



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:42 AM UTC

PDB ID : 8QH2 / pdb_00008qh2
Title : Crystal structure of chimeric UAP1L1
Authors : Chen, X.; Yan, K.; Bartual, S.; van Aalten, D.
Deposited on : 2023-09-06
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

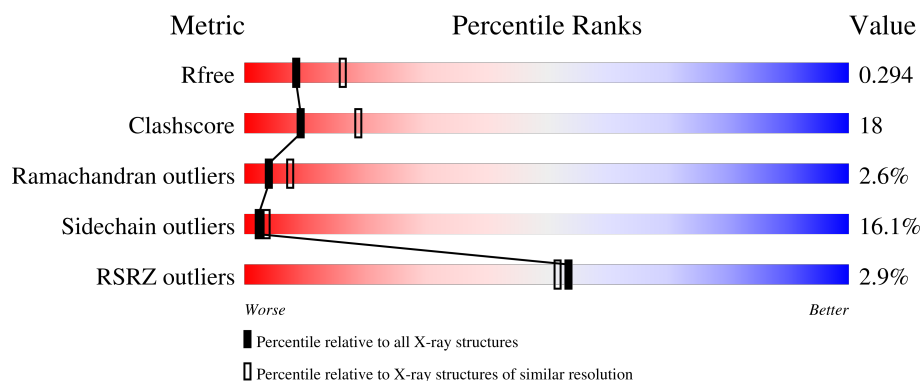
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	510	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3810 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 UDP-N-acetylhexosamine pyrophosphorylase,UDP-N-acetylhexosamine pyrophosphorylase-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	89	0	0
			3810	2437	664	693	16			

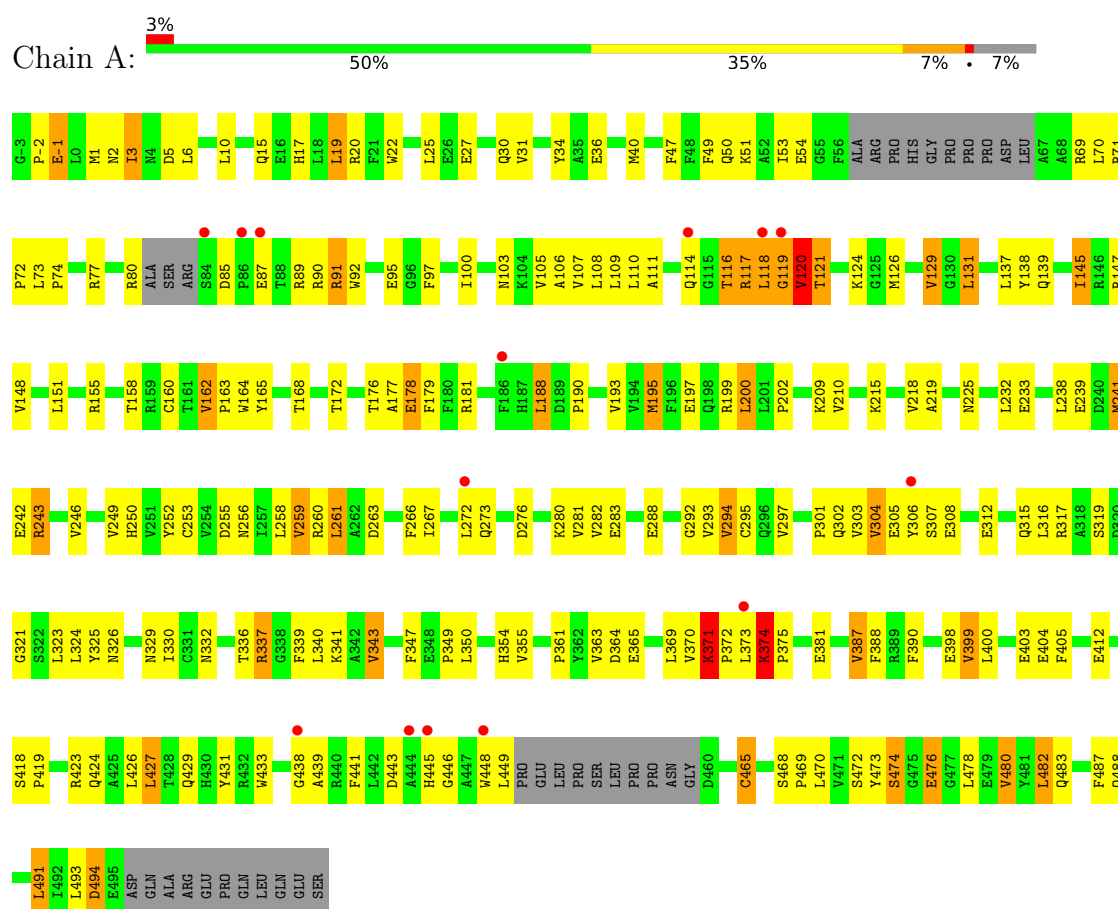
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q16222
A	-2	PRO	-	expression tag	UNP Q16222
A	-1	GLU	-	expression tag	UNP Q16222
A	0	LEU	-	expression tag	UNP Q16222

3 Residue-property plots

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 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 ($RSRZ > 2$). 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: UDP-N-acetylhexosamine pyrophosphorylase, UDP-N-acetylhexosamine pyrophosphorylase-like protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	113.52Å 113.52Å 43.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.00 – 2.75 41.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (41.00-2.75) 100.0 (41.00-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.19.2-4158-000	Depositor
R, R_{free}	0.233 , 0.283 (Not available) , 0.294	Depositor DCC
R_{free} test set	672 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	77.3	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 104.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3810	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.40	0/3898	0.62	0/5274

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	3810	0	3772	133	0
All	All	3810	0	3772	133	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 18.

All (133) 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:19:LEU:HD12	1:A:22:TRP:HB2	1.59	0.84
1:A:256:ASN:HB3	1:A:259:VAL:HG22	1.63	0.80
1:A:40:MET:HE3	1:A:202:PRO:HG3	1.69	0.75
1:A:100:ILE:HG13	1:A:105:VAL:HG21	1.69	0.74
1:A:172:THR:O	1:A:176:THR:HG22	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:TRP:O	1:A:30:GLN:NE2	2.23	0.71
1:A:250:HIS:NE2	1:A:332:ASN:OD1	2.21	0.70
1:A:176:THR:HG23	1:A:195:MET:HE1	1.72	0.70
1:A:238:LEU:HA	1:A:241:MET:HE2	1.72	0.70
1:A:110:LEU:HD13	1:A:225:ASN:HB2	1.75	0.69
1:A:304:VAL:HG23	1:A:308:GLU:HB2	1.75	0.69
1:A:323:LEU:O	1:A:326:ASN:ND2	2.27	0.67
1:A:336:THR:HG23	1:A:339:PHE:H	1.59	0.67
1:A:72:PRO:HG3	1:A:324:LEU:HD11	1.75	0.66
1:A:474:SER:N	1:A:476:GLU:OE2	2.25	0.66
1:A:73:LEU:HG	1:A:295:CYS:HA	1.76	0.66
1:A:443:ASP:N	1:A:446:GLY:O	2.20	0.66
1:A:363:VAL:HG23	1:A:369:LEU:HG	1.79	0.65
1:A:116:THR:O	1:A:118:LEU:N	2.30	0.65
1:A:258:LEU:HD23	1:A:426:LEU:HD12	1.77	0.64
1:A:343:VAL:HA	1:A:347:PHE:HD2	1.62	0.63
1:A:439:ALA:HB2	1:A:482:LEU:HB3	1.82	0.61
1:A:178:GLU:HA	1:A:181:ARG:HD2	1.82	0.61
1:A:17:HIS:CE1	1:A:209:LYS:HB3	2.36	0.61
1:A:47:PHE:CZ	1:A:51:LYS:HE2	2.36	0.60
1:A:246:VAL:O	1:A:337:ARG:NH1	2.34	0.60
1:A:119:GLY:O	1:A:121:THR:N	2.35	0.60
1:A:423:ARG:HG2	1:A:427:LEU:HD23	1.82	0.60
1:A:294:VAL:HG22	1:A:303:VAL:HG22	1.85	0.59
1:A:145:ILE:HD12	1:A:162:VAL:HG13	1.84	0.58
1:A:242:GLU:HG3	1:A:337:ARG:NH2	2.19	0.58
1:A:10:LEU:HD22	1:A:15:GLN:HB2	1.86	0.57
1:A:70:LEU:HG	1:A:297:VAL:HG22	1.88	0.56
1:A:263:ASP:O	1:A:267:ILE:HD12	2.05	0.56
1:A:190:PRO:HA	1:A:193:VAL:HG12	1.86	0.56
1:A:106:ALA:HA	1:A:162:VAL:HG23	1.88	0.55
1:A:20:ARG:HE	1:A:20:ARG:HA	1.71	0.55
1:A:253:CYS:HB3	1:A:255:ASP:OD1	2.07	0.55
1:A:317:ARG:NH1	1:A:321:GLY:O	2.39	0.55
1:A:27:GLU:O	1:A:31:VAL:HG23	2.07	0.55
1:A:145:ILE:HA	1:A:148:VAL:HG12	1.90	0.54
1:A:108:LEU:HA	1:A:165:TYR:O	2.08	0.53
1:A:343:VAL:HA	1:A:347:PHE:CD2	2.43	0.53
1:A:339:PHE:O	1:A:343:VAL:HG13	2.09	0.52
1:A:118:LEU:HG	1:A:124:LYS:HD3	1.90	0.52
1:A:1:MET:HB2	1:A:34:TYR:CE2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:LEU:HD11	1:A:215:LYS:HB3	1.93	0.51
1:A:90:ARG:HG2	1:A:91:ARG:HH21	1.75	0.51
1:A:177:ALA:HB2	1:A:195:MET:HE2	1.93	0.51
1:A:239:GLU:O	1:A:243:ARG:N	2.43	0.51
1:A:280:LYS:HB2	1:A:332:ASN:HB3	1.91	0.51
1:A:1:MET:HE1	1:A:31:VAL:HG13	1.92	0.51
1:A:85:ASP:C	1:A:87:GLU:H	2.19	0.50
1:A:494:ASP:N	1:A:494:ASP:OD1	2.43	0.50
1:A:147:ARG:HD3	1:A:470:LEU:O	2.11	0.50
1:A:337:ARG:HD2	1:A:341:LYS:HE2	1.93	0.50
1:A:349:PRO:HB2	1:A:350:LEU:HD12	1.94	0.50
1:A:292:GLY:HA2	1:A:305:GLU:HA	1.94	0.50
1:A:108:LEU:HG	1:A:165:TYR:HB2	1.94	0.49
1:A:288:GLU:OE1	1:A:288:GLU:N	2.41	0.49
1:A:361:PRO:HA	1:A:371:LYS:HA	1.94	0.49
1:A:129:VAL:HG12	1:A:131:LEU:HD13	1.94	0.49
1:A:145:ILE:HD11	1:A:188:LEU:HD21	1.94	0.49
1:A:242:GLU:HA	1:A:337:ARG:CZ	2.42	0.49
1:A:282:VAL:HG12	1:A:399:VAL:O	2.12	0.49
1:A:95:GLU:OE2	1:A:272:LEU:HB2	2.13	0.49
1:A:2:ASN:ND2	1:A:5:ASP:HB2	2.28	0.48
1:A:111:ALA:O	1:A:168:THR:HA	2.13	0.48
1:A:429:GLN:OE1	1:A:433:TRP:NE1	2.45	0.48
1:A:20:ARG:HA	1:A:20:ARG:NE	2.29	0.48
1:A:25:LEU:HD21	1:A:215:LYS:HB2	1.95	0.48
1:A:40:MET:HE1	1:A:210:VAL:HG21	1.96	0.47
1:A:293:VAL:HG11	1:A:325:TYR:HB3	1.97	0.47
1:A:482:LEU:HD12	1:A:487:PHE:CZ	2.49	0.47
1:A:17:HIS:HE1	1:A:209:LYS:HB3	1.79	0.47
1:A:118:LEU:O	1:A:120:VAL:N	2.48	0.47
1:A:1:MET:HE2	1:A:1:MET:HB3	1.73	0.47
1:A:232:LEU:HD13	1:A:238:LEU:HD11	1.97	0.47
1:A:305:GLU:O	1:A:308:GLU:N	2.39	0.47
1:A:109:LEU:HD21	1:A:138:TYR:HE1	1.79	0.47
1:A:92:TRP:N	1:A:92:TRP:CD1	2.80	0.46
1:A:106:ALA:HA	1:A:163:PRO:HD2	1.97	0.46
1:A:129:VAL:HG22	1:A:419:PRO:HA	1.98	0.46
1:A:364:ASP:OD1	1:A:365:GLU:N	2.44	0.46
1:A:91:ARG:O	1:A:95:GLU:HG2	2.15	0.46
1:A:276:ASP:OD1	1:A:276:ASP:N	2.44	0.46
1:A:340:LEU:HD23	1:A:340:LEU:HA	1.66	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:LYS:HD2	1:A:375:PRO:O	2.15	0.45
1:A:252:TYR:CZ	1:A:261:LEU:HD21	2.51	0.45
1:A:126:MET:HG2	1:A:179:PHE:CD2	2.52	0.45
1:A:478:LEU:HB3	1:A:482:LEU:HD23	1.97	0.45
1:A:347:PHE:HB3	1:A:390:PHE:CE2	2.53	0.44
1:A:323:LEU:HA	1:A:323:LEU:HD23	1.64	0.43
1:A:49:PHE:CE2	1:A:53:ILE:HD11	2.53	0.43
1:A:476:GLU:CD	1:A:476:GLU:H	2.24	0.43
1:A:19:LEU:HD13	1:A:19:LEU:HA	1.72	0.43
1:A:107:VAL:HG23	1:A:164:TRP:HA	2.01	0.43
1:A:465:CYS:HB2	1:A:491:LEU:HD23	2.00	0.43
1:A:50:GLN:O	1:A:54:GLU:HG3	2.18	0.43
1:A:151:LEU:O	1:A:155:ARG:HB2	2.19	0.43
1:A:232:LEU:HB3	1:A:238:LEU:HD13	2.01	0.43
1:A:70:LEU:O	1:A:71:ARG:HD2	2.19	0.43
1:A:20:ARG:HE	1:A:20:ARG:CA	2.32	0.43
1:A:168:THR:O	1:A:197:GLU:HA	2.18	0.43
1:A:312:GLU:N	1:A:312:GLU:OE1	2.52	0.43
1:A:103:ASN:OD1	1:A:160:CYS:HB2	2.20	0.42
1:A:276:ASP:O	1:A:336:THR:HG22	2.19	0.42
1:A:77:ARG:CD	1:A:301:PRO:HG3	2.49	0.42
1:A:263:ASP:O	1:A:266:PHE:N	2.53	0.42
1:A:339:PHE:HE2	1:A:387:VAL:HG13	1.83	0.42
1:A:90:ARG:NE	1:A:90:ARG:HA	2.34	0.42
1:A:116:THR:C	1:A:118:LEU:H	2.23	0.42
1:A:241:MET:HE2	1:A:241:MET:HB2	1.93	0.42
1:A:97:PHE:CZ	1:A:151:LEU:HB3	2.55	0.42
1:A:354:HIS:N	1:A:381:GLU:O	2.33	0.42
1:A:371:LYS:H	1:A:371:LYS:HG2	1.62	0.41
1:A:85:ASP:O	1:A:87:GLU:N	2.53	0.41
1:A:261:LEU:HD23	1:A:261:LEU:HA	1.58	0.41
1:A:469:PRO:HA	1:A:472:SER:O	2.20	0.41
1:A:199:ARG:C	1:A:200:LEU:HD13	2.45	0.41
1:A:260:ARG:HB3	1:A:263:ASP:HB2	2.02	0.41
1:A:431:TYR:HE1	1:A:441:PHE:CD1	2.37	0.41
1:A:473:TYR:CD1	1:A:473:TYR:C	2.98	0.41
1:A:302:GLN:HB2	1:A:388:PHE:HB2	2.02	0.41
1:A:36:GLU:OE2	1:A:218:VAL:HB	2.20	0.41
1:A:117:ARG:NE	1:A:412:GLU:OE1	2.43	0.41
1:A:330:ILE:HB	1:A:405:PHE:CD1	2.56	0.41
1:A:480:VAL:H	1:A:480:VAL:HG12	1.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:VAL:CG2	1:A:398:GLU:HG2	2.51	0.40
1:A:343:VAL:HA	1:A:347:PHE:HB2	2.03	0.40
1:A:438:GLY:HA3	1:A:483:GLN:HA	2.03	0.40
1:A:97:PHE:HA	1:A:100:ILE:HD13	2.02	0.40
1:A:3:ILE:HD13	1:A:6:LEU:HB3	2.02	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	468/510 (92%)	406 (87%)	50 (11%)	12 (3%)	4 7

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-2	PRO
1	A	372	PRO
1	A	117	ARG
1	A	119	GLY
1	A	120	VAL
1	A	306	TYR
1	A	219	ALA
1	A	-1	GLU
1	A	74	PRO
1	A	448	TRP
1	A	371	LYS
1	A	374	LYS

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	403/432 (93%)	338 (84%)	65 (16%)	2 4

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	GLU
1	A	3	ILE
1	A	19	LEU
1	A	69	ARG
1	A	80	ARG
1	A	89	ARG
1	A	91	ARG
1	A	114	GLN
1	A	116	THR
1	A	118	LEU
1	A	120	VAL
1	A	121	THR
1	A	129	VAL
1	A	131	LEU
1	A	137	LEU
1	A	139	GLN
1	A	145	ILE
1	A	158	THR
1	A	162	VAL
1	A	178	GLU
1	A	188	LEU
1	A	195	MET
1	A	200	LEU
1	A	233	GLU
1	A	241	MET
1	A	243	ARG
1	A	249	VAL
1	A	259	VAL
1	A	261	LEU
1	A	273	GLN

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Mol	Chain	Res	Type
1	A	283	GLU
1	A	294	VAL
1	A	304	VAL
1	A	307	SER
1	A	315	GLN
1	A	316	LEU
1	A	319	SER
1	A	329	ASN
1	A	337	ARG
1	A	343	VAL
1	A	355	VAL
1	A	370	VAL
1	A	371	LYS
1	A	373	LEU
1	A	374	LYS
1	A	387	VAL
1	A	399	VAL
1	A	400	LEU
1	A	403	GLU
1	A	404	GLU
1	A	418	SER
1	A	424	GLN
1	A	427	LEU
1	A	445	HIS
1	A	449	LEU
1	A	465	CYS
1	A	468	SER
1	A	474	SER
1	A	476	GLU
1	A	480	VAL
1	A	482	LEU
1	A	488	GLN
1	A	491	LEU
1	A	493	LEU
1	A	494	ASP

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	198	GLN

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.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	476/510 (93%)	0.34	14 (2%)	53 52	48, 79, 122, 165	76 (15%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	445	HIS	6.4
1	A	87	GLU	5.2
1	A	119	GLY	4.0
1	A	186	PHE	3.3
1	A	438	GLY	3.2
1	A	373	LEU	2.9
1	A	118	LEU	2.8
1	A	84	SER	2.3
1	A	272	LEU	2.2
1	A	448	TRP	2.2
1	A	444	ALA	2.1
1	A	306	TYR	2.1
1	A	86	PRO	2.0
1	A	114	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.