



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:01 AM UTC

PDB ID : 7PVJ / pdb_00007pvj
Title : Crystal structure of Thioredoxin Reductase from Brugia Malayi in complex with auranofin
Authors : Fata, F.; Ardini, M.; Silvestri, I.; Gabriele, F.; Ippoliti, R.; Gencheva, R.; Cheng, Q.; Arner, E.S.J.; Angelucci, F.; Williams, D.L.
Deposited on : 2021-10-04
Resolution : 3.10 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

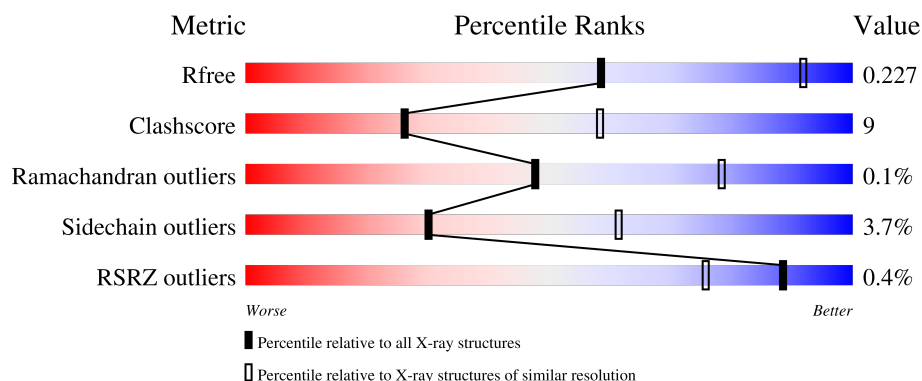
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	598	
1	B	598	
1	C	598	

2 Entry composition [i](#)

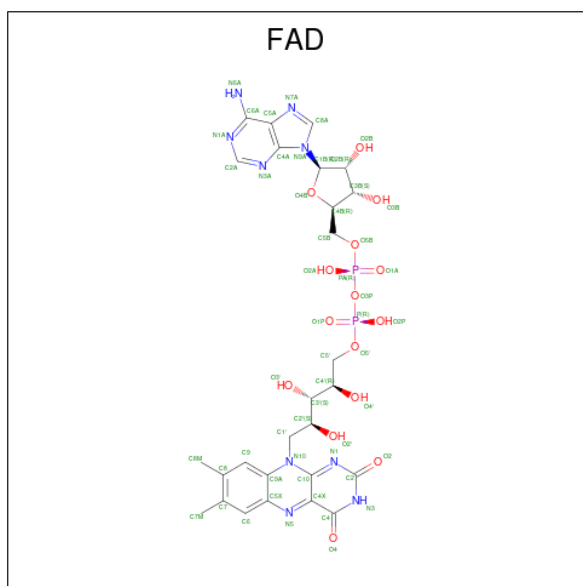
There are 6 unique types of molecules in this entry. The entry contains 13416 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutaredoxin domain-containing protein.

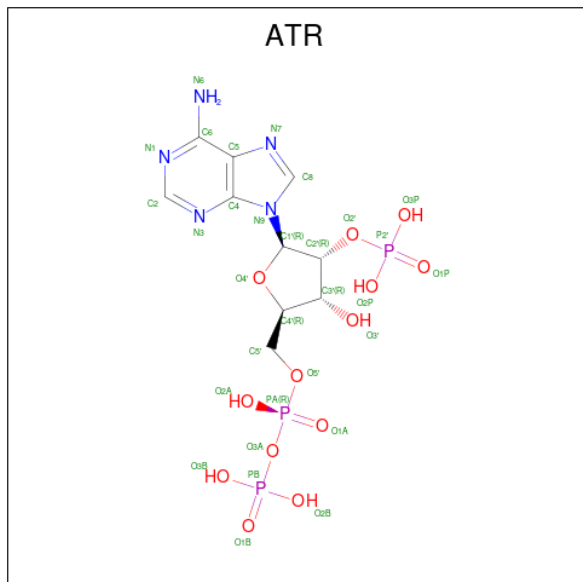
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	577	Total	C	N	O	S	0	0	0
			4484	2861	752	842	29			
1	B	577	Total	C	N	O	S	0	0	0
			4482	2860	752	841	29			
1	C	541	Total	C	N	O	S	0	0	0
			4218	2698	711	782	27			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



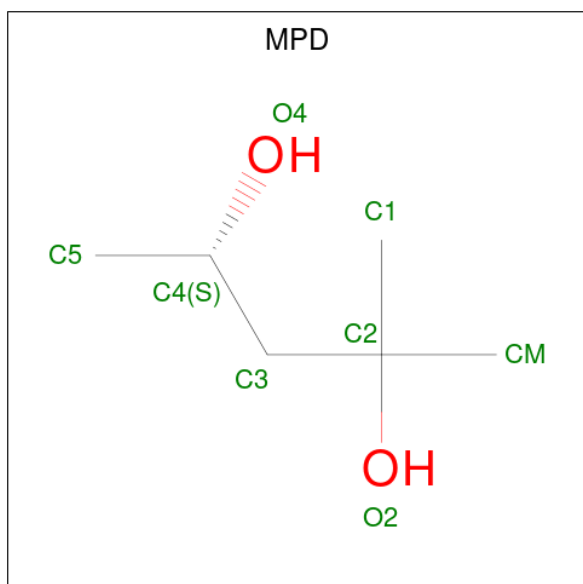
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 2'-MONOPHOSPHOADENOSINE-5'-DIPHOSPHATE (CCD ID: ATR) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	31	10	5	13	3	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	8	6	2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

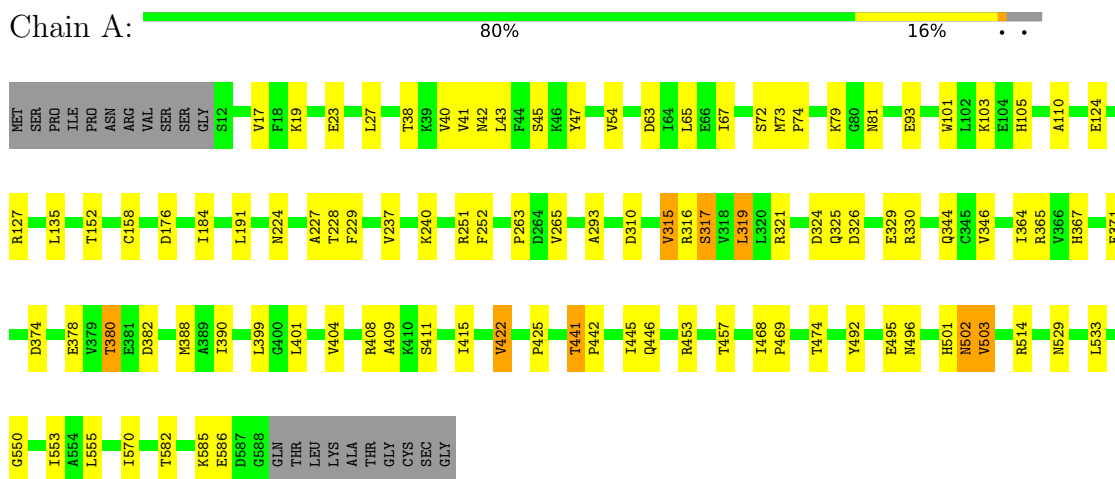
- Molecule 6 is GOLD ION (CCD ID: AU) (formula: Au) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Au	0	0
			1	1		

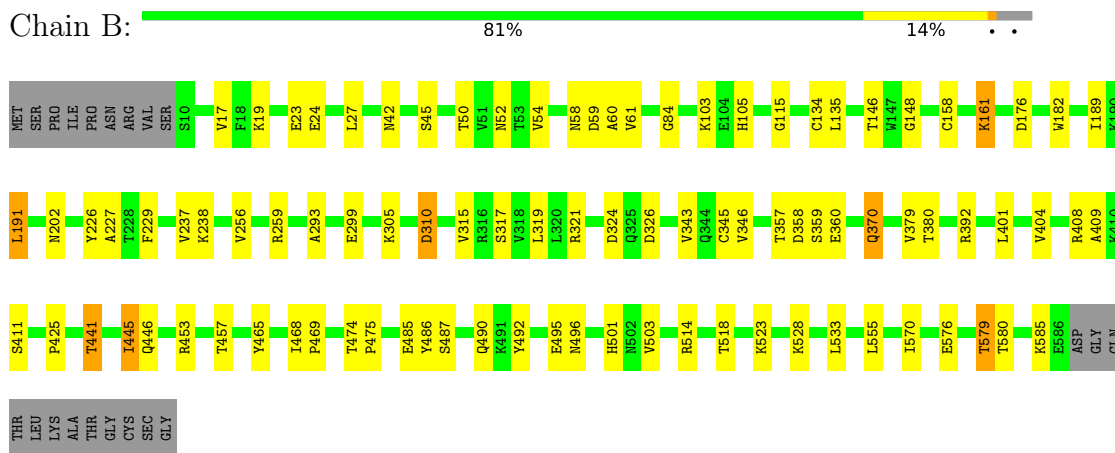
3 Residue-property plots

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: Glutaredoxin domain-containing protein

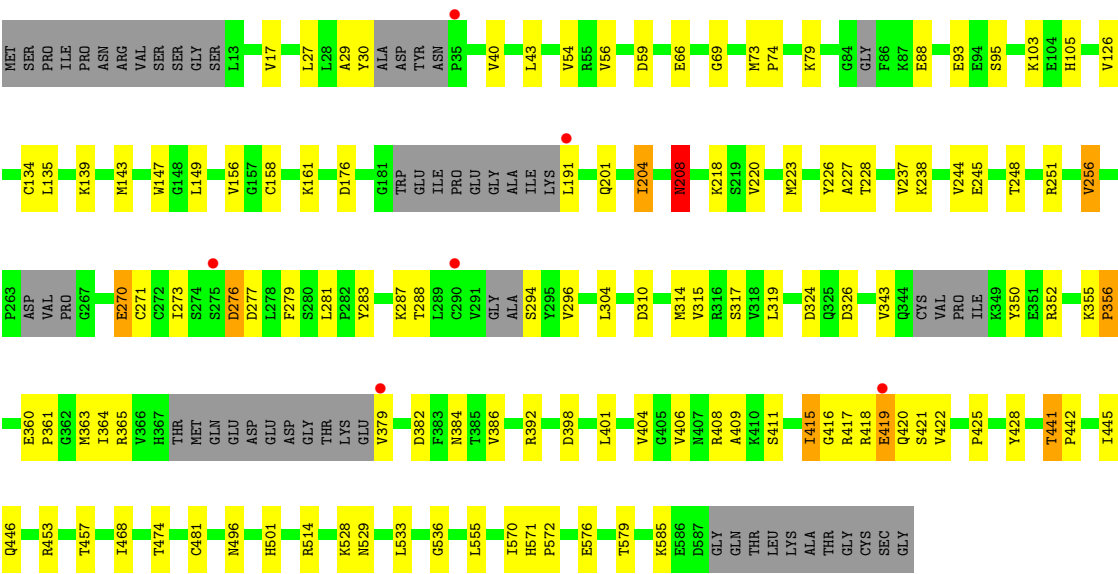


- Molecule 1: Glutaredoxin domain-containing protein



- Molecule 1: Glutaredoxin domain-containing protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	147.33Å 261.66Å 130.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.36 – 3.10 49.36 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.36-3.10) 99.9 (49.36-3.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.186 , 0.229 0.190 , 0.227	Depositor DCC
R_{free} test set	2303 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	100.7	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.012 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.014 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13416	wwPDB-VP
Average B, all atoms (Å ²)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: NA, AU, FAD, ATR, MPD

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	1.10	1/4573 (0.0%)	1.43	5/6182 (0.1%)
1	B	1.09	3/4571 (0.1%)	1.43	6/6179 (0.1%)
1	C	1.08	0/4294	1.45	9/5789 (0.2%)
All	All	1.09	4/13438 (0.0%)	1.44	20/18150 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	84	GLY	C-O	6.37	1.27	1.24
1	B	148	GLY	C-O	5.27	1.28	1.23
1	B	115	GLY	C-O	5.26	1.28	1.23
1	A	502	ASN	C-N	-5.17	1.25	1.33

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	374	ASP	CA-CB-CG	9.37	121.97	112.60
1	C	276	ASP	CA-CB-CG	7.90	120.50	112.60
1	B	343	VAL	CA-C-N	7.44	132.51	120.60
1	B	343	VAL	C-N-CA	7.44	132.51	120.60
1	C	425	PRO	N-CA-C	6.56	122.09	114.03
1	C	208	ASN	CA-CB-CG	6.45	119.05	112.60
1	C	528	LYS	CA-C-N	6.26	129.52	120.38
1	C	528	LYS	C-N-CA	6.26	129.52	120.38
1	B	425	PRO	N-CA-C	6.22	121.68	114.03
1	A	425	PRO	N-CA-C	6.13	121.56	114.03
1	C	356	PRO	N-CA-C	5.92	123.21	114.80
1	B	61	VAL	N-CA-CB	5.89	117.05	110.51
1	A	315	VAL	N-CA-C	-5.70	99.57	108.86
1	A	152	THR	CA-CB-OG1	-5.53	101.30	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	208	ASN	N-CA-CB	5.29	118.37	110.22
1	A	317	SER	N-CA-C	-5.13	99.88	110.80
1	C	256	VAL	CA-C-O	-5.10	118.19	122.63
1	B	370	GLN	CB-CG-CD	5.06	121.20	112.60
1	B	425	PRO	CB-CA-C	-5.03	103.85	111.44
1	C	317	SER	N-CA-C	5.00	121.46	110.80

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	4484	0	4511	77	0
1	B	4482	0	4512	54	0
1	C	4218	0	4261	120	0
2	A	53	0	31	0	0
2	B	53	0	31	4	0
2	C	53	0	31	4	0
3	A	31	0	11	6	0
4	A	16	0	28	3	0
4	B	24	0	42	3	0
5	A	1	0	0	0	0
6	B	1	0	0	0	0
All	All	13416	0	13458	253	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 9.

All (253) 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:390:ILE:O	3:A:602:ATR:H5'1	1.21	1.36
1:C:417:ARG:HG2	1:C:422:VAL:HG23	1.25	1.12
1:C:356:PRO:O	1:C:360:GLU:O	1.70	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:409:ALA:CA	1:C:415:ILE:HD11	1.87	1.04
1:C:501:HIS:CD2	1:C:585:LYS:HB2	1.94	1.02
1:C:156:VAL:CG1	1:C:276:ASP:HA	1.88	1.01
1:C:417:ARG:CG	1:C:422:VAL:HG23	1.92	1.00
1:A:502:ASN:OD1	1:A:503:VAL:O	1.79	0.98
1:C:156:VAL:HG11	1:C:276:ASP:HA	1.45	0.98
1:A:390:ILE:O	3:A:602:ATR:C5'	2.12	0.97
1:C:409:ALA:HA	1:C:415:ILE:HD11	1.46	0.95
1:C:501:HIS:HD2	1:C:585:LYS:HB2	1.32	0.93
1:C:126:VAL:HG21	1:C:220:VAL:HG22	1.47	0.92
1:A:293:ALA:HB2	1:A:315:VAL:HG13	1.52	0.89
1:C:156:VAL:HG11	1:C:276:ASP:CA	2.04	0.87
1:C:93:GLU:O	1:C:95:SER:O	1.95	0.85
1:A:501:HIS:CD2	1:A:585:LYS:HB2	2.12	0.85
1:C:416:GLY:HA2	1:C:422:VAL:H	1.44	0.81
1:A:501:HIS:HD2	1:A:585:LYS:HB2	1.46	0.81
1:C:273:ILE:HG22	1:C:386:VAL:O	1.81	0.81
1:C:417:ARG:HG2	1:C:422:VAL:CG2	2.08	0.80
1:C:156:VAL:HG12	1:C:156:VAL:O	1.82	0.80
1:C:294:SER:OG	1:C:296:VAL:HG22	1.81	0.79
1:A:293:ALA:HB2	1:A:315:VAL:CG1	2.10	0.79
1:C:66:GLU:O	1:C:418:ARG:NH1	2.16	0.79
1:A:317:SER:HB2	3:A:602:ATR:O2P	1.84	0.77
1:C:419:GLU:O	1:C:428:TYR:HD2	1.66	0.77
1:C:161:LYS:CD	2:C:601:FAD:H6	2.15	0.76
1:C:287:LYS:H	1:C:384:ASN:HD22	1.31	0.76
1:C:73:MET:HB2	1:C:74:PRO:HD3	1.67	0.75
1:C:416:GLY:HA2	1:C:422:VAL:N	2.00	0.75
1:C:156:VAL:HG13	1:C:276:ASP:HA	1.69	0.75
1:C:270:GLU:HB2	1:C:350:TYR:HB2	1.68	0.75
1:C:143:MET:HG2	1:C:277:ASP:CG	2.13	0.74
1:C:161:LYS:HD3	2:C:601:FAD:H6	1.68	0.74
1:C:270:GLU:OE1	1:C:271:CYS:SG	2.45	0.73
1:A:502:ASN:OD1	1:A:503:VAL:N	2.21	0.73
1:C:238:LYS:HG2	1:C:244:VAL:HG22	1.70	0.72
1:C:143:MET:HG2	1:C:277:ASP:OD1	1.90	0.72
1:C:409:ALA:N	1:C:415:ILE:HD11	2.06	0.71
1:C:404:VAL:HG13	1:C:406:VAL:HG23	1.72	0.70
1:C:415:ILE:O	1:C:422:VAL:HG12	1.92	0.70
1:B:161:LYS:HD2	1:B:299:GLU:OE1	1.92	0.70
1:B:501:HIS:NE2	1:B:585:LYS:HA	2.07	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:MET:HB2	1:A:74:PRO:HD3	1.72	0.69
1:B:357:THR:CG2	1:B:360:GLU:HB2	2.24	0.68
1:C:105:HIS:ND1	1:C:248:THR:CG2	2.57	0.68
4:A:604:MPD:H52	4:A:604:MPD:HM1	1.75	0.68
1:C:256:VAL:HG23	1:C:256:VAL:O	1.93	0.68
1:C:223:MET:HE1	1:C:245:GLU:OE2	1.95	0.67
1:C:191:LEU:HD12	1:C:191:LEU:O	1.94	0.67
1:C:201:GLN:NE2	1:C:279:PHE:O	2.28	0.67
1:A:415:ILE:HG22	1:A:422:VAL:HG21	1.75	0.66
1:C:156:VAL:CG1	1:C:276:ASP:CA	2.69	0.66
1:C:156:VAL:HG11	1:C:276:ASP:CB	2.24	0.66
1:B:453:ARG:O	1:B:457:THR:HB	1.97	0.65
4:A:604:MPD:HM1	4:A:604:MPD:C5	2.26	0.65
1:C:453:ARG:O	1:C:457:THR:HB	1.97	0.65
1:B:495:GLU:CD	1:B:495:GLU:H	2.04	0.64
1:C:139:LYS:CE	1:C:226:TYR:CE2	2.80	0.64
1:C:139:LYS:HE3	1:C:226:TYR:CE2	2.33	0.64
1:C:149:LEU:O	1:C:208:ASN:HB3	1.98	0.64
1:A:316:ARG:HB3	1:A:346:VAL:HG23	1.79	0.64
1:A:453:ARG:O	1:A:457:THR:HB	1.97	0.63
1:B:576:GLU:O	1:B:579:THR:HB	1.98	0.63
1:A:415:ILE:HG22	1:A:422:VAL:CG2	2.29	0.63
1:B:579:THR:HG22	1:B:580:THR:HG23	1.81	0.63
1:A:317:SER:CB	3:A:602:ATR:O2P	2.48	0.62
1:B:486:TYR:CE2	1:B:490:GLN:OE1	2.53	0.62
1:C:360:GLU:CD	1:C:361:PRO:HD2	2.25	0.62
1:C:419:GLU:O	1:C:428:TYR:CD2	2.52	0.62
1:C:29:ALA:O	1:C:30:TYR:HB3	2.00	0.61
1:C:501:HIS:CD2	1:C:585:LYS:CB	2.79	0.61
1:C:156:VAL:HG11	1:C:276:ASP:HB3	1.81	0.61
1:B:161:LYS:CD	1:B:299:GLU:OE1	2.50	0.60
1:A:495:GLU:OE2	1:A:495:GLU:N	2.26	0.60
1:C:135:LEU:HD23	1:C:227:ALA:HB2	1.81	0.60
1:B:27:LEU:HD11	1:B:54:VAL:HG23	1.85	0.59
1:C:416:GLY:HA3	1:C:421:SER:HA	1.84	0.59
1:A:316:ARG:HB3	1:A:346:VAL:CG2	2.32	0.59
1:A:365:ARG:NH2	1:A:382:ASP:OD1	2.36	0.59
1:B:358:ASP:O	1:B:359:SER:HB2	2.03	0.59
1:B:495:GLU:O	1:B:528:LYS:NZ	2.27	0.59
1:C:126:VAL:HG21	1:C:220:VAL:CG2	2.27	0.58
1:B:135:LEU:HD23	1:B:227:ALA:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:392:ARG:HD2	2:C:601:FAD:HM82	1.85	0.58
1:C:287:LYS:N	1:C:384:ASN:HD22	2.02	0.58
1:C:27:LEU:HD11	1:C:54:VAL:HG23	1.87	0.57
1:C:126:VAL:HG11	1:C:218:LYS:O	2.04	0.57
1:B:465:TYR:CE1	4:B:603:MPD:H11	2.39	0.57
1:A:321:ARG:HD2	3:A:602:ATR:O2B	2.06	0.56
1:C:409:ALA:CB	1:C:415:ILE:HD11	2.34	0.56
1:B:135:LEU:HD12	1:B:135:LEU:N	2.21	0.56
4:B:603:MPD:HM1	4:B:603:MPD:O4	2.06	0.56
1:B:357:THR:HG23	1:B:360:GLU:HB2	1.88	0.56
1:C:147:TRP:CH2	1:C:276:ASP:HB2	2.41	0.56
1:C:288:THR:HG21	1:C:304:LEU:HD13	1.86	0.55
1:B:134:CYS:C	1:B:135:LEU:HD12	2.31	0.55
1:B:161:LYS:NZ	2:B:601:FAD:O4	2.40	0.55
1:A:315:VAL:O	1:A:344:GLN:HA	2.06	0.55
1:C:156:VAL:CG1	1:C:156:VAL:O	2.51	0.55
4:B:604:MPD:O4	4:B:604:MPD:O2	2.22	0.54
1:A:184:ILE:HD13	1:B:189:ILE:HG21	1.89	0.54
1:A:124:GLU:OE2	1:A:127:ARG:NH1	2.41	0.54
1:A:325:GLN:O	1:A:329:GLU:HG2	2.07	0.54
1:C:126:VAL:CG2	1:C:220:VAL:HG22	2.30	0.54
1:C:73:MET:HB2	1:C:74:PRO:CD	2.37	0.54
1:A:492:TYR:O	1:A:496:ASN:ND2	2.39	0.53
1:A:27:LEU:HD11	1:A:54:VAL:HG23	1.90	0.53
1:C:270:GLU:CB	1:C:350:TYR:HB2	2.39	0.53
1:A:315:VAL:HG21	1:A:319:LEU:HD21	1.91	0.53
1:B:486:TYR:O	1:B:487:SER:C	2.50	0.53
1:C:294:SER:OG	1:C:296:VAL:CG2	2.53	0.53
1:C:134:CYS:C	1:C:135:LEU:HD12	2.33	0.53
1:B:317:SER:OG	1:B:321:ARG:NH2	2.42	0.53
1:A:365:ARG:HD3	1:A:382:ASP:OD1	2.08	0.52
1:C:415:ILE:O	1:C:422:VAL:CG1	2.57	0.52
1:A:570:ILE:HD11	1:B:469:PRO:HB2	1.90	0.52
1:C:496:ASN:HB3	1:C:529:ASN:ND2	2.25	0.52
1:A:124:GLU:CD	1:A:127:ARG:NH1	2.68	0.52
1:A:229:PHE:CD1	1:A:404:VAL:HG21	2.44	0.52
1:A:364:ILE:HD12	1:A:364:ILE:N	2.25	0.52
1:B:486:TYR:HE2	1:B:490:GLN:OE1	1.92	0.51
1:B:345:CYS:SG	1:B:370:GLN:OE1	2.64	0.51
1:A:40:VAL:O	1:A:43:LEU:HB3	2.11	0.51
1:B:404:VAL:HG12	1:B:404:VAL:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ASP:O	1:A:67:ILE:HD12	2.10	0.50
1:A:42:ASN:O	1:A:45:SER:OG	2.24	0.50
1:C:135:LEU:HD12	1:C:135:LEU:N	2.26	0.50
1:A:326:ASP:OD1	1:A:330:ARG:NH2	2.44	0.50
1:C:364:ILE:HD12	1:C:364:ILE:N	2.26	0.50
1:B:226:TYR:CE1	2:B:601:FAD:N6A	2.80	0.50
1:C:417:ARG:CD	1:C:422:VAL:HG23	2.41	0.50
1:A:228:THR:HG23	1:A:399:LEU:C	2.36	0.50
1:C:69:GLY:O	1:C:417:ARG:NH1	2.44	0.50
1:C:287:LYS:H	1:C:384:ASN:ND2	2.06	0.50
1:A:73:MET:HB2	1:A:74:PRO:CD	2.40	0.49
1:A:19:LYS:HE3	1:A:23:GLU:HG3	1.94	0.49
1:A:310:ASP:OD1	1:A:310:ASP:C	2.56	0.48
1:C:139:LYS:HE2	1:C:226:TYR:CE2	2.48	0.48
1:C:416:GLY:C	1:C:420:GLN:O	2.56	0.48
1:C:273:ILE:HG22	1:C:386:VAL:C	2.38	0.48
1:B:191:LEU:C	1:B:191:LEU:HD13	2.38	0.48
1:C:363:MET:CE	1:C:382:ASP:HB3	2.43	0.48
1:B:310:ASP:OD1	1:B:310:ASP:C	2.57	0.48
1:C:496:ASN:ND2	1:C:529:ASN:OD1	2.47	0.48
1:B:229:PHE:CD1	1:B:404:VAL:HG21	2.48	0.48
1:B:401:LEU:HD12	1:B:408:ARG:HD3	1.96	0.48
1:A:43:LEU:HD11	1:A:93:GLU:CD	2.39	0.48
1:C:310:ASP:C	1:C:310:ASP:OD1	2.57	0.47
1:C:66:GLU:C	1:C:418:ARG:NH1	2.63	0.47
1:C:161:LYS:HG3	1:C:279:PHE:CZ	2.49	0.47
1:C:105:HIS:ND1	1:C:248:THR:HG22	2.28	0.47
1:A:404:VAL:HG12	1:A:404:VAL:O	2.13	0.47
1:B:305:LYS:HD2	1:B:305:LYS:O	2.15	0.47
1:B:293:ALA:HB1	1:B:319:LEU:HA	1.97	0.46
1:C:43:LEU:HD11	1:C:93:GLU:HG3	1.97	0.46
1:C:576:GLU:O	1:C:579:THR:OG1	2.31	0.46
1:B:485:GLU:OE1	1:B:523:LYS:NZ	2.39	0.46
1:A:47:TYR:OH	1:A:93:GLU:OE1	2.14	0.46
1:A:124:GLU:CD	1:A:127:ARG:HH11	2.23	0.46
1:A:365:ARG:HD2	1:A:380:THR:CB	2.46	0.46
1:A:555:LEU:N	1:A:555:LEU:HD12	2.31	0.46
1:C:191:LEU:HD12	1:C:191:LEU:C	2.40	0.46
1:A:38:THR:O	1:A:41:VAL:HG12	2.15	0.46
1:A:401:LEU:HD12	1:A:408:ARG:HD3	1.98	0.46
1:A:409:ALA:O	1:A:411:SER:O	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:586:GLU:OE1	1:A:586:GLU:HA	2.16	0.46
1:C:315:VAL:HG11	1:C:319:LEU:CD2	2.46	0.45
1:C:355:LYS:HG2	1:C:356:PRO:HD2	1.98	0.45
1:B:24:GLU:OE2	1:B:52:ASN:OD1	2.35	0.45
1:C:555:LEU:HD12	1:C:555:LEU:N	2.32	0.45
1:B:555:LEU:N	1:B:555:LEU:HD12	2.32	0.45
1:C:401:LEU:HD12	1:C:408:ARG:HD3	1.97	0.45
1:A:469:PRO:HB2	1:B:570:ILE:HD11	1.99	0.45
1:C:409:ALA:O	1:C:411:SER:O	2.35	0.45
1:C:271:CYS:SG	1:C:352:ARG:NH1	2.90	0.45
1:A:330:ARG:HH21	1:A:330:ARG:HG3	1.82	0.44
1:A:501:HIS:CD2	1:A:585:LYS:CB	2.93	0.44
1:B:259:ARG:HA	1:B:259:ARG:HD2	1.85	0.44
1:C:363:MET:HE1	1:C:382:ASP:HB3	1.98	0.44
1:A:365:ARG:HD2	1:A:380:THR:HB	2.00	0.44
1:A:441:THR:HB	1:A:442:PRO:HD3	2.00	0.44
1:B:135:LEU:N	1:B:135:LEU:CD1	2.81	0.44
1:C:571:HIS:HA	1:C:572:PRO:HA	1.85	0.44
1:B:409:ALA:O	1:B:411:SER:O	2.36	0.43
1:A:293:ALA:HB2	1:A:315:VAL:HG11	1.96	0.43
1:A:415:ILE:O	1:A:422:VAL:HG23	2.19	0.43
1:C:40:VAL:O	1:C:43:LEU:HB3	2.19	0.43
1:B:324:ASP:OD1	1:B:326:ASP:HB3	2.19	0.43
1:C:17:VAL:HG22	1:C:54:VAL:HG11	2.00	0.43
1:C:409:ALA:N	1:C:415:ILE:CD1	2.78	0.43
1:A:293:ALA:HB1	1:A:319:LEU:HA	2.00	0.43
1:A:388:MET:HE2	1:A:388:MET:HB2	1.85	0.43
1:B:465:TYR:HA	1:B:468:ILE:HD11	2.00	0.43
1:C:270:GLU:OE2	1:C:352:ARG:NH1	2.52	0.43
1:C:409:ALA:HB2	1:C:415:ILE:HG12	2.01	0.43
1:C:446:GLN:OE1	1:C:446:GLN:HA	2.19	0.43
1:A:316:ARG:NH1	3:A:602:ATR:N7	2.67	0.42
1:B:19:LYS:HE2	1:B:23:GLU:OE2	2.18	0.42
1:B:492:TYR:O	1:B:496:ASN:ND2	2.39	0.42
1:B:446:GLN:HA	1:B:446:GLN:OE1	2.19	0.42
2:B:601:FAD:H9	2:B:601:FAD:H1'1	1.86	0.42
1:A:367:HIS:HD2	1:A:378:GLU:OE1	2.02	0.42
1:B:42:ASN:O	1:B:45:SER:HB2	2.19	0.42
1:C:139:LYS:HE2	1:C:226:TYR:CD2	2.55	0.42
1:C:283:TYR:CE2	1:C:356:PRO:HB2	2.54	0.42
1:C:441:THR:HB	1:C:442:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:441:THR:O	1:C:445:ILE:HG13	2.19	0.42
1:C:481:CYS:O	1:C:536:GLY:HA2	2.20	0.42
1:A:65:LEU:HD21	1:A:72:SER:O	2.20	0.42
1:B:441:THR:O	1:B:445:ILE:HG13	2.20	0.42
1:C:103:LYS:O	1:C:105:HIS:CD2	2.73	0.42
1:C:176:ASP:OD2	1:C:514:ARG:NH1	2.48	0.42
2:C:601:FAD:H9	2:C:601:FAD:H1'1	1.85	0.42
1:A:191:LEU:HD12	1:B:182:TRP:CE2	2.55	0.42
1:A:496:ASN:HB3	1:A:529:ASN:ND2	2.35	0.42
1:A:103:LYS:O	1:A:105:HIS:CD2	2.73	0.42
1:A:176:ASP:OD2	1:A:514:ARG:NH1	2.49	0.42
1:B:392:ARG:HD2	2:B:601:FAD:HM82	2.02	0.42
1:C:204:ILE:HG22	1:C:208:ASN:ND2	2.35	0.42
1:A:17:VAL:HG22	1:A:54:VAL:HG11	2.02	0.41
4:A:603:MPD:H52	4:A:603:MPD:O2	2.20	0.41
1:A:81:ASN:HD21	1:A:251:ARG:HD3	1.85	0.41
1:A:135:LEU:HD13	1:A:227:ALA:HB2	2.01	0.41
1:C:256:VAL:O	1:C:256:VAL:CG2	2.63	0.41
1:A:79:LYS:NZ	1:A:101:TRP:O	2.50	0.41
1:B:176:ASP:OD2	1:B:514:ARG:NH1	2.46	0.41
1:A:550:GLY:O	1:A:553:ILE:HG22	2.20	0.41
1:C:30:TYR:H	1:C:56:VAL:HG23	1.85	0.41
1:A:191:LEU:HD12	1:B:182:TRP:CD1	2.56	0.41
1:C:147:TRP:CZ3	1:C:276:ASP:HB2	2.56	0.41
1:C:409:ALA:HB2	1:C:415:ILE:HD11	2.01	0.41
1:C:417:ARG:N	1:C:420:GLN:O	2.54	0.41
1:A:110:ALA:HB3	1:A:252:PHE:CD1	2.56	0.41
1:A:446:GLN:HA	1:A:446:GLN:OE1	2.20	0.41
1:A:224:ASN:OD1	1:A:240:LYS:HE3	2.20	0.41
1:B:495:GLU:OE1	1:B:495:GLU:N	2.37	0.41
1:C:79:LYS:O	1:C:251:ARG:NH1	2.54	0.41
1:C:135:LEU:N	1:C:135:LEU:CD1	2.84	0.41
1:C:417:ARG:HG2	1:C:422:VAL:CB	2.50	0.41
1:C:501:HIS:NE2	1:C:585:LYS:HA	2.35	0.41
1:C:501:HIS:NE2	1:C:585:LYS:CA	2.84	0.41
1:A:324:ASP:OD1	1:A:326:ASP:HB3	2.20	0.41
1:A:365:ARG:HD2	1:A:380:THR:OG1	2.20	0.41
1:B:58:ASN:C	1:B:60:ALA:H	2.28	0.41
1:C:228:THR:OG1	1:C:398:ASP:O	2.38	0.41
1:B:103:LYS:O	1:B:105:HIS:CD2	2.75	0.40
1:C:324:ASP:OD1	1:C:326:ASP:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:PRO:HB2	1:A:265:VAL:HG23	2.04	0.40
1:C:273:ILE:CD1	1:C:281:LEU:HD11	2.51	0.40
1:C:496:ASN:HB3	1:C:529:ASN:HD21	1.86	0.40
1:B:17:VAL:HG22	1:B:54:VAL:HG11	2.03	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	575/598 (96%)	549 (96%)	26 (4%)	0	100	100
1	B	575/598 (96%)	542 (94%)	32 (6%)	1 (0%)	43	73
1	C	525/598 (88%)	498 (95%)	27 (5%)	0	100	100
All	All	1675/1794 (93%)	1589 (95%)	85 (5%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	59	ASP

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	488/504 (97%)	475 (97%)	13 (3%)	39	67
1	B	488/504 (97%)	466 (96%)	22 (4%)	24	56
1	C	460/504 (91%)	442 (96%)	18 (4%)	28	60
All	All	1436/1512 (95%)	1383 (96%)	53 (4%)	30	61

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	CYS
1	A	237	VAL
1	A	319	LEU
1	A	371	GLU
1	A	380	THR
1	A	422	VAL
1	A	441	THR
1	A	445	ILE
1	A	468	ILE
1	A	474	THR
1	A	503	VAL
1	A	533	LEU
1	A	582	THR
1	B	50	THR
1	B	146	THR
1	B	158	CYS
1	B	161	LYS
1	B	191	LEU
1	B	202	ASN
1	B	237	VAL
1	B	238	LYS
1	B	256	VAL
1	B	310	ASP
1	B	315	VAL
1	B	346	VAL
1	B	379	VAL
1	B	380	THR
1	B	441	THR
1	B	445	ILE
1	B	474	THR
1	B	475	PRO
1	B	503	VAL
1	B	518	THR
1	B	533	LEU

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Mol	Chain	Res	Type
1	B	579	THR
1	C	59	ASP
1	C	88	GLU
1	C	158	CYS
1	C	204	ILE
1	C	208	ASN
1	C	237	VAL
1	C	270	GLU
1	C	314	MET
1	C	343	VAL
1	C	365	ARG
1	C	379	VAL
1	C	415	ILE
1	C	419	GLU
1	C	441	THR
1	C	468	ILE
1	C	474	THR
1	C	533	LEU
1	C	570	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	99	ASN
1	A	105	HIS
1	A	155	ASN
1	A	195	GLN
1	A	208	ASN
1	A	367	HIS
1	A	529	ASN
1	B	34	ASN
1	B	91	GLN
1	B	99	ASN
1	B	105	HIS
1	B	203	HIS
1	B	208	ASN
1	C	48	ASN
1	C	58	ASN
1	C	99	ASN
1	C	208	ASN
1	C	384	ASN

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Mol	Chain	Res	Type
1	C	496	ASN
1	C	529	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 ⓘ

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPD	A	604	-	7,7,7	0.15	0	9,10,10	0.30	0
4	MPD	B	605	-	7,7,7	0.15	0	9,10,10	0.30	0
4	MPD	B	603	-	7,7,7	0.15	0	9,10,10	0.30	0
2	FAD	C	601	-	58,58,58	0.55	0	85,89,89	0.55	1 (1%)
4	MPD	B	604	-	7,7,7	0.15	0	9,10,10	0.30	0
2	FAD	B	601	-	58,58,58	0.61	0	85,89,89	0.64	1 (1%)
3	ATR	A	602	-	32,33,33	1.34	4 (12%)	50,52,52	1.79	10 (20%)
4	MPD	A	603	-	7,7,7	0.16	0	9,10,10	0.30	0
2	FAD	A	601	5	58,58,58	0.63	0	85,89,89	0.62	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
4	MPD	A	604	-	-	1/5/5/5	-
4	MPD	B	605	-	-	1/5/5/5	-
4	MPD	B	603	-	-	3/5/5/5	-
2	FAD	C	601	-	-	5/34/50/50	0/6/6/6
4	MPD	B	604	-	-	0/5/5/5	-
2	FAD	B	601	-	-	2/34/50/50	0/6/6/6
3	ATR	A	602	-	-	6/21/37/37	0/3/3/3
4	MPD	A	603	-	-	3/5/5/5	-
2	FAD	A	601	5	-	5/34/50/50	0/6/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	ATR	C5-C4	4.63	1.47	1.39
3	A	602	ATR	C5-C6	2.68	1.48	1.41
3	A	602	ATR	C5-N7	-2.31	1.34	1.39
3	A	602	ATR	C8-N7	2.20	1.35	1.31

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	ATR	C5-C4-N3	-5.73	118.82	126.72
3	A	602	ATR	N3-C4-N9	4.54	134.89	127.17
3	A	602	ATR	C2-N3-C4	3.63	120.70	111.83
3	A	602	ATR	C4-C5-N7	-3.48	106.61	110.58
3	A	602	ATR	C2'-C1'-N9	-3.29	108.34	113.75
3	A	602	ATR	N3-C2-N1	-3.20	123.75	128.58
3	A	602	ATR	C4-N9-C8	2.67	108.54	105.74
3	A	602	ATR	C5-N7-C8	2.51	107.40	103.45
3	A	602	ATR	C6-C5-N7	2.10	136.14	132.09
2	B	601	FAD	C4-N3-C2	-2.08	121.94	125.64
2	C	601	FAD	C4-N3-C2	-2.08	121.94	125.64
3	A	602	ATR	P2'-O2'-C2'	-2.07	117.90	123.43

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	PA-O3P-P-O5'
2	C	601	FAD	N10-C1'-C2'-O2'
2	C	601	FAD	PA-O3P-P-O5'
3	A	602	ATR	C5'-O5'-PA-O2A
3	A	602	ATR	C5'-O5'-PA-O3A
4	A	603	MPD	C2-C3-C4-O4
4	A	603	MPD	C2-C3-C4-C5
4	B	603	MPD	O2-C2-C3-C4
4	B	603	MPD	CM-C2-C3-C4
2	A	601	FAD	O4B-C4B-C5B-O5B
2	C	601	FAD	O4B-C4B-C5B-O5B
2	A	601	FAD	C3B-C4B-C5B-O5B
2	C	601	FAD	C3B-C4B-C5B-O5B
2	B	601	FAD	PA-O3P-P-O5'
2	A	601	FAD	C2B-C1B-N9A-C8A
4	A	604	MPD	C1-C2-C3-C4
3	A	602	ATR	C5'-O5'-PA-O1A
3	A	602	ATR	O4'-C4'-C5'-O5'
3	A	602	ATR	C3'-C4'-C5'-O5'
2	C	601	FAD	C2B-C1B-N9A-C8A
4	A	603	MPD	O2-C2-C3-C4
4	B	605	MPD	O2-C2-C3-C4
3	A	602	ATR	C2'-O2'-P2'-O3P
2	B	601	FAD	C2B-C1B-N9A-C8A
4	B	603	MPD	C1-C2-C3-C4
2	A	601	FAD	P-O3P-PA-O2A

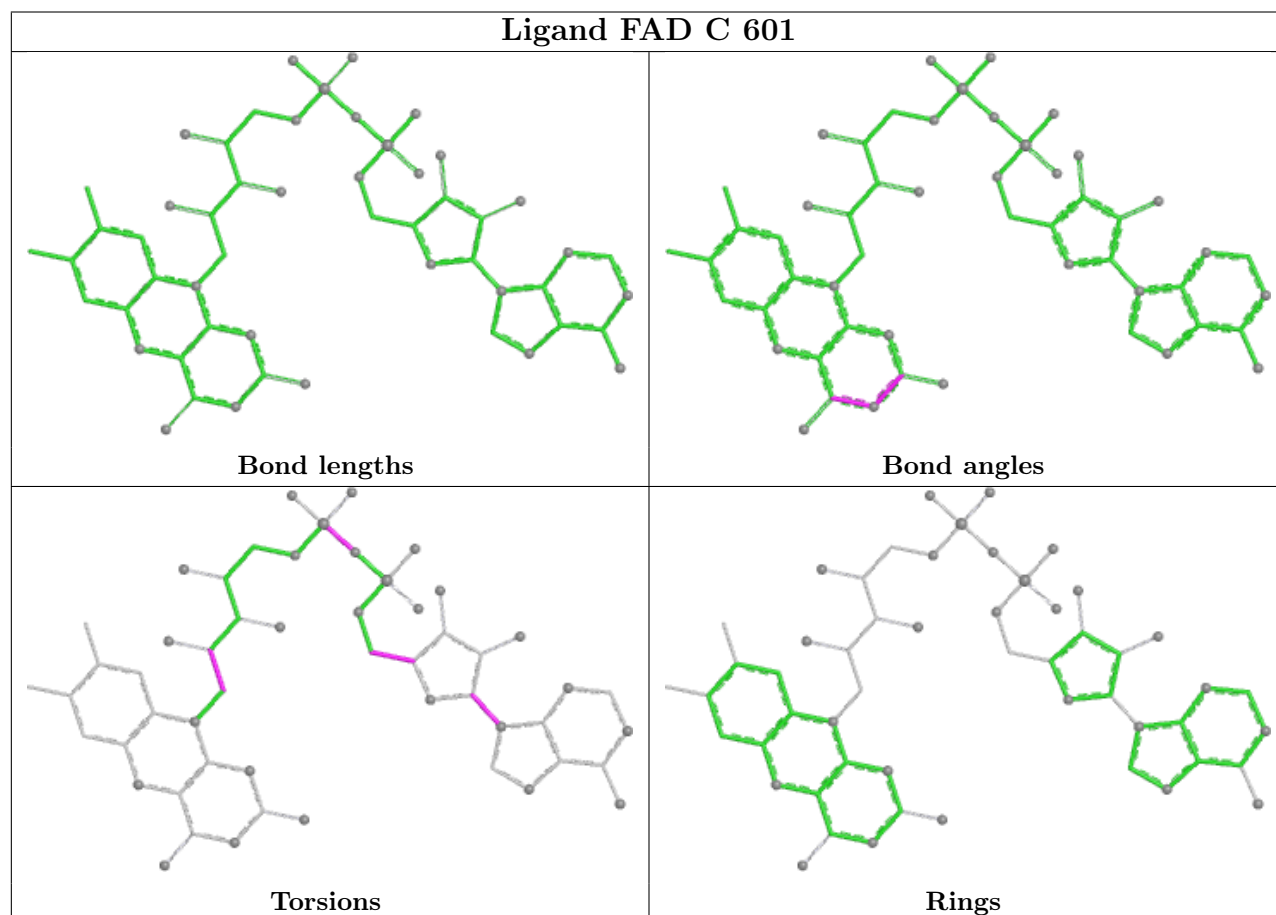
There are no ring outliers.

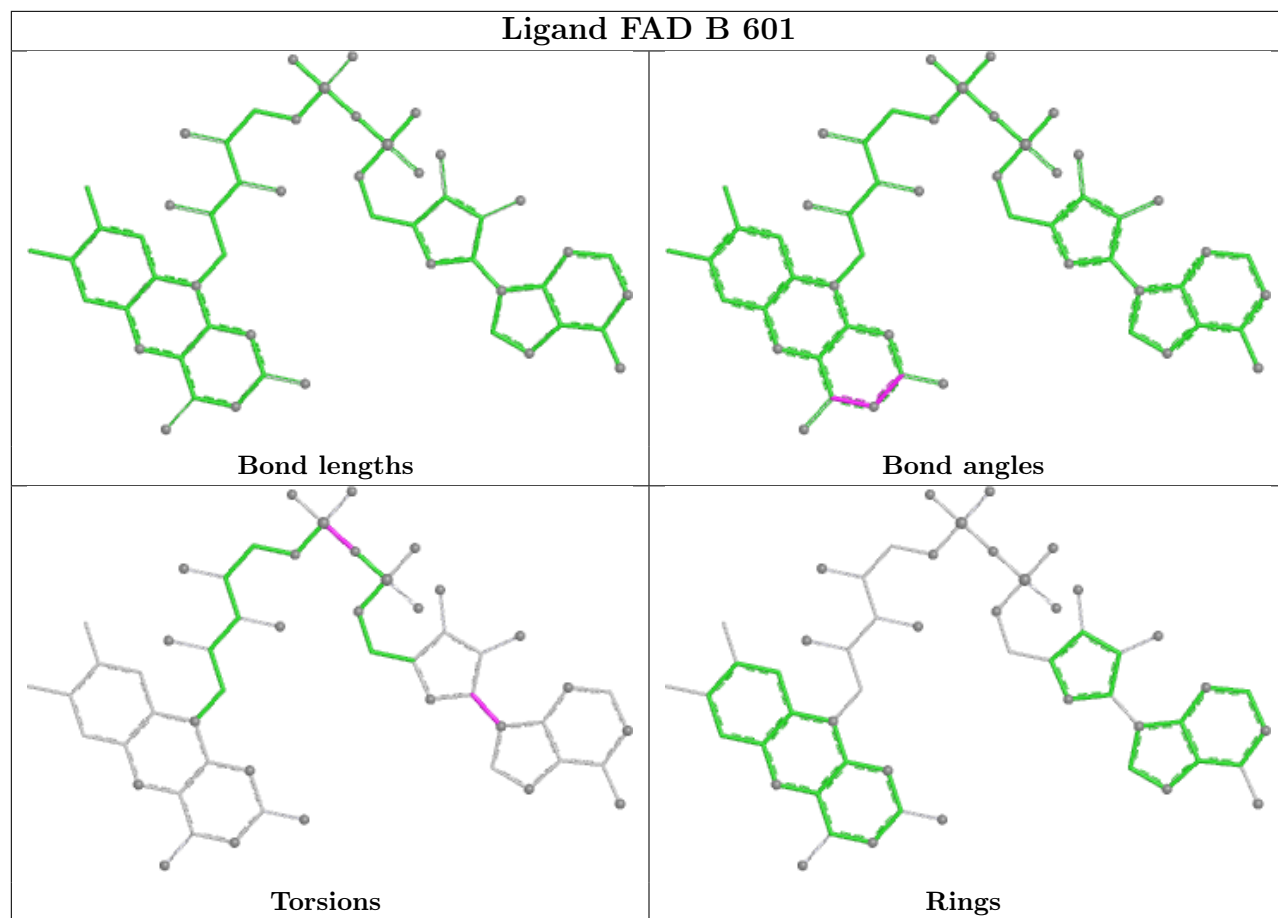
7 monomers are involved in 20 short contacts:

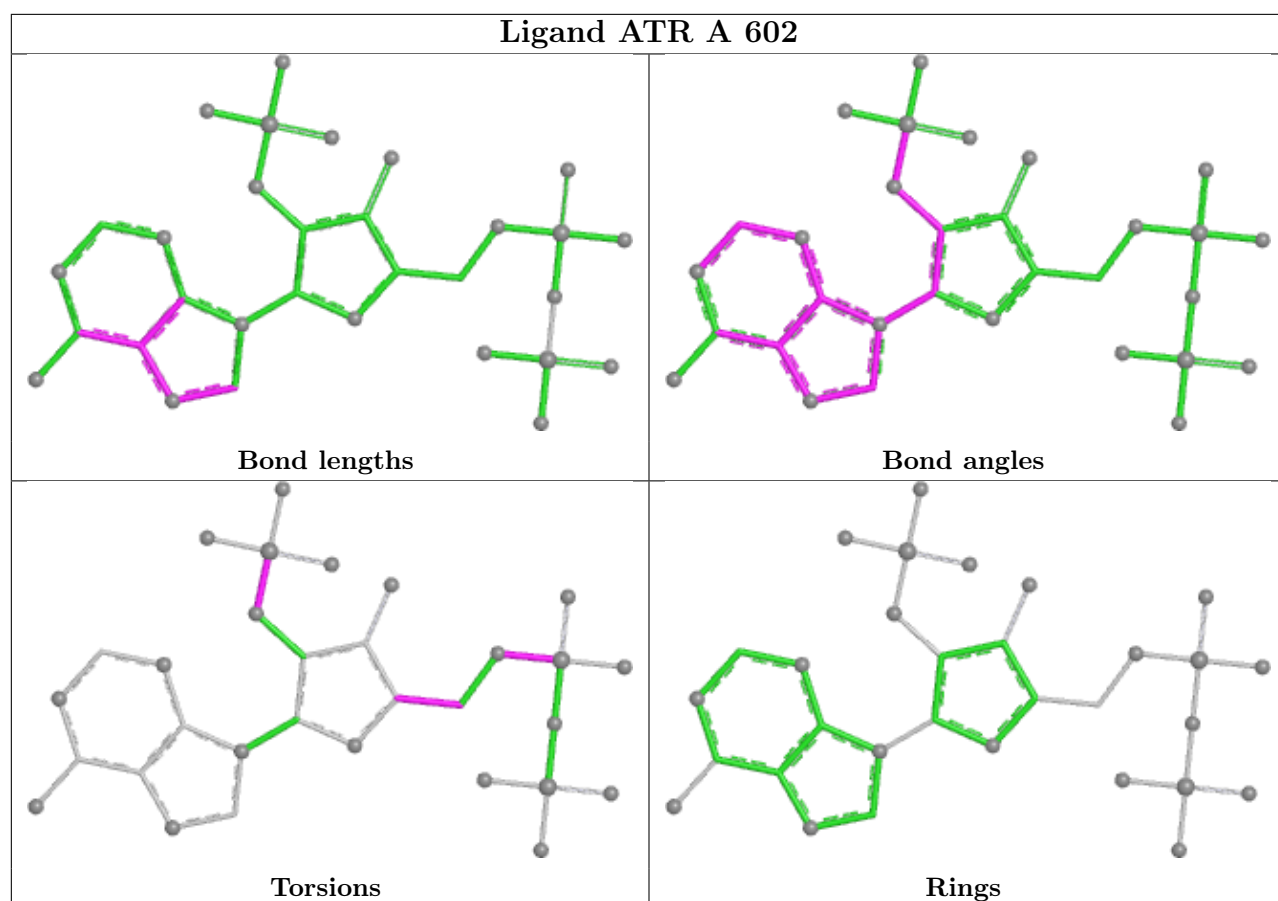
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	604	MPD	2	0
4	B	603	MPD	2	0
2	C	601	FAD	4	0
4	B	604	MPD	1	0
2	B	601	FAD	4	0
3	A	602	ATR	6	0
4	A	603	MPD	1	0

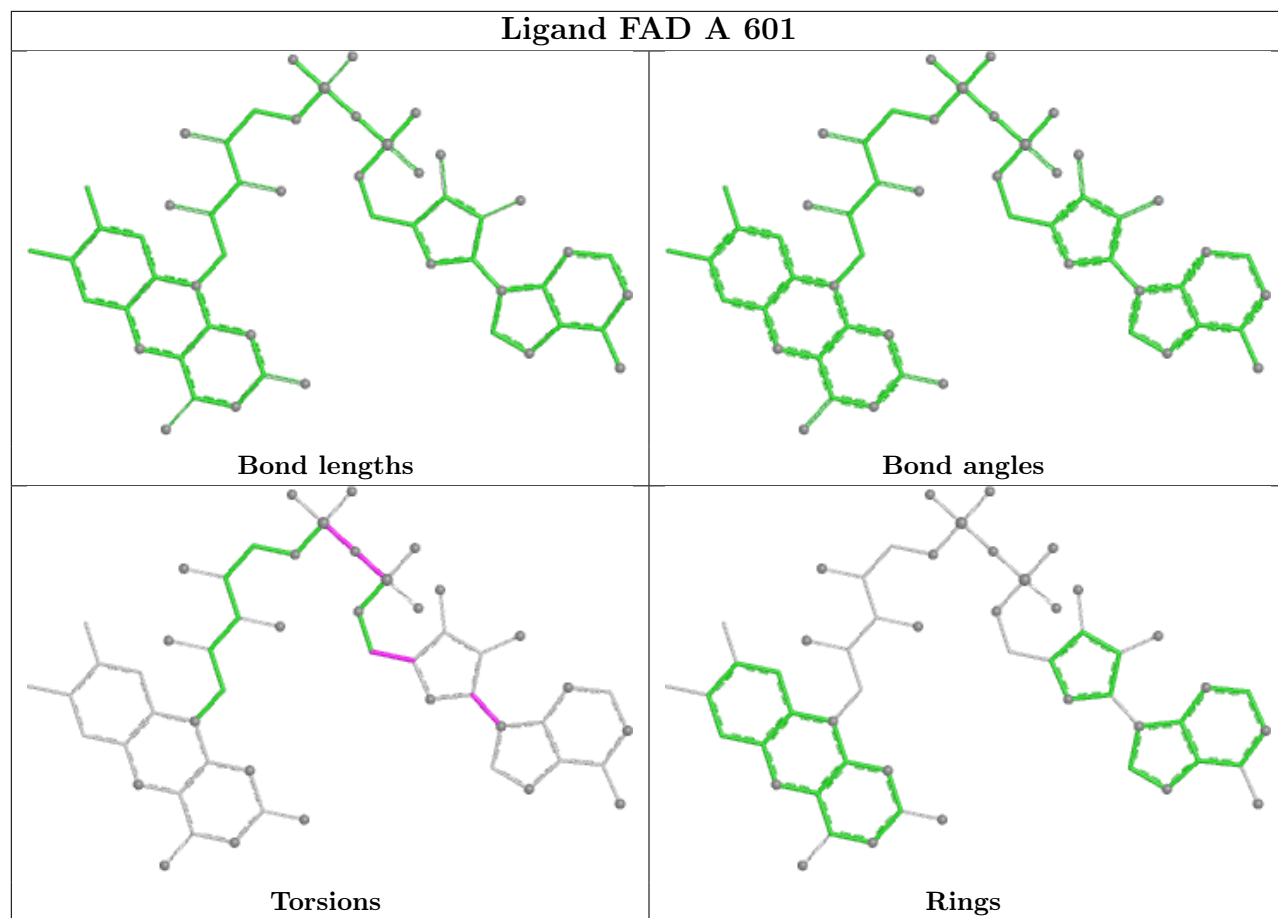
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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









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	577/598 (96%)	-0.58	0	100 100	61, 93, 160, 231	1 (0%)
1	B	577/598 (96%)	-0.63	0	100 100	61, 101, 153, 214	1 (0%)
1	C	541/598 (90%)	-0.10	6 (1%)	78 59	80, 172, 274, 338	1 (0%)
All	All	1695/1794 (94%)	-0.44	6 (0%)	88 76	61, 112, 226, 338	3 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	35	PRO	3.7
1	C	191	LEU	2.8
1	C	379	VAL	2.3
1	C	290	CYS	2.2
1	C	419	GLU	2.2
1	C	275	SER	2.1

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](#)

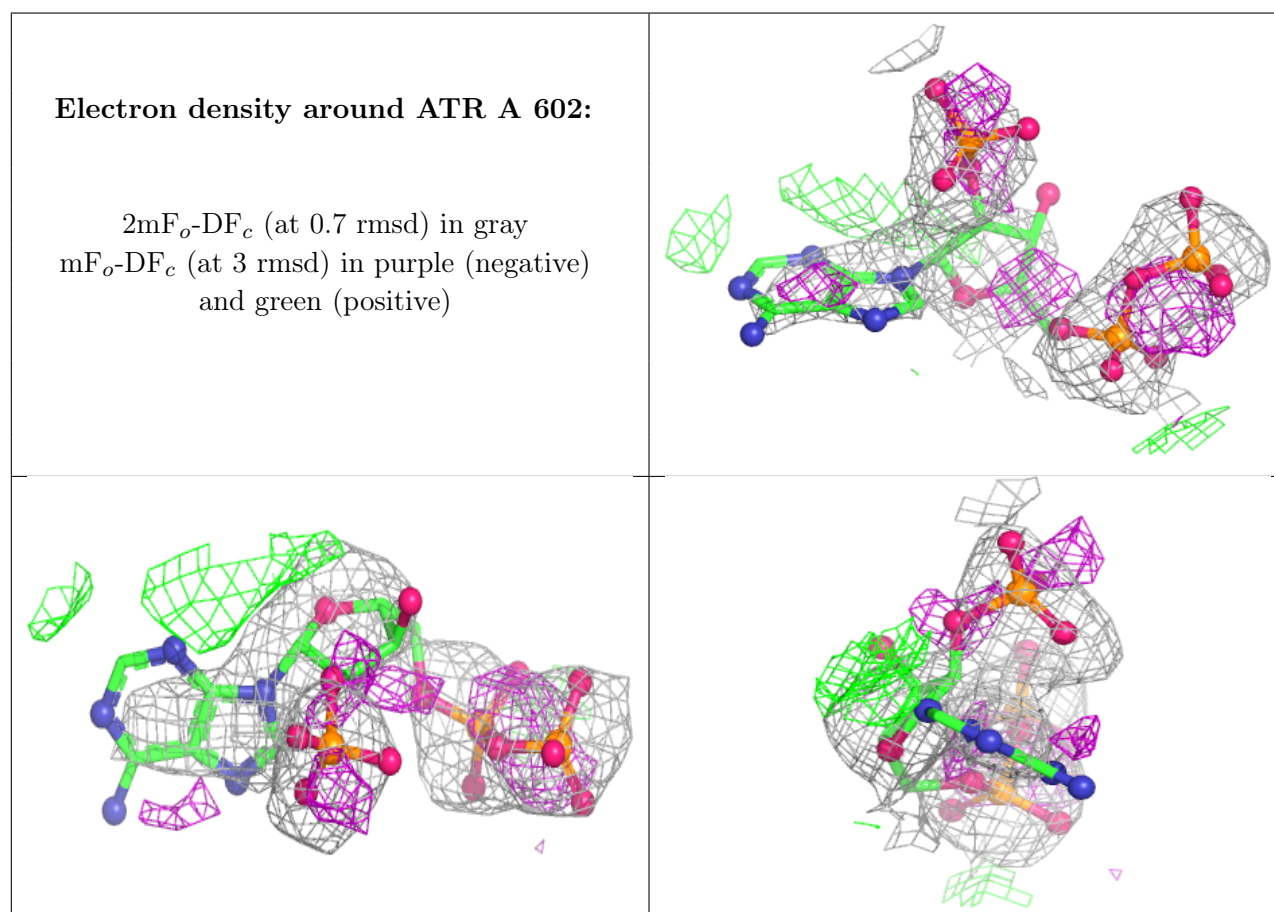
There are no oligosaccharides in this entry.

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.

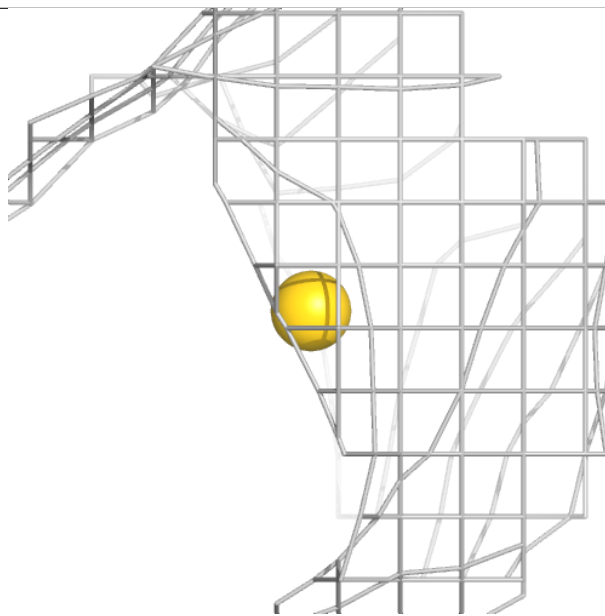
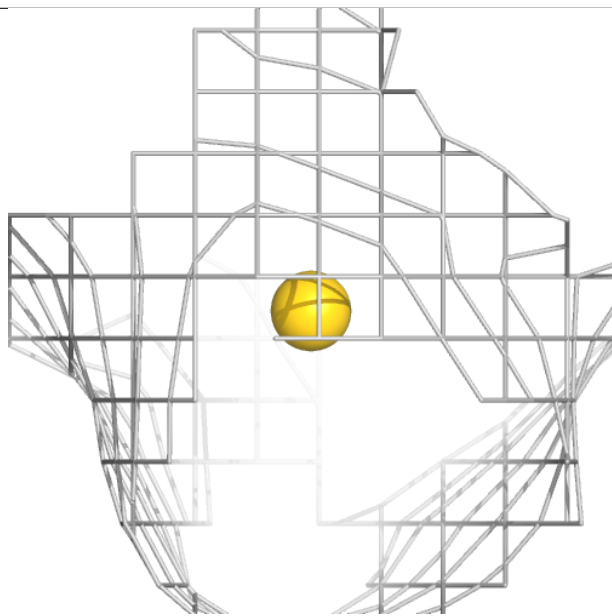
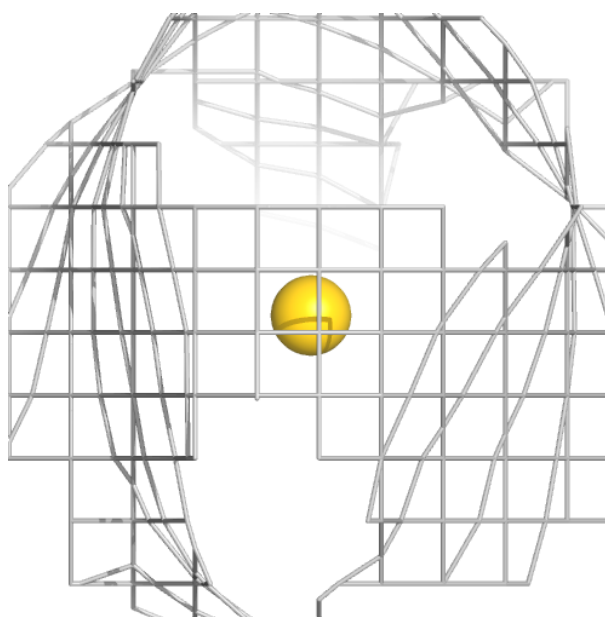
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NA	A	605	1/1	0.74	0.18	79,79,79,79	1
3	ATR	A	602	31/31	0.76	0.13	90,108,121,124	31
4	MPD	B	604	8/8	0.83	0.17	124,132,139,141	0
4	MPD	B	605	8/8	0.88	0.24	104,115,124,130	0
6	AU	B	602	1/1	0.88	0.07	175,175,175,175	1
4	MPD	B	603	8/8	0.90	0.16	102,120,128,129	0
4	MPD	A	604	8/8	0.91	0.14	127,135,147,155	0
2	FAD	C	601	53/53	0.92	0.09	98,144,170,172	0
4	MPD	A	603	8/8	0.93	0.14	89,107,122,122	0
2	FAD	B	601	53/53	0.96	0.07	65,80,99,102	0
2	FAD	A	601	53/53	0.98	0.05	60,73,94,100	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.



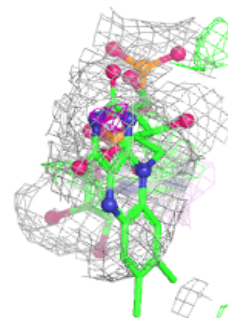
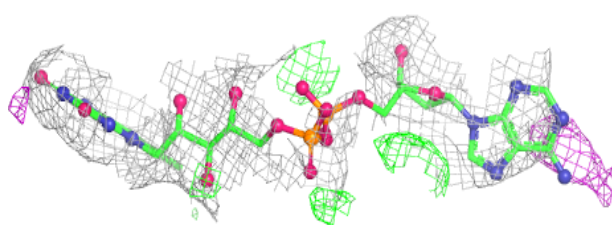
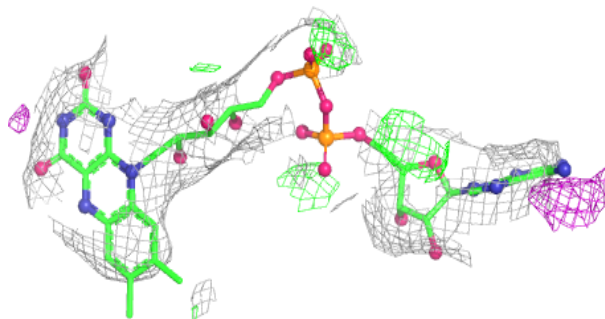
Electron density around AU B 602:

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

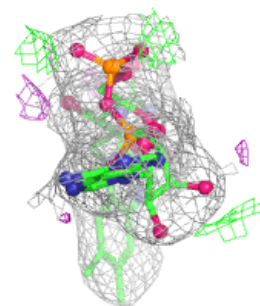
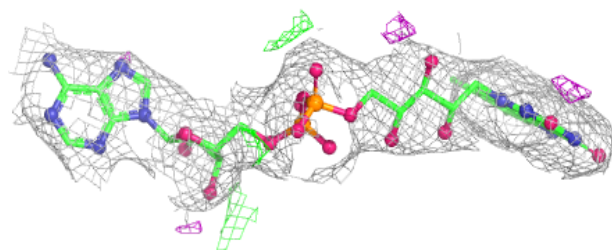
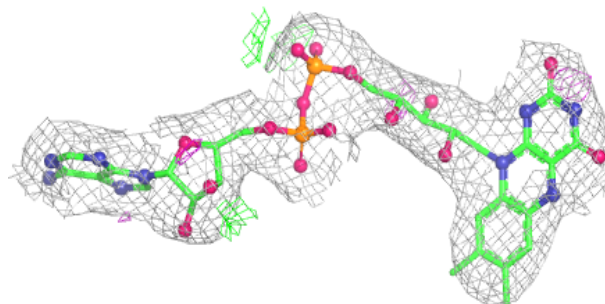


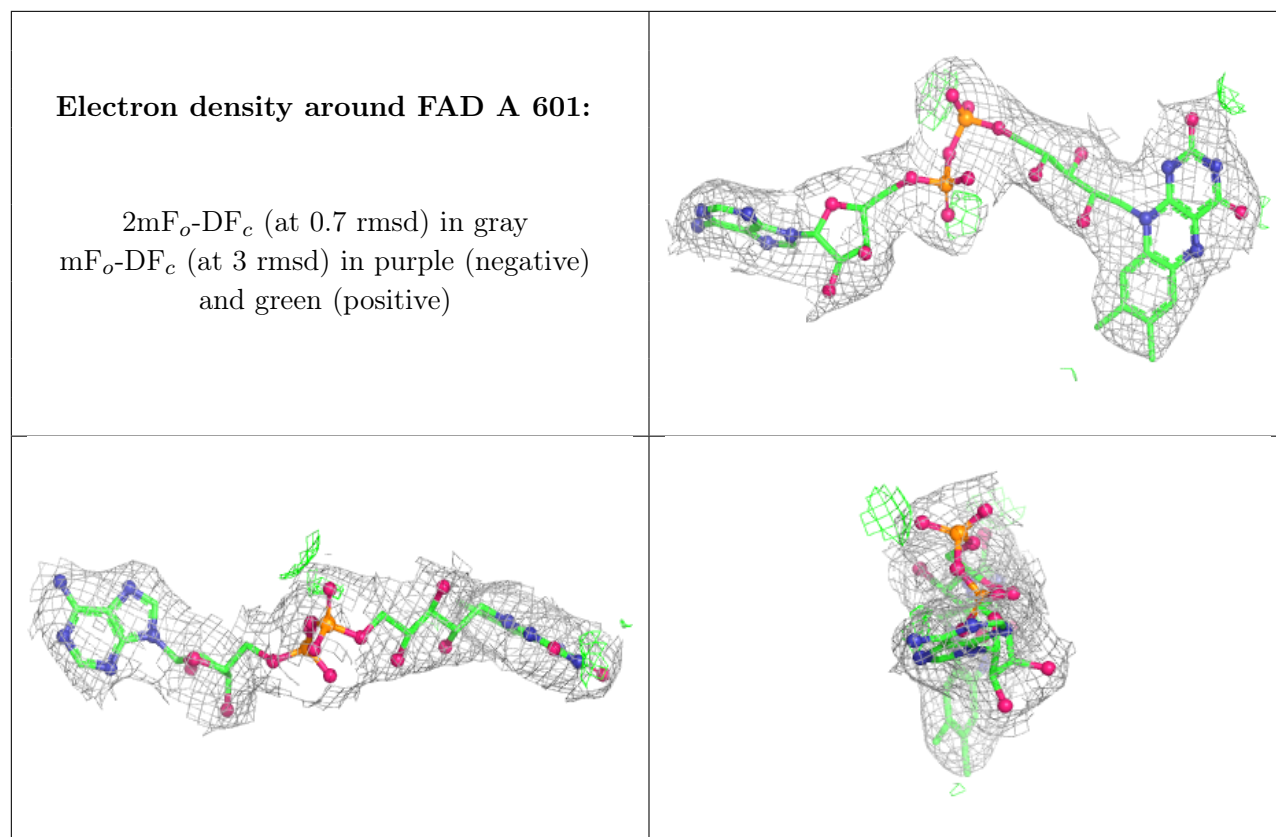
Electron density around FAD C 601:

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

**Electron density around FAD B 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.