



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:55 PM UTC

PDB ID : 6EIV / pdb_00006eiv
Title : DYRK1A in complex with JWD-065
Authors : Rothweiler, U.
Deposited on : 2017-09-19
Resolution : 2.68 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

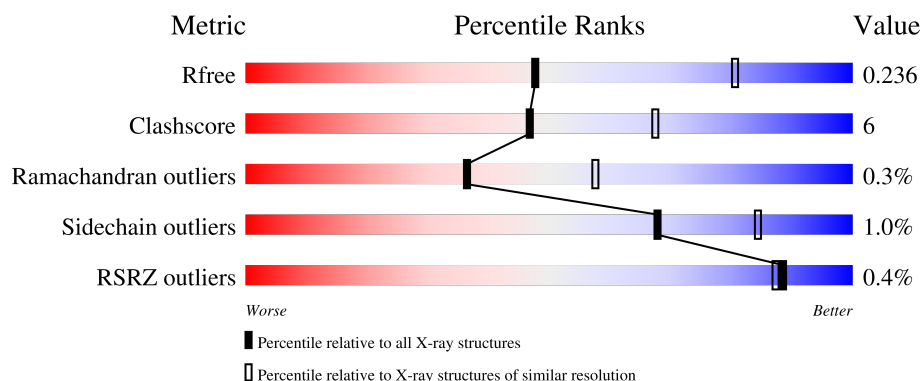
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

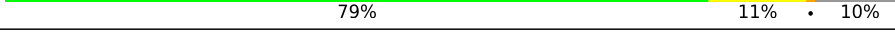
The reported resolution of this entry is 2.68 Å.

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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	368	 81% 12% 7%
1	B	368	 75% 17% 8%
1	C	368	 73% 16% 10%
1	D	368	 79% 11% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11167 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.

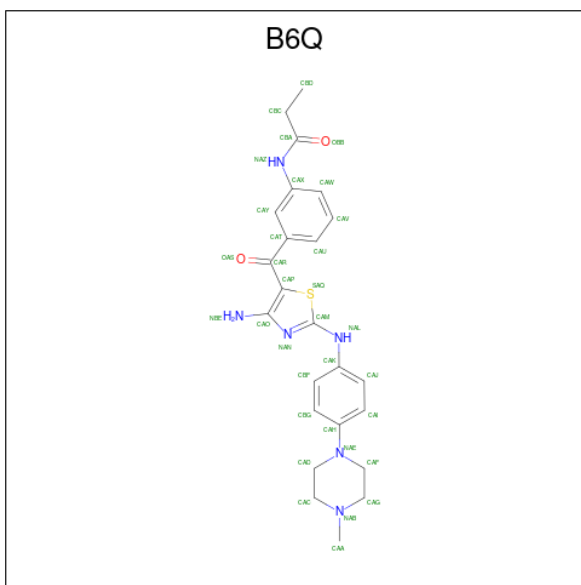
- Molecule 1 is a protein called Dual specificity tyrosine-phosphorylation-regulated kinase 1A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	P	S	0	0	0
			2794	1798	477	501	1	17			
1	B	340	Total	C	N	O	P	S	0	0	0
			2757	1774	470	495	1	17			
1	C	333	Total	C	N	O	P	S	0	0	0
			2649	1697	452	482	1	17			
1	D	333	Total	C	N	O	P	S	0	0	0
			2711	1752	455	486	1	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	GLY	-	expression tag	UNP Q13627
A	124	ALA	-	expression tag	UNP Q13627
A	125	SER	-	expression tag	UNP Q13627
B	123	GLY	-	expression tag	UNP Q13627
B	124	ALA	-	expression tag	UNP Q13627
B	125	SER	-	expression tag	UNP Q13627
C	123	GLY	-	expression tag	UNP Q13627
C	124	ALA	-	expression tag	UNP Q13627
C	125	SER	-	expression tag	UNP Q13627
D	123	GLY	-	expression tag	UNP Q13627
D	124	ALA	-	expression tag	UNP Q13627
D	125	SER	-	expression tag	UNP Q13627

- Molecule 2 is {N}-[3-[[4-azanyl-2-[[4-(4-methylpiperazin-1-yl)phenyl]amino]-1,3-thiazol-5-yl]carbonyl]phenyl]propanamide (CCD ID: B6Q) (formula: C₂₄H₂₈N₆O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 33	C 24	N 6	O 2	S 1	0	0
2	B	1	Total 33	C 24	N 6	O 2	S 1	0	0
2	C	1	Total 33	C 24	N 6	O 2	S 1	0	0
2	D	1	Total 33	C 24	N 6	O 2	S 1	0	0

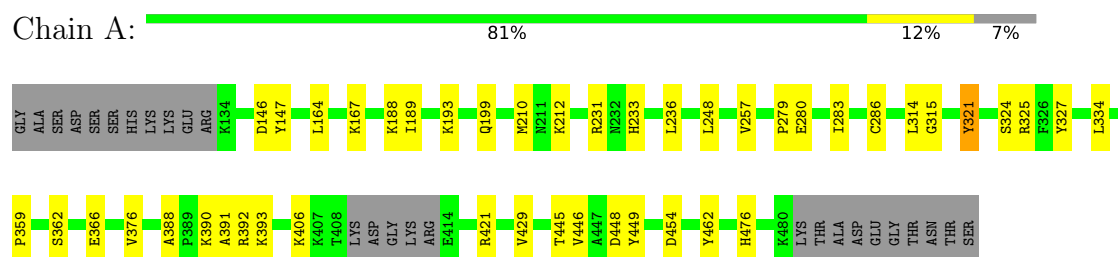
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	38	Total O 38 38	0	0
3	B	20	Total O 20 20	0	0
3	C	28	Total O 28 28	0	0
3	D	38	Total O 38 38	0	0

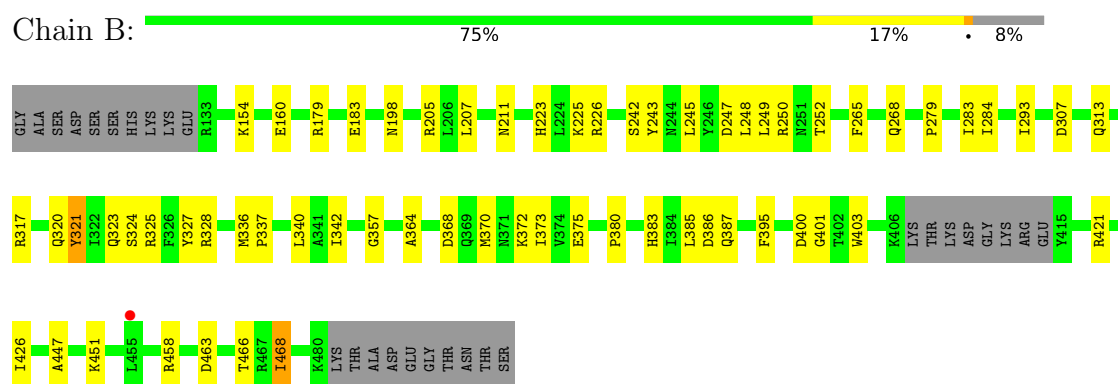
3 Residue-property plots

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.

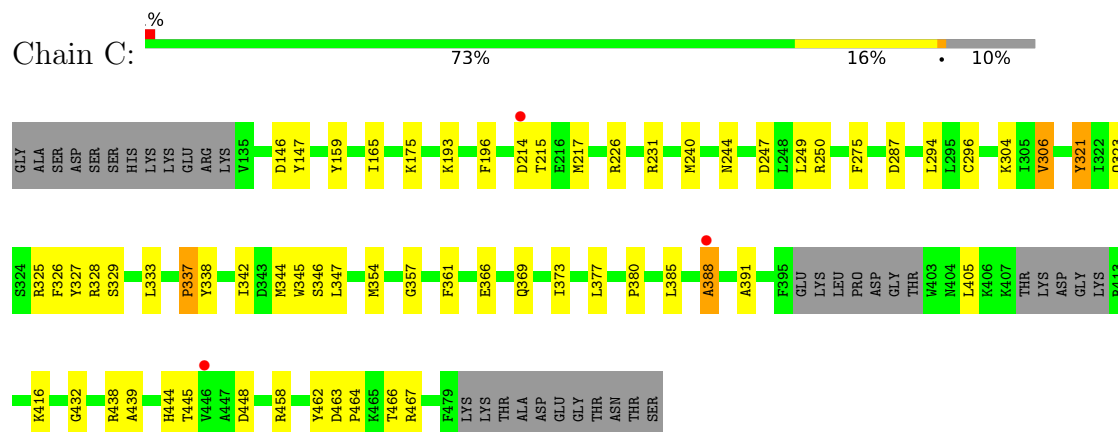
- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



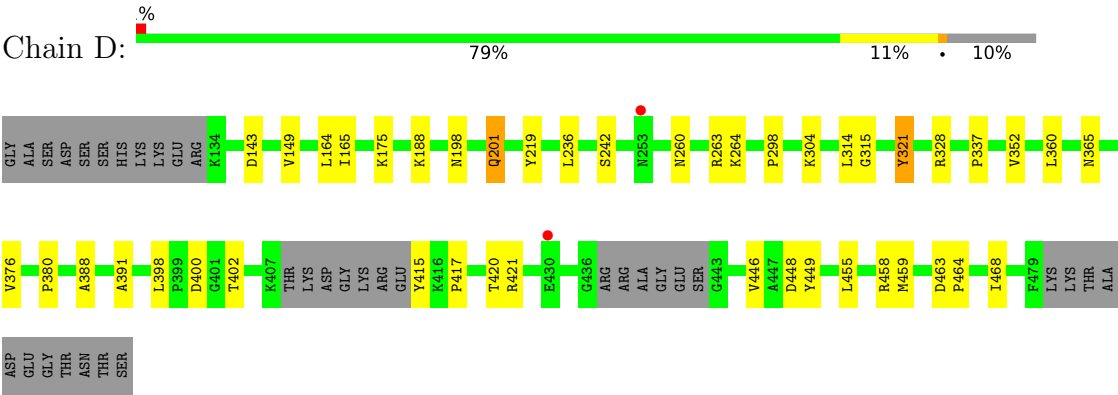
- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



● Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.76Å 88.54Å 231.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.62 – 2.68 48.62 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.1 (48.62-2.68) 98.9 (48.62-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.197 , 0.235 0.195 , 0.236	Depositor DCC
R_{free} test set	2639 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 37.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.368 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11167	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PTR, B6Q

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	0.26	0/2842	0.69	0/3833
1	B	0.27	0/2804	0.75	0/3785
1	C	0.28	0/2692	0.76	5/3643 (0.1%)
1	D	0.28	0/2758	0.72	1/3723 (0.0%)
All	All	0.27	0/11096	0.73	6/14984 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	416	LYS	CA-C-N	6.40	124.27	119.66
1	C	416	LYS	C-N-CA	6.40	124.27	119.66
1	C	388	ALA	CA-C-N	5.91	125.61	119.05
1	C	388	ALA	C-N-CA	5.91	125.61	119.05
1	C	377	LEU	N-CA-C	5.52	116.93	110.41
1	D	380	PRO	O-C-N	5.06	123.64	121.31

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	2794	0	2787	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2757	0	2734	39	0
1	C	2649	0	2551	37	0
1	D	2711	0	2693	26	0
2	A	33	0	0	0	0
2	B	33	0	0	0	0
2	C	33	0	0	2	0
2	D	33	0	0	0	0
3	A	38	0	0	1	0
3	B	20	0	0	1	0
3	C	28	0	0	0	0
3	D	38	0	0	2	0
All	All	11167	0	10765	125	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 6.

All (125) 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:C:338:TYR:HB2	1:C:342:ILE:HD11	1.59	0.85
1:A:279:PRO:HG3	1:D:337:PRO:HG3	1.63	0.79
1:C:323:GLN:NE2	1:C:346:SER:OG	2.21	0.73
1:B:183:GLU:OE1	1:B:225:LYS:NZ	2.25	0.69
1:C:249:LEU:HD22	1:C:357:GLY:HA2	1.76	0.68
1:A:321:PTR:O1P	1:A:325:ARG:NH2	2.27	0.67
1:D:415:TYR:N	3:D:601:HOH:O	2.29	0.66
1:D:398:LEU:HB2	1:D:402:THR:HG22	1.78	0.65
1:C:294:LEU:HD13	1:C:306:VAL:HG21	1.79	0.64
1:C:458:ARG:NH2	1:C:466:THR:O	2.31	0.64
1:C:388:ALA:HB3	1:C:391:ALA:HB2	1.80	0.63
1:D:352:VAL:HG11	1:D:360:LEU:HD13	1.82	0.62
1:B:226:ARG:NH2	3:B:601:HOH:O	2.32	0.62
1:A:388:ALA:HB3	1:A:391:ALA:HB2	1.81	0.62
1:B:207:LEU:O	1:B:211:ASN:ND2	2.25	0.62
1:C:380:PRO:HB2	1:C:385:LEU:HD11	1.80	0.62
1:C:247:ASP:OD1	1:C:250:ARG:NH2	2.34	0.61
1:C:463:ASP:O	1:C:467:ARG:N	2.33	0.61
1:C:296:CYS:SG	1:C:304:LYS:NZ	2.68	0.59
1:B:321:PTR:O1P	1:B:328:ARG:NH2	2.30	0.58
1:A:147:TYR:OH	1:A:231:ARG:NH1	2.37	0.58
1:B:324:SER:HB2	1:B:327:TYR:HD2	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:458:ARG:C	1:B:468:ILE:HD11	2.31	0.56
1:B:336:MET:HE2	1:B:337:PRO:HD2	1.87	0.56
1:B:386:ASP:OD1	1:B:403:TRP:NE1	2.27	0.56
1:B:247:ASP:OD1	1:B:250:ARG:NH2	2.36	0.56
1:C:244:ASN:HA	1:C:294:LEU:HA	1.88	0.56
1:B:380:PRO:HB2	1:B:385:LEU:HD13	1.89	0.55
1:B:325:ARG:NH1	1:B:364:ALA:O	2.41	0.53
1:A:210:MET:HE2	1:A:283:ILE:HD13	1.90	0.53
1:B:154:LYS:HE2	1:B:160:GLU:HG2	1.91	0.53
1:A:324:SER:HB2	1:A:327:TYR:HD2	1.73	0.52
1:C:321:PTR:O1P	1:C:328:ARG:NH2	2.32	0.52
1:B:400:ASP:OD1	1:B:401:GLY:N	2.43	0.52
1:A:454:ASP:OD1	1:A:476:HIS:NE2	2.36	0.52
1:D:321:PTR:O3P	1:D:328:ARG:NH2	2.32	0.52
1:C:215:THR:HG22	1:C:217:MET:H	1.74	0.52
1:C:159:TYR:OH	1:C:226:ARG:NH1	2.42	0.52
1:D:458:ARG:HB3	1:D:468:ILE:HB	1.90	0.51
1:C:326:PHE:HA	1:C:369:GLN:NE2	2.26	0.51
1:B:421:ARG:HG2	1:B:426:ILE:HD11	1.93	0.50
1:B:283:ILE:HG12	1:B:313:GLN:HG2	1.92	0.50
1:C:165:ILE:HD13	1:C:175:LYS:HB2	1.93	0.50
1:B:205:ARG:NH2	1:B:313:GLN:OE1	2.41	0.50
1:B:249:LEU:HD22	1:B:357:GLY:HA2	1.92	0.50
1:D:417:PRO:O	1:D:420:THR:OG1	2.26	0.50
1:B:284:ILE:HG12	1:B:340:LEU:HG	1.94	0.49
1:B:323:GLN:NE2	1:B:342:ILE:O	2.44	0.49
1:C:369:GLN:O	1:C:373:ILE:HG12	2.12	0.49
1:A:314:LEU:O	1:D:315:GLY:HA2	2.13	0.49
1:C:147:TYR:OH	1:C:231:ARG:NH1	2.46	0.49
1:C:438:ARG:HH11	1:C:444:HIS:CE1	2.31	0.49
1:D:388:ALA:HB3	1:D:391:ALA:HB2	1.95	0.49
1:C:462:TYR:O	1:C:464:PRO:HD3	2.14	0.48
1:A:334:LEU:HB3	1:A:388:ALA:HB1	1.95	0.48
1:C:249:LEU:HD21	1:C:354:MET:HA	1.96	0.48
1:C:361:PHE:O	1:C:369:GLN:NE2	2.47	0.48
1:A:188:LYS:HB3	1:A:236:LEU:HB2	1.95	0.48
1:D:352:VAL:HG21	1:D:360:LEU:HD22	1.96	0.48
1:A:445:THR:OG1	1:A:448:ASP:OD2	2.32	0.48
1:A:146:ASP:HA	1:A:233:HIS:HE1	1.78	0.47
1:D:421:ARG:NH1	3:D:604:HOH:O	2.45	0.47
1:B:223:HIS:CD2	1:B:225:LYS:HG2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:446:VAL:HA	1:D:449:TYR:HB2	1.95	0.47
1:B:317:ARG:NH2	1:B:320:GLN:OE1	2.48	0.47
1:B:447:ALA:O	1:B:451:LYS:HG2	2.15	0.47
1:D:198:ASN:HA	1:D:201:GLN:HG3	1.96	0.46
1:B:385:LEU:HD21	1:B:395:PHE:CD2	2.50	0.46
1:C:321:PTR:O3P	1:C:325:ARG:NH2	2.49	0.46
1:D:143:ASP:OD1	1:D:143:ASP:N	2.45	0.46
1:A:193:LYS:HE2	1:A:193:LYS:HB3	1.69	0.45
1:A:164:LEU:HD21	1:A:167:LYS:HE3	1.98	0.45
1:C:217:MET:HB3	1:C:275:PHE:HB2	1.96	0.45
1:C:342:ILE:O	1:C:345:TRP:HB3	2.16	0.45
1:A:390:LYS:NZ	3:A:604:HOH:O	2.43	0.45
1:B:279:PRO:HG3	1:C:337:PRO:HD3	1.98	0.45
1:D:219:TYR:O	1:D:304:LYS:NZ	2.32	0.45
1:A:199:GLN:OE1	1:B:198:ASN:ND2	2.50	0.45
1:B:160:GLU:OE2	1:B:179:ARG:NE	2.40	0.45
1:D:376:VAL:HA	1:D:421:ARG:O	2.17	0.45
1:B:223:HIS:HD2	1:B:225:LYS:HG2	1.82	0.44
1:B:370:MET:HE3	1:B:373:ILE:HB	1.99	0.44
1:C:329:SER:HA	1:C:345:TRP:CD1	2.52	0.44
1:D:165:ILE:HD13	1:D:175:LYS:HB2	1.99	0.44
1:A:280:GLU:OE1	1:A:280:GLU:N	2.50	0.44
1:B:458:ARG:HB3	1:B:468:ILE:HD11	2.00	0.44
1:D:365:ASN:N	1:D:365:ASN:OD1	2.50	0.43
1:B:245:LEU:HG	1:B:293:ILE:HB	1.99	0.43
1:A:248:LEU:HD21	1:A:257:VAL:HG11	1.99	0.43
1:D:242:SER:OG	1:D:298:PRO:HD3	2.18	0.43
1:B:243:TYR:HE1	1:B:248:LEU:HB2	1.82	0.43
1:A:315:GLY:HA2	1:D:314:LEU:O	2.19	0.43
1:C:244:ASN:ND2	1:C:247:ASP:OD2	2.51	0.43
1:B:372:LYS:O	1:B:375:GLU:HB2	2.18	0.42
1:C:445:THR:N	1:C:448:ASP:OD2	2.52	0.42
1:C:175:LYS:HE3	1:C:240:MET:HE2	2.00	0.42
1:D:260:ASN:O	1:D:264:LYS:HG2	2.19	0.42
1:B:265:PHE:HA	1:B:268:GLN:HE21	1.84	0.42
1:D:260:ASN:HA	1:D:263:ARG:NH1	2.34	0.42
1:A:366:GLU:OE1	1:A:393:LYS:NZ	2.50	0.42
1:D:188:LYS:HB3	1:D:236:LEU:HB2	2.01	0.42
1:B:242:SER:OG	1:B:243:TYR:N	2.48	0.42
1:D:463:ASP:HA	1:D:464:PRO:HD3	1.93	0.42
1:B:248:LEU:O	1:B:252:THR:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASP:OD1	1:A:147:TYR:N	2.48	0.42
1:B:368:ASP:O	1:B:372:LYS:HG3	2.20	0.42
1:C:165:ILE:HD12	2:C:501:B6Q:CBF	2.49	0.42
1:A:359:PRO:HB2	1:A:362:SER:HB3	2.01	0.41
1:B:463:ASP:HB3	1:B:466:THR:OG1	2.21	0.41
1:C:432:GLY:HA3	1:C:439:ALA:HB2	2.01	0.41
1:A:376:VAL:HG22	1:A:421:ARG:HD3	2.03	0.41
1:C:327:TYR:O	1:C:346:SER:OG	2.35	0.41
1:D:455:LEU:HD11	1:D:459:MET:HE3	2.02	0.41
1:C:193:LYS:HB3	1:C:196:PHE:HD2	1.86	0.41
1:C:345:TRP:HZ2	1:C:373:ILE:HD12	1.86	0.41
1:B:317:ARG:HH21	1:B:320:GLN:HG2	1.86	0.41
1:C:344:MET:O	1:C:347:LEU:HB3	2.20	0.41
1:A:167:LYS:HE3	1:A:167:LYS:HB2	1.85	0.41
1:C:328:ARG:HG2	1:C:333:LEU:HG	2.02	0.41
1:B:383:HIS:O	1:B:387:GLN:NE2	2.54	0.40
2:C:501:B6Q:NBE	2:C:501:B6Q:OAS	2.52	0.40
1:A:429:VAL:HG12	1:A:446:VAL:HG13	2.04	0.40
1:A:429:VAL:HG22	1:A:449:TYR:HB3	2.03	0.40
1:A:392:ARG:O	1:A:406:LYS:NZ	2.33	0.40
1:D:149:VAL:HG21	1:D:164:LEU:HD13	2.04	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	337/368 (92%)	321 (95%)	16 (5%)	0	100	100
1	B	335/368 (91%)	315 (94%)	19 (6%)	1 (0%)	36	57
1	C	326/368 (89%)	297 (91%)	26 (8%)	3 (1%)	14	31
1	D	326/368 (89%)	306 (94%)	20 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1324/1472 (90%)	1239 (94%)	81 (6%)	4 (0%)	36 57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	405	LEU
1	C	287	ASP
1	B	307	ASP
1	C	337	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	A	301/324 (93%)	297 (99%)	4 (1%)	61 81
1	B	295/324 (91%)	294 (100%)	1 (0%)	86 94
1	C	276/324 (85%)	272 (99%)	4 (1%)	59 80
1	D	292/324 (90%)	289 (99%)	3 (1%)	68 84
All	All	1164/1296 (90%)	1152 (99%)	12 (1%)	68 84

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	189	ILE
1	A	212	LYS
1	A	286	CYS
1	A	462	TYR
1	B	468	ILE
1	C	146	ASP
1	C	214	ASP
1	C	306	VAL
1	C	366	GLU
1	D	201	GLN
1	D	400	ASP

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Mol	Chain	Res	Type
1	D	448	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	HIS
1	A	260	ASN
1	A	371	ASN
1	B	244	ASN
1	B	424	HIS
1	C	211	ASN
1	C	223	HIS
1	C	227	HIS
1	C	253	ASN
1	C	365	ASN
1	C	425	ASN
1	D	198	ASN
1	D	201	GLN
1	D	323	GLN

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	$\# Z > 2$
1	PTR	B	321	1	15,16,17	1.33	1 (6%)	17,22,24	0.50	0
1	PTR	C	321	1	15,16,17	1.33	1 (6%)	17,22,24	0.47	0
1	PTR	D	321	1	15,16,17	1.32	1 (6%)	17,22,24	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PTR	A	321	1	15,16,17	1.32	1 (6%)	17,22,24	0.57	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
1	PTR	B	321	1	-	0/10/11/13	0/1/1/1
1	PTR	C	321	1	-	0/10/11/13	0/1/1/1
1	PTR	D	321	1	-	0/10/11/13	0/1/1/1
1	PTR	A	321	1	-	0/10/11/13	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	321	PTR	OH-CZ	-4.47	1.30	1.40
1	B	321	PTR	OH-CZ	-4.45	1.30	1.40
1	A	321	PTR	OH-CZ	-4.36	1.30	1.40
1	D	321	PTR	OH-CZ	-4.35	1.30	1.40

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	321	PTR	1	0
1	C	321	PTR	2	0
1	D	321	PTR	1	0
1	A	321	PTR	1	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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	B6Q	D	501	-	35,36,36	1.90	6 (17%)	48,50,50	2.44	7 (14%)
2	B6Q	A	501	-	35,36,36	1.87	6 (17%)	48,50,50	2.59	7 (14%)
2	B6Q	C	501	-	35,36,36	1.85	6 (17%)	48,50,50	2.41	9 (18%)
2	B6Q	B	501	-	35,36,36	1.84	6 (17%)	48,50,50	2.45	9 (18%)

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	B6Q	D	501	-	-	10/22/32/32	0/4/4/4
2	B6Q	A	501	-	-	8/22/32/32	0/4/4/4
2	B6Q	C	501	-	-	4/22/32/32	0/4/4/4
2	B6Q	B	501	-	-	8/22/32/32	0/4/4/4

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	B6Q	CAM-SAQ	-4.90	1.65	1.73
2	A	501	B6Q	CAM-SAQ	-4.88	1.65	1.73
2	D	501	B6Q	CAP-SAQ	-4.74	1.64	1.74
2	D	501	B6Q	CAP-CAR	-4.71	1.33	1.47
2	D	501	B6Q	CAT-CAR	-4.66	1.39	1.49
2	A	501	B6Q	CAT-CAR	-4.65	1.39	1.49
2	B	501	B6Q	CAM-SAQ	-4.61	1.66	1.73
2	C	501	B6Q	CAP-CAR	-4.57	1.34	1.47
2	B	501	B6Q	CAT-CAR	-4.54	1.40	1.49
2	B	501	B6Q	CAP-CAR	-4.53	1.34	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	B6Q	CAP-CAR	-4.52	1.34	1.47
2	C	501	B6Q	CAT-CAR	-4.48	1.40	1.49
2	D	501	B6Q	CAM-SAQ	-4.45	1.66	1.73
2	A	501	B6Q	CAP-SAQ	-4.25	1.65	1.74
2	B	501	B6Q	CAX-NAZ	-3.85	1.33	1.41
2	C	501	B6Q	CAP-SAQ	-3.82	1.66	1.74
2	B	501	B6Q	CAP-SAQ	-3.78	1.66	1.74
2	C	501	B6Q	CAK-NAL	-3.74	1.34	1.41
2	A	501	B6Q	CAX-NAZ	-3.74	1.34	1.41
2	D	501	B6Q	CAX-NAZ	-3.71	1.34	1.41
2	B	501	B6Q	CAK-NAL	-3.62	1.34	1.41
2	D	501	B6Q	CAK-NAL	-3.52	1.34	1.41
2	A	501	B6Q	CAK-NAL	-3.42	1.34	1.41
2	C	501	B6Q	CAX-NAZ	-3.42	1.34	1.41

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	B6Q	CAM-SAQ-CAP	12.66	95.46	88.34
2	D	501	B6Q	CAM-SAQ-CAP	11.61	94.87	88.34
2	B	501	B6Q	CAM-SAQ-CAP	11.25	94.67	88.34
2	C	501	B6Q	CAM-SAQ-CAP	11.21	94.65	88.34
2	A	501	B6Q	SAQ-CAM-NAN	-8.88	107.83	115.78
2	D	501	B6Q	SAQ-CAM-NAN	-8.41	108.25	115.78
2	B	501	B6Q	SAQ-CAM-NAN	-7.98	108.64	115.78
2	C	501	B6Q	SAQ-CAM-NAN	-7.78	108.82	115.78
2	B	501	B6Q	CAX-NAZ-CBA	-4.70	119.20	127.52
2	A	501	B6Q	CAP-CAO-NAN	-4.62	113.76	117.84
2	D	501	B6Q	CAP-CAO-NAN	-4.55	113.82	117.84
2	B	501	B6Q	CAP-CAO-NAN	-4.27	114.07	117.84
2	C	501	B6Q	CAP-CAO-NAN	-4.21	114.13	117.84
2	D	501	B6Q	CAF-NAE-CAD	3.94	120.44	111.57
2	C	501	B6Q	CAK-NAL-CAM	-3.92	121.08	129.46
2	A	501	B6Q	CAF-NAE-CAD	3.90	120.33	111.57
2	B	501	B6Q	CAK-NAL-CAM	-3.72	121.52	129.46
2	A	501	B6Q	CAX-NAZ-CBA	-3.42	121.47	127.52
2	D	501	B6Q	CAX-NAZ-CBA	-3.07	122.08	127.52
2	D	501	B6Q	CAK-NAL-CAM	-2.95	123.17	129.46
2	A	501	B6Q	CAK-NAL-CAM	-2.91	123.25	129.46
2	C	501	B6Q	CAO-CAP-CAR	-2.87	122.00	124.52
2	B	501	B6Q	CAF-NAE-CAD	2.82	117.91	111.57
2	C	501	B6Q	CAF-NAE-CAD	2.79	117.85	111.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	B6Q	CAR-CAP-SAQ	2.62	129.42	117.41
2	B	501	B6Q	CAR-CAP-SAQ	2.47	128.73	117.41
2	B	501	B6Q	CAG-NAB-CAC	2.41	113.37	109.54
2	C	501	B6Q	CAX-NAZ-CBA	-2.29	123.47	127.52
2	C	501	B6Q	CAG-NAB-CAC	2.27	113.15	109.54
2	A	501	B6Q	CAR-CAP-SAQ	2.18	127.40	117.41
2	B	501	B6Q	CAO-CAP-CAR	-2.11	122.67	124.52
2	D	501	B6Q	CAR-CAP-SAQ	2.11	127.06	117.41

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	B6Q	OBB-CBA-NAZ-CAX
2	B	501	B6Q	OBB-CBA-NAZ-CAX
2	C	501	B6Q	OBB-CBA-NAZ-CAX
2	D	501	B6Q	OBB-CBA-NAZ-CAX
2	A	501	B6Q	CBC-CBA-NAZ-CAX
2	B	501	B6Q	CBC-CBA-NAZ-CAX
2	C	501	B6Q	CBC-CBA-NAZ-CAX
2	D	501	B6Q	CBC-CBA-NAZ-CAX
2	B	501	B6Q	CAW-CAX-NAZ-CBA
2	B	501	B6Q	CAY-CAX-NAZ-CBA
2	D	501	B6Q	CAY-CAX-NAZ-CBA
2	D	501	B6Q	CAW-CAX-NAZ-CBA
2	A	501	B6Q	CAY-CAX-NAZ-CBA
2	A	501	B6Q	CAW-CAX-NAZ-CBA
2	A	501	B6Q	OAS-CAR-CAT-CAU
2	A	501	B6Q	OAS-CAR-CAT-CAY
2	C	501	B6Q	CAY-CAX-NAZ-CBA
2	A	501	B6Q	CAP-CAR-CAT-CAY
2	C	501	B6Q	CAW-CAX-NAZ-CBA
2	A	501	B6Q	CAP-CAR-CAT-CAU
2	D	501	B6Q	OAS-CAR-CAT-CAU
2	D	501	B6Q	CAP-CAR-CAT-CAU
2	D	501	B6Q	OAS-CAR-CAT-CAY
2	B	501	B6Q	CAI-CAH-NAE-CAD
2	D	501	B6Q	CAP-CAR-CAT-CAY
2	B	501	B6Q	CBG-CAH-NAE-CAF
2	D	501	B6Q	SAQ-CAM-NAL-CAK
2	D	501	B6Q	NAN-CAM-NAL-CAK
2	B	501	B6Q	CBG-CAH-NAE-CAD

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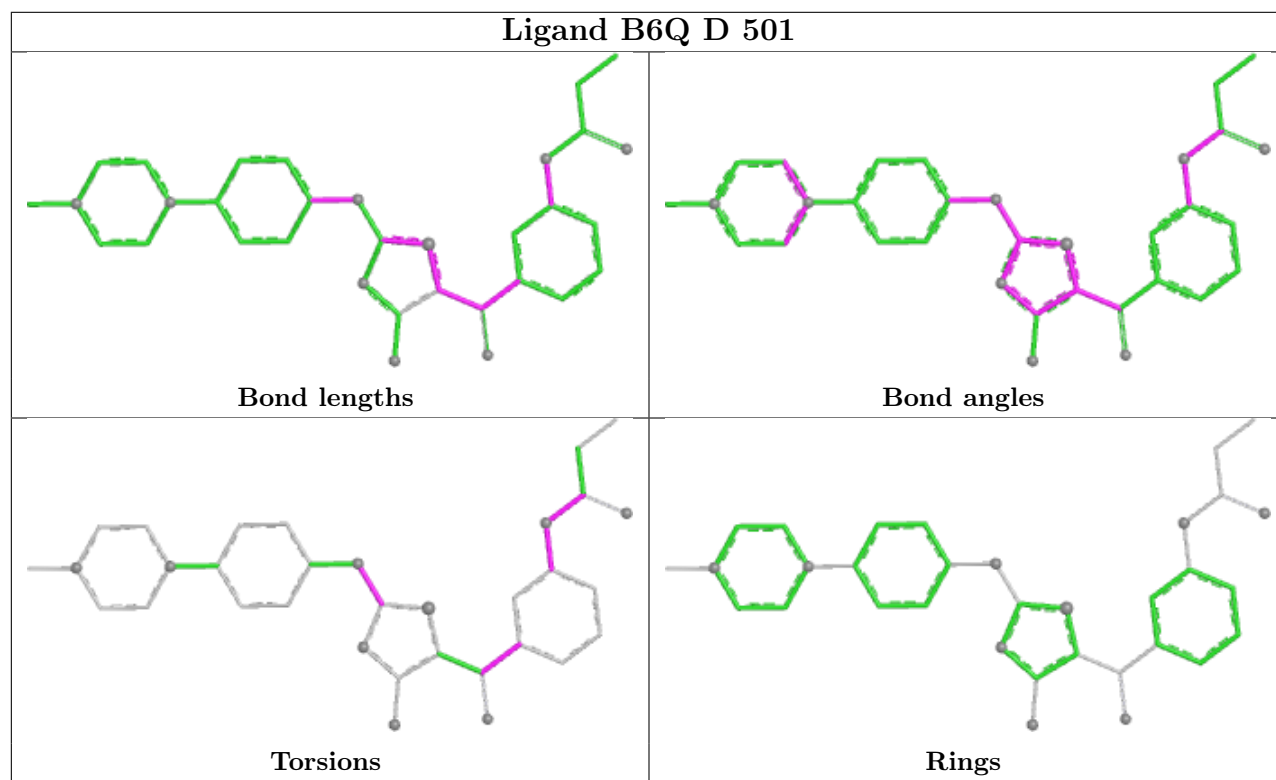
Mol	Chain	Res	Type	Atoms
2	B	501	B6Q	CAI-CAH-NAE-CAF

There are no ring outliers.

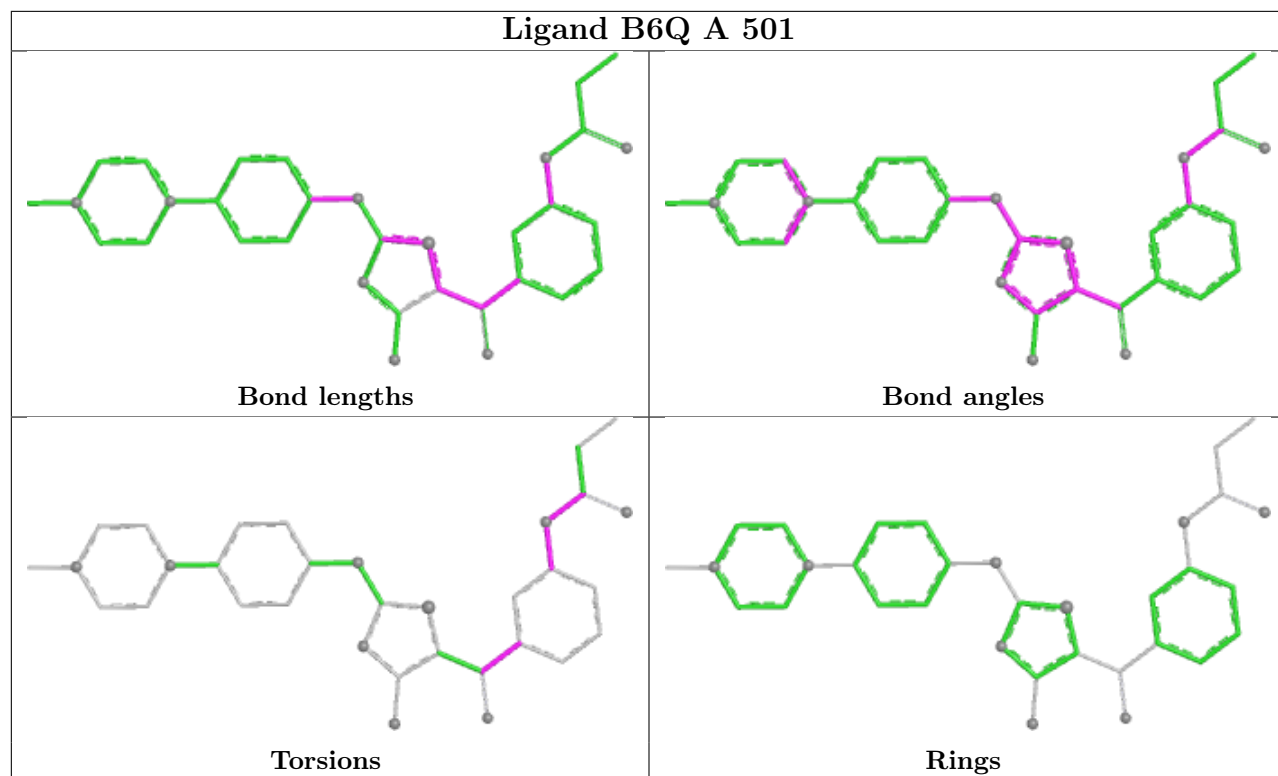
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	B6Q	2	0

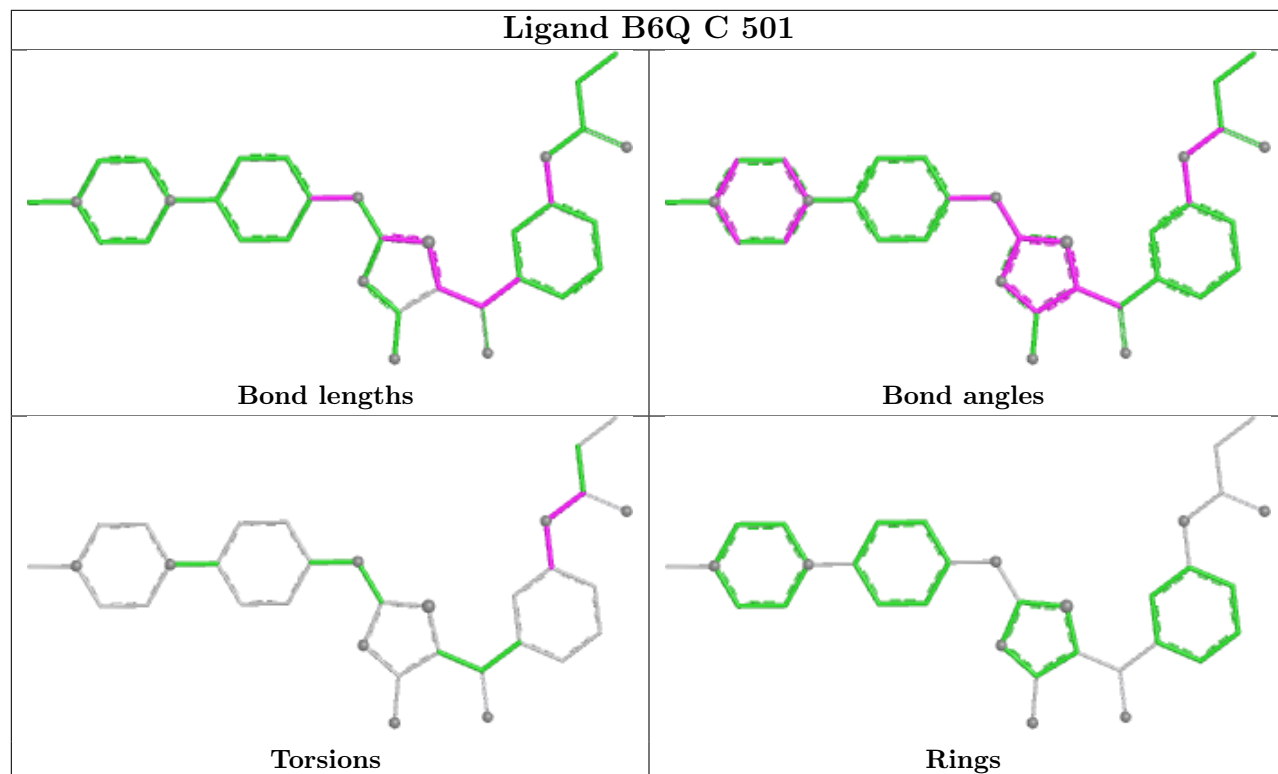
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

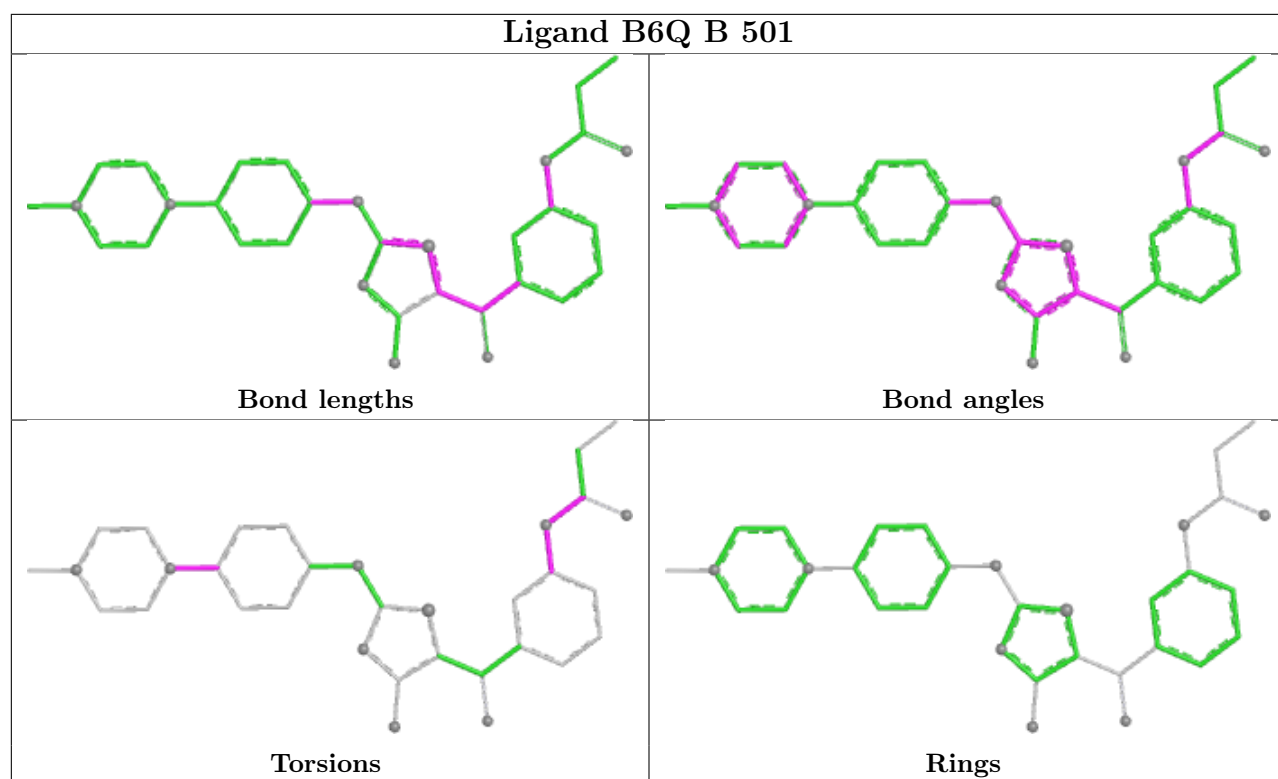


Ligand B6Q A 501



Ligand B6Q C 501





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	341/368 (92%)	-1.00	0 100 100	27, 45, 67, 87	0
1	B	339/368 (92%)	-0.70	1 (0%) 90 89	37, 65, 108, 144	0
1	C	332/368 (90%)	-0.60	3 (0%) 81 79	31, 74, 120, 146	0
1	D	332/368 (90%)	-0.81	2 (0%) 85 84	32, 55, 88, 111	0
All	All	1344/1472 (91%)	-0.78	6 (0%) 88 87	27, 57, 109, 146	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	446	VAL	2.4
1	C	214	ASP	2.4
1	D	253	ASN	2.2
1	C	388	ALA	2.2
1	D	430	GLU	2.1
1	B	455	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	PTR	B	321	16/17	0.98	0.05	52,67,80,80	0
1	PTR	A	321	16/17	0.99	0.05	26,39,53,56	0
1	PTR	C	321	16/17	0.99	0.04	61,70,105,119	0
1	PTR	D	321	16/17	0.99	0.04	37,47,65,68	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

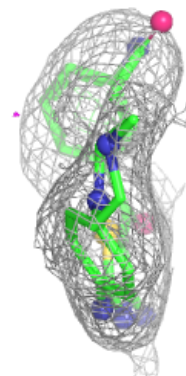
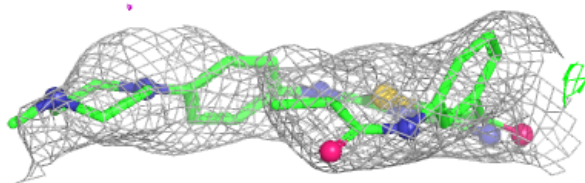
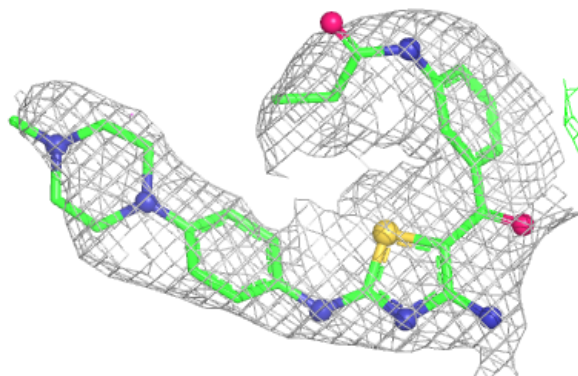
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	B6Q	A	501	33/33	0.98	0.07	44,64,72,77	0
2	B6Q	B	501	33/33	0.98	0.08	63,74,95,100	0
2	B6Q	C	501	33/33	0.98	0.06	62,69,76,82	0
2	B6Q	D	501	33/33	0.99	0.06	52,63,83,92	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.

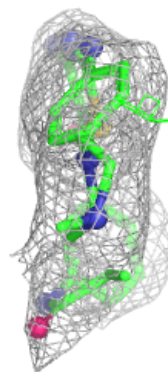
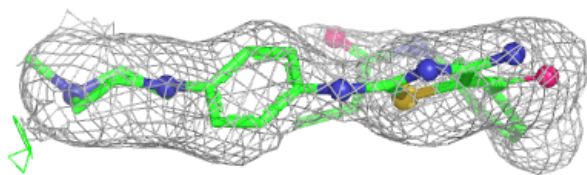
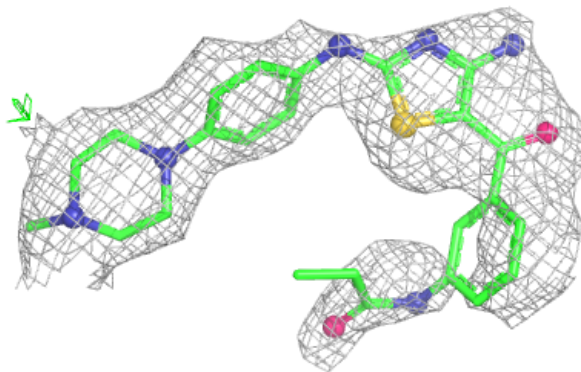
Electron density around B6Q A 501:

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



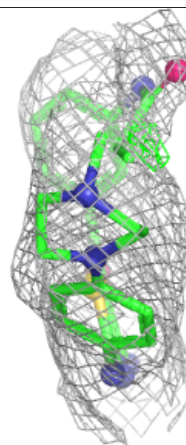
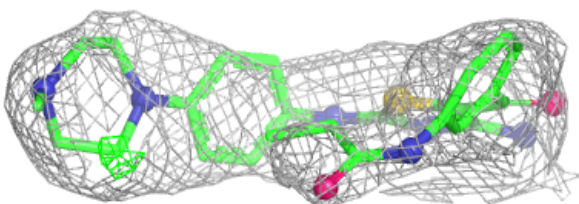
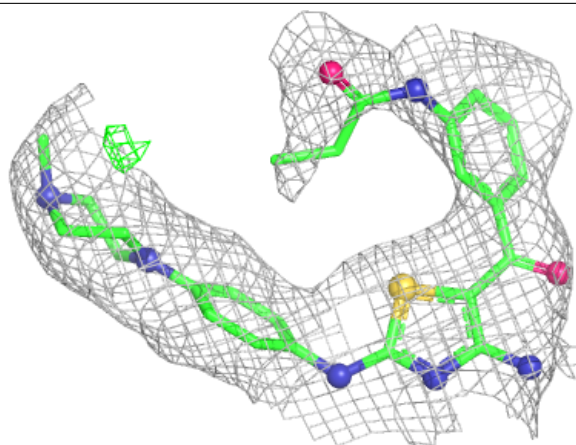
Electron density around B6Q B 501:

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



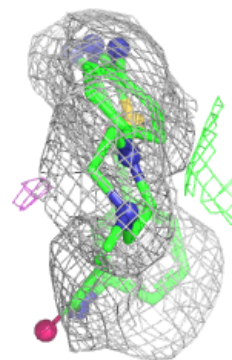
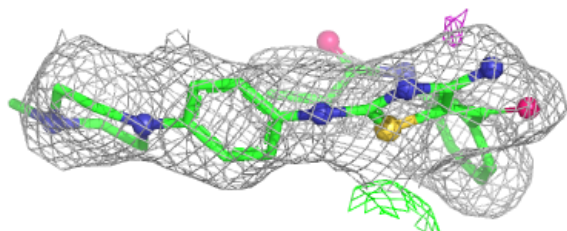
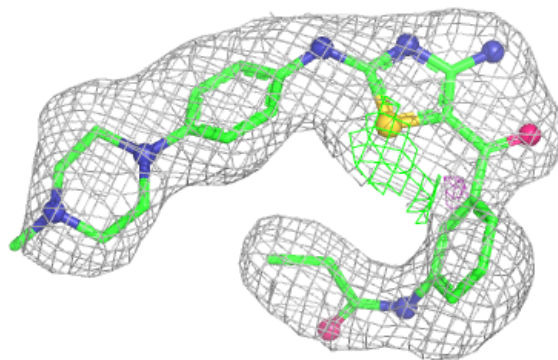
Electron density around B6Q C 501:

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



Electron density around B6Q D 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.