



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:08 PM UTC

PDB ID : 4TMN / pdb_00004tmn
Title : SLOW-AND FAST-BINDING INHIBITORS OF THERMOLYSIN DISPLAY
DIFFERENT MODES OF BINDING. CRYSTALLOGRAPHIC ANALY-
SIS OF EXTENDED PHOSPHONAMIDATE TRANSITION-STATE ANA-
LOGUES
Authors : Holden, H.M.; Tronrud, D.E.; Monzingo, A.F.; Weaver, L.H.; Matthews, B.W.
Deposited on : 1987-06-29
Resolution : 1.70 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](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

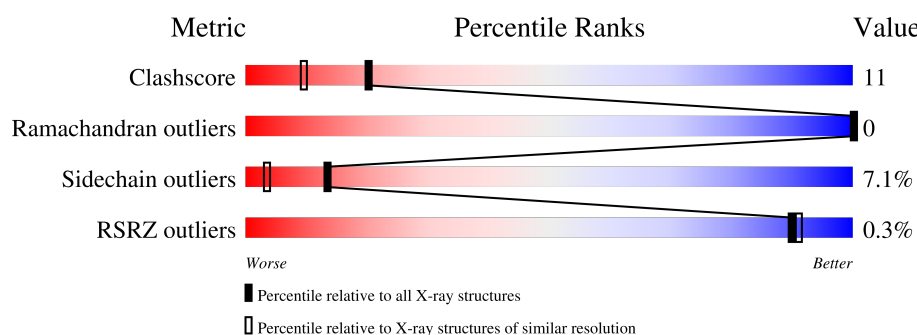
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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))
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	E	316	

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	OPK	E	317	-	-	X	-

2 Entry composition [i](#)

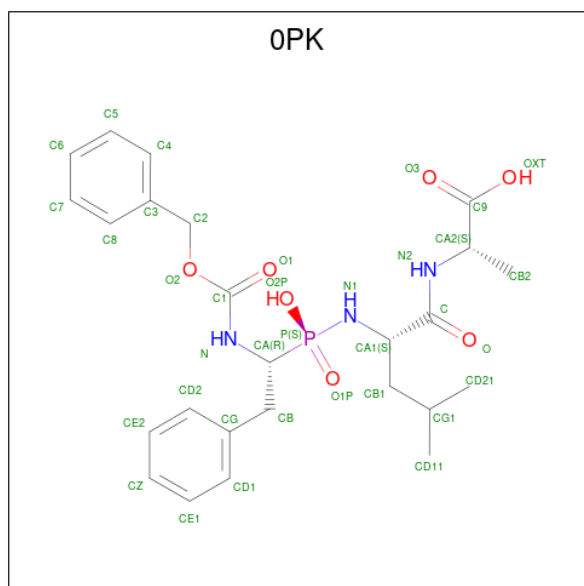
There are 5 unique types of molecules in this entry. The entry contains 2635 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 THERMOLYSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	316	Total	C	N	O	S	0	0	0
			2432	1528	408	494	2			

- Molecule 2 is N-[(S)-[(1R)-1-{[(benzyloxy)carbonyl]amino}-2-phenylethyl](hydroxy)phosphoryl]-L-leucyl-L-alanine (CCD ID: 0PK) (formula: C₂₅H₃₄N₃O₇P).



- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	1	Total 1	Zn 1	0	0

- Molecule 5 is water.

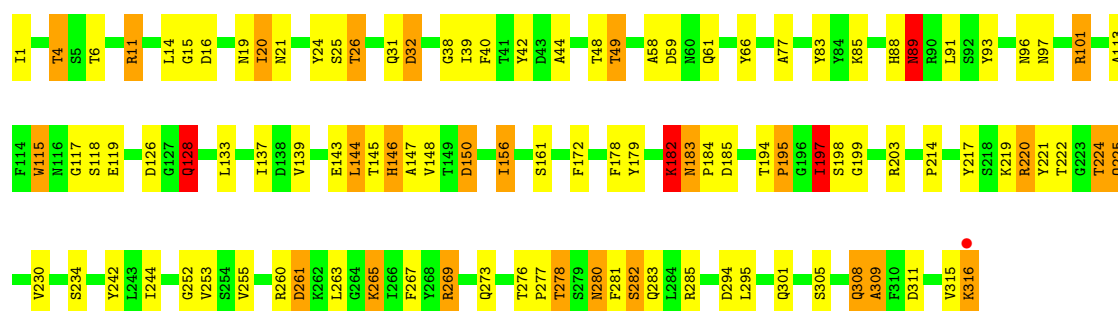
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	162	Total 162	O 162	0	0

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: THERMOLYSIN

Chain E: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	94.10Å 94.10Å 131.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 1.70 81.49 – 1.71	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-1.70) 75.8 (81.49-1.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 1.71Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.170 , (Not available) 0.157 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	15.4	Xtriage
Anisotropy	0.548	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 73.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2635	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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: 0PK, ZN, CA

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	E	1.80	33/2491 (1.3%)	2.01	86/3391 (2.5%)

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	E	2	0

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	203	ARG	NE-CZ	7.77	1.41	1.33
1	E	261	ASP	CG-OD2	7.05	1.38	1.25
1	E	150	ASP	CG-OD2	6.84	1.38	1.25
1	E	183	ASN	C-O	-6.59	1.20	1.23
1	E	145	THR	C-O	-6.35	1.16	1.24
1	E	242	TYR	CA-C	6.24	1.60	1.52
1	E	252	GLY	C-N	6.17	1.41	1.33
1	E	308	GLN	CD-OE1	6.14	1.35	1.23
1	E	139	VAL	CA-C	5.89	1.60	1.52
1	E	15	GLY	CA-C	5.83	1.59	1.51
1	E	119	GLU	CD-OE2	5.69	1.36	1.25
1	E	316	LYS	C-OXT	5.67	1.34	1.23
1	E	267	PHE	CA-C	5.58	1.60	1.52
1	E	260	ARG	C-N	-5.56	1.26	1.33
1	E	88	HIS	CA-C	-5.55	1.45	1.52
1	E	128	GLN	C-N	-5.51	1.25	1.33
1	E	172	PHE	C-O	-5.41	1.17	1.24
1	E	263	LEU	CA-C	5.40	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	96	ASN	CG-OD1	5.38	1.33	1.23
1	E	260	ARG	CA-CB	-5.35	1.44	1.53
1	E	261	ASP	CA-CB	-5.32	1.45	1.53
1	E	253	VAL	N-CA	5.25	1.52	1.46
1	E	147	ALA	N-CA	5.23	1.52	1.46
1	E	39	ILE	C-O	-5.23	1.18	1.24
1	E	146	HIS	CG-ND1	5.22	1.44	1.38
1	E	267	PHE	C-O	5.22	1.30	1.24
1	E	198	SER	C-O	-5.18	1.17	1.23
1	E	11	ARG	CA-C	-5.17	1.46	1.52
1	E	19	ASN	CA-C	-5.13	1.46	1.52
1	E	91	LEU	C-O	5.13	1.29	1.23
1	E	83	TYR	CA-CB	5.12	1.61	1.53
1	E	185	ASP	CA-C	-5.11	1.46	1.52
1	E	77	ALA	C-N	5.07	1.40	1.33

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	269	ARG	CD-NE-CZ	-13.22	105.89	124.40
1	E	260	ARG	N-CA-C	12.65	125.15	111.36
1	E	101	ARG	CG-CD-NE	-10.70	88.46	112.00
1	E	269	ARG	NE-CZ-NH1	-10.44	111.06	121.50
1	E	133	LEU	N-CA-C	9.13	122.21	111.71
1	E	316	LYS	CB-CA-C	8.83	126.88	110.10
1	E	269	ARG	NE-CZ-NH2	8.62	126.95	119.20
1	E	6	THR	OG1-CB-CG2	-8.35	92.61	109.30
1	E	118	SER	N-CA-CB	8.07	124.12	110.49
1	E	58	ALA	N-CA-C	8.04	120.12	111.36
1	E	278	THR	N-CA-CB	-8.00	98.13	110.92
1	E	261	ASP	N-CA-CB	-7.96	98.53	110.07
1	E	61	GLN	CA-CB-CG	-7.63	98.85	114.10
1	E	49	THR	N-CA-CB	-7.53	98.27	109.95
1	E	183	ASN	CB-CA-C	-7.42	104.55	111.00
1	E	265	LYS	CA-CB-CG	-7.41	99.28	114.10
1	E	96	ASN	CB-CG-ND2	-7.30	105.45	116.40
1	E	156	ILE	CA-CB-CG2	7.26	122.84	110.50
1	E	148	VAL	N-CA-CB	7.00	117.98	110.62
1	E	49	THR	CB-CA-C	6.94	121.15	109.84
1	E	26	THR	CA-CB-OG1	6.91	119.97	109.60
1	E	89	ASN	CA-CB-CG	-6.83	105.77	112.60
1	E	126	ASP	CB-CA-C	-6.78	99.11	110.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	224	THR	CA-CB-CG2	6.71	121.91	110.50
1	E	234	SER	N-CA-C	-6.46	104.66	112.54
1	E	220	ARG	NE-CZ-NH2	-6.43	113.41	119.20
1	E	222	THR	CA-C-O	6.42	126.12	119.05
1	E	295	LEU	N-CA-C	6.35	119.50	111.69
1	E	278	THR	CA-CB-CG2	6.34	121.27	110.50
1	E	93	TYR	N-CA-C	6.26	119.08	111.82
1	E	261	ASP	CB-CA-C	-6.21	101.09	110.90
1	E	182	LYS	CG-CD-CE	6.11	125.34	111.30
1	E	197	ILE	CA-CB-CG1	6.04	120.67	110.40
1	E	224	THR	OG1-CB-CG2	6.01	121.31	109.30
1	E	282	SER	CB-CA-C	5.99	121.04	110.85
1	E	198	SER	CB-CA-C	-5.96	98.50	109.46
1	E	4	THR	OG1-CB-CG2	-5.93	97.43	109.30
1	E	224	THR	N-CA-CB	-5.91	102.45	110.67
1	E	66	TYR	CA-C-N	5.81	128.54	120.29
1	E	66	TYR	C-N-CA	5.81	128.54	120.29
1	E	26	THR	OG1-CB-CG2	5.76	120.83	109.30
1	E	144	LEU	CD1-CG-CD2	-5.76	98.12	110.80
1	E	20	ILE	CA-CB-CG1	-5.73	100.66	110.40
1	E	308	GLN	O-C-N	5.73	127.97	122.07
1	E	96	ASN	OD1-CG-ND2	5.71	128.31	122.60
1	E	278	THR	OG1-CB-CG2	5.70	120.71	109.30
1	E	316	LYS	N-CA-CB	5.69	120.17	110.50
1	E	21	ASN	OD1-CG-ND2	-5.67	116.93	122.60
1	E	301	GLN	N-CA-C	-5.66	105.26	111.82
1	E	179	TYR	CB-CA-C	-5.65	101.42	110.79
1	E	101	ARG	CD-NE-CZ	5.63	132.28	124.40
1	E	260	ARG	NE-CZ-NH1	5.63	127.13	121.50
1	E	273	GLN	CB-CA-C	-5.61	99.17	109.15
1	E	26	THR	N-CA-CB	5.55	119.87	110.49
1	E	309	ALA	O-C-N	-5.53	116.37	122.07
1	E	273	GLN	CB-CG-CD	5.50	121.95	112.60
1	E	220	ARG	CD-NE-CZ	5.45	132.03	124.40
1	E	101	ARG	CB-CA-C	-5.42	99.01	109.37
1	E	16	ASP	CA-CB-CG	5.39	117.99	112.60
1	E	281	PHE	CA-CB-CG	5.34	119.14	113.80
1	E	115	TRP	N-CA-C	-5.34	99.81	108.52
1	E	14	LEU	CD1-CG-CD2	-5.33	99.06	110.80
1	E	32	ASP	CA-CB-CG	5.32	117.92	112.60
1	E	185	ASP	O-C-N	5.29	128.54	123.04
1	E	195	PRO	CA-C-N	-5.28	111.06	121.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	195	PRO	C-N-CA	-5.28	111.06	121.41
1	E	48	THR	CA-C-O	5.26	125.58	119.32
1	E	183	ASN	OD1-CG-ND2	5.26	127.86	122.60
1	E	230	VAL	CA-CB-CG1	5.25	119.32	110.40
1	E	161	SER	CA-C-O	5.24	126.00	119.97
1	E	44	ALA	O-C-N	-5.24	115.14	122.37
1	E	214	PRO	N-CA-CB	5.21	107.94	103.31
1	E	59	ASP	N-CA-C	5.17	120.23	112.94
1	E	25	SER	CA-CB-OG	-5.16	100.78	111.10
1	E	315	VAL	CA-C-N	5.15	130.98	121.70
1	E	315	VAL	C-N-CA	5.15	130.98	121.70
1	E	199	GLY	CA-C-O	-5.12	113.34	119.02
1	E	311	ASP	N-CA-C	-5.12	105.70	111.28
1	E	278	THR	CA-CB-OG1	5.12	117.27	109.60
1	E	26	THR	CA-CB-CG2	5.09	119.16	110.50
1	E	305	SER	CA-CB-OG	-5.05	101.00	111.10
1	E	101	ARG	CA-CB-CG	-5.02	104.06	114.10
1	E	113	ALA	N-CA-C	-5.02	101.11	108.99
1	E	117	GLY	N-CA-C	-5.01	108.70	115.21
1	E	197	ILE	CB-CA-C	-5.01	103.66	110.77
1	E	145	THR	N-CA-C	-5.01	106.01	111.82

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	E	26	THR	CB
1	E	278	THR	CB

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	E	2432	0	2267	31	0
2	E	36	0	32	24	0
3	E	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	1	0	0	0	0
5	E	162	0	0	4	2
All	All	2635	0	2299	54	2

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 11.

All (54) 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)
2:E:317:0PK:HD23	2:E:317:0PK:CD21	0.97	1.15
2:E:317:0PK:CD21	2:E:317:0PK:HD21	0.97	1.14
2:E:317:0PK:H71	5:E:481:HOH:O	1.47	1.13
2:E:317:0PK:CD11	2:E:317:0PK:HD13	0.97	1.12
2:E:317:0PK:CD11	2:E:317:0PK:HD11	0.97	1.09
2:E:317:0PK:CD11	2:E:317:0PK:HD12	0.97	1.09
2:E:317:0PK:CD21	2:E:317:0PK:HD22	0.97	1.04
1:E:285:ARG:HD3	1:E:316:LYS:HD3	1.43	0.97
2:E:317:0PK:HD22	2:E:317:0PK:CG1	2.05	0.86
2:E:317:0PK:HD23	2:E:317:0PK:CG1	2.05	0.86
2:E:317:0PK:HD21	2:E:317:0PK:CG1	2.05	0.85
2:E:317:0PK:HD23	2:E:317:0PK:HD22	1.58	0.85
2:E:317:0PK:HD23	2:E:317:0PK:HD21	1.58	0.85
2:E:317:0PK:HD13	2:E:317:0PK:CG1	2.08	0.84
2:E:317:0PK:HD11	2:E:317:0PK:CG1	2.08	0.84
2:E:317:0PK:HD11	2:E:317:0PK:HD12	1.58	0.83
2:E:317:0PK:HD13	2:E:317:0PK:HD12	1.58	0.83
2:E:317:0PK:HD13	2:E:317:0PK:HD11	1.58	0.83
2:E:317:0PK:HD21	2:E:317:0PK:HD22	1.58	0.83
2:E:317:0PK:HD12	2:E:317:0PK:CG1	2.08	0.82
1:E:269:ARG:NH1	1:E:294:ASP:OD2	2.18	0.75
1:E:4:THR:HG22	1:E:24:TYR:HB3	1.68	0.73
1:E:137:ILE:H	1:E:182:LYS:HZ2	1.42	0.66
1:E:221:TYR:OH	1:E:225:GLN:HG3	1.97	0.65
1:E:280:ASN:ND2	1:E:283:GLN:H	1.97	0.63
1:E:42:TYR:HE2	1:E:101:ARG:HG2	1.66	0.61
2:E:317:0PK:C7	5:E:481:HOH:O	2.25	0.57
1:E:280:ASN:HD22	1:E:280:ASN:C	2.17	0.53
1:E:280:ASN:HD22	1:E:283:GLN:H	1.56	0.52
1:E:115:TRP:NE1	1:E:150:ASP:OD2	2.38	0.51
1:E:31:GLN:HG3	1:E:40:PHE:CE1	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:282:SER:HA	1:E:316:LYS:HD2	1.93	0.49
1:E:143:GLU:OE1	2:E:317:OPK:O2P	2.29	0.49
1:E:137:ILE:H	1:E:182:LYS:NZ	2.08	0.49
1:E:32:ASP:O	1:E:38:GLY:HA2	2.13	0.48
1:E:217:TYR:O	1:E:220:ARG:HB2	2.13	0.48
1:E:42:TYR:CE2	1:E:101:ARG:HG2	2.48	0.48
1:E:269:ARG:HH11	1:E:269:ARG:HD3	1.04	0.46
2:E:317:OPK:HD22	2:E:317:OPK:CB1	2.47	0.45
2:E:317:OPK:HD23	2:E:317:OPK:CB1	2.47	0.45
1:E:194:THR:HA	1:E:195:PRO:HD2	1.73	0.45
1:E:156:ILE:HD13	1:E:156:ILE:HG21	1.75	0.44
1:E:128:GLN:O	1:E:195:PRO:HD2	2.17	0.44
1:E:137:ILE:CG2	1:E:182:LYS:HD3	2.49	0.43
1:E:244:ILE:HG13	1:E:309:ALA:CB	2.49	0.42
1:E:276:THR:HB	1:E:277:PRO:CD	2.49	0.42
1:E:85:LYS:O	1:E:89:ASN:HA	2.20	0.42
1:E:276:THR:HB	1:E:277:PRO:HD2	2.01	0.42
1:E:255:VAL:HG22	1:E:308:GLN:HB3	2.02	0.42
2:E:317:OPK:C8	5:E:481:HOH:O	2.62	0.41
1:E:197:ILE:N	1:E:197:ILE:CD1	2.84	0.41
1:E:219:LYS:HE2	5:E:370:HOH:O	2.21	0.41
1:E:178:PHE:CE1	1:E:184:PRO:HB2	2.56	0.40
1:E:143:GLU:O	1:E:146:HIS:HB2	2.21	0.40

All (2) 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)
5:E:433:HOH:O	5:E:433:HOH:O[7_555]	1.59	0.61
5:E:385:HOH:O	5:E:385:HOH:O[12_565]	2.02	0.18

5.3 Torsion angles

5.3.1 Protein backbone

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	E	314/316 (99%)	301 (96%)	13 (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	E	252/252 (100%)	234 (93%)	18 (7%)	13	3

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	1	ILE
1	E	11	ARG
1	E	20	ILE
1	E	26	THR
1	E	49	THR
1	E	89	ASN
1	E	97	ASN
1	E	128	GLN
1	E	144	LEU
1	E	182	LYS
1	E	183	ASN
1	E	197	ILE
1	E	224	THR
1	E	225	GLN
1	E	261	ASP
1	E	265	LYS
1	E	278	THR
1	E	280	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	21	ASN
1	E	97	ASN
1	E	158	GLN
1	E	159	ASN
1	E	183	ASN
1	E	246	GLN
1	E	280	ASN
1	E	290	GLN
1	E	301	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](#)

Of 6 ligands modelled in this entry, 5 are monoatomic - leaving 1 for Mogul analysis.

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	OPK	E	317	4	35,37,37	1.55	9 (25%)	42,50,50	2.86	17 (40%)

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
2	0PK	E	317	4	-	1/35/40/40	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	317	0PK	P-O1P	3.63	1.58	1.46
2	E	317	0PK	O1-C1	3.59	1.28	1.21
2	E	317	0PK	CB-CA	-2.96	1.49	1.54
2	E	317	0PK	CA-N	2.91	1.49	1.46
2	E	317	0PK	C6-C5	-2.26	1.33	1.38
2	E	317	0PK	C8-C3	-2.19	1.34	1.38
2	E	317	0PK	CZ-CE2	2.14	1.42	1.38
2	E	317	0PK	CA2-C9	2.11	1.55	1.52
2	E	317	0PK	C5-C4	2.05	1.42	1.38

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	317	0PK	C6-C7-C8	9.88	132.43	120.24
2	E	317	0PK	C7-C6-C5	-7.19	110.03	119.87
2	E	317	0PK	CB-CA-N	-6.16	104.18	111.36
2	E	317	0PK	OXT-C9-O3	-4.73	113.35	124.08
2	E	317	0PK	C2-C3-C4	3.67	129.11	120.64
2	E	317	0PK	OXT-C9-CA2	3.65	125.88	113.84
2	E	317	0PK	CG-CB-CA	-3.35	106.03	113.34
2	E	317	0PK	O2-C2-C3	-3.19	101.66	109.40
2	E	317	0PK	C7-C8-C3	-3.13	116.21	120.61
2	E	317	0PK	CD21-CG1-CB1	-3.10	99.95	111.08
2	E	317	0PK	O2P-P-CA	2.69	112.54	105.06
2	E	317	0PK	CD2-CG-CD1	2.41	121.82	118.23
2	E	317	0PK	CB-CG-CD1	-2.31	116.60	120.90
2	E	317	0PK	C8-C3-C4	-2.24	114.90	118.23
2	E	317	0PK	C5-C4-C3	2.21	123.73	120.61
2	E	317	0PK	C2-O2-C1	2.17	120.81	115.93
2	E	317	0PK	C2-C3-C8	-2.06	115.90	120.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

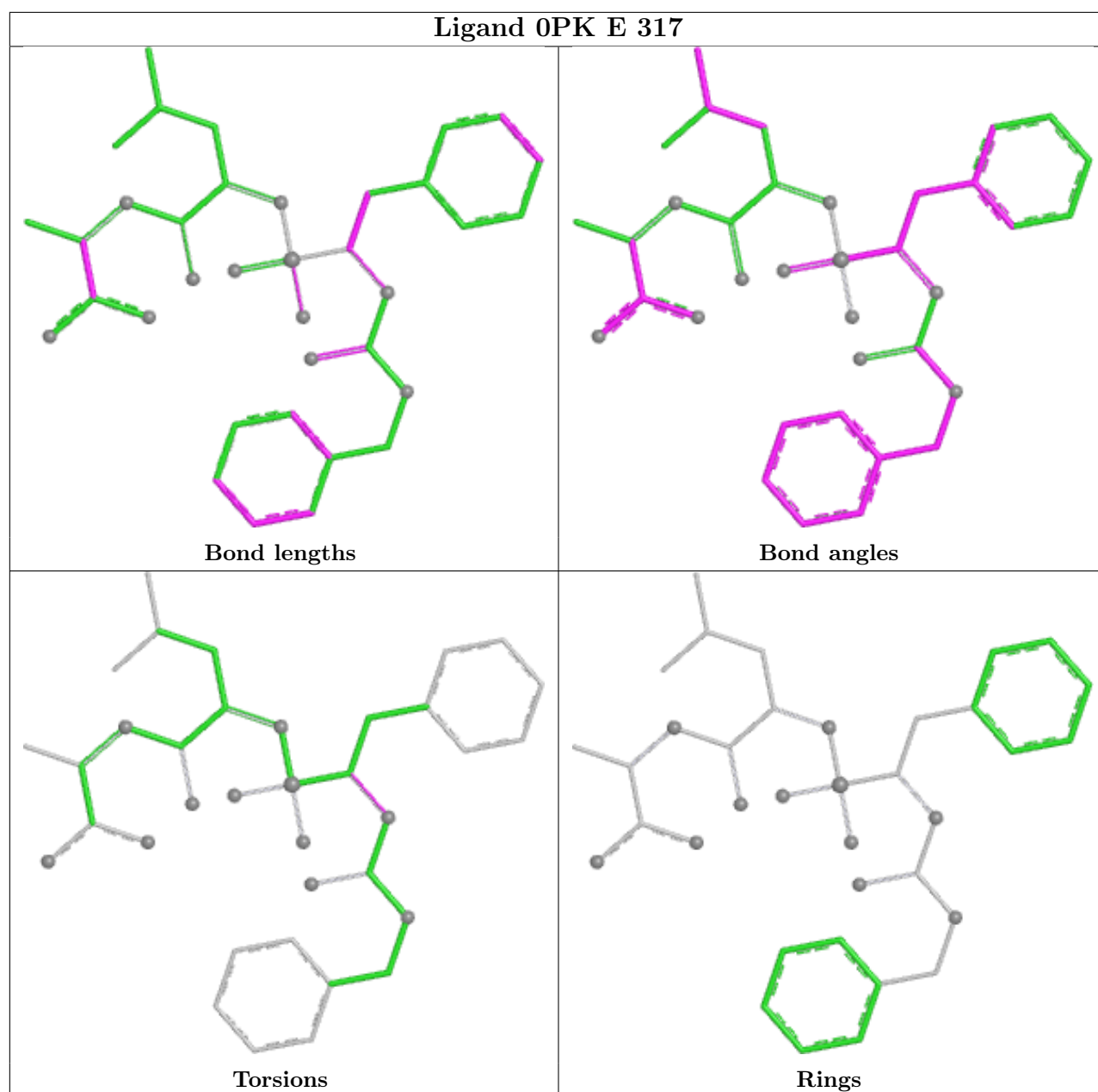
Mol	Chain	Res	Type	Atoms
2	E	317	0PK	P-CA-N-C1

There are no ring outliers.

1 monomer is involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	317	0PK	24	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	E	316/316 (100%)	-0.57	1 (0%) 90 91	7, 14, 27, 38	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	316	LYS	2.2

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](#)

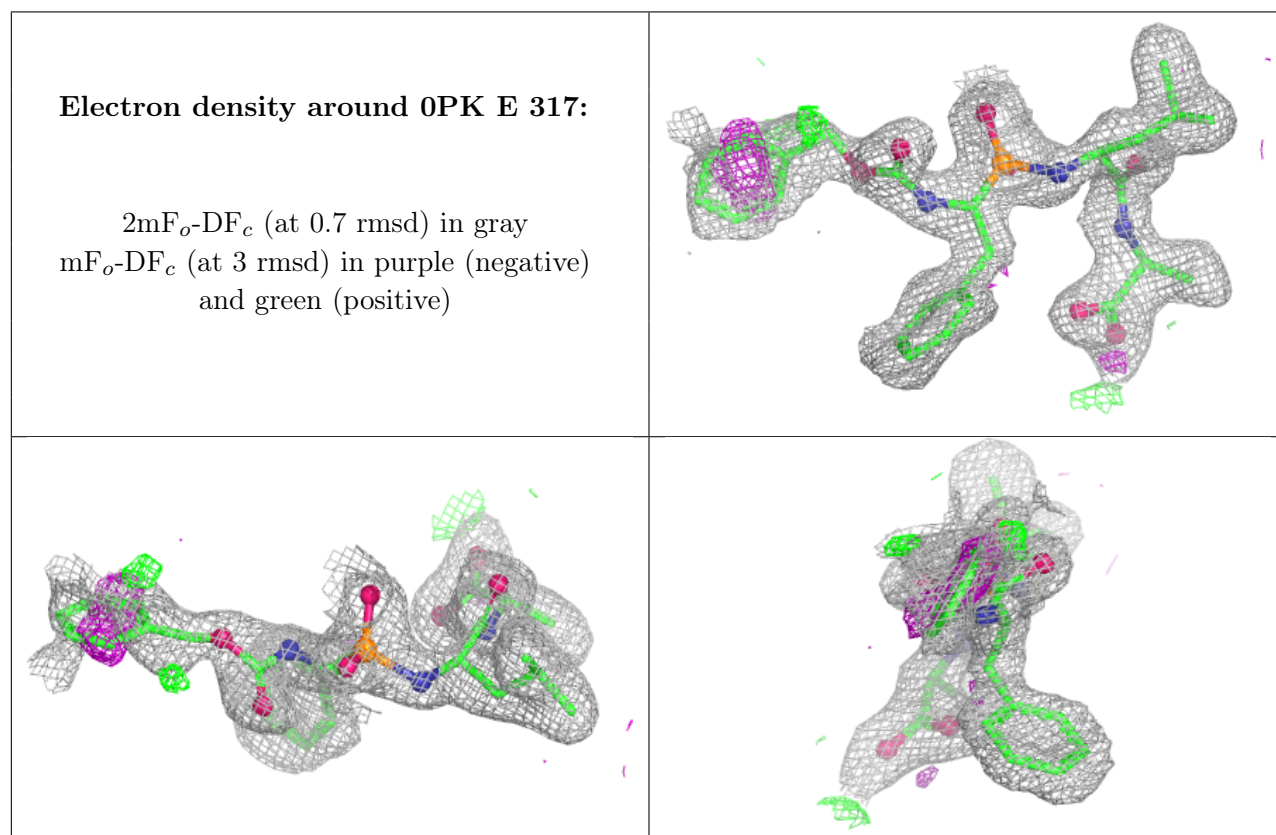
There are no oligosaccharides in this entry.

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	OPK	E	317	36/36	0.94	0.09	10,21,37,45	0
3	CA	E	318	1/1	0.99	0.02	12,12,12,12	0
3	CA	E	319	1/1	0.99	0.01	14,14,14,14	0
3	CA	E	320	1/1	0.99	0.05	11,11,11,11	0
3	CA	E	321	1/1	0.99	0.04	17,17,17,17	0
4	ZN	E	322	1/1	1.00	0.01	12,12,12,12	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.