



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:56 AM UTC

PDB ID : 3H85 / pdb_00003h85
Title : Molecular basis for the association of PIPKI gamma-p90 with the clathrin adaptor AP-2
Authors : Vahedi-Faridi, A.; Kahlfeldt, N.; Schaefer, J.G.; Haucke, V.
Deposited on : 2009-04-28
Resolution : 2.60 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

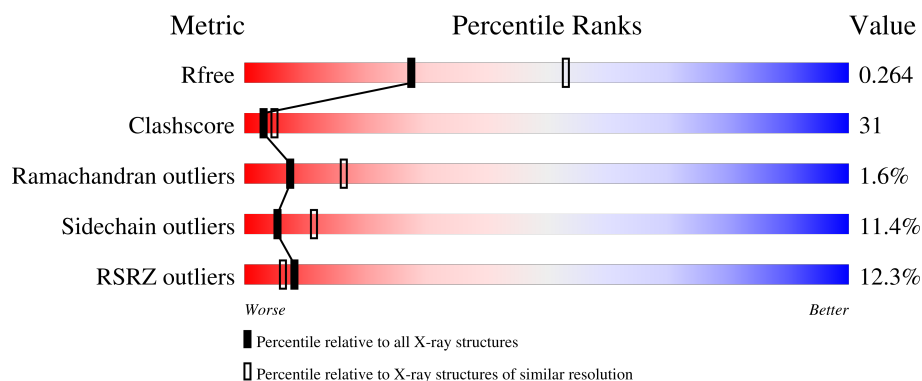
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	8	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2142 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	253	2039	1311	359	355	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 Phosphatidylinositol-4-phosphate 5-kinase type-1 gamma.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	8	Total	C	N	O	0	0	0
			71	48	11	12			

- Molecule 3 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ni	0	0
			1	1		

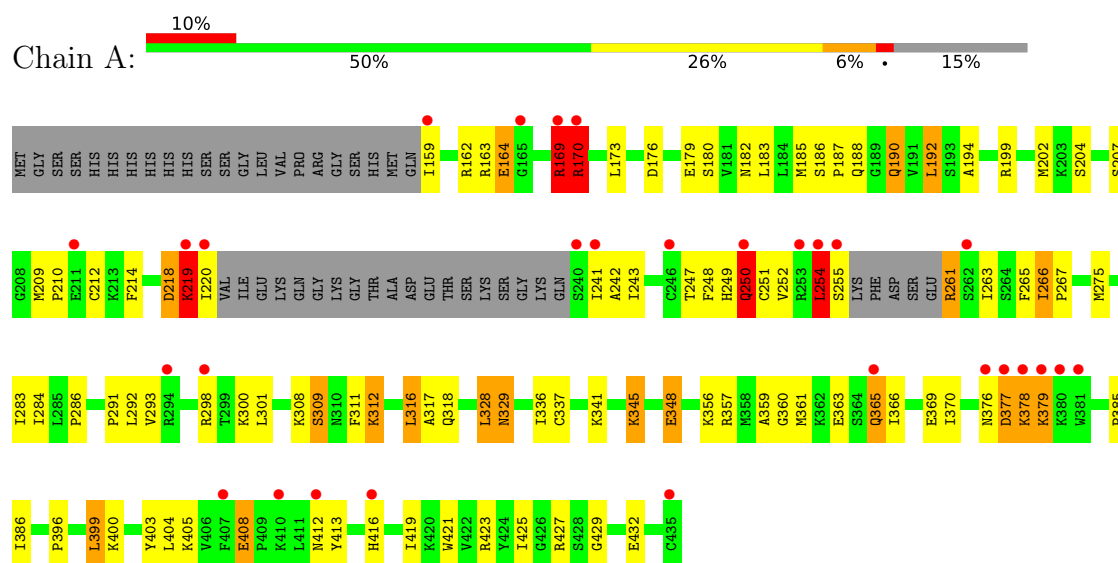
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total	O	0	0
			31	31		

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: AP-2 complex subunit mu-1



• Molecule 2: Phosphatidylinositol-4-phosphate 5-kinase type-1 gamma



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	125.30Å 125.30Å 74.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.01 – 2.60 41.01 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.6 (41.01-2.60) 92.6 (41.01-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.231 , 0.263 (Not available) , 0.264	Depositor DCC
R_{free} test set	792 reflections (4.15%)	wwPDB-VP
Wilson B-factor (Å ²)	67.5	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2142	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NI

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	0.52	0/2080	0.89	5/2796 (0.2%)
2	P	0.74	0/75	0.82	1/102 (1.0%)
All	All	0.53	0/2155	0.89	6/2898 (0.2%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	ARG	N-CA-C	-7.05	99.06	110.20
1	A	219	LYS	N-CA-C	-6.02	105.04	112.38
1	A	266	ILE	CA-C-N	5.72	123.78	119.66
1	A	266	ILE	C-N-CA	5.72	123.78	119.66
2	P	2	TRP	N-CA-C	5.55	118.30	109.65
1	A	164	GLU	N-CA-C	5.30	116.87	111.14

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	2039	0	2132	132	0
2	P	71	0	65	5	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	31	0	0	9	0
All	All	2142	0	2197	134	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 31.

All (134) 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:261:ARG:O	1:A:261:ARG:HD3	1.45	1.16
1:A:248:PHE:HB3	1:A:252:VAL:HG11	1.19	1.14
1:A:399:LEU:C	1:A:399:LEU:HD23	1.77	1.10
1:A:399:LEU:HD23	1:A:400:LYS:N	1.67	1.09
1:A:425:ILE:HG12	2:P:1:SER:HA	1.35	1.05
1:A:312:LYS:N	1:A:312:LYS:HD3	1.69	1.04
1:A:248:PHE:CB	1:A:252:VAL:HG11	1.88	1.03
1:A:312:LYS:H	1:A:312:LYS:CD	1.71	1.00
1:A:248:PHE:HB3	1:A:252:VAL:CG1	1.91	0.99
1:A:169:ARG:CB	1:A:169:ARG:HH11	1.76	0.98
1:A:169:ARG:HH11	1:A:169:ARG:CG	1.78	0.95
1:A:214:PHE:HE2	1:A:275:MET:CE	1.80	0.94
1:A:379:LYS:H	1:A:379:LYS:HD3	1.34	0.91
1:A:214:PHE:HE2	1:A:275:MET:HE2	1.33	0.91
1:A:254:LEU:HD23	1:A:254:LEU:H	1.36	0.90
1:A:214:PHE:CE2	1:A:275:MET:HE2	2.09	0.87
1:A:312:LYS:N	1:A:312:LYS:CD	2.34	0.84
4:A:31:HOH:O	2:P:8:HIS:HE1	1.60	0.84
1:A:400:LYS:CE	4:A:24:HOH:O	2.24	0.84
1:A:399:LEU:C	1:A:399:LEU:CD2	2.50	0.83
1:A:261:ARG:HD3	1:A:261:ARG:C	2.03	0.82
1:A:329:ASN:HD22	1:A:329:ASN:H	1.28	0.81
1:A:186:SER:HB3	1:A:192:LEU:HD11	1.65	0.79
1:A:218:ASP:CG	1:A:218:ASP:O	2.25	0.79
1:A:328:LEU:H	1:A:328:LEU:CD2	1.96	0.78
1:A:400:LYS:HE2	4:A:24:HOH:O	1.83	0.78
1:A:169:ARG:HH11	1:A:169:ARG:HG2	1.47	0.78
1:A:180:SER:HA	1:A:427:ARG:O	1.84	0.77
1:A:328:LEU:H	1:A:328:LEU:HD22	1.49	0.77
1:A:311:PHE:HB2	1:A:312:LYS:HE3	1.67	0.76
1:A:328:LEU:HD22	1:A:328:LEU:N	2.02	0.75
1:A:163:ARG:O	1:A:209:MET:HE1	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:GLU:HA	1:A:209:MET:CE	2.17	0.74
1:A:255:SER:CB	1:A:261:ARG:HB2	2.18	0.74
1:A:400:LYS:HE3	4:A:24:HOH:O	1.87	0.74
1:A:169:ARG:CB	1:A:169:ARG:NH1	2.50	0.74
1:A:292:LEU:HD23	1:A:292:LEU:C	2.15	0.72
1:A:219:LYS:NZ	1:A:219:LYS:HB3	2.07	0.69
1:A:169:ARG:O	1:A:170:ARG:CB	2.41	0.69
1:A:249:HIS:O	1:A:252:VAL:HG12	1.94	0.67
1:A:219:LYS:NZ	1:A:219:LYS:CB	2.57	0.67
1:A:308:LYS:NZ	1:A:363:GLU:OE1	2.28	0.67
1:A:169:ARG:NH1	1:A:169:ARG:HB2	2.10	0.66
1:A:309:SER:HB2	1:A:360:GLY:HA2	1.78	0.66
1:A:219:LYS:HB3	1:A:219:LYS:HZ3	1.59	0.66
1:A:169:ARG:HG2	1:A:169:ARG:NH1	2.11	0.65
1:A:312:LYS:H	1:A:312:LYS:CE	2.09	0.65
1:A:254:LEU:HD23	1:A:254:LEU:N	2.08	0.64
1:A:248:PHE:CB	1:A:252:VAL:CG1	2.64	0.64
1:A:210:PRO:O	1:A:266:ILE:HG12	1.98	0.63
1:A:399:LEU:HD23	1:A:400:LYS:CA	2.28	0.63
1:A:170:ARG:HH21	1:A:419:ILE:HD13	1.62	0.62
1:A:377:ASP:OD1	1:A:377:ASP:N	2.33	0.61
1:A:291:PRO:HB3	1:A:386:ILE:HD12	1.83	0.60
1:A:163:ARG:O	1:A:209:MET:CE	2.49	0.60
1:A:249:HIS:O	1:A:251:CYS:N	2.34	0.60
1:A:365:GLN:O	1:A:365:GLN:HG3	2.01	0.60
1:A:252:VAL:HG21	1:A:263:ILE:HG23	1.85	0.59
1:A:248:PHE:CG	1:A:252:VAL:HG11	2.40	0.57
1:A:220:ILE:HG12	1:A:242:ALA:HA	1.88	0.56
1:A:179:GLU:OE1	1:A:396:PRO:HD2	2.06	0.56
1:A:427:ARG:HD3	4:A:25:HOH:O	2.05	0.56
1:A:261:ARG:C	1:A:261:ARG:CD	2.72	0.55
1:A:255:SER:HB3	1:A:261:ARG:HB2	1.87	0.55
1:A:311:PHE:C	1:A:361:MET:HE2	2.32	0.54
1:A:404:LEU:C	1:A:404:LEU:HD23	2.32	0.54
1:A:345:LYS:HD3	1:A:348:GLU:OE1	2.07	0.54
1:A:214:PHE:CE2	1:A:275:MET:CE	2.71	0.54
1:A:284:ILE:O	1:A:286:PRO:HD3	2.08	0.53
1:A:218:ASP:O	1:A:220:ILE:HG13	2.09	0.53
1:A:249:HIS:ND1	1:A:251:CYS:HB3	2.23	0.53
1:A:164:GLU:HA	1:A:209:MET:HE2	1.90	0.52
1:A:173:LEU:C	1:A:173:LEU:HD12	2.35	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:LEU:CD2	1:A:328:LEU:N	2.62	0.51
1:A:218:ASP:C	1:A:220:ILE:N	2.67	0.50
1:A:169:ARG:O	1:A:170:ARG:HB3	2.12	0.50
1:A:170:ARG:HE	1:A:419:ILE:HD12	1.77	0.50
1:A:421:TRP:HB3	2:P:4:TYR:HB3	1.93	0.50
1:A:159:ILE:HG23	1:A:159:ILE:O	2.11	0.50
1:A:190:GLN:O	1:A:192:LEU:HD13	2.11	0.50
1:A:292:LEU:C	1:A:292:LEU:CD2	2.84	0.50
1:A:318:GLN:NE2	2:P:2:TRP:CZ2	2.79	0.50
1:A:241:ILE:HD11	1:A:243:ILE:HD11	1.93	0.50
1:A:300:LYS:HD2	1:A:369:GLU:HG2	1.93	0.50
1:A:163:ARG:C	1:A:209:MET:HE1	2.37	0.49
1:A:376:ASN:OD1	1:A:378:LYS:HB2	2.12	0.49
1:A:219:LYS:CB	1:A:219:LYS:HZ2	2.23	0.49
1:A:311:PHE:O	1:A:361:MET:HE2	2.12	0.49
1:A:379:LYS:HD3	1:A:379:LYS:N	2.16	0.49
1:A:317:ALA:O	1:A:357:ARG:HA	2.13	0.48
1:A:183:LEU:HD11	1:A:185:MET:HE3	1.97	0.47
1:A:212:CYS:HA	1:A:405:LYS:O	2.14	0.47
1:A:329:ASN:H	1:A:329:ASN:ND2	2.05	0.47
1:A:412:ASN:OD1	1:A:413:TYR:N	2.48	0.47
1:A:164:GLU:HA	1:A:209:MET:HE3	1.96	0.46
1:A:376:ASN:HB3	1:A:379:LYS:NZ	2.31	0.46
1:A:169:ARG:C	1:A:169:ARG:HD3	2.41	0.46
1:A:292:LEU:HD23	1:A:293:VAL:N	2.31	0.46
1:A:252:VAL:HA	1:A:265:PHE:HB3	1.98	0.45
1:A:241:ILE:HD12	1:A:243:ILE:HD12	1.98	0.45
1:A:219:LYS:HZ2	1:A:219:LYS:HB2	1.81	0.45
1:A:408:GLU:O	4:A:3:HOH:O	2.21	0.45
1:A:190:GLN:O	1:A:192:LEU:CD1	2.65	0.45
1:A:192:LEU:CD1	1:A:192:LEU:N	2.80	0.44
1:A:298:ARG:NH1	1:A:377:ASP:HB3	2.32	0.44
1:A:301:LEU:HB2	1:A:370:ILE:HB	1.98	0.44
1:A:316:LEU:HD21	1:A:357:ARG:HD2	2.00	0.44
1:A:207:SER:O	1:A:413:TYR:OH	2.34	0.44
1:A:169:ARG:O	1:A:170:ARG:HB2	2.18	0.43
1:A:311:PHE:HB2	1:A:312:LYS:CE	2.42	0.43
1:A:202:MET:HE2	1:A:204:SER:HB3	1.99	0.43
1:A:316:LEU:HD12	1:A:359:ALA:HA	2.01	0.43
1:A:329:ASN:HD22	1:A:329:ASN:N	2.04	0.43
1:A:173:LEU:HD12	1:A:173:LEU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:GLY:HA2	4:A:26:HOH:O	2.18	0.43
1:A:162:ARG:HD2	1:A:267:PRO:O	2.18	0.42
1:A:283:ILE:CG2	1:A:284:ILE:N	2.81	0.42
1:A:190:GLN:HG2	1:A:192:LEU:CD1	2.49	0.42
1:A:385:PRO:HB2	1:A:432:GLU:HB3	2.02	0.42
1:A:376:ASN:HB3	1:A:379:LYS:HZ1	1.85	0.42
1:A:316:LEU:HD12	1:A:316:LEU:HA	1.79	0.42
1:A:247:THR:HG22	4:A:436:HOH:O	2.20	0.42
1:A:283:ILE:HG22	1:A:284:ILE:N	2.33	0.42
1:A:188:GLN:CD	1:A:188:GLN:N	2.78	0.41
1:A:176:ASP:CG	1:A:423:ARG:HE	2.29	0.41
1:A:255:SER:OG	1:A:261:ARG:HB2	2.19	0.41
1:A:199:ARG:HA	1:A:275:MET:O	2.21	0.41
1:A:337:CYS:HB3	1:A:366:ILE:HG13	2.03	0.41
1:A:312:LYS:H	1:A:312:LYS:HE3	1.84	0.41
1:A:182:ASN:O	1:A:194:ALA:HA	2.20	0.40
1:A:187:PRO:HD2	1:A:188:GLN:OE1	2.21	0.40
1:A:250:GLN:N	4:A:12:HOH:O	2.36	0.40
2:P:7:LEU:HA	2:P:7:LEU:HD12	1.74	0.40
1:A:403:TYR:C	1:A:403:TYR:CD1	3.00	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	247/299 (83%)	229 (93%)	14 (6%)	4 (2%)	7	16
2	P	6/8 (75%)	5 (83%)	1 (17%)	0	100	100
All	All	253/307 (82%)	234 (92%)	15 (6%)	4 (2%)	7	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	254	LEU
1	A	250	GLN
1	A	378	LYS
1	A	170	ARG

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	229/268 (85%)	204 (89%)	25 (11%)	6	13
2	P	8/8 (100%)	6 (75%)	2 (25%)	0	1
All	All	237/276 (86%)	210 (89%)	27 (11%)	5	11

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	ARG
1	A	170	ARG
1	A	190	GLN
1	A	192	LEU
1	A	218	ASP
1	A	219	LYS
1	A	250	GLN
1	A	254	LEU
1	A	261	ARG
1	A	309	SER
1	A	312	LYS
1	A	316	LEU
1	A	328	LEU
1	A	329	ASN
1	A	336	ILE
1	A	341	LYS
1	A	345	LYS
1	A	348	GLU
1	A	356	LYS
1	A	365	GLN

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Mol	Chain	Res	Type
1	A	377	ASP
1	A	379	LYS
1	A	399	LEU
1	A	408	GLU
1	A	416	HIS
2	P	2	TRP
2	P	3	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	195	HIS
1	A	329	ASN
2	P	8	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	253/299 (84%)	0.58	29 (11%) 9 7	46, 62, 82, 92	0
2	P	8/8 (100%)	1.02	3 (37%) 1 0	57, 63, 67, 67	0
All	All	261/307 (85%)	0.59	32 (12%) 8 6	46, 63, 82, 92	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	254	LEU	7.5
1	A	220	ILE	6.3
1	A	219	LYS	4.7
1	A	255	SER	4.4
1	A	379	LYS	4.4
1	A	240	SER	4.2
1	A	412	ASN	4.1
1	A	410	LYS	3.3
1	A	298	ARG	3.1
2	P	8	HIS	2.9
1	A	416	HIS	2.9
1	A	381	TRP	2.9
1	A	253	ARG	2.9
1	A	159	ILE	2.9
1	A	211	GLU	2.8
1	A	365	GLN	2.8
1	A	246	CYS	2.7
1	A	378	LYS	2.6
1	A	294	ARG	2.6
1	A	435	CYS	2.6
2	P	2	TRP	2.5
1	A	165	GLY	2.4
1	A	380	LYS	2.3
1	A	170	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	376	ASN	2.2
1	A	169	ARG	2.2
1	A	377	ASP	2.1
1	A	262	SER	2.1
1	A	241	ILE	2.1
1	A	407	PHE	2.1
1	A	250	GLN	2.1
2	P	1	SER	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](#)

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. The B-factors column lists the minimum, median, 95th 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	B-factors($\text{\AA}^2$)	Q<0.9
3	NI	A	1	1/1	0.95	0.13	98,98,98,98	1

6.5 Other polymers [i](#)

There are no such residues in this entry.