



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:10 PM GMT

PDB ID : 3WD9
Title : Crystal structure of phosphodiesterase 4B in complex with compound 10f
Authors : Takahashi, M.; Hanzawa, H.
Deposited on : 2013-06-11
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

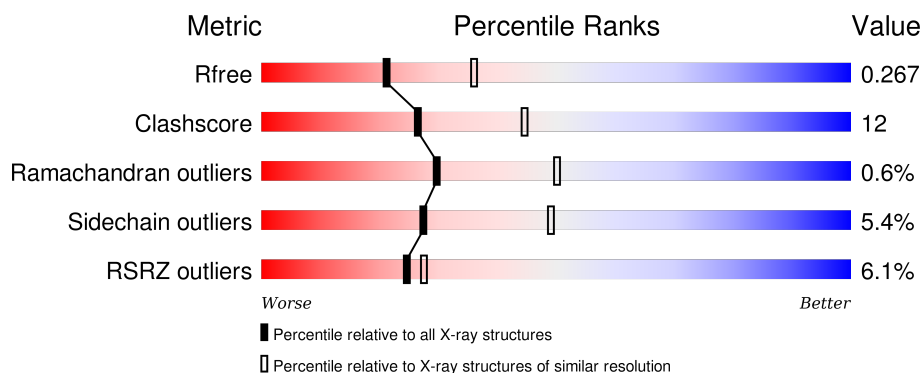
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


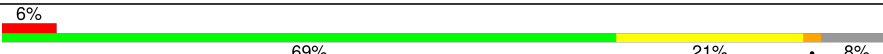
The reported resolution of this entry is 2.50 Å.

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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	377	
1	B	377	

2 Entry composition [i](#)

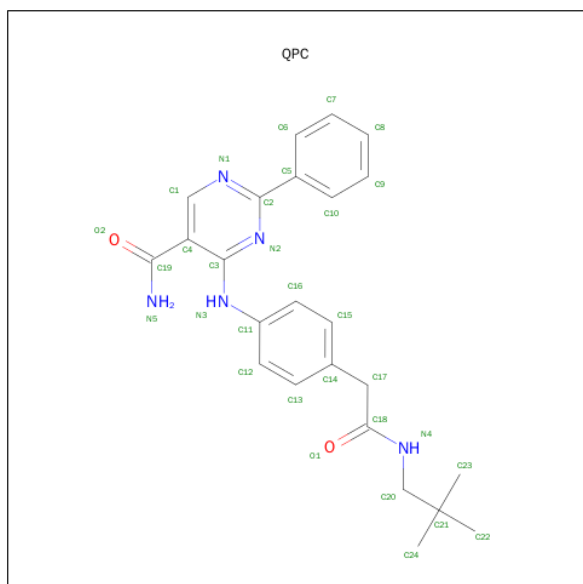
There are 5 unique types of molecules in this entry. The entry contains 5991 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-specific 3',5'-cyclic phosphodiesterase 4B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2792	1763	471	537	21			
1	B	345	Total	C	N	O	S	0	0	0
			2792	1763	471	537	21			

- Molecule 2 is 4-[(4-{2-[(2,2-DIMETHYLPROPYL)AMINO]-2-OXOETHYL}PHENYL)AMINO]-2-PHENYLPYRIMIDINE-5-CARBOXAMIDE (three-letter code: QPC) (formula: C₂₄H₂₇N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			31	24	5	2		
2	B	1	Total	C	N	O	0	0
			31	24	5	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

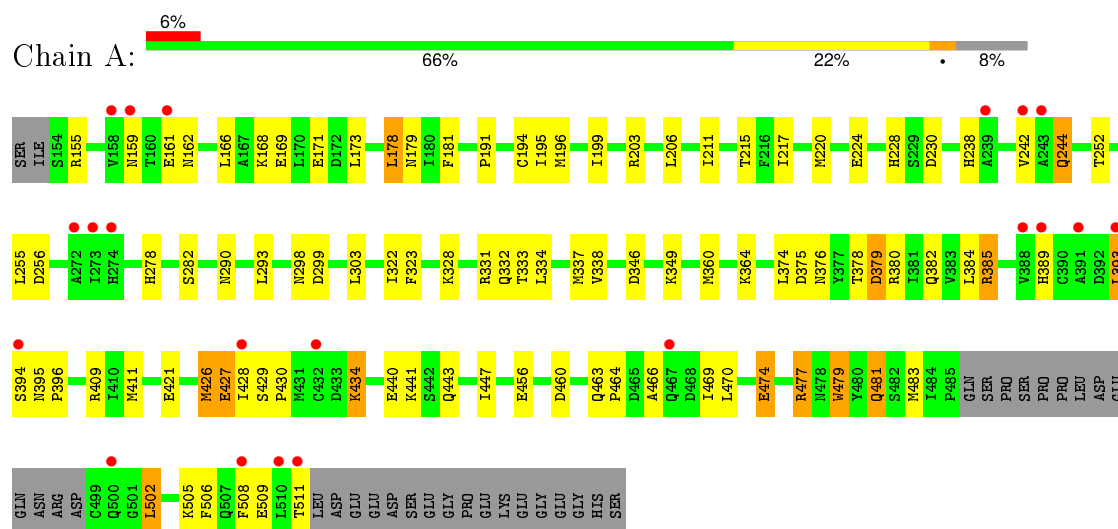
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	159	Total 159	O 159	0	0
5	B	182	Total 182	O 182	0	0

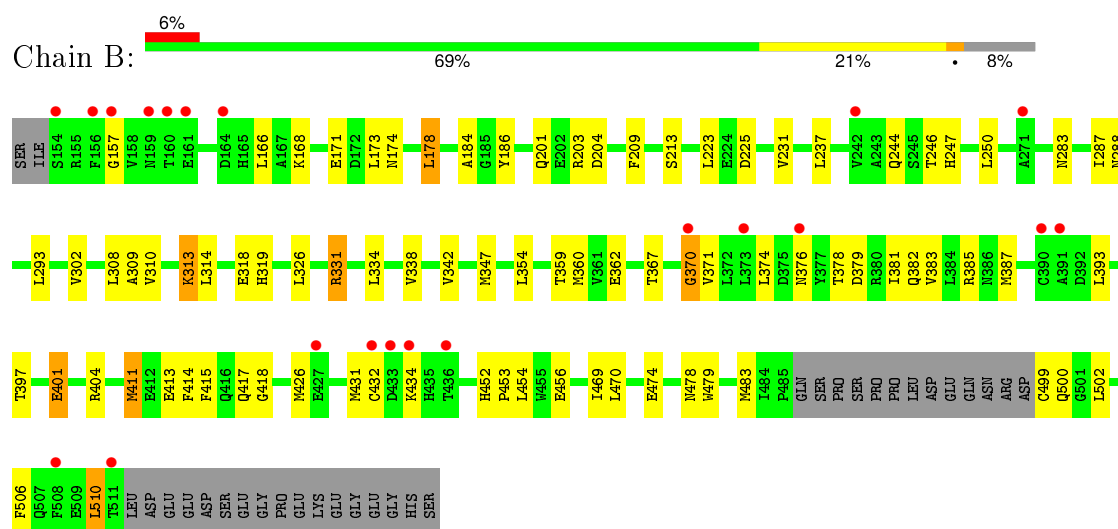
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4B



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	81.31Å 157.38Å 58.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.81 – 2.50 32.57 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.5 (19.81-2.50) 90.6 (32.57-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.174 , 0.267 0.174 , 0.267	Depositor DCC
R_{free} test set	2400 reflections (11.11%)	DCC
Wilson B-factor (Å ²)	40.5	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 24067 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5991	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: QPC, ZN, 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	1/2849 (0.0%)	0.69	0/3859
1	B	0.62	0/2849	0.67	1/3859 (0.0%)
All	All	0.63	1/5698 (0.0%)	0.68	1/7718 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	479	TRP	CD2-CE2	5.09	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	204	ASP	CB-CG-OD1	5.42	123.18	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2792	0	2710	69	0
1	B	2792	0	2710	65	0
2	A	31	0	27	2	0
2	B	31	0	27	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	159	0	0	6	0
5	B	182	0	0	5	0
All	All	5991	0	5474	136	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 12.

All (136) 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:331:ARG:HB3	1:B:331:ARG:HH11	1.37	0.87
1:B:411:MET:HE1	1:B:502:LEU:HD13	1.58	0.86
1:B:411:MET:CE	1:B:502:LEU:HD13	2.06	0.85
1:B:510:LEU:HD23	5:B:1098:HOH:O	1.77	0.84
1:B:203:ARG:HE	1:B:247:HIS:HD2	1.30	0.78
1:B:367:THR:HG23	1:B:370:GLY:H	1.52	0.75
1:A:411:MET:HE1	1:A:502:LEU:HD11	1.68	0.72
1:B:318:GLU:HG3	1:B:319:HIS:CE1	2.26	0.70
1:B:367:THR:HG22	1:B:371:VAL:H	1.55	0.70
1:B:367:THR:HG22	1:B:371:VAL:N	2.07	0.69
1:B:367:THR:CG2	1:B:370:GLY:H	2.06	0.68
1:A:376:ASN:ND2	1:A:378:THR:HB	2.09	0.68
1:B:173:LEU:HA	1:B:178:LEU:HD13	1.75	0.67
1:A:421:GLU:OE2	5:A:1039:HOH:O	2.11	0.67
1:B:173:LEU:HA	1:B:178:LEU:CD1	2.28	0.64
1:A:427:GLU:HG3	1:A:428:ILE:N	2.13	0.63
1:B:283:ASN:O	1:B:287:ILE:HG13	1.99	0.63
1:B:381:ILE:O	1:B:385:ARG:HG3	1.98	0.62
1:B:411:MET:CE	1:B:502:LEU:CD1	2.78	0.62
1:A:203:ARG:NH1	5:A:1019:HOH:O	2.28	0.61
1:B:374:LEU:HD11	1:B:383:VAL:HG21	1.81	0.61
1:A:427:GLU:HG3	1:A:429:SER:H	1.67	0.60
1:A:505:LYS:O	1:A:509:GLU:HG2	2.02	0.59
1:A:427:GLU:HG3	1:A:429:SER:N	2.18	0.58
1:A:434:LYS:H	1:A:434:LYS:HD3	1.67	0.58
1:B:318:GLU:HG3	1:B:319:HIS:ND1	2.19	0.58
1:A:168:LYS:HD3	1:A:171:GLU:OE2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:GLU:OE1	1:A:429:SER:HB3	2.05	0.57
1:B:397:THR:HB	1:B:469:ILE:HG23	1.87	0.57
1:A:364:LYS:HA	5:A:1072:HOH:O	2.02	0.57
1:A:346:ASP:HB3	1:A:349:LYS:HD2	1.85	0.56
1:B:302:VAL:HG23	5:B:1048:HOH:O	2.05	0.56
1:B:418:GLY:HA3	1:B:431:MET:CE	2.35	0.56
1:B:382:GLN:HG3	1:B:385:ARG:HH11	1.71	0.56
1:B:203:ARG:HE	1:B:247:HIS:CD2	2.18	0.55
1:B:411:MET:HE1	1:B:502:LEU:CD1	2.33	0.55
1:B:246:THR:HG23	1:B:387:MET:CE	2.36	0.55
1:A:426:MET:O	1:A:427:GLU:C	2.45	0.55
1:A:508:PHE:O	1:A:511:THR:HG23	2.07	0.55
1:A:389:HIS:CE1	1:A:393:LEU:HD22	2.43	0.54
1:B:288:ASN:HD22	1:B:426:MET:HE1	1.73	0.54
1:A:333:THR:O	1:A:337:MET:HG3	2.07	0.53
1:A:282:SER:HB3	2:A:901:QPC:H17	1.90	0.53
1:B:331:ARG:HH11	1:B:331:ARG:CB	2.16	0.53
1:B:347:MET:HE1	1:B:393:LEU:HD11	1.91	0.53
2:B:901:QPC:H11	2:B:901:QPC:N2	2.24	0.53
1:A:173:LEU:HA	1:A:178:LEU:HD13	1.90	0.52
1:A:334:LEU:O	1:A:338:VAL:HG23	2.09	0.52
1:A:506:PHE:HB2	2:A:901:QPC:H5	1.91	0.51
1:A:427:GLU:CG	1:A:428:ILE:N	2.73	0.51
1:B:414:PHE:CE2	2:B:901:QPC:H17	2.45	0.51
1:A:443:GLN:O	1:A:447:ILE:HG13	2.12	0.50
1:A:206:LEU:HB2	5:A:1014:HOH:O	2.11	0.50
1:A:178:LEU:HD22	1:A:244:GLN:CD	2.33	0.49
1:B:418:GLY:HA3	1:B:431:MET:HE1	1.95	0.49
1:B:474:GLU:OE2	1:B:478:ASN:ND2	2.46	0.49
1:B:417:GLN:OE1	2:B:901:QPC:H27	2.13	0.48
1:A:155:ARG:HG2	1:A:217:ILE:HG21	1.95	0.48
1:B:246:THR:HG23	1:B:387:MET:HE3	1.97	0.47
1:A:470:LEU:O	1:A:474:GLU:HG2	2.14	0.47
1:B:415:PHE:HA	1:B:431:MET:HE1	1.96	0.47
1:A:376:ASN:HD22	1:A:378:THR:HB	1.75	0.47
1:A:196:MET:HG3	1:A:220:MET:CE	2.45	0.47
1:A:323:PHE:O	1:A:331:ARG:HD3	2.15	0.47
1:A:393:LEU:HD12	1:A:393:LEU:HA	1.67	0.47
1:B:246:THR:HG23	1:B:387:MET:HE2	1.97	0.46
1:B:479:TRP:CH2	1:B:483:MET:HG3	2.49	0.46
1:B:231:VAL:CG1	1:B:413:GLU:OE1	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:ASP:O	1:A:256:ASP:OD2	2.34	0.46
1:A:252:THR:HG21	5:A:1107:HOH:O	2.14	0.46
1:B:168:LYS:HA	1:B:171:GLU:HG3	1.98	0.46
1:A:199:ILE:O	1:A:203:ARG:HG3	2.15	0.46
1:A:211:ILE:HG12	1:A:322:ILE:O	2.15	0.46
1:A:440:GLU:HG2	1:A:483:MET:HB2	1.98	0.45
1:A:426:MET:O	1:A:427:GLU:O	2.33	0.45
1:A:429:SER:HA	1:A:430:PRO:HD3	1.78	0.45
1:B:354:LEU:HD13	1:B:454:LEU:HA	1.98	0.45
1:A:460:ASP:O	1:A:463:GLN:HG3	2.16	0.45
1:A:382:GLN:HA	1:A:385:ARG:HD2	1.98	0.45
1:B:331:ARG:NE	5:B:1014:HOH:O	2.19	0.44
1:B:415:PHE:HD1	1:B:431:MET:HE3	1.82	0.44
1:A:477:ARG:HH11	1:A:481:GLN:HG3	1.82	0.44
1:A:179:ASN:OD1	1:A:181:PHE:HB2	2.17	0.44
1:B:385:ARG:NH1	5:B:1056:HOH:O	2.42	0.44
1:B:334:LEU:O	1:B:338:VAL:HG23	2.17	0.44
1:A:411:MET:HE1	1:A:502:LEU:CD1	2.43	0.44
1:B:502:LEU:H	1:B:502:LEU:HG	1.65	0.44
1:A:228:HIS:HB3	1:A:230:ASP:OD1	2.17	0.44
1:B:174:ASN:OD1	1:B:247:HIS:HE1	2.01	0.43
1:A:303:LEU:HA	1:A:303:LEU:HD23	1.75	0.43
1:B:401:GLU:H	1:B:401:GLU:HG3	1.34	0.43
1:B:223:LEU:HA	1:B:314:LEU:CD1	2.49	0.43
1:B:456:GLU:HA	1:B:470:LEU:HD11	1.99	0.43
1:A:293:LEU:HA	1:A:293:LEU:HD23	1.81	0.43
1:A:328:LYS:O	1:A:332:GLN:HG2	2.19	0.43
1:B:411:MET:HE2	1:B:502:LEU:CD1	2.48	0.43
1:A:427:GLU:HG3	1:A:428:ILE:H	1.81	0.43
1:B:250:LEU:HG	1:B:387:MET:HE3	2.00	0.43
1:A:244:GLN:HE21	1:A:244:GLN:HB3	1.61	0.43
1:A:395:ASN:HB2	1:A:396:PRO:HD3	2.01	0.43
1:B:231:VAL:HG12	1:B:413:GLU:OE1	2.19	0.42
1:B:452:HIS:HB3	1:B:453:PRO:HD3	2.01	0.42
1:B:331:ARG:NH2	5:B:1014:HOH:O	2.41	0.42
1:B:506:PHE:O	1:B:510:LEU:HB2	2.19	0.42
1:B:415:PHE:HA	1:B:431:MET:CE	2.50	0.42
2:B:901:QPC:C16	2:B:901:QPC:N2	2.83	0.42
1:B:308:LEU:HD21	1:B:342:VAL:HB	2.00	0.42
1:B:184:ALA:HB2	1:B:237:LEU:HD13	2.00	0.42
1:B:168:LYS:HG3	1:B:171:GLU:OE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:VAL:HA	1:B:313:LYS:HD3	2.02	0.42
1:A:159:ASN:C	1:A:161:GLU:N	2.72	0.42
1:A:166:LEU:HD21	1:A:194:CYS:SG	2.59	0.42
1:A:215:THR:HG22	1:A:322:ILE:HG22	2.02	0.41
1:A:466:ALA:O	1:A:469:ILE:HB	2.20	0.41
1:B:359:THR:HA	1:B:362:GLU:OE1	2.21	0.41
1:A:228:HIS:HD2	1:A:278:HIS:O	2.02	0.41
1:B:376:ASN:HD21	1:B:378:THR:HB	1.85	0.41
1:A:374:LEU:HD13	1:A:380:ARG:HA	2.02	0.41
1:A:298:ASN:O	1:A:299:ASP:HB2	2.21	0.41
1:A:409:ARG:NH1	5:A:1057:HOH:O	2.53	0.41
1:B:360:MET:HE1	1:B:382:GLN:HB3	2.01	0.41
1:A:238:HIS:O	1:A:242:VAL:HG23	2.20	0.41
1:B:209:PHE:CD1	1:B:326:LEU:HD11	2.55	0.41
1:B:244:GLN:O	1:B:247:HIS:HB3	2.21	0.41
1:A:428:ILE:HD12	1:A:429:SER:N	2.36	0.41
1:A:255:LEU:HD23	1:A:255:LEU:HA	1.83	0.41
1:A:474:GLU:H	1:A:474:GLU:HG2	1.54	0.41
1:A:159:ASN:HB3	1:A:161:GLU:HG3	2.02	0.41
1:A:380:ARG:O	1:A:384:LEU:HG	2.21	0.41
1:A:463:GLN:HA	1:A:464:PRO:HA	1.88	0.41
1:A:394:SER:O	1:A:395:ASN:C	2.59	0.41
1:A:191:PRO:HD2	1:A:224:GLU:OE2	2.21	0.41
1:A:374:LEU:HD22	1:A:379:ASP:HB3	2.02	0.40
1:B:166:LEU:HB2	1:B:186:TYR:HB3	2.03	0.40
1:A:195:ILE:HA	1:A:195:ILE:HD12	1.92	0.40
1:B:309:ALA:O	1:B:313:LYS:HB3	2.22	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	341/377 (90%)	330 (97%)	10 (3%)	1 (0%)	46	68
1	B	341/377 (90%)	325 (95%)	13 (4%)	3 (1%)	21	37
All	All	682/754 (90%)	655 (96%)	23 (3%)	4 (1%)	30	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	427	GLU
1	B	370	GLY
1	B	510	LEU
1	B	157	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	313/342 (92%)	294 (94%)	19 (6%)	23	42
1	B	313/342 (92%)	298 (95%)	15 (5%)	31	55
All	All	626/684 (92%)	592 (95%)	34 (5%)	27	49

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	169	GLU
1	A	178	LEU
1	A	244	GLN
1	A	290	ASN
1	A	360	MET
1	A	375	ASP
1	A	379	ASP
1	A	385	ARG
1	A	393	LEU
1	A	426	MET
1	A	434	LYS

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Mol	Chain	Res	Type
1	A	441	LYS
1	A	456	GLU
1	A	474	GLU
1	A	477	ARG
1	A	479	TRP
1	A	481	GLN
1	A	502	LEU
1	B	178	LEU
1	B	201	GLN
1	B	213	SER
1	B	225	ASP
1	B	293	LEU
1	B	313	LYS
1	B	331	ARG
1	B	379	ASP
1	B	401	GLU
1	B	404	ARG
1	B	411	MET
1	B	432	CYS
1	B	434	LYS
1	B	499	CYS
1	B	500	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	ASN
1	A	228	HIS
1	A	244	GLN
1	A	284	GLN
1	A	288	ASN
1	A	376	ASN
1	B	159	ASN
1	B	182	ASN
1	B	201	GLN
1	B	226	HIS
1	B	228	HIS
1	B	244	GLN
1	B	247	HIS
1	B	284	GLN
1	B	288	ASN
1	B	305	ASN

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Mol	Chain	Res	Type
1	B	376	ASN
1	B	386	ASN
1	B	481	GLN
1	B	500	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
2	QPC	A	901	-	33,33,33	0.94	2 (6%)	44,46,46	1.14	2 (4%)
2	QPC	B	901	-	33,33,33	1.09	4 (12%)	44,46,46	1.08	4 (9%)

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	QPC	A	901	-	-	0/22/22/22	0/3/3/3
2	QPC	B	901	-	-	0/22/22/22	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	QPC	C4-C3	-2.80	1.38	1.42
2	A	901	QPC	C18-N4	2.19	1.38	1.33
2	B	901	QPC	C19-N5	2.38	1.37	1.33
2	B	901	QPC	C18-N4	2.75	1.40	1.33
2	A	901	QPC	C3-N3	3.11	1.42	1.36
2	B	901	QPC	C3-N3	3.18	1.42	1.36

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	QPC	C4-C3-N2	-2.42	119.25	122.65
2	A	901	QPC	C11-N3-C3	-2.31	122.90	128.96
2	A	901	QPC	C1-C4-C3	2.11	116.62	114.51
2	B	901	QPC	C15-C16-C11	2.12	122.66	120.28
2	B	901	QPC	C1-C4-C3	2.34	116.84	114.51
2	B	901	QPC	C21-C20-N4	2.69	118.13	113.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	QPC	2	0
2	B	901	QPC	4	0

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	345/377 (91%)	0.05	21 (6%) 25 27	33, 48, 85, 123	0
1	B	345/377 (91%)	-0.03	21 (6%) 25 27	34, 46, 90, 122	0
All	All	690/754 (91%)	0.01	42 (6%) 25 27	33, 47, 89, 123	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	THR	6.4
1	A	511	THR	6.3
1	B	157	GLY	5.1
1	B	508	PHE	4.3
1	B	161	GLU	3.9
1	A	508	PHE	3.6
1	A	242	VAL	3.5
1	B	511	THR	3.2
1	B	159	ASN	3.1
1	A	428	ILE	3.1
1	A	158	VAL	3.1
1	B	436	THR	3.0
1	A	159	ASN	2.9
1	B	391	ALA	2.9
1	B	433	ASP	2.8
1	A	273	ILE	2.8
1	A	239	ALA	2.7
1	B	390	CYS	2.7
1	B	242	VAL	2.7
1	B	154	SER	2.6
1	A	161	GLU	2.6
1	A	388	VAL	2.5
1	A	274	HIS	2.5
1	A	391	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	164	ASP	2.4
1	A	510	LEU	2.4
1	A	389	HIS	2.4
1	A	393	LEU	2.4
1	B	370	GLY	2.3
1	A	432	CYS	2.3
1	A	394	SER	2.3
1	B	376	ASN	2.3
1	B	271	ALA	2.2
1	B	432	CYS	2.2
1	B	156	PHE	2.2
1	B	373	LEU	2.1
1	B	434	LYS	2.1
1	A	243	ALA	2.1
1	A	500	GLN	2.1
1	A	272	ALA	2.0
1	B	427	GLU	2.0
1	A	467	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
2	QPC	B	901	31/31	0.95	0.18	0.53	38,49,64,70	0
4	CA	A	903	1/1	0.96	0.23	0.40	52,52,52,52	0
2	QPC	A	901	31/31	0.95	0.18	0.26	43,52,71,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	B	903	1/1	1.00	0.18	0.10	45,45,45,45	0
3	ZN	B	902	1/1	1.00	0.13	-1.38	47,47,47,47	0
3	ZN	A	902	1/1	1.00	0.13	-1.95	51,51,51,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.