



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:27 AM UTC

PDB ID : 8FHF / pdb_00008fhf
Title : Crystal structure of PPARgamma ligand-binding domain in complex with ZINC5672437
Authors : Shang, J.; Kojetin, D.J.
Deposited on : 2022-12-14
Resolution : 2.10 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

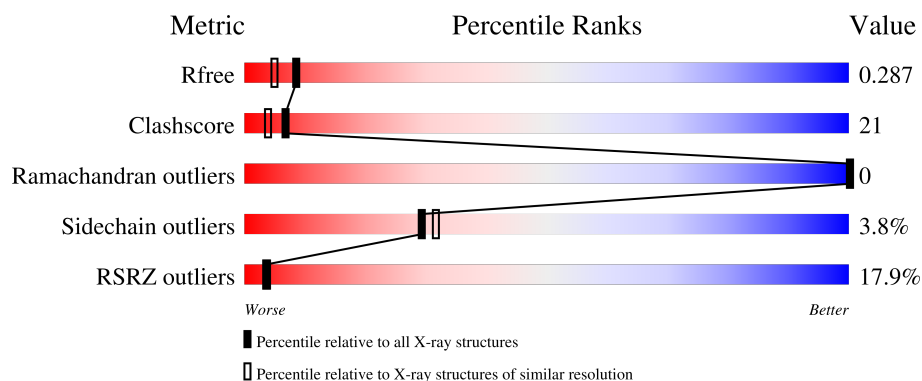
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	275	20% 67% 24% 6%
1	B	275	13% 65% 23% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	KNA	A	502	-	-	X	-
3	KNA	B	502	-	-	X	-

2 Entry composition [i](#)

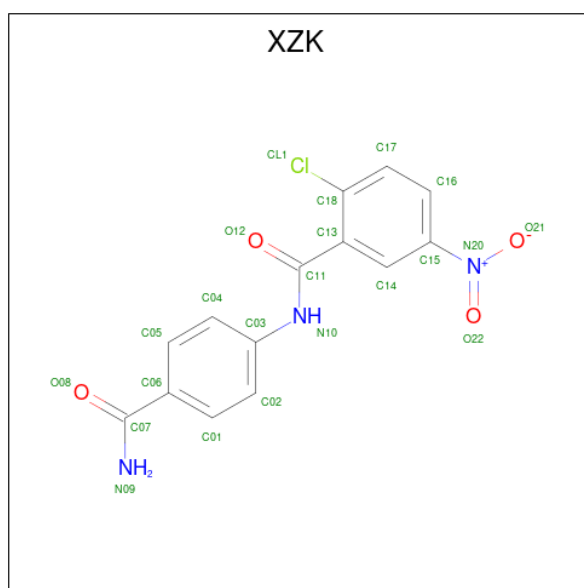
There are 4 unique types of molecules in this entry. The entry contains 4355 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Peroxisome proliferator-activated receptor gamma.

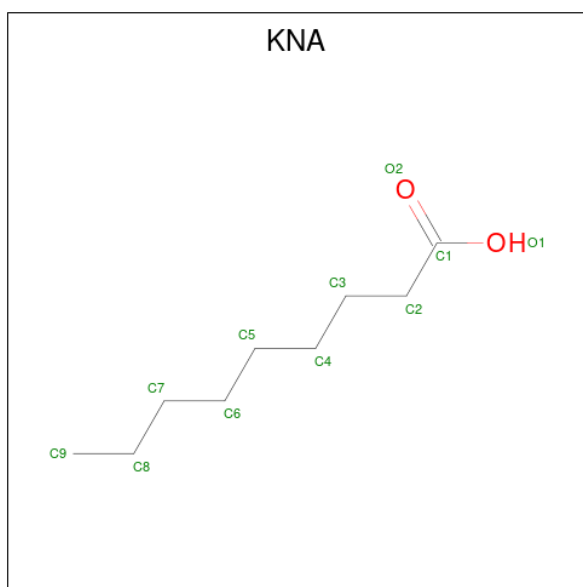
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			2063	1331	336	386	10			
1	B	246	Total	C	N	O	S	0	0	0
			1970	1273	322	366	9			

- Molecule 2 is N-(4-carbamoylphenyl)-2-chloro-5-nitrobenzamide (CCD ID: XZK) (formula: $C_{14}H_{10}ClN_3O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	14	3	4		
2	B	1	Total	C	N	O	0	0
			21	14	3	4		

- Molecule 3 is nonanoic acid (CCD ID: KNA) (formula: $C_9H_{18}O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	9	2		
3	B	1	Total	C	O	0	0
			11	9	2		

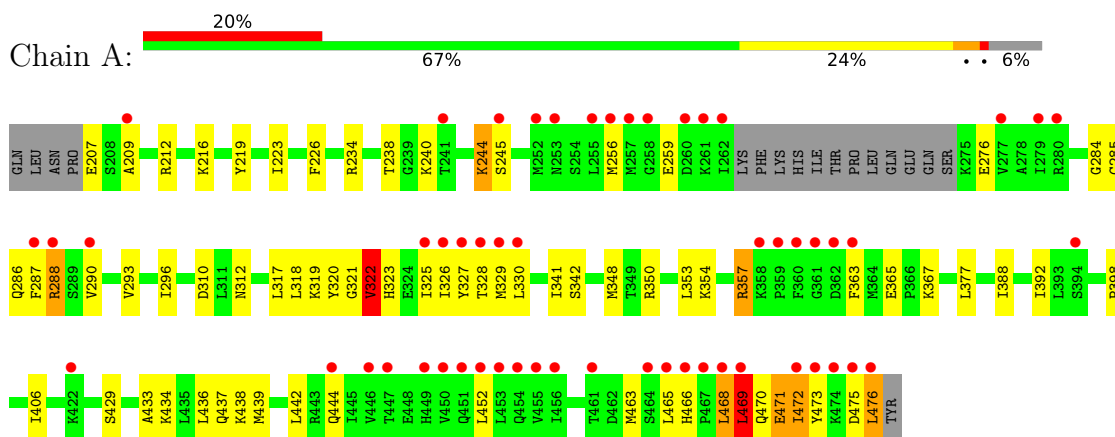
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	140	Total	O	0	0
			140	140		
4	B	118	Total	O	0	0
			118	118		

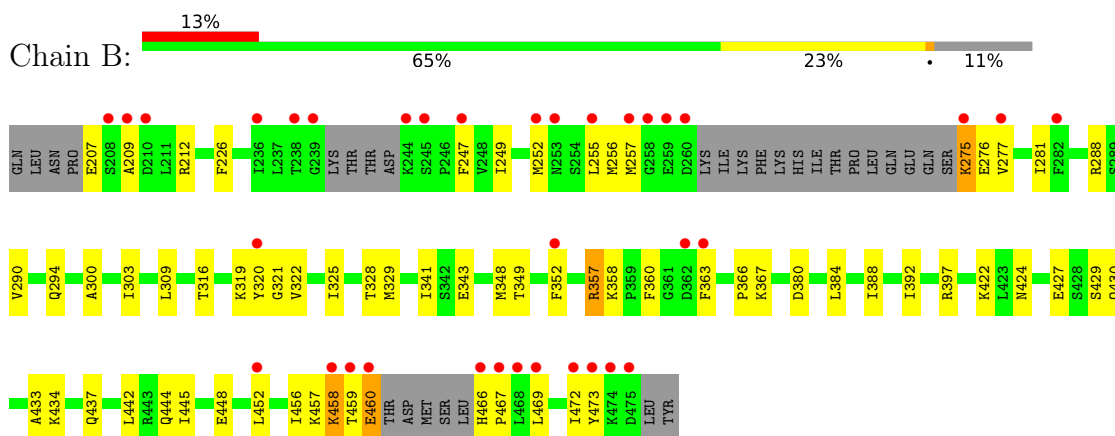
3 Residue-property plots [i](#)

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: Peroxisome proliferator-activated receptor gamma



- Molecule 1: Peroxisome proliferator-activated receptor gamma



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	91.64Å 60.70Å 117.08Å 90.00° 102.51° 90.00°	Depositor
Resolution (Å)	38.89 – 2.10 38.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.5 (38.89-2.10) 97.4 (38.89-2.10)	Depositor EDS
R_{merge}	0.02	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.233 , 0.285 0.244 , 0.287	Depositor DCC
R_{free} test set	2000 reflections (5.42%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4355	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.63	6/2096 (0.3%)	0.67	1/2823 (0.0%)
1	B	0.60	3/2001 (0.1%)	0.71	1/2692 (0.0%)
All	All	0.62	9/4097 (0.2%)	0.69	2/5515 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	322	VAL	C-O	-8.69	1.14	1.24
1	A	321	GLY	C-O	-7.14	1.15	1.24
1	B	322	VAL	C-O	-6.97	1.16	1.24
1	A	472	ILE	N-CA	-6.15	1.39	1.46
1	B	309	LEU	C-O	-5.78	1.16	1.23
1	A	468	LEU	C-O	-5.74	1.17	1.24
1	B	319	LYS	C-O	-5.37	1.17	1.24
1	A	322	VAL	N-CA	-5.24	1.40	1.46
1	A	469	LEU	C-O	-5.14	1.17	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	360	PHE	N-CA-C	5.72	118.46	111.82
1	A	322	VAL	N-CA-C	5.39	116.12	110.62

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	2063	0	2125	113	0
1	B	1970	0	2021	63	0
2	A	21	0	0	3	0
2	B	21	0	0	0	0
3	A	11	0	17	29	0
3	B	11	0	17	10	0
4	A	140	0	0	9	0
4	B	118	0	0	11	0
All	All	4355	0	4180	177	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 21.

All (177) 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:288:ARG:CZ	3:A:502:KNA:H6A	1.52	1.36
1:A:327:TYR:CZ	1:A:367:LYS:HE2	1.84	1.11
1:A:288:ARG:CZ	3:A:502:KNA:C6	2.33	1.06
1:A:288:ARG:NE	3:A:502:KNA:H6A	1.71	1.04
1:A:290:VAL:HG12	1:A:469:LEU:CD2	1.91	1.01
1:A:327:TYR:OH	1:A:367:LYS:HE2	1.65	0.94
1:A:325:ILE:CD1	1:A:392:ILE:HG12	1.99	0.92
1:A:290:VAL:HG12	1:A:469:LEU:HD21	1.53	0.89
1:A:296:ILE:CD1	1:A:326:ILE:CG1	2.52	0.88
1:A:296:ILE:HD11	1:A:326:ILE:HG12	1.54	0.87
1:A:288:ARG:NH2	3:A:502:KNA:H6A	1.90	0.86
1:A:288:ARG:NH1	3:A:502:KNA:C4	2.38	0.86
1:A:296:ILE:CD1	1:A:326:ILE:HG12	2.05	0.86
1:B:316:THR:HG23	1:B:320:TYR:HE1	1.42	0.85
1:A:288:ARG:NH1	3:A:502:KNA:C5	2.40	0.84
1:B:316:THR:HG23	1:B:320:TYR:CE1	2.15	0.81
1:A:327:TYR:CZ	1:A:367:LYS:CE	2.63	0.79
1:A:288:ARG:HG2	3:A:502:KNA:H8A	1.63	0.79
1:A:288:ARG:HG2	3:A:502:KNA:H7	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ARG:HG2	3:A:502:KNA:C7	2.14	0.78
1:A:296:ILE:CD1	1:A:326:ILE:HG13	2.15	0.77
1:A:288:ARG:NH1	3:A:502:KNA:H4	1.98	0.77
1:A:288:ARG:NH1	3:A:502:KNA:H5A	2.01	0.76
1:A:288:ARG:HH11	3:A:502:KNA:H4	1.49	0.76
1:A:259:GLU:OE1	1:A:259:GLU:N	2.18	0.76
1:A:296:ILE:HD11	1:A:326:ILE:CG1	2.17	0.72
1:A:288:ARG:CZ	3:A:502:KNA:C5	2.66	0.72
1:A:471:GLU:O	1:A:471:GLU:HG3	1.90	0.72
1:B:341:ILE:HA	3:B:502:KNA:H7A	1.71	0.71
1:A:468:LEU:HD12	1:A:468:LEU:O	1.91	0.71
1:A:363:PHE:HD2	1:A:452:LEU:HD12	1.57	0.70
1:A:288:ARG:HG2	3:A:502:KNA:C8	2.21	0.70
1:A:320:TYR:CZ	1:A:398:PRO:HG2	2.26	0.70
1:A:288:ARG:HB3	2:A:501:XZK:C02	2.21	0.70
1:A:240:LYS:HG3	4:A:601:HOH:O	1.91	0.69
1:A:290:VAL:HG12	1:A:469:LEU:HD22	1.74	0.69
1:A:325:ILE:HG23	1:A:388:ILE:HD12	1.74	0.69
1:A:325:ILE:CD1	1:A:392:ILE:CG1	2.72	0.68
1:A:325:ILE:HD11	1:A:392:ILE:HG12	1.74	0.68
1:A:290:VAL:CG1	1:A:469:LEU:HD21	2.25	0.67
1:A:256:MET:HB3	4:A:661:HOH:O	1.94	0.67
1:B:288:ARG:HE	3:B:502:KNA:H5A	1.61	0.65
1:A:323:HIS:CE1	1:A:472:ILE:HG21	2.30	0.65
1:A:325:ILE:HG12	1:A:388:ILE:HG23	1.79	0.64
1:A:288:ARG:NH2	3:A:502:KNA:C6	2.55	0.63
1:A:288:ARG:HH11	3:A:502:KNA:C4	2.06	0.63
1:B:458:LYS:O	1:B:458:LYS:HE3	1.98	0.62
1:B:276:GLU:OE2	1:B:357:ARG:HD2	1.98	0.62
1:A:219:TYR:CZ	1:A:223:ILE:HD11	2.34	0.62
1:B:357:ARG:HG2	4:B:625:HOH:O	1.98	0.62
1:A:323:HIS:NE2	1:A:473:TYR:OH	2.21	0.62
1:B:247:PHE:HE2	1:B:257:MET:HE2	1.64	0.62
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.30	0.61
1:A:287:PHE:O	1:A:290:VAL:HG22	2.01	0.60
1:A:466:HIS:HB3	1:A:469:LEU:HD23	1.82	0.60
1:A:240:LYS:N	4:A:601:HOH:O	2.34	0.60
1:B:367:LYS:NZ	4:B:607:HOH:O	2.32	0.60
1:A:296:ILE:HD13	1:A:326:ILE:HG13	1.84	0.60
1:A:290:VAL:CG1	1:A:469:LEU:CD2	2.75	0.60
1:B:252:MET:SD	1:B:277:VAL:HG21	2.42	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:ARG:NH2	1:B:460:GLU:OE2	2.35	0.59
1:A:325:ILE:HD12	1:A:392:ILE:HG12	1.80	0.59
1:B:444:GLN:HE21	1:B:444:GLN:HA	1.67	0.59
1:A:288:ARG:CD	3:A:502:KNA:H6A	2.32	0.59
1:A:473:TYR:OH	2:A:501:XZK:O22	2.21	0.59
1:A:341:ILE:HB	3:A:502:KNA:H6	1.84	0.58
1:B:357:ARG:CG	4:B:625:HOH:O	2.50	0.58
1:A:348:MET:HE3	1:A:353:LEU:HD21	1.86	0.57
1:A:293:VAL:HG22	1:A:326:ILE:HD11	1.87	0.57
1:A:328:THR:OG1	1:A:442:LEU:HD11	2.05	0.57
1:B:459:THR:HG21	4:B:682:HOH:O	2.05	0.56
1:B:430:GLN:O	1:B:434:LYS:HG3	2.06	0.56
3:A:502:KNA:H2A	3:A:502:KNA:H7A	1.88	0.55
1:B:363:PHE:HD2	1:B:452:LEU:HD22	1.71	0.55
4:A:632:HOH:O	1:B:444:GLN:HG2	2.05	0.55
1:B:321:GLY:O	1:B:325:ILE:HD12	2.06	0.55
1:A:444:GLN:HE22	1:B:444:GLN:NE2	2.05	0.54
1:B:349:THR:HG23	1:B:352:PHE:H	1.72	0.54
1:B:341:ILE:HB	3:B:502:KNA:H5	1.90	0.54
1:A:388:ILE:O	1:A:392:ILE:HG13	2.08	0.53
1:A:433:ALA:O	1:A:437:GLN:HG3	2.09	0.53
1:B:247:PHE:CE2	1:B:257:MET:HE2	2.43	0.53
1:A:365:GLU:HG2	4:A:685:HOH:O	2.08	0.53
2:A:501:XZK:C01	3:A:502:KNA:H9A	2.39	0.52
1:A:436:LEU:O	1:A:439:MET:HB2	2.10	0.52
1:A:325:ILE:CG1	1:A:388:ILE:HG23	2.39	0.52
1:B:320:TYR:O	1:B:397:ARG:CZ	2.57	0.52
1:A:293:VAL:CG2	1:A:326:ILE:HD11	2.40	0.52
1:B:288:ARG:HG3	3:B:502:KNA:H6	1.91	0.52
1:A:276:GLU:OE2	1:A:357:ARG:NE	2.40	0.51
1:B:249:ILE:HD12	1:B:255:LEU:HA	1.91	0.51
1:B:328:THR:OG1	1:B:442:LEU:HD11	2.11	0.51
1:B:207:GLU:HG3	1:B:209:ALA:H	1.76	0.51
1:B:276:GLU:OE2	1:B:357:ARG:CD	2.59	0.50
1:B:343:GLU:H	3:B:502:KNA:H9B	1.76	0.50
1:B:466:HIS:N	1:B:467:PRO:HD3	2.26	0.50
1:A:327:TYR:CE2	1:A:367:LYS:CE	2.94	0.50
1:A:317:LEU:HD21	1:A:406:ILE:HD13	1.94	0.50
1:A:328:THR:HG1	1:A:442:LEU:HD11	1.76	0.50
1:A:284:GLY:O	1:A:288:ARG:HB2	2.11	0.50
1:A:207:GLU:HG3	1:A:209:ALA:H	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:ARG:O	1:A:216:LYS:HG2	2.13	0.49
1:A:377:LEU:HD11	1:A:438:LYS:HG3	1.95	0.49
1:A:286:GLN:HE22	1:A:463:MET:HE3	1.78	0.49
3:A:502:KNA:C7	3:A:502:KNA:H2A	2.42	0.49
1:A:342:SER:H	3:A:502:KNA:H6	1.77	0.48
1:A:325:ILE:HD11	1:A:388:ILE:O	2.14	0.48
1:A:465:LEU:HG	1:A:470:GLN:HG3	1.95	0.48
1:A:319:LYS:NZ	1:A:471:GLU:O	2.47	0.48
1:B:444:GLN:O	1:B:448:GLU:HG3	2.13	0.48
1:B:472:ILE:HG22	1:B:473:TYR:CD1	2.49	0.48
1:B:427:GLU:OE2	4:B:601:HOH:O	2.20	0.47
1:B:349:THR:CG2	1:B:352:PHE:H	2.25	0.47
1:A:437:GLN:HG2	4:B:687:HOH:O	2.14	0.47
1:A:437:GLN:HA	4:B:687:HOH:O	2.14	0.47
1:B:366:PRO:HB3	4:B:605:HOH:O	2.13	0.47
1:A:363:PHE:CD2	1:A:452:LEU:HD12	2.44	0.47
1:B:380:ASP:OD2	1:B:424:ASN:ND2	2.35	0.47
1:A:341:ILE:HB	3:A:502:KNA:C6	2.45	0.47
1:B:288:ARG:HG3	3:B:502:KNA:C6	2.45	0.46
1:A:296:ILE:HD12	1:A:326:ILE:CG1	2.42	0.46
1:B:275:LYS:HE2	4:B:709:HOH:O	2.14	0.46
1:A:323:HIS:NE2	1:A:473:TYR:CZ	2.83	0.46
1:B:325:ILE:HD11	1:B:392:ILE:HG13	1.98	0.46
1:B:226:PHE:CG	1:B:329:MET:HE1	2.50	0.46
1:B:357:ARG:HH22	1:B:460:GLU:CD	2.24	0.46
1:A:226:PHE:CE1	1:A:329:MET:HE1	2.51	0.45
1:A:429:SER:O	1:A:434:LYS:HE3	2.16	0.45
1:A:244:LYS:HA	1:A:244:LYS:HD3	1.66	0.45
1:A:326:ILE:O	4:A:602:HOH:O	2.20	0.45
1:A:476:LEU:HD12	1:A:476:LEU:HA	1.75	0.45
1:B:422:LYS:NZ	4:B:608:HOH:O	2.33	0.45
1:B:429:SER:O	1:B:434:LYS:NZ	2.44	0.45
1:B:384:LEU:O	1:B:388:ILE:HG23	2.15	0.45
1:B:343:GLU:H	3:B:502:KNA:C9	2.30	0.45
1:A:327:TYR:OH	1:A:367:LYS:CE	2.51	0.44
1:B:288:ARG:HG3	3:B:502:KNA:H5A	1.99	0.44
1:B:458:LYS:HE3	1:B:458:LYS:HB3	1.79	0.44
1:A:348:MET:HE1	3:A:502:KNA:H3	2.00	0.44
1:B:433:ALA:O	1:B:437:GLN:HG3	2.17	0.44
1:A:323:HIS:CE1	1:A:472:ILE:CG2	3.01	0.44
1:B:320:TYR:HB2	1:B:397:ARG:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:LEU:O	1:A:322:VAL:HB	2.17	0.43
1:A:322:VAL:O	1:A:326:ILE:HG13	2.19	0.43
1:B:348:MET:HE3	1:B:348:MET:HB2	1.80	0.43
1:B:212:ARG:HA	1:B:212:ARG:HD2	1.76	0.43
1:A:212:ARG:HA	1:A:212:ARG:HD2	1.67	0.43
1:A:319:LYS:HE3	1:A:476:LEU:O	2.18	0.43
1:A:323:HIS:HE1	1:A:472:ILE:HG21	1.80	0.43
1:A:325:ILE:CD1	1:A:388:ILE:HG23	2.48	0.43
1:A:310:ASP:OD2	1:A:312:ASN:ND2	2.52	0.43
1:B:288:ARG:CD	3:B:502:KNA:H8A	2.49	0.42
1:B:442:LEU:O	1:B:445:ILE:HB	2.19	0.42
1:A:341:ILE:HA	3:A:502:KNA:H8	2.01	0.42
1:B:276:GLU:CD	1:B:357:ARG:HH11	2.27	0.42
1:A:354:LYS:HB2	1:A:354:LYS:HE3	1.84	0.42
1:A:288:ARG:CG	3:A:502:KNA:H7	2.41	0.42
1:A:350:ARG:NH2	4:A:603:HOH:O	2.22	0.42
1:A:444:GLN:NE2	1:B:444:GLN:NE2	2.67	0.42
1:B:277:VAL:O	1:B:281:ILE:HG12	2.19	0.42
1:B:466:HIS:O	1:B:469:LEU:HB2	2.19	0.42
1:A:288:ARG:NE	3:A:502:KNA:C6	2.61	0.42
1:A:234:ARG:HA	1:A:234:ARG:HD3	1.79	0.41
1:B:288:ARG:HG3	3:B:502:KNA:C5	2.50	0.41
1:B:452:LEU:HD23	1:B:456:ILE:HD11	2.01	0.41
1:A:238:THR:OG1	4:A:601:HOH:O	2.20	0.41
1:A:288:ARG:CZ	3:A:502:KNA:C4	2.97	0.41
1:B:226:PHE:CD1	1:B:329:MET:HE1	2.56	0.41
1:A:398:PRO:O	4:A:604:HOH:O	2.22	0.41
1:B:300:ALA:HA	1:B:303:ILE:HD12	2.02	0.41
1:A:286:GLN:NE2	1:A:463:MET:HE3	2.36	0.41
1:A:323:HIS:CE1	1:A:473:TYR:CE2	3.08	0.41
1:A:326:ILE:O	1:A:330:LEU:HB2	2.21	0.41
1:B:255:LEU:HD21	1:B:281:ILE:HD11	2.02	0.41
1:B:294:GLN:HB3	4:B:692:HOH:O	2.20	0.41
1:A:285:CYS:CB	3:A:502:KNA:H2	2.52	0.40
1:A:296:ILE:HD12	1:A:326:ILE:HG12	1.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	254/275 (92%)	246 (97%)	8 (3%)	0	100	100
1	B	238/275 (86%)	232 (98%)	6 (2%)	0	100	100
All	All	492/550 (90%)	478 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	231/248 (93%)	222 (96%)	9 (4%)	28	31
1	B	219/248 (88%)	211 (96%)	8 (4%)	30	33
All	All	450/496 (91%)	433 (96%)	17 (4%)	29	32

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	244	LYS
1	A	245	SER
1	A	288	ARG
1	A	322	VAL
1	A	357	ARG
1	A	469	LEU
1	A	471	GLU
1	A	475	ASP

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Mol	Chain	Res	Type
1	A	476	LEU
1	B	256	MET
1	B	275	LYS
1	B	290	VAL
1	B	357	ARG
1	B	358	LYS
1	B	457	LYS
1	B	458	LYS
1	B	460	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	294	GLN
1	A	345	GLN
1	A	410	GLN
1	A	444	GLN
1	B	437	GLN
1	B	444	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	XZK	B	501	1	22,22,23	4.22	10 (45%)	28,30,32	1.08	2 (7%)
3	KNA	A	502	-	10,10,10	0.66	0	10,10,10	0.97	0
2	XZK	A	501	1	22,22,23	4.11	10 (45%)	28,30,32	1.22	3 (10%)
3	KNA	B	502	-	10,10,10	0.83	0	10,10,10	1.24	2 (20%)

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	XZK	B	501	1	-	6/14/16/16	0/2/2/2
3	KNA	A	502	-	-	5/8/8/8	-
2	XZK	A	501	1	-	4/14/16/16	0/2/2/2
3	KNA	B	502	-	-	4/8/8/8	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	XZK	O22-N20	15.01	1.48	1.22
2	A	501	XZK	O22-N20	14.62	1.48	1.22
2	B	501	XZK	C07-N09	8.41	1.48	1.33
2	A	501	XZK	C07-N09	8.32	1.48	1.33
2	A	501	XZK	O08-C07	-5.20	1.14	1.24
2	B	501	XZK	O08-C07	-4.92	1.15	1.24
2	B	501	XZK	C05-C04	3.51	1.44	1.38
2	A	501	XZK	C05-C04	3.39	1.44	1.38
2	A	501	XZK	C17-C16	3.35	1.44	1.38
2	B	501	XZK	C02-C01	3.12	1.43	1.38
2	B	501	XZK	C17-C16	3.05	1.44	1.38
2	A	501	XZK	C06-C07	2.83	1.54	1.50
2	B	501	XZK	C06-C07	2.83	1.54	1.50
2	A	501	XZK	C02-C01	2.67	1.43	1.38
2	B	501	XZK	C18-C13	-2.60	1.35	1.39
2	B	501	XZK	C16-C15	-2.52	1.34	1.38
2	B	501	XZK	C11-N10	2.39	1.42	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	XZK	C18-C13	-2.37	1.35	1.39
2	A	501	XZK	C16-C15	-2.32	1.34	1.38
2	A	501	XZK	C11-N10	2.27	1.42	1.35

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	XZK	C06-C07-N09	3.14	121.61	117.74
2	A	501	XZK	C16-C15-N20	2.77	121.76	119.34
2	A	501	XZK	C05-C06-C01	2.49	121.74	118.57
3	B	502	KNA	O1-C1-C2	2.31	121.30	114.00
2	A	501	XZK	C04-C03-C02	2.24	122.01	119.04
3	B	502	KNA	C3-C2-C1	-2.23	108.68	114.51
2	B	501	XZK	C05-C06-C01	2.03	121.15	118.57

There are no chirality outliers.

All (19) torsion outliers are listed below:

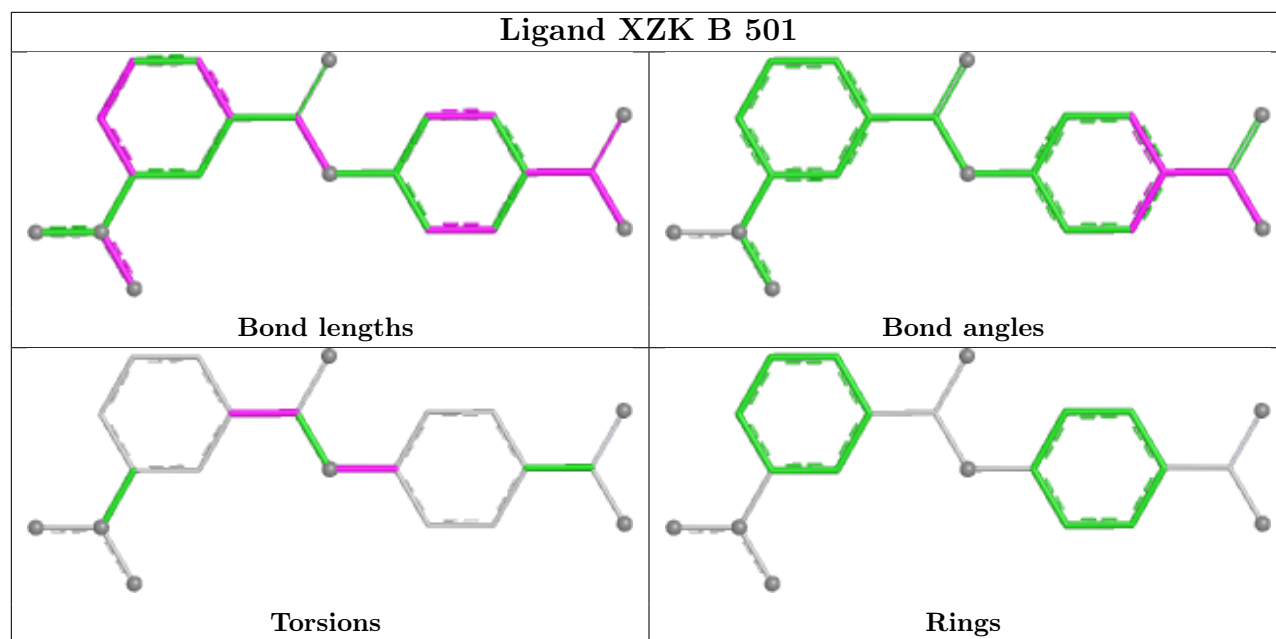
Mol	Chain	Res	Type	Atoms
2	B	501	XZK	N10-C11-C13-C14
2	B	501	XZK	O12-C11-C13-C14
2	B	501	XZK	N10-C11-C13-C18
2	B	501	XZK	O12-C11-C13-C18
2	A	501	XZK	N10-C11-C13-C14
3	B	502	KNA	C3-C4-C5-C6
2	A	501	XZK	O12-C11-C13-C14
3	B	502	KNA	C1-C2-C3-C4
2	A	501	XZK	N10-C11-C13-C18
3	B	502	KNA	C5-C6-C7-C8
2	A	501	XZK	O12-C11-C13-C18
3	B	502	KNA	C2-C3-C4-C5
3	A	502	KNA	C3-C4-C5-C6
3	A	502	KNA	C2-C3-C4-C5
3	A	502	KNA	C1-C2-C3-C4
2	B	501	XZK	C04-C03-N10-C11
3	A	502	KNA	C5-C6-C7-C8
3	A	502	KNA	C6-C7-C8-C9
2	B	501	XZK	C02-C03-N10-C11

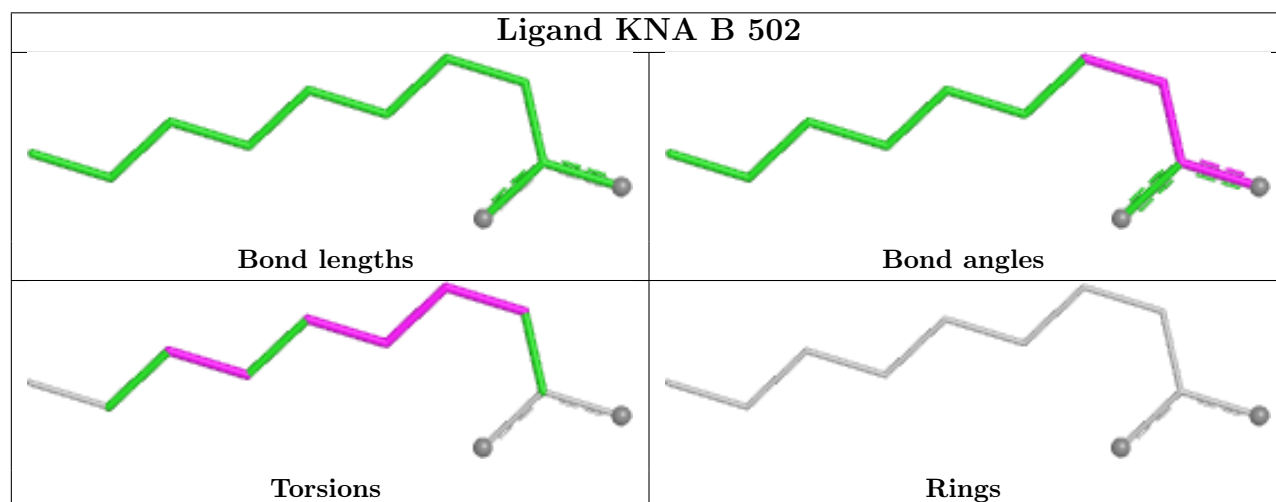
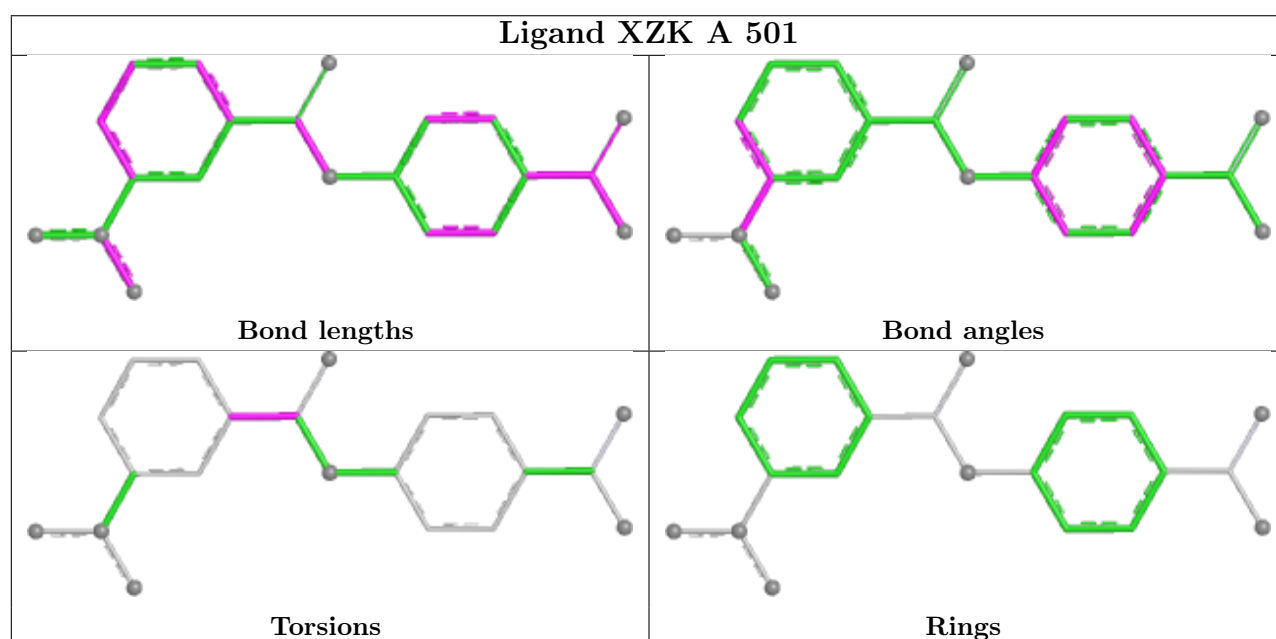
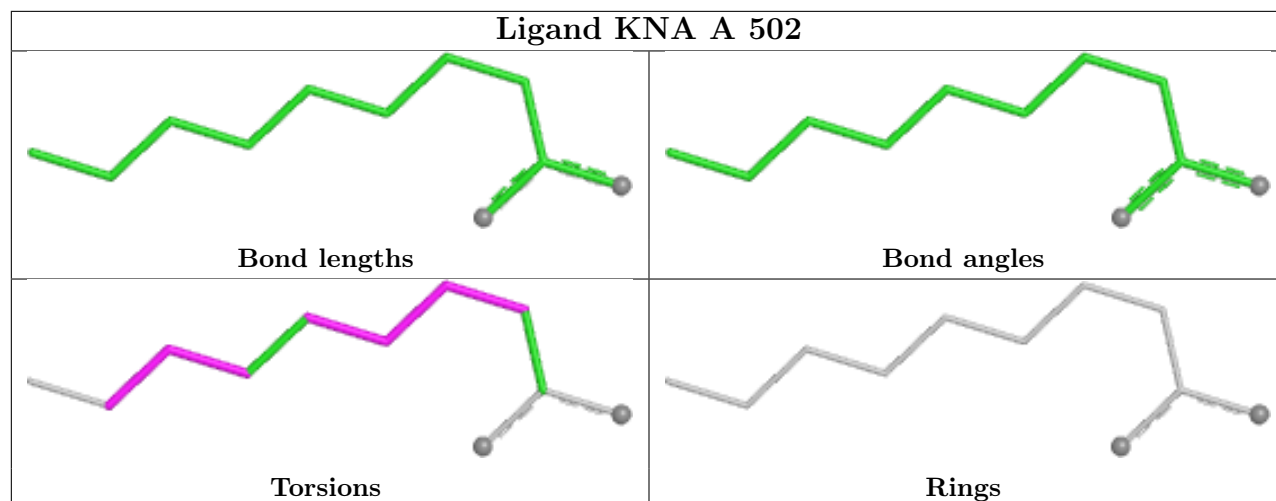
There are no ring outliers.

3 monomers are involved in 41 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	KNA	29	0
2	A	501	XZK	3	0
3	B	502	KNA	10	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

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	258/275 (93%)	1.03	55 (21%) 2 2	10, 26, 54, 81	0
1	B	246/275 (89%)	0.90	35 (14%) 6 6	10, 27, 56, 76	0
All	All	504/550 (91%)	0.97	90 (17%) 3 4	10, 26, 56, 81	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	476	LEU	8.7
1	A	327	TYR	6.8
1	A	326	ILE	6.1
1	B	239	GLY	5.5
1	A	329	MET	5.0
1	B	473	TYR	4.9
1	A	245	SER	4.8
1	A	262	ILE	4.6
1	A	325	ILE	4.4
1	B	320	TYR	4.0
1	A	330	LEU	4.0
1	B	209	ALA	3.9
1	A	261	LYS	3.9
1	A	465	LEU	3.9
1	B	460	GLU	3.7
1	B	247	PHE	3.7
1	B	475	ASP	3.6
1	B	252	MET	3.6
1	A	252	MET	3.5
1	B	244	LYS	3.5
1	A	473	TYR	3.5
1	A	253	ASN	3.3
1	B	258	GLY	3.2
1	B	259	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	475	ASP	3.1
1	A	363	PHE	3.1
1	A	452	LEU	3.1
1	B	458	LYS	3.1
1	A	447	THR	3.1
1	A	287	PHE	3.1
1	B	466	HIS	3.1
1	A	359	PRO	3.0
1	B	275	LYS	2.9
1	B	474	LYS	2.9
1	A	288	ARG	2.9
1	B	459	THR	2.9
1	B	452	LEU	2.8
1	A	467	PRO	2.8
1	A	328	THR	2.8
1	A	358	LYS	2.7
1	A	453	LEU	2.7
1	A	279	ILE	2.7
1	B	472	ILE	2.7
1	B	362	ASP	2.7
1	B	468	LEU	2.7
1	B	277	VAL	2.7
1	A	362	ASP	2.7
1	B	260	ASP	2.7
1	B	352	PHE	2.6
1	B	469	LEU	2.6
1	B	245	SER	2.6
1	A	468	LEU	2.6
1	B	238	THR	2.6
1	A	469	LEU	2.6
1	B	208	SER	2.6
1	B	363	PHE	2.5
1	A	461	THR	2.5
1	A	422	LYS	2.5
1	A	394	SER	2.4
1	A	474	LYS	2.4
1	B	210	ASP	2.4
1	A	472	ILE	2.4
1	A	241	THR	2.4
1	B	257	MET	2.4
1	A	256	MET	2.3
1	A	360	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	282	PHE	2.3
1	A	277	VAL	2.3
1	A	290	VAL	2.2
1	A	446	VAL	2.2
1	B	467	PRO	2.2
1	A	464	SER	2.2
1	A	280	ARG	2.2
1	A	260	ASP	2.2
1	A	456	ILE	2.2
1	A	450	VAL	2.2
1	A	455	VAL	2.2
1	A	444	GLN	2.2
1	A	257	MET	2.1
1	A	209	ALA	2.1
1	B	236	ILE	2.1
1	B	255	LEU	2.1
1	A	449	HIS	2.1
1	A	361	GLY	2.1
1	B	253	ASN	2.0
1	A	451	GLN	2.0
1	A	466	HIS	2.0
1	A	255	LEU	2.0
1	A	454	GLN	2.0
1	A	258	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

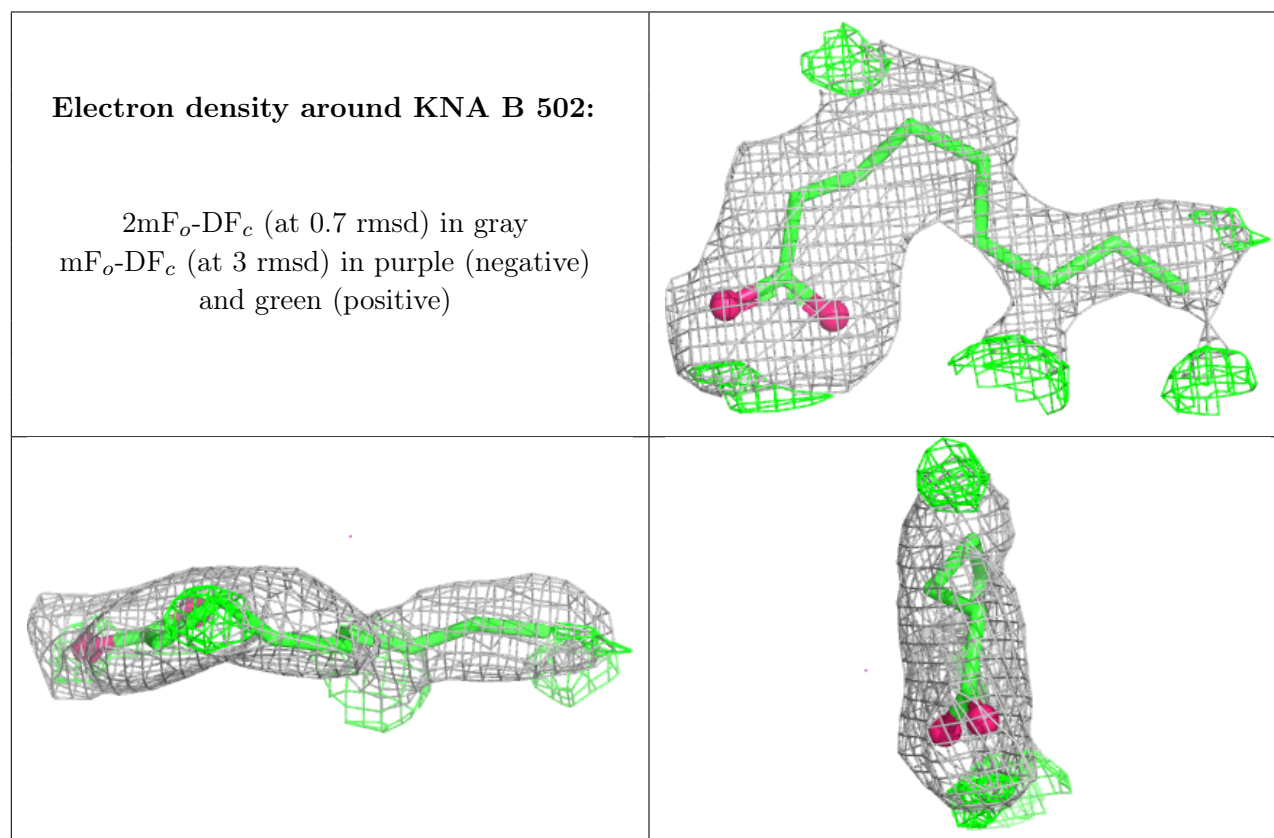
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

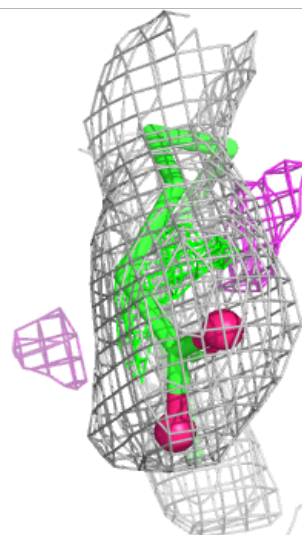
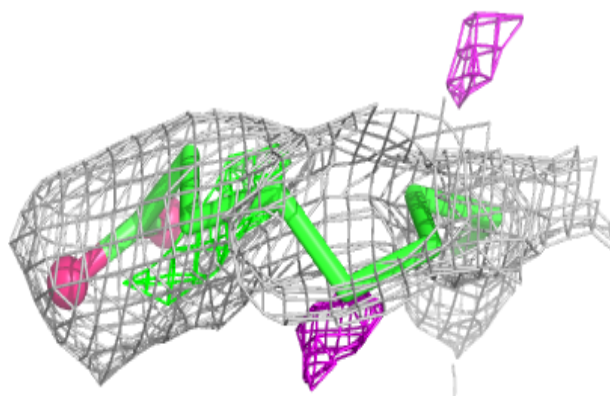
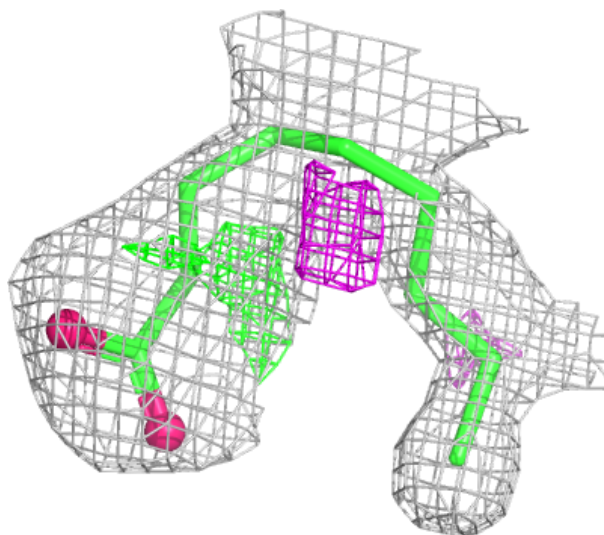
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	KNA	B	502	11/11	0.71	0.19	30,33,39,43	0
3	KNA	A	502	11/11	0.75	0.22	20,40,45,46	0
2	XZK	A	501	21/22	0.86	0.13	23,34,56,66	0
2	XZK	B	501	21/22	0.88	0.12	20,31,42,53	0

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.



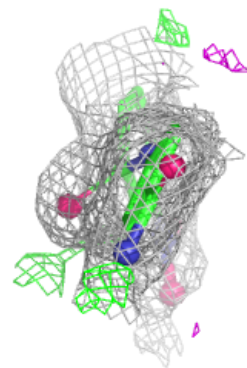
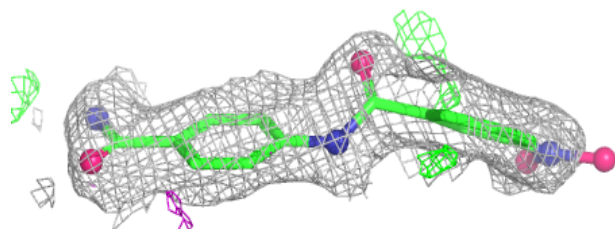
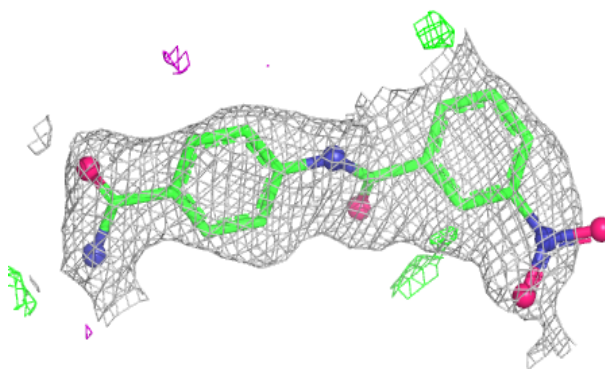
Electron density around KNA A 502:

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

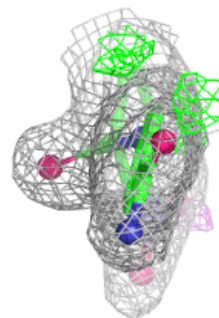
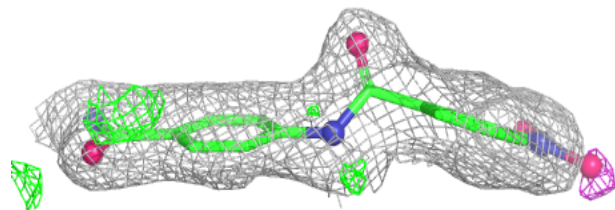
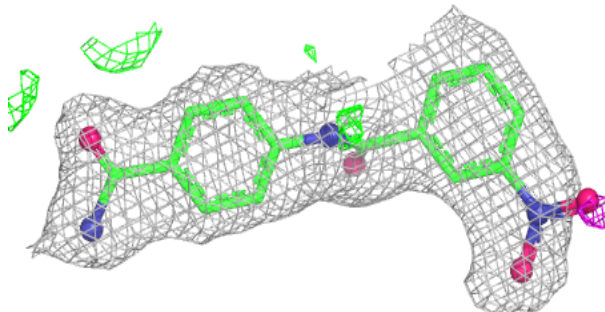


Electron density around XZK A 501:

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

**Electron density around XZK B 501:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.