



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:03 AM UTC

PDB ID : 7LQU / pdb_00007lqu
Title : Crystal Structure of HIV-1 RT in Complex with NBD-14075
Authors : Losada, N.; Ruiz, F.X.; Gruber, K.; Das, K.; Arnold, E.
Deposited on : 2021-02-15
Resolution : 2.60 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

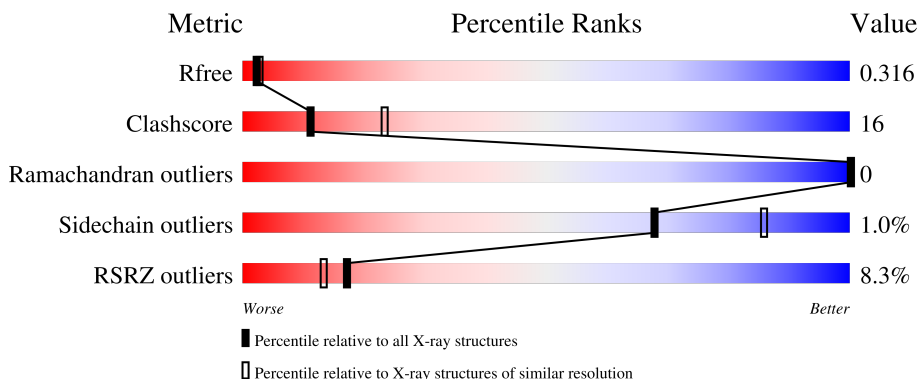
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	557	10% 71% 28% .
2	B	429	6% 67% 28% ..

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
4	SO4	B	501	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8072 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 Reverse transcriptase p66.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	1	0
			4521	2926	751	836	8			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	172	ALA	LYS	engineered mutation	UNP P03366
A	173	ALA	LYS	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366

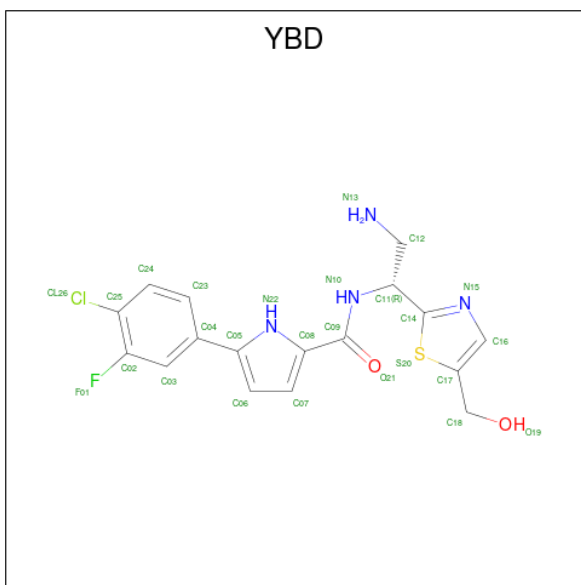
- Molecule 2 is a protein called Reverse transcriptase p51.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	412	Total	C	N	O	S	0	0	0
			3409	2220	564	618	7			

There are 2 discrepancies between the modelled and reference sequences:

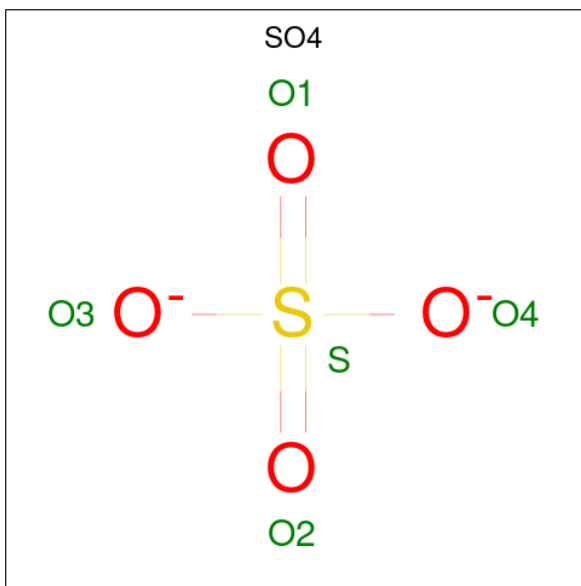
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP P03366
B	280	SER	CYS	engineered mutation	UNP P03366

- Molecule 3 is {N}-[(1 {R})-2-azanyl-1-[5-(hydroxymethyl)-1,3-thiazol-2-yl]ethyl]-5-(4-chloranyl-3-fluoranyl-phenyl)-1 {H}-pyrrole-2-carboxamide (CCD ID: YBD) (formula: C₁₇H₁₆ClFN₄O₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	S	0	0
			26	17	1	1	4	2	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	63	Total 63	O 63	0	0
5	B	38	Total 38	O 38	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.68Å 73.05Å 107.29Å 90.00° 99.58° 90.00°	Depositor
Resolution (Å)	47.90 – 2.60 47.90 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.3 (47.90-2.60) 95.3 (47.90-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.61Å)	Xtriage
Refinement program	PHENIX dev_3051	Depositor
R, R_{free}	0.248 , 0.310 0.252 , 0.316	Depositor DCC
R_{free} test set	1797 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	65.0	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8072	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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: SO4, YBD

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	0.43	0/4643	0.64	2/6311 (0.0%)
2	B	0.43	0/3506	0.66	6/4762 (0.1%)
All	All	0.43	0/8149	0.65	8/11073 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	419	THR	CA-C-N	8.18	128.81	120.38
1	A	419	THR	C-N-CA	8.18	128.81	120.38
2	B	270	ILE	CA-C-N	7.18	131.22	121.74
2	B	270	ILE	C-N-CA	7.18	131.22	121.74
2	B	138	GLU	N-CA-C	5.76	123.06	110.80
2	B	272	PRO	N-CA-C	5.72	121.49	113.53
2	B	271	TYR	CB-CA-C	-5.29	104.00	110.08
2	B	358	ARG	N-CA-C	5.14	116.98	107.73

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	4521	0	4576	154	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3409	0	3441	118	0
3	A	26	0	0	2	0
4	A	10	0	0	1	0
4	B	5	0	0	2	0
5	A	63	0	0	21	0
5	B	38	0	0	39	0
All	All	8072	0	8017	260	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 16.

All (260) 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:61:PHE:CZ	2:B:74:LEU:HD23	1.69	1.26
1:A:544:GLY:O	1:A:548:VAL:HG12	1.31	1.25
2:B:374:LYS:HE3	5:B:635:HOH:O	1.40	1.17
1:A:325:LEU:HG	1:A:387:PRO:HB3	1.38	1.01
2:B:95:PRO:HD2	5:B:628:HOH:O	1.60	1.01
2:B:315:HIS:HA	5:B:606:HOH:O	1.60	1.00
2:B:61:PHE:CZ	2:B:74:LEU:CD2	2.45	0.99
2:B:354:TYR:HB2	5:B:632:HOH:O	1.62	0.98
1:A:110:ASP:HB2	5:A:706:HOH:O	1.62	0.97
1:A:542:ILE:O	1:A:542:ILE:HG22	1.63	0.97
2:B:61:PHE:CE2	2:B:74:LEU:HD23	2.03	0.93
1:A:424:LYS:HB2	5:A:717:HOH:O	1.69	0.93
2:B:241:VAL:HB	5:B:629:HOH:O	1.68	0.92
1:A:137:ASN:HB2	5:A:711:HOH:O	1.71	0.91
1:A:441:TYR:CD2	1:A:544:GLY:HA3	2.05	0.91
2:B:315:HIS:HE1	5:B:633:HOH:O	1.54	0.90
1:A:116:PHE:HZ	5:A:707:HOH:O	1.58	0.87
2:B:94:ILE:HG13	2:B:161:GLN:NE2	1.89	0.87
2:B:94:ILE:HG23	5:B:628:HOH:O	1.74	0.87
1:A:441:TYR:CG	1:A:544:GLY:HA3	2.11	0.86
2:B:94:ILE:HD11	2:B:161:GLN:HG3	1.56	0.84
1:A:536:VAL:HB	1:A:542:ILE:HD13	1.59	0.84
2:B:357:MET:HE2	5:B:623:HOH:O	1.76	0.84
1:A:94:ILE:HG12	1:A:183:TYR:OH	1.77	0.84
1:A:441:TYR:HB3	1:A:544:GLY:HA3	1.60	0.83
1:A:441:TYR:CB	1:A:544:GLY:HA3	2.09	0.83
2:B:126:LYS:HB3	5:B:616:HOH:O	1.77	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:93:GLY:HA2	5:B:611:HOH:O	1.79	0.82
2:B:238:LYS:HE2	5:B:609:HOH:O	1.79	0.81
2:B:257:ILE:HG22	2:B:283:LEU:HD11	1.63	0.80
1:A:105:SER:HG	1:A:198:HIS:CE1	2.03	0.76
1:A:210:LEU:HD12	5:A:754:HOH:O	1.85	0.75
1:A:537:PRO:O	1:A:542:ILE:HD12	1.86	0.75
1:A:37:ILE:HD13	1:A:73:LYS:HB3	1.69	0.74
2:B:93:GLY:CA	5:B:611:HOH:O	2.36	0.73
1:A:109:LEU:HD22	1:A:217:PRO:HG2	1.72	0.72
1:A:229:TRP:HB2	3:A:601:YBD:C07	2.19	0.72
2:B:315:HIS:CE1	5:B:633:HOH:O	2.35	0.72
1:A:105:SER:OG	1:A:198:HIS:ND1	2.21	0.72
1:A:253:THR:HG22	1:A:292:VAL:HG22	1.72	0.71
1:A:547:GLN:HB3	5:A:729:HOH:O	1.91	0.70
1:A:92:LEU:HB3	5:B:625:HOH:O	1.90	0.70
1:A:203:GLU:HA	1:A:206:ARG:HD3	1.74	0.70
1:A:183:TYR:CE2	1:A:184:MET:HG2	2.26	0.70
2:B:235:HIS:HB3	2:B:238:LYS:HG2	1.73	0.70
2:B:66:LYS:NZ	2:B:67:ASP:OD2	2.24	0.68
1:A:86:ASP:HB2	5:B:604:HOH:O	1.95	0.67
2:B:282:LEU:HD21	2:B:296:THR:HG23	1.75	0.67
1:A:210:LEU:CD1	5:A:754:HOH:O	2.42	0.66
2:B:60:VAL:HG23	2:B:75:VAL:HG22	1.78	0.65
2:B:104:LYS:HA	2:B:237:ASP:HB2	1.78	0.65
2:B:354:TYR:HE1	5:B:635:HOH:O	1.79	0.65
1:A:544:GLY:O	1:A:548:VAL:CG1	2.26	0.65
1:A:210:LEU:HB2	5:A:754:HOH:O	1.97	0.64
1:A:548:VAL:CG2	5:A:734:HOH:O	2.45	0.64
1:A:90:VAL:HG11	2:B:140:PRO:HB3	1.78	0.64
1:A:441:TYR:HB3	1:A:544:GLY:CA	2.28	0.64
1:A:542:ILE:O	1:A:542:ILE:CG2	2.37	0.64
1:A:191:SER:OG	1:A:198:HIS:ND1	2.21	0.63
2:B:228:LEU:CD2	5:B:630:HOH:O	2.47	0.62
1:A:441:TYR:CB	1:A:544:GLY:CA	2.78	0.62
1:A:223:LYS:HB3	1:A:227:PHE:HE1	1.64	0.61
1:A:40:GLU:HA	1:A:43:LYS:HD2	1.82	0.61
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.00	0.61
1:A:325:LEU:CG	1:A:387:PRO:HB3	2.22	0.60
1:A:449:GLU:OE2	1:A:449:GLU:N	2.33	0.60
1:A:490:GLY:O	1:A:528:LYS:NZ	2.32	0.60
2:B:356:ARG:HG2	2:B:357:MET:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:270:ILE:HG12	2:B:346:PHE:HB3	1.83	0.59
1:A:542:ILE:HG23	2:B:283:LEU:HD23	1.85	0.59
1:A:551:LEU:HG	5:A:710:HOH:O	2.02	0.59
2:B:46:LYS:HE2	2:B:116:PHE:HB3	1.85	0.59
2:B:238:LYS:CD	5:B:609:HOH:O	2.50	0.58
1:A:56:TYR:O	1:A:143:ARG:NH2	2.30	0.58
1:A:543:GLY:O	1:A:544:GLY:C	2.46	0.58
2:B:95:PRO:CD	5:B:628:HOH:O	2.33	0.58
2:B:94:ILE:CG1	2:B:161:GLN:NE2	2.65	0.58
1:A:266:TRP:O	1:A:269:GLN:HG2	2.03	0.58
2:B:154:LYS:CD	5:B:613:HOH:O	2.51	0.58
2:B:136:ASN:HB2	5:B:610:HOH:O	2.05	0.57
2:B:360:ALA:HB2	5:B:608:HOH:O	2.03	0.57
1:A:277:ARG:HH12	1:A:356:ARG:NH2	2.03	0.57
2:B:94:ILE:CD1	2:B:161:GLN:HG3	2.33	0.57
2:B:350:LYS:HE3	2:B:378:GLU:OE1	2.05	0.56
1:A:388:LYS:HE3	5:A:740:HOH:O	2.05	0.56
2:B:354:TYR:CB	5:B:632:HOH:O	2.37	0.56
1:A:543:GLY:HA3	2:B:285:GLY:O	2.05	0.56
2:B:360:ALA:HB1	2:B:367:GLN:NE2	2.21	0.56
1:A:183:TYR:CD2	1:A:184:MET:HG2	2.41	0.56
1:A:552:VAL:CG1	5:A:734:HOH:O	2.52	0.55
2:B:354:TYR:CA	5:B:632:HOH:O	2.53	0.55
2:B:89:GLU:HG3	2:B:90:VAL:H	1.72	0.55
1:A:116:PHE:CZ	5:A:707:HOH:O	2.43	0.55
1:A:356:ARG:HD3	1:A:367:GLN:NE2	2.22	0.55
1:A:404:GLU:OE2	1:A:431:LYS:NZ	2.32	0.55
2:B:421:PRO:HB2	2:B:423:VAL:HG12	1.87	0.55
1:A:516:GLU:O	1:A:520:GLN:HG3	2.07	0.55
1:A:540:LYS:O	1:A:541:GLY:C	2.50	0.55
1:A:257:ILE:HB	1:A:283:LEU:HD21	1.89	0.54
2:B:393:ILE:HD13	2:B:398:TRP:HB2	1.89	0.54
1:A:325:LEU:HG	1:A:387:PRO:CB	2.26	0.54
1:A:50:ILE:CG1	1:A:143:ARG:HB3	2.38	0.53
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.89	0.53
1:A:24:TRP:HE3	1:A:25:PRO:HD2	1.73	0.53
1:A:57:ASN:OD1	1:A:143:ARG:NH1	2.42	0.53
2:B:73:LYS:NZ	2:B:146:TYR:OH	2.42	0.53
1:A:7:THR:HG21	1:A:121:ASP:HA	1.91	0.53
1:A:17:ASP:O	1:A:83:ARG:HD3	2.09	0.52
2:B:20:LYS:NZ	5:B:604:HOH:O	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:237:ASP:C	5:B:609:HOH:O	2.51	0.52
1:A:543:GLY:HA3	2:B:283:LEU:O	2.09	0.52
1:A:324:ASP:O	1:A:343:GLN:HG2	2.10	0.52
1:A:441:TYR:CG	1:A:544:GLY:CA	2.89	0.52
2:B:324:ASP:O	2:B:343:GLN:HG2	2.09	0.52
2:B:320:ASP:OD2	2:B:323:LYS:HG3	2.10	0.52
1:A:223:LYS:HB3	1:A:227:PHE:CE1	2.44	0.52
1:A:548:VAL:HA	1:A:551:LEU:HB2	1.92	0.52
2:B:20:LYS:HE2	2:B:55:PRO:HB2	1.91	0.52
1:A:167:ILE:HD13	1:A:214:LEU:HD11	1.90	0.52
1:A:458:VAL:CG2	2:B:286:THR:HG21	2.40	0.52
2:B:94:ILE:CG2	5:B:628:HOH:O	2.47	0.52
1:A:547:GLN:NE2	5:A:708:HOH:O	2.41	0.51
2:B:136:ASN:HB3	2:B:138:GLU:HG3	1.92	0.51
1:A:252:TRP:CH2	1:A:260:LEU:HD22	2.45	0.51
1:A:268:SER:HB3	1:A:274:ILE:HB	1.92	0.51
1:A:364:ASP:HB3	1:A:423:VAL:HG13	1.93	0.51
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.45	0.51
1:A:39:THR:HG22	1:A:43:LYS:HE3	1.93	0.51
2:B:61:PHE:HZ	2:B:74:LEU:CD2	2.17	0.51
2:B:360:ALA:O	2:B:362:THR:N	2.38	0.51
1:A:376:THR:O	1:A:380:ILE:HG12	2.10	0.51
1:A:254:VAL:HG23	1:A:291:GLU:O	2.11	0.51
2:B:94:ILE:HG13	2:B:161:GLN:CD	2.34	0.51
1:A:199:ARG:O	1:A:203:GLU:HG2	2.11	0.51
1:A:207:GLN:HA	1:A:210:LEU:HD12	1.93	0.51
1:A:325:LEU:HD11	1:A:341:ILE:CG2	2.41	0.51
1:A:541:GLY:HA2	1:A:546:GLU:HB2	1.92	0.51
1:A:57:ASN:OD1	1:A:131:THR:OG1	2.23	0.50
1:A:458:VAL:HG23	2:B:286:THR:HG21	1.93	0.50
1:A:23:GLN:HG2	1:A:133:PRO:HD3	1.93	0.50
1:A:540:LYS:HB2	1:A:542:ILE:HG13	1.92	0.50
1:A:7:THR:CG2	1:A:121:ASP:HA	2.41	0.50
1:A:219:LYS:HG2	1:A:220:LYS:H	1.77	0.50
1:A:267:ALA:HB1	1:A:310:LEU:HD21	1.94	0.50
2:B:275:LYS:HD2	4:B:501:SO4:O4	2.12	0.50
1:A:115:TYR:OH	1:A:150:PRO:O	2.30	0.49
1:A:93:GLY:HA3	2:B:137:ASN:CG	2.37	0.49
2:B:194:GLU:O	2:B:198:HIS:N	2.44	0.49
1:A:159:ILE:O	1:A:163:SER:OG	2.24	0.49
1:A:171:PHE:CD2	1:A:205:LEU:HD13	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:423:VAL:HA	2:B:426:TRP:CD1	2.47	0.49
2:B:335:GLY:HA3	2:B:356:ARG:HD3	1.94	0.48
1:A:203:GLU:HA	1:A:206:ARG:HH11	1.77	0.48
1:A:114:ALA:HA	5:A:726:HOH:O	2.14	0.48
2:B:354:TYR:CD1	2:B:374:LYS:HD2	2.48	0.48
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.53	0.48
1:A:379:SER:CB	1:A:387:PRO:HD3	2.44	0.47
2:B:174:GLN:HG3	2:B:175:ASN:OD1	2.14	0.47
2:B:359:GLY:HA3	2:B:366:LYS:NZ	2.29	0.47
1:A:441:TYR:CD2	1:A:544:GLY:CA	2.88	0.47
2:B:239:TRP:CZ3	2:B:350:LYS:HE2	2.49	0.47
1:A:50:ILE:HG12	1:A:143:ARG:HB3	1.96	0.47
2:B:86:ASP:HB3	2:B:88:TRP:CZ2	2.50	0.47
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.55	0.47
1:A:356:ARG:HG2	1:A:357:MET:O	2.15	0.47
1:A:380:ILE:HD12	2:B:27:THR:HG22	1.97	0.47
1:A:443:ASP:HB2	1:A:548:VAL:HG13	1.97	0.47
1:A:551:LEU:N	5:A:710:HOH:O	2.47	0.47
2:B:154:LYS:HD3	5:B:613:HOH:O	2.13	0.46
1:A:547:GLN:C	5:A:710:HOH:O	2.58	0.46
1:A:547:GLN:CB	5:A:729:HOH:O	2.53	0.46
2:B:94:ILE:HD11	2:B:161:GLN:CG	2.38	0.46
1:A:443:ASP:OD1	1:A:444:GLY:N	2.44	0.46
2:B:238:LYS:HD3	5:B:609:HOH:O	2.14	0.46
2:B:354:TYR:HA	5:B:632:HOH:O	2.14	0.46
1:A:35:VAL:HG22	1:A:132:ILE:HG21	1.98	0.46
1:A:447:ASN:O	1:A:451:LYS:N	2.48	0.46
2:B:317:VAL:HG22	2:B:347:LYS:HD2	1.98	0.46
1:A:105:SER:HG	1:A:198:HIS:CG	2.26	0.45
1:A:447:ASN:OD1	1:A:450:THR:HG23	2.16	0.45
1:A:419:THR:HA	1:A:420:PRO:HD3	1.72	0.45
1:A:537:PRO:O	1:A:542:ILE:CD1	2.60	0.45
1:A:424:LYS:HD3	1:A:426:TRP:CH2	2.51	0.45
1:A:544:GLY:HA2	2:B:286:THR:HG22	1.98	0.45
2:B:76:ASP:OD2	2:B:78:ARG:NH2	2.47	0.45
1:A:229:TRP:CB	3:A:601:YBD:C07	2.92	0.45
2:B:103:LYS:O	2:B:236:PRO:HG2	2.16	0.45
2:B:278:GLN:HG3	2:B:298:GLU:HB3	1.97	0.45
1:A:252:TRP:CE3	1:A:256:ASP:HB3	2.52	0.45
1:A:405:TYR:CE2	1:A:407:GLN:HB2	2.52	0.45
1:A:92:LEU:CG	5:B:625:HOH:O	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ASP:O	1:A:115:TYR:N	2.50	0.45
1:A:543:GLY:CA	2:B:285:GLY:O	2.65	0.45
1:A:115:TYR:CE2	1:A:156:SER:HB3	2.52	0.44
1:A:427:TYR:HB2	4:A:602:SO4:O3	2.18	0.44
2:B:228:LEU:HD21	5:B:630:HOH:O	2.16	0.44
1:A:206:ARG:HH21	1:A:218:ASP:CG	2.25	0.44
1:A:228:LEU:HD22	1:A:233:GLU:CG	2.48	0.44
1:A:228:LEU:HD22	1:A:233:GLU:HG2	1.99	0.44
1:A:363:ASN:OD1	1:A:363:ASN:C	2.60	0.44
1:A:277:ARG:HH12	1:A:356:ARG:HH21	1.65	0.44
2:B:237:ASP:CB	5:B:609:HOH:O	2.65	0.44
2:B:336:GLN:HB3	2:B:353:LYS:NZ	2.33	0.44
1:A:269:GLN:HA	1:A:351:THR:O	2.18	0.44
2:B:28:GLU:HB2	2:B:135:ILE:HD11	2.00	0.44
2:B:84:THR:HB	2:B:154:LYS:HE2	1.99	0.44
2:B:277:ARG:N	4:B:501:SO4:O2	2.51	0.44
1:A:101:LYS:HE2	1:A:101:LYS:H	1.82	0.43
2:B:274:ILE:O	2:B:275:LYS:HD3	2.17	0.43
1:A:548:VAL:HG22	5:A:734:HOH:O	2.16	0.43
2:B:160:PHE:HD2	2:B:164:MET:HB2	1.83	0.43
2:B:236:PRO:HA	2:B:239:TRP:NE1	2.33	0.43
1:A:257:ILE:O	1:A:261:VAL:HG23	2.18	0.43
1:A:130:PHE:CE1	1:A:144:TYR:HB3	2.53	0.43
2:B:212:TRP:O	2:B:215:THR:HG22	2.19	0.43
2:B:257:ILE:CG2	2:B:283:LEU:HD11	2.41	0.43
1:A:88:TRP:HB3	2:B:52:PRO:HA	2.01	0.43
2:B:238:LYS:CE	5:B:609:HOH:O	2.47	0.43
2:B:254:VAL:O	2:B:258:GLN:HG3	2.18	0.43
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.88	0.43
2:B:42:GLU:OE2	2:B:49:LYS:HG2	2.19	0.43
2:B:282:LEU:CD2	2:B:296:THR:HG23	2.44	0.43
1:A:231:GLY:C	1:A:242:GLN:HB2	2.44	0.42
2:B:236:PRO:HA	2:B:239:TRP:CE2	2.55	0.42
2:B:238:LYS:N	5:B:609:HOH:O	2.51	0.42
2:B:289:LEU:HD23	2:B:289:LEU:HA	1.81	0.42
1:A:122:GLU:OE1	1:A:125:ARG:HD2	2.20	0.42
1:A:311:LYS:HG3	1:A:312:GLU:N	2.35	0.42
1:A:369:THR:O	1:A:373[B]:GLN:HG2	2.20	0.42
1:A:122:GLU:HA	1:A:125:ARG:HG3	2.02	0.42
1:A:88:TRP:HB2	2:B:54:ASN:O	2.20	0.42
2:B:410:TRP:O	2:B:411:ILE:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LEU:HB3	1:A:214:LEU:HB2	2.02	0.42
1:A:456:GLY:C	5:A:734:HOH:O	2.63	0.42
2:B:305:GLU:O	2:B:308:GLU:HB2	2.19	0.42
2:B:63:ILE:O	2:B:71:TRP:HA	2.20	0.42
2:B:175:ASN:HB3	2:B:178:ILE:HG12	2.02	0.41
1:A:92:LEU:CB	5:B:625:HOH:O	2.61	0.41
1:A:210:LEU:CB	5:A:754:HOH:O	2.64	0.41
2:B:136:ASN:ND2	5:B:610:HOH:O	2.51	0.41
1:A:42:GLU:OE2	1:A:49:LYS:HG3	2.20	0.41
1:A:233:GLU:O	1:A:239:TRP:HA	2.20	0.41
2:B:258:GLN:HG2	2:B:283:LEU:CD2	2.50	0.41
2:B:260:LEU:HD21	2:B:303:LEU:HD13	2.02	0.41
2:B:316:GLY:N	5:B:606:HOH:O	2.45	0.41
1:A:50:ILE:HG13	1:A:143:ARG:HB3	2.02	0.41
2:B:136:ASN:CB	5:B:610:HOH:O	2.66	0.41
1:A:357:MET:O	1:A:359:GLY:N	2.51	0.41
1:A:452:LEU:HD23	1:A:470:THR:HA	2.03	0.41
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.84	0.41
2:B:104:LYS:HB2	2:B:192:ASP:HA	2.02	0.41
2:B:165:THR:O	2:B:168:LEU:N	2.50	0.41
2:B:424:LYS:HG2	2:B:425:LEU:HD23	2.02	0.41
1:A:233:GLU:HG3	1:A:242:GLN:HA	2.02	0.41
2:B:61:PHE:CZ	2:B:74:LEU:HD21	2.49	0.41
2:B:105:SER:O	2:B:190:GLY:HA2	2.21	0.41
2:B:120:LEU:HD23	2:B:125:ARG:HG2	2.02	0.41
1:A:116:PHE:CZ	1:A:151:GLN:HG2	2.57	0.40
1:A:164:MET:HE1	1:A:209:LEU:CD2	2.51	0.40
1:A:282:LEU:HD21	1:A:296:THR:HG23	2.02	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	555/557 (100%)	506 (91%)	49 (9%)	0	100	100
2	B	408/429 (95%)	374 (92%)	34 (8%)	0	100	100
All	All	963/986 (98%)	880 (91%)	83 (9%)	0	100	100

There are no Ramachandran outliers to report.

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	496/495 (100%)	491 (99%)	5 (1%)	68	86
2	B	374/390 (96%)	369 (99%)	5 (1%)	61	82
All	All	870/885 (98%)	860 (99%)	10 (1%)	68	84

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	363	ASN
1	A	373[A]	GLN
1	A	373[B]	GLN
1	A	399	GLU
2	B	58	THR
2	B	185	ASP
2	B	280	SER
2	B	287	LYS
2	B	357	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	147	ASN
1	A	182	GLN

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Mol	Chain	Res	Type
1	A	242	GLN
1	A	306	ASN
1	A	340	GLN
1	A	474	ASN
1	A	520	GLN
2	B	161	GLN
2	B	255	ASN
2	B	258	GLN
2	B	407	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	YBD	A	601	-	28,28,28	2.38	10 (35%)	37,39,39	1.70	9 (24%)
4	SO4	A	602	-	4,4,4	0.42	0	6,6,6	0.07	0
4	SO4	B	501	-	4,4,4	0.42	0	6,6,6	0.07	0
4	SO4	A	603	-	4,4,4	0.43	0	6,6,6	0.07	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
3	YBD	A	601	-	-	6/19/20/20	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	YBD	C09-N10	5.46	1.45	1.34
3	A	601	YBD	C14-S20	-5.29	1.63	1.75
3	A	601	YBD	C16-N15	5.11	1.47	1.37
3	A	601	YBD	C17-S20	-3.96	1.60	1.73
3	A	601	YBD	C04-C05	3.67	1.53	1.47
3	A	601	YBD	C18-C17	3.43	1.53	1.49
3	A	601	YBD	O21-C09	-2.41	1.19	1.23
3	A	601	YBD	C25-CL26	2.25	1.78	1.73
3	A	601	YBD	C08-C09	2.05	1.53	1.48
3	A	601	YBD	C14-N15	2.05	1.39	1.30

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	YBD	C17-S20-C14	4.63	98.82	90.69
3	A	601	YBD	S20-C14-N15	-4.33	110.07	115.98
3	A	601	YBD	C17-C16-N15	-3.64	110.15	117.00
3	A	601	YBD	C04-C05-N22	2.56	126.00	122.76
3	A	601	YBD	C07-C08-C09	-2.31	126.37	132.26
3	A	601	YBD	C08-C09-N10	2.26	119.88	116.86
3	A	601	YBD	C04-C05-C06	-2.25	126.31	130.58
3	A	601	YBD	C11-C14-S20	2.17	124.90	121.80
3	A	601	YBD	C18-C17-S20	2.17	124.98	120.20

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	YBD	C12-C11-N10-C09
3	A	601	YBD	C14-C11-N10-C09
3	A	601	YBD	N22-C08-C09-N10
3	A	601	YBD	N22-C08-C09-O21

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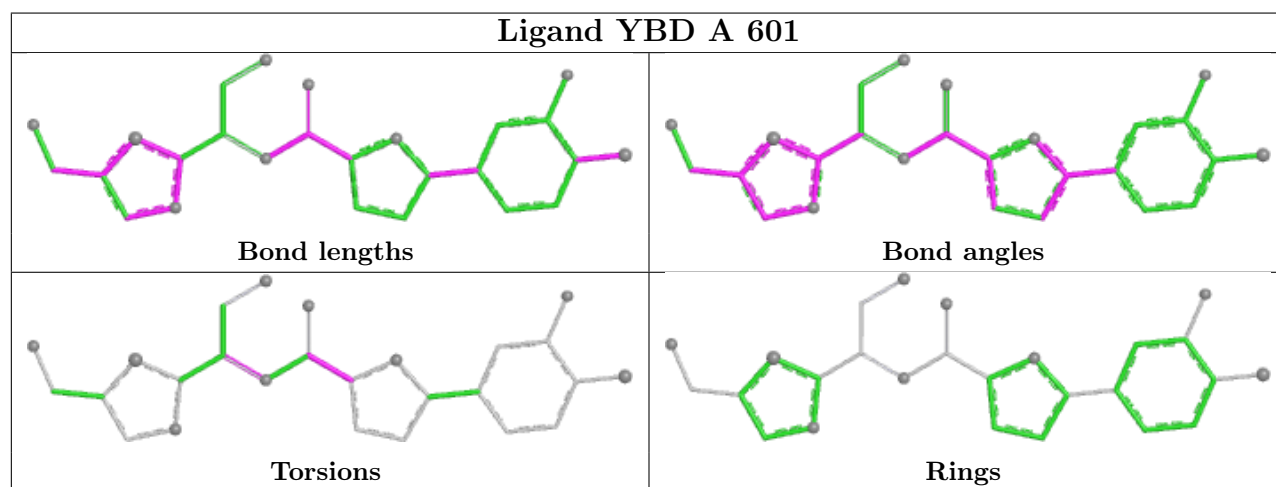
Mol	Chain	Res	Type	Atoms
3	A	601	YBD	C07-C08-C09-O21
3	A	601	YBD	C07-C08-C09-N10

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	YBD	2	0
4	A	602	SO4	1	0
4	B	501	SO4	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.



5.7 Other polymers [i](#)

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	556/557 (99%)	0.83	53 (9%)	14 10	35, 95, 130, 145	1 (0%)
2	B	412/429 (96%)	0.66	27 (6%)	24 19	41, 78, 116, 140	0
All	All	968/986 (98%)	0.76	80 (8%)	17 13	35, 85, 128, 145	1 (0%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	551	LEU	4.7
1	A	552	VAL	4.6
1	A	92	LEU	4.5
1	A	214	LEU	4.5
2	B	241	VAL	4.4
1	A	189	VAL	4.3
1	A	544	GLY	4.1
2	B	230	MET	3.9
1	A	279	LEU	3.9
2	B	232	TYR	3.9
1	A	289	LEU	3.8
2	B	240	THR	3.7
1	A	90	VAL	3.6
1	A	554	ALA	3.5
2	B	237	ASP	3.5
1	A	294	PRO	3.4
1	A	138	GLU	3.3
2	B	359	GLY	3.3
1	A	288	ALA	3.2
2	B	94	ILE	3.2
1	A	325	LEU	3.2
1	A	227	PHE	3.2
1	A	180	ILE	3.0
1	A	245	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	115	TYR	3.0
2	B	212	TRP	3.0
2	B	239	TRP	3.0
2	B	229	TRP	2.9
1	A	548	VAL	2.9
1	A	202	ILE	2.9
2	B	228	LEU	2.9
1	A	91	GLN	2.9
1	A	246	LEU	2.8
1	A	309	ILE	2.8
2	B	5	ILE	2.8
1	A	301	LEU	2.8
1	A	184	MET	2.8
1	A	303	LEU	2.7
1	A	2	ILE	2.7
1	A	205	LEU	2.7
2	B	428	GLN	2.7
2	B	90	VAL	2.6
1	A	187	LEU	2.6
1	A	247	PRO	2.6
1	A	553	SER	2.5
1	A	100	LEU	2.5
1	A	318	TYR	2.5
1	A	545	ASN	2.5
1	A	244	ILE	2.5
2	B	214	LEU	2.5
1	A	313	PRO	2.5
1	A	65	LYS	2.4
1	A	542	ILE	2.4
1	A	290	THR	2.4
2	B	14	PRO	2.4
2	B	358	ARG	2.4
1	A	292	VAL	2.4
1	A	24	TRP	2.3
1	A	61	PHE	2.3
1	A	108	VAL	2.3
1	A	547	GLN	2.3
2	B	93	GLY	2.2
2	B	74	LEU	2.2
1	A	225	PRO	2.2
2	B	354	TYR	2.2
2	B	238	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	420	PRO	2.1
1	A	546	GLU	2.1
2	B	383	TRP	2.1
2	B	360	ALA	2.1
1	A	283	LEU	2.1
1	A	136	ASN	2.1
1	A	221	HIS	2.1
1	A	130	PHE	2.1
1	A	127	TYR	2.1
1	A	218	ASP	2.0
1	A	299	ALA	2.0
2	B	149	LEU	2.0
2	B	213	GLY	2.0
2	B	60	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

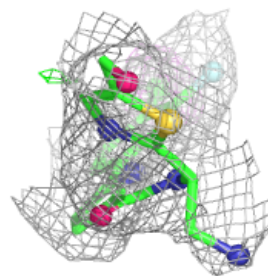
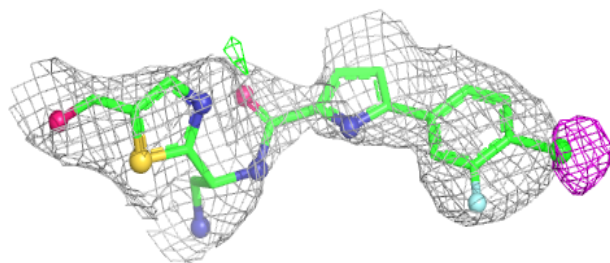
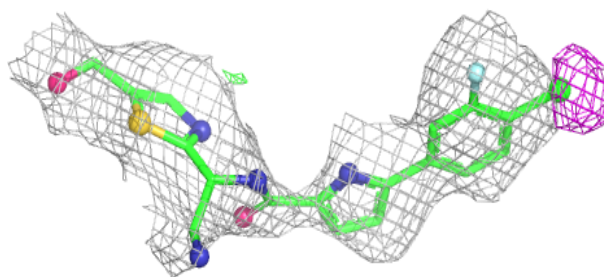
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	YBD	A	601	26/26	0.65	0.16	97,107,111,114	0
4	SO4	B	501	5/5	0.74	0.19	94,95,97,99	0
4	SO4	A	603	5/5	0.75	0.13	96,98,99,99	0
4	SO4	A	602	5/5	0.84	0.28	91,92,92,93	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 YBD A 601:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.