



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:31 AM UTC

PDB ID : 6ZRC / pdb_00006zrc
Title : Structure of the human RBAP48 in complex with a macrocyclic peptide cyclized via a xylene linker attached to two cysteines
Authors : Vetter, I.R.; Porfetye, A.T.
Deposited on : 2020-07-13
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
Mogul	:	2022.3.0, CSD as543be (2022)
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

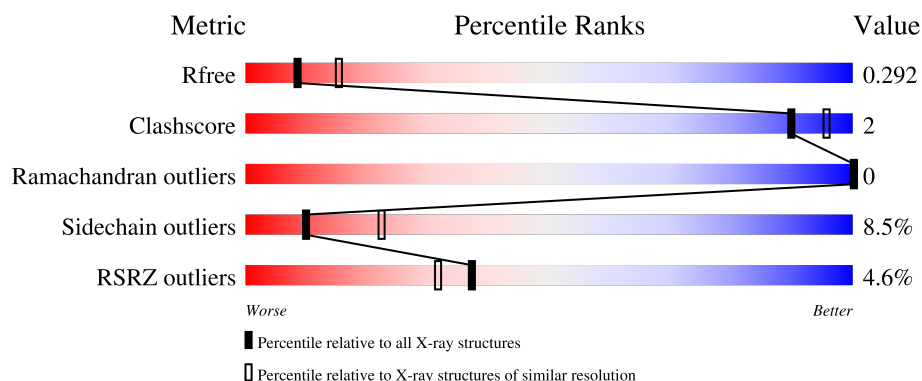
1 Overall quality at a glance

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	427	4% 78% 11% • 11%
1	B	427	3% 78% 10% • 11%
2	P	16	12% 88% 12%
2	Q	16	6% 88% 12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6335 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 Histone-binding protein RBBP4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	0
			3042	1917	519	596	10			
1	B	381	Total	C	N	O	S	0	0	0
			3046	1919	519	598	10			

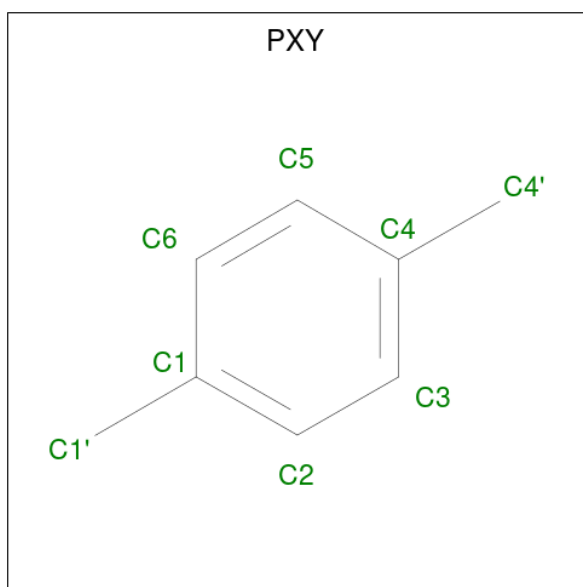
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q09028
A	0	PRO	-	expression tag	UNP Q09028
B	-1	GLY	-	expression tag	UNP Q09028
B	0	PRO	-	expression tag	UNP Q09028

- Molecule 2 is a protein called macrocyclic peptide based on residues 659-672 of the metastasis-associated protein MTA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	16	Total	C	N	O	S	0	0	1
			115	70	26	17	2			
2	Q	16	Total	C	N	O	S	0	0	1
			115	70	26	17	2			

- Molecule 3 is PARA-XYLENE (CCD ID: PXY) (formula: C₈H₁₀).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	P	1	Total C 8 8	0	0
3	Q	1	Total C 8 8	0	0

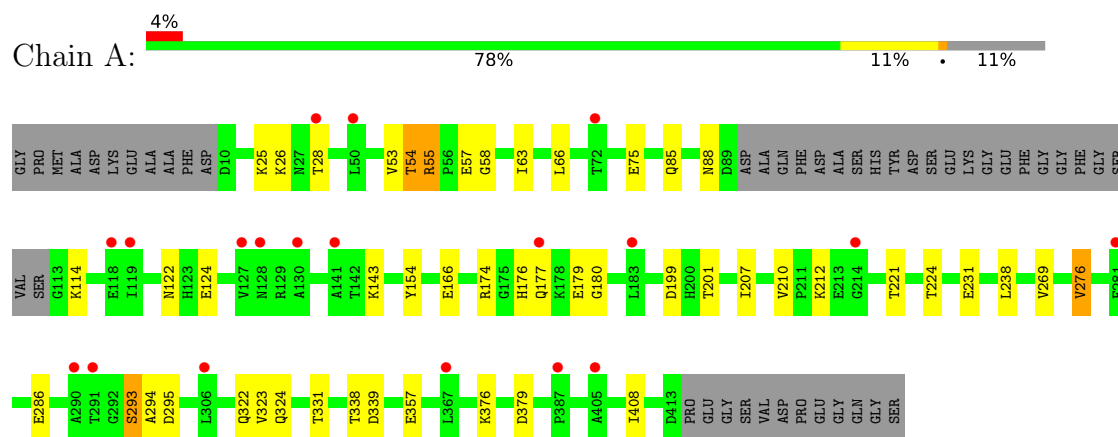
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O 1 1	0	0

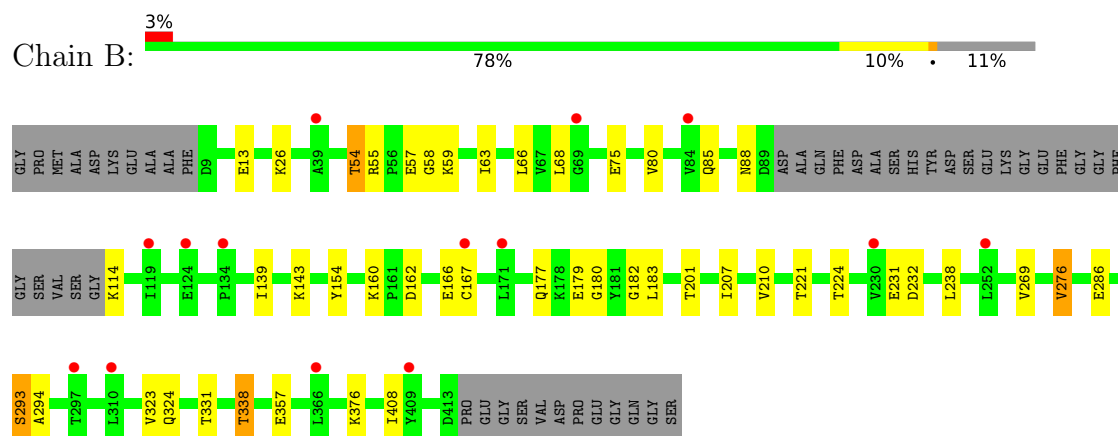
3 Residue-property plots [i](#)

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.

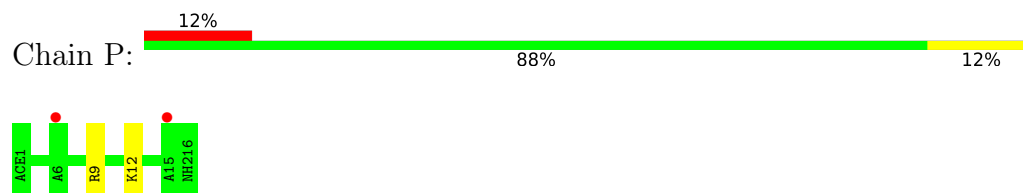
- Molecule 1: Histone-binding protein RBBP4



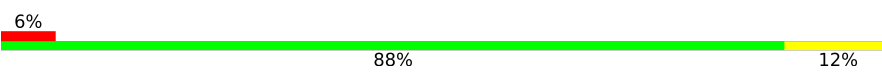
- Molecule 1: Histone-binding protein RBBP4

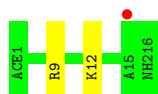


- Molecule 2: macrocyclic peptide based on residues 659-672 of the metastasis-associated protein MTA1



- Molecule 2: macrocyclic peptide based on residues 659-672 of the metastasis-associated protein MTA1

Chain Q: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	71.96Å 96.33Å 149.30Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	48.16 – 2.60 48.16 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.16-2.60) 99.5 (48.16-2.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.18rc7_3834	Depositor
R, R_{free}	0.263 , 0.293 0.265 , 0.292	Depositor DCC
R_{free} test set	1577 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	64.0	Xtriage
Anisotropy	0.930	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 177.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.458 for -h,-k,l	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	6335	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACE, NH2, PXY

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	0/3125	0.85	7/4261 (0.2%)
1	B	0.52	0/3129	0.83	5/4267 (0.1%)
2	P	0.75	0/114	0.95	0/152
2	Q	0.66	0/114	0.96	0/152
All	All	0.53	0/6482	0.84	12/8832 (0.1%)

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	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	58	GLY	CA-C-N	7.08	138.98	126.45
1	A	58	GLY	C-N-CA	7.08	138.98	126.45
1	B	58	GLY	CA-C-N	6.61	138.14	126.45
1	B	58	GLY	C-N-CA	6.61	138.14	126.45
1	A	295	ASP	CA-C-N	5.48	131.66	123.05
1	A	295	ASP	C-N-CA	5.48	131.66	123.05
1	B	154	TYR	CA-CB-CG	5.25	123.35	113.90
1	B	59	LYS	CA-C-N	5.11	130.57	122.61
1	B	59	LYS	C-N-CA	5.11	130.57	122.61
1	A	339	ASP	CA-C-N	5.05	132.74	124.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	339	ASP	C-N-CA	5.05	132.74	124.31
1	A	154	TYR	CA-CB-CG	5.01	122.91	113.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	55	ARG	Peptide
1	B	162	ASP	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	3042	0	2895	12	0
1	B	3046	0	2896	10	1
2	P	115	0	121	0	0
2	Q	115	0	121	0	0
3	P	8	0	7	0	0
3	Q	8	0	7	0	0
4	B	1	0	0	0	0
All	All	6335	0	6047	22	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 2.

All (22) 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:B:54:THR:HG23	1:B:63:ILE:HB	1.93	0.50
1:A:276:VAL:HA	1:A:293:SER:HB2	1.94	0.50
1:B:180:GLY:HA3	1:B:231:GLU:HA	1.95	0.49
1:B:276:VAL:HA	1:B:293:SER:HB2	1.95	0.49
1:A:176:HIS:O	1:A:177:GLN:NE2	2.45	0.49
1:A:338:THR:HG23	1:A:376:LYS:HD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:THR:HG23	1:B:376:LYS:HD2	1.97	0.47
1:B:323:VAL:O	1:B:324:GLN:NE2	2.48	0.46
1:A:122:ASN:OD1	1:A:122:ASN:N	2.49	0.45
1:A:293:SER:OG	1:A:294:ALA:N	2.49	0.45
1:A:54:THR:HG23	1:A:63:ILE:HB	1.99	0.44
1:A:63:ILE:HD13	1:A:85:GLN:HG2	2.00	0.43
1:B:63:ILE:HD13	1:B:85:GLN:HG2	2.00	0.43
1:A:238:LEU:HD12	1:A:286:GLU:HB3	2.00	0.43
1:B:68:LEU:HB2	1:B:80:VAL:HB	2.00	0.43
1:A:199:ASP:OD1	1:A:199:ASP:N	2.50	0.42
1:B:238:LEU:HD12	1:B:286:GLU:HB3	2.00	0.42
1:A:180:GLY:HA3	1:A:231:GLU:HA	2.01	0.42
1:A:322:GLN:NE2	1:A:379:ASP:OD1	2.48	0.42
1:B:293:SER:OG	1:B:294:ALA:N	2.52	0.41
1:B:182:GLY:HA3	1:B:232:ASP:HA	2.03	0.40
1:A:323:VAL:O	1:A:324:GLN:NE2	2.54	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)
1:B:167:CYS:CB	1:B:167:CYS:SG[2_556]	2.18	0.02

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	377/427 (88%)	363 (96%)	14 (4%)	0	100	100
1	B	377/427 (88%)	363 (96%)	14 (4%)	0	100	100
2	P	14/16 (88%)	13 (93%)	1 (7%)	0	100	100
2	Q	14/16 (88%)	14 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	782/886 (88%)	753 (96%)	29 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	343/376 (91%)	315 (92%)	28 (8%)	10	24
1	B	344/376 (92%)	316 (92%)	28 (8%)	11	24
2	P	11/11 (100%)	9 (82%)	2 (18%)	2	3
2	Q	11/11 (100%)	9 (82%)	2 (18%)	2	3
All	All	709/774 (92%)	649 (92%)	60 (8%)	10	22

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LYS
1	A	26	LYS
1	A	28	THR
1	A	53	VAL
1	A	54	THR
1	A	55	ARG
1	A	57	GLU
1	A	66	LEU
1	A	75	GLU
1	A	88	ASN
1	A	114	LYS
1	A	124	GLU
1	A	143	LYS
1	A	166	GLU
1	A	174	ARG
1	A	179	GLU
1	A	201	THR

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Mol	Chain	Res	Type
1	A	207	ILE
1	A	210	VAL
1	A	212	LYS
1	A	221	THR
1	A	224	THR
1	A	269	VAL
1	A	276	VAL
1	A	293	SER
1	A	331	THR
1	A	357	GLU
1	A	408	ILE
2	P	9	ARG
2	P	12	LYS
2	Q	9	ARG
2	Q	12	LYS
1	B	13	GLU
1	B	26	LYS
1	B	54	THR
1	B	55	ARG
1	B	57	GLU
1	B	66	LEU
1	B	75	GLU
1	B	88	ASN
1	B	114	LYS
1	B	139	ILE
1	B	143	LYS
1	B	160	LYS
1	B	166	GLU
1	B	177	GLN
1	B	179	GLU
1	B	183	LEU
1	B	201	THR
1	B	207	ILE
1	B	210	VAL
1	B	221	THR
1	B	224	THR
1	B	269	VAL
1	B	276	VAL
1	B	293	SER
1	B	331	THR
1	B	338	THR
1	B	357	GLU

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Mol	Chain	Res	Type
1	B	408	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	A	136	ASN
1	A	328	HIS
1	B	136	ASN
1	B	239	HIS
1	B	250	GLN
1	B	322	GLN
1	B	328	HIS

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](#)

2 ligands are modelled in this entry.

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
3	PXY	Q	101	2	8,8,8	0.77	0	10,10,10	1.34	1 (10%)
3	PXY	P	101	2	8,8,8	0.56	0	10,10,10	1.46	0

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
3	PXY	Q	101	2	-	-	0/1/1/1
3	PXY	P	101	2	-	-	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	101	PXY	C3-C2-C1	-2.02	119.05	121.37

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	381/427 (89%)	0.53	19 (4%) 34 28	92, 114, 141, 175	0
1	B	381/427 (89%)	0.51	14 (3%) 45 39	92, 112, 143, 168	0
2	P	14/16 (87%)	0.87	2 (14%) 6 5	109, 114, 119, 120	0
2	Q	14/16 (87%)	0.72	1 (7%) 22 17	110, 116, 120, 125	0
All	All	790/886 (89%)	0.53	36 (4%) 37 32	92, 114, 142, 175	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	177	GLN	3.9
1	A	290	ALA	3.8
1	A	306	LEU	3.5
2	P	15	ALA	3.1
1	B	252	LEU	3.0
1	B	69	GLY	2.9
1	B	119	ILE	2.9
1	A	119	ILE	2.8
1	A	128	ASN	2.7
1	A	72	THR	2.5
1	A	118	GLU	2.5
1	A	28	THR	2.5
1	A	281	PHE	2.4
1	A	291	THR	2.4
1	B	171	LEU	2.3
1	A	387	PRO	2.3
1	B	39	ALA	2.3
1	A	214	GLY	2.3
1	A	367	LEU	2.2
1	A	50	LEU	2.2
1	B	124	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	130	ALA	2.1
1	B	297	THR	2.1
1	A	141	ALA	2.1
2	P	6	ALA	2.1
2	Q	15	ALA	2.1
1	A	127	VAL	2.1
1	B	84	VAL	2.1
1	A	405	ALA	2.1
1	B	310	LEU	2.1
1	B	230	VAL	2.1
1	B	366	LEU	2.0
1	B	167	CYS	2.0
1	B	409	TYR	2.0
1	B	134	PRO	2.0
1	A	183	LEU	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	PXY	Q	101	8/8	0.94	0.10	125,127,129,130	0
3	PXY	P	101	8/8	0.96	0.16	118,121,125,125	0

6.5 Other polymers [i](#)

There are no such residues in this entry.