



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:04 PM UTC

PDB ID : 7JVX / pdb_00007jvx
Title : Crystal structure of PTEN (aa 7-353 followed by spacer TGGGSGGTGGGSG-GTGGGCY ligated to peptide pSDpTpTDpSDPENEPFDED)
Authors : Dempsey, D.; Phan, K.; Cole, P.; Gabelli, S.B.
Deposited on : 2020-08-24
Resolution : 3.20 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

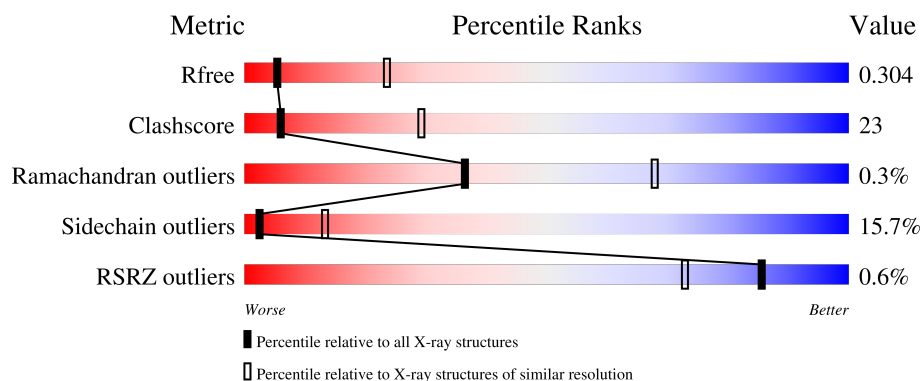
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	422	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	501	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2639 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.

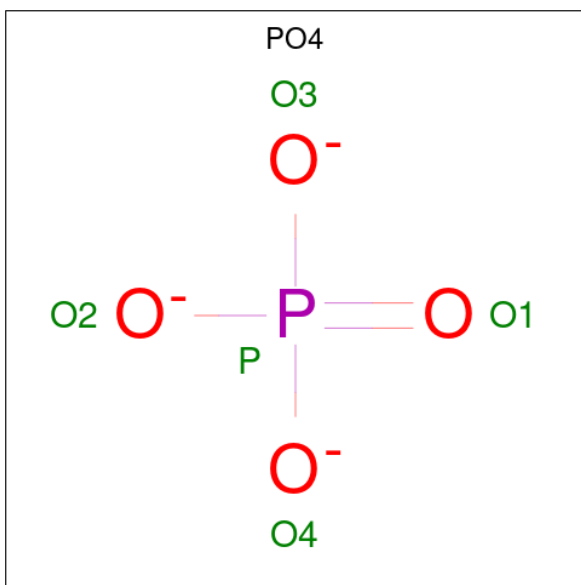
- Molecule 1 is a protein called Phosphatidylinositol 3,4,5-trisphosphate 3-phosphatase and dual-specificity protein phosphatase PTEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2629	1703	450	460	16			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	381	THR	-	insertion	UNP P60484
A	382	GLY	-	insertion	UNP P60484
A	383	GLY	-	insertion	UNP P60484
A	384	GLY	-	insertion	UNP P60484
A	385	SER	-	insertion	UNP P60484
A	386	GLY	-	insertion	UNP P60484
A	387	GLY	-	insertion	UNP P60484
A	388	THR	-	insertion	UNP P60484
A	389	GLY	-	insertion	UNP P60484
A	390	GLY	-	insertion	UNP P60484
A	391	GLY	-	insertion	UNP P60484
A	392	SER	-	insertion	UNP P60484
A	393	GLY	-	insertion	UNP P60484
A	394	GLY	-	insertion	UNP P60484
A	395	THR	-	insertion	UNP P60484
A	396	GLY	-	insertion	UNP P60484
A	397	GLY	-	insertion	UNP P60484
A	398	GLY	-	insertion	UNP P60484
A	399	CYS	-	insertion	UNP P60484

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

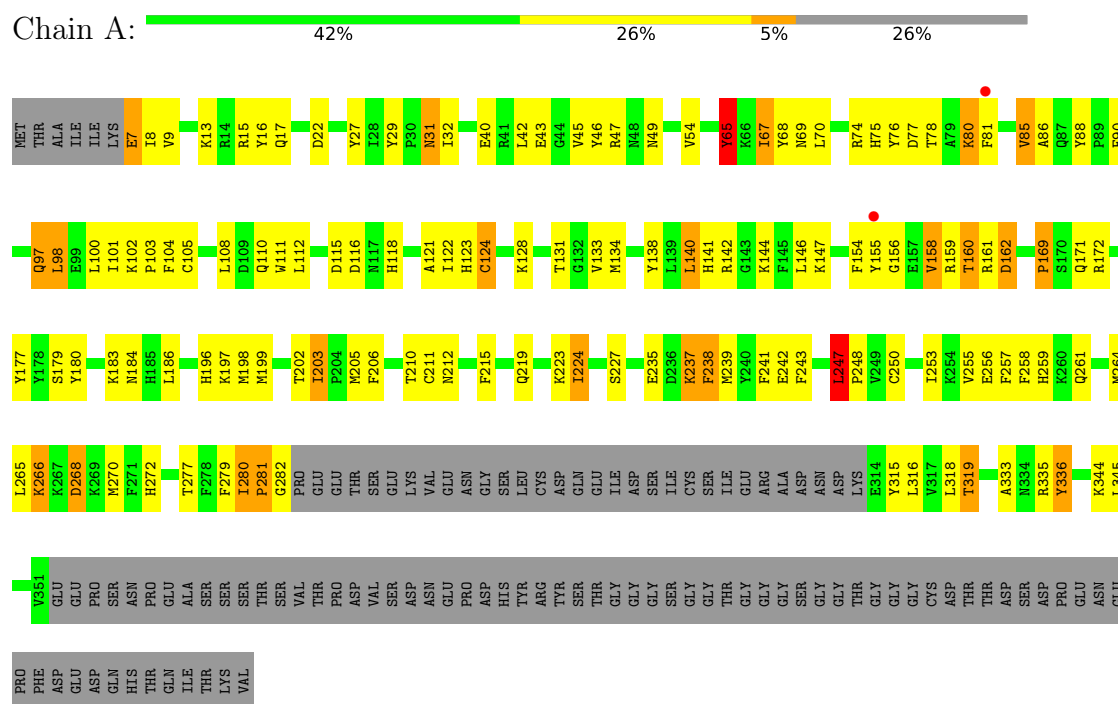


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

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: Phosphatidylinositol 3,4,5-trisphosphate 3-phosphatase and dual-specificity protein phosphatase PTEN



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	113.39Å 113.39Å 57.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.80 – 3.20 19.80 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.3 (19.80-3.20) 94.3 (19.80-3.20)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 3.22Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.244 , 0.296 0.252 , 0.304	Depositor DCC
R_{free} test set	314 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	81.7	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.056 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2639	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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: PO4

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	0.91	1/2703 (0.0%)	1.45	15/3644 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	169	PRO	C-O	-6.51	1.16	1.24

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	ASN	CA-C-N	9.69	130.07	119.90
1	A	212	ASN	C-N-CA	9.69	130.07	119.90
1	A	247	LEU	CA-C-N	8.95	131.03	119.84
1	A	247	LEU	C-N-CA	8.95	131.03	119.84
1	A	202	THR	CA-C-N	-7.57	116.75	123.33
1	A	202	THR	C-N-CA	-7.57	116.75	123.33
1	A	160	THR	CB-CA-C	-6.88	108.63	116.54
1	A	202	THR	CA-CB-OG1	-6.80	99.40	109.60
1	A	212	ASN	CA-C-O	-6.31	113.87	121.00
1	A	27	TYR	CB-CA-C	-5.71	101.48	110.79
1	A	65	TYR	CB-CA-C	5.67	119.67	110.37
1	A	315	TYR	CB-CA-C	-5.66	100.54	113.33
1	A	336	TYR	CB-CA-C	5.65	119.68	110.24
1	A	40	GLU	N-CA-C	-5.18	105.72	111.36
1	A	196	HIS	CA-CB-CG	-5.01	108.79	113.80

There are no chirality outliers.

There are no planarity outliers.

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	2629	0	2589	119	1
2	A	10	0	0	3	0
All	All	2639	0	2589	119	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 23.

All (119) 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:237:LYS:HE3	1:A:237:LYS:HA	1.19	1.08
1:A:54:VAL:HG22	1:A:81:PHE:CE1	1.92	1.05
1:A:67:ILE:HD12	1:A:67:ILE:H	1.24	1.02
1:A:154:PHE:O	1:A:158:VAL:HG23	1.62	0.99
1:A:43:GLU:H	1:A:46:TYR:HB2	1.38	0.88
1:A:237:LYS:HE3	1:A:237:LYS:CA	2.02	0.86
1:A:7:GLU:N	1:A:7:GLU:OE1	2.13	0.81
1:A:162:ASP:OD1	1:A:162:ASP:N	2.15	0.80
1:A:68:TYR:CD2	1:A:86:ALA:HB3	2.20	0.77
1:A:154:PHE:CE1	1:A:158:VAL:HG21	2.22	0.75
1:A:88:TYR:HD2	1:A:100:LEU:HD11	1.51	0.73
1:A:54:VAL:HG22	1:A:81:PHE:CD1	2.23	0.73
1:A:282:GLY:H	1:A:316:LEU:HD22	1.52	0.73
1:A:154:PHE:O	1:A:158:VAL:CG2	2.37	0.72
1:A:124:CYS:HG	1:A:131:THR:HG1	1.31	0.72
1:A:140:LEU:HD23	1:A:140:LEU:N	2.06	0.71
1:A:17:GLN:HA	1:A:22:ASP:HA	1.74	0.70
1:A:198:MET:HE1	1:A:257:PHE:CZ	2.29	0.67
1:A:177:TYR:CD1	1:A:279:PHE:CD2	2.82	0.67
1:A:54:VAL:CG2	1:A:81:PHE:CE1	2.75	0.65
1:A:261:GLN:HB3	1:A:264:MET:HB2	1.79	0.64
1:A:67:ILE:HD12	1:A:67:ILE:N	2.07	0.62
1:A:80:LYS:O	1:A:80:LYS:HE3	1.99	0.62
1:A:319:THR:HB	1:A:344:LYS:HG2	1.82	0.62
1:A:266:LYS:N	1:A:266:LYS:HD3	2.15	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:TYR:OH	1:A:85:VAL:HG11	2.01	0.60
1:A:133:VAL:CG2	1:A:171:GLN:HB3	2.32	0.60
1:A:197:LYS:HD3	1:A:242:GLU:HG2	1.84	0.60
1:A:146:LEU:HD12	1:A:146:LEU:N	2.15	0.59
1:A:197:LYS:HB3	1:A:242:GLU:HA	1.84	0.59
1:A:198:MET:HE1	1:A:257:PHE:CE2	2.37	0.59
1:A:247:LEU:HD23	1:A:248:PRO:HD2	1.85	0.58
1:A:80:LYS:HE3	1:A:80:LYS:HA	1.86	0.58
1:A:76:TYR:OH	1:A:85:VAL:CG1	2.51	0.58
1:A:197:LYS:HB2	1:A:241:PHE:O	2.04	0.57
1:A:100:LEU:HD12	1:A:100:LEU:O	2.04	0.57
1:A:67:ILE:H	1:A:67:ILE:CD1	2.03	0.57
1:A:88:TYR:HD2	1:A:100:LEU:CD1	2.18	0.57
1:A:68:TYR:CD2	1:A:86:ALA:CB	2.88	0.57
1:A:102:LYS:HB3	1:A:103:PRO:HD3	1.87	0.56
1:A:75:HIS:ND1	2:A:502:PO4:O3	2.36	0.56
1:A:54:VAL:HG22	1:A:81:PHE:CZ	2.40	0.56
1:A:65:TYR:CD2	1:A:67:ILE:HD11	2.41	0.56
1:A:237:LYS:HA	1:A:237:LYS:CE	2.13	0.55
1:A:264:MET:HB3	1:A:266:LYS:HE2	1.87	0.55
1:A:65:TYR:CD1	1:A:65:TYR:N	2.75	0.55
1:A:203:ILE:HD12	1:A:239:MET:HE2	1.88	0.55
1:A:215:PHE:HB3	1:A:257:PHE:CE1	2.43	0.54
1:A:279:PHE:CD1	1:A:279:PHE:N	2.75	0.54
1:A:90:PHE:CE1	1:A:134:MET:HE2	2.43	0.54
1:A:235:GLU:O	1:A:235:GLU:HG2	2.07	0.54
1:A:238:PHE:CD1	1:A:238:PHE:N	2.75	0.54
1:A:224:ILE:HG13	1:A:247:LEU:HD21	1.90	0.53
1:A:76:TYR:HE2	1:A:123:HIS:NE2	2.06	0.52
1:A:67:ILE:HA	1:A:121:ALA:O	2.10	0.52
1:A:156:GLY:O	1:A:159:ARG:O	2.28	0.52
1:A:219:GLN:NE2	1:A:250:CYS:O	2.42	0.52
1:A:133:VAL:HG21	1:A:171:GLN:HB3	1.92	0.52
1:A:256:GLU:HG2	1:A:258:PHE:CZ	2.46	0.51
1:A:74:ARG:NH1	2:A:501:PO4:O4	2.42	0.51
1:A:105:CYS:HB3	1:A:141:HIS:ND1	2.27	0.50
1:A:177:TYR:HB3	1:A:279:PHE:CZ	2.46	0.50
1:A:203:ILE:HD11	1:A:210:THR:HB	1.93	0.50
1:A:97:GLN:O	1:A:100:LEU:HB3	2.11	0.50
1:A:197:LYS:HA	1:A:243:PHE:CE1	2.46	0.50
1:A:256:GLU:HG2	1:A:258:PHE:CE1	2.46	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:LYS:HE3	1:A:80:LYS:CA	2.41	0.49
1:A:88:TYR:CD2	1:A:100:LEU:HD11	2.40	0.49
1:A:266:LYS:HD3	1:A:266:LYS:H	1.76	0.49
1:A:211:CYS:HA	1:A:259:HIS:CE1	2.47	0.49
1:A:250:CYS:HA	1:A:277:THR:HG21	1.93	0.49
1:A:280:ILE:CD1	1:A:280:ILE:N	2.76	0.49
1:A:268:ASP:N	1:A:268:ASP:OD1	2.43	0.48
1:A:13:LYS:HA	1:A:160:THR:HA	1.95	0.48
1:A:70:LEU:HD11	1:A:134:MET:HG3	1.96	0.48
1:A:115:ASP:HB3	1:A:118:HIS:CG	2.49	0.47
1:A:333:ALA:HB3	1:A:335:ARG:HG3	1.97	0.47
1:A:122:ILE:HG22	1:A:131:THR:HG23	1.97	0.47
1:A:128:LYS:O	1:A:155:TYR:OH	2.31	0.47
1:A:76:TYR:C	1:A:76:TYR:CD1	2.93	0.46
1:A:318:LEU:C	1:A:318:LEU:HD13	2.41	0.46
1:A:76:TYR:HE2	1:A:123:HIS:CE1	2.34	0.45
1:A:206:PHE:CE1	1:A:336:TYR:HB3	2.51	0.45
1:A:140:LEU:CD1	1:A:179:SER:HB2	2.45	0.45
1:A:255:VAL:O	1:A:272:HIS:HA	2.17	0.45
1:A:105:CYS:HB3	1:A:141:HIS:CG	2.51	0.45
1:A:238:PHE:H	1:A:238:PHE:HD1	1.65	0.45
1:A:243:PHE:CD1	1:A:243:PHE:N	2.84	0.45
1:A:65:TYR:N	1:A:65:TYR:HD1	2.13	0.45
1:A:198:MET:HE1	1:A:257:PHE:HZ	1.78	0.45
1:A:203:ILE:HD11	1:A:210:THR:HA	1.99	0.45
1:A:177:TYR:CD1	1:A:279:PHE:CE2	3.05	0.44
1:A:88:TYR:CE2	1:A:104:PHE:HA	2.52	0.44
1:A:256:GLU:HA	1:A:272:HIS:HB3	2.00	0.44
1:A:154:PHE:C	1:A:154:PHE:CD1	2.97	0.43
1:A:180:TYR:O	1:A:184:ASN:ND2	2.52	0.43
1:A:280:ILE:N	1:A:280:ILE:HD12	2.32	0.43
1:A:98:LEU:HD13	1:A:98:LEU:C	2.44	0.43
1:A:280:ILE:N	1:A:281:PRO:HD2	2.34	0.43
1:A:203:ILE:HD12	1:A:239:MET:CE	2.49	0.42
1:A:29:TYR:HB2	1:A:32:ILE:HD12	2.00	0.42
1:A:280:ILE:N	1:A:281:PRO:CD	2.82	0.42
1:A:184:ASN:CB	1:A:186:LEU:HG	2.51	0.41
1:A:16:TYR:O	1:A:17:GLN:C	2.63	0.41
1:A:31:ASN:HB2	1:A:112:LEU:HD13	2.02	0.41
1:A:74:ARG:HH11	2:A:501:PO4:P	2.43	0.41
1:A:177:TYR:HB3	1:A:279:PHE:CE2	2.56	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:LYS:O	1:A:223:LYS:HG2	2.19	0.41
1:A:69:ASN:OD1	1:A:69:ASN:C	2.63	0.41
1:A:112:LEU:HD23	1:A:112:LEU:HA	1.84	0.41
1:A:80:LYS:HE3	1:A:80:LYS:C	2.45	0.41
1:A:140:LEU:HD12	1:A:179:SER:HB2	2.02	0.41
1:A:169:PRO:HA	1:A:172:ARG:HB2	2.01	0.41
1:A:177:TYR:O	1:A:180:TYR:HB3	2.21	0.41
1:A:248:PRO:HG2	1:A:248:PRO:O	2.22	0.40
1:A:78:THR:OG1	1:A:85:VAL:HG21	2.22	0.40
1:A:76:TYR:CD1	1:A:76:TYR:O	2.75	0.40
1:A:138:TYR:CD1	1:A:138:TYR:O	2.75	0.40
1:A:335:ARG:HE	1:A:335:ARG:HB3	1.50	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:A:47:ARG:NH1	1:A:319:THR:O[6_445]	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	310/422 (74%)	281 (91%)	28 (9%)	1 (0%)	36 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	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	286/377 (76%)	241 (84%)	45 (16%)	2 13

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	8	ILE
1	A	9	VAL
1	A	15	ARG
1	A	31	ASN
1	A	42	LEU
1	A	45	VAL
1	A	49	ASN
1	A	65	TYR
1	A	67	ILE
1	A	77	ASP
1	A	80	LYS
1	A	85	VAL
1	A	97	GLN
1	A	98	LEU
1	A	101	ILE
1	A	108	LEU
1	A	110	GLN
1	A	111	TRP
1	A	116	ASP
1	A	124	CYS
1	A	140	LEU
1	A	142	ARG
1	A	144	LYS
1	A	147	LYS
1	A	158	VAL
1	A	162	ASP
1	A	183	LYS
1	A	199	MET
1	A	203	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	205	MET
1	A	224	ILE
1	A	227	SER
1	A	237	LYS
1	A	238	PHE
1	A	247	LEU
1	A	253	ILE
1	A	265	LEU
1	A	266	LYS
1	A	268	ASP
1	A	270	MET
1	A	280	ILE
1	A	281	PRO
1	A	319	THR
1	A	345	LEU

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	110	GLN
1	A	184	ASN
1	A	261	GLN

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 ⓘ

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	$\# Z > 2$
2	PO4	A	502	-	4,4,4	0.84	0	6,6,6	0.44	0
2	PO4	A	501	-	4,4,4	0.75	0	6,6,6	0.45	0

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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	PO4	1	0
2	A	501	PO4	2	0

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	314/422 (74%)	-0.10	2 (0%) 85 73	47, 83, 138, 174	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	PHE	3.2
1	A	155	TYR	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 [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(Å ²)	Q<0.9
2	PO4	A	502	5/5	0.71	0.15	110,110,110,110	0
2	PO4	A	501	5/5	0.74	0.10	123,129,134,135	0

6.5 Other polymers [i](#)

There are no such residues in this entry.