



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:29 AM UTC

PDB ID : 8F5S / pdb_00008f5s
Title : human branched chain ketoacid dehydrogenase kinase in complex with inhibitors
Authors : Liu, S.; Roth Flach, R.; Bollinger, E.; Filipski, K.
Deposited on : 2022-11-15
Resolution : 2.79 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

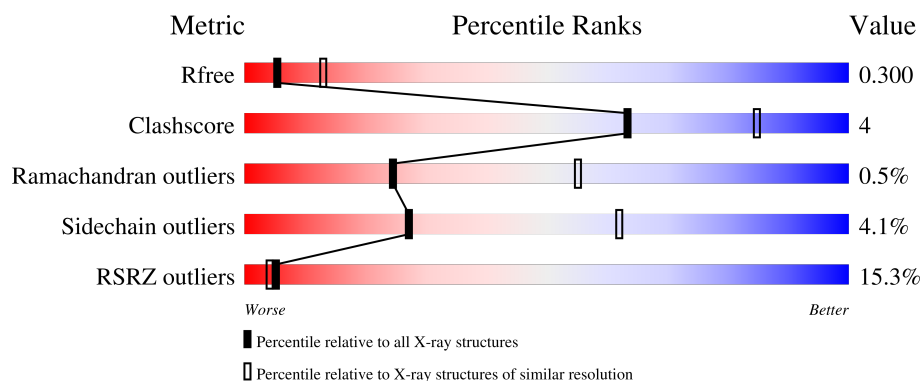
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	388	
1	B	388	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5204 atoms, of which 24 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.

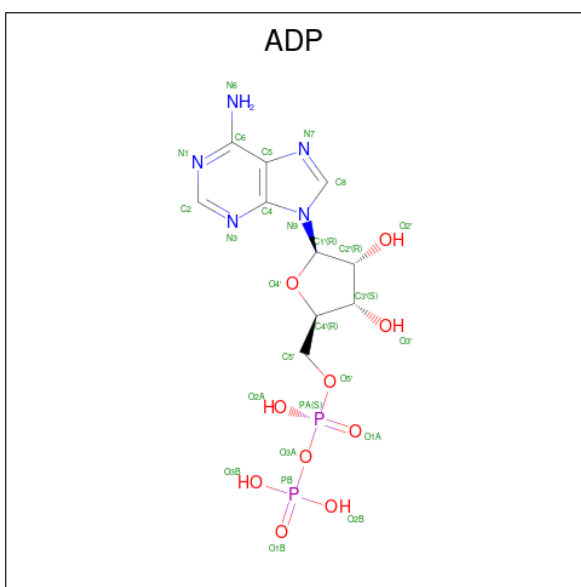
- Molecule 1 is a protein called [3-methyl-2-oxobutanoate dehydrogenase [lipoamide]] kinase, mitochondrial.

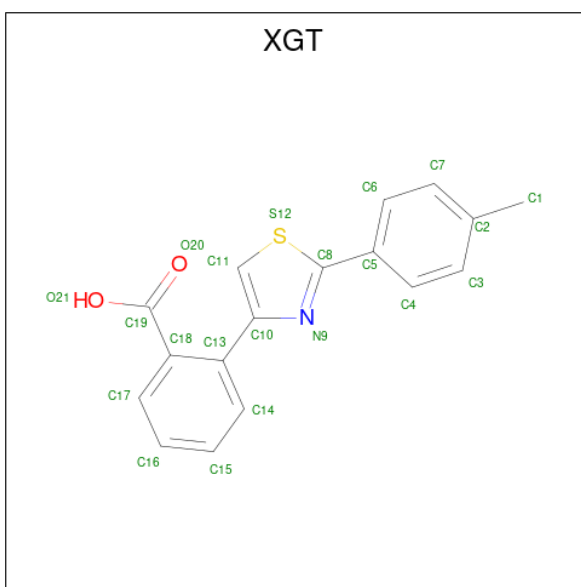
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2554	1630	459	452	13			
1	B	314	Total	C	N	O	S	0	0	0
			2490	1592	447	439	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	383	HIS	-	expression tag	UNP O14874
A	384	HIS	-	expression tag	UNP O14874
A	385	HIS	-	expression tag	UNP O14874
A	386	HIS	-	expression tag	UNP O14874
A	387	HIS	-	expression tag	UNP O14874
A	388	HIS	-	expression tag	UNP O14874
B	383	HIS	-	expression tag	UNP O14874
B	384	HIS	-	expression tag	UNP O14874
B	385	HIS	-	expression tag	UNP O14874
B	386	HIS	-	expression tag	UNP O14874
B	387	HIS	-	expression tag	UNP O14874
B	388	HIS	-	expression tag	UNP O14874

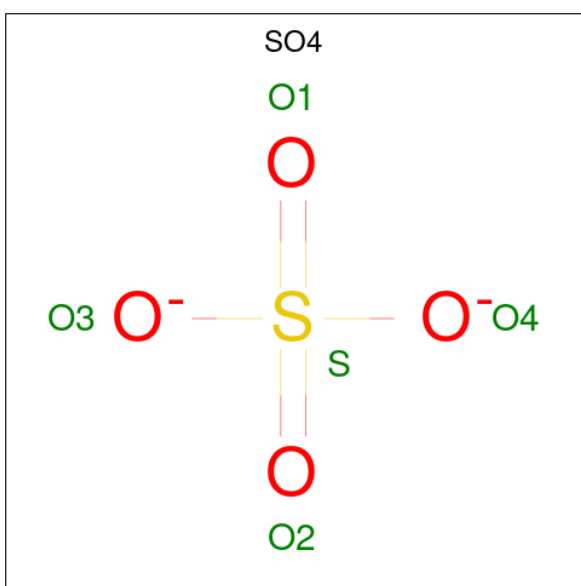
- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).





Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	S	12	0
			33	17	12	1	2	1		
5	B	1	Total	C	H	N	O	S	12	0
			33	17	12	1	2	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

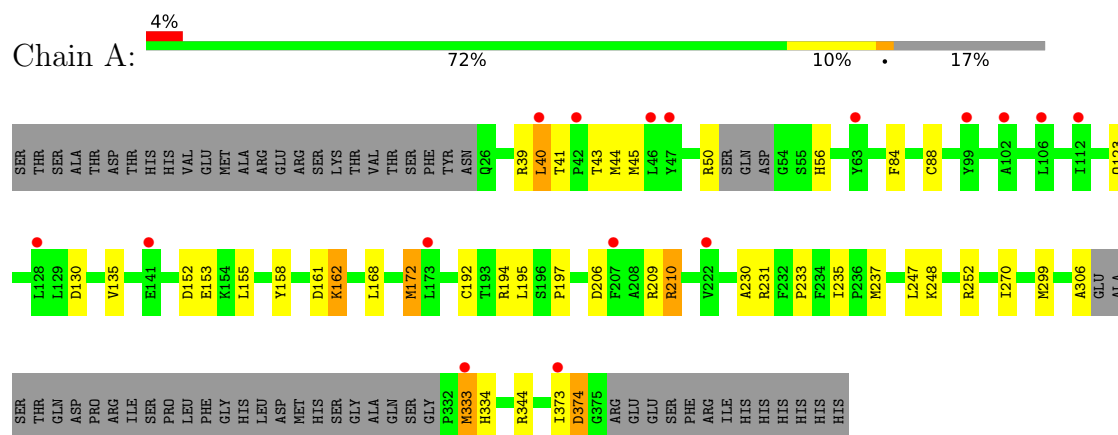
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	11	Total	O	0	0
			11	11		
7	B	1	Total	O	0	0
			1	1		

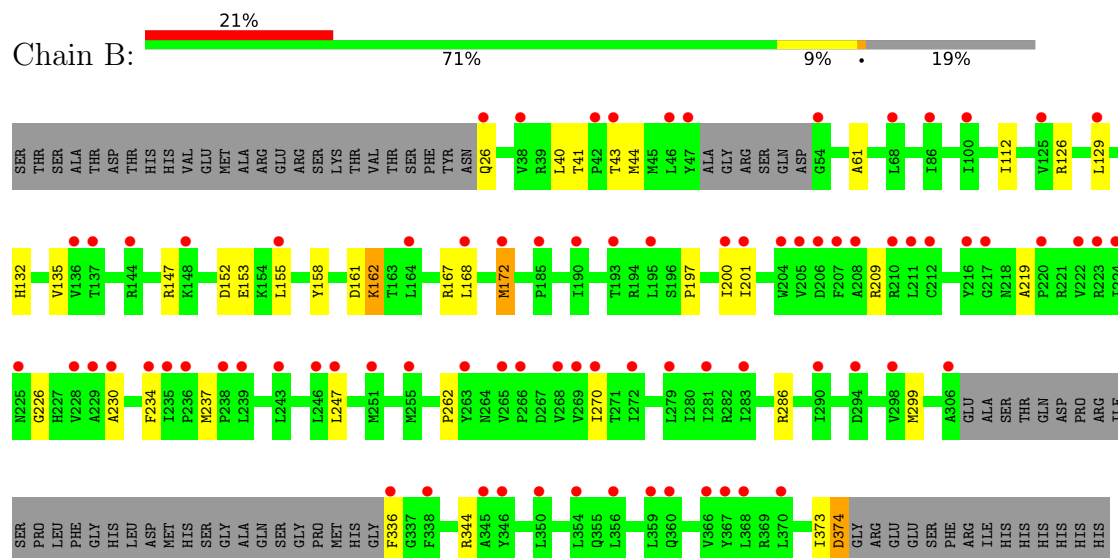
3 Residue-property plots [i](#)

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: [3-methyl-2-oxobutanoate dehydrogenase [lipoamide]] kinase, mitochondrial



- Molecule 1: [3-methyl-2-oxobutanoate dehydrogenase [lipoamide]] kinase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	117.77Å 117.77Å 154.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	117.77 – 2.79 117.77 – 2.79	Depositor EDS
% Data completeness (in resolution range)	63.1 (117.77-2.79) 63.1 (117.77-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.77Å)	Xtriage
Refinement program	BUSTER 2.11.8 (8-JUN-2022)	Depositor
R, R_{free}	0.257 , 0.297 0.252 , 0.300	Depositor DCC
R_{free} test set	843 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	88.4	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 103.9	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	0.92	EDS
Total number of atoms	5204	wwPDB-VP
Average B, all atoms (Å ²)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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: MG, SO4, XGT, ADP, K

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	0.68	0/2610	1.10	3/3532 (0.1%)
1	B	0.62	0/2544	1.10	6/3446 (0.2%)
All	All	0.65	0/5154	1.10	9/6978 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	201	ILE	N-CA-C	-6.19	104.81	110.82
1	A	40	LEU	CA-C-N	5.76	128.86	120.51
1	A	40	LEU	C-N-CA	5.76	128.86	120.51
1	B	161	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	26	GLN	CA-C-N	5.19	127.75	120.28
1	B	26	GLN	C-N-CA	5.19	127.75	120.28
1	B	262	PRO	CA-C-N	5.18	130.21	121.14
1	B	262	PRO	C-N-CA	5.18	130.21	121.14
1	A	161	ASP	CA-CB-CG	5.14	117.74	112.60

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	2554	0	2576	24	0
1	B	2490	0	2510	18	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
5	A	21	12	0	0	0
5	B	21	12	0	0	0
6	A	20	0	0	0	0
6	B	5	0	0	0	0
7	A	11	0	0	0	0
7	B	1	0	0	0	0
All	All	5180	24	5110	41	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 4.

All (41) 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:248:LYS:HE3	1:A:333:MET:HA	1.56	0.85
1:B:41:THR:HB	1:B:44:MET:HG3	1.62	0.79
1:B:209:ARG:HG2	1:B:219:ALA:HB3	1.64	0.78
1:A:41:THR:HB	1:A:44:MET:HG3	1.65	0.76
1:A:158:TYR:HB3	1:A:162:LYS:NZ	2.11	0.65
1:B:158:TYR:HB3	1:B:162:LYS:NZ	2.12	0.64
1:A:158:TYR:HB3	1:A:162:LYS:HZ2	1.65	0.61
1:B:158:TYR:HB3	1:B:162:LYS:HZ1	1.68	0.58
1:B:132:HIS:HB3	1:B:167:ARG:HH12	1.69	0.56
1:A:41:THR:HB	1:A:44:MET:CG	2.37	0.54
1:A:299:MET:HE2	1:A:344:ARG:HD2	1.88	0.54
1:B:299:MET:HE2	1:B:344:ARG:HD2	1.90	0.53
1:A:247:LEU:HD21	1:A:270:ILE:HD11	1.91	0.52
1:B:41:THR:HB	1:B:44:MET:CG	2.35	0.52
1:B:209:ARG:HG2	1:B:219:ALA:CB	2.38	0.51
1:A:50:ARG:CB	1:A:56:HIS:ND1	2.75	0.50
1:A:194:ARG:HE	1:B:209:ARG:HH12	1.59	0.50
1:B:373:ILE:O	1:B:374:ASP:HB2	2.13	0.49
1:A:248:LYS:CE	1:A:333:MET:HA	2.38	0.47
1:B:247:LEU:HD21	1:B:270:ILE:HD11	1.96	0.47
1:A:168:LEU:HD12	1:A:237:MET:HE3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:ILE:O	1:A:374:ASP:CB	2.63	0.46
1:B:126:ARG:HA	1:B:129:LEU:HD12	1.99	0.45
1:B:168:LEU:HD12	1:B:237:MET:HE3	1.99	0.45
1:A:192:CYS:O	1:A:233:PRO:HA	2.17	0.44
1:A:40:LEU:HG	1:A:45:MET:HG2	1.99	0.44
1:A:252:ARG:HD2	1:A:334:HIS:CG	2.52	0.43
1:B:172:MET:HE3	1:B:237:MET:HE2	2.01	0.43
1:B:197:PRO:HD2	1:B:230:ALA:O	2.19	0.42
1:A:197:PRO:HD2	1:A:230:ALA:O	2.19	0.42
1:B:61:ALA:HB1	1:B:112:ILE:HD13	2.02	0.42
1:B:153:GLU:H	1:B:153:GLU:CD	2.28	0.42
1:A:172:MET:HE3	1:A:237:MET:HE2	2.02	0.41
1:A:153:GLU:CD	1:A:153:GLU:H	2.28	0.41
1:A:172:MET:HE2	1:A:235:ILE:HG21	2.02	0.41
1:A:152:ASP:HB3	1:A:155:LEU:HB2	2.03	0.41
1:B:152:ASP:HB3	1:B:155:LEU:HB2	2.02	0.41
1:A:209:ARG:NH1	1:A:210:ARG:HH12	2.19	0.40
1:A:84:PHE:CZ	1:A:88:CYS:SG	3.14	0.40
1:A:306:ALA:CB	1:A:334:HIS:ND1	2.85	0.40
1:A:235:ILE:HG13	1:A:373:ILE:HD11	2.04	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	316/388 (81%)	305 (96%)	9 (3%)	2 (1%)	21	51
1	B	308/388 (79%)	295 (96%)	12 (4%)	1 (0%)	36	66
All	All	624/776 (80%)	600 (96%)	21 (3%)	3 (0%)	24	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	ASP
1	B	226	GLY
1	A	333	MET

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	273/337 (81%)	262 (96%)	11 (4%)	28	63
1	B	266/337 (79%)	255 (96%)	11 (4%)	27	62
All	All	539/674 (80%)	517 (96%)	22 (4%)	27	62

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ARG
1	A	43	THR
1	A	123	GLN
1	A	130	ASP
1	A	135	VAL
1	A	162	LYS
1	A	172	MET
1	A	195	LEU
1	A	206	ASP
1	A	210	ARG
1	A	231	ARG
1	B	40	LEU
1	B	43	THR
1	B	135	VAL
1	B	147	ARG
1	B	162	LYS
1	B	172	MET
1	B	200	ILE
1	B	234	PHE
1	B	286	ARG
1	B	336	PHE

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Mol	Chain	Res	Type
1	B	374	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	127	GLN
1	A	214	HIS
1	A	292	HIS
1	B	214	HIS
1	B	249	ASN
1	B	292	HIS

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](#)

Of 12 ligands modelled in this entry, 3 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	$\# Z > 2$
5	XGT	A	404	-	23,23,23	0.24	0	32,32,32	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	A	407	-	4,4,4	0.34	0	6,6,6	0.77	0
6	SO4	A	405	-	4,4,4	0.36	0	6,6,6	0.22	0
2	ADP	B	401	3	28,29,29	0.41	0	43,45,45	0.47	0
6	SO4	B	404	-	4,4,4	0.35	0	6,6,6	0.43	0
6	SO4	A	406	-	4,4,4	0.30	0	6,6,6	0.45	0
5	XGT	B	403	-	23,23,23	0.83	2 (8%)	32,32,32	0.78	2 (6%)
2	ADP	A	401	4,3	28,29,29	0.42	0	43,45,45	0.47	0
6	SO4	A	408	-	4,4,4	0.36	0	6,6,6	0.38	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
5	XGT	A	404	-	-	7/12/12/12	0/3/3/3
5	XGT	B	403	-	-	7/12/12/12	0/3/3/3
2	ADP	A	401	4,3	-	2/16/32/32	0/3/3/3
2	ADP	B	401	3	-	1/16/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	403	XGT	O20-C19	2.96	1.31	1.22
5	B	403	XGT	O21-C19	-2.55	1.22	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	403	XGT	O20-C19-C18	-2.93	114.97	121.97
5	B	403	XGT	O21-C19-C18	2.11	121.29	115.28

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ADP	PA-O3A-PB-O3B
2	B	401	ADP	PA-O3A-PB-O2B
5	A	404	XGT	C11-C10-C13-C18
5	B	403	XGT	N9-C10-C13-C18

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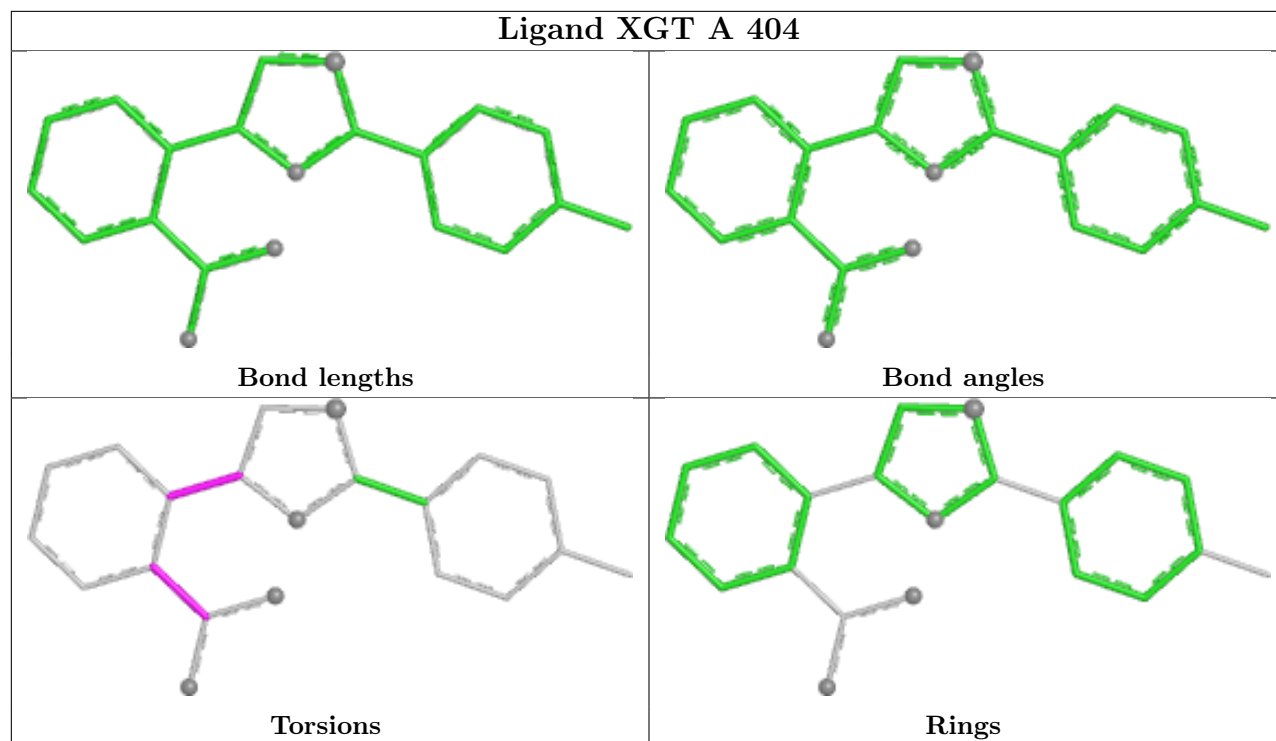
Mol	Chain	Res	Type	Atoms
5	B	403	XGT	C11-C10-C13-C18
5	A	404	XGT	N9-C10-C13-C18
5	B	403	XGT	C17-C18-C19-O20
5	A	404	XGT	C17-C18-C19-O21
5	A	404	XGT	C17-C18-C19-O20
5	B	403	XGT	C17-C18-C19-O21
2	A	401	ADP	PA-O3A-PB-O1B
5	B	403	XGT	C13-C18-C19-O20
5	A	404	XGT	C13-C18-C19-O21
5	A	404	XGT	N9-C10-C13-C14
5	B	403	XGT	N9-C10-C13-C14
5	B	403	XGT	C13-C18-C19-O21
5	A	404	XGT	C13-C18-C19-O20

There are no ring outliers.

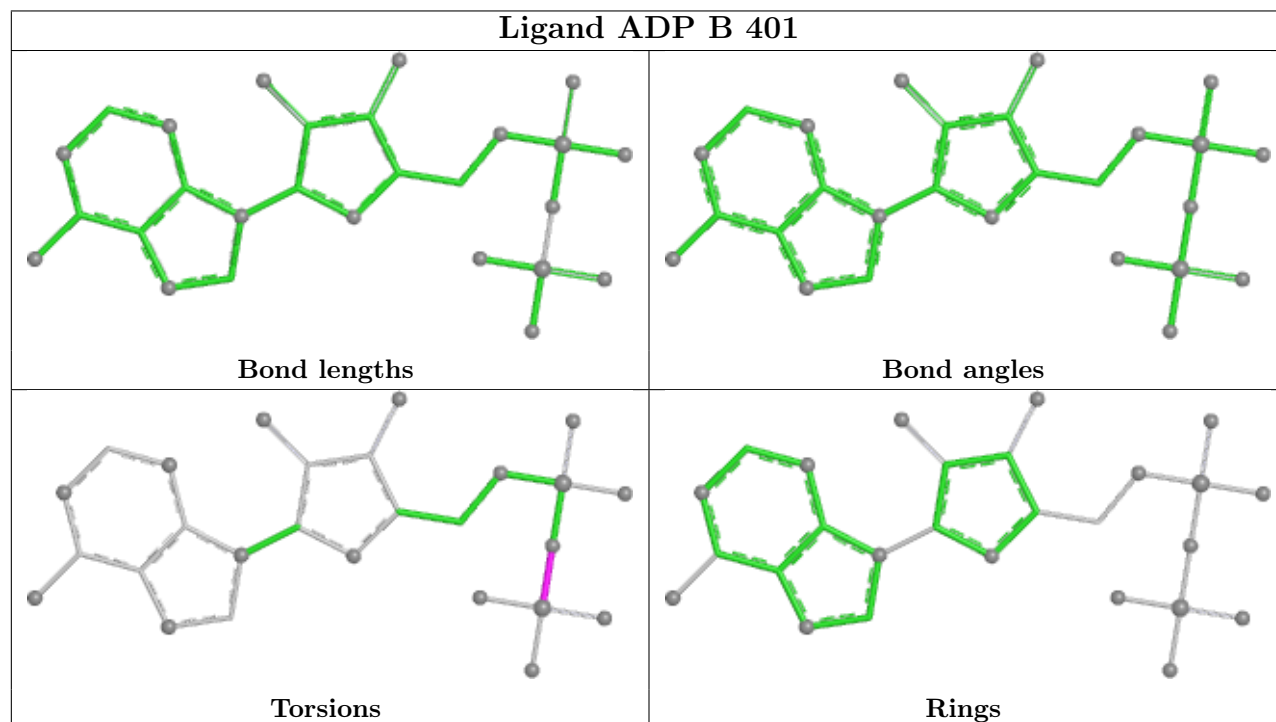
No monomer is involved in short contacts.

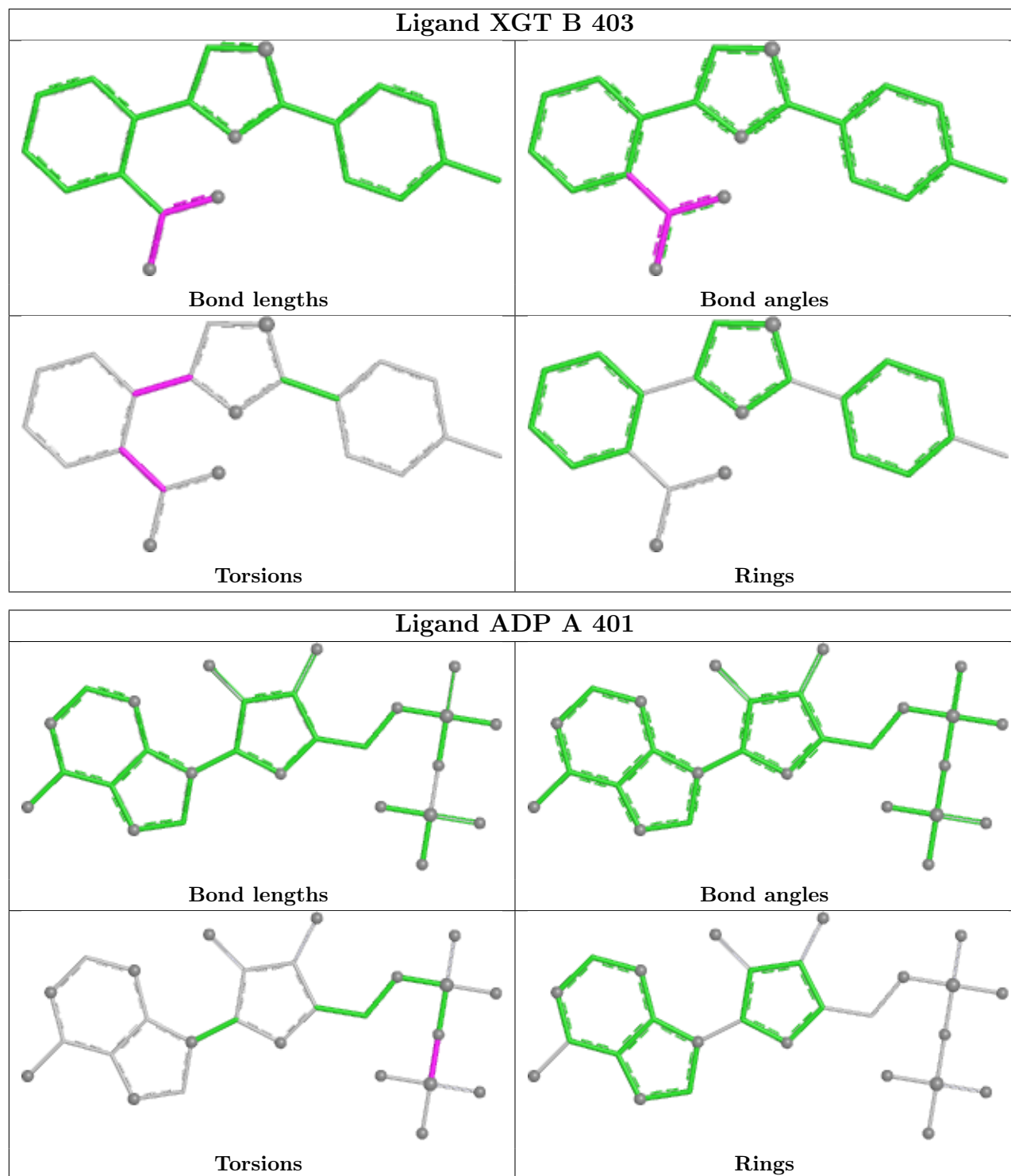
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 XGT A 404



Ligand ADP B 401





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	322/388 (82%)	0.58	16 (4%) 34 26	59, 87, 150, 163	0
1	B	314/388 (80%)	1.47	81 (25%) 1 1	119, 163, 187, 196	0
All	All	636/776 (81%)	1.02	97 (15%) 5 4	59, 143, 183, 196	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	281	ILE	6.1
1	B	217	GLY	5.7
1	B	251	MET	5.6
1	B	247	LEU	5.6
1	B	211	LEU	5.3
1	B	205	VAL	5.3
1	B	47	TYR	5.2
1	B	239	LEU	5.2
1	B	246	LEU	4.8
1	B	368	LEU	4.5
1	B	42	PRO	4.4
1	B	155	LEU	4.3
1	B	224	ILE	4.2
1	B	346	TYR	4.1
1	B	283	ILE	3.9
1	B	238	PRO	3.9
1	B	200	ILE	3.9
1	B	356	LEU	3.5
1	B	207	PHE	3.5
1	B	216	TYR	3.4
1	B	234	PHE	3.4
1	B	265	VAL	3.4
1	B	243	LEU	3.4
1	B	269	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	294	ASP	3.4
1	B	193	THR	3.3
1	B	206	ASP	3.3
1	B	268	VAL	3.3
1	B	222	VAL	3.2
1	B	204	TRP	3.2
1	B	225	ASN	3.2
1	A	106	LEU	3.0
1	A	47	TYR	3.0
1	B	43	THR	2.9
1	B	236	PRO	2.9
1	A	99	TYR	2.9
1	B	350	LEU	2.9
1	B	366	VAL	2.9
1	B	208	ALA	2.8
1	B	235	ILE	2.7
1	B	148	LYS	2.7
1	A	63	TYR	2.7
1	A	128	LEU	2.7
1	B	228	VAL	2.7
1	B	298	VAL	2.7
1	B	345	ALA	2.7
1	A	42	PRO	2.6
1	A	112	ILE	2.6
1	A	40	LEU	2.6
1	B	338	PHE	2.6
1	B	263	TYR	2.5
1	B	201	ILE	2.5
1	A	46	LEU	2.5
1	B	279	LEU	2.5
1	B	136	VAL	2.5
1	B	255	MET	2.5
1	B	46	LEU	2.5
1	B	220	PRO	2.5
1	B	266	PRO	2.4
1	B	26	GLN	2.4
1	B	354	LEU	2.4
1	A	102	ALA	2.4
1	B	229	ALA	2.4
1	B	270	ILE	2.3
1	B	360	GLN	2.3
1	B	168	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	86	ILE	2.3
1	B	230	ALA	2.3
1	B	336	PHE	2.3
1	B	100	ILE	2.2
1	B	164	LEU	2.2
1	B	38	VAL	2.2
1	A	373	ILE	2.2
1	B	190	ILE	2.2
1	B	367	TYR	2.2
1	B	185	PRO	2.2
1	B	54	GLY	2.2
1	B	68	LEU	2.2
1	B	129	LEU	2.2
1	B	172	MET	2.2
1	A	207	PHE	2.2
1	B	306	ALA	2.2
1	A	141	GLU	2.2
1	B	212	CYS	2.2
1	A	173	LEU	2.2
1	B	125	VAL	2.2
1	B	272	ILE	2.2
1	B	195	LEU	2.1
1	B	144	ARG	2.1
1	A	333	MET	2.1
1	B	223	ARG	2.1
1	B	370	LEU	2.1
1	B	210	ARG	2.0
1	B	359	LEU	2.0
1	A	222	VAL	2.0
1	B	137	THR	2.0
1	B	290	ILE	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

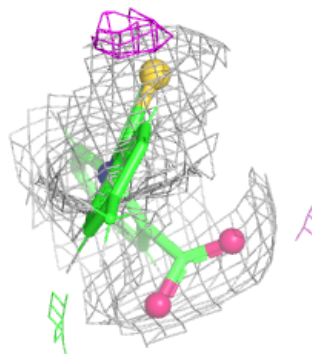
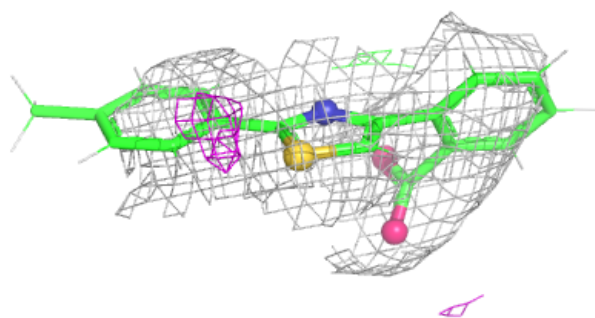
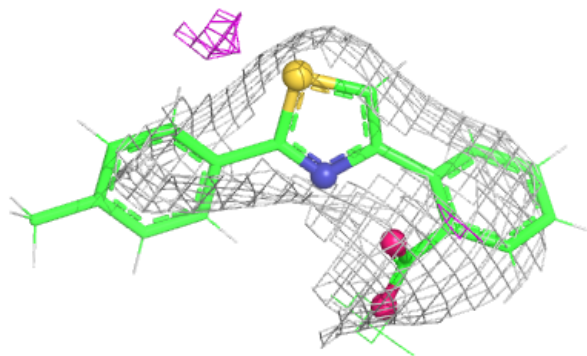
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SO4	A	407	5/5	0.58	0.15	174,174,174,174	0
6	SO4	A	408	5/5	0.70	0.11	150,150,150,151	0
6	SO4	B	404	5/5	0.76	0.11	173,173,173,173	0
3	MG	B	402	1/1	0.77	0.27	153,153,153,153	0
6	SO4	A	405	5/5	0.82	0.08	188,188,188,188	0
3	MG	A	402	1/1	0.86	0.29	64,64,64,64	0
5	XGT	B	403	21/21	0.90	0.12	145,147,177,179	12
2	ADP	B	401	27/27	0.91	0.12	163,163,165,165	0
6	SO4	A	406	5/5	0.94	0.07	139,139,139,139	0
4	K	A	403	1/1	0.94	0.11	184,184,184,184	0
2	ADP	A	401	27/27	0.95	0.09	74,76,79,79	0
5	XGT	A	404	21/21	0.96	0.09	54,86,88,88	12

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.

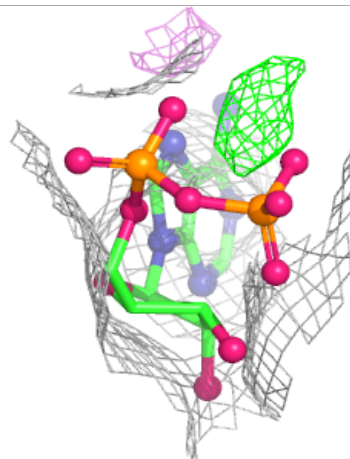
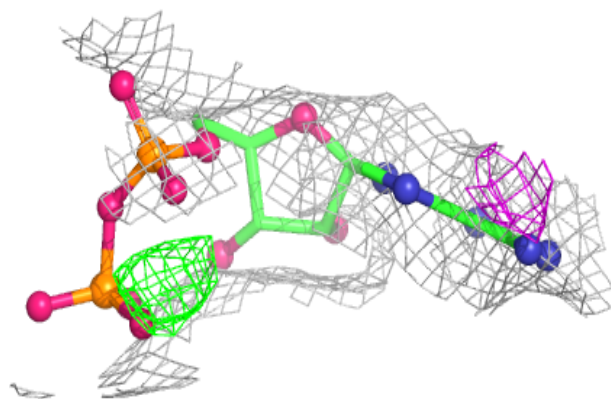
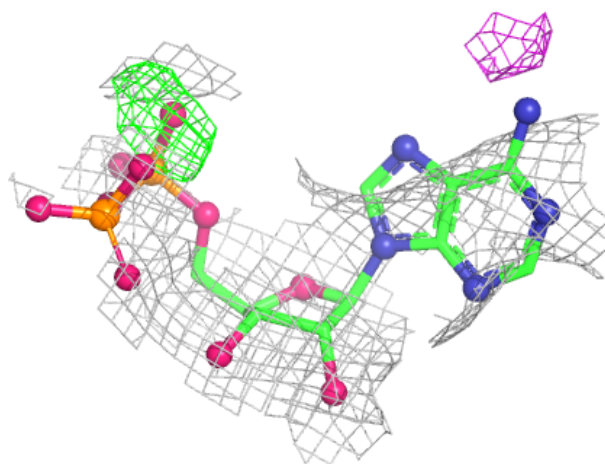
Electron density around XGT B 403:

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



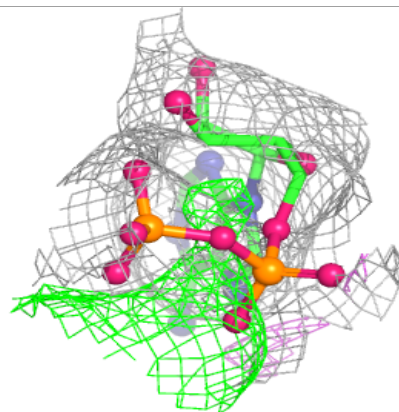
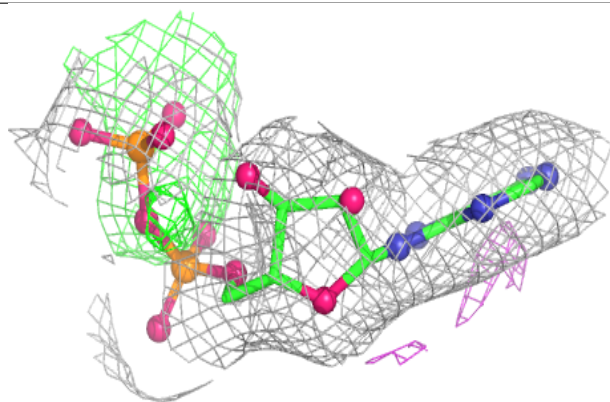
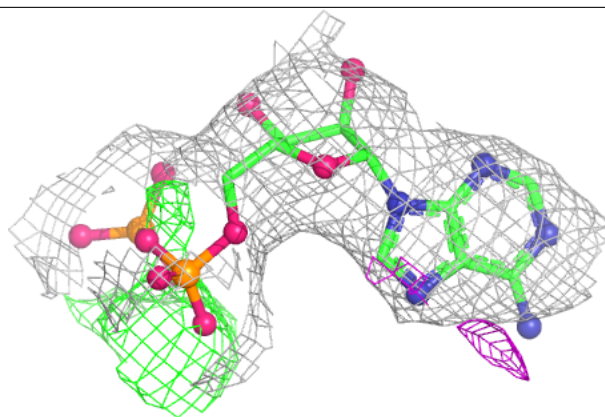
Electron density around ADP B 401:

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

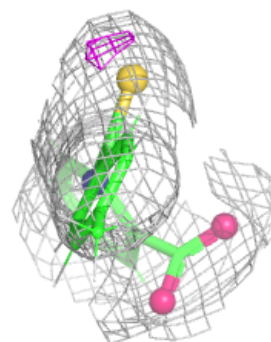
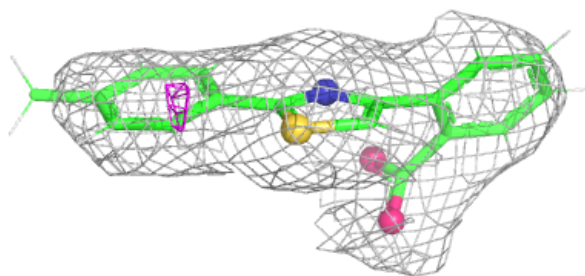
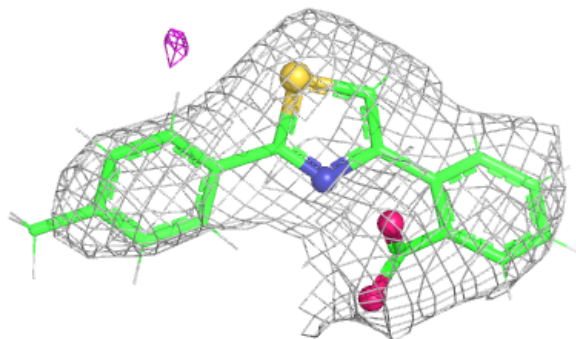


Electron density around ADP A 401:

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

**Electron density around XGT A 404:**

$2mF_o-DF_c$ (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.