



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:34 PM UTC

PDB ID : 2PR9 / pdb_00002pr9
Title : Mu2 adaptin subunit (AP50) of AP2 adaptor (second domain), complexed with GABAA receptor-gamma2 subunit-derived internalization peptide DEEY-GYECL
Authors : Vahedi-Faridi, A.; Haucke, V.; Kittler, J.T.; Kukhtina, V.; Moss, S.J.; Saenger, W.; Chen, G.-J.; Tretter, V.; Smith, K.; Yan, Z.; McAinsh, K.; Arancibia-Carcamo, L.
Deposited on : 2007-05-04
Resolution : 2.51 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
Xtrriage (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

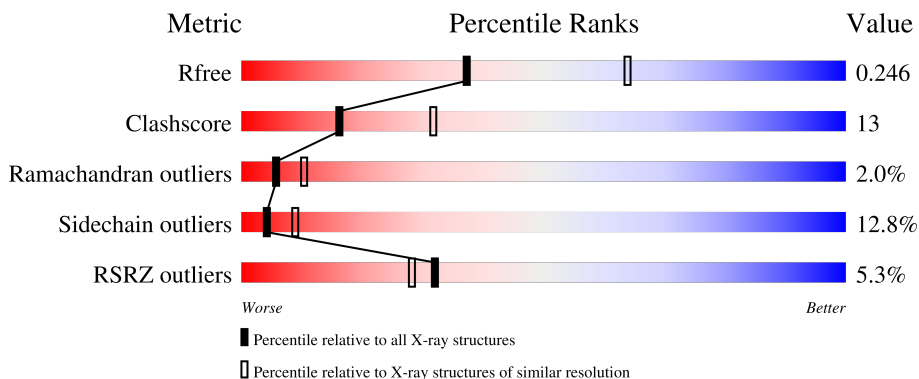
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	299	
2	P	10	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2174 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 AP-2 complex subunit mu-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	254	2029	1305	353	357	14	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	MET	-	expression tag	UNP P84092
A	138	GLY	-	expression tag	UNP P84092
A	139	SER	-	expression tag	UNP P84092
A	140	SER	-	expression tag	UNP P84092
A	141	HIS	-	expression tag	UNP P84092
A	142	HIS	-	expression tag	UNP P84092
A	143	HIS	-	expression tag	UNP P84092
A	144	HIS	-	expression tag	UNP P84092
A	145	HIS	-	expression tag	UNP P84092
A	146	HIS	-	expression tag	UNP P84092
A	147	SER	-	expression tag	UNP P84092
A	148	SER	-	expression tag	UNP P84092
A	149	GLY	-	expression tag	UNP P84092
A	150	LEU	-	expression tag	UNP P84092
A	151	VAL	-	expression tag	UNP P84092
A	152	PRO	-	expression tag	UNP P84092
A	153	ARG	-	expression tag	UNP P84092
A	154	GLY	-	expression tag	UNP P84092
A	155	SER	-	expression tag	UNP P84092
A	156	HIS	-	expression tag	UNP P84092
A	157	MET	-	expression tag	UNP P84092

- Molecule 2 is a protein called GABA(A) receptor subunit gamma-2 peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	P	10	86	52	10	23	1	0	0	0

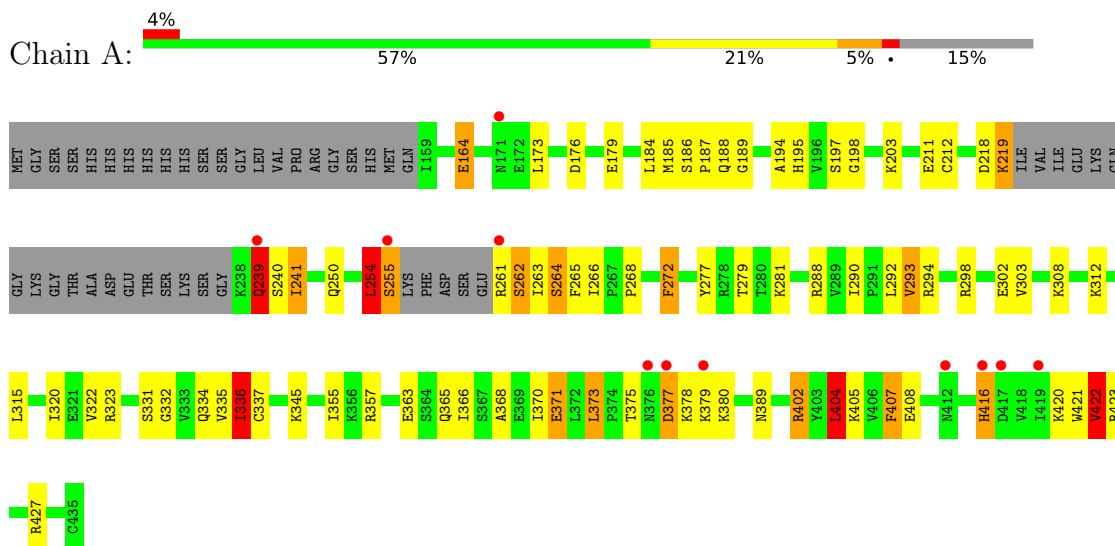
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total	O	0	0
			54	54		
3	P	5	Total	O	0	0
			5	5		

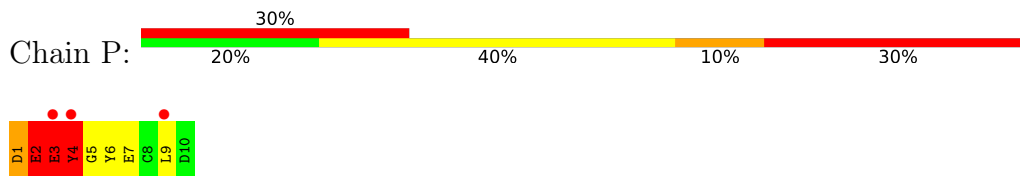
3 Residue-property plots [i](#)

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 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: AP-2 complex subunit mu-1



- Molecule 2: GABA(A) receptor subunit gamma-2 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	126.30Å 126.30Å 74.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	9.97 – 2.51 9.97 – 2.51	Depositor EDS
% Data completeness (in resolution range)	100.0 (9.97-2.51) 98.0 (9.97-2.51)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 2.50Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.203 , 0.240 (Not available) , 0.246	Depositor DCC
R_{free} test set	913 reflections (4.00%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtrriage
Anisotropy	0.030	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.044 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2174	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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	1.97	30/2070 (1.4%)	1.61	28/2786 (1.0%)
2	P	1.93	3/87 (3.4%)	1.93	3/115 (2.6%)
All	All	1.96	33/2157 (1.5%)	1.63	31/2901 (1.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
2	P	0	2
All	All	0	4

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	255	SER	C-O	30.05	1.83	1.23
1	A	219	LYS	C-O	19.90	1.63	1.23
1	A	261	ARG	NE-CZ	13.64	1.48	1.33
1	A	255	SER	CA-C	11.73	1.77	1.52
1	A	290	ILE	CA-CB	-10.95	1.44	1.54
1	A	261	ARG	CZ-NH2	10.23	1.46	1.33
1	A	254	LEU	C-O	9.30	1.35	1.24
1	A	261	ARG	CZ-NH1	7.67	1.43	1.32
1	A	239	GLN	CA-C	7.65	1.63	1.52
1	A	262	SER	CA-C	7.55	1.62	1.53
1	A	254	LEU	C-N	7.23	1.43	1.33
1	A	261	ARG	CD-NE	7.22	1.56	1.46
1	A	261	ARG	CG-CD	7.06	1.73	1.52
2	P	3	GLU	CA-C	6.96	1.62	1.52
1	A	315	LEU	N-CA	6.86	1.53	1.45
1	A	375	THR	CA-C	-6.46	1.44	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	427	ARG	NE-CZ	6.34	1.40	1.33
1	A	322	VAL	CA-CB	-6.14	1.47	1.54
1	A	293	VAL	C-O	-6.11	1.17	1.24
1	A	323	ARG	CB-CG	-6.05	1.34	1.52
1	A	239	GLN	N-CA	5.81	1.53	1.46
2	P	3	GLU	N-CA	5.76	1.53	1.46
1	A	323	ARG	CA-CB	-5.71	1.45	1.53
1	A	370	ILE	CA-CB	-5.68	1.47	1.54
1	A	176	ASP	C-O	5.63	1.30	1.24
1	A	254	LEU	CA-C	5.52	1.60	1.52
1	A	254	LEU	N-CA	5.44	1.53	1.46
1	A	272	PHE	N-CA	5.42	1.52	1.46
1	A	422	VAL	C-O	5.30	1.29	1.24
2	P	4	TYR	CA-C	5.27	1.64	1.52
1	A	241	ILE	CA-CB	-5.23	1.46	1.54
1	A	389	ASN	CG-OD1	5.16	1.33	1.23
1	A	427	ARG	CZ-NH1	5.08	1.39	1.32

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	SER	CA-C-O	-16.12	93.40	120.80
1	A	241	ILE	N-CA-C	13.69	128.49	109.45
1	A	219	LYS	CA-C-O	-10.96	102.17	120.80
2	P	3	GLU	N-CA-C	10.57	133.32	110.80
1	A	263	ILE	N-CA-C	8.49	120.16	107.51
1	A	189	GLY	N-CA-C	8.45	127.18	115.27
1	A	265	PHE	CA-C-N	-8.30	116.61	122.59
1	A	265	PHE	C-N-CA	-8.30	116.61	122.59
1	A	422	VAL	CB-CA-C	-7.99	97.91	110.69
1	A	312	LYS	CA-C-N	7.53	128.20	119.47
1	A	312	LYS	C-N-CA	7.53	128.20	119.47
2	P	3	GLU	CA-C-N	6.94	134.19	121.70
2	P	3	GLU	C-N-CA	6.94	134.19	121.70
1	A	164	GLU	N-CA-C	6.92	121.97	112.90
1	A	254	LEU	CA-C-N	-6.61	109.81	121.70
1	A	254	LEU	C-N-CA	-6.61	109.81	121.70
1	A	240	SER	N-CA-C	6.45	118.83	107.61
1	A	179	GLU	CA-C-N	-6.26	114.16	122.99
1	A	179	GLU	C-N-CA	-6.26	114.16	122.99
1	A	365	GLN	N-CA-C	5.79	118.99	109.72
1	A	323	ARG	N-CA-CB	-5.72	101.72	110.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	303	VAL	N-CA-C	5.68	116.04	107.80
1	A	265	PHE	N-CA-C	5.54	116.94	108.52
1	A	264	SER	N-CA-C	5.49	118.41	109.24
1	A	336	ILE	CG1-CB-CG2	-5.41	94.47	110.70
1	A	404	LEU	CA-CB-CG	-5.41	97.36	116.30
1	A	261	ARG	NE-CZ-NH2	-5.39	114.35	119.20
1	A	335	VAL	N-CA-C	5.11	117.09	108.81
1	A	373	LEU	CA-C-N	-5.08	114.55	119.78
1	A	373	LEU	C-N-CA	-5.08	114.55	119.78
1	A	266	ILE	CA-C-O	5.01	122.11	119.15

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	218	ASP	Peptide
1	A	262	SER	Peptide
2	P	2	GLU	Peptide
2	P	3	GLU	Peptide

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	2029	0	2095	45	0
2	P	86	0	65	15	0
3	A	54	0	0	13	1
3	P	5	0	0	0	0
All	All	2174	0	2160	56	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 13.

All (56) 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:255:SER:CA	1:A:255:SER:C	1.77	1.54
1:A:255:SER:C	1:A:255:SER:O	1.83	1.18
1:A:185:MET:HE2	3:A:479:HOH:O	1.64	0.94
2:P:5:GLY:H	2:P:6:TYR:HA	1.37	0.90
1:A:402:ARG:HD2	3:A:474:HOH:O	1.78	0.80
2:P:2:GLU:O	2:P:3:GLU:HG3	1.87	0.75
2:P:4:TYR:HB3	2:P:5:GLY:HA3	1.67	0.74
1:A:254:LEU:H	1:A:254:LEU:HD22	1.53	0.72
1:A:345:LYS:HD2	3:A:453:HOH:O	1.88	0.72
1:A:173:LEU:HD22	1:A:404:LEU:HD22	1.69	0.71
2:P:5:GLY:N	2:P:6:TYR:HA	2.02	0.71
1:A:173:LEU:CD2	1:A:404:LEU:HD22	2.22	0.70
1:A:416:HIS:HB3	3:A:446:HOH:O	1.92	0.69
2:P:5:GLY:H	2:P:6:TYR:CA	2.09	0.65
1:A:421:TRP:HB3	2:P:6:TYR:HB3	1.79	0.65
1:A:185:MET:CE	3:A:479:HOH:O	2.32	0.65
1:A:255:SER:C	1:A:255:SER:CB	2.72	0.62
2:P:5:GLY:N	2:P:6:TYR:CA	2.62	0.62
1:A:294:ARG:NH1	1:A:302:GLU:HG3	2.15	0.61
1:A:212:CYS:HA	1:A:405:LYS:O	2.00	0.60
2:P:4:TYR:CB	2:P:5:GLY:HA3	2.33	0.57
1:A:422:VAL:HG22	2:P:9:LEU:HB2	1.86	0.57
1:A:308:LYS:HG2	1:A:363:GLU:HG3	1.92	0.51
1:A:377:ASP:CB	3:A:472:HOH:O	2.59	0.51
1:A:281:LYS:CG	3:A:466:HOH:O	2.58	0.51
2:P:1:ASP:O	2:P:2:GLU:HG3	2.11	0.51
1:A:194:ALA:O	1:A:195:HIS:HB3	2.11	0.50
1:A:332:GLY:O	1:A:371:GLU:HG3	2.12	0.50
2:P:1:ASP:O	2:P:2:GLU:CG	2.60	0.49
1:A:268:PRO:HD2	1:A:272:PHE:CE1	2.48	0.49
1:A:239:GLN:CG	1:A:402:ARG:HH22	2.25	0.49
1:A:239:GLN:HG2	1:A:402:ARG:HH22	1.78	0.48
1:A:337:CYS:HB3	1:A:366:ILE:HG13	1.94	0.48
1:A:407:PHE:HB3	3:A:455:HOH:O	2.12	0.48
1:A:186:SER:HB2	1:A:187:PRO:CD	2.44	0.47
1:A:255:SER:CA	1:A:255:SER:O	2.62	0.47
1:A:334:GLN:O	1:A:368:ALA:HA	2.15	0.47
1:A:377:ASP:HB3	3:A:472:HOH:O	2.14	0.47
1:A:255:SER:C	1:A:255:SER:N	2.65	0.47
1:A:423:ARG:HG2	2:P:6:TYR:CE1	2.50	0.46
1:A:423:ARG:HD3	2:P:5:GLY:HA2	1.97	0.46
1:A:254:LEU:C	1:A:255:SER:C	2.84	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:478:HOH:O	2:P:7:GLU:HG3	2.17	0.45
1:A:173:LEU:C	1:A:173:LEU:HD12	2.43	0.44
1:A:377:ASP:HB2	3:A:472:HOH:O	2.19	0.43
1:A:239:GLN:HG2	1:A:402:ARG:HH12	1.84	0.43
1:A:404:LEU:O	1:A:420:LYS:HE2	2.19	0.43
1:A:173:LEU:HD21	1:A:404:LEU:HD22	1.98	0.42
1:A:198:GLY:HA3	1:A:277:TYR:CZ	2.54	0.42
1:A:357:ARG:HB3	3:A:485:HOH:O	2.19	0.42
1:A:331:SER:HB3	1:A:373:LEU:HG	2.01	0.42
1:A:250:GLN:HB2	3:A:442:HOH:O	2.19	0.42
1:A:268:PRO:HD2	1:A:272:PHE:CD1	2.56	0.41
1:A:336:ILE:HD13	1:A:336:ILE:HG21	1.64	0.41
2:P:2:GLU:C	2:P:3:GLU:HG3	2.45	0.41
1:A:320:ILE:HB	1:A:355:ILE:HB	2.03	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)
3:A:463:HOH:O	3:A:463:HOH:O[4_565]	2.08	0.12

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	248/299 (83%)	237 (96%)	8 (3%)	3 (1%)	10	20
2	P	8/10 (80%)	4 (50%)	2 (25%)	2 (25%)	0	0
All	All	256/309 (83%)	241 (94%)	10 (4%)	5 (2%)	6	10

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	P	3	GLU
1	A	254	LEU
1	A	239	GLN
1	A	378	LYS
2	P	2	GLU

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	225/268 (84%)	198 (88%)	27 (12%)	5	10
2	P	9/9 (100%)	6 (67%)	3 (33%)	0	0
All	All	234/277 (84%)	204 (87%)	30 (13%)	4	9

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	164	GLU
1	A	184	LEU
1	A	188	GLN
1	A	197	SER
1	A	203	LYS
1	A	211	GLU
1	A	219	LYS
1	A	239	GLN
1	A	241	ILE
1	A	254	LEU
1	A	264	SER
1	A	279	THR
1	A	288	ARG
1	A	292	LEU
1	A	293	VAL
1	A	298	ARG
1	A	336	ILE
1	A	371	GLU
1	A	377	ASP

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Mol	Chain	Res	Type
1	A	379	LYS
1	A	380	LYS
1	A	402	ARG
1	A	404	LEU
1	A	407	PHE
1	A	408	GLU
1	A	416	HIS
1	A	422	VAL
2	P	1	ASP
2	P	3	GLU
2	P	4	TYR

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

Mol	Chain	Res	Type
1	A	190	GLN
1	A	250	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

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	254/299 (84%)	0.01	11 (4%) 40 35	36, 49, 67, 94	0
2	P	10/10 (100%)	1.63	3 (30%) 1 1	46, 56, 70, 72	0
All	All	264/309 (85%)	0.07	14 (5%) 32 28	36, 50, 67, 94	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	239	GLN	3.6
2	P	3	GLU	2.8
1	A	255	SER	2.8
1	A	261	ARG	2.6
2	P	4	TYR	2.5
1	A	419	ILE	2.4
1	A	412	ASN	2.4
1	A	379	LYS	2.3
1	A	416	HIS	2.3
1	A	377	ASP	2.3
2	P	9	LEU	2.1
1	A	171	ASN	2.1
1	A	376	ASN	2.1
1	A	417	ASP	2.1

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.