



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:53 AM UTC

PDB ID : 3KC1 / pdb_00003kc1
Title : Crystal structure of human liver FBPase in complex with tricyclic inhibitor 19a
Authors : Takahashi, M.; Sone, J.; Hanzawa, H.
Deposited on : 2009-10-20
Resolution : 2.25 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

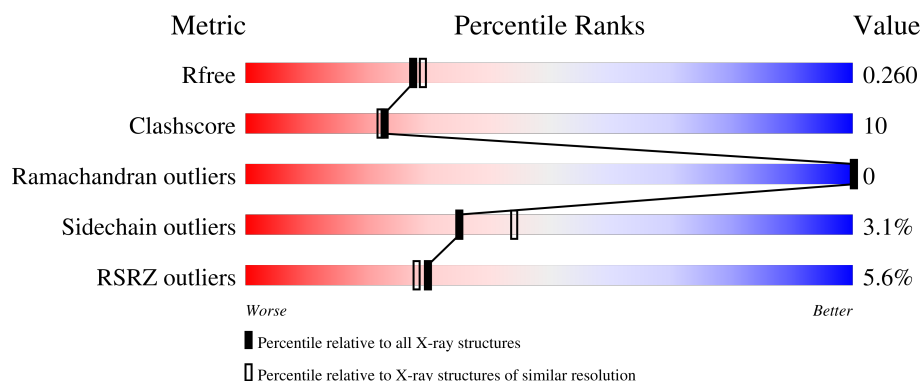
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	337	 7% 74% 18% • 5%
1	B	337	 5% 72% 20% • 6%
1	C	337	 5% 76% 17% • 6%
1	D	337	 3% 74% 18% • 6%

2 Entry composition [i](#)

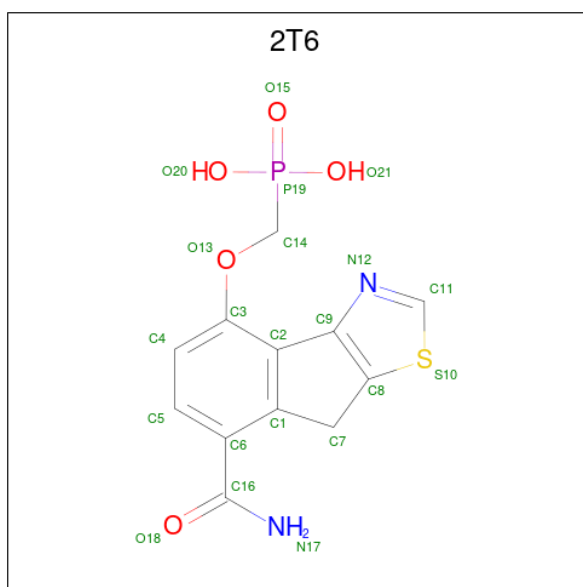
There are 3 unique types of molecules in this entry. The entry contains 10173 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2441	1556	409	459	17			
1	B	318	Total	C	N	O	S	0	0	0
			2434	1551	408	458	17			
1	C	318	Total	C	N	O	S	0	0	0
			2434	1551	408	458	17			
1	D	317	Total	C	N	O	S	0	0	0
			2425	1545	406	457	17			

- Molecule 2 is {[(7-carbamoyl-8H-indeno[1,2-d][1,3]thiazol-4-yl)oxy]methyl}phosphonic acid (CCD ID: 2T6) (formula: C₁₂H₁₁N₂O₅PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			21	12	2	5	1	1		
2	B	1	Total	C	N	O	P	S	0	0
			21	12	2	5	1	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	S	0	0
			21	12	2	5	1	1		
2	D	1	Total	C	N	O	P	S	0	0
			21	12	2	5	1	1		

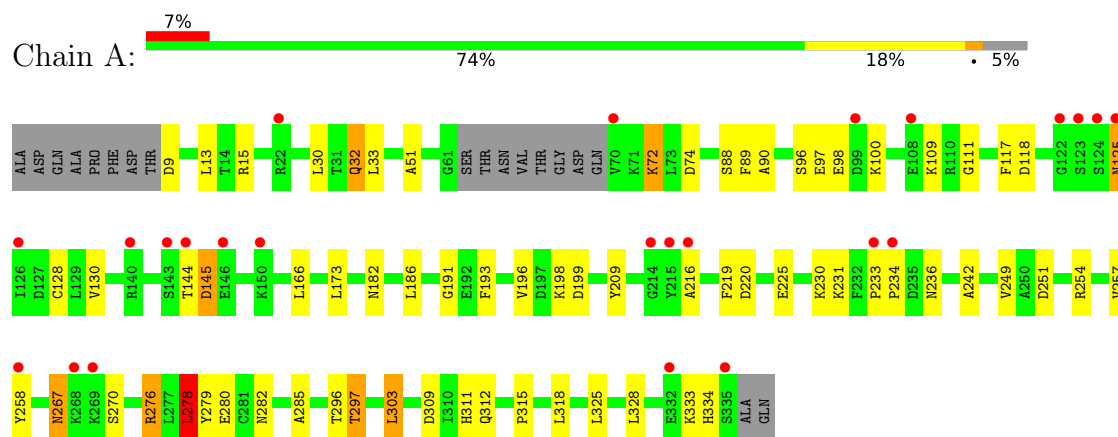
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	60	Total	O	0	0
			60	60		
3	B	72	Total	O	0	0
			72	72		
3	C	102	Total	O	0	0
			102	102		
3	D	121	Total	O	0	0
			121	121		

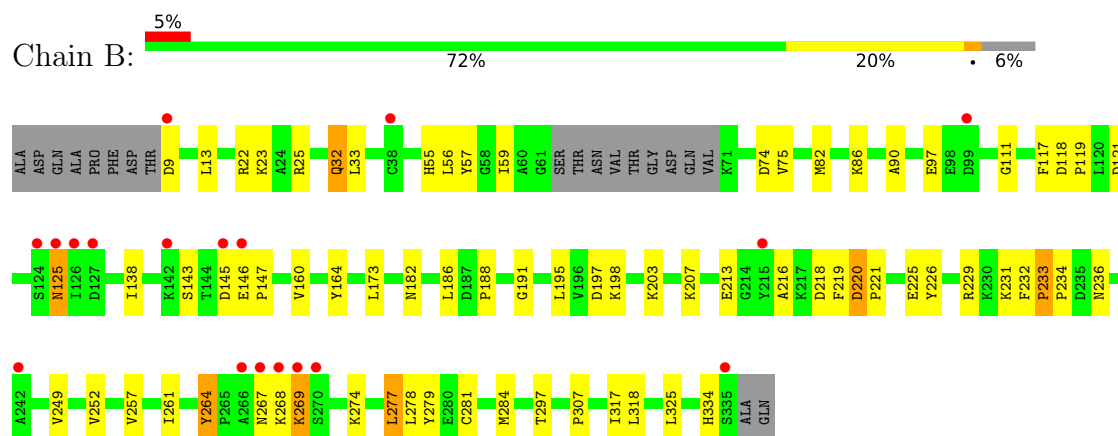
3 Residue-property plots [i](#)

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.

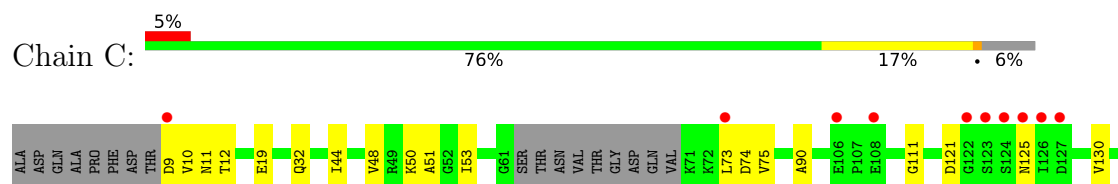
• Molecule 1: Fructose-1,6-bisphosphatase 1

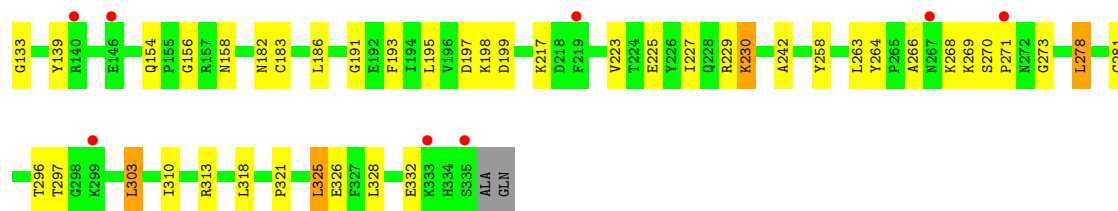


• Molecule 1: Fructose-1,6-bisphosphatase 1

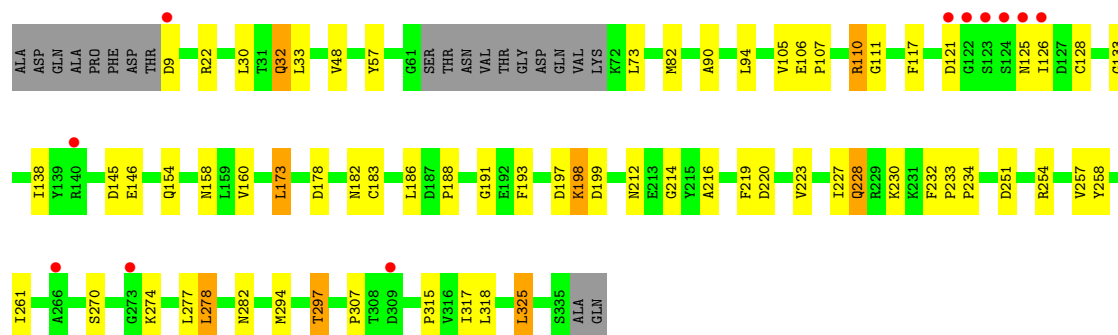
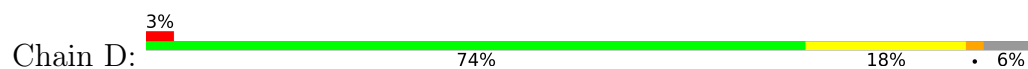


• Molecule 1: Fructose-1,6-bisphosphatase 1





● Molecule 1: Fructose-1,6-bisphosphatase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.36Å 83.21Å 278.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 2.25 19.88 – 2.25	Depositor EDS
% Data completeness (in resolution range)	95.9 (19.88-2.25) 95.8 (19.88-2.25)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.18 (at 2.24Å)	Xtriage
Refinement program	CNS, CNX 2000.1	Depositor
R, R_{free}	0.228 , 0.262 0.226 , 0.260	Depositor DCC
R_{free} test set	7157 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10173	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: 2T6

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	0.40	0/2484	0.90	5/3354 (0.1%)
1	B	0.40	0/2477	0.94	15/3344 (0.4%)
1	C	0.41	0/2477	0.95	10/3344 (0.3%)
1	D	0.42	0/2468	0.92	6/3333 (0.2%)
All	All	0.41	0/9906	0.93	36/13375 (0.3%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	278	LEU	N-CA-C	7.57	119.17	111.07
1	B	278	LEU	N-CA-C	7.48	119.43	111.28
1	C	270	SER	CA-C-N	7.40	128.05	119.47
1	C	270	SER	C-N-CA	7.40	128.05	119.47
1	D	278	LEU	N-CA-C	7.27	118.84	111.07
1	A	278	LEU	N-CA-C	7.09	118.66	111.07
1	D	297	THR	N-CA-C	-6.91	104.85	113.28
1	A	297	THR	N-CA-C	-6.88	104.89	113.28
1	C	297	THR	N-CA-C	-6.86	104.91	113.28
1	B	297	THR	N-CA-C	-6.85	104.92	113.28
1	B	232	PHE	CA-C-N	6.72	127.30	120.38
1	B	232	PHE	C-N-CA	6.72	127.30	120.38
1	C	264	TYR	CA-C-N	6.21	126.67	120.52
1	C	264	TYR	C-N-CA	6.21	126.67	120.52
1	A	144	THR	N-CA-C	-6.16	105.56	114.12
1	B	233	PRO	N-CA-C	6.01	118.03	110.70
1	B	264	TYR	CA-C-N	5.99	126.45	120.52
1	B	264	TYR	C-N-CA	5.99	126.45	120.52
1	B	233	PRO	CA-C-N	5.99	125.54	119.19
1	B	233	PRO	C-N-CA	5.99	125.54	119.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	220	ASP	N-CA-C	-5.98	101.83	110.08
1	B	118	ASP	CA-C-N	5.72	125.19	119.24
1	B	118	ASP	C-N-CA	5.72	125.19	119.24
1	B	220	ASP	N-CA-C	-5.67	102.26	110.08
1	C	268	LYS	N-CA-C	-5.55	105.38	111.82
1	A	220	ASP	N-CA-C	-5.52	102.47	110.08
1	B	257	VAL	N-CA-C	5.27	115.48	110.53
1	D	270	SER	CA-C-N	5.27	125.58	119.47
1	D	270	SER	C-N-CA	5.27	125.58	119.47
1	C	50	LYS	N-CA-C	5.23	118.96	112.58
1	A	257	VAL	N-CA-C	5.12	115.35	110.53
1	C	281	CYS	N-CA-C	5.08	116.67	111.03
1	B	281	CYS	N-CA-C	5.07	116.65	111.03
1	D	257	VAL	N-CA-C	5.03	115.26	110.53
1	C	264	TYR	N-CA-C	-5.02	101.53	108.76
1	B	277	LEU	N-CA-C	5.01	116.43	111.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2441	0	2491	59	0
1	B	2434	0	2482	56	0
1	C	2434	0	2482	49	0
1	D	2425	0	2469	62	0
2	A	21	0	9	1	0
2	B	21	0	9	0	0
2	C	21	0	9	0	0
2	D	21	0	9	1	0
3	A	60	0	0	2	0
3	B	72	0	0	0	0
3	C	102	0	0	0	0
3	D	121	0	0	2	0
All	All	10173	0	9960	206	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 10.

All (206) 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:A:267:ASN:ND2	1:A:270:SER:H	1.53	1.06
1:B:252:VAL:HG11	1:B:284:MET:HE2	1.33	1.04
1:B:284:MET:HA	1:B:284:MET:HE3	1.43	0.98
1:D:82:MET:HA	1:D:82:MET:HE2	1.52	0.92
1:A:32:GLN:HE21	1:A:32:GLN:HA	1.35	0.91
1:A:267:ASN:HD21	1:A:270:SER:H	1.16	0.88
1:B:32:GLN:HE21	1:B:32:GLN:HA	1.41	0.84
1:C:125:ASN:HA	1:D:258:TYR:OH	1.80	0.81
1:D:32:GLN:HE21	1:D:32:GLN:HA	1.46	0.80
1:C:230:LYS:HE3	1:C:230:LYS:HA	1.64	0.80
1:D:82:MET:HE3	1:D:117:PHE:CZ	2.16	0.80
1:A:72:LYS:HA	1:A:72:LYS:HE3	1.65	0.79
1:D:82:MET:HE1	1:D:94:LEU:HB3	1.63	0.78
1:B:164:TYR:CE1	1:B:284:MET:HE1	2.21	0.76
1:B:252:VAL:CG1	1:B:284:MET:HE2	2.15	0.75
1:D:105:VAL:O	1:D:110:ARG:HD3	1.88	0.74
1:A:267:ASN:C	1:A:267:ASN:HD22	1.95	0.73
1:D:82:MET:HE3	1:D:117:PHE:HZ	1.53	0.72
1:B:216:ALA:HA	1:B:219:PHE:CD2	2.29	0.68
1:C:48:VAL:HA	1:C:73:LEU:HD21	1.76	0.67
1:A:230:LYS:HA	1:A:230:LYS:HE2	1.77	0.67
1:D:212:ASN:HD22	1:D:214:GLY:H	1.44	0.66
1:B:121:ASP:OD2	1:B:249:VAL:HG23	1.98	0.63
1:A:267:ASN:HD21	1:A:270:SER:N	1.94	0.62
1:A:118:ASP:HA	3:A:359:HOH:O	1.98	0.62
1:B:191:GLY:HA3	1:D:191:GLY:HA3	1.81	0.62
1:B:207:LYS:HB2	1:B:207:LYS:NZ	2.15	0.62
1:D:212:ASN:ND2	1:D:214:GLY:H	1.96	0.62
1:D:82:MET:CE	1:D:94:LEU:HD13	2.31	0.61
1:A:279:TYR:OH	1:A:311:HIS:HD2	1.84	0.61
1:D:48:VAL:HG22	1:D:73:LEU:HD21	1.82	0.61
1:D:274:LYS:NZ	1:D:274:LYS:HB3	2.15	0.60
1:B:13:LEU:HD22	1:B:173:LEU:HD23	1.83	0.60
1:D:82:MET:HE1	1:D:94:LEU:CB	2.29	0.59
1:D:107:PRO:HA	1:D:110:ARG:HG3	1.82	0.59
1:B:267:ASN:OD1	1:B:269:LYS:HG3	2.01	0.59
1:C:266:ALA:HA	1:C:273:GLY:HA2	1.85	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:ASN:HD21	1:B:269:LYS:NZ	2.00	0.59
1:D:228:GLN:HE21	1:D:228:GLN:HA	1.68	0.59
1:B:117:PHE:CE2	1:B:119:PRO:HG3	2.38	0.58
1:D:216:ALA:HA	1:D:219:PHE:CD2	2.38	0.58
1:C:10:VAL:CG1	1:D:57:TYR:O	2.51	0.58
1:A:100:LYS:O	1:A:311:HIS:HE1	1.87	0.58
1:C:242:ALA:H	1:D:212:ASN:HD21	1.51	0.57
1:A:15:ARG:HG3	1:C:32:GLN:HE22	1.69	0.57
1:B:221:PRO:HB2	1:B:334:HIS:CD2	2.40	0.57
1:A:267:ASN:ND2	1:A:267:ASN:C	2.61	0.57
1:B:164:TYR:HE1	1:B:284:MET:HE1	1.69	0.57
1:C:10:VAL:HG12	1:D:57:TYR:O	2.05	0.56
1:B:143:SER:HB2	1:B:145:ASP:OD1	2.05	0.56
1:B:195:LEU:HD21	1:B:198:LYS:HG2	1.88	0.56
1:D:251:ASP:OD1	1:D:254:ARG:NH2	2.38	0.56
1:D:318:LEU:HD12	1:D:318:LEU:C	2.31	0.55
1:C:229:ARG:HB3	1:C:229:ARG:NH1	2.20	0.55
1:D:212:ASN:HD22	1:D:212:ASN:C	2.14	0.55
1:D:254:ARG:HD2	3:D:367:HOH:O	2.05	0.55
1:B:218:ASP:HA	1:B:268:LYS:HE2	1.89	0.55
1:A:216:ALA:HA	1:A:219:PHE:CD2	2.42	0.54
1:B:117:PHE:HE2	1:B:119:PRO:HG3	1.72	0.54
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.90	0.54
1:A:32:GLN:HA	1:A:32:GLN:NE2	2.16	0.54
1:A:166:LEU:HD13	1:A:249:VAL:HG12	1.88	0.54
1:C:321:PRO:O	1:C:325:LEU:HD22	2.08	0.54
1:B:220:ASP:OD1	1:B:221:PRO:HD2	2.08	0.54
1:B:218:ASP:O	1:B:267:ASN:HB2	2.08	0.54
1:A:318:LEU:C	1:A:318:LEU:HD12	2.33	0.54
1:A:30:LEU:HD23	2:A:401:2T6:C9	2.38	0.53
1:D:294:MET:SD	1:D:325:LEU:HD13	2.49	0.53
1:D:278:LEU:HD22	3:D:354:HOH:O	2.08	0.53
1:B:221:PRO:HB2	1:B:334:HIS:HD2	1.73	0.52
1:C:182:ASN:ND2	1:C:199:ASP:H	2.07	0.52
1:B:284:MET:HA	1:B:284:MET:CE	2.27	0.52
1:C:229:ARG:HB3	1:C:229:ARG:HH11	1.74	0.52
1:C:90:ALA:C	1:C:111:GLY:HA3	2.35	0.52
1:D:173:LEU:C	1:D:173:LEU:HD13	2.35	0.52
1:B:318:LEU:C	1:B:318:LEU:HD12	2.35	0.51
1:C:230:LYS:NZ	1:C:326:GLU:OE2	2.43	0.51
1:C:182:ASN:HD22	1:C:198:LYS:HA	1.75	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:GLN:HA	1:B:32:GLN:NE2	2.18	0.51
1:D:9:ASP:OD1	1:D:9:ASP:N	2.43	0.51
1:D:145:ASP:O	1:D:146:GLU:C	2.52	0.51
1:A:125:ASN:O	1:A:128:CYS:SG	2.69	0.51
1:C:125:ASN:O	1:C:130:VAL:HG12	2.10	0.51
1:D:82:MET:HE1	1:D:94:LEU:HD13	1.92	0.51
1:C:229:ARG:NH1	1:C:326:GLU:OE2	2.44	0.51
1:D:182:ASN:HD22	1:D:198:LYS:HA	1.76	0.50
1:D:82:MET:HE3	1:D:117:PHE:CE1	2.46	0.50
1:A:297:THR:HB	1:A:315:PRO:HG2	1.93	0.50
1:D:121:ASP:N	1:D:133:GLY:O	2.39	0.50
1:A:13:LEU:HD22	1:A:173:LEU:HD13	1.93	0.50
1:C:269:LYS:C	1:C:271:PRO:HD3	2.37	0.50
1:C:195:LEU:HD21	1:C:198:LYS:HG2	1.93	0.49
1:C:11:ASN:HD22	1:C:12:THR:H	1.58	0.49
1:D:230:LYS:HA	1:D:230:LYS:HE2	1.94	0.49
1:B:90:ALA:C	1:B:111:GLY:HA3	2.38	0.49
1:B:186:LEU:O	1:B:188:PRO:HD3	2.13	0.49
1:A:182:ASN:HD22	1:A:198:LYS:HA	1.77	0.48
1:A:100:LYS:O	1:A:311:HIS:CE1	2.65	0.48
1:A:225:GLU:OE2	1:A:334:HIS:HE1	1.97	0.48
1:A:267:ASN:HD22	1:A:270:SER:H	1.50	0.48
1:A:13:LEU:CD2	1:A:173:LEU:HD13	2.42	0.48
1:D:297:THR:HB	1:D:315:PRO:HG2	1.95	0.48
1:C:278:LEU:HD22	1:C:310:ILE:HA	1.96	0.48
1:A:72:LYS:HA	1:A:72:LYS:CE	2.41	0.48
1:A:267:ASN:ND2	1:A:270:SER:N	2.38	0.48
1:C:154:GLN:HE21	1:C:158:ASN:HD22	1.62	0.47
1:C:318:LEU:C	1:C:318:LEU:HD12	2.39	0.47
1:C:296:THR:HG21	1:C:328:LEU:HD21	1.96	0.47
1:A:182:ASN:ND2	1:A:199:ASP:H	2.11	0.47
1:D:82:MET:HE2	1:D:94:LEU:HD13	1.95	0.47
1:B:182:ASN:HD22	1:B:198:LYS:HA	1.79	0.47
1:C:242:ALA:H	1:D:212:ASN:ND2	2.13	0.47
1:D:32:GLN:HA	1:D:32:GLN:NE2	2.24	0.47
1:C:229:ARG:HH11	1:C:229:ARG:CB	2.28	0.47
1:A:72:LYS:HB3	1:A:74:ASP:OD2	2.15	0.47
1:D:90:ALA:C	1:D:111:GLY:HA3	2.39	0.47
1:D:138:ILE:O	1:D:160:VAL:HG22	2.14	0.47
1:B:55:HIS:HA	1:B:59:ILE:HG22	1.96	0.46
1:A:97:GLU:O	1:A:97:GLU:HG2	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:LEU:HD23	1:D:126:ILE:CD1	2.45	0.46
1:D:183:CYS:HB2	1:D:197:ASP:HB2	1.97	0.46
1:B:226:TYR:HA	1:B:229:ARG:HG3	1.98	0.46
1:A:251:ASP:OD1	1:A:254:ARG:NH2	2.47	0.46
1:C:121:ASP:N	1:C:133:GLY:O	2.42	0.46
1:D:106:GLU:O	1:D:110:ARG:HG2	2.16	0.46
1:D:82:MET:HA	1:D:82:MET:CE	2.33	0.46
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.51	0.46
1:D:223:VAL:O	1:D:227:ILE:HG12	2.16	0.46
1:D:261:ILE:HD11	1:D:317:ILE:CG2	2.46	0.46
1:D:182:ASN:ND2	1:D:199:ASP:H	2.14	0.45
1:C:258:TYR:OH	1:D:125:ASN:HA	2.16	0.45
1:C:139:TYR:OH	1:C:156:GLY:HA2	2.16	0.45
1:A:145:ASP:OD2	1:A:145:ASP:N	2.49	0.45
1:C:10:VAL:HG13	1:D:57:TYR:O	2.17	0.45
1:C:223:VAL:O	1:C:227:ILE:HG12	2.17	0.45
1:A:230:LYS:HE2	1:A:230:LYS:CA	2.46	0.45
1:C:53:ILE:HG12	1:C:53:ILE:O	2.16	0.45
1:D:228:GLN:HA	1:D:228:GLN:NE2	2.30	0.45
1:A:32:GLN:HE21	1:A:32:GLN:CA	2.16	0.45
1:C:328:LEU:O	1:C:332:GLU:HG3	2.17	0.45
1:A:231:LYS:HE2	1:B:216:ALA:HB2	1.99	0.44
1:A:296:THR:HG21	1:A:328:LEU:HD21	1.99	0.44
1:B:97:GLU:HB2	1:B:279:TYR:CE1	2.52	0.44
1:C:263:LEU:HD12	1:C:263:LEU:N	2.32	0.44
1:B:138:ILE:O	1:B:160:VAL:HG22	2.16	0.44
1:A:15:ARG:HG3	1:C:32:GLN:NE2	2.32	0.44
1:A:90:ALA:C	1:A:111:GLY:HA3	2.42	0.44
1:B:33:LEU:C	1:B:33:LEU:HD23	2.43	0.44
1:B:261:ILE:HD11	1:B:317:ILE:CG2	2.47	0.44
1:A:276:ARG:N	1:A:276:ARG:HD2	2.33	0.44
1:C:186:LEU:HB2	1:C:193:PHE:CE1	2.53	0.44
1:A:225:GLU:OE2	1:A:334:HIS:CE1	2.71	0.44
1:B:23:LYS:HA	1:B:25:ARG:NH1	2.33	0.44
1:A:231:LYS:HE2	1:B:216:ALA:CB	2.48	0.43
1:A:125:ASN:O	1:A:130:VAL:HG12	2.18	0.43
1:B:233:PRO:HA	1:B:234:PRO:HD3	1.67	0.43
1:C:74:ASP:OD1	1:C:75:VAL:N	2.50	0.43
1:A:88:SER:O	1:A:89:PHE:HB2	2.19	0.43
1:B:74:ASP:OD1	1:B:75:VAL:N	2.52	0.43
1:B:225:GLU:OE1	1:B:334:HIS:HE1	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:PRO:HA	1:A:234:PRO:HD3	1.88	0.43
1:B:213:GLU:OE2	1:B:231:LYS:NZ	2.43	0.43
1:C:156:GLY:HA3	1:C:303:LEU:HD12	2.01	0.43
1:A:96:SER:HB2	1:A:117:PHE:CZ	2.53	0.43
1:A:109:LYS:HZ3	1:C:9:ASP:CG	2.27	0.43
1:C:51:ALA:HA	1:D:188:PRO:HD2	2.01	0.43
1:D:154:GLN:HE21	1:D:158:ASN:HD22	1.65	0.43
1:B:267:ASN:HD21	1:B:269:LYS:HZ2	1.66	0.43
1:A:51:ALA:HA	1:B:188:PRO:HD2	2.00	0.42
1:D:33:LEU:C	1:D:33:LEU:HD23	2.44	0.42
1:A:90:ALA:HA	1:A:111:GLY:HA3	2.01	0.42
1:B:13:LEU:CD2	1:B:173:LEU:HD23	2.48	0.42
1:D:30:LEU:HD23	2:D:401:2T6:C9	2.49	0.42
1:A:209:TYR:CZ	1:A:242:ALA:HB2	2.54	0.42
1:A:309:ASP:HB3	1:A:312:GLN:HB3	2.00	0.42
1:C:44:ILE:O	1:C:48:VAL:HG23	2.20	0.42
1:D:186:LEU:HB2	1:D:193:PHE:CE1	2.54	0.42
1:A:258:TYR:OH	1:B:125:ASN:HA	2.19	0.42
1:B:203:LYS:HE3	1:B:203:LYS:HB2	1.79	0.42
1:B:22:ARG:HA	1:B:22:ARG:HD3	1.88	0.42
1:C:183:CYS:HB2	1:C:197:ASP:HB2	2.02	0.42
1:D:22:ARG:HG2	1:D:22:ARG:HH11	1.84	0.42
1:A:278:LEU:HD13	3:A:339:HOH:O	2.20	0.42
1:B:264:TYR:CZ	1:B:274:LYS:HD3	2.55	0.42
1:C:258:TYR:HE1	1:D:128:CYS:HB3	1.85	0.42
1:A:196:VAL:HG12	1:B:57:TYR:CZ	2.55	0.42
1:B:164:TYR:CD1	1:B:284:MET:HE1	2.55	0.42
1:B:82:MET:O	1:B:86:LYS:HG3	2.20	0.41
1:C:217:LYS:HA	1:D:232:PHE:CE2	2.55	0.41
1:B:236:ASN:O	1:B:236:ASN:CG	2.62	0.41
1:D:325:LEU:HD12	1:D:325:LEU:HA	1.90	0.41
1:B:146:GLU:HG2	1:B:147:PRO:HD2	2.03	0.41
1:C:230:LYS:HE3	1:C:230:LYS:CA	2.43	0.41
1:C:198:LYS:O	1:C:199:ASP:C	2.64	0.41
1:B:216:ALA:HA	1:B:219:PHE:CE2	2.56	0.41
1:B:277:LEU:HD23	1:B:307:PRO:HB3	2.03	0.40
1:A:33:LEU:C	1:A:33:LEU:HD23	2.46	0.40
1:A:225:GLU:OE1	1:A:333:LYS:HE2	2.21	0.40
1:B:207:LYS:HB2	1:B:207:LYS:HZ2	1.86	0.40
1:D:198:LYS:O	1:D:199:ASP:C	2.65	0.40
1:D:277:LEU:HD23	1:D:307:PRO:HB3	2.03	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:GLU:O	1:C:229:ARG:HG3	2.21	0.40
1:C:230:LYS:HA	1:C:230:LYS:CE	2.44	0.40
1:A:276:ARG:O	1:A:280:GLU:HB2	2.21	0.40
1:A:285:ALA:CB	1:A:303:LEU:HD13	2.52	0.40
1:D:233:PRO:HA	1:D:234:PRO:HD2	1.97	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	315/337 (94%)	302 (96%)	13 (4%)	0	100	100
1	B	314/337 (93%)	307 (98%)	7 (2%)	0	100	100
1	C	314/337 (93%)	307 (98%)	7 (2%)	0	100	100
1	D	313/337 (93%)	305 (97%)	8 (3%)	0	100	100
All	All	1256/1348 (93%)	1221 (97%)	35 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	266/280 (95%)	253 (95%)	13 (5%)	22	24

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	265/280 (95%)	258 (97%)	7 (3%)	40	50
1	C	265/280 (95%)	260 (98%)	5 (2%)	50	61
1	D	264/280 (94%)	256 (97%)	8 (3%)	36	45
All	All	1060/1120 (95%)	1027 (97%)	33 (3%)	35	44

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASP
1	A	32	GLN
1	A	72	LYS
1	A	98	GLU
1	A	125	ASN
1	A	145	ASP
1	A	236	ASN
1	A	267	ASN
1	A	276	ARG
1	A	278	LEU
1	A	282	ASN
1	A	303	LEU
1	A	325	LEU
1	B	9	ASP
1	B	32	GLN
1	B	56	LEU
1	B	125	ASN
1	B	197	ASP
1	B	269	LYS
1	B	325	LEU
1	C	19	GLU
1	C	230	LYS
1	C	303	LEU
1	C	313	ARG
1	C	325	LEU
1	D	32	GLN
1	D	110	ARG
1	D	173	LEU
1	D	178	ASP
1	D	198	LYS
1	D	228	GLN
1	D	282	ASN
1	D	325	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	35	ASN
1	A	125	ASN
1	A	154	GLN
1	A	182	ASN
1	A	267	ASN
1	A	282	ASN
1	A	311	HIS
1	B	32	GLN
1	B	35	ASN
1	B	83	ASN
1	B	101	HIS
1	B	125	ASN
1	B	154	GLN
1	B	158	ASN
1	B	182	ASN
1	B	272	ASN
1	B	282	ASN
1	B	334	HIS
1	C	11	ASN
1	C	35	ASN
1	C	83	ASN
1	C	101	HIS
1	C	125	ASN
1	C	154	GLN
1	C	182	ASN
1	C	228	GLN
1	C	267	ASN
1	C	282	ASN
1	D	32	GLN
1	D	35	ASN
1	D	83	ASN
1	D	154	GLN
1	D	182	ASN
1	D	212	ASN
1	D	228	GLN
1	D	236	ASN
1	D	282	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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	2T6	C	401	-	22,23,23	1.69	5 (22%)	23,35,35	1.72	6 (26%)
2	2T6	A	401	-	22,23,23	1.73	7 (31%)	23,35,35	1.70	6 (26%)
2	2T6	B	401	-	22,23,23	1.72	5 (22%)	23,35,35	1.71	6 (26%)
2	2T6	D	401	-	22,23,23	1.71	3 (13%)	23,35,35	1.67	6 (26%)

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	2T6	C	401	-	-	2/10/18/18	0/3/3/3
2	2T6	A	401	-	-	3/10/18/18	0/3/3/3
2	2T6	B	401	-	-	4/10/18/18	0/3/3/3
2	2T6	D	401	-	-	1/10/18/18	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	2T6	C11-N12	3.60	1.39	1.30
2	D	401	2T6	C7-C1	3.53	1.56	1.50
2	A	401	2T6	C11-N12	3.46	1.39	1.30
2	D	401	2T6	C11-N12	3.46	1.39	1.30
2	B	401	2T6	C11-N12	3.43	1.39	1.30
2	D	401	2T6	C8-C9	-3.36	1.33	1.36
2	B	401	2T6	C8-C9	-3.36	1.33	1.36
2	A	401	2T6	C7-C1	3.33	1.56	1.50
2	C	401	2T6	C7-C1	3.30	1.56	1.50
2	A	401	2T6	C8-C9	-3.30	1.33	1.36
2	B	401	2T6	C7-C1	3.22	1.55	1.50
2	C	401	2T6	C8-C9	-3.10	1.33	1.36
2	A	401	2T6	C6-C16	-2.31	1.46	1.50
2	B	401	2T6	C6-C16	-2.18	1.47	1.50
2	A	401	2T6	P19-O20	-2.14	1.50	1.55
2	C	401	2T6	P19-O21	-2.06	1.50	1.55
2	A	401	2T6	P19-O21	-2.06	1.50	1.55
2	A	401	2T6	C2-C3	-2.05	1.37	1.40
2	B	401	2T6	P19-O21	-2.03	1.50	1.55
2	C	401	2T6	P19-O20	-2.03	1.50	1.55

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	2T6	S10-C11-N12	-4.35	112.34	116.08
2	C	401	2T6	S10-C11-N12	-4.30	112.39	116.08
2	A	401	2T6	S10-C11-N12	-4.23	112.45	116.08
2	D	401	2T6	S10-C11-N12	-4.13	112.53	116.08
2	C	401	2T6	C9-C8-S10	3.19	111.19	109.47
2	D	401	2T6	C9-C8-S10	3.18	111.19	109.47
2	C	401	2T6	O18-C16-C6	3.11	123.91	120.24
2	A	401	2T6	C9-C8-S10	3.05	111.12	109.47
2	B	401	2T6	O18-C16-C6	3.02	123.80	120.24
2	A	401	2T6	O18-C16-C6	2.98	123.75	120.24
2	D	401	2T6	O18-C16-C6	2.91	123.67	120.24
2	B	401	2T6	C9-C8-S10	2.90	111.04	109.47
2	A	401	2T6	C3-C2-C9	-2.65	129.10	133.41
2	B	401	2T6	C3-C2-C9	-2.60	129.19	133.41
2	D	401	2T6	C3-C2-C9	-2.58	129.22	133.41
2	A	401	2T6	C6-C16-N17	-2.48	114.43	118.29
2	B	401	2T6	C6-C16-N17	-2.44	114.50	118.29
2	C	401	2T6	C3-C2-C9	-2.38	129.54	133.41

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	2T6	C6-C16-N17	-2.37	114.61	118.29
2	D	401	2T6	C6-C16-N17	-2.22	114.85	118.29
2	A	401	2T6	C2-C9-C8	-2.19	107.74	111.44
2	B	401	2T6	C2-C9-C8	-2.18	107.77	111.44
2	C	401	2T6	C2-C9-C8	-2.12	107.86	111.44
2	D	401	2T6	C2-C9-C8	-2.04	107.99	111.44

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	2T6	O13-C14-P19-O20
2	B	401	2T6	O13-C14-P19-O15
2	B	401	2T6	C2-C3-O13-C14
2	C	401	2T6	C2-C3-O13-C14
2	D	401	2T6	C2-C3-O13-C14
2	A	401	2T6	C2-C3-O13-C14
2	A	401	2T6	N17-C16-C6-C1
2	A	401	2T6	O18-C16-C6-C1
2	C	401	2T6	O18-C16-C6-C1
2	B	401	2T6	O13-C14-P19-O21

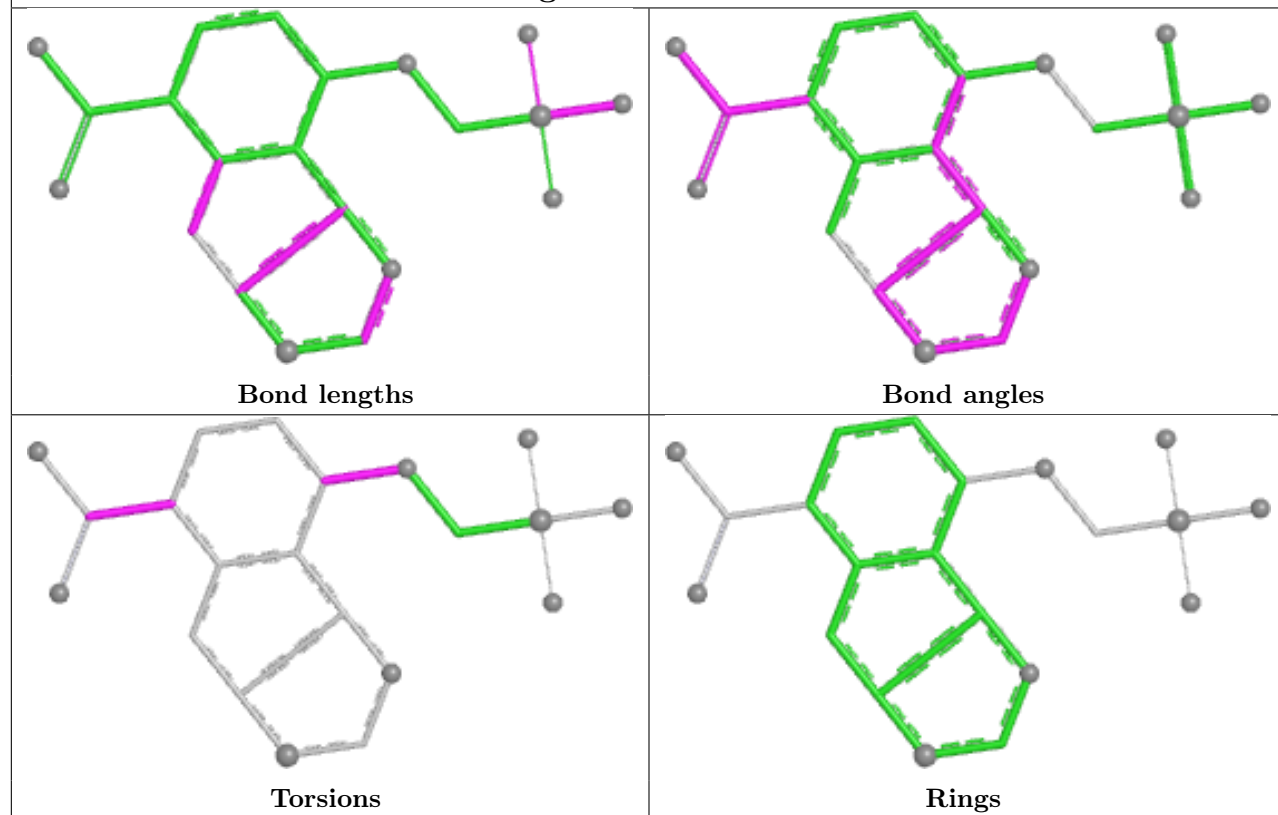
There are no ring outliers.

2 monomers are involved in 2 short contacts:

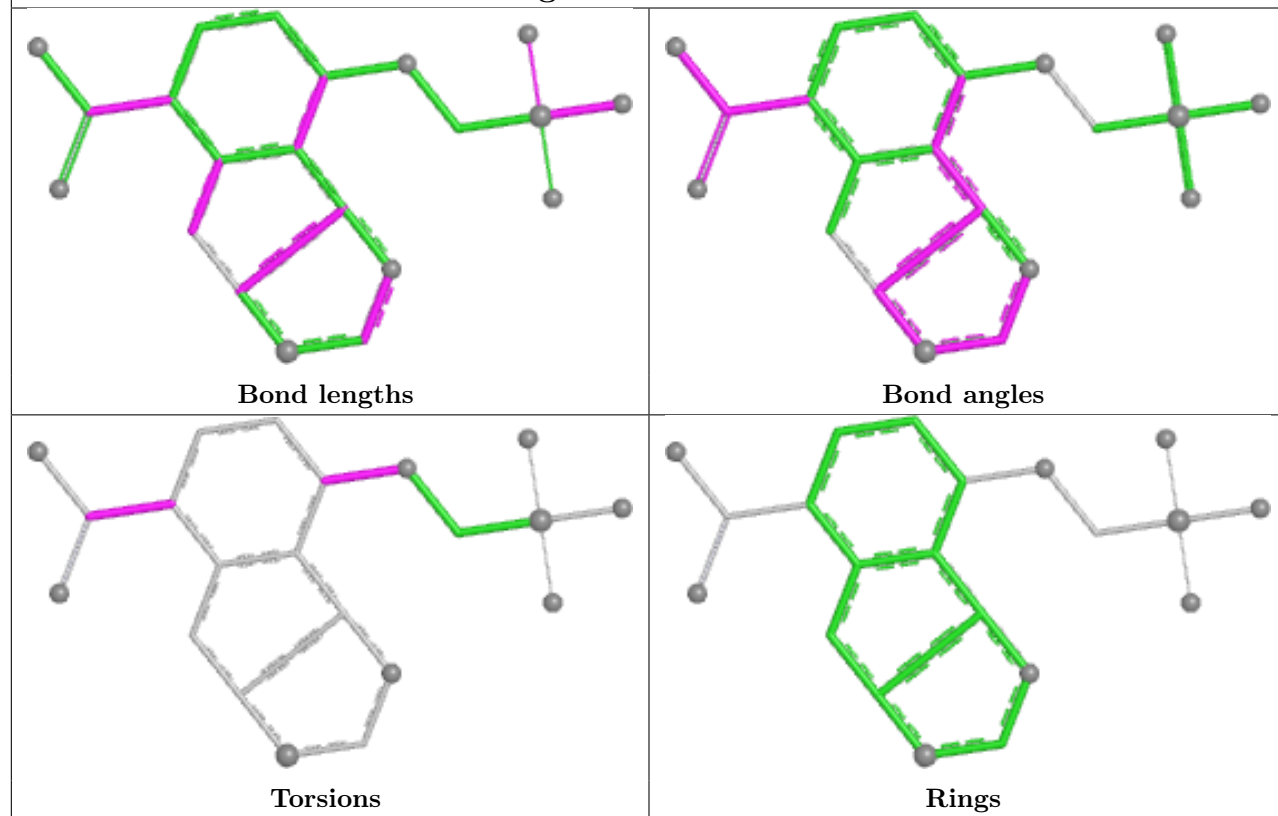
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	2T6	1	0
2	D	401	2T6	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

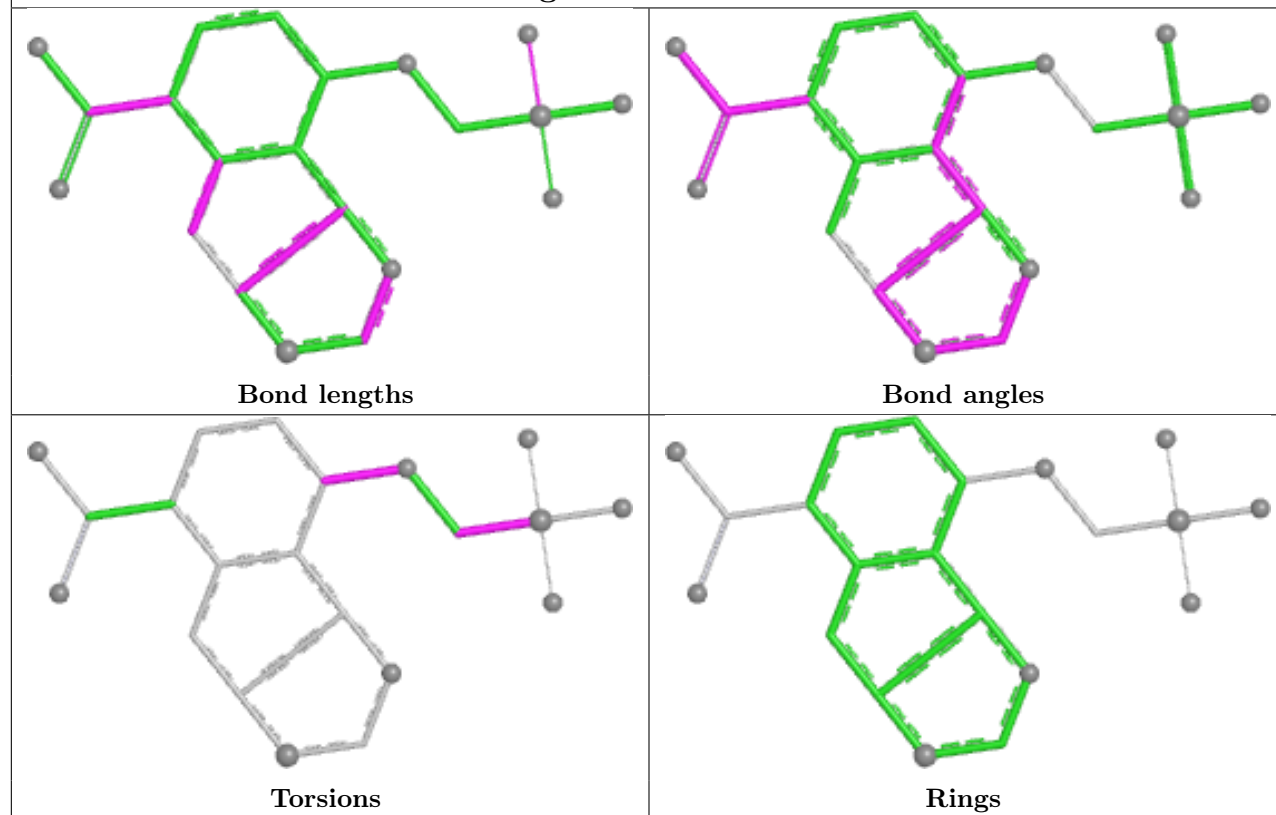
Ligand 2T6 C 401



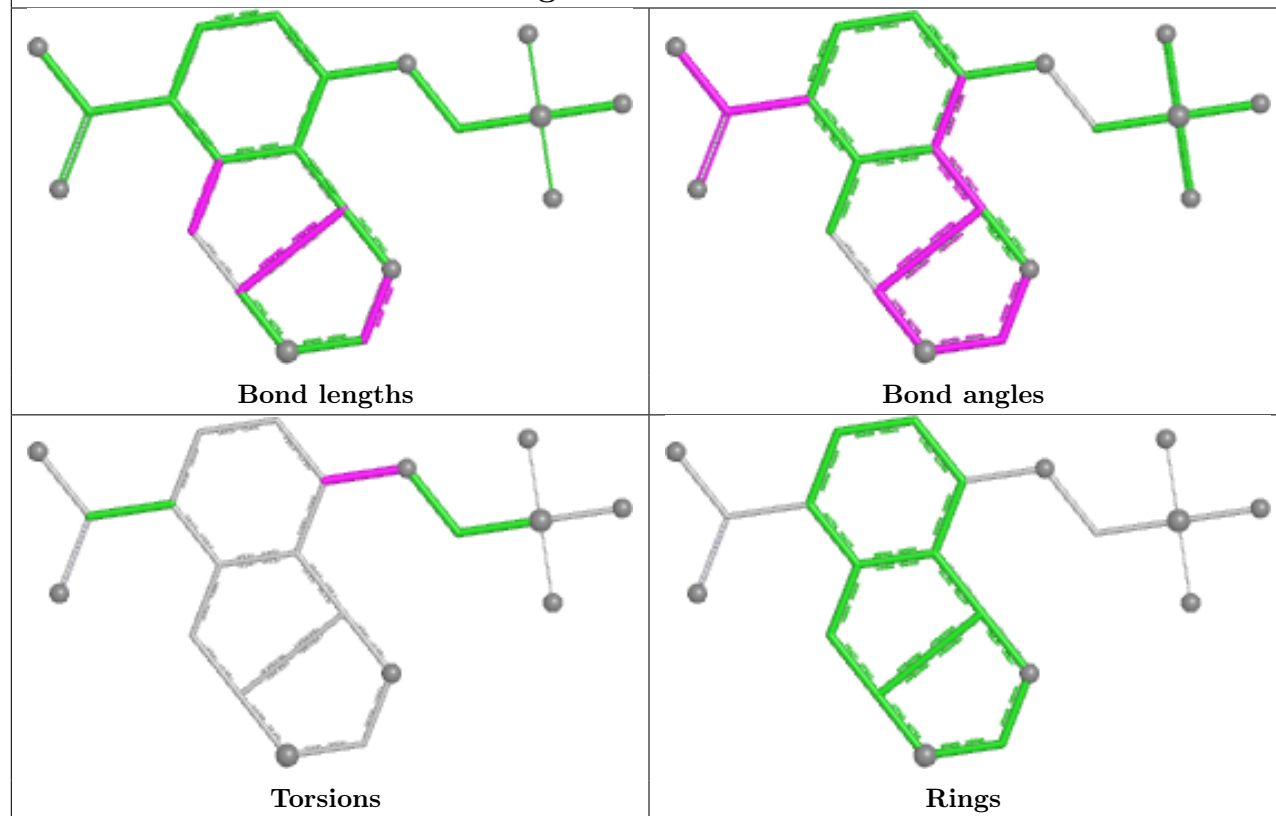
Ligand 2T6 A 401



Ligand 2T6 B 401



Ligand 2T6 D 401



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	319/337 (94%)	0.53	24 (7%)	20 18	16, 25, 42, 53	0
1	B	318/337 (94%)	0.47	18 (5%)	29 27	16, 24, 40, 50	0
1	C	318/337 (94%)	0.34	18 (5%)	29 27	15, 22, 37, 50	0
1	D	317/337 (94%)	0.21	11 (3%)	47 47	15, 22, 35, 48	0
All	All	1272/1348 (94%)	0.39	71 (5%)	30 28	15, 24, 39, 53	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	125	ASN	6.5
1	A	124	SER	5.8
1	B	126	ILE	5.8
1	C	125	ASN	5.7
1	B	125	ASN	5.6
1	B	124	SER	5.1
1	A	126	ILE	4.7
1	B	335	SER	4.6
1	A	215	TYR	4.6
1	C	124	SER	4.4
1	C	267	ASN	4.2
1	C	126	ILE	4.0
1	A	108	GLU	3.9
1	B	9	ASP	3.8
1	D	124	SER	3.7
1	D	121	ASP	3.6
1	C	127	ASP	3.5
1	B	269	LYS	3.4
1	C	146	GLU	3.2
1	D	122	GLY	3.2
1	D	9	ASP	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	99	ASP	3.0
1	A	335	SER	3.0
1	B	270	SER	2.9
1	A	144	THR	2.9
1	C	140	ARG	2.9
1	A	123	SER	2.9
1	C	271	PRO	2.9
1	A	22	ARG	2.8
1	B	242	ALA	2.8
1	B	145	ASP	2.8
1	D	125	ASN	2.8
1	A	216	ALA	2.8
1	A	234	PRO	2.7
1	B	268	LYS	2.6
1	D	123	SER	2.6
1	A	70	VAL	2.6
1	A	214	GLY	2.6
1	B	215	TYR	2.6
1	B	146	GLU	2.5
1	C	333	LYS	2.5
1	A	258	TYR	2.5
1	C	73	LEU	2.5
1	A	268	LYS	2.5
1	B	142	LYS	2.5
1	C	219	PHE	2.5
1	C	122	GLY	2.5
1	B	127	ASP	2.5
1	A	122	GLY	2.4
1	C	335	SER	2.4
1	C	123	SER	2.3
1	A	332	GLU	2.3
1	B	266	ALA	2.2
1	D	309	ASP	2.2
1	A	140	ARG	2.2
1	C	108	GLU	2.2
1	B	267	ASN	2.2
1	C	9	ASP	2.2
1	D	273	GLY	2.2
1	C	299	LYS	2.2
1	B	38	CYS	2.1
1	D	140	ARG	2.1
1	A	150	LYS	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	143	SER	2.1
1	C	106	GLU	2.1
1	B	99	ASP	2.1
1	A	146	GLU	2.0
1	D	266	ALA	2.0
1	A	233	PRO	2.0
1	D	126	ILE	2.0
1	A	269	LYS	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](#)

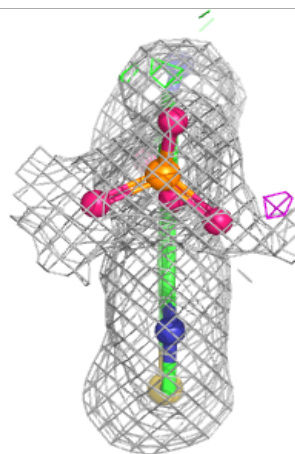
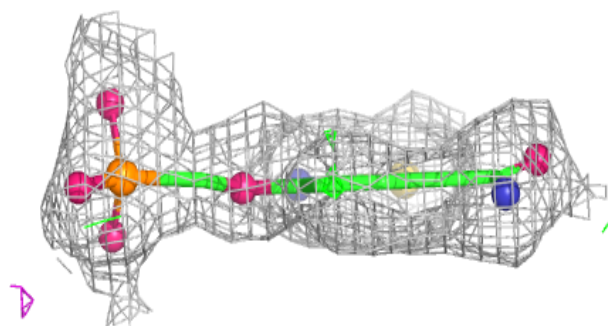
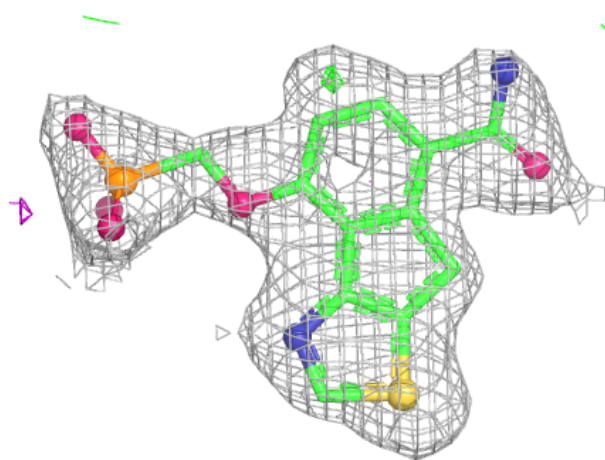
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	2T6	A	401	21/21	0.95	0.10	26,28,31,32	0
2	2T6	B	401	21/21	0.95	0.08	20,23,25,28	0
2	2T6	C	401	21/21	0.97	0.06	19,22,26,29	0
2	2T6	D	401	21/21	0.98	0.05	18,20,22,24	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.

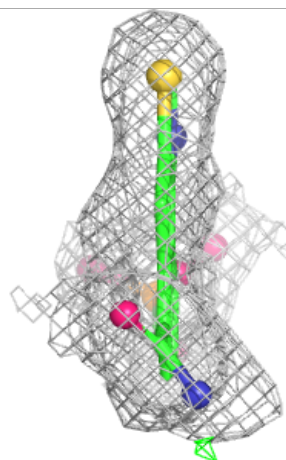
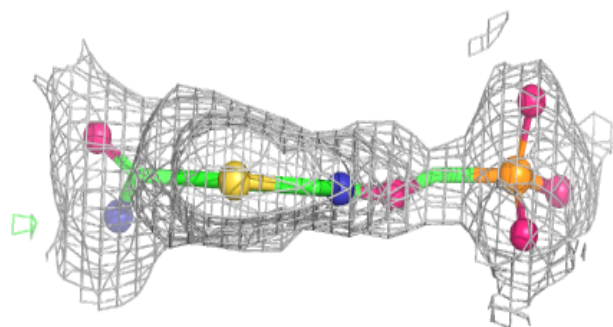
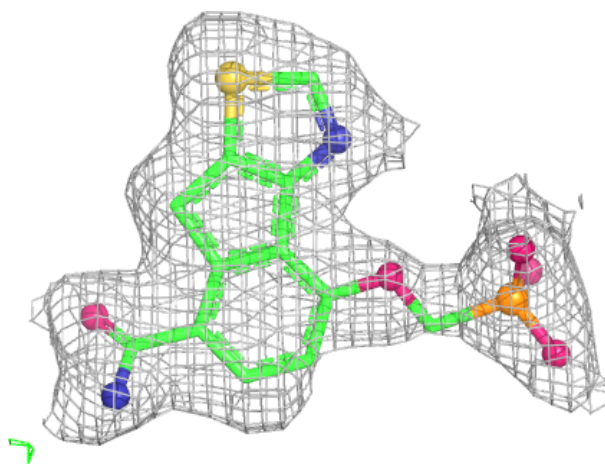
Electron density around 2T6 A 401:

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



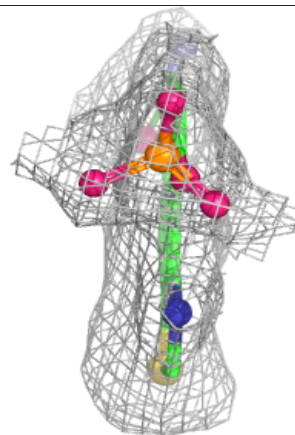
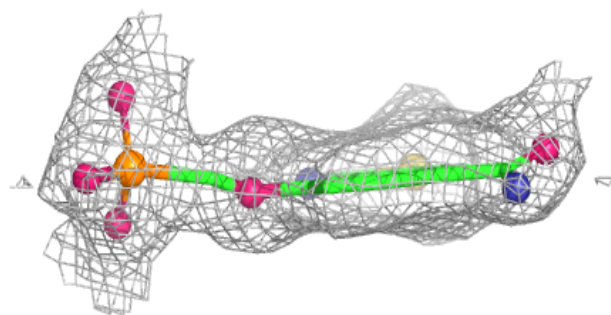
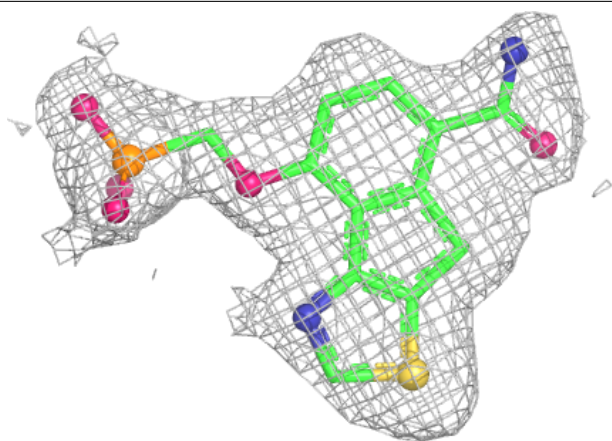
Electron density around 2T6 B 401:

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



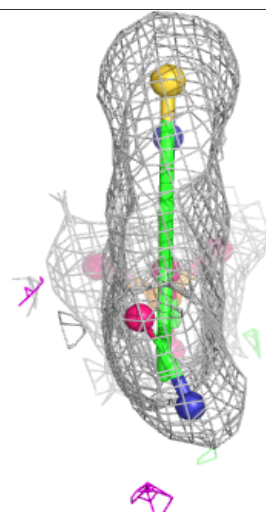
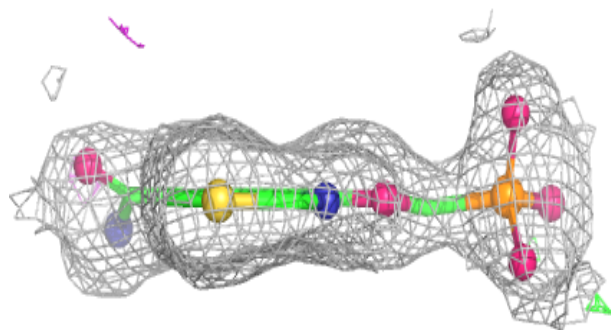
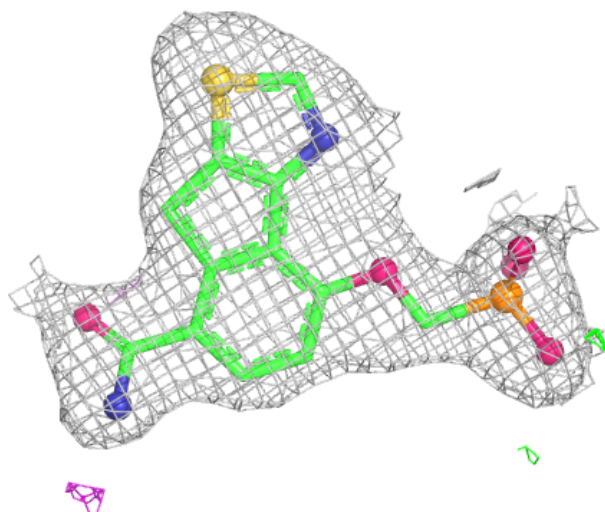
Electron density around 2T6 C 401:

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



Electron density around 2T6 D 401:

$2mF_o-DF_c$ (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.