



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:22 AM UTC

PDB ID : 4WBB / pdb_00004wbb
Title : Single Turnover Autophosphorylation Cycle of the PKA RIIB Holoenzyme
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Deposited on : 2014-09-02
Resolution : 2.80 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

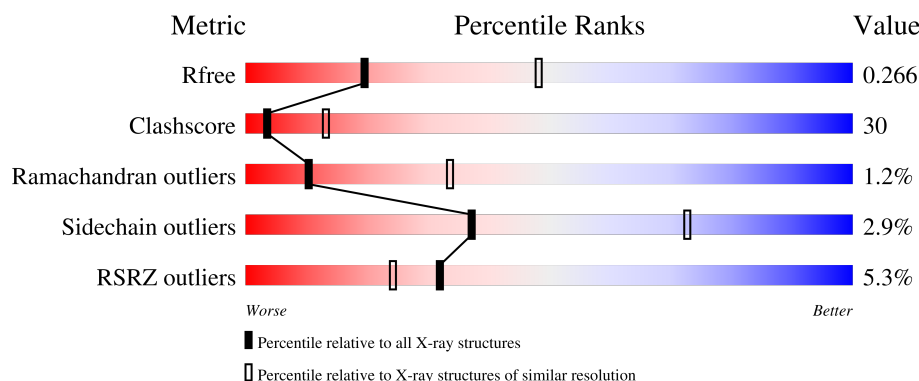
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	416	5% 32% 28% • 35%
2	B	350	4% 60% 33% • •

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	SEP	B	338	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4986 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 cAMP-dependent protein kinase type II-beta regulatory subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	P	S	0	0	0
			2128	1340	372	400	1	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	230	LYS	ARG	engineered mutation	UNP P31324

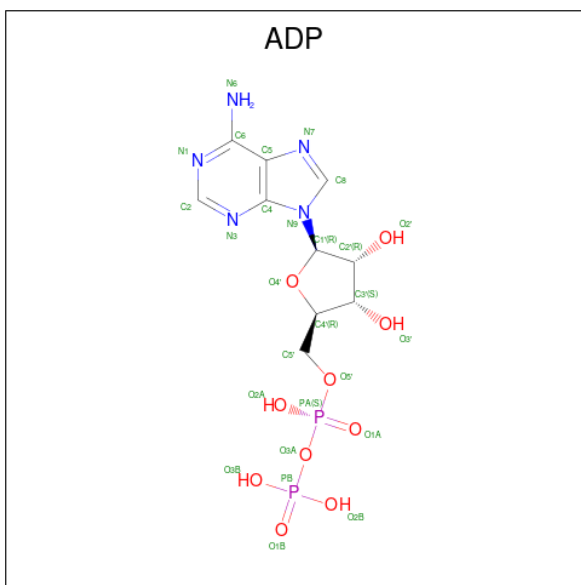
- Molecule 2 is a protein called cAMP-dependent protein kinase catalytic subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	337	Total	C	N	O	P	S	0	0	0
			2787	1803	466	507	3	8			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

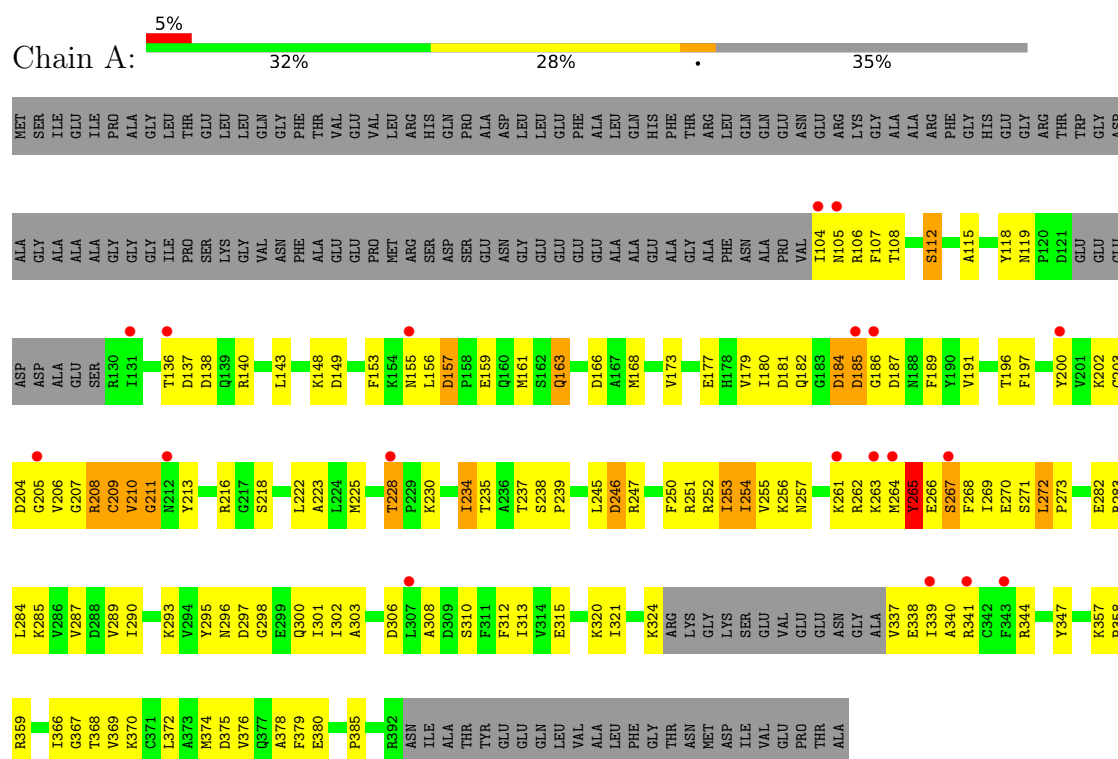
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	24	Total O 24 24	0	0
5	B	18	Total O 18 18	0	0

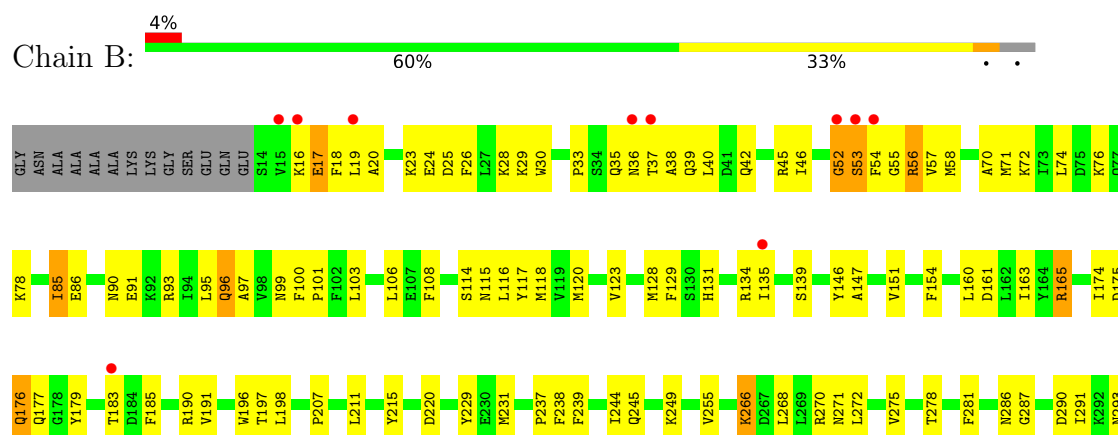
3 Residue-property plots

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: cAMP-dependent protein kinase type II-beta regulatory subunit



- Molecule 2: cAMP-dependent protein kinase catalytic subunit alpha





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	150.38Å 213.95Å 61.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.0 (50.00-2.80) 91.6 (50.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.65Å)	Xtriage
Refinement program		Depositor
R, R_{free}	0.232 , 0.279 (Not available) , 0.266	Depositor DCC
R_{free} test set	1271 reflections (4.36%)	wwPDB-VP
Wilson B-factor (Å ²)	56.6	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4986	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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: ADP, SEP, TPO, 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	A	0.63	1/2149 (0.0%)	1.13	18/2886 (0.6%)
2	B	0.58	0/2824	0.99	12/3801 (0.3%)
All	All	0.60	1/4973 (0.0%)	1.05	30/6687 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	157	ASP	CA-C	-6.15	1.47	1.53

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	357	LYS	CA-C-N	9.75	132.03	119.84
1	A	357	LYS	C-N-CA	9.75	132.03	119.84
1	A	209	CYS	N-CA-C	-9.48	94.91	110.17
2	B	46	ILE	N-CA-C	9.34	119.24	110.74
1	A	157	ASP	N-CA-C	-7.36	101.14	109.60
1	A	228	THR	CA-C-N	7.35	129.03	119.84
1	A	228	THR	C-N-CA	7.35	129.03	119.84
2	B	183	THR	N-CA-C	6.62	121.98	113.18
2	B	16	LYS	N-CA-C	6.60	119.76	111.24
1	A	272	LEU	CA-C-N	6.60	126.84	119.32
1	A	272	LEU	C-N-CA	6.60	126.84	119.32
2	B	165	ARG	N-CA-C	6.58	121.08	113.38
2	B	191	VAL	N-CA-C	6.30	116.69	107.37
1	A	208	ARG	CA-C-N	-6.29	111.94	122.64
1	A	208	ARG	C-N-CA	-6.29	111.94	122.64
1	A	265	TYR	CA-C-N	6.02	129.16	120.79
1	A	265	TYR	C-N-CA	6.02	129.16	120.79
2	B	46	ILE	CB-CA-C	-5.93	104.11	111.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	293	ASN	N-CA-C	-5.76	106.87	112.97
1	A	254	ILE	N-CA-C	5.67	116.40	110.62
2	B	17	GLU	N-CA-C	-5.62	104.79	111.03
1	A	203	CYS	N-CA-C	5.54	118.83	111.75
2	B	52	GLY	CA-C-N	5.49	131.85	123.23
2	B	52	GLY	C-N-CA	5.49	131.85	123.23
2	B	315	ILE	CA-C-N	-5.23	114.39	119.78
2	B	315	ILE	C-N-CA	-5.23	114.39	119.78
1	A	119	ASN	CB-CA-C	5.22	116.08	110.08
1	A	267	SER	N-CA-C	-5.21	102.78	111.37
1	A	157	ASP	N-CA-CB	5.11	117.27	109.75
1	A	378	ALA	N-CA-C	-5.05	105.67	111.07

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	2128	0	2132	163	0
2	B	2787	0	2767	135	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	12	1	0
5	A	24	0	0	35	0
5	B	18	0	0	9	0
All	All	4986	0	4911	291	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 30.

All (291) 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:B:53:SER:HA	2:B:54:PHE:CD1	1.42	1.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:33:PRO:HB3	2:B:96:GLN:NE2	1.34	1.36
1:A:368:THR:HA	5:A:623:HOH:O	1.21	1.27
1:A:337:VAL:HG12	5:A:602:HOH:O	1.37	1.21
1:A:257:ASN:N	5:A:618:HOH:O	1.76	1.18
1:A:204:ASP:HB3	1:A:205:GLY:HA3	1.29	1.15
1:A:200:TYR:HB2	1:A:208:ARG:O	1.46	1.14
2:B:53:SER:CA	2:B:54:PHE:CD1	2.30	1.13
2:B:33:PRO:CB	2:B:96:GLN:NE2	2.12	1.13
2:B:318:PHE:HD1	2:B:319:LYS:O	1.35	1.09
1:A:204:ASP:CB	1:A:205:GLY:HA3	1.83	1.07
2:B:37:THR:OG1	2:B:108:PHE:HB3	1.52	1.07
2:B:146:TYR:HB2	2:B:231:MET:HE1	1.38	1.05
1:A:181:ASP:O	1:A:184:ASP:HB2	1.60	1.01
2:B:53:SER:HB3	2:B:54:PHE:CE1	1.99	0.97
1:A:186:GLY:O	1:A:223:ALA:HB1	1.65	0.97
2:B:318:PHE:CD1	2:B:319:LYS:O	2.17	0.97
2:B:85:ILE:HB	5:B:506:HOH:O	1.64	0.95
1:A:230:LYS:HG2	5:A:611:HOH:O	1.66	0.94
2:B:33:PRO:HB3	2:B:96:GLN:HE22	0.88	0.94
1:A:263:LYS:HE2	5:A:610:HOH:O	1.68	0.94
1:A:168:MET:HE3	1:A:245:LEU:HB2	1.49	0.93
2:B:29:LYS:N	5:B:517:HOH:O	1.97	0.93
1:A:337:VAL:CG1	5:A:602:HOH:O	2.01	0.93
1:A:155:ASN:HD21	2:B:211:LEU:HD22	1.31	0.92
1:A:197:PHE:HB2	1:A:213:TYR:HB2	1.48	0.92
1:A:263:LYS:HG2	1:A:266:GLU:OE2	1.68	0.92
1:A:155:ASN:HD21	2:B:211:LEU:CD2	1.83	0.90
1:A:186:GLY:O	1:A:223:ALA:CB	2.20	0.89
1:A:254:ILE:C	5:A:618:HOH:O	2.15	0.89
1:A:210:VAL:HG12	1:A:211:GLY:H	1.37	0.89
2:B:91:GLU:HG2	2:B:118:MET:HE1	1.55	0.89
1:A:185:ASP:HA	5:A:616:HOH:O	1.74	0.87
2:B:25:ASP:O	5:B:517:HOH:O	1.91	0.87
1:A:209:CYS:O	5:A:601:HOH:O	1.92	0.87
1:A:204:ASP:HB3	1:A:205:GLY:CA	2.04	0.85
1:A:337:VAL:O	5:A:602:HOH:O	1.95	0.84
1:A:104:ILE:HG22	1:A:105:ASN:H	1.40	0.84
2:B:146:TYR:CB	2:B:231:MET:HE1	2.08	0.83
1:A:185:ASP:CG	5:A:616:HOH:O	2.21	0.83
2:B:268:LEU:HB2	2:B:294:HIS:CE1	2.15	0.82
1:A:191:VAL:HG12	1:A:218:SER:HB3	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ASP:O	5:A:609:HOH:O	1.97	0.81
1:A:143:LEU:CD1	1:A:191:VAL:HG21	2.11	0.80
2:B:33:PRO:CB	2:B:96:GLN:HE21	1.90	0.80
1:A:180:ILE:HD13	1:A:230:LYS:HD2	1.65	0.79
1:A:295:TYR:HB2	1:A:369:VAL:HG22	1.64	0.78
1:A:155:ASN:OD1	1:A:155:ASN:O	2.01	0.78
1:A:266:GLU:O	1:A:269:ILE:N	2.16	0.78
2:B:96:GLN:O	2:B:96:GLN:HG2	1.84	0.77
1:A:155:ASN:ND2	2:B:211:LEU:HD22	1.99	0.76
2:B:281:PHE:HE2	2:B:294:HIS:HD2	1.30	0.76
2:B:52:GLY:C	2:B:55:GLY:HA3	2.11	0.76
1:A:367:GLY:O	5:A:623:HOH:O	2.01	0.76
2:B:53:SER:HA	2:B:54:PHE:HD1	0.88	0.76
1:A:184:ASP:O	5:A:611:HOH:O	2.02	0.75
1:A:254:ILE:O	5:A:618:HOH:O	2.01	0.75
2:B:20:ALA:HA	2:B:23:LYS:HE2	1.68	0.75
1:A:302:ILE:HG21	1:A:359:ARG:HD2	1.67	0.75
1:A:301:ILE:HG22	1:A:302:ILE:HG13	1.67	0.74
2:B:20:ALA:O	2:B:24:GLU:HG3	1.87	0.74
1:A:118:TYR:OH	1:A:256:LYS:HE2	1.88	0.74
1:A:191:VAL:HG12	1:A:218:SER:CB	2.18	0.74
2:B:53:SER:CA	2:B:54:PHE:HD1	1.82	0.74
1:A:107:PHE:O	5:A:603:HOH:O	2.06	0.73
1:A:208:ARG:NH2	5:A:605:HOH:O	2.21	0.72
2:B:238:PHE:CE1	2:B:255:VAL:HG22	2.24	0.72
1:A:209:CYS:C	5:A:601:HOH:O	2.31	0.71
1:A:272:LEU:HD12	1:A:273:PRO:HD2	1.70	0.71
1:A:295:TYR:O	5:A:623:HOH:O	2.08	0.71
2:B:96:GLN:HB2	2:B:106:LEU:HD23	1.72	0.71
1:A:297:ASP:OD2	5:A:621:HOH:O	2.08	0.71
2:B:54:PHE:HB3	2:B:78:LYS:HD3	1.73	0.71
1:A:204:ASP:CB	1:A:205:GLY:CA	2.66	0.71
1:A:138:ASP:OD1	5:A:604:HOH:O	2.10	0.70
2:B:271:ASN:HB3	2:B:281:PHE:CD1	2.26	0.69
1:A:196:THR:HB	5:A:624:HOH:O	1.92	0.69
1:A:255:VAL:C	5:A:618:HOH:O	2.35	0.69
1:A:298:GLY:N	5:A:621:HOH:O	2.25	0.68
2:B:24:GLU:O	2:B:28:LYS:HD3	1.93	0.68
1:A:310:SER:HA	1:A:376:VAL:HG23	1.75	0.68
1:A:143:LEU:HD12	1:A:191:VAL:HG21	1.75	0.68
1:A:153:PHE:HB3	1:A:161:MET:HE1	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:ARG:O	1:A:265:TYR:N	2.27	0.68
1:A:338:GLU:OE2	1:A:341:ARG:CD	2.41	0.68
2:B:55:GLY:O	2:B:56:ARG:HB2	1.94	0.68
2:B:268:LEU:HD13	2:B:294:HIS:CD2	2.30	0.67
2:B:53:SER:CB	2:B:54:PHE:CD1	2.78	0.67
1:A:143:LEU:CD1	1:A:191:VAL:CG2	2.72	0.67
2:B:33:PRO:HG3	2:B:93:ARG:NH1	2.09	0.67
2:B:53:SER:CB	2:B:54:PHE:CE1	2.76	0.67
2:B:37:THR:OG1	2:B:108:PHE:CB	2.36	0.66
1:A:210:VAL:HG12	1:A:211:GLY:N	2.11	0.66
1:A:337:VAL:CB	5:A:602:HOH:O	2.36	0.66
1:A:112:SEP:O3P	5:A:607:HOH:O	2.13	0.65
1:A:320:LYS:HB2	1:A:366:ILE:HD11	1.78	0.65
1:A:283:ARG:O	1:A:287:VAL:HG23	1.96	0.65
1:A:197:PHE:N	5:A:624:HOH:O	2.14	0.65
1:A:108:THR:O	5:A:613:HOH:O	2.14	0.64
2:B:76:LYS:HG2	2:B:116:LEU:CD1	2.28	0.64
2:B:294:HIS:CG	2:B:295:LYS:H	2.15	0.64
1:A:157:ASP:OD2	1:A:261:LYS:CE	2.45	0.64
2:B:207:PRO:HG3	2:B:275:VAL:HG12	1.78	0.64
2:B:38:ALA:HB1	2:B:42:GLN:OE1	1.98	0.63
1:A:310:SER:HB3	1:A:375:ASP:HA	1.79	0.63
2:B:29:LYS:HB2	5:B:517:HOH:O	1.99	0.63
1:A:295:TYR:HB2	1:A:369:VAL:CG2	2.28	0.62
1:A:245:LEU:HD21	1:A:250:PHE:CD1	2.34	0.62
1:A:315:GLU:CD	1:A:370:LYS:HZ1	2.07	0.62
2:B:336:ARG:NH2	2:B:338:SEP:O2P	2.33	0.61
1:A:262:ARG:HD2	2:B:196:TRP:CH2	2.36	0.61
1:A:338:GLU:OE2	1:A:341:ARG:HD3	2.01	0.61
1:A:155:ASN:ND2	2:B:211:LEU:CD2	2.58	0.60
2:B:86:GLU:H	2:B:86:GLU:CD	2.09	0.60
1:A:230:LYS:CG	5:A:611:HOH:O	2.33	0.60
1:A:261:LYS:O	1:A:265:TYR:CG	2.55	0.60
1:A:320:LYS:HB2	1:A:366:ILE:CD1	2.32	0.60
2:B:270:ARG:NH2	5:B:501:HOH:O	2.36	0.59
1:A:264:MET:O	5:A:612:HOH:O	2.16	0.59
1:A:310:SER:HB2	1:A:374:MET:O	2.03	0.59
2:B:96:GLN:C	2:B:96:GLN:OE1	2.46	0.58
1:A:261:LYS:O	1:A:265:TYR:CD2	2.57	0.58
1:A:262:ARG:O	1:A:265:TYR:HB2	2.03	0.58
2:B:100:PHE:HB3	2:B:103:LEU:HD12	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:96:GLN:O	2:B:96:GLN:CG	2.52	0.58
2:B:272:LEU:HD21	2:B:291:ILE:HD11	1.86	0.58
2:B:103:LEU:HD22	2:B:185:PHE:HZ	1.69	0.57
1:A:267:SER:HA	1:A:270:GLU:HG2	1.86	0.57
2:B:115:ASN:HB2	2:B:117:TYR:CE2	2.38	0.57
1:A:196:THR:OG1	1:A:237:THR:OG1	2.06	0.57
1:A:143:LEU:HD11	1:A:191:VAL:CG2	2.35	0.57
1:A:268:PHE:HB3	1:A:313:ILE:HG21	1.86	0.57
1:A:180:ILE:HD11	1:A:234:ILE:HG13	1.86	0.57
2:B:268:LEU:HD13	2:B:294:HIS:CG	2.40	0.56
2:B:30:TRP:CZ3	2:B:93:ARG:HG2	2.40	0.56
1:A:104:ILE:HG22	1:A:105:ASN:N	2.17	0.56
2:B:72:LYS:HG2	2:B:74:LEU:CD1	2.35	0.56
2:B:266:LYS:O	2:B:270:ARG:HG3	2.06	0.56
1:A:181:ASP:HB2	1:A:184:ASP:CG	2.30	0.56
1:A:163:GLN:HE22	1:A:257:ASN:HD22	1.54	0.55
1:A:272:LEU:HD13	1:A:347:TYR:CD2	2.42	0.55
2:B:229:TYR:CG	2:B:237:PRO:HG3	2.41	0.55
2:B:24:GLU:HA	2:B:28:LYS:NZ	2.20	0.55
1:A:181:ASP:O	1:A:184:ASP:CB	2.46	0.55
1:A:320:LYS:HE3	1:A:338:GLU:HB2	1.89	0.55
2:B:198:LEU:HD23	2:B:198:LEU:C	2.31	0.55
2:B:238:PHE:CE1	2:B:255:VAL:CG2	2.90	0.55
1:A:303:ALA:O	1:A:306:ASP:HB2	2.06	0.54
2:B:52:GLY:O	2:B:55:GLY:HA3	2.06	0.54
2:B:76:LYS:HG2	2:B:116:LEU:HD12	1.89	0.54
2:B:336:ARG:HH21	2:B:338:SEP:HB3	1.71	0.54
2:B:314:PHE:C	2:B:315:ILE:HD12	2.33	0.54
2:B:179:TYR:CE1	2:B:308:ARG:HA	2.42	0.54
1:A:296:ASN:HA	5:A:623:HOH:O	2.08	0.54
1:A:186:GLY:O	1:A:223:ALA:HB2	2.05	0.53
1:A:337:VAL:HB	5:A:602:HOH:O	2.01	0.53
2:B:177:GLN:O	2:B:177:GLN:HG2	2.09	0.53
2:B:146:TYR:HD1	2:B:231:MET:CE	2.22	0.53
1:A:338:GLU:OE2	1:A:341:ARG:HG3	2.09	0.53
1:A:143:LEU:HD11	1:A:191:VAL:HG21	1.89	0.53
2:B:286:ASN:HB2	2:B:290:ASP:OD2	2.09	0.53
1:A:148:LYS:NZ	1:A:149:ASP:OD2	2.39	0.52
1:A:206:VAL:HG12	1:A:208:ARG:HG3	1.91	0.52
1:A:339:ILE:O	1:A:340:ALA:HB2	2.08	0.52
2:B:160:LEU:O	2:B:161:ASP:HB2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:281:PHE:CE2	2:B:294:HIS:HD2	2.17	0.52
1:A:140:ARG:NH1	1:A:166:ASP:OD1	2.39	0.52
2:B:163:ILE:HG13	2:B:165:ARG:HG3	1.92	0.52
2:B:245:GLN:O	2:B:249:LYS:HG3	2.10	0.52
2:B:349:GLU:O	2:B:349:GLU:HG2	2.08	0.52
1:A:155:ASN:HD21	2:B:211:LEU:HD21	1.69	0.52
2:B:93:ARG:O	2:B:96:GLN:OE1	2.27	0.52
1:A:256:LYS:N	5:A:618:HOH:O	2.42	0.51
1:A:238:SER:HB2	1:A:239:PRO:HD2	1.93	0.51
1:A:315:GLU:CD	1:A:370:LYS:NZ	2.69	0.51
1:A:272:LEU:HD13	1:A:347:TYR:CG	2.45	0.51
2:B:52:GLY:C	2:B:55:GLY:CA	2.82	0.51
1:A:115:ALA:HB2	1:A:225:MET:O	2.11	0.51
1:A:173:VAL:HA	1:A:177:GLU:OE1	2.11	0.51
2:B:268:LEU:CB	2:B:294:HIS:CE1	2.91	0.50
1:A:196:THR:CB	1:A:237:THR:HG1	2.25	0.50
1:A:302:ILE:HD13	1:A:359:ARG:HH11	1.75	0.50
2:B:17:GLU:HG2	2:B:18:PHE:N	2.26	0.50
1:A:282:GLU:O	1:A:285:LYS:HB2	2.10	0.50
1:A:157:ASP:OD2	1:A:261:LYS:HE2	2.12	0.50
2:B:90:ASN:O	2:B:91:GLU:C	2.55	0.50
2:B:128:MET:O	2:B:129:PHE:C	2.55	0.50
1:A:104:ILE:C	1:A:106:ARG:H	2.19	0.50
1:A:376:VAL:O	1:A:379:PHE:HB3	2.12	0.50
1:A:136:THR:HG22	1:A:137:ASP:N	2.26	0.49
1:A:216:ARG:HH11	1:A:216:ARG:HG3	1.76	0.49
1:A:284:LEU:HA	1:A:287:VAL:HG23	1.95	0.49
2:B:95:LEU:CD2	2:B:120:MET:HE2	2.42	0.49
2:B:294:HIS:CG	2:B:295:LYS:N	2.81	0.49
2:B:35:GLN:O	2:B:37:THR:N	2.46	0.49
2:B:147:ALA:O	2:B:151:VAL:HG23	2.13	0.49
1:A:182:GLN:C	1:A:184:ASP:H	2.20	0.49
1:A:181:ASP:C	1:A:184:ASP:HB2	2.35	0.48
1:A:338:GLU:OE2	1:A:341:ARG:CG	2.62	0.48
1:A:155:ASN:OD1	1:A:155:ASN:C	2.57	0.48
1:A:204:ASP:HB2	1:A:205:GLY:HA3	1.83	0.48
2:B:33:PRO:CA	2:B:96:GLN:NE2	2.77	0.48
1:A:186:GLY:CA	1:A:230:LYS:HZ2	2.26	0.48
2:B:215:TYR:CD1	2:B:215:TYR:C	2.91	0.48
2:B:20:ALA:O	2:B:23:LYS:HG2	2.14	0.48
2:B:281:PHE:HE2	2:B:294:HIS:CD2	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ASP:HB2	1:A:184:ASP:OD2	2.14	0.47
1:A:222:LEU:HD23	1:A:228:THR:HG21	1.95	0.47
1:A:200:TYR:CB	1:A:208:ARG:O	2.39	0.47
1:A:186:GLY:HA3	1:A:230:LYS:HZ2	1.79	0.47
4:A:502:ADP:N6	2:B:70:ALA:HB3	2.30	0.47
1:A:266:GLU:O	1:A:267:SER:C	2.58	0.47
2:B:338:SEP:O1P	2:B:340:ASN:HB2	2.15	0.47
1:A:187:ASP:O	1:A:246:ASP:HA	2.15	0.46
1:A:191:VAL:HG12	1:A:218:SER:HB2	1.95	0.46
1:A:290:ILE:HD12	1:A:374:MET:HB2	1.97	0.46
1:A:293:LYS:HD3	1:A:295:TYR:OH	2.15	0.46
2:B:55:GLY:O	2:B:56:ARG:CB	2.62	0.46
2:B:76:LYS:HG2	2:B:116:LEU:HD11	1.97	0.46
1:A:157:ASP:OD2	1:A:261:LYS:NZ	2.48	0.46
1:A:271:SER:O	1:A:273:PRO:HD3	2.16	0.46
1:A:302:ILE:HD13	1:A:359:ARG:NH1	2.29	0.46
2:B:24:GLU:HA	2:B:28:LYS:HZ3	1.80	0.46
1:A:156:LEU:O	1:A:157:ASP:C	2.56	0.46
1:A:180:ILE:HD11	1:A:234:ILE:CG1	2.45	0.46
2:B:301:ASP:CG	2:B:304:ALA:HB3	2.41	0.46
1:A:252:ARG:O	1:A:253:ILE:C	2.57	0.46
1:A:300:GLN:OE1	1:A:303:ALA:HB2	2.16	0.46
2:B:58:MET:HE2	2:B:58:MET:HB3	1.85	0.46
1:A:312:PHE:HA	1:A:372:LEU:O	2.16	0.46
2:B:291:ILE:O	2:B:297:PHE:HD1	1.99	0.46
1:A:181:ASP:O	1:A:184:ASP:N	2.46	0.45
2:B:146:TYR:CD1	2:B:231:MET:CE	2.99	0.45
1:A:184:ASP:HB3	5:A:609:HOH:O	2.15	0.45
2:B:96:GLN:OE1	2:B:97:ALA:HB2	2.16	0.45
1:A:136:THR:HG22	1:A:138:ASP:H	1.82	0.45
2:B:239:PHE:CD1	2:B:239:PHE:C	2.94	0.45
2:B:312:ALA:C	2:B:314:PHE:H	2.24	0.45
1:A:379:PHE:C	1:A:379:PHE:CD2	2.95	0.45
2:B:135:ILE:O	2:B:135:ILE:HG22	2.16	0.45
2:B:315:ILE:HD12	2:B:315:ILE:N	2.32	0.45
1:A:324:LYS:O	1:A:324:LYS:HG2	2.17	0.45
2:B:20:ALA:HA	2:B:23:LYS:HG2	1.99	0.44
2:B:26:PHE:HE1	2:B:190:ARG:NH1	2.14	0.44
2:B:86:GLU:CD	2:B:86:GLU:N	2.76	0.44
2:B:116:LEU:HD23	2:B:350:PHE:CD2	2.52	0.44
1:A:136:THR:CG2	1:A:137:ASP:N	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ASP:HA	1:A:247:ARG:HH12	1.82	0.44
1:A:302:ILE:CG2	1:A:359:ARG:HD2	2.41	0.44
2:B:134:ARG:NH1	5:B:502:HOH:O	2.48	0.44
2:B:335:ILE:O	2:B:335:ILE:HG22	2.18	0.44
2:B:176:GLN:HG2	5:B:507:HOH:O	2.17	0.44
2:B:36:ASN:OD1	2:B:38:ALA:O	2.35	0.44
2:B:54:PHE:HA	2:B:55:GLY:HA3	1.64	0.44
2:B:154:PHE:CE1	2:B:220:ASP:HB3	2.53	0.44
1:A:186:GLY:HA2	1:A:230:LYS:NZ	2.33	0.43
2:B:72:LYS:HG2	2:B:74:LEU:HD11	2.00	0.43
1:A:251:ARG:HG2	1:A:251:ARG:HH21	1.83	0.43
1:A:156:LEU:HD23	1:A:156:LEU:HA	1.81	0.42
1:A:157:ASP:HB3	1:A:159:GLU:H	1.83	0.42
2:B:101:PRO:HG2	2:B:306:TYR:CE2	2.54	0.42
2:B:28:LYS:HD3	2:B:28:LYS:N	2.34	0.42
2:B:229:TYR:CD2	2:B:229:TYR:C	2.97	0.42
1:A:189:PHE:CE2	1:A:191:VAL:HG13	2.55	0.42
2:B:281:PHE:CE2	2:B:294:HIS:CD2	3.02	0.42
1:A:310:SER:CB	1:A:375:ASP:HA	2.48	0.42
2:B:95:LEU:HD21	2:B:120:MET:HE2	2.01	0.42
1:A:202:LYS:HA	1:A:207:GLY:HA2	2.01	0.42
2:B:116:LEU:HD23	2:B:350:PHE:CG	2.54	0.42
1:A:185:ASP:CA	5:A:616:HOH:O	2.50	0.42
2:B:114:SER:OG	2:B:338:SEP:HB2	2.20	0.42
2:B:39:GLN:HG2	2:B:42:GLN:NE2	2.35	0.42
2:B:175:ASP:C	2:B:175:ASP:OD1	2.61	0.42
1:A:308:ALA:CA	1:A:359:ARG:HH21	2.32	0.41
2:B:33:PRO:HG3	2:B:93:ARG:HH12	1.83	0.41
2:B:131:HIS:CE1	2:B:314:PHE:HZ	2.38	0.41
2:B:40:LEU:HD11	2:B:45:ARG:HH21	1.86	0.41
2:B:123:VAL:HG12	2:B:174:ILE:O	2.20	0.41
2:B:301:ASP:OD1	2:B:304:ALA:HB3	2.20	0.41
2:B:52:GLY:C	5:B:511:HOH:O	2.64	0.41
1:A:196:THR:CB	1:A:237:THR:OG1	2.68	0.41
1:A:315:GLU:O	1:A:344:ARG:HG3	2.20	0.40
1:A:380:GLU:HG2	1:A:385:PRO:HA	2.03	0.40
1:A:163:GLN:NE2	1:A:257:ASN:HD22	2.17	0.40
1:A:206:VAL:HG12	1:A:208:ARG:CG	2.51	0.40
1:A:289:VAL:HG21	2:B:278:THR:CG2	2.52	0.40
2:B:57:VAL:HA	2:B:71:MET:O	2.21	0.40
2:B:134:ARG:NH2	5:B:503:HOH:O	2.41	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:286:ASN:O	2:B:287:GLY:C	2.63	0.40
1:A:251:ARG:HG2	1:A:251:ARG:NH2	2.37	0.40
2:B:268:LEU:HA	2:B:268:LEU:HD12	1.78	0.40

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	A	262/416 (63%)	235 (90%)	23 (9%)	4 (2%)	8	28
2	B	332/350 (95%)	291 (88%)	38 (11%)	3 (1%)	14	41
All	All	594/766 (78%)	526 (89%)	61 (10%)	7 (1%)	10	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	56	ARG
1	A	211	GLY
1	A	210	VAL
2	B	19	LEU
2	B	85	ILE
1	A	253	ILE
1	A	358	PRO

5.3.2 Protein sidechains [i](#)

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	227/339 (67%)	218 (96%)	9 (4%)	28	63
2	B	295/302 (98%)	289 (98%)	6 (2%)	48	80
All	All	522/641 (81%)	507 (97%)	15 (3%)	37	73

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	GLN
1	A	179	VAL
1	A	184	ASP
1	A	185	ASP
1	A	234	ILE
1	A	235	THR
1	A	246	ASP
1	A	265	TYR
1	A	321	ILE
2	B	53	SER
2	B	96	GLN
2	B	99	ASN
2	B	176	GLN
2	B	244	ILE
2	B	266	LYS

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

Mol	Chain	Res	Type
1	A	119	ASN
1	A	155	ASN
1	A	163	GLN
1	A	300	GLN
1	A	364	HIS
2	B	68	HIS
2	B	115	ASN
2	B	274	GLN
2	B	326	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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
2	SEP	B	338	2	8,9,10	1.52	1 (12%)	7,12,14	2.25	2 (28%)
1	SEP	A	112	3,1	8,9,10	1.65	2 (25%)	7,12,14	0.93	0
2	SEP	B	139	2	8,9,10	1.70	1 (12%)	7,12,14	1.58	1 (14%)
2	TPO	B	197	2	8,10,11	1.77	2 (25%)	10,14,16	2.29	2 (20%)

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	SEP	B	338	2	-	3/6/8/10	-
1	SEP	A	112	3,1	-	1/6/8/10	-
2	SEP	B	139	2	-	3/6/8/10	-
2	TPO	B	197	2	-	0/9/11/13	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	139	SEP	P-O1P	3.67	1.61	1.50
1	A	112	SEP	P-O1P	3.53	1.61	1.50
2	B	197	TPO	P-O1P	3.45	1.61	1.50
2	B	338	SEP	P-O1P	3.35	1.60	1.50
2	B	197	TPO	P-O3P	-2.24	1.46	1.54
1	A	112	SEP	P-O3P	2.14	1.62	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	197	TPO	P-OG1-CB	-6.26	106.31	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	338	SEP	OG-CB-CA	4.88	112.89	108.14
2	B	139	SEP	OG-CB-CA	3.71	111.76	108.14
2	B	338	SEP	O2P-P-OG	2.89	114.20	106.67
2	B	197	TPO	CG2-CB-CA	-2.71	107.98	113.26

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	139	SEP	CB-OG-P-O1P
2	B	139	SEP	CB-OG-P-O2P
2	B	338	SEP	N-CA-CB-OG
2	B	338	SEP	C-CA-CB-OG
2	B	338	SEP	CA-CB-OG-P
2	B	139	SEP	CB-OG-P-O3P
1	A	112	SEP	CA-CB-OG-P

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	338	SEP	4	0
1	A	112	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 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
4	ADP	A	502	3	28,29,29	1.42	4 (14%)	43,45,45	1.85	9 (20%)

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
4	ADP	A	502	3	-	5/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	502	ADP	C5-C4	4.74	1.47	1.39
4	A	502	ADP	C5-C6	2.77	1.48	1.41
4	A	502	ADP	C8-N7	2.37	1.36	1.31
4	A	502	ADP	C5-N7	-2.21	1.35	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	502	ADP	C5-C4-N3	-5.85	118.67	126.72
4	A	502	ADP	N3-C4-N9	4.61	135.00	127.17
4	A	502	ADP	C2-N3-C4	3.70	120.87	111.83
4	A	502	ADP	C4-C5-N7	-3.48	106.60	110.58
4	A	502	ADP	N3-C2-N1	-3.23	123.70	128.58
4	A	502	ADP	C4-N9-C8	2.58	108.45	105.74
4	A	502	ADP	C3'-C2'-C1'	2.53	106.25	101.46
4	A	502	ADP	C5-N7-C8	2.52	107.41	103.45
4	A	502	ADP	C6-C5-N7	2.08	136.09	132.09

There are no chirality outliers.

All (5) torsion outliers are listed below:

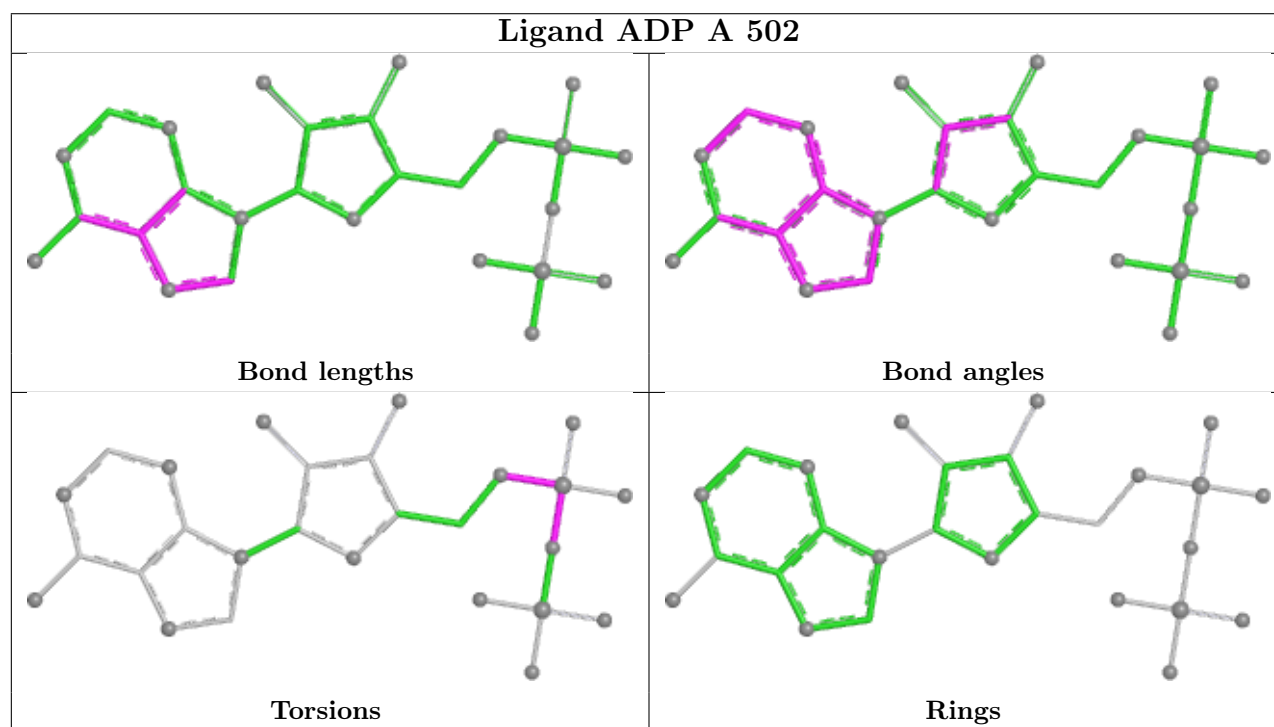
Mol	Chain	Res	Type	Atoms
4	A	502	ADP	C5'-O5'-PA-O2A
4	A	502	ADP	C5'-O5'-PA-O1A
4	A	502	ADP	C5'-O5'-PA-O3A
4	A	502	ADP	PB-O3A-PA-O2A
4	A	502	ADP	PB-O3A-PA-O1A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	502	ADP	1	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 ⓘ

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, 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	268/416 (64%)	0.43	19 (7%) 22 16	32, 56, 82, 110	0
2	B	334/350 (95%)	0.22	13 (3%) 43 34	29, 52, 84, 120	0
All	All	602/766 (78%)	0.31	32 (5%) 32 24	29, 54, 83, 120	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	104	ILE	6.4
1	A	343	PHE	5.5
1	A	186	GLY	5.5
1	A	264	MET	5.5
2	B	295	LYS	5.0
2	B	52	GLY	4.3
2	B	37	THR	3.9
1	A	263	LYS	3.8
1	A	261	LYS	3.8
2	B	53	SER	3.4
2	B	36	ASN	3.3
1	A	185	ASP	3.3
1	A	105	ASN	3.2
1	A	136	THR	3.1
1	A	307	LEU	3.0
2	B	183	THR	3.0
1	A	131	ILE	3.0
1	A	155	ASN	2.9
1	A	212	ASN	2.9
2	B	54	PHE	2.9
1	A	200	TYR	2.9
2	B	16	LYS	2.9
2	B	135	ILE	2.7
2	B	15	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	339	ILE	2.7
2	B	318	PHE	2.7
2	B	294	HIS	2.6
2	B	19	LEU	2.6
1	A	341	ARG	2.4
1	A	205	GLY	2.1
1	A	267	SER	2.1
1	A	228	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	SEP	B	139	10/11	0.90	0.13	53,58,75,81	0
1	SEP	A	112	10/11	0.95	0.08	29,37,50,52	0
2	SEP	B	338	10/11	0.95	0.08	50,59,67,68	0
2	TPO	B	197	11/12	0.98	0.06	31,33,36,37	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

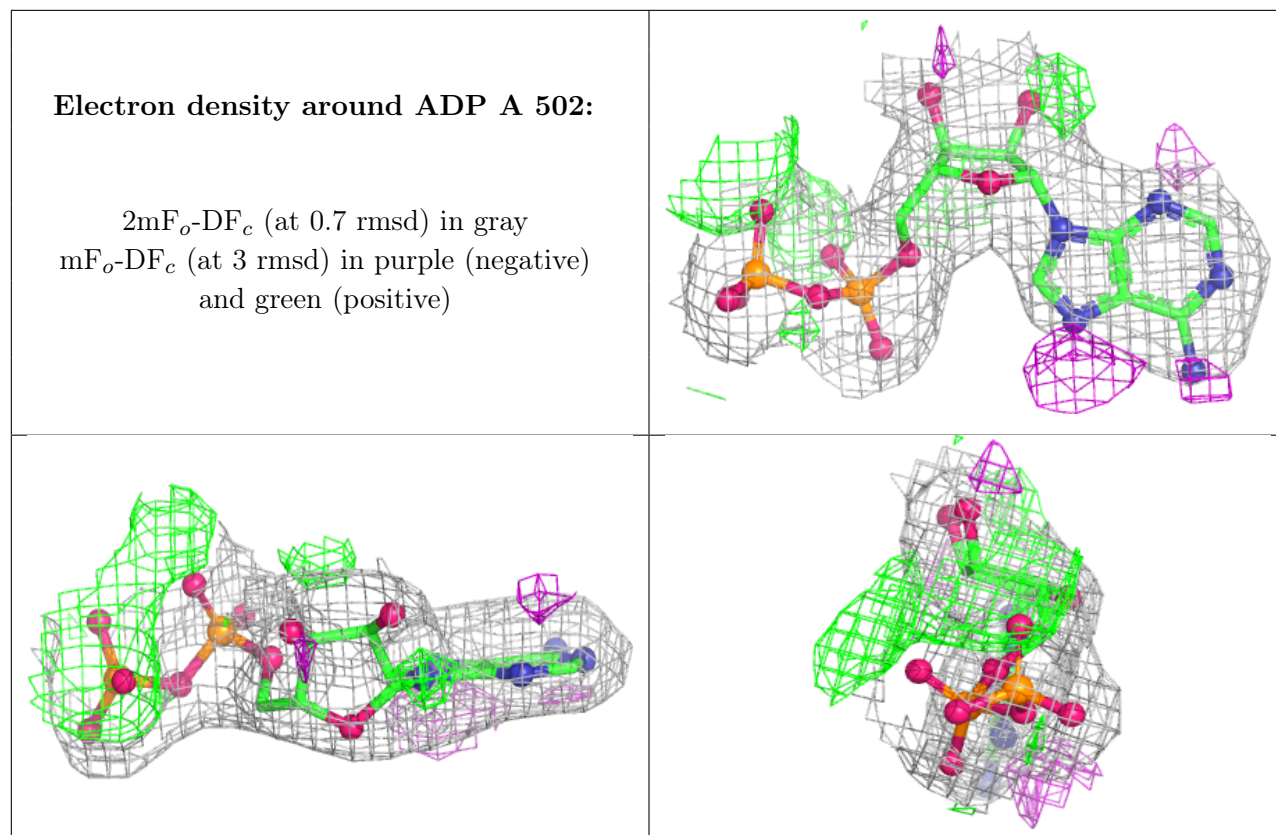
6.4 Ligands [i](#)

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
4	ADP	A	502	27/27	0.90	0.12	32,45,48,50	0
3	CA	B	401	1/1	0.95	0.12	43,43,43,43	0
3	CA	A	501	1/1	0.96	0.05	44,44,44,44	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 ⓘ

There are no such residues in this entry.