208 (45%)	916 (92%)	73 (7%)	1 (0%)	50	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	LYS

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	446/955 (47%)	446 (100%)	0	100	100
2	B	433/941 (46%)	431 (100%)	2 (0%)	86	93
All	All	879/1896 (46%)	877 (100%)	2 (0%)	91	97

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	469	HIS
2	B	529	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN
1	A	143	ASN
1	A	266	GLN
1	A	286	ASN
1	A	291	GLN
1	A	397	HIS
1	A	419	GLN
2	B	193	GLN
2	B	202	GLN
2	B	317	GLN
2	B	326	GLN
2	B	341	HIS
2	B	532	ASN

5.3.3 RNA [i](#)

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](#)

There are no ligands in this entry.

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.