



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:55 PM GMT

PDB ID : 1YDT  
Title : STRUCTURE OF CAMP-DEPENDENT PROTEIN KINASE, ALPHA-CATALYTIC SUBUNIT IN COMPLEX WITH H89 PROTEIN KINASE INHIBITOR N-[2-(4-BROMOCINNAMYLAMINO)ETHYL]-5-ISOQUINOLIN E  
Authors : Engh, R.A.; Girod, A.; Kinzel, V.; Huber, R.; Bossemeyer, D.  
Deposited on : 1996-07-24  
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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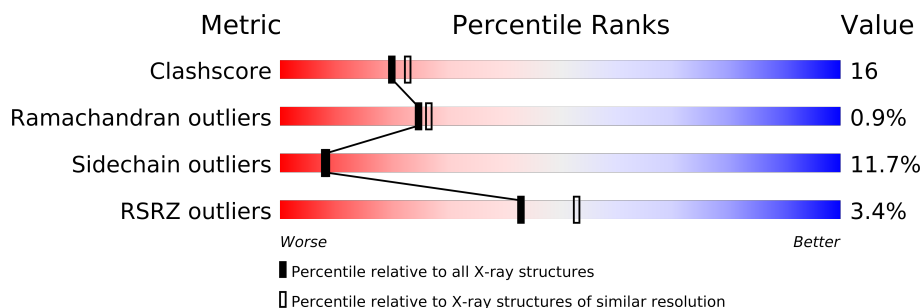
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.30 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	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	E	350	
2	I	20	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2983 atoms, of which 0 are hydrogen and 0 are deuterium.

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 C-AMP-DEPENDENT PROTEIN KINASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	E	336	2778	1801	465	501	2	9	0	0	0

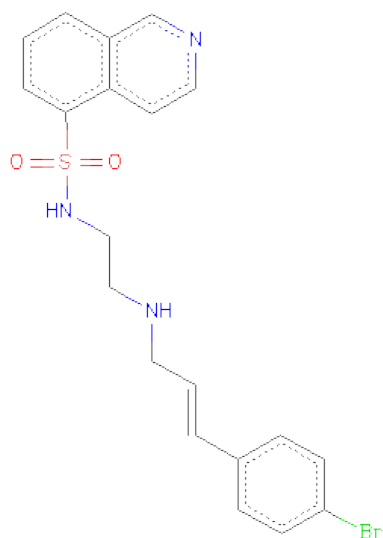
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	124	ALA	PRO	CONFLICT	UNP P00517
E	197	TPO	THR	MODIFIED RESIDUE	UNP P00517
E	286	ASP	ASN	CONFLICT	UNP P00517
E	338	SEP	SER	MODIFIED RESIDUE	UNP P00517

- Molecule 2 is a protein called PROTEIN KINASE INHIBITOR PEPTIDE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	I	20	157	94	32	31	0	0	0

- Molecule 3 is N-[2-(4-BROMOCINNAMYLAMINO)ETHYL]-5-ISOQUINOLINESULFONAMIDE (three-letter code: IQB) (formula: C<sub>20</sub>H<sub>20</sub>BrN<sub>3</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	N	O	S		
3	E	1	27	1	20	3	2	1	0	0

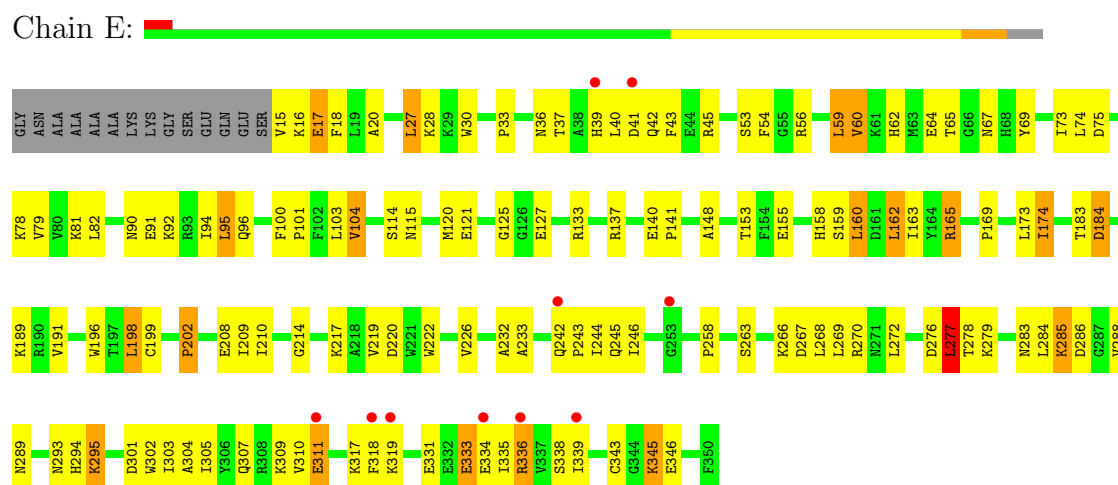
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	18	Total	O	0	0
			18	18		
4	I	3	Total	O	0	0
			3	3		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: C-AMP-DEPENDENT PROTEIN KINASE



- Molecule 2: PROTEIN KINASE INHIBITOR PEPTIDE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.58Å 76.28Å 80.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.30 19.48 – 1.99	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.30) 62.7 (19.48-1.99)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.98 (at 1.99Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.194 , (Not available) 0.228 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	26.7	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 37.5	EDS
Estimated twinning fraction	0.026 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 20066 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2983	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, IQB, SEP

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	0.67	1/2827 (0.0%)	0.86	6/3807 (0.2%)
2	I	0.65	0/159	0.82	0/212
All	All	0.66	1/2986 (0.0%)	0.86	6/4019 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	199	CYS	CB-SG	-8.22	1.68	1.82

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	59	LEU	N-CA-C	-5.76	95.45	111.00
1	E	277	LEU	CA-CB-CG	5.17	127.19	115.30
1	E	165	ARG	N-CA-C	5.15	124.91	111.00
1	E	114	SER	N-CA-C	5.08	124.71	111.00
1	E	162	LEU	CA-CB-CG	5.07	126.96	115.30
1	E	196	TRP	N-CA-C	5.02	124.56	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2778	0	2757	86	0
2	I	157	0	146	7	0
3	E	27	0	20	4	0
4	E	18	0	0	3	0
4	I	3	0	0	0	0
All	All	2983	0	2923	94	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 16.

All (94) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:184:ASP:HB2	4:E:413:HOH:O	1.74	0.87
1:E:242:GLN:HG3	1:E:244:ILE:HG12	1.63	0.81
3:E:351:IQB:H3'1	3:E:351:IQB:O2	1.80	0.80
1:E:242:GLN:HG2	1:E:245:GLN:HG3	1.67	0.76
1:E:243:PRO:HA	1:E:246:ILE:HD12	1.70	0.73
1:E:54:PHE:HD2	1:E:82:LEU:HD12	1.54	0.73
1:E:75:ASP:O	1:E:79:VAL:HG23	1.89	0.72
1:E:295:LYS:H	1:E:295:LYS:HD3	1.54	0.71
1:E:104:VAL:HG21	1:E:183:THR:HG22	1.77	0.67
3:E:351:IQB:H3'2	4:E:413:HOH:O	1.94	0.66
1:E:304:ALA:HA	1:E:309:LYS:HE2	1.78	0.65
1:E:319:LYS:HD2	1:E:319:LYS:N	2.11	0.65
1:E:319:LYS:HD2	1:E:319:LYS:H	1.62	0.64
1:E:242:GLN:HG2	1:E:245:GLN:OE1	2.01	0.60
1:E:266:LYS:O	1:E:270:ARG:HG3	2.02	0.60
1:E:345:LYS:H	1:E:345:LYS:HD2	1.66	0.60
1:E:125:GLY:HA3	1:E:174:ILE:O	2.02	0.60
1:E:303:ILE:HG22	1:E:307:GLN:NE2	2.18	0.59
1:E:285:LYS:N	1:E:285:LYS:HD2	2.18	0.58
1:E:78:LYS:HA	1:E:81:LYS:HE3	1.84	0.57
1:E:30:TRP:O	1:E:33:PRO:HD3	2.05	0.57
1:E:277:LEU:O	1:E:283:ASN:ND2	2.38	0.56
1:E:140:GLU:HB2	1:E:141:PRO:HD3	1.88	0.56
1:E:100:PHE:CD1	1:E:101:PRO:HD2	2.41	0.55
1:E:56:ARG:NH2	1:E:331:GLU:O	2.39	0.55
1:E:90:ASN:O	1:E:94:ILE:HG13	2.06	0.55
1:E:189:LYS:HD3	1:E:191:VAL:HG21	1.89	0.54
1:E:209:ILE:HD11	1:E:219:VAL:HG11	1.89	0.54
1:E:304:ALA:CA	1:E:309:LYS:HE2	2.38	0.53
1:E:304:ALA:CB	1:E:309:LYS:HE2	2.38	0.53
1:E:158:HIS:HE1	1:E:220:ASP:OD2	1.90	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:242:GLN:HG2	1:E:245:GLN:CG	2.38	0.53
1:E:69:TYR:HB3	1:E:120:MET:O	2.09	0.53
1:E:45:ARG:HH12	1:E:334:GLU:HA	1.74	0.52
2:I:22:ILE:O	2:I:22:ILE:HG13	2.08	0.52
1:E:301:ASP:OD2	1:E:304:ALA:HB2	2.10	0.51
1:E:295:LYS:H	1:E:295:LYS:CD	2.20	0.51
1:E:294:HIS:CE1	1:E:295:LYS:HE3	2.45	0.51
1:E:333:GLU:HG2	1:E:336:ARG:HH21	1.75	0.51
1:E:73:ILE:HD13	1:E:335:ILE:HD11	1.92	0.50
1:E:244:ILE:HG13	1:E:245:GLN:N	2.26	0.50
1:E:65:THR:HB	1:E:67:ASN:ND2	2.26	0.50
1:E:33:PRO:HA	1:E:96:GLN:OE1	2.12	0.49
1:E:121:GLU:O	3:E:351:IQB:H6	2.12	0.48
1:E:202:PRO:HD3	2:I:22:ILE:HG23	1.94	0.48
3:E:351:IQB:C3'	3:E:351:IQB:O2	2.57	0.48
1:E:74:LEU:O	1:E:115:ASN:HB3	2.14	0.48
2:I:23:HIS:CD2	2:I:24:ASP:H	2.32	0.48
1:E:43:PHE:HB3	1:E:60:VAL:HG13	1.97	0.47
1:E:242:GLN:CG	1:E:245:GLN:HG3	2.42	0.47
1:E:284:LEU:HB3	1:E:285:LYS:HD2	1.97	0.47
1:E:295:LYS:N	1:E:295:LYS:HD3	2.24	0.47
1:E:288:VAL:O	1:E:288:VAL:HG22	2.14	0.46
1:E:285:LYS:H	1:E:285:LYS:HD2	1.79	0.46
1:E:189:LYS:HD3	1:E:191:VAL:CG2	2.45	0.46
1:E:15:VAL:O	1:E:18:PHE:N	2.48	0.46
1:E:62:HIS:NE2	1:E:64:GLU:HG2	2.31	0.46
1:E:91:GLU:O	1:E:95:LEU:HB2	2.16	0.46
1:E:54:PHE:CD2	1:E:82:LEU:HD12	2.42	0.45
1:E:338:SEP:O3P	1:E:339:ILE:HG13	2.16	0.45
1:E:343:CYS:HB3	1:E:346:GLU:OE1	2.16	0.45
1:E:158:HIS:CG	1:E:217:LYS:HD2	2.52	0.45
1:E:103:LEU:HD21	1:E:153:THR:CG2	2.47	0.45
1:E:127:GLU:HA	1:E:173:LEU:HD23	1.99	0.44
1:E:159:SER:OG	1:E:160:LEU:HD13	2.18	0.44
1:E:39:HIS:HB2	1:E:42:GLN:NE2	2.33	0.44
1:E:304:ALA:HB2	1:E:309:LYS:HE2	1.99	0.44
2:I:23:HIS:CD2	2:I:24:ASP:N	2.86	0.43
1:E:148:ALA:HA	1:E:302:TRP:CZ3	2.53	0.43
1:E:222:TRP:O	1:E:226:VAL:HG23	2.19	0.43
2:I:11:ILE:HD12	2:I:11:ILE:HG23	1.76	0.43
1:E:233:ALA:HB2	1:E:258:PRO:HG2	2.00	0.43
1:E:310:VAL:CG1	1:E:311:GLU:N	2.82	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:163:ILE:HG22	1:E:165:ARG:HG3	2.01	0.42
1:E:310:VAL:HG12	1:E:311:GLU:N	2.35	0.42
1:E:155:GLU:HG3	1:E:288:VAL:HG11	2.01	0.42
1:E:317:LYS:HA	1:E:317:LYS:HD2	1.72	0.42
1:E:305:ILE:HG12	1:E:310:VAL:HG21	2.01	0.42
1:E:198:LEU:HD23	1:E:209:ILE:HG22	2.02	0.42
1:E:285:LYS:HG2	1:E:286:ASP:OD2	2.20	0.42
1:E:305:ILE:CG1	1:E:310:VAL:HG21	2.50	0.42
1:E:137:ARG:NH1	1:E:232:ALA:O	2.52	0.42
1:E:148:ALA:HA	1:E:302:TRP:HZ3	1.85	0.41
2:I:20:ASN:HD22	2:I:20:ASN:N	2.18	0.41
1:E:345:LYS:HB3	1:E:345:LYS:HE3	1.79	0.41
1:E:289:ASN:O	1:E:293:ASN:HB2	2.20	0.41
1:E:285:LYS:HZ3	1:E:285:LYS:H	1.68	0.41
1:E:133:ARG:NH1	4:E:400:HOH:O	2.38	0.41
1:E:100:PHE:CG	1:E:101:PRO:HD2	2.55	0.41
1:E:17:GLU:O	1:E:20:ALA:HB3	2.21	0.41
1:E:345:LYS:CD	1:E:345:LYS:H	2.30	0.40
1:E:276:ASP:OD1	1:E:279:LYS:HG3	2.20	0.40
1:E:169:PRO:HD2	2:I:19:ARG:NH2	2.37	0.40
1:E:27:LEU:HA	1:E:27:LEU:HD12	1.91	0.40

There are no symmetry-related clashes.

## 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	332/350 (95%)	310 (93%)	19 (6%)	3 (1%)	25	26
2	I	18/20 (90%)	16 (89%)	2 (11%)	0	100	100
All	All	350/370 (95%)	326 (93%)	21 (6%)	3 (1%)	25	26

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	41	ASP
1	E	184	ASP
1	E	214	GLY

### 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	294/302 (97%)	260 (88%)	34 (12%)	8	8
2	I	15/15 (100%)	13 (87%)	2 (13%)	6	6
All	All	309/317 (98%)	273 (88%)	36 (12%)	8	8

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	16	LYS
1	E	17	GLU
1	E	27	LEU
1	E	28	LYS
1	E	36	ASN
1	E	37	THR
1	E	40	LEU
1	E	53	SER
1	E	59	LEU
1	E	60	VAL
1	E	92	LYS
1	E	95	LEU
1	E	104	VAL
1	E	160	LEU
1	E	162	LEU
1	E	174	ILE
1	E	198	LEU
1	E	202	PRO
1	E	208	GLU
1	E	210	ILE
1	E	263	SER
1	E	267	ASP
1	E	268	LEU

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Mol	Chain	Res	Type
1	E	269	LEU
1	E	272	LEU
1	E	277	LEU
1	E	278	THR
1	E	285	LYS
1	E	295	LYS
1	E	311	GLU
1	E	318	PHE
1	E	333	GLU
1	E	336	ARG
1	E	345	LYS
2	I	15	ARG
2	I	20	ASN

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

Mol	Chain	Res	Type
1	E	42	GLN
1	E	62	HIS
1	E	67	ASN
1	E	77	GLN
1	E	113	ASN
1	E	158	HIS
1	E	307	GLN
2	I	20	ASN
2	I	23	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	TPO	E	197	1	10,10,11	6.48	5 (50%)	12,14,16	1.37	2 (16%)
1	SEP	E	338	1	9,9,10	5.66	2 (22%)	10,12,14	1.96	3 (30%)

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
1	TPO	E	197	1	-	0/9/11/13	0/0/0/0
1	SEP	E	338	1	-	0/6/8/10	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	197	TPO	O-C	19.64	1.24	1.11
1	E	338	SEP	O-C	16.64	1.22	1.11
1	E	197	TPO	CA-C	2.90	1.53	1.48
1	E	197	TPO	P-O2P	-2.47	1.45	1.54
1	E	197	TPO	OG1-CB	-2.46	1.40	1.45
1	E	197	TPO	P-O1P	-2.44	1.43	1.51
1	E	338	SEP	CA-C	2.04	1.52	1.48

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	338	SEP	C-CA-N	3.53	117.35	113.83
1	E	338	SEP	OG-CB-CA	-3.16	104.22	108.69
1	E	197	TPO	O3P-P-O2P	2.27	116.44	107.61
1	E	338	SEP	O2P-P-O1P	2.07	117.22	110.44
1	E	197	TPO	OG1-P-O1P	-2.04	101.09	106.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

1 ligand is 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	IQB	E	351	-	29,29,29	1.22	3 (10%)	39,39,39	1.40	3 (7%)

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	IQB	E	351	-	-	0/16/17/17	0/1/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	351	IQB	C9-C8	3.05	1.40	1.36
3	E	351	IQB	C2-C1	2.74	1.40	1.37
3	E	351	IQB	C6-N7	2.15	1.36	1.32

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	351	IQB	C1-S-N1'	-4.74	100.60	106.85
3	E	351	IQB	C9-C8-N7	-3.64	120.04	123.84
3	E	351	IQB	C1-C10-C5	2.59	119.97	117.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	E	336/350 (96%)	0.09	10 (2%) 48 58	10, 24, 52, 69	0
2	I	20/20 (100%)	0.06	2 (10%) 8 13	12, 21, 65, 77	0
All	All	356/370 (96%)	0.09	12 (3%) 43 53	10, 24, 53, 77	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	23	HIS	4.4
2	I	24	ASP	4.1
1	E	318	PHE	3.7
1	E	311	GLU	3.2
1	E	253	GLY	2.9
1	E	39	HIS	2.8
1	E	336	ARG	2.7
1	E	334	GLU	2.6
1	E	339	ILE	2.4
1	E	242	GLN	2.3
1	E	319	LYS	2.1
1	E	41	ASP	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	TPO	E	197	11/12	0.10	-0.69	14,19,21,21	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	SEP	E	338	10/11	0.20	-0.70	49,52,55,55	0

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	IQB	E	351	27/27	0.12	-0.33	20,25,39,49	0

### 6.5 Other polymers

There are no such residues in this entry.