



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:36 AM UTC

PDB ID : 3TPV / pdb_00003tpv
Title : Structure of pHipA bound to ADP
Authors : Schumacher, M.A.; Link, T.M.; Brennan, R.G.
Deposited on : 2011-09-08
Resolution : 2.30 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
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

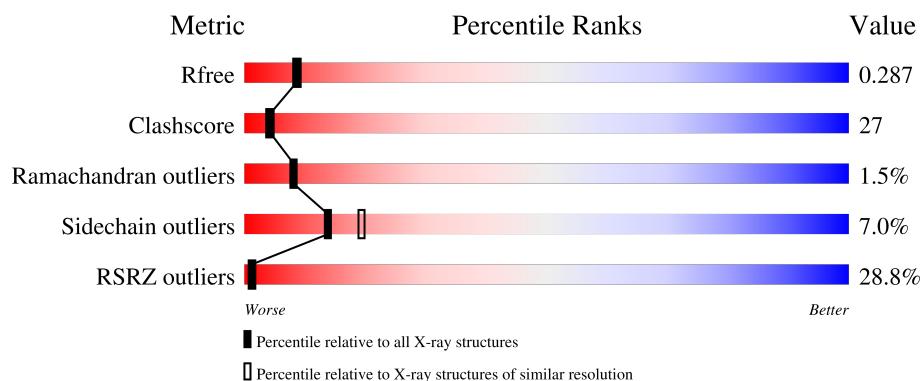
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	B	440	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3500 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.

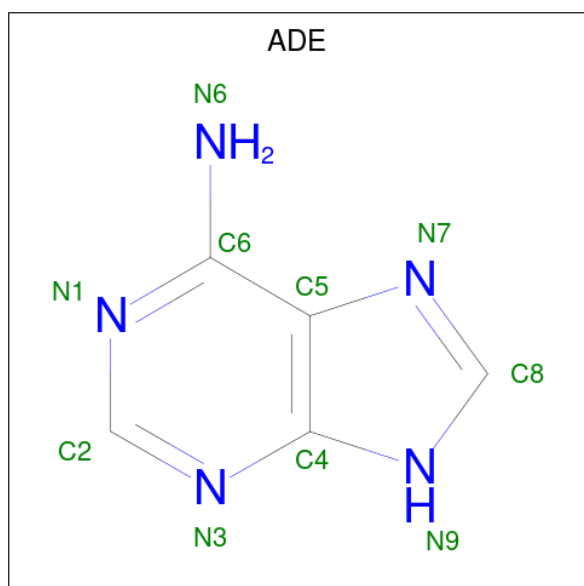
- Molecule 1 is a protein called Serine/threonine-protein kinase HipA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	B	421	3329	2132	582	600	1	14	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	243	ARG	GLU	conflict	UNP P23874

- Molecule 2 is ADENINE (CCD ID: ADE) (formula: C₅H₅N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	N		
2	B	1	10	5	5	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

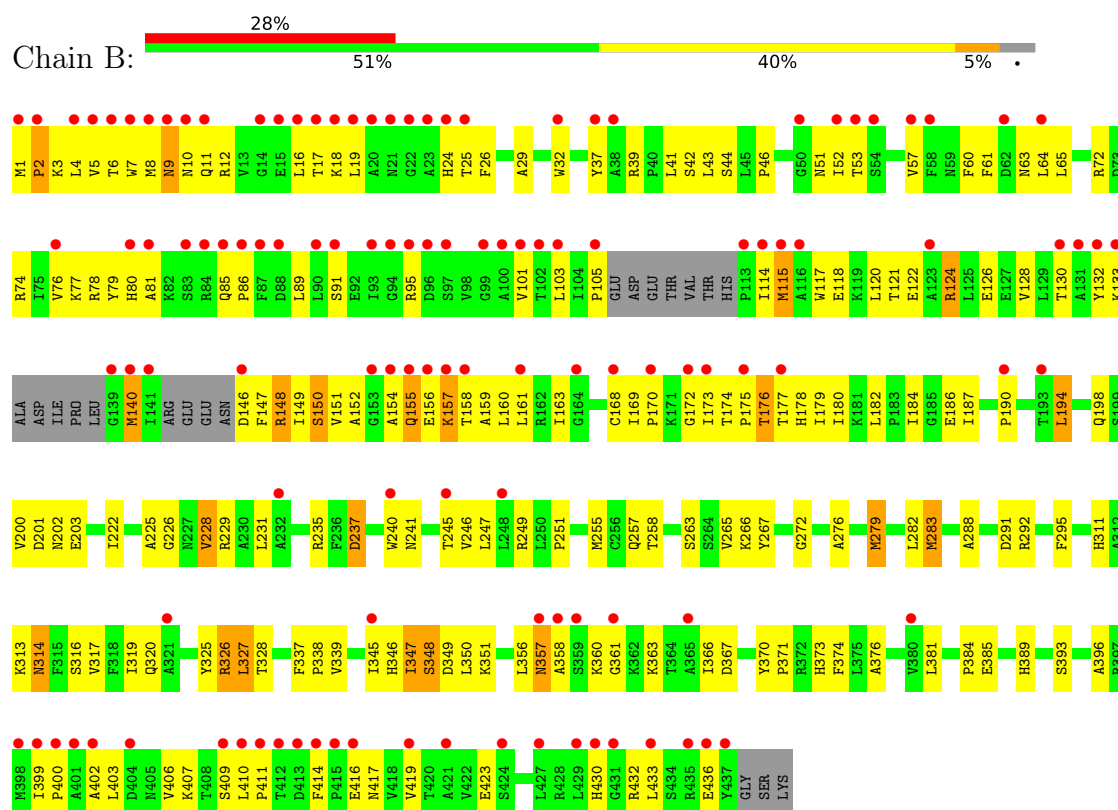
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	156	Total	O	0	0
			156	156		

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: Serine/threonine-protein kinase HipA



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	133.70Å 133.70Å 48.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.60 – 2.30 38.60 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (38.60-2.30) 99.0 (38.60-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 1.89Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.239 , 0.289 0.257 , 0.287	Depositor DCC
R_{free} test set	2853 reflections (7.30%)	wwPDB-VP
Wilson B-factor (Å ²)	36.2	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3500	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: ADE, SO4, SEP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	B	0.94	2/3390 (0.1%)	0.86	1/4587 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	279	MET	CG-SD	8.42	2.01	1.80
1	B	279	MET	SD-CE	-7.32	1.61	1.79

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	237	ASP	N-CA-C	-5.22	106.76	112.72

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	B	3329	0	3383	181	0
2	B	10	0	4	0	0
3	B	5	0	0	0	0
4	B	156	0	0	15	0
All	All	3500	0	3387	181	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 27.

All (181) 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:B:279:MET:SD	1:B:279:MET:CG	2.01	1.46
1:B:159:ALA:HB1	1:B:176:THR:HG21	1.35	1.04
1:B:279:MET:HE1	1:B:292:ARG:HA	1.44	0.98
1:B:198:GLN:NE2	1:B:417:ASN:HD21	1.77	0.82
1:B:198:GLN:HE22	1:B:417:ASN:HD21	1.30	0.78
1:B:241:ASN:ND2	1:B:246:VAL:H	1.81	0.77
1:B:279:MET:SD	1:B:279:MET:CB	2.75	0.74
1:B:326:ARG:HD3	4:B:444:HOH:O	1.90	0.71
1:B:155:GLN:HG3	1:B:184:ILE:HG12	1.73	0.70
1:B:114:ILE:O	1:B:115:MET:HB2	1.90	0.70
1:B:279:MET:HE3	1:B:295:PHE:CD2	2.28	0.69
1:B:241:ASN:HD21	1:B:246:VAL:H	1.40	0.68
1:B:41:LEU:HD11	1:B:101:VAL:HG12	1.75	0.68
1:B:345:ILE:HG22	1:B:346:HIS:N	2.10	0.66
1:B:1:MET:HB3	1:B:18:LYS:H	1.61	0.65
1:B:95:ARG:HG3	1:B:173:ILE:HD12	1.77	0.65
1:B:279:MET:CG	1:B:279:MET:CE	2.75	0.65
1:B:235:ARG:HB3	1:B:237:ASP:OD1	1.97	0.64
1:B:345:ILE:CG2	1:B:349:ASP:HB2	2.27	0.64
1:B:114:ILE:HG13	1:B:115:MET:H	1.62	0.64
1:B:148:ARG:HD2	4:B:546:HOH:O	1.99	0.63
1:B:279:MET:HA	1:B:279:MET:HE2	1.80	0.62
1:B:8:MET:HE2	1:B:11:GLN:HE21	1.64	0.62
1:B:9:ASN:C	1:B:10:ASN:HD22	2.08	0.62
1:B:1:MET:HE2	1:B:17:THR:HA	1.82	0.62
1:B:85:GLN:O	1:B:89:LEU:HD23	2.00	0.61
1:B:346:HIS:HD2	4:B:459:HOH:O	1.84	0.61
1:B:160:LEU:O	1:B:176:THR:HG23	2.01	0.61
1:B:202:ASN:ND2	1:B:338:PRO:HG2	2.16	0.60
1:B:200:VAL:HG13	1:B:231:LEU:HB2	1.84	0.60
1:B:241:ASN:HD21	1:B:246:VAL:N	2.00	0.59
1:B:407:LYS:HE3	1:B:419:VAL:HG11	1.86	0.58
1:B:18:LYS:HG2	1:B:19:LEU:N	2.19	0.58
1:B:345:ILE:HG22	1:B:346:HIS:H	1.68	0.58
1:B:3:LYS:HG3	1:B:17:THR:CG2	2.35	0.57
1:B:114:ILE:HG21	1:B:173:ILE:HG22	1.84	0.57
1:B:1:MET:HB3	1:B:18:LYS:N	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:LEU:HB3	1:B:366:ILE:HD12	1.87	0.56
1:B:357:ASN:ND2	4:B:596:HOH:O	2.39	0.55
1:B:155:GLN:HB2	4:B:540:HOH:O	2.06	0.55
1:B:39:ARG:HD3	1:B:319:ILE:HG22	1.88	0.54
1:B:257:GLN:HA	4:B:467:HOH:O	2.07	0.54
1:B:43:LEU:HD21	1:B:325:TYR:CD1	2.43	0.54
1:B:187:ILE:HB	1:B:194:LEU:HB2	1.89	0.54
1:B:132:TYR:O	1:B:182:LEU:HD13	2.07	0.54
1:B:65:LEU:HD11	1:B:86:PRO:HA	1.90	0.54
1:B:399:ILE:O	1:B:403:LEU:HG	2.08	0.53
1:B:241:ASN:ND2	1:B:246:VAL:N	2.54	0.53
1:B:117:TRP:NE1	1:B:163:ILE:HG13	2.22	0.53
1:B:8:MET:HE2	1:B:11:GLN:NE2	2.23	0.53
1:B:283:MET:HA	1:B:283:MET:HE2	1.91	0.53
1:B:147:PHE:C	1:B:148:ARG:HG2	2.35	0.52
1:B:198:GLN:NE2	1:B:417:ASN:ND2	2.52	0.52
1:B:148:ARG:HD3	1:B:267:TYR:CZ	2.45	0.52
1:B:161:LEU:HD21	1:B:163:ILE:HD11	1.91	0.52
1:B:4:LEU:HD11	1:B:91:SER:HB2	1.92	0.51
1:B:283:MET:HE3	1:B:292:ARG:CD	2.40	0.51
1:B:276:ALA:HA	1:B:381:LEU:HD21	1.93	0.51
1:B:358:ALA:HB2	1:B:363:LYS:HG2	1.92	0.51
1:B:399:ILE:HB	1:B:430:HIS:CD2	2.45	0.51
1:B:158:THR:CG2	1:B:180:ILE:HB	2.40	0.51
1:B:279:MET:HE2	1:B:282:LEU:HD12	1.93	0.51
1:B:407:LYS:HA	1:B:410:LEU:HD12	1.93	0.51
1:B:363:LYS:HG3	1:B:373:HIS:CD2	2.46	0.50
1:B:407:LYS:C	1:B:409:SER:H	2.19	0.50
1:B:2:PRO:O	1:B:17:THR:HG22	2.10	0.50
1:B:18:LYS:HG2	1:B:19:LEU:H	1.75	0.50
1:B:44:SER:C	1:B:46:PRO:HD3	2.37	0.50
1:B:291:ASP:CG	1:B:327:LEU:HB2	2.37	0.50
1:B:399:ILE:HD13	1:B:430:HIS:HB2	1.94	0.50
1:B:152:ALA:HB3	4:B:524:HOH:O	2.12	0.49
1:B:396:ALA:HB2	1:B:433:LEU:HB3	1.95	0.49
1:B:8:MET:O	1:B:9:ASN:HB2	2.14	0.48
1:B:399:ILE:HB	1:B:400:PRO:HD3	1.95	0.48
1:B:6:THR:HA	1:B:103:LEU:HD23	1.95	0.48
1:B:72:ARG:O	1:B:76:VAL:HG23	2.14	0.48
1:B:118:GLU:HG2	1:B:169:ILE:O	2.14	0.48
1:B:194:LEU:CD1	1:B:345:ILE:HD11	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:GLU:CD	1:B:226:GLY:HA3	2.39	0.48
1:B:283:MET:HE3	1:B:292:ARG:HD3	1.95	0.48
1:B:26:PHE:O	1:B:51:ASN:HA	2.14	0.48
1:B:337:PHE:HB2	1:B:338:PRO:HD3	1.96	0.48
1:B:345:ILE:CG2	1:B:346:HIS:N	2.76	0.48
1:B:179:ILE:HG13	1:B:235:ARG:HG2	1.96	0.47
1:B:222:ILE:HD13	1:B:229:ARG:HH12	1.79	0.47
1:B:201:ASP:OD2	1:B:229:ARG:NH1	2.48	0.47
1:B:12:ARG:NH2	4:B:529:HOH:O	2.48	0.47
1:B:126:GLU:O	1:B:130:THR:HG23	2.14	0.47
1:B:265:VAL:O	1:B:272:GLY:HA3	2.14	0.47
1:B:411:PRO:HG2	1:B:414:PHE:HB2	1.97	0.47
1:B:203:GLU:HG2	1:B:231:LEU:HD21	1.97	0.47
1:B:229:ARG:HH11	1:B:229:ARG:HG3	1.80	0.47
1:B:385:GLU:HG2	1:B:389:HIS:CD2	2.50	0.47
1:B:225:ALA:O	1:B:228:VAL:HG13	2.15	0.46
1:B:370:TYR:HB3	1:B:371:PRO:HD2	1.97	0.46
1:B:120:LEU:HD21	1:B:169:ILE:HD12	1.96	0.46
1:B:132:TYR:HB2	1:B:182:LEU:CD1	2.45	0.46
1:B:25:THR:HG22	1:B:53:THR:HG22	1.96	0.46
1:B:78:ARG:HH12	1:B:132:TYR:HD2	1.63	0.46
1:B:349:ASP:HA	1:B:351:LYS:NZ	2.31	0.46
1:B:255:MET:O	1:B:258:THR:HG22	2.16	0.46
1:B:130:THR:HA	1:B:133:LYS:HE3	1.96	0.46
1:B:370:TYR:CD2	1:B:432:ARG:HD3	2.50	0.46
1:B:120:LEU:HD22	1:B:124:ARG:HD2	1.98	0.45
1:B:198:GLN:HA	4:B:487:HOH:O	2.15	0.45
1:B:152:ALA:HB2	1:B:187:ILE:HD12	1.98	0.45
1:B:4:LEU:HD23	1:B:105:PRO:HA	1.97	0.45
1:B:347:ILE:CG2	1:B:348:SER:N	2.79	0.45
1:B:363:LYS:HD3	4:B:523:HOH:O	2.15	0.45
1:B:76:VAL:HG13	1:B:81:ALA:HB3	1.98	0.45
1:B:370:TYR:CE2	1:B:432:ARG:HD3	2.52	0.45
1:B:117:TRP:CD1	1:B:163:ILE:HG13	2.51	0.45
1:B:358:ALA:HB3	1:B:361:GLY:C	2.42	0.45
1:B:385:GLU:HG2	1:B:389:HIS:HD2	1.82	0.45
1:B:25:THR:HA	1:B:52:ILE:O	2.16	0.44
1:B:279:MET:HE3	1:B:295:PHE:HB3	1.98	0.44
1:B:29:ALA:O	1:B:32:TRP:HB3	2.18	0.44
1:B:176:THR:CG2	1:B:178:HIS:H	2.29	0.44
1:B:126:GLU:OE2	1:B:226:GLY:HA3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:GLU:HG3	4:B:591:HOH:O	2.17	0.44
1:B:114:ILE:HG13	1:B:115:MET:N	2.31	0.44
1:B:245:THR:HG22	1:B:246:VAL:HG23	2.00	0.44
1:B:152:ALA:C	1:B:154:ALA:N	2.76	0.44
1:B:345:ILE:HD12	1:B:345:ILE:N	2.33	0.44
1:B:25:THR:CG2	1:B:53:THR:HG22	2.47	0.44
1:B:57:VAL:O	1:B:61:PHE:HD1	2.01	0.43
1:B:186:GLU:O	1:B:186:GLU:HG3	2.16	0.43
1:B:345:ILE:CG2	1:B:346:HIS:H	2.30	0.43
1:B:376:ALA:HA	4:B:512:HOH:O	2.17	0.43
1:B:255:MET:HA	1:B:258:THR:HG22	2.00	0.43
1:B:65:LEU:CD1	1:B:86:PRO:HA	2.48	0.43
1:B:175:PRO:HB2	1:B:247:LEU:HD23	2.00	0.43
1:B:316:SER:C	1:B:328:THR:HG23	2.42	0.43
1:B:320:GLN:HE21	1:B:320:GLN:HB2	1.65	0.43
1:B:16:LEU:HD12	1:B:25:THR:O	2.17	0.43
1:B:74:ARG:HH21	1:B:156:GLU:HB2	1.83	0.43
1:B:176:THR:CG2	1:B:177:THR:N	2.81	0.43
1:B:25:THR:HG22	1:B:53:THR:HA	2.00	0.43
1:B:95:ARG:CG	1:B:173:ILE:HD12	2.46	0.43
1:B:177:THR:HG22	1:B:247:LEU:HD21	2.00	0.43
1:B:194:LEU:N	1:B:194:LEU:HD13	2.34	0.43
1:B:240:TRP:CE2	1:B:247:LEU:HD13	2.54	0.43
1:B:346:HIS:O	1:B:347:ILE:C	2.62	0.43
1:B:356:LEU:HD11	1:B:374:PHE:CE2	2.54	0.43
1:B:132:TYR:HB2	1:B:182:LEU:HD13	2.01	0.42
1:B:174:THR:HA	1:B:175:PRO:HD3	1.86	0.42
1:B:151:VAL:HG13	1:B:152:ALA:N	2.34	0.42
1:B:276:ALA:HA	1:B:381:LEU:CD2	2.48	0.42
1:B:279:MET:CE	1:B:295:PHE:HB3	2.49	0.42
1:B:367:ASP:HB2	4:B:520:HOH:O	2.19	0.42
1:B:416:GLU:N	1:B:416:GLU:OE2	2.52	0.42
1:B:402:ALA:O	1:B:406:VAL:HG23	2.19	0.42
1:B:279:MET:SD	1:B:279:MET:HB3	2.59	0.42
1:B:311:HIS:CE1	1:B:313:LYS:HB2	2.55	0.42
1:B:157:LYS:O	1:B:157:LYS:HD2	2.20	0.41
1:B:407:LYS:C	1:B:409:SER:N	2.77	0.41
1:B:419:VAL:O	1:B:423:GLU:HB2	2.20	0.41
1:B:61:PHE:O	1:B:64:LEU:HB2	2.20	0.41
1:B:283:MET:HE2	1:B:288:ALA:HB1	2.00	0.41
1:B:5:VAL:HB	1:B:7:TRP:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:ASN:OD1	1:B:263:SER:HB3	2.20	0.41
1:B:176:THR:HG22	1:B:178:HIS:H	1.85	0.41
1:B:77:LYS:C	1:B:79:TYR:H	2.29	0.41
1:B:194:LEU:CD1	1:B:194:LEU:N	2.83	0.41
1:B:317:VAL:CA	1:B:328:THR:HG23	2.51	0.41
1:B:140:MET:HE1	1:B:190:PRO:HA	2.01	0.41
1:B:170:PRO:HB2	1:B:174:THR:O	2.21	0.41
1:B:117:TRP:HB3	1:B:161:LEU:HD23	2.03	0.41
1:B:339:VAL:HG12	1:B:339:VAL:O	2.21	0.41
1:B:74:ARG:O	1:B:77:LYS:HB2	2.21	0.41
1:B:42:SER:HB2	1:B:60:PHE:CE1	2.55	0.41
1:B:158:THR:HG22	1:B:180:ILE:HB	2.02	0.40
1:B:150:SEP:O3P	1:B:266:LYS:NZ	2.53	0.40
1:B:132:TYR:HD1	1:B:182:LEU:HD11	1.86	0.40
1:B:157:LYS:H	1:B:157:LYS:HG3	1.48	0.40
1:B:175:PRO:HG3	1:B:249:ARG:CZ	2.51	0.40
1:B:42:SER:HB2	1:B:60:PHE:CD1	2.57	0.40
1:B:95:ARG:HD3	1:B:172:GLY:O	2.22	0.40
1:B:132:TYR:O	1:B:133:LYS:HB3	2.20	0.40
1:B:251:PRO:HB3	4:B:583:HOH:O	2.22	0.40
1:B:314:ASN:ND2	4:B:568:HOH:O	2.55	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	B	412/440 (94%)	382 (93%)	24 (6%)	6 (2%)	8 8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	PRO
1	B	115	MET
1	B	80	HIS
1	B	436	GLU
1	B	9	ASN
1	B	384	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	B	356/373 (95%)	331 (93%)	25 (7%)	14	19

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	24	HIS
1	B	37	TYR
1	B	121	THR
1	B	122	GLU
1	B	124	ARG
1	B	128	VAL
1	B	140	MET
1	B	146	ASP
1	B	148	ARG
1	B	149	ILE
1	B	155	GLN
1	B	157	LYS
1	B	168	CYS
1	B	176	THR
1	B	194	LEU
1	B	228	VAL
1	B	283	MET
1	B	314	ASN
1	B	326	ARG
1	B	327	LEU
1	B	347	ILE

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Mol	Chain	Res	Type
1	B	348	SER
1	B	357	ASN
1	B	360	LYS
1	B	393	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	10	ASN
1	B	24	HIS
1	B	59	ASN
1	B	85	GLN
1	B	165	ASN
1	B	191	ASN
1	B	198	GLN
1	B	320	GLN
1	B	346	HIS
1	B	357	ASN
1	B	387	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	SEP	B	150	1	8,9,10	1.13	0	7,12,14	1.57	1 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SEP	B	150	1	-	1/6/8/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	SEP	O3P-P-O1P	3.15	123.12	110.83

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	150	SEP	CB-OG-P-O1P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	150	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	ADE	B	500	-	11,11,11	2.07	4 (36%)	15,15,15	1.01	1 (6%)
3	SO4	B	833	-	4,4,4	0.56	0	6,6,6	0.17	0

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	ADE	B	500	-	-	-	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	ADE	C4-N3	4.69	1.42	1.35
2	B	500	ADE	C8-N9	2.56	1.41	1.35
2	B	500	ADE	C2-N3	2.53	1.38	1.33
2	B	500	ADE	C2-N1	2.01	1.37	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	ADE	N3-C2-N1	-2.37	125.00	128.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	B	420/440 (95%)	1.38	121 (28%) 1 1	17, 43, 76, 84	1 (0%)

All (121) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	87	PHE	6.3
1	B	37	TYR	5.9
1	B	17	THR	5.5
1	B	100	ALA	5.4
1	B	115	MET	5.3
1	B	2	PRO	5.3
1	B	76	VAL	5.2
1	B	24	HIS	4.7
1	B	4	LEU	4.5
1	B	175	PRO	4.3
1	B	437	TYR	4.2
1	B	50	GLY	4.1
1	B	113	PRO	4.1
1	B	99	GLY	4.0
1	B	86	PRO	4.0
1	B	20	ALA	4.0
1	B	411	PRO	3.9
1	B	132	TYR	3.9
1	B	90	LEU	3.9
1	B	25	THR	3.8
1	B	81	ALA	3.8
1	B	415	PRO	3.8
1	B	410	LEU	3.8
1	B	23	ALA	3.7
1	B	7	TRP	3.6
1	B	9	ASN	3.6
1	B	93	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	139	GLY	3.6
1	B	401	ALA	3.5
1	B	103	LEU	3.5
1	B	431	GLY	3.4
1	B	435	ARG	3.4
1	B	114	ILE	3.3
1	B	105	PRO	3.3
1	B	18	LYS	3.3
1	B	153	GLY	3.3
1	B	57	VAL	3.3
1	B	19	LEU	3.3
1	B	94	GLY	3.2
1	B	404	ASP	3.2
1	B	91	SER	3.2
1	B	399	ILE	3.1
1	B	102	THR	3.1
1	B	22	GLY	3.1
1	B	11	GLN	3.1
1	B	157	LYS	3.1
1	B	398	MET	3.1
1	B	80	HIS	3.0
1	B	245	THR	3.0
1	B	155	GLN	3.0
1	B	345	ILE	3.0
1	B	164	GLY	3.0
1	B	14	GLY	2.9
1	B	400	PRO	2.9
1	B	424	SER	2.9
1	B	16	LEU	2.8
1	B	156	GLU	2.8
1	B	158	THR	2.8
1	B	88	ASP	2.8
1	B	141	ILE	2.8
1	B	133	LYS	2.8
1	B	6	THR	2.7
1	B	97	SER	2.7
1	B	8	MET	2.7
1	B	232	ALA	2.7
1	B	21	ASN	2.7
1	B	54	SER	2.7
1	B	359	SER	2.7
1	B	358	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	85	GLN	2.6
1	B	409	SER	2.6
1	B	15	GLU	2.6
1	B	436	GLU	2.6
1	B	402	ALA	2.6
1	B	58	PHE	2.6
1	B	5	VAL	2.6
1	B	131	ALA	2.5
1	B	101	VAL	2.5
1	B	172	GLY	2.5
1	B	419	VAL	2.5
1	B	140	MET	2.5
1	B	95	ARG	2.4
1	B	240	TRP	2.4
1	B	361	GLY	2.4
1	B	427	LEU	2.4
1	B	84	ARG	2.4
1	B	321	ALA	2.4
1	B	130	THR	2.4
1	B	177	THR	2.4
1	B	83	SER	2.4
1	B	248	LEU	2.4
1	B	32	TRP	2.4
1	B	365	ALA	2.3
1	B	161	LEU	2.3
1	B	10	ASN	2.3
1	B	116	ALA	2.3
1	B	53	THR	2.3
1	B	429	LEU	2.2
1	B	123	ALA	2.2
1	B	154	ALA	2.2
1	B	193	THR	2.2
1	B	433	LEU	2.2
1	B	96	ASP	2.2
1	B	190	PRO	2.2
1	B	168	CYS	2.2
1	B	38	ALA	2.2
1	B	380	VAL	2.2
1	B	62	ASP	2.1
1	B	170	PRO	2.1
1	B	421	ALA	2.1
1	B	1	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	357	ASN	2.1
1	B	173	ILE	2.1
1	B	414	PHE	2.1
1	B	412	THR	2.1
1	B	430	HIS	2.1
1	B	416	GLU	2.0
1	B	146	ASP	2.0
1	B	413	ASP	2.0
1	B	52	ILE	2.0
1	B	64	LEU	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
1	SEP	B	150	10/11	0.94	0.10	31,35,37,38	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
3	SO4	B	833	5/5	0.67	0.23	67,67,69,69	0
2	ADE	B	500	10/10	0.89	0.10	31,36,45,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.