



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:39 AM UTC

PDB ID : 2ZDX / pdb_00002zdx
Title : Inhibitor-bound structures of human pyruvate dehydrogenase kinase 4
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Deposited on : 2007-11-30
Resolution : 2.54 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

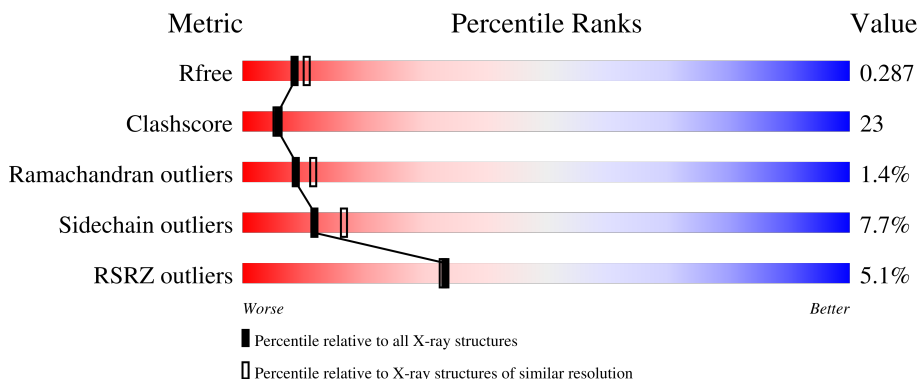
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	394	4% 54% 28% • 14%
1	B	394	5% 45% 36% 6% 13%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5623 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.

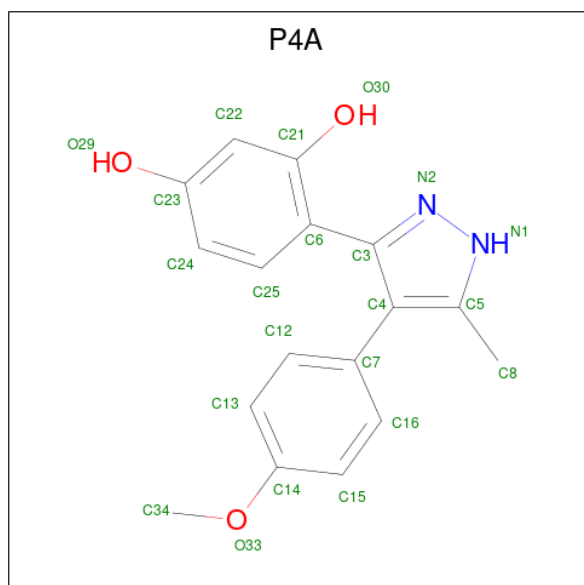
- Molecule 1 is a protein called Pyruvate dehydrogenase kinase isozyme 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2720	1743	460	504	13			
1	B	344	Total	C	N	O	S	0	1	0
			2780	1779	470	516	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	GLY	-	expression tag	UNP Q16654
A	19	PRO	-	expression tag	UNP Q16654
B	18	GLY	-	expression tag	UNP Q16654
B	19	PRO	-	expression tag	UNP Q16654

- Molecule 2 is 4-[4-(4-methoxyphenyl)-5-methyl-1H-pyrazol-3-yl]benzene-1,3-diol (CCD ID: P4A) (formula: C₁₇H₁₆N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			22	17	2	3		
2	B	1	Total	C	N	O	0	0
			22	17	2	3		

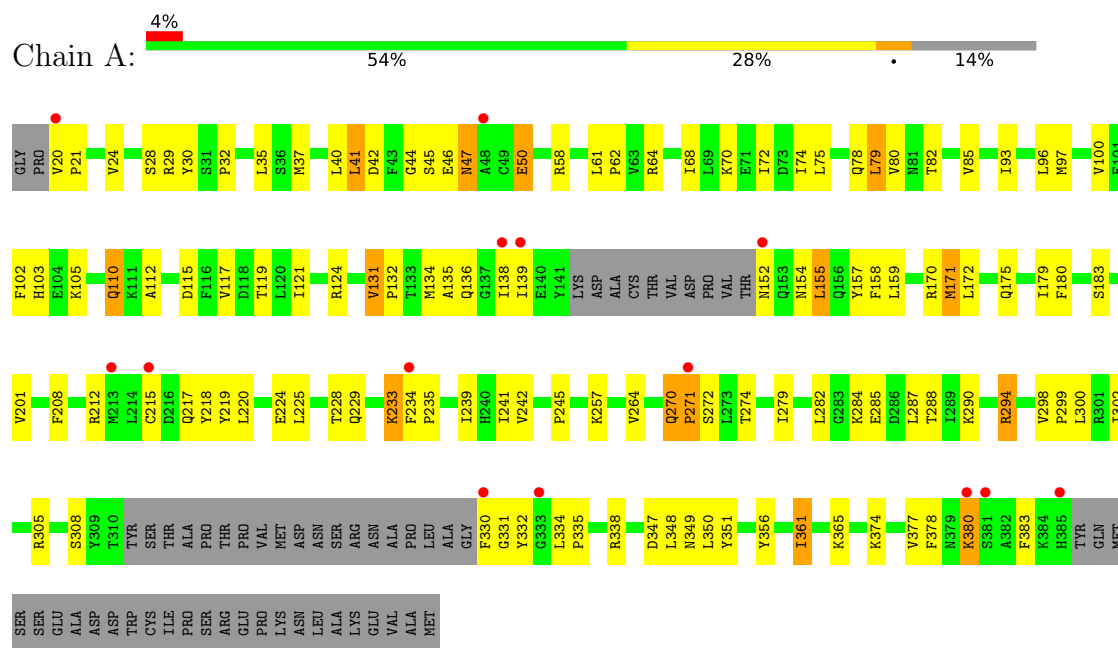
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	43	Total	O	0	0
			43	43		
3	B	36	Total	O	0	0
			36	36		

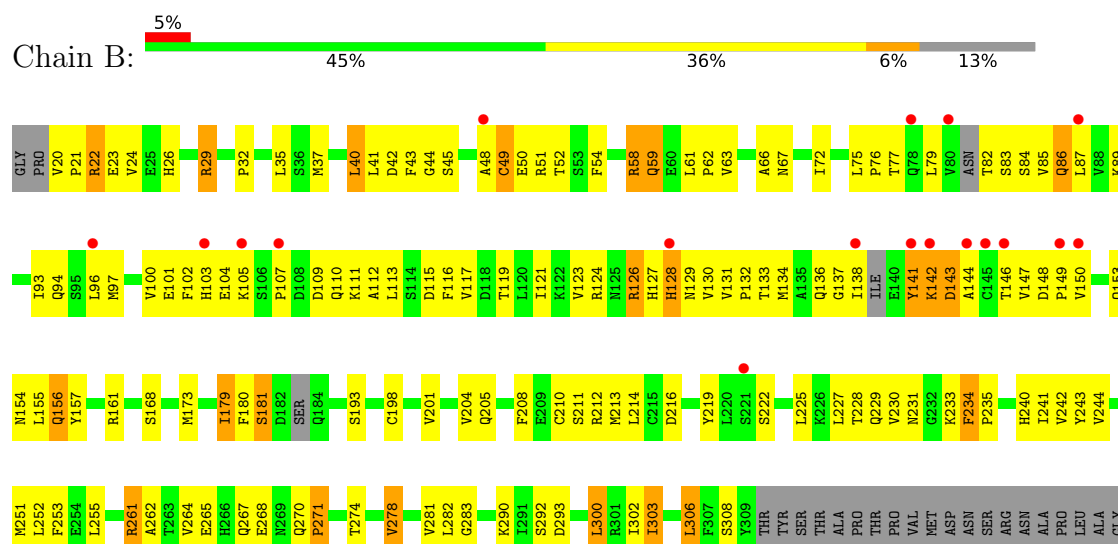
3 Residue-property plots

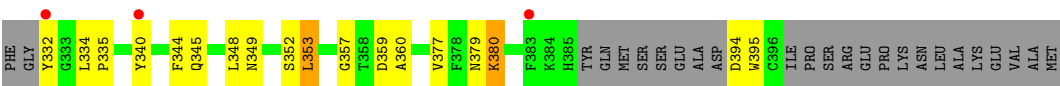
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Pyruvate dehydrogenase kinase isozyme 4



• Molecule 1: Pyruvate dehydrogenase kinase isozyme 4





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.08Å 69.08Å 85.63Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	48.68 – 2.54 48.68 – 2.54	Depositor EDS
% Data completeness (in resolution range)	87.0 (48.68-2.54) 87.2 (48.68-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.99 (at 2.54Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.222 , 0.290 0.220 , 0.287	Depositor DCC
R_{free} test set	1144 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5623	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.60% 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: P4A

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	0.44	0/2782	0.78	3/3768 (0.1%)
1	B	0.42	0/2847	0.79	2/3855 (0.1%)
All	All	0.43	0/5629	0.79	5/7623 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	PHE	CA-C-N	5.51	126.73	119.84
1	A	234	PHE	C-N-CA	5.51	126.73	119.84
1	A	50	GLU	N-CA-C	-5.38	106.67	113.18
1	B	234	PHE	CA-C-N	5.11	126.23	119.84
1	B	234	PHE	C-N-CA	5.11	126.23	119.84

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	2720	0	2706	99	0
1	B	2780	0	2754	156	0
2	A	22	0	14	0	0
2	B	22	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	43	0	0	2	0
3	B	36	0	0	0	0
All	All	5623	0	5488	253	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 23.

All (253) 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:270:GLN:H	1:A:270:GLN:NE2	1.58	1.00
1:B:32:PRO:HD2	1:B:377:VAL:HG22	1.48	0.93
1:A:288:THR:HG21	1:B:353:LEU:HD22	1.53	0.89
1:A:131:VAL:HG13	1:A:132:PRO:HD3	1.56	0.85
1:B:233:LYS:HB2	1:B:282:LEU:HD12	1.57	0.84
1:A:124:ARG:HH21	1:A:170:ARG:HE	1.25	0.84
1:B:156:GLN:HE22	1:B:308:SER:HA	1.45	0.80
1:B:270:GLN:HG3	1:B:271:PRO:HD2	1.64	0.79
1:A:24:VAL:HG12	1:A:96:LEU:HD13	1.64	0.78
1:A:270:GLN:H	1:A:270:GLN:HE21	1.29	0.78
1:B:332:TYR:C	1:B:335:PRO:HD2	2.09	0.78
1:A:134:MET:HE2	1:A:158:PHE:HE2	1.49	0.77
1:A:212:ARG:HH12	1:A:224:GLU:HG2	1.49	0.77
1:B:129:ASN:C	1:B:132:PRO:HD2	2.12	0.75
1:A:138:ILE:HD13	1:A:159:LEU:HD11	1.69	0.74
1:B:72:ILE:HA	1:B:75:LEU:HD23	1.70	0.73
1:B:142:LYS:HD2	1:B:147:VAL:HG21	1.68	0.73
1:B:117:VAL:HG11	1:B:181:SER:HB2	1.71	0.73
1:B:101:GLU:O	1:B:104:GLU:HG2	1.89	0.72
1:B:21:PRO:C	1:B:23:GLU:H	1.97	0.71
1:A:349:ASN:HB2	1:A:361:ILE:HD11	1.73	0.71
1:B:93:ILE:O	1:B:97:MET:HG2	1.90	0.71
1:B:233:LYS:HD2	1:B:282:LEU:HD11	1.72	0.71
1:A:349:ASN:HB2	1:A:361:ILE:CD1	2.21	0.70
1:B:96:LEU:O	1:B:100:VAL:HG23	1.91	0.70
1:A:124:ARG:HH21	1:A:170:ARG:NE	1.91	0.69
1:B:59:GLN:O	1:B:62:PRO:HD2	1.93	0.68
1:A:257:LYS:HZ3	1:A:330:PHE:HE1	1.41	0.68
1:A:115:ASP:O	1:A:119:THR:HG23	1.93	0.68
1:B:201:VAL:O	1:B:205:GLN:HG3	1.94	0.67
1:A:21:PRO:O	1:A:24:VAL:HG22	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:THR:HG21	1:A:294:ARG:HG3	1.75	0.67
1:B:212:ARG:HD3	1:B:222:SER:OG	1.94	0.66
1:A:241:ILE:HD13	1:A:287:LEU:HG	1.77	0.66
1:A:75:LEU:HD23	1:A:79:LEU:HD13	1.78	0.66
1:B:50:GLU:HB3	1:B:180:PHE:CD1	2.31	0.66
1:A:37:MET:HE1	1:A:242:VAL:O	1.96	0.65
1:B:141:TYR:HD2	1:B:142:LYS:H	1.42	0.65
1:B:59:GLN:HE22	1:B:63:VAL:CG2	2.09	0.65
1:A:45:SER:HB3	1:A:46:GLU:OE1	1.95	0.65
1:B:24:VAL:HG23	1:B:66:ALA:HA	1.78	0.65
1:B:89:LYS:O	1:B:93:ILE:HG13	1.98	0.64
1:A:105:LYS:HD3	1:A:112:ALA:HB2	1.80	0.64
1:A:138:ILE:HD11	1:A:332:TYR:OH	1.97	0.64
1:B:23:GLU:HG3	1:B:97:MET:SD	2.38	0.63
1:B:352:SER:C	1:B:353:LEU:HD23	2.23	0.63
1:B:303:ILE:O	1:B:306:LEU:HB2	1.98	0.63
1:B:241:ILE:CD1	1:B:243:TYR:HB2	2.28	0.63
1:B:282:LEU:HD13	1:B:283:GLY:N	2.14	0.62
1:A:102:PHE:CE1	1:A:119:THR:HG21	2.34	0.62
1:B:265:GLU:O	1:B:268:GLU:HG3	1.99	0.62
1:B:142:LYS:CD	1:B:147:VAL:HG21	2.29	0.61
1:A:45:SER:C	1:A:47:ASN:H	2.09	0.61
1:B:41:LEU:O	1:B:45:SER:HB3	2.00	0.61
1:B:251:MET:HE2	1:B:340:TYR:HB3	1.83	0.61
1:B:124:ARG:HH22	1:B:128:HIS:CE1	2.19	0.61
1:B:82:THR:O	1:B:86:GLN:HG2	2.02	0.60
1:A:61:LEU:HB2	1:A:62:PRO:HD3	1.82	0.60
1:B:208:PHE:CZ	1:B:225:LEU:HB2	2.37	0.60
1:B:241:ILE:HD11	1:B:243:TYR:HB2	1.83	0.60
1:A:270:GLN:NE2	1:A:270:GLN:N	2.42	0.59
1:B:37:MET:O	1:B:41:LEU:HB2	2.02	0.59
1:A:37:MET:O	1:A:41:LEU:HB2	2.03	0.58
1:A:124:ARG:NH2	1:A:170:ARG:HE	1.97	0.58
1:A:212:ARG:NH1	1:A:224:GLU:HG2	2.20	0.57
1:A:282:LEU:HD13	1:A:287:LEU:HD23	1.87	0.56
1:B:150:VAL:O	1:B:153:GLN:HG2	2.05	0.56
1:B:26:HIS:O	1:B:29:ARG:HB2	2.06	0.56
1:A:100:VAL:HA	1:A:103:HIS:CE1	2.40	0.55
1:A:175:GLN:O	1:A:179:ILE:HG13	2.06	0.55
1:B:59:GLN:HE22	1:B:63:VAL:HG23	1.71	0.55
1:B:86:GLN:HG3	1:B:87:LEU:N	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:VAL:HG13	1:A:132:PRO:CD	2.31	0.55
1:B:334:LEU:HB3	1:B:335:PRO:HD3	1.87	0.55
1:B:270:GLN:HG3	1:B:271:PRO:CD	2.33	0.54
1:B:49:CYS:SG	1:B:51:ARG:HB2	2.48	0.54
1:B:82:THR:HG21	1:B:137:GLY:HA2	1.89	0.54
1:B:255:LEU:HD13	1:B:360:ALA:HB1	1.89	0.54
1:B:214:LEU:HD13	1:B:253:PHE:HZ	1.73	0.54
1:B:117:VAL:O	1:B:121:ILE:HG13	2.07	0.53
1:B:233:LYS:HD2	1:B:282:LEU:CD1	2.37	0.53
1:A:45:SER:C	1:A:47:ASN:N	2.65	0.53
1:B:82:THR:HG22	1:B:84:SER:H	1.73	0.53
1:B:142:LYS:HB2	1:B:147:VAL:HG21	1.91	0.53
1:B:227:LEU:C	1:B:227:LEU:HD23	2.33	0.53
1:A:102:PHE:HE1	1:A:119:THR:HG21	1.74	0.53
1:A:46:GLU:CD	1:A:46:GLU:H	2.17	0.53
1:A:135:ALA:O	1:A:138:ILE:HG12	2.09	0.53
1:B:332:TYR:O	1:B:335:PRO:HD2	2.07	0.53
1:B:134:MET:O	1:B:138:ILE:HG12	2.10	0.52
1:B:61:LEU:HB2	1:B:62:PRO:HD3	1.91	0.52
1:B:157:TYR:O	1:B:161:ARG:HG3	2.08	0.52
1:A:219:TYR:CD1	1:A:264:VAL:HG11	2.44	0.52
1:A:270:GLN:HB3	1:A:271:PRO:CD	2.39	0.52
1:B:23:GLU:HG3	1:B:97:MET:HE1	1.90	0.52
1:A:139:ILE:HG22	1:A:139:ILE:O	2.09	0.52
1:B:216:ASP:OD2	1:B:222:SER:HB2	2.09	0.52
1:B:292:SER:HB3	1:B:359:ASP:OD1	2.10	0.52
1:B:117:VAL:CG1	1:B:181:SER:HB2	2.38	0.52
1:A:96:LEU:C	1:A:96:LEU:HD23	2.35	0.51
1:A:217:GLN:HB3	1:A:218:TYR:CD2	2.46	0.51
1:B:37:MET:HE1	1:B:243:TYR:CD1	2.45	0.51
1:B:50:GLU:CD	1:B:113:LEU:HD23	2.36	0.51
1:A:50:GLU:HB3	1:A:180:PHE:CE1	2.45	0.51
1:B:37:MET:HE1	1:B:243:TYR:HA	1.91	0.51
1:B:193:SER:O	1:B:244:VAL:HG13	2.11	0.51
1:B:22:ARG:O	1:B:22:ARG:HG3	2.10	0.51
1:B:129:ASN:O	1:B:132:PRO:HD2	2.10	0.51
1:B:105:LYS:HD3	1:B:112:ALA:HB2	1.93	0.50
1:B:124:ARG:HH12	1:B:128:HIS:CE1	2.29	0.50
1:A:68:ILE:O	1:A:72:ILE:HG13	2.11	0.50
1:A:138:ILE:HG13	1:A:139:ILE:N	2.26	0.50
1:B:48:ALA:O	1:B:49:CYS:C	2.54	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:ASP:HB2	1:B:149:PRO:CD	2.41	0.50
1:B:219:TYR:CD1	1:B:264:VAL:HG11	2.47	0.50
1:B:293:ASP:O	1:B:357:GLY:HA3	2.12	0.50
1:B:353:LEU:HD23	1:B:353:LEU:N	2.27	0.50
1:A:380:LYS:HA	1:A:383:PHE:HB2	1.94	0.50
1:B:59:GLN:C	1:B:59:GLN:HE21	2.20	0.50
1:B:261:ARG:O	1:B:265:GLU:HB2	2.12	0.50
1:A:233:LYS:O	1:A:235:PRO:HD3	2.12	0.49
1:B:261:ARG:HG3	1:B:262:ALA:N	2.24	0.49
1:A:100:VAL:HA	1:A:103:HIS:ND1	2.27	0.49
1:B:54:PHE:CZ	1:B:102:PHE:HB3	2.47	0.49
1:B:84:SER:HA	1:B:133:THR:CG2	2.43	0.49
1:A:302:ILE:HG13	1:A:305:ARG:NE	2.28	0.49
1:B:394:ASP:OD2	1:B:395:TRP:HD1	1.95	0.49
1:A:28:SER:C	1:A:30:TYR:H	2.19	0.49
1:A:229:GLN:HB2	1:A:239:ILE:HD11	1.93	0.49
1:A:208:PHE:CZ	1:A:225:LEU:HB2	2.48	0.49
1:A:351:TYR:CE1	1:A:361:ILE:HD11	2.47	0.49
1:B:241:ILE:HD12	1:B:242:VAL:C	2.37	0.49
1:B:300:LEU:O	1:B:300:LEU:HD22	2.13	0.48
1:A:302:ILE:HG12	1:A:305:ARG:HG2	1.95	0.48
1:B:127:HIS:HA	1:B:130:VAL:HG23	1.96	0.48
1:A:42:ASP:C	1:A:44:GLY:H	2.21	0.48
1:B:75:LEU:HD12	1:B:79:LEU:HD23	1.95	0.48
1:A:32:PRO:HD2	1:A:377:VAL:HG22	1.94	0.47
1:B:179:ILE:O	1:B:179:ILE:CG2	2.59	0.47
1:B:107:PRO:O	1:B:113:LEU:HD11	2.14	0.47
1:A:270:GLN:HE21	1:A:270:GLN:N	2.04	0.47
1:B:42:ASP:C	1:B:44:GLY:H	2.22	0.47
1:B:198:CYS:O	1:B:240:HIS:HA	2.14	0.47
1:B:208:PHE:CE2	1:B:225:LEU:HB2	2.50	0.47
1:B:20:VAL:HG11	1:B:24:VAL:HG11	1.97	0.47
1:B:59:GLN:C	1:B:62:PRO:HD2	2.40	0.47
1:B:97:MET:HA	1:B:100:VAL:HG23	1.95	0.47
1:B:113:LEU:N	1:B:113:LEU:HD12	2.30	0.47
1:B:116:PHE:HA	1:B:119:THR:OG1	2.14	0.47
1:B:131:VAL:N	1:B:132:PRO:CD	2.78	0.47
1:B:379:ASN:HB2	1:B:380:LYS:HD2	1.96	0.47
1:B:148:ASP:HB2	1:B:149:PRO:HD2	1.97	0.47
1:B:334:LEU:N	1:B:335:PRO:CD	2.78	0.47
1:B:282:LEU:HD13	1:B:282:LEU:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:GLN:NE2	1:B:63:VAL:HG23	2.30	0.46
1:A:270:GLN:HB3	1:A:271:PRO:HD2	1.97	0.46
1:A:294:ARG:HD2	1:A:356:TYR:CZ	2.50	0.46
1:A:80:VAL:HA	1:A:85:VAL:HG11	1.98	0.46
1:A:215:CYS:HG	1:A:219:TYR:HD2	1.61	0.46
1:A:24:VAL:HG12	1:A:96:LEU:CD1	2.41	0.46
1:A:331:GLY:HA3	1:A:332:TYR:HA	1.70	0.46
1:B:21:PRO:C	1:B:23:GLU:N	2.64	0.46
1:B:21:PRO:O	1:B:23:GLU:N	2.48	0.46
1:B:116:PHE:HA	1:B:119:THR:HG1	1.81	0.46
1:B:211:SER:HA	1:B:253:PHE:HE1	1.81	0.46
1:A:37:MET:HE3	1:A:242:VAL:HG12	1.98	0.46
1:B:20:VAL:CG1	1:B:24:VAL:HG11	2.45	0.46
1:A:58:ARG:O	1:A:103:HIS:CE1	2.69	0.45
1:B:37:MET:HA	1:B:40:LEU:HB2	1.98	0.45
1:A:157:TYR:OH	1:B:394:ASP:OD1	2.29	0.45
1:B:43:PHE:HE2	1:B:52:THR:HG22	1.80	0.45
1:A:334:LEU:HD12	1:A:334:LEU:HA	1.79	0.45
1:B:76:PRO:O	1:B:77:THR:C	2.59	0.45
1:A:284:LYS:HD2	1:A:285:GLU:HG3	1.98	0.45
1:B:82:THR:HG22	1:B:83:SER:N	2.32	0.45
1:B:133:THR:O	1:B:136:GLN:HB3	2.17	0.45
1:B:292:SER:CB	1:B:359:ASP:OD1	2.65	0.45
1:B:96:LEU:O	1:B:96:LEU:HD23	2.17	0.45
1:B:23:GLU:HG3	1:B:97:MET:CE	2.47	0.44
1:A:351:TYR:HE1	1:A:361:ILE:HD11	1.82	0.44
1:B:142:LYS:HD3	1:B:155:LEU:HD11	1.98	0.44
1:B:267:GLN:HB3	1:B:270:GLN:HB3	2.00	0.44
1:A:50:GLU:HB3	1:A:180:PHE:CD1	2.52	0.44
1:B:105:LYS:HD3	1:B:112:ALA:CB	2.47	0.44
1:B:231:ASN:O	1:B:235:PRO:HA	2.18	0.44
1:A:201:VAL:HG13	3:A:520:HOH:O	2.16	0.44
1:B:96:LEU:HD23	1:B:96:LEU:C	2.41	0.44
1:B:144:ALA:C	1:B:146:THR:H	2.25	0.44
1:A:380:LYS:O	1:A:383:PHE:N	2.50	0.44
1:B:230:VAL:O	1:B:281:VAL:HA	2.18	0.44
1:B:82:THR:O	1:B:86:GLN:HB3	2.18	0.44
1:B:124:ARG:O	1:B:127:HIS:N	2.50	0.43
1:A:271:PRO:HB2	1:A:272:SER:H	1.64	0.43
1:B:85:VAL:O	1:B:86:GLN:C	2.60	0.43
1:B:59:GLN:HE21	1:B:59:GLN:CA	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ILE:O	1:A:97:MET:HG2	2.18	0.43
1:A:134:MET:HE2	1:A:158:PHE:CE2	2.40	0.43
1:B:54:PHE:HD1	1:B:116:PHE:CD2	2.35	0.43
1:B:97:MET:HA	1:B:100:VAL:CG2	2.48	0.43
1:B:210:CYS:HA	1:B:213:MET:HE3	2.01	0.43
1:A:217:GLN:HE21	1:A:217:GLN:HB2	1.63	0.43
1:B:123:VAL:HG23	1:B:124:ARG:N	2.33	0.43
1:B:42:ASP:C	1:B:44:GLY:N	2.76	0.43
1:B:101:GLU:C	1:B:103:HIS:H	2.26	0.43
1:B:204:VAL:HA	1:B:252:LEU:HD13	2.01	0.43
1:B:228:THR:HG22	1:B:229:GLN:N	2.33	0.43
1:A:299:PRO:O	1:A:302:ILE:HG22	2.19	0.42
1:A:61:LEU:HD23	1:A:61:LEU:HA	1.86	0.42
1:A:334:LEU:HB3	1:A:335:PRO:HD3	2.02	0.42
1:A:171:MET:O	1:A:175:GLN:HB2	2.19	0.42
1:B:144:ALA:C	1:B:146:THR:N	2.77	0.42
1:A:74:ILE:HD12	1:A:378:PHE:HE1	1.84	0.42
1:A:155:LEU:HD22	1:A:155:LEU:HA	1.62	0.42
1:A:79:LEU:O	1:A:82:THR:HG23	2.20	0.42
1:A:228:THR:O	1:A:279:ILE:HA	2.19	0.42
1:B:58:ARG:O	1:B:103:HIS:HE1	2.02	0.42
1:A:132:PRO:O	1:A:136:GLN:HG2	2.20	0.42
1:B:79:LEU:HA	1:B:82:THR:OG1	2.20	0.42
1:B:348:LEU:HD23	1:B:349:ASN:N	2.35	0.42
1:A:28:SER:C	1:A:30:TYR:N	2.77	0.42
1:B:156:GLN:HE21	1:B:156:GLN:HB2	1.66	0.42
1:A:41:LEU:HD12	1:A:41:LEU:HA	1.84	0.42
1:B:267:GLN:HE22	1:B:274:THR:H	1.68	0.42
1:A:245:PRO:HD2	3:A:502:HOH:O	2.19	0.42
1:B:115:ASP:O	1:B:119:THR:HG23	2.20	0.42
1:A:42:ASP:C	1:A:44:GLY:N	2.78	0.42
1:A:305:ARG:O	1:A:308:SER:HB3	2.20	0.42
1:B:124:ARG:CD	1:B:173:MET:HE3	2.49	0.42
1:B:241:ILE:HD12	1:B:242:VAL:O	2.20	0.42
1:B:344:PHE:O	1:B:345:GLN:HB2	2.19	0.42
1:A:64:ARG:NH2	1:A:172:LEU:HD11	2.35	0.41
1:A:380:LYS:HA	1:A:380:LYS:HE3	2.02	0.41
1:B:82:THR:C	1:B:84:SER:H	2.28	0.41
1:B:302:ILE:O	1:B:303:ILE:C	2.63	0.41
1:A:46:GLU:OE1	1:A:46:GLU:N	2.52	0.41
1:B:63:VAL:O	1:B:67:ASN:ND2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:GLN:H	1:A:110:GLN:NE2	2.18	0.41
1:A:290:LYS:HG3	1:A:361:ILE:HG23	2.02	0.41
1:B:109:ASP:O	1:B:113:LEU:HD13	2.21	0.41
1:A:302:ILE:O	1:A:305:ARG:HG2	2.20	0.41
1:B:76:PRO:HB2	1:B:143:ASP:HA	2.01	0.41
1:B:94:GLN:HB3	1:B:126:ARG:NH1	2.36	0.41
1:B:241:ILE:HD12	1:B:243:TYR:HB2	2.00	0.41
1:A:117:VAL:O	1:A:121:ILE:HG13	2.20	0.41
1:A:134:MET:CE	1:A:158:PHE:HE2	2.23	0.41
1:B:113:LEU:N	1:B:113:LEU:CD1	2.83	0.41
1:B:84:SER:OG	1:B:137:GLY:N	2.53	0.41
1:B:20:VAL:HA	1:B:21:PRO:HD3	1.89	0.41
1:B:37:MET:HE2	1:B:344:PHE:CE2	2.56	0.40
1:B:101:GLU:C	1:B:103:HIS:N	2.80	0.40
1:B:109:ASP:OD1	1:B:111:LYS:HB3	2.22	0.40
1:B:278:VAL:HA	1:B:290:LYS:O	2.21	0.40
1:A:347:ASP:OD1	1:A:348:LEU:N	2.45	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	A	331/394 (84%)	313 (95%)	15 (4%)	3 (1%)	14	20
1	B	333/394 (84%)	301 (90%)	26 (8%)	6 (2%)	6	7
All	All	664/788 (84%)	614 (92%)	41 (6%)	9 (1%)	9	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	LYS

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Mol	Chain	Res	Type
1	A	271	PRO
1	A	29	ARG
1	B	22	ARG
1	B	142	LYS
1	B	271	PRO
1	B	49	CYS
1	B	143	ASP
1	B	303	ILE

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	310/358 (87%)	284 (92%)	26 (8%)	10	14
1	B	318/358 (89%)	296 (93%)	22 (7%)	14	20
All	All	628/716 (88%)	580 (92%)	48 (8%)	12	17

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	35	LEU
1	A	40	LEU
1	A	41	LEU
1	A	47	ASN
1	A	70	LYS
1	A	78	GLN
1	A	79	LEU
1	A	110	GLN
1	A	131	VAL
1	A	152	ASN
1	A	154	ASN
1	A	155	LEU
1	A	171	MET
1	A	183	SER

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Mol	Chain	Res	Type
1	A	220	LEU
1	A	270	GLN
1	A	294	ARG
1	A	298	VAL
1	A	300	LEU
1	A	338	ARG
1	A	350	LEU
1	A	361	ILE
1	A	365	LYS
1	A	374	LYS
1	A	380	LYS
1	B	29	ARG
1	B	35	LEU
1	B	40	LEU
1	B	58	ARG
1	B	59	GLN
1	B	86	GLN
1	B	110	GLN
1	B	126	ARG
1	B	128	HIS
1	B	141	TYR
1	B	154	ASN
1	B	156	GLN
1	B	168	SER
1	B	179	ILE
1	B	181	SER
1	B	234	PHE
1	B	261	ARG
1	B	278	VAL
1	B	300	LEU
1	B	306	LEU
1	B	353	LEU
1	B	380	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	HIS
1	A	47	ASN
1	A	59	GLN
1	A	78	GLN
1	A	81	ASN

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Mol	Chain	Res	Type
1	A	94	GLN
1	A	110	GLN
1	A	136	GLN
1	A	153	GLN
1	A	174	ASN
1	A	197	ASN
1	A	217	GLN
1	A	237	GLN
1	A	250	HIS
1	A	270	GLN
1	B	59	GLN
1	B	78	GLN
1	B	86	GLN
1	B	103	HIS
1	B	129	ASN
1	B	136	GLN
1	B	153	GLN
1	B	156	GLN
1	B	174	ASN
1	B	190	HIS
1	B	205	GLN
1	B	267	GLN
1	B	270	GLN
1	B	349	ASN

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	P4A	B	1500	-	24,24,24	3.42	19 (79%)	30,34,34	1.02	3 (10%)
2	P4A	A	500	-	24,24,24	3.48	18 (75%)	30,34,34	1.02	2 (6%)

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	P4A	B	1500	-	-	5/10/10/10	0/3/3/3
2	P4A	A	500	-	-	5/10/10/10	0/3/3/3

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	P4A	C6-C3	5.66	1.57	1.48
2	A	500	P4A	C16-C7	5.62	1.47	1.39
2	B	1500	P4A	C16-C7	5.60	1.47	1.39
2	A	500	P4A	C6-C21	5.09	1.48	1.40
2	B	1500	P4A	C6-C3	4.87	1.56	1.48
2	A	500	P4A	C25-C6	4.84	1.47	1.39
2	B	1500	P4A	C25-C6	4.64	1.47	1.39
2	B	1500	P4A	C12-C7	4.31	1.45	1.39
2	B	1500	P4A	C3-N2	4.25	1.40	1.34
2	B	1500	P4A	C7-C4	4.20	1.57	1.49
2	A	500	P4A	C13-C14	4.19	1.46	1.38
2	A	500	P4A	C12-C7	4.18	1.45	1.39
2	A	500	P4A	C7-C4	4.18	1.57	1.49
2	B	1500	P4A	C6-C21	4.17	1.47	1.40
2	B	1500	P4A	C15-C14	4.00	1.46	1.38
2	B	1500	P4A	C13-C14	3.96	1.46	1.38
2	A	500	P4A	C22-C23	3.94	1.45	1.39
2	B	1500	P4A	C22-C21	3.93	1.44	1.38
2	B	1500	P4A	C16-C15	3.91	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	P4A	C22-C21	3.85	1.44	1.38
2	A	500	P4A	C15-C14	3.84	1.45	1.38
2	A	500	P4A	C16-C15	3.78	1.44	1.38
2	B	1500	P4A	C22-C23	3.69	1.44	1.39
2	A	500	P4A	C3-N2	3.66	1.39	1.34
2	A	500	P4A	C13-C12	3.38	1.44	1.38
2	B	1500	P4A	C5-N1	3.14	1.40	1.34
2	B	1500	P4A	C13-C12	2.93	1.43	1.38
2	A	500	P4A	C25-C24	2.80	1.43	1.38
2	B	1500	P4A	C25-C24	2.72	1.43	1.38
2	A	500	P4A	C4-C5	2.64	1.44	1.39
2	B	1500	P4A	C4-C5	2.62	1.44	1.39
2	B	1500	P4A	C8-C5	2.60	1.54	1.49
2	A	500	P4A	C8-C5	2.59	1.54	1.49
2	A	500	P4A	C5-N1	2.56	1.39	1.34
2	B	1500	P4A	C24-C23	2.47	1.43	1.39
2	A	500	P4A	O33-C14	2.16	1.41	1.37
2	B	1500	P4A	O33-C14	2.16	1.41	1.37

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	P4A	C34-O33-C14	-2.96	111.14	117.50
2	B	1500	P4A	C34-O33-C14	-2.85	111.38	117.50
2	B	1500	P4A	C7-C4-C5	2.50	130.40	126.78
2	B	1500	P4A	C4-C5-N1	2.24	108.65	106.70
2	A	500	P4A	C4-C5-N1	2.22	108.64	106.70

There are no chirality outliers.

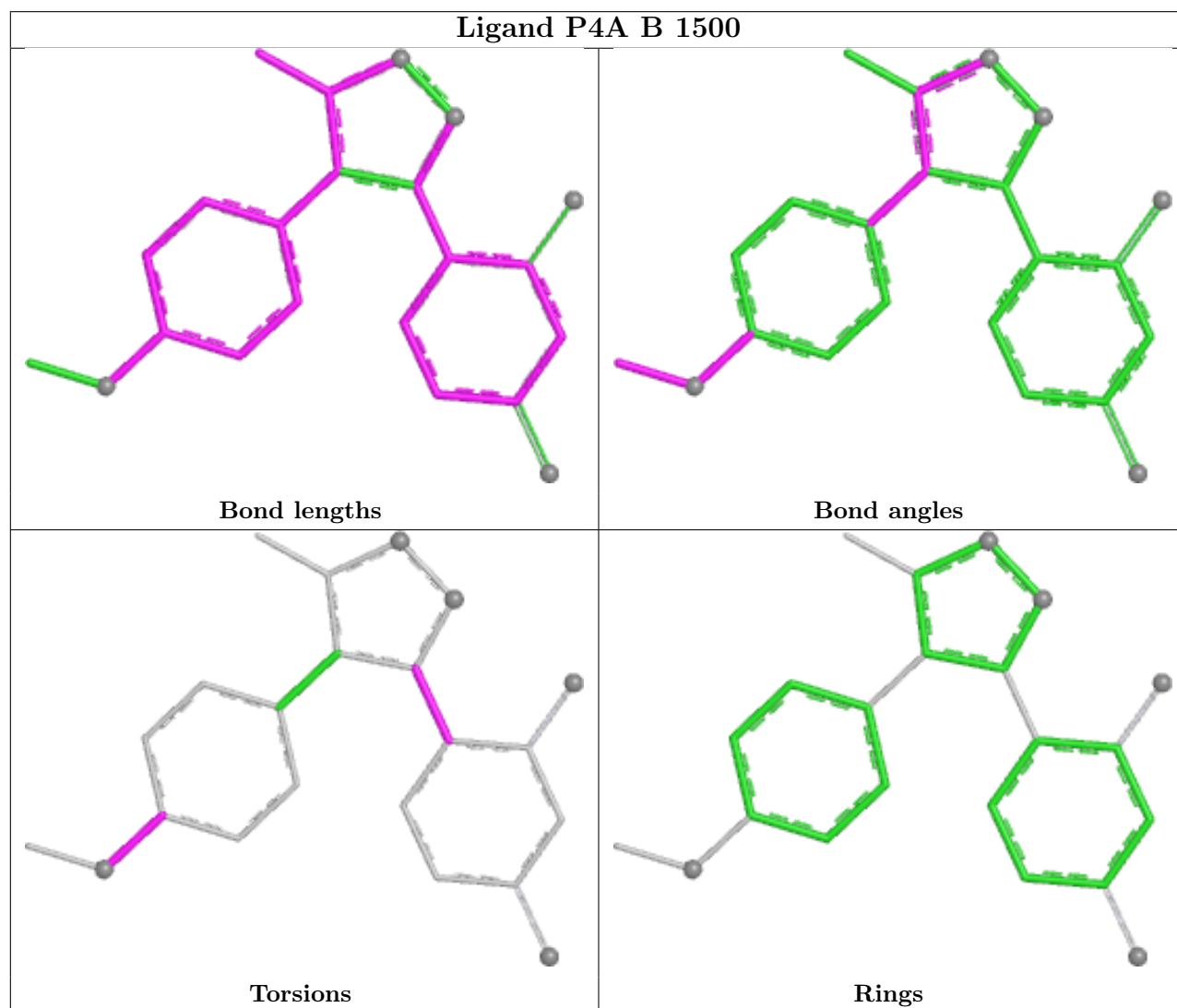
All (10) torsion outliers are listed below:

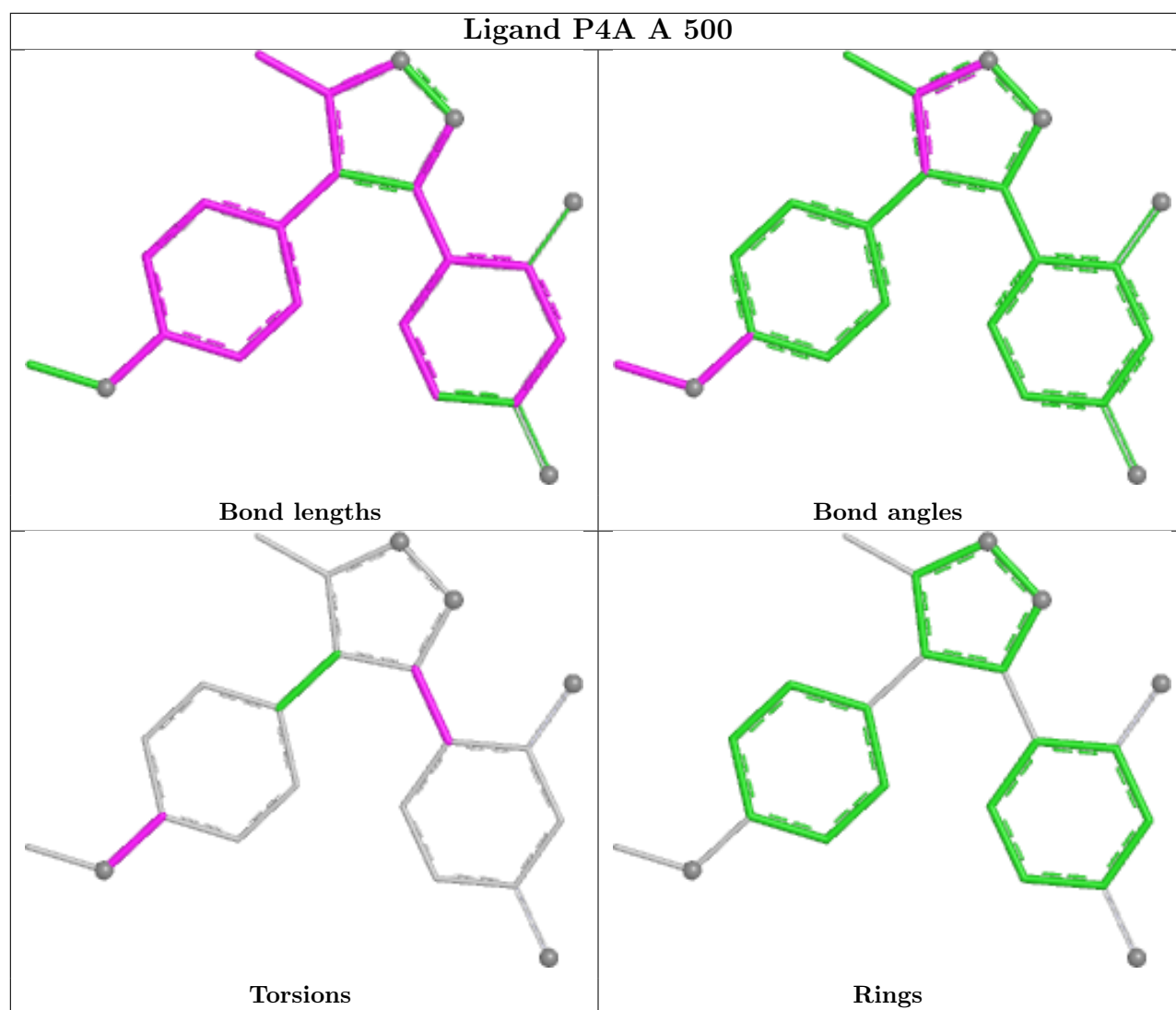
Mol	Chain	Res	Type	Atoms
2	A	500	P4A	C13-C14-O33-C34
2	B	1500	P4A	C15-C14-O33-C34
2	A	500	P4A	C15-C14-O33-C34
2	B	1500	P4A	C13-C14-O33-C34
2	A	500	P4A	N2-C3-C6-C21
2	B	1500	P4A	N2-C3-C6-C21
2	A	500	P4A	C4-C3-C6-C25
2	B	1500	P4A	C4-C3-C6-C25
2	A	500	P4A	C4-C3-C6-C21
2	B	1500	P4A	C4-C3-C6-C21

There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	337/394 (85%)	0.22	14 (4%)	40 41	25, 42, 71, 86	1 (0%)
1	B	344/394 (87%)	0.47	21 (6%)	27 26	22, 50, 81, 99	2 (0%)
All	All	681/788 (86%)	0.35	35 (5%)	33 33	22, 45, 78, 99	3 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	138	ILE	3.5
1	B	78	GLN	3.4
1	B	128	HIS	3.2
1	A	139	ILE	3.2
1	B	107	PRO	3.2
1	B	141	TYR	3.0
1	B	105	LYS	2.8
1	A	385	HIS	2.8
1	B	142	LYS	2.7
1	A	234	PHE	2.6
1	B	150	VAL	2.6
1	B	144	ALA	2.6
1	A	333	GLY	2.6
1	B	383	PHE	2.6
1	A	20	VAL	2.5
1	B	221	SER	2.5
1	B	48	ALA	2.5
1	B	145	CYS	2.5
1	B	103	HIS	2.5
1	A	215	CYS	2.5
1	B	146	THR	2.5
1	A	213	MET	2.4
1	B	87	LEU	2.4
1	A	330	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	80	VAL	2.4
1	A	152	ASN	2.4
1	A	381	SER	2.4
1	B	138	ILE	2.4
1	A	48	ALA	2.3
1	B	149	PRO	2.3
1	A	271	PRO	2.2
1	B	340	TYR	2.2
1	B	332	TYR	2.2
1	A	380	LYS	2.1
1	B	96	LEU	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](#)

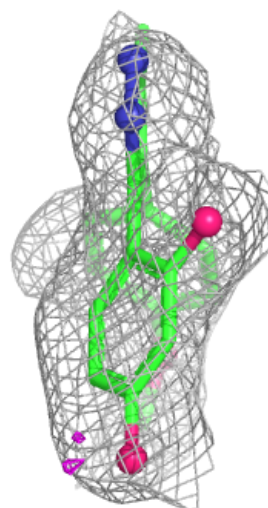
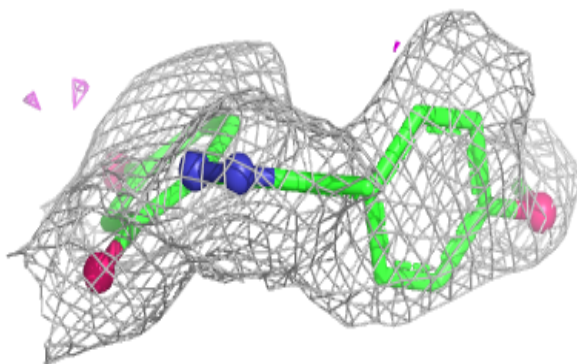
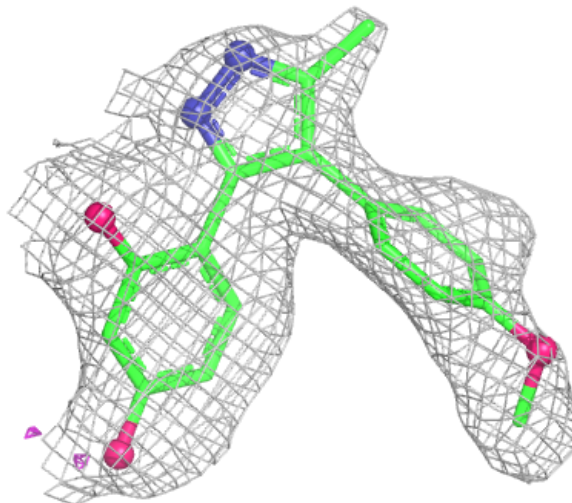
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

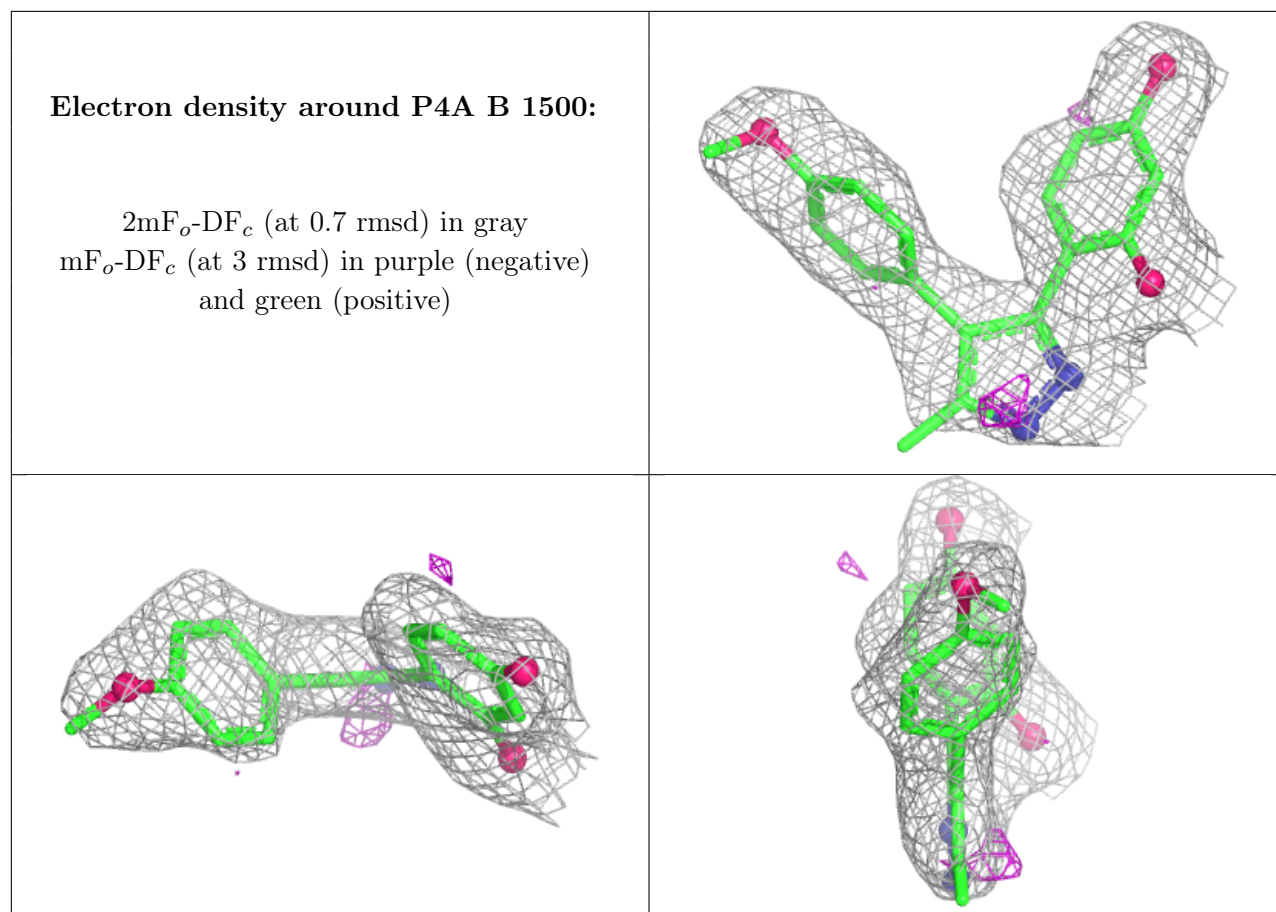
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	P4A	A	500	22/22	0.92	0.10	27,40,52,56	0
2	P4A	B	1500	22/22	0.93	0.10	32,38,47,53	0

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 P4A A 500:

$2mF_o-DF_c$ (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.