Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2020 – 03:13 PM GMT

PDB ID	:	6SMO
Title	:	AntDE:AntF (apo): type II PKS acyl-carrier protein in complex with its ketosynthase bound to the hexaketide
Authors	:	Braeuer, A.; Zhou, Q.; Grammbitter, G.L.C.; Schmalhofer, M.; Ruehl, M.; Kaila, V.R.I.; Bode, H.; Groll, M.
Deposited on	:	2019-08-22
Resolution	:	2.70 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

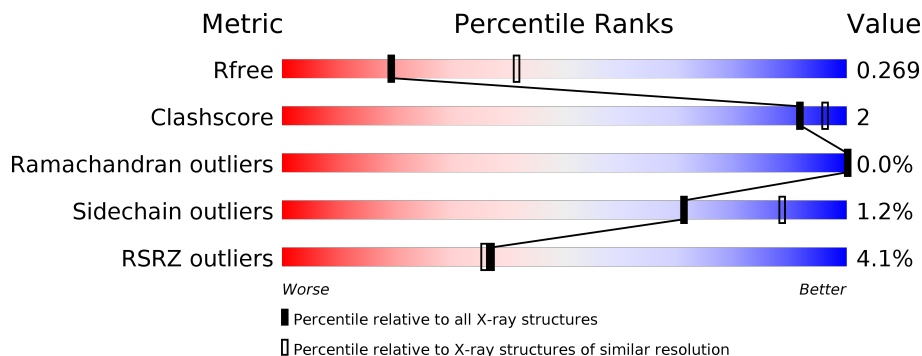|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.14.6

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.
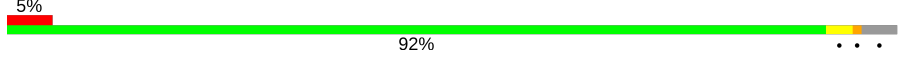
Metric	Percentile Ranks	Value
Rfree		0.269
Clashscore		2
Ramachandran outliers		0.0%
Sidechain outliers		1.2%
RSRZ outliers		4.1%

Worse | | *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on 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 <=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	97	14% / 69% · 27%
2	B	443	2% / 88% · · 8%
2	D	443	2% / 90% · 7%
2	F	443	2% / 88% 5% 8%
3	C	371	5% / 87% 6% 6%
3	E	371	5% / 92% · 5%

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Mol	Chain	Length	Quality of chain
3	G	371	

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 18074 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 Acyl carrier protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	71	Total	C	N	O	S	0	0	0
			562	357	87	116	2			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	GLY	-	expression tag	UNP A0A2S8QL96
A	-13	SER	-	expression tag	UNP A0A2S8QL96
A	-12	SER	-	expression tag	UNP A0A2S8QL96
A	-11	HIS	-	expression tag	UNP A0A2S8QL96
A	-10	HIS	-	expression tag	UNP A0A2S8QL96
A	-9	HIS	-	expression tag	UNP A0A2S8QL96
A	-8	HIS	-	expression tag	UNP A0A2S8QL96
A	-7	HIS	-	expression tag	UNP A0A2S8QL96
A	-6	HIS	-	expression tag	UNP A0A2S8QL96
A	-5	SER	-	expression tag	UNP A0A2S8QL96
A	-4	GLY	-	expression tag	UNP A0A2S8QL96
A	-3	ASP	-	expression tag	UNP A0A2S8QL96
A	-2	PRO	-	expression tag	UNP A0A2S8QL96
A	-1	ALA	-	expression tag	UNP A0A2S8QL96
A	0	SER	-	expression tag	UNP A0A2S8QL96

- Molecule 2 is a protein called PKS_KS domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	408	Total	C	N	O	S	0	0	0
			3053	1913	525	594	21			
2	D	410	Total	C	N	O	S	0	0	0
			3069	1921	528	599	21			
2	F	408	Total	C	N	O	S	0	0	0
			3050	1908	525	596	21			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	GLY	-	expression tag	UNP Q7MZT3
B	-13	SER	-	expression tag	UNP Q7MZT3
B	-12	SER	-	expression tag	UNP Q7MZT3
B	-11	HIS	-	expression tag	UNP Q7MZT3
B	-10	HIS	-	expression tag	UNP Q7MZT3
B	-9	HIS	-	expression tag	UNP Q7MZT3
B	-8	HIS	-	expression tag	UNP Q7MZT3
B	-7	HIS	-	expression tag	UNP Q7MZT3
B	-6	HIS	-	expression tag	UNP Q7MZT3
B	-5	SER	-	expression tag	UNP Q7MZT3
B	-4	GLY	-	expression tag	UNP Q7MZT3
B	-3	ASP	-	expression tag	UNP Q7MZT3
B	-2	PRO	-	expression tag	UNP Q7MZT3
B	-1	ALA	-	expression tag	UNP Q7MZT3
B	0	SER	-	expression tag	UNP Q7MZT3
D	-14	GLY	-	expression tag	UNP Q7MZT3
D	-13	SER	-	expression tag	UNP Q7MZT3
D	-12	SER	-	expression tag	UNP Q7MZT3
D	-11	HIS	-	expression tag	UNP Q7MZT3
D	-10	HIS	-	expression tag	UNP Q7MZT3
D	-9	HIS	-	expression tag	UNP Q7MZT3
D	-8	HIS	-	expression tag	UNP Q7MZT3
D	-7	HIS	-	expression tag	UNP Q7MZT3
D	-6	HIS	-	expression tag	UNP Q7MZT3
D	-5	SER	-	expression tag	UNP Q7MZT3
D	-4	GLY	-	expression tag	UNP Q7MZT3
D	-3	ASP	-	expression tag	UNP Q7MZT3
D	-2	PRO	-	expression tag	UNP Q7MZT3
D	-1	ALA	-	expression tag	UNP Q7MZT3
D	0	SER	-	expression tag	UNP Q7MZT3
F	-14	GLY	-	expression tag	UNP Q7MZT3
F	-13	SER	-	expression tag	UNP Q7MZT3
F	-12	SER	-	expression tag	UNP Q7MZT3
F	-11	HIS	-	expression tag	UNP Q7MZT3
F	-10	HIS	-	expression tag	UNP Q7MZT3
F	-9	HIS	-	expression tag	UNP Q7MZT3
F	-8	HIS	-	expression tag	UNP Q7MZT3
F	-7	HIS	-	expression tag	UNP Q7MZT3
F	-6	HIS	-	expression tag	UNP Q7MZT3
F	-5	SER	-	expression tag	UNP Q7MZT3
F	-4	GLY	-	expression tag	UNP Q7MZT3
F	-3	ASP	-	expression tag	UNP Q7MZT3

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-2	PRO	-	expression tag	UNP Q7MZT3
F	-1	ALA	-	expression tag	UNP Q7MZT3
F	0	SER	-	expression tag	UNP Q7MZT3

- Molecule 3 is a protein called Ketoacyl_synth_N domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	347	Total 2682	C 1697	N 456	O 520	S 9	0	0	0
3	E	352	Total 2713	C 1712	N 463	O 528	S 10	0	0	0
3	G	355	Total 2740	C 1730	N 465	O 535	S 10	0	0	0

- Molecule 4 is 3,5,7,9,11-pentakis(oxidanylidene)dodecanal (three-letter code: LKZ) (formula: $C_{12}H_{14}O_6$).
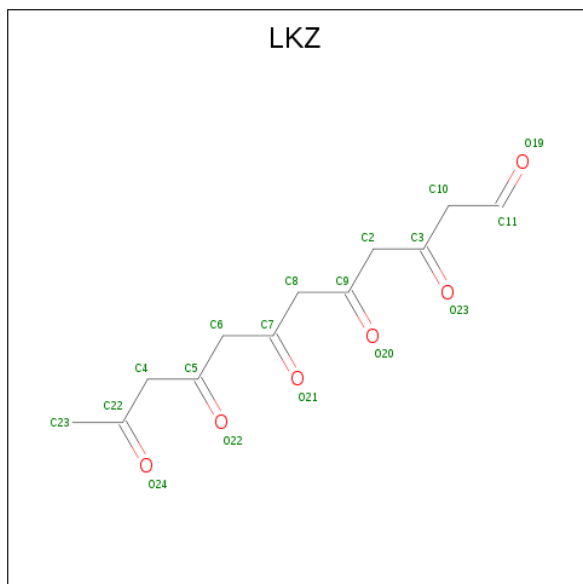
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total 18	C 12	O 6	0	0
4	D	1	Total 18	C 12	O 6	0	0
4	F	1	Total 18	C 12	O 6	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	2	Total Na 2 2		0	0

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O_4S).
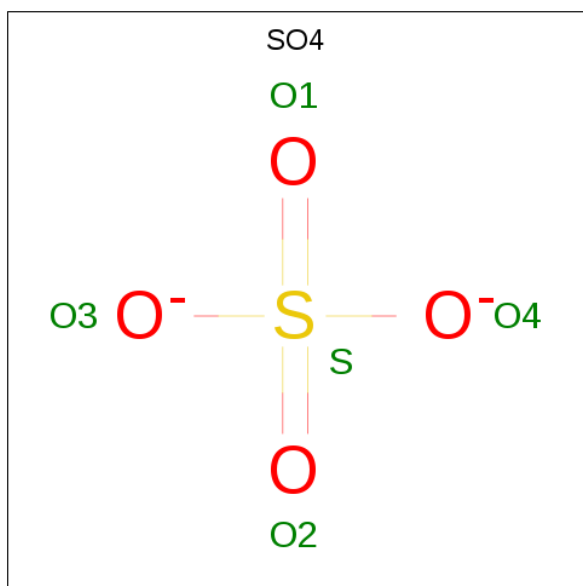
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	G	1	Total 5	O 4	S 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	3	Total 3	O 3	0	0
7	B	23	Total 23	O 23	0	0
7	C	15	Total 15	O 15	0	0
7	D	32	Total 32	O 32	0	0
7	E	20	Total 20	O 20	0	0
7	F	30	Total 30	O 30	0	0
7	G	21	Total 21	O 21	0	0

# 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.
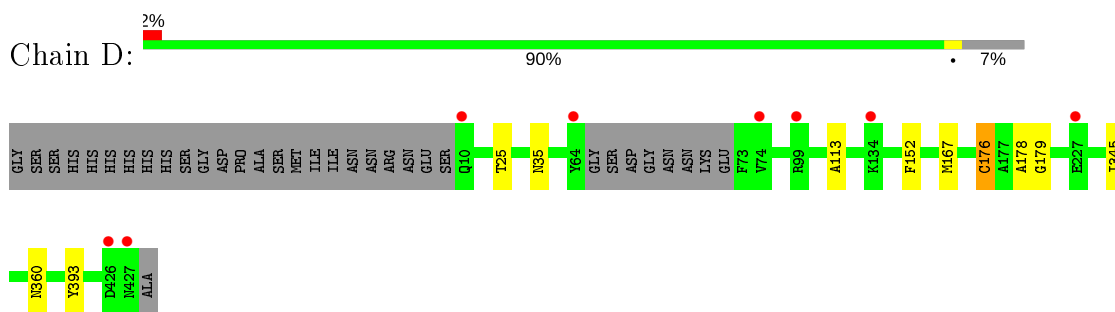
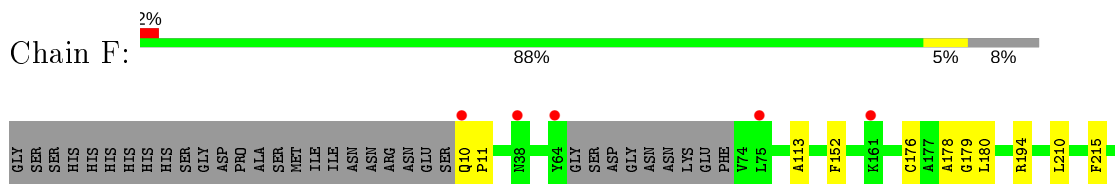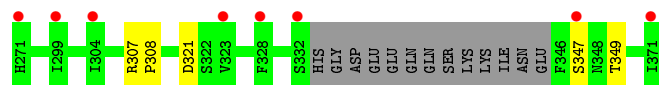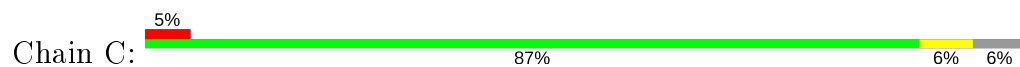- Molecule 1: Acyl carrier protein
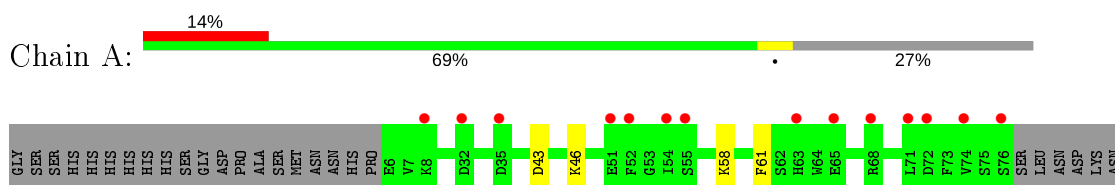
- Molecule 2: PKS_KS domain-containing protein
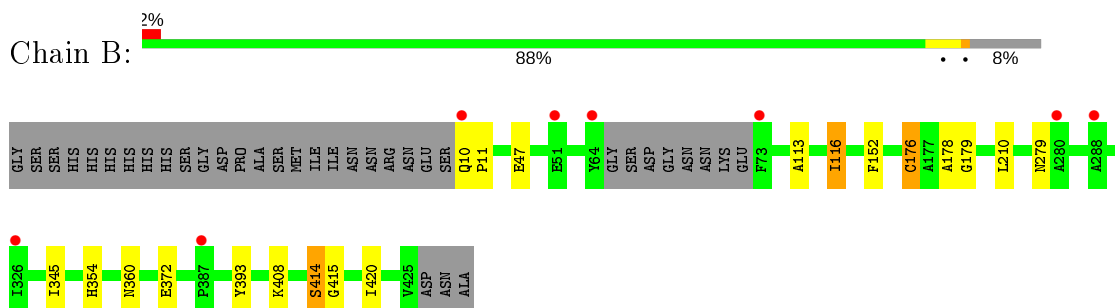
- Molecule 2: PKS_KS domain-containing protein

- Molecule 2: PKS_KS domain-containing protein
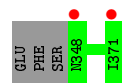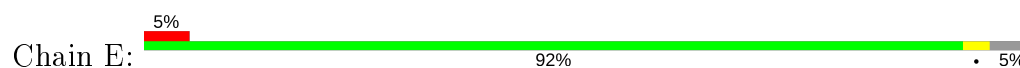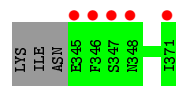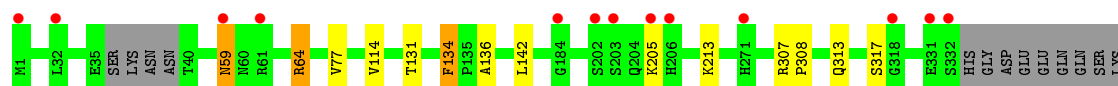
• Molecule 3: Ketoacyl_synth_N domain-containing protein

Chain C:

• Molecule 3: Ketoacyl_synth_N domain-containing protein

Chain E:

• Molecule 3: Ketoacyl_synth_N domain-containing protein

Chain G:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α, β, γ	107.65Å 107.65Å 201.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 29.86 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.2 (30.00-2.70) 98.2 (29.86-2.70)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.226 , 0.265 0.226 , 0.269	Depositor DCC
R_{free} test set	3077 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å²)	41.2	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.32 , 22.0	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	18074	wwPDB-VP
Average B, all atoms (Å²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.53% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, LKZ, SO4

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	0/570	0.68	0/769
2	B	0.68	0/3109	0.73	1/4206 (0.0%)
2	D	0.67	0/3125	0.73	1/4228 (0.0%)
2	F	0.68	0/3105	0.73	1/4201 (0.0%)
3	C	0.69	0/2729	0.72	0/3682
3	E	0.68	0/2759	0.71	0/3723
3	G	0.68	0/2787	0.71	1/3760 (0.0%)
All	All	0.68	0/18184	0.72	4/24569 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	176	CYS	CA-CB-SG	-5.76	103.62	114.00
2	B	176	CYS	CA-CB-SG	-5.41	104.26	114.00
3	G	59	ASN	CB-CA-C	5.38	121.16	110.40
2	F	176	CYS	CA-CB-SG	-5.20	104.63	114.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	562	0	544	4	0
2	B	3053	0	2992	10	0
2	D	3069	0	3001	5	0
2	F	3050	0	2987	11	0
3	C	2682	0	2645	18	0
3	E	2713	0	2679	6	0
3	G	2740	0	2699	8	0
4	B	18	0	0	2	0
4	D	18	0	0	1	0
4	F	18	0	0	0	0
5	D	2	0	0	0	0
6	G	5	0	0	0	0
7	A	3	0	0	0	0
7	B	23	0	0	0	0
7	C	15	0	0	0	0
7	D	32	0	0	0	0
7	E	20	0	0	0	0
7	F	30	0	0	0	0
7	G	21	0	0	0	0
All	All	18074	0	17547	55	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 2.

All (55) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:414:SER:HB3	3:C:140:GLY:HA3	1.36	1.07
2:B:414:SER:HB3	3:C:140:GLY:CA	2.03	0.87
3:C:98:ILE:HG21	3:C:142:LEU:HD12	1.76	0.66
3:C:14:HIS:CD2	3:C:15:PRO:HD2	2.30	0.66
2:F:113:ALA:HB1	2:F:178:ALA:HB1	1.80	0.64
3:C:14:HIS:CG	3:C:15:PRO:HD2	2.33	0.63
3:C:139:GLN:NE2	3:C:152:SER:O	2.32	0.63
3:E:54:GLY:H	3:E:63:ARG:HH12	1.49	0.60
1:A:46:LYS:NZ	3:C:122:ILE:O	2.33	0.59
1:A:46:LYS:NZ	3:C:125:MET:O	2.37	0.57
2:D:113:ALA:HB1	2:D:178:ALA:HB1	1.86	0.57
2:F:215:PHE:HB3	2:F:221:LEU:HD13	1.85	0.57
2:B:113:ALA:HB1	2:B:178:ALA:HB1	1.88	0.56
2:B:176:CYS:SG	4:B:501:LKZ:C2	2.94	0.55
3:C:139:GLN:NE2	3:C:152:SER:C	2.60	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:307:ARG:N	3:E:308:PRO:HD2	2.23	0.53
2:B:116:ILE:HG22	4:B:501:LKZ:C6	2.39	0.53
3:C:139:GLN:HE22	3:C:152:SER:C	2.12	0.53
3:E:134:PHE:CE2	3:E:136:ALA:HB3	2.46	0.51
3:G:134:PHE:CE2	3:G:136:ALA:HB3	2.46	0.51
2:B:345:ILE:O	2:B:393:TYR:HA	2.11	0.51
2:D:179:GLY:HA3	2:D:360:ASN:HB2	1.92	0.50
2:F:215:PHE:CB	2:F:221:LEU:HD13	2.43	0.48
2:F:356:LEU:N	2:F:356:LEU:HD23	2.29	0.48
2:F:210:LEU:HD13	3:G:114:VAL:HG22	1.96	0.47
1:A:58:LYS:HA	1:A:61:PHE:CE2	2.50	0.46
3:C:16:ASP:OD1	3:C:47:ARG:NH1	2.45	0.46
2:B:408:LYS:HB3	2:B:420:ILE:HG13	1.98	0.46
3:G:313:GLN:H	3:G:317:SER:HG	1.64	0.46
2:B:279:ASN:HD22	2:B:415:GLY:HA2	1.81	0.46
2:D:25:THR:O	2:D:35:ASN:ND2	2.48	0.45
3:C:98:ILE:HG21	3:C:142:LEU:CD1	2.44	0.45
2:D:176:CYS:SG	4:D:501:LKZ:C2	2.95	0.45
2:F:179:GLY:HA3	2:F:360:ASN:HB2	1.99	0.45
3:C:134:PHE:HA	3:C:135:PRO:HD3	1.87	0.45
3:E:134:PHE:HA	3:E:135:PRO:HD3	1.83	0.45
3:G:64:ARG:HD3	3:G:64:ARG:HA	1.53	0.44
3:C:307:ARG:N	3:C:308:PRO:CD	2.81	0.44
3:C:141:GLN:OE1	3:C:141:GLN:HA	2.18	0.44
2:D:345:ILE:O	2:D:393:TYR:HA	2.18	0.44
2:F:210:LEU:CD1	3:G:114:VAL:HG22	2.47	0.44
2:F:180:LEU:HD12	2:F:410:SER:HB2	1.99	0.44
3:G:307:ARG:N	3:G:308:PRO:CD	2.81	0.43
2:B:179:GLY:HA3	2:B:360:ASN:HB2	2.01	0.42
3:C:73:GLY:HA3	3:C:138:LEU:HD21	2.00	0.42
3:C:110:ILE:HD13	3:C:133:TRP:HB2	2.02	0.42
2:B:10:GLN:N	2:B:11:PRO:HD2	2.35	0.42
3:E:54:GLY:N	3:E:63:ARG:HH12	2.15	0.41
2:F:10:GLN:N	2:F:11:PRO:HD2	2.36	0.41
2:F:313:TYR:HE2	2:F:367:LEU:HD13	1.85	0.41
3:E:270:ASN:O	3:E:270:ASN:ND2	2.53	0.41
2:F:311:VAL:HG21	2:F:407:LEU:HG	2.03	0.41
3:G:205:LYS:HD3	3:G:213:LYS:HD2	2.03	0.40
1:A:43:ASP:OD2	3:C:64:ARG:NH1	2.55	0.40
3:G:77:VAL:HG21	3:G:142:LEU:HD22	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	69/97 (71%)	65 (94%)	4 (6%)	0	100	100
2	B	404/443 (91%)	384 (95%)	20 (5%)	0	100	100
2	D	406/443 (92%)	384 (95%)	22 (5%)	0	100	100
2	F	404/443 (91%)	388 (96%)	16 (4%)	0	100	100
3	C	341/371 (92%)	328 (96%)	12 (4%)	1 (0%)	41	66
3	E	346/371 (93%)	331 (96%)	15 (4%)	0	100	100
3	G	349/371 (94%)	334 (96%)	15 (4%)	0	100	100
All	All	2319/2539 (91%)	2214 (96%)	104 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	347	SER

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	63/86 (73%)	63 (100%)	0	100	100
2	B	319/348 (92%)	312 (98%)	7 (2%)	52	79
2	D	321/348 (92%)	319 (99%)	2 (1%)	86	95
2	F	319/348 (92%)	316 (99%)	3 (1%)	78	92
3	C	291/312 (93%)	288 (99%)	3 (1%)	76	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	E	294/312 (94%)	290 (99%)	4 (1%)	67	86
3	G	297/312 (95%)	293 (99%)	4 (1%)	69	87
All	All	1904/2066 (92%)	1881 (99%)	23 (1%)	71	88

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	47	GLU
2	B	116	ILE
2	B	152	PHE
2	B	210	LEU
2	B	354	HIS
2	B	372	GLU
2	B	414	SER
3	C	134	PHE
3	C	321	ASP
3	C	349	THR
2	D	152	PHE
2	D	167	MET
3	E	58	ILE
3	E	59	ASN
3	E	69	PHE
3	E	321	ASP
2	F	152	PHE
2	F	194	ARG
2	F	372	GLU
3	G	59	ASN
3	G	64	ARG
3	G	131	THR
3	G	134	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	E	141	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	LKZ	D	501	-	16,17,17	0.52	0	17,21,21	1.02	1 (5%)
4	LKZ	B	501	-	16,17,17	0.54	0	17,21,21	0.76	0
6	SO4	G	401	-	4,4,4	0.39	0	6,6,6	0.05	0
4	LKZ	F	501	-	16,17,17	0.38	0	17,21,21	1.10	2 (11%)

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	LKZ	D	501	-	-	10/18/19/19	-
4	LKZ	B	501	-	-	8/18/19/19	-
4	LKZ	F	501	-	-	10/18/19/19	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	F	501	LKZ	O20-C9-C2	2.17	124.50	120.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	501	LKZ	C11-C10-C3	-2.10	101.79	111.68
4	F	501	LKZ	C11-C10-C3	-2.02	102.21	111.68

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	501	LKZ	O24-C22-C4-C5
4	F	501	LKZ	C23-C22-C4-C5
4	B	501	LKZ	C22-C4-C5-C6
4	B	501	LKZ	O24-C22-C4-C5
4	B	501	LKZ	C23-C22-C4-C5
4	F	501	LKZ	C5-C6-C7-O21
4	F	501	LKZ	C7-C8-C9-O20
4	B	501	LKZ	C22-C4-C5-O22
4	F	501	LKZ	C5-C6-C7-C8
4	F	501	LKZ	C9-C2-C3-C10
4	D	501	LKZ	C4-C5-C6-C7
4	D	501	LKZ	C5-C6-C7-C8
4	D	501	LKZ	C7-C8-C9-C2
4	D	501	LKZ	C23-C22-C4-C5
4	F	501	LKZ	C11-C10-C3-C2
4	D	501	LKZ	C11-C10-C3-C2
4	B	501	LKZ	C11-C10-C3-C2
4	D	501	LKZ	C22-C4-C5-O22
4	D	501	LKZ	C7-C8-C9-O20
4	B	501	LKZ	O22-C5-C6-C7
4	F	501	LKZ	C6-C7-C8-C9
4	F	501	LKZ	C7-C8-C9-C2
4	F	501	LKZ	C3-C2-C9-C8
4	D	501	LKZ	C22-C4-C5-C6
4	D	501	LKZ	C3-C2-C9-C8
4	B	501	LKZ	C4-C5-C6-C7
4	B	501	LKZ	C7-C8-C9-C2
4	D	501	LKZ	O24-C22-C4-C5

There are no ring outliers.

2 monomers are involved in 3 short contacts:

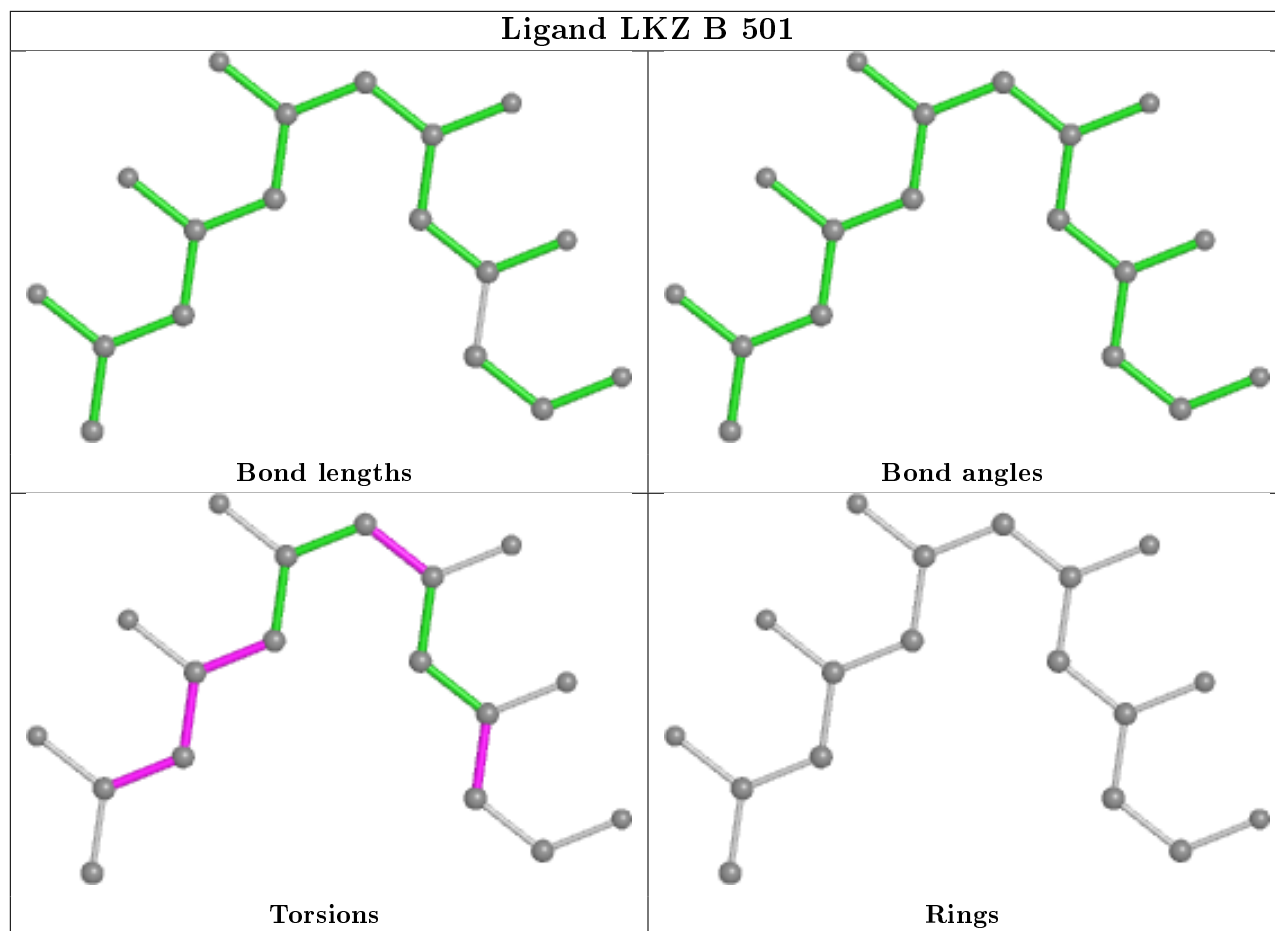| Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	501	LKZ	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	501	LKZ	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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 LKZ D 501

Bond lengths

Bond angles

Torsions

Rings

Ligand LKZ B 501

Bond lengths

Bond angles

Torsions

Rings

Ligand LKZ F 501

Bond lengths

Bond angles

Torsions

Rings
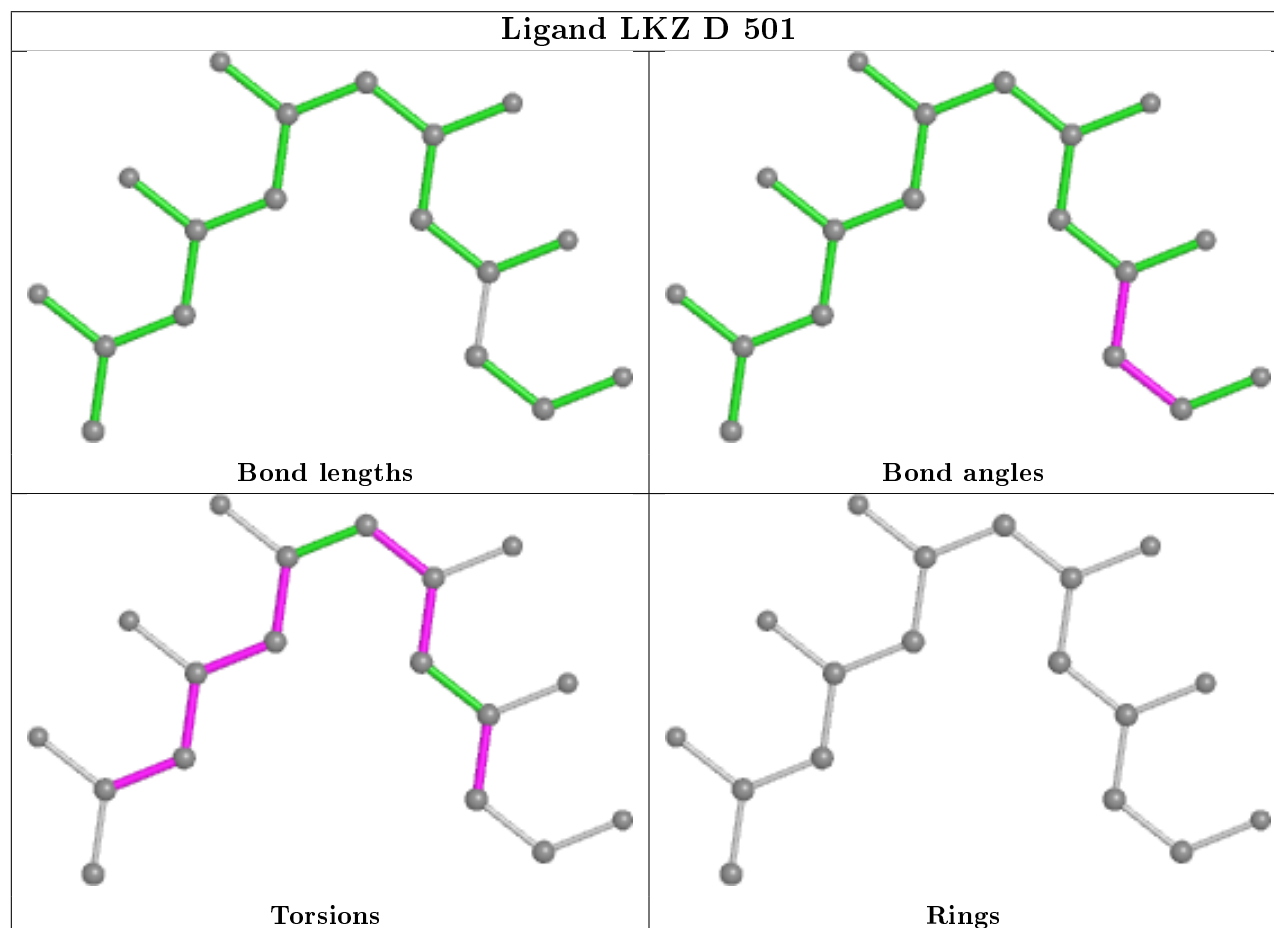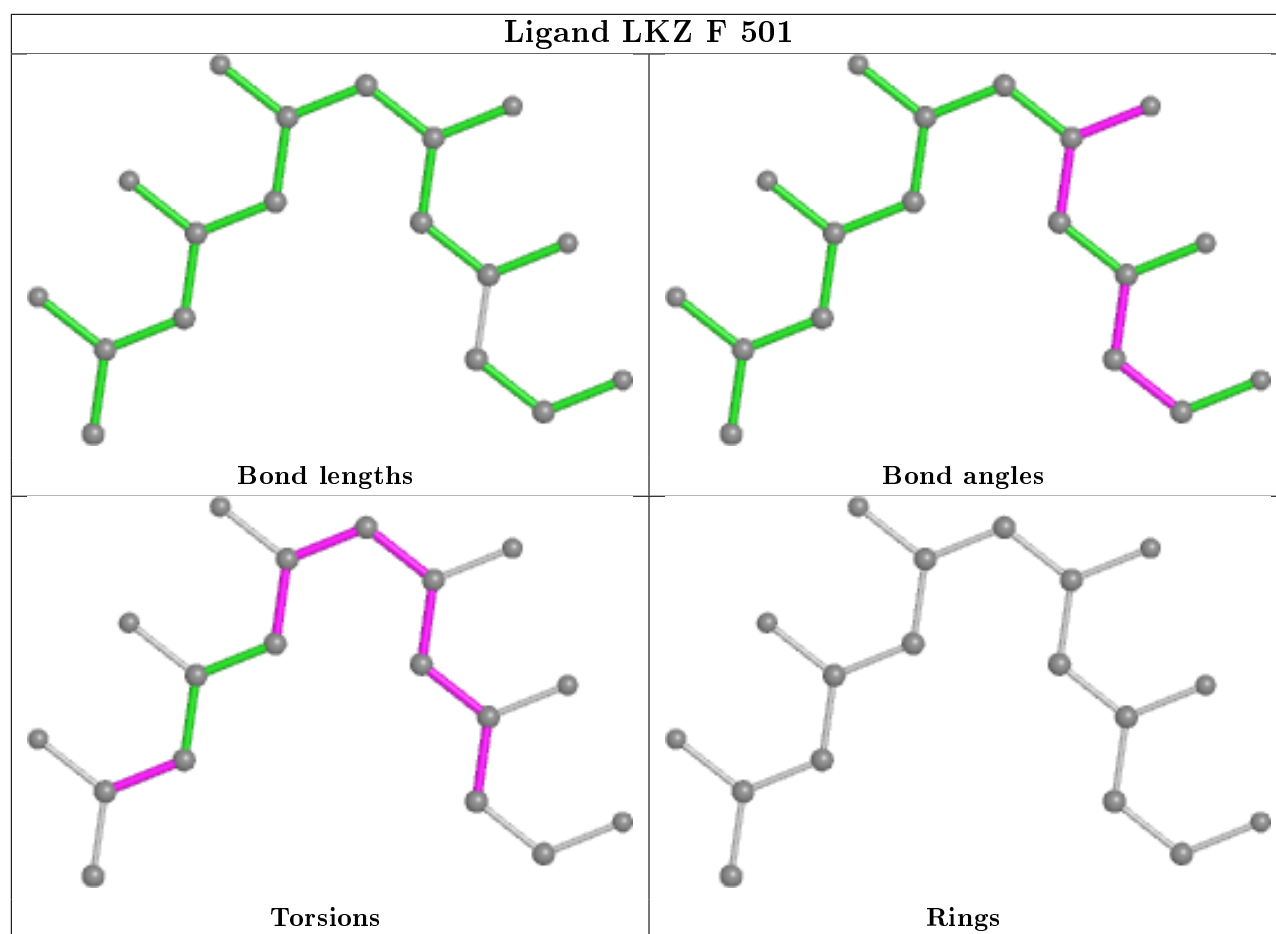
5.7 Other polymers (i)

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(Å²)	Q<0.9
1	A	71/97 (73%)	1.12	14 (19%)	1	0	62, 94, 114, 130	0
2	B	408/443 (92%)	0.22	8 (1%)	65	67	30, 40, 65, 85	0
2	D	410/443 (92%)	0.15	8 (1%)	65	67	28, 38, 61, 82	0
2	F	408/443 (92%)	0.19	9 (2%)	62	63	31, 42, 63, 85	0
3	C	347/371 (93%)	0.40	20 (5%)	23	22	31, 47, 71, 83	0
3	E	352/371 (94%)	0.34	19 (5%)	25	24	27, 42, 68, 84	0
3	G	355/371 (95%)	0.37	18 (5%)	28	26	30, 43, 65, 80	0
All	All	2351/2539 (92%)	0.30	96 (4%)	37	36	27, 42, 72, 130	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	52	PHE	5.6
3	G	332	SER	5.1
3	E	54	GLY	4.9
3	G	1	MET	4.7
3	G	348	ASN	4.7
2	D	10	GLN	4.4
3	G	205	LYS	4.4
3	E	332	SER	4.3
3	E	1	MET	4.3
3	C	56	LYS	4.2
3	E	201	THR	4.2
2	F	262	GLN	4.2
3	C	202	SER	4.1
3	E	42	ALA	4.1
3	E	371	ILE	4.0
3	G	346	PHE	4.0
3	E	33	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
3	G	203	SER	3.9
3	E	53	ASP	3.8
3	C	201	THR	3.6
3	C	32	LEU	3.6
2	B	73	PHE	3.6
2	D	74	VAL	3.6
3	E	41	THR	3.6
2	F	64	TYR	3.5
2	D	64	TYR	3.5
3	G	345	GLU	3.5
3	G	271	HIS	3.5
2	B	51	GLU	3.5
3	G	371	ILE	3.5
1	A	51	GLU	3.4
3	E	51	ASP	3.4
3	C	45	ILE	3.3
1	A	76	SER	3.3
3	C	43	SER	3.3
1	A	32	ASP	3.3
2	D	427	ASN	3.3
1	A	72	ASP	3.3
3	C	29	GLN	3.3
3	E	50	SER	3.2
3	C	371	ILE	3.2
3	C	271	HIS	3.1
3	C	31	LEU	3.1
3	G	32	LEU	3.0
2	F	10	GLN	2.9
2	F	75	LEU	2.9
3	E	203	SER	2.9
3	C	53	ASP	2.8
2	B	326	ILE	2.8
2	D	227	GLU	2.8
3	E	204	GLN	2.7
3	E	39	ASN	2.7
3	C	2	ARG	2.6
2	F	258	GLU	2.6
3	E	32	LEU	2.6
3	G	59	ASN	2.5
3	C	332	SER	2.5
3	C	347	SER	2.5
3	G	347	SER	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	387	PRO	2.5
1	A	54	ILE	2.5
1	A	8	LYS	2.5
1	A	74	VAL	2.5
3	C	203	SER	2.5
1	A	63	HIS	2.5
3	G	202	SER	2.4
3	G	61	ARG	2.4
3	E	348	ASN	2.4
3	G	318	GLY	2.4
3	E	328	PHE	2.4
3	E	56	LYS	2.3
2	D	134	LYS	2.3
2	F	38	ASN	2.3
3	C	204	GLN	2.3
3	E	34	GLN	2.3
2	D	99	ARG	2.3
3	G	184	GLY	2.3
1	A	71	LEU	2.3
2	F	227	GLU	2.3
2	B	64	TYR	2.2
3	G	206	HIS	2.2
3	G	331	GLU	2.2
1	A	55	SER	2.2
3	C	328	PHE	2.2
2	B	288	ALA	2.2
3	C	323	VAL	2.1
2	F	261	LYS	2.1
3	C	304	ILE	2.1
1	A	68	ARG	2.1
2	D	426	ASP	2.1
1	A	35	ASP	2.1
1	A	65	GLU	2.0
2	B	280	ALA	2.0
2	F	161	LYS	2.0
2	B	10	GLN	2.0
3	C	299	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.
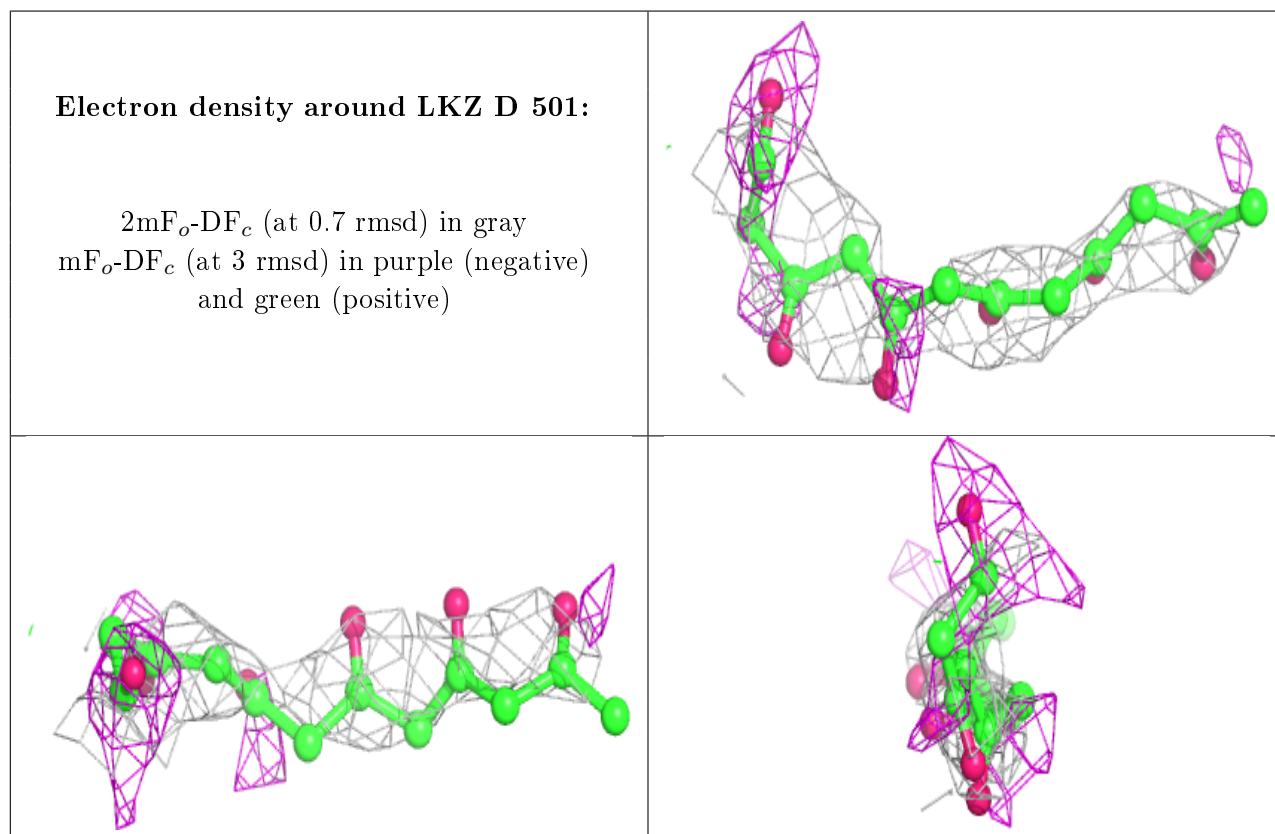
6.3 Carbohydrates (i)

There are no monosaccharides 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(Å^2)	Q<0.9
5	NA	D	503	1/1	0.82	0.42	38,38,38,38	0
4	LKZ	D	501	18/18	0.86	0.35	51,54,55,56	0
4	LKZ	F	501	18/18	0.86	0.29	48,50,56,57	0
4	LKZ	B	501	18/18	0.87	0.33	49,53,57,58	0
5	NA	D	502	1/1	0.91	0.13	22,22,22,22	0
6	SO4	G	401	5/5	0.94	0.23	69,69,72,73	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.
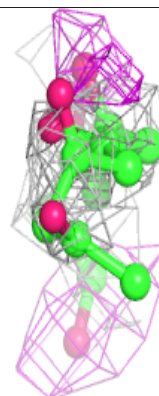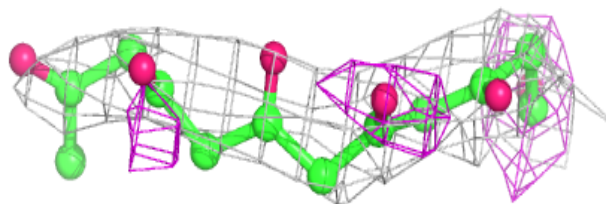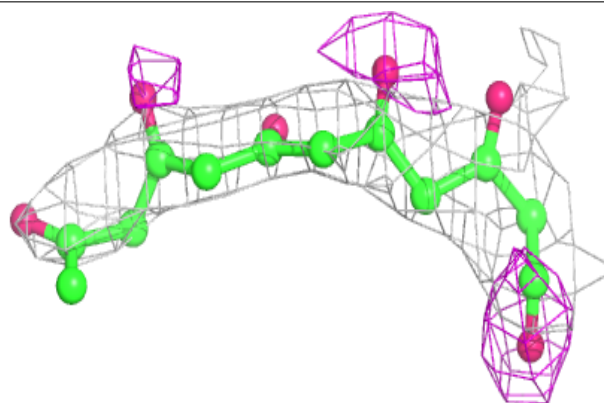
Electron density around LKZ D 501:

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

Electron density around LKZ F 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
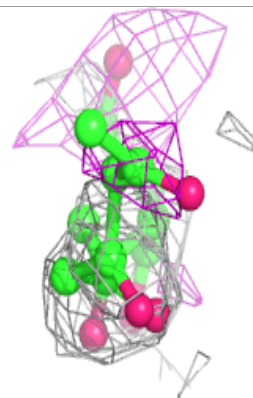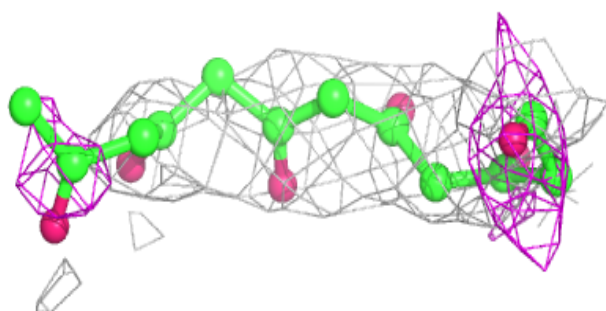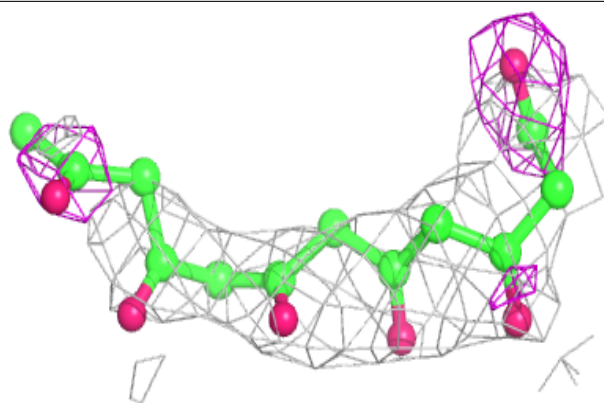
Electron density around LKZ B 501:

$2mF_o-DF_c$ (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.