



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:28 PM UTC

PDB ID : 2Q7O / pdb_00002q7o
Title : Structure of human purine nucleoside phosphorylase in complex with L-Immucillin-H
Authors : Rinaldo-Matthis, A.; Almo, S.C.; Schramm, V.L.
Deposited on : 2007-06-07
Resolution : 2.90 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
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

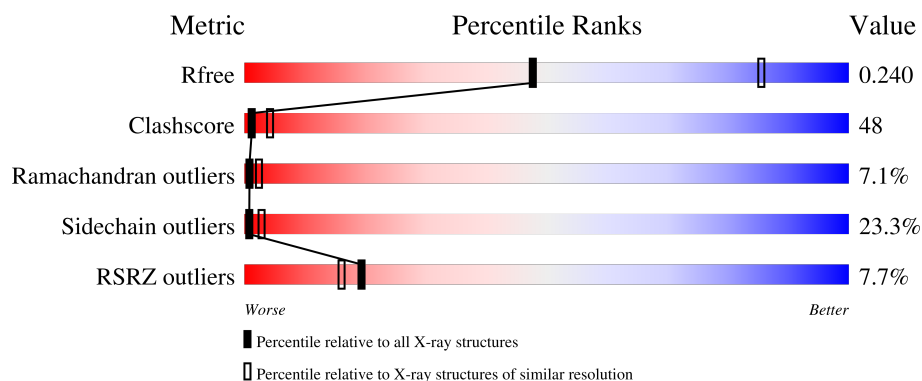
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	289	

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	E	290	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IMH	E	291	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2268 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.

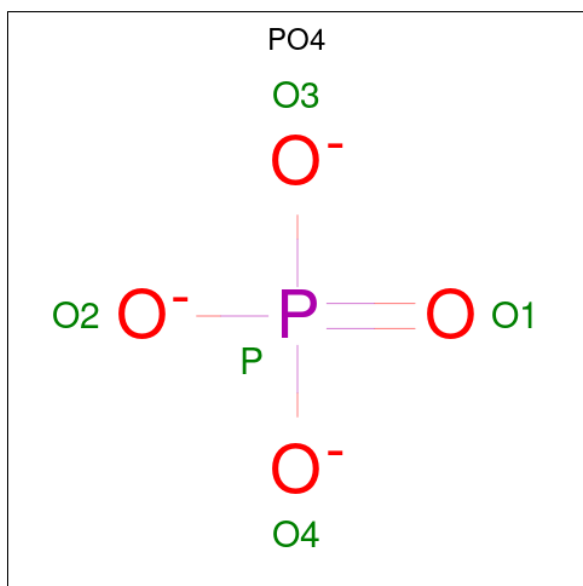
- Molecule 1 is a protein called Purine nucleoside phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	284	2223	1413	389	405	16	0	0	0

There is a discrepancy between the modelled and reference sequences:

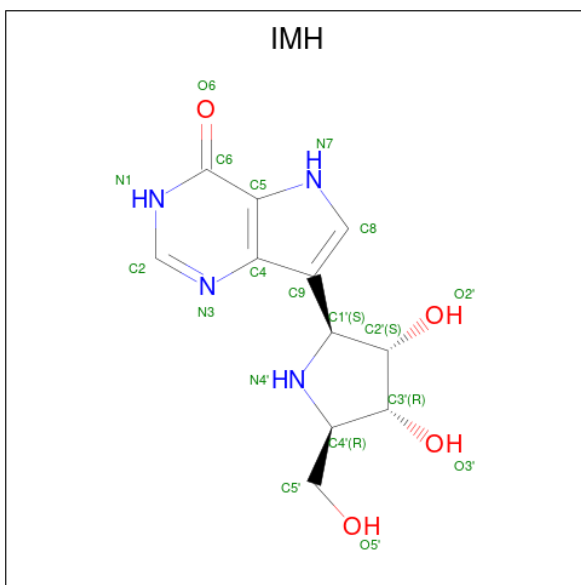
Chain	Residue	Modelled	Actual	Comment	Reference
E	51	SER	GLY	conflict	UNP P00491

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

- Molecule 3 is 1,4-DIDEOXY-4-AZA-1-(S)-(9-DEAZAHYPOXANTHIN-9-YL)-D-RIBITOL (CCD ID: IMH) (formula: $C_{11}H_{14}N_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	N	O	0	0
			19	11	4	4		

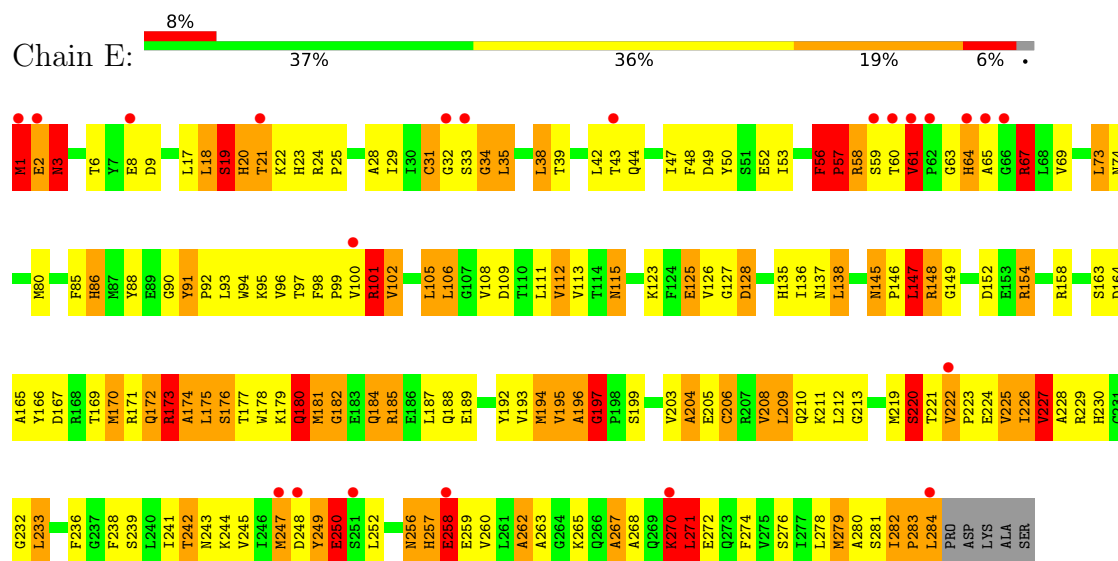
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	21	Total	O	0	0
			21	21		

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: Purine nucleoside phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	143.46Å 143.46Å 166.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.54 – 2.90 34.54 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (34.54-2.90) 99.9 (34.54-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.90Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.216 , 0.262 (Not available) , 0.240	Depositor DCC
R_{free} test set	745 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	77.0	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for -2/3*h-1/3*k+2/3*l,-1/3*h-2/3*k-2/3*l,2/3*h-2/3*k+1/3*l 0.017 for -h,1/3*h-1/3*k+2/3*l,2/3*h+4/3*k+1/3*l 0.012 for -1/3*h+1/3*k-2/3*l,-k,-4/3*h-2/3*k+1/3*l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2268	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: PO4, IMH

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	1.39	18/2274 (0.8%)	1.52	34/3076 (1.1%)

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	E	0	14

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	91	TYR	CA-C	10.82	1.65	1.52
1	E	147	LEU	CA-CB	-9.70	1.38	1.53
1	E	225	VAL	CA-CB	-8.13	1.45	1.54
1	E	96	VAL	C-O	7.93	1.33	1.24
1	E	96	VAL	CA-CB	-7.32	1.45	1.54
1	E	91	TYR	N-CA	6.99	1.57	1.46
1	E	227	VAL	N-CA	6.87	1.55	1.46
1	E	56	PHE	CA-C	-6.42	1.46	1.53
1	E	226	ILE	C-O	6.12	1.31	1.24
1	E	226	ILE	N-CA	-5.98	1.39	1.46
1	E	25	PRO	CA-C	-5.94	1.47	1.53
1	E	28	ALA	CA-CB	-5.88	1.44	1.53
1	E	92	PRO	CA-CB	-5.47	1.45	1.53
1	E	102	VAL	N-CA	5.39	1.52	1.46
1	E	95	LYS	C-O	-5.27	1.17	1.24
1	E	226	ILE	CB-CG1	5.11	1.63	1.53
1	E	61	VAL	CA-CB	5.03	1.60	1.54
1	E	228	ALA	N-CA	5.01	1.52	1.46

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	282	ILE	CA-C-N	10.22	132.61	119.84
1	E	282	ILE	C-N-CA	10.22	132.61	119.84
1	E	226	ILE	CA-C-N	-8.33	108.44	121.02
1	E	226	ILE	C-N-CA	-8.33	108.44	121.02
1	E	225	VAL	CB-CA-C	-8.27	101.48	111.81
1	E	101	ARG	CA-C-N	-8.17	108.82	120.42
1	E	101	ARG	C-N-CA	-8.17	108.82	120.42
1	E	225	VAL	O-C-N	-7.77	113.94	121.94
1	E	226	ILE	CA-C-O	7.76	129.12	121.05
1	E	96	VAL	CA-C-O	7.17	128.50	121.05
1	E	138	LEU	N-CA-C	6.70	122.76	113.45
1	E	91	TYR	CA-C-N	-6.69	113.78	120.21
1	E	91	TYR	C-N-CA	-6.69	113.78	120.21
1	E	226	ILE	CB-CA-C	-6.36	103.71	112.04
1	E	67	ARG	N-CA-C	6.10	119.42	109.24
1	E	203	VAL	N-CA-C	-6.08	104.37	111.00
1	E	147	LEU	N-CA-CB	6.03	119.37	110.33
1	E	96	VAL	CB-CA-C	-6.02	104.15	112.04
1	E	249	TYR	N-CA-C	-5.99	101.18	110.10
1	E	3	ASN	N-CA-CB	5.83	120.35	110.49
1	E	270	LYS	CA-C-N	5.82	132.17	121.70
1	E	270	LYS	C-N-CA	5.82	132.17	121.70
1	E	102	VAL	N-CA-CB	5.76	119.22	110.58
1	E	95	LYS	CA-C-O	-5.62	114.53	120.55
1	E	173	ARG	N-CA-C	5.61	122.74	110.80
1	E	248	ASP	N-CA-C	5.39	117.97	108.75
1	E	270	LYS	N-CA-C	5.38	122.26	110.80
1	E	173	ARG	CA-C-N	5.22	131.52	121.54
1	E	173	ARG	C-N-CA	5.22	131.52	121.54
1	E	204	ALA	N-CA-C	-5.21	105.69	111.36
1	E	145	ASN	CB-CA-C	5.05	117.60	110.02
1	E	206	CYS	CA-CB-SG	-5.04	102.81	114.40
1	E	135	HIS	N-CA-C	5.02	117.01	109.23
1	E	73	LEU	N-CA-C	-5.01	100.28	108.76

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	1	MET	Peptide
1	E	173	ARG	Peptide

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Mol	Chain	Res	Type	Group
1	E	19	SER	Peptide
1	E	194	MET	Peptide
1	E	196	ALA	Peptide
1	E	197	GLY	Peptide
1	E	220	SER	Peptide
1	E	256	ASN	Peptide
1	E	257	HIS	Peptide
1	E	267	ALA	Peptide
1	E	270	LYS	Peptide
1	E	32	GLY	Peptide
1	E	34	GLY	Peptide
1	E	64	HIS	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	E	2223	0	2201	214	0
2	E	5	0	0	6	0
3	E	19	0	13	0	0
4	E	21	0	0	10	0
All	All	2268	0	2214	214	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 48.

All (214) 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:E:220:SER:HB3	2:E:290:PO4:O3	1.38	1.23
1:E:220:SER:HB3	2:E:290:PO4:P	1.91	1.10
1:E:256:ASN:HD22	1:E:258:GLU:CB	1.64	1.09
1:E:23:HIS:O	1:E:24:ARG:HD3	1.50	1.09
1:E:194:MET:HB2	1:E:222:VAL:HG21	1.32	1.08
1:E:256:ASN:ND2	1:E:258:GLU:HB2	1.72	1.03
1:E:256:ASN:HD22	1:E:258:GLU:HB2	0.86	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:58:ARG:HH11	1:E:58:ARG:HG3	1.25	0.99
1:E:35:LEU:HD11	1:E:38:LEU:HD22	1.44	0.95
1:E:31:CYS:SG	1:E:35:LEU:CD2	2.56	0.94
1:E:100:VAL:HG21	1:E:224:GLU:HB2	1.49	0.93
1:E:265:LYS:O	1:E:268:ALA:HB2	1.70	0.92
1:E:35:LEU:HD21	1:E:38:LEU:HD23	1.51	0.91
1:E:31:CYS:SG	1:E:35:LEU:HD23	2.12	0.90
1:E:283:PRO:O	1:E:284:LEU:HD22	1.76	0.86
1:E:172:GLN:HA	1:E:175:LEU:HD23	1.57	0.84
1:E:59:SER:HA	1:E:91:TYR:OH	1.78	0.83
1:E:23:HIS:HB3	4:E:307:HOH:O	1.79	0.83
1:E:35:LEU:O	1:E:35:LEU:HG	1.80	0.82
1:E:19:SER:O	1:E:20:HIS:ND1	2.12	0.82
1:E:222:VAL:HB	1:E:223:PRO:HD3	1.64	0.79
1:E:180:GLN:CG	1:E:180:GLN:O	2.30	0.78
1:E:35:LEU:HD11	1:E:38:LEU:CD2	2.14	0.78
1:E:256:ASN:ND2	1:E:258:GLU:CB	2.39	0.77
1:E:194:MET:HB2	1:E:222:VAL:CG2	2.13	0.77
1:E:256:ASN:HB2	1:E:258:GLU:HG2	1.68	0.76
1:E:179:LYS:C	1:E:181:MET:H	1.95	0.74
1:E:19:SER:C	1:E:20:HIS:ND1	2.45	0.74
1:E:1:MET:SD	1:E:1:MET:C	2.71	0.73
1:E:194:MET:SD	1:E:223:PRO:HD3	2.30	0.72
1:E:98:PHE:HB3	1:E:99:PRO:HD3	1.71	0.71
1:E:35:LEU:HA	4:E:308:HOH:O	1.90	0.71
1:E:170:MET:HA	1:E:170:MET:CE	2.20	0.71
1:E:249:TYR:O	1:E:250:GLU:HB2	1.88	0.71
1:E:35:LEU:CD1	1:E:38:LEU:HD22	2.21	0.70
1:E:170:MET:O	1:E:174:ALA:HB2	1.92	0.70
1:E:179:LYS:C	1:E:181:MET:N	2.49	0.70
1:E:137:ASN:ND2	1:E:226:ILE:HD11	2.07	0.69
1:E:17:LEU:C	1:E:19:SER:H	2.00	0.69
1:E:220:SER:CB	2:E:290:PO4:O3	2.31	0.68
1:E:257:HIS:H	1:E:258:GLU:HG2	1.60	0.67
1:E:100:VAL:CG2	1:E:224:GLU:HB2	2.23	0.67
1:E:50:TYR:HA	1:E:53:ILE:HD12	1.76	0.67
1:E:145:ASN:HB3	4:E:304:HOH:O	1.93	0.66
1:E:258:GLU:CD	1:E:258:GLU:H	2.03	0.66
1:E:100:VAL:HG21	1:E:224:GLU:CB	2.25	0.66
1:E:220:SER:CB	2:E:290:PO4:P	2.78	0.66
1:E:242:THR:HG22	1:E:243:ASN:OD1	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:179:LYS:O	1:E:181:MET:N	2.29	0.65
1:E:222:VAL:HB	1:E:223:PRO:CD	2.27	0.65
1:E:47:ILE:HD13	1:E:67:ARG:HH12	1.59	0.65
1:E:93:LEU:O	1:E:97:THR:HG23	1.96	0.65
1:E:179:LYS:O	1:E:182:GLY:N	2.29	0.64
1:E:58:ARG:HG3	1:E:58:ARG:NH1	1.99	0.64
1:E:232:GLY:O	1:E:233:LEU:C	2.37	0.64
1:E:63:GLY:HA3	1:E:65:ALA:H	1.62	0.64
1:E:63:GLY:HA2	1:E:64:HIS:HB2	1.79	0.64
1:E:59:SER:CA	1:E:91:TYR:OH	2.45	0.63
1:E:180:GLN:O	1:E:180:GLN:HG3	1.99	0.63
1:E:137:ASN:HD22	1:E:226:ILE:HD11	1.64	0.62
1:E:258:GLU:CD	1:E:258:GLU:N	2.57	0.62
1:E:170:MET:O	1:E:174:ALA:CB	2.48	0.62
1:E:20:HIS:ND1	1:E:20:HIS:N	2.45	0.61
1:E:180:GLN:O	1:E:180:GLN:HG2	1.99	0.61
1:E:47:ILE:HD13	1:E:67:ARG:NH1	2.15	0.61
1:E:194:MET:CB	1:E:222:VAL:HG21	2.21	0.59
1:E:210:GLN:HE22	1:E:247:MET:HE3	1.68	0.59
1:E:109:ASP:HA	1:E:233:LEU:HD22	1.84	0.58
1:E:206:CYS:SG	1:E:245:VAL:CG1	2.92	0.58
1:E:88:TYR:HB3	1:E:197:GLY:HA2	1.84	0.58
1:E:258:GLU:CA	1:E:258:GLU:OE1	2.50	0.58
1:E:172:GLN:HA	1:E:175:LEU:CD2	2.29	0.58
1:E:233:LEU:HD23	1:E:233:LEU:N	2.18	0.58
1:E:170:MET:HA	1:E:170:MET:HE3	1.85	0.58
1:E:249:TYR:O	1:E:250:GLU:CB	2.52	0.58
1:E:35:LEU:O	1:E:35:LEU:CG	2.50	0.58
1:E:31:CYS:HB3	1:E:35:LEU:CD2	2.34	0.58
1:E:112:VAL:HA	1:E:236:PHE:O	2.03	0.58
1:E:220:SER:CB	2:E:290:PO4:O4	2.52	0.57
1:E:220:SER:O	1:E:222:VAL:N	2.37	0.57
1:E:97:THR:HG21	1:E:146:PRO:HB3	1.87	0.57
1:E:241:ILE:HG22	1:E:242:THR:N	2.19	0.56
1:E:17:LEU:C	1:E:19:SER:N	2.63	0.56
1:E:86:HIS:N	1:E:86:HIS:CD2	2.74	0.56
1:E:194:MET:SD	1:E:222:VAL:HB	2.45	0.56
1:E:272:GLU:O	1:E:276:SER:HB2	2.06	0.56
1:E:24:ARG:N	4:E:307:HOH:O	2.39	0.55
1:E:195:VAL:O	1:E:219:MET:HA	2.06	0.55
1:E:31:CYS:SG	1:E:35:LEU:HD22	2.46	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:179:LYS:C	1:E:182:GLY:H	2.15	0.55
1:E:206:CYS:SG	1:E:245:VAL:HG12	2.47	0.55
1:E:63:GLY:CA	1:E:65:ALA:H	2.19	0.54
1:E:33:SER:O	1:E:35:LEU:N	2.40	0.54
1:E:31:CYS:CB	1:E:35:LEU:HD23	2.39	0.53
1:E:97:THR:O	1:E:100:VAL:HB	2.08	0.53
1:E:21:THR:O	1:E:21:THR:CG2	2.57	0.53
1:E:98:PHE:N	1:E:99:PRO:HD2	2.24	0.53
1:E:171:ARG:O	1:E:175:LEU:HD22	2.07	0.53
1:E:172:GLN:C	1:E:174:ALA:HB3	2.33	0.53
1:E:61:VAL:HG13	4:E:300:HOH:O	2.08	0.53
1:E:61:VAL:O	1:E:63:GLY:HA2	2.08	0.53
1:E:222:VAL:CB	1:E:223:PRO:HD3	2.38	0.53
1:E:2:GLU:OE2	1:E:2:GLU:HA	2.09	0.52
1:E:271:LEU:H	1:E:274:PHE:H	1.58	0.52
1:E:60:THR:HG23	4:E:300:HOH:O	2.08	0.52
1:E:109:ASP:OD1	1:E:109:ASP:C	2.52	0.52
1:E:171:ARG:O	1:E:174:ALA:HB3	2.09	0.52
1:E:65:ALA:HB1	1:E:85:PHE:HE2	1.75	0.52
1:E:206:CYS:SG	1:E:245:VAL:HG11	2.50	0.52
1:E:222:VAL:O	1:E:223:PRO:C	2.53	0.52
1:E:18:LEU:O	1:E:19:SER:HB3	2.10	0.52
1:E:31:CYS:CB	1:E:35:LEU:CD2	2.88	0.52
1:E:53:ILE:O	1:E:58:ARG:NH2	2.42	0.52
1:E:258:GLU:OE1	1:E:258:GLU:HA	2.10	0.52
1:E:35:LEU:HD21	1:E:38:LEU:CD2	2.32	0.51
1:E:221:THR:O	1:E:224:GLU:N	2.42	0.51
1:E:39:THR:HA	1:E:42:LEU:HD12	1.93	0.51
1:E:241:ILE:CG2	1:E:242:THR:N	2.74	0.50
1:E:88:TYR:CB	1:E:197:GLY:HA2	2.42	0.50
1:E:258:GLU:O	1:E:259:GLU:C	2.54	0.50
1:E:57:PRO:HB2	1:E:85:PHE:CZ	2.46	0.50
1:E:148:ARG:HG3	1:E:149:GLY:N	2.26	0.50
1:E:256:ASN:HB2	1:E:258:GLU:CG	2.39	0.50
1:E:270:LYS:CA	1:E:271:LEU:HB2	2.41	0.50
1:E:21:THR:O	1:E:24:ARG:NH1	2.44	0.50
1:E:205:GLU:O	1:E:208:VAL:HG13	2.11	0.49
1:E:220:SER:HB2	2:E:290:PO4:O4	2.12	0.49
1:E:192:TYR:CE2	1:E:222:VAL:HG23	2.47	0.49
1:E:152:ASP:OD1	1:E:154:ARG:N	2.44	0.49
1:E:170:MET:HA	1:E:170:MET:HE2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:43:THR:O	1:E:44:GLN:HB2	2.12	0.49
1:E:164:ASP:CG	1:E:164:ASP:O	2.53	0.49
1:E:88:TYR:HB3	1:E:197:GLY:CA	2.43	0.49
1:E:221:THR:O	1:E:222:VAL:C	2.56	0.49
1:E:47:ILE:HG22	1:E:48:PHE:N	2.28	0.49
1:E:171:ARG:C	1:E:174:ALA:HB3	2.38	0.49
1:E:173:ARG:N	1:E:174:ALA:HB3	2.28	0.49
1:E:187:LEU:HD12	1:E:187:LEU:HA	1.40	0.48
1:E:98:PHE:HB3	1:E:99:PRO:CD	2.43	0.47
1:E:137:ASN:ND2	1:E:226:ILE:CD1	2.78	0.47
1:E:196:ALA:O	1:E:197:GLY:O	2.32	0.47
1:E:225:VAL:O	1:E:226:ILE:C	2.57	0.47
1:E:242:THR:CG2	1:E:243:ASN:OD1	2.63	0.47
1:E:99:PRO:O	1:E:102:VAL:HG22	2.15	0.47
1:E:204:ALA:O	1:E:208:VAL:HG12	2.14	0.47
1:E:184:GLN:HG2	1:E:185:ARG:N	2.29	0.47
1:E:220:SER:O	1:E:221:THR:HB	2.15	0.47
1:E:210:GLN:OE1	1:E:247:MET:HG3	2.16	0.46
1:E:105:LEU:HD23	1:E:105:LEU:HA	1.72	0.46
1:E:31:CYS:HB3	1:E:35:LEU:HD22	1.98	0.46
1:E:136:ILE:HB	1:E:193:VAL:HG23	1.98	0.46
1:E:222:VAL:CB	1:E:223:PRO:CD	2.90	0.46
1:E:271:LEU:HD23	1:E:271:LEU:HA	1.89	0.46
1:E:170:MET:CE	1:E:170:MET:CA	2.94	0.46
1:E:49:ASP:O	1:E:52:GLU:HG2	2.16	0.46
1:E:125:GLU:OE2	1:E:244:LYS:HE3	2.15	0.46
1:E:211:LYS:C	1:E:213:GLY:H	2.24	0.46
1:E:169:THR:O	1:E:173:ARG:HB2	2.16	0.45
1:E:21:THR:O	1:E:21:THR:HG22	2.16	0.45
1:E:63:GLY:HA2	1:E:64:HIS:CB	2.46	0.45
1:E:265:LYS:C	1:E:267:ALA:H	2.24	0.45
1:E:279:MET:HA	1:E:282:ILE:HD12	1.99	0.45
1:E:74:ASN:HD21	1:E:280:ALA:HB2	1.80	0.45
1:E:173:ARG:HD3	1:E:281:SER:HB2	1.99	0.45
1:E:61:VAL:CG1	4:E:300:HOH:O	2.65	0.45
1:E:98:PHE:N	1:E:99:PRO:CD	2.80	0.45
1:E:33:SER:C	1:E:35:LEU:H	2.25	0.44
1:E:73:LEU:HD23	1:E:279:MET:HG2	1.98	0.44
1:E:6:THR:N	1:E:9:ASP:OD2	2.42	0.44
1:E:98:PHE:O	1:E:99:PRO:C	2.59	0.44
1:E:1:MET:SD	1:E:1:MET:O	2.76	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:258:GLU:N	1:E:258:GLU:OE1	2.51	0.44
1:E:88:TYR:HB3	1:E:197:GLY:H	1.83	0.43
1:E:3:ASN:HA	1:E:94:TRP:CD2	2.54	0.43
1:E:229:ARG:O	1:E:230:HIS:C	2.57	0.43
1:E:211:LYS:C	1:E:213:GLY:N	2.76	0.43
1:E:258:GLU:O	1:E:260:VAL:N	2.51	0.43
1:E:115:ASN:ND2	1:E:239:SER:OG	2.52	0.43
1:E:137:ASN:C	1:E:137:ASN:OD1	2.59	0.43
1:E:111:LEU:HD11	1:E:113:VAL:HG12	2.01	0.43
1:E:238:PHE:CD1	1:E:238:PHE:C	2.97	0.43
1:E:94:TRP:CD1	1:E:94:TRP:H	2.37	0.43
1:E:47:ILE:CG2	1:E:48:PHE:N	2.81	0.42
1:E:192:TYR:HE2	1:E:222:VAL:HG23	1.82	0.42
1:E:19:SER:O	1:E:19:SER:OG	2.35	0.42
1:E:65:ALA:HB1	1:E:85:PHE:CE2	2.54	0.42
1:E:171:ARG:O	1:E:174:ALA:CB	2.66	0.42
1:E:268:ALA:O	1:E:272:GLU:HG3	2.20	0.42
1:E:29:ILE:HD12	1:E:80:MET:HE2	2.01	0.42
1:E:50:TYR:CD2	1:E:56:PHE:CE2	3.08	0.42
1:E:74:ASN:ND2	1:E:280:ALA:HB2	2.35	0.42
1:E:127:GLY:O	1:E:128:ASP:O	2.37	0.42
1:E:88:TYR:C	1:E:90:GLY:H	2.28	0.42
1:E:165:ALA:HB3	1:E:166:TYR:CD2	2.54	0.42
1:E:167:ASP:HB3	1:E:170:MET:HB2	2.01	0.42
1:E:152:ASP:OD1	1:E:152:ASP:C	2.62	0.42
1:E:245:VAL:HG13	4:E:301:HOH:O	2.19	0.42
1:E:61:VAL:O	1:E:64:HIS:HB2	2.20	0.41
1:E:176:SER:O	1:E:178:TRP:N	2.52	0.41
1:E:270:LYS:HB3	1:E:271:LEU:HB2	2.01	0.41
1:E:31:CYS:HB3	1:E:35:LEU:HD23	1.99	0.41
1:E:101:ARG:HD2	4:E:311:HOH:O	2.21	0.41
1:E:106:LEU:HA	1:E:106:LEU:HD12	1.68	0.41
1:E:172:GLN:O	1:E:175:LEU:N	2.53	0.41
1:E:193:VAL:HG11	1:E:209:LEU:HG	2.02	0.41
1:E:42:LEU:HD11	1:E:80:MET:HE3	2.02	0.41
1:E:111:LEU:HD21	1:E:113:VAL:CG1	2.50	0.41
1:E:123:LYS:HE3	1:E:188:GLN:HE22	1.85	0.41
1:E:147:LEU:HD12	1:E:147:LEU:HA	1.73	0.41
1:E:166:TYR:O	1:E:167:ASP:C	2.63	0.41
1:E:39:THR:HG22	1:E:69:VAL:HG21	2.01	0.41
1:E:262:ALA:O	1:E:263:ALA:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:18:LEU:HD13	1:E:106:LEU:HD12	2.02	0.41
1:E:227:VAL:HG23	4:E:311:HOH:O	2.20	0.41

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	E	282/289 (98%)	225 (80%)	37 (13%)	20 (7%)	1 2

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	19	SER
1	E	173	ARG
1	E	174	ALA
1	E	176	SER
1	E	250	GLU
1	E	271	LEU
1	E	283	PRO
1	E	34	GLY
1	E	128	ASP
1	E	177	THR
1	E	180	GLN
1	E	197	GLY
1	E	258	GLU
1	E	18	LEU
1	E	262	ALA
1	E	270	LYS
1	E	222	VAL
1	E	3	ASN
1	E	182	GLY

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Mol	Chain	Res	Type
1	E	57	PRO

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	236/240 (98%)	181 (77%)	55 (23%)	1 3

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	1	MET
1	E	2	GLU
1	E	8	GLU
1	E	19	SER
1	E	20	HIS
1	E	21	THR
1	E	22	LYS
1	E	31	CYS
1	E	35	LEU
1	E	38	LEU
1	E	56	PHE
1	E	57	PRO
1	E	58	ARG
1	E	61	VAL
1	E	67	ARG
1	E	86	HIS
1	E	101	ARG
1	E	105	LEU
1	E	106	LEU
1	E	108	VAL
1	E	112	VAL
1	E	115	ASN
1	E	125	GLU
1	E	126	VAL
1	E	138	LEU

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Mol	Chain	Res	Type
1	E	147	LEU
1	E	148	ARG
1	E	154	ARG
1	E	158	ARG
1	E	163	SER
1	E	170	MET
1	E	172	GLN
1	E	175	LEU
1	E	180	GLN
1	E	181	MET
1	E	184	GLN
1	E	185	ARG
1	E	189	GLU
1	E	195	VAL
1	E	199	SER
1	E	208	VAL
1	E	209	LEU
1	E	212	LEU
1	E	220	SER
1	E	227	VAL
1	E	233	LEU
1	E	242	THR
1	E	247	MET
1	E	250	GLU
1	E	252	LEU
1	E	258	GLU
1	E	271	LEU
1	E	278	LEU
1	E	279	MET
1	E	284	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	23	HIS
1	E	74	ASN
1	E	104	HIS
1	E	115	ASN
1	E	172	GLN
1	E	256	ASN
1	E	269	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$
3	IMH	E	291	-	21,21,21	2.79	5 (23%)	22,31,31	2.90	12 (54%)
2	PO4	E	290	-	4,4,4	1.88	1 (25%)	6,6,6	1.45	1 (16%)

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	IMH	E	291	-	4/4/4/5	3/6/22/22	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	291	IMH	C5-C4	7.06	1.45	1.38
3	E	291	IMH	O6-C6	6.94	1.36	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	291	IMH	C3'-C4'	-5.37	1.48	1.53
3	E	291	IMH	C5-N7	-3.64	1.32	1.38
2	E	290	PO4	P-O1	3.48	1.58	1.50
3	E	291	IMH	C2-N3	2.50	1.35	1.30

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	291	IMH	C4-C5-C6	-6.80	115.43	121.41
3	E	291	IMH	C5'-C4'-C3'	-6.46	102.54	113.68
3	E	291	IMH	O3'-C3'-C4'	-4.01	102.19	112.92
3	E	291	IMH	O5'-C5'-C4'	-3.43	102.79	111.10
3	E	291	IMH	O6-C6-C5	-3.41	119.04	127.26
3	E	291	IMH	C5-C6-N1	3.32	117.54	110.26
2	E	290	PO4	O4-P-O3	3.09	117.54	107.91
3	E	291	IMH	C2'-C3'-C4'	2.86	106.89	102.55
3	E	291	IMH	C2'-C1'-N4'	-2.78	100.94	104.81
3	E	291	IMH	C5'-C4'-N4'	2.49	115.85	111.47
3	E	291	IMH	C6-C5-N7	2.41	135.75	131.14
3	E	291	IMH	C4-C5-N7	2.23	108.81	105.57
3	E	291	IMH	C2-N1-C6	-2.12	122.18	125.09

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	291	IMH	C3'
3	E	291	IMH	C1'
3	E	291	IMH	C2'
3	E	291	IMH	C4'

All (3) torsion outliers are listed below:

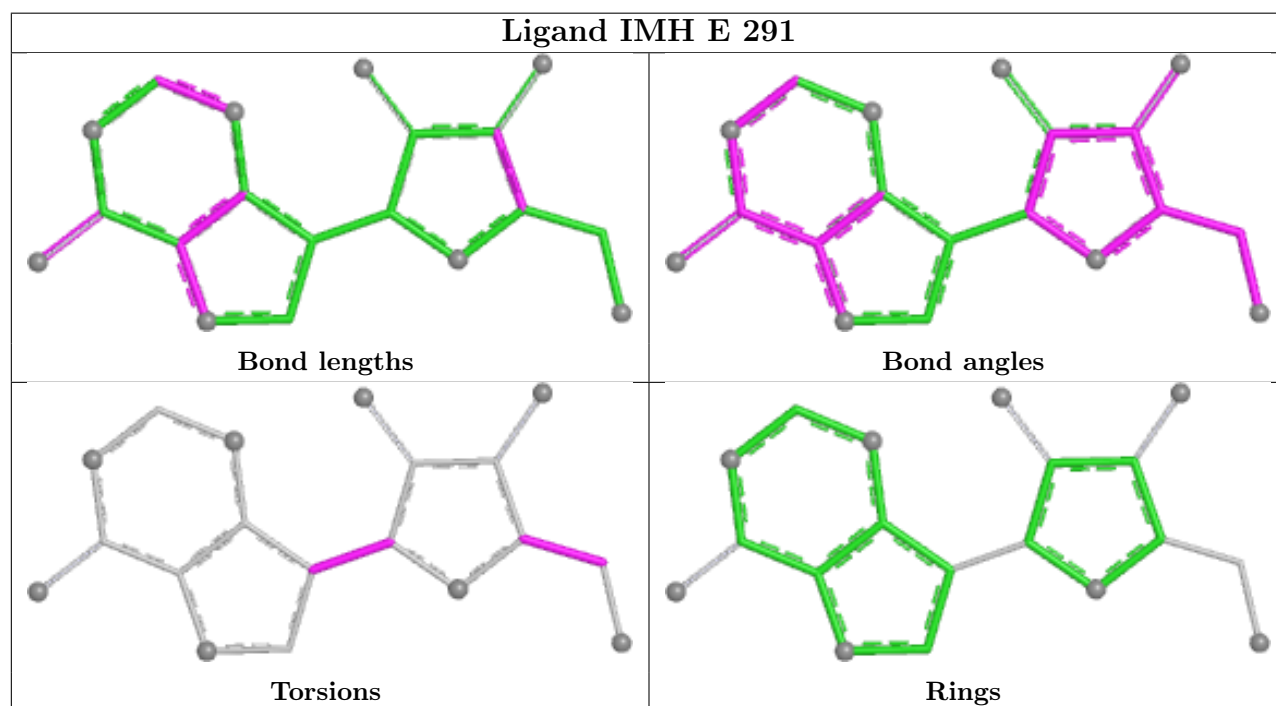
Mol	Chain	Res	Type	Atoms
3	E	291	IMH	N4'-C4'-C5'-O5'
3	E	291	IMH	C3'-C4'-C5'-O5'
3	E	291	IMH	C2'-C1'-C9-C8

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	290	PO4	6	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	284/289 (98%)	0.54	22 (7%) 19 16	44, 68, 99, 109	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	65	ALA	6.1
1	E	62	PRO	4.8
1	E	61	VAL	4.0
1	E	222	VAL	3.7
1	E	59	SER	3.6
1	E	247	MET	3.6
1	E	258	GLU	3.3
1	E	32	GLY	3.0
1	E	284	LEU	2.9
1	E	66	GLY	2.8
1	E	64	HIS	2.8
1	E	60	THR	2.8
1	E	33	SER	2.4
1	E	248	ASP	2.3
1	E	21	THR	2.3
1	E	43	THR	2.3
1	E	2	GLU	2.3
1	E	270	LYS	2.2
1	E	100	VAL	2.2
1	E	1	MET	2.1
1	E	8	GLU	2.0
1	E	251	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](#)

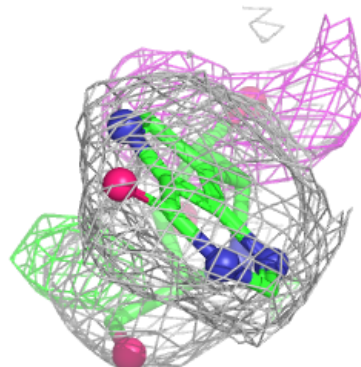
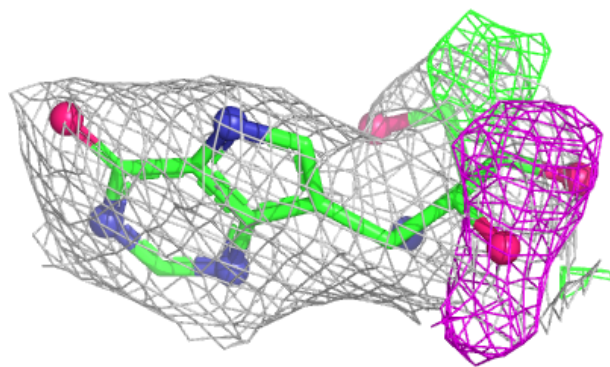
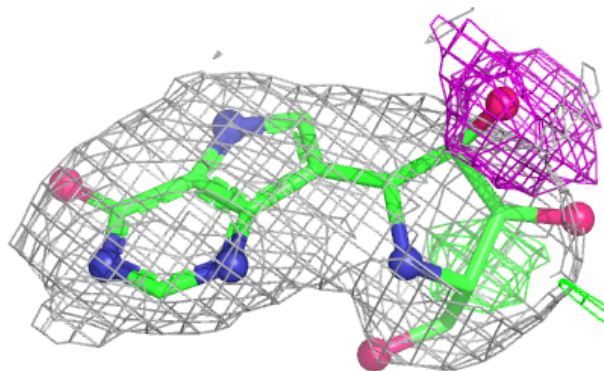
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	E	290	5/5	0.83	0.21	133,135,136,137	0
3	IMH	E	291	19/19	0.91	0.14	68,71,89,90	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.

Electron density around IMH E 291:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.