wwPDB X-ray Structure Validation Summary Report (i)

Nov 5, 2023 – 01:59 PM EST

PDB ID	:	4IY7
Title	:	crystal structure of cystathionine gamma lyase (XometC) from Xanthomonas oryzae pv. oryzae in complex with E-site serine, A-site external aldimine structure with serine and A-site external aldimine structure with aminoacrylate intermediates
Authors	:	Ngo, H.P.T.; Kim, J.K.; Kang, L.W.
Deposited on	:	2013-01-28
Resolution	:	1.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

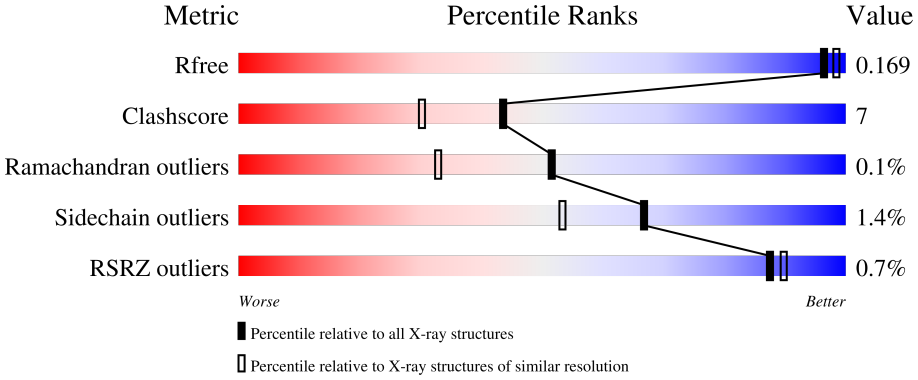
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.70 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 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 <=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	397	.% 79% 15% • •
1	B	397	82% 13% • •
1	C	397	.% 83% 13% •
1	D	397	.% 84% 12% • •

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
6	PYR	D	402	-	X	-	-
7	KOU	B	403	-	-	X	-
7	KOU	C	402	-	-	X	-
7	KOU	D	403	-	-	X	-

2 Entry composition (i)

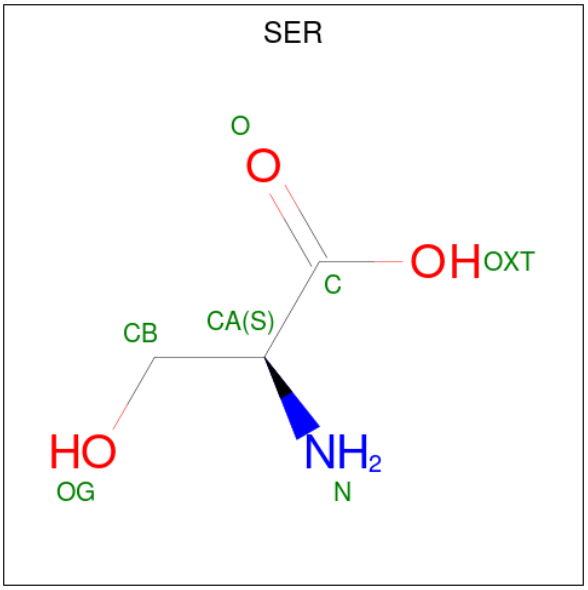
There are 8 unique types of molecules in this entry. The entry contains 13157 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cystathionine gamma-lyase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	381	2889	1830	509	535	15	0	6	0
1	B	381	2892	1832	509	536	15	0	7	0
1	C	381	2886	1828	509	534	15	0	5	0
1	D	383	2918	1849	515	539	15	0	8	0

- Molecule 2 is SERINE (three-letter code: SER) (formula: $C_3H_7NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	7	3	1	3	0	0
2	B	1	7	3	1	3	0	0

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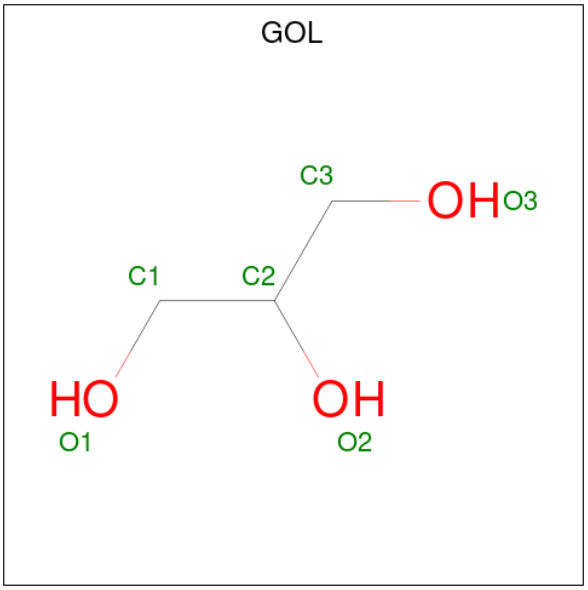
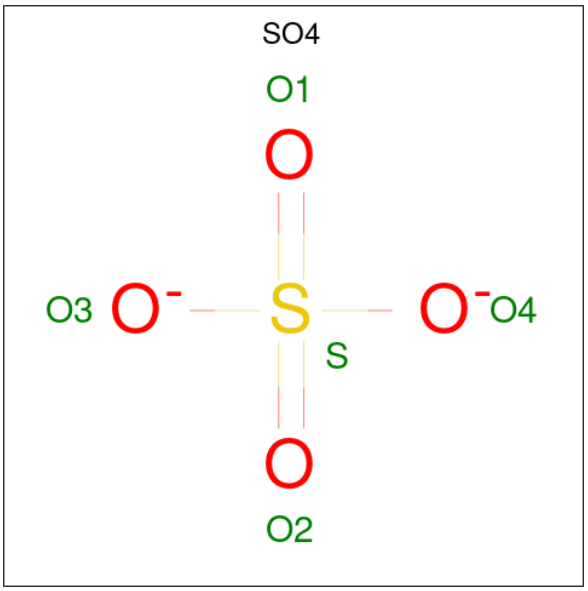
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	C	1	7	3	1	3	0	0
			Total	C	N	O		
2	D	1	7	3	1	3	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



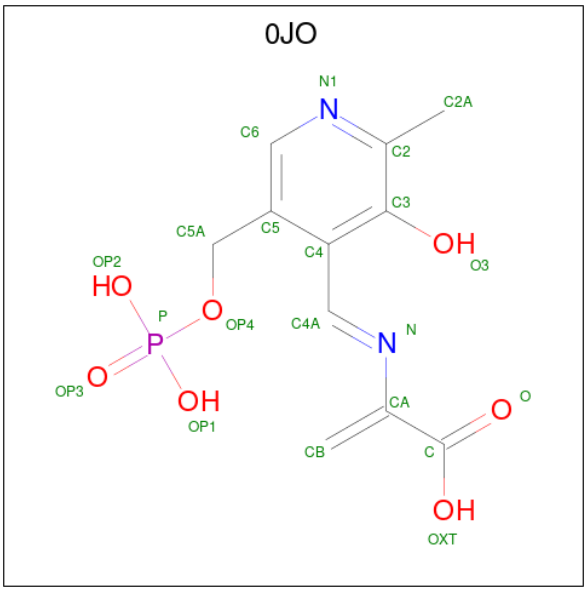
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
3	A	1	6	3	3		0	0
			Total	C	O			
3	A	1	6	3	3		0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



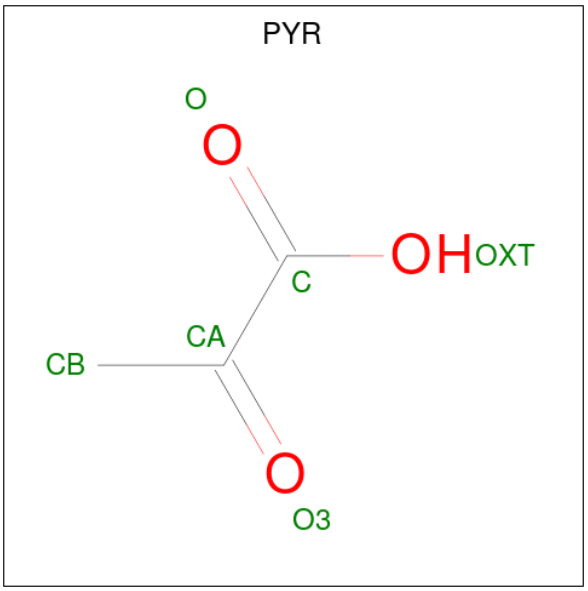
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	A	1	5	4	1	0	0

- Molecule 5 is 2-{[(E)-{3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridin-4-yl}methylidene]amino}prop-2-enoic acid (three-letter code: 0JO) (formula: $C_{11}H_{13}N_2O_7P$).



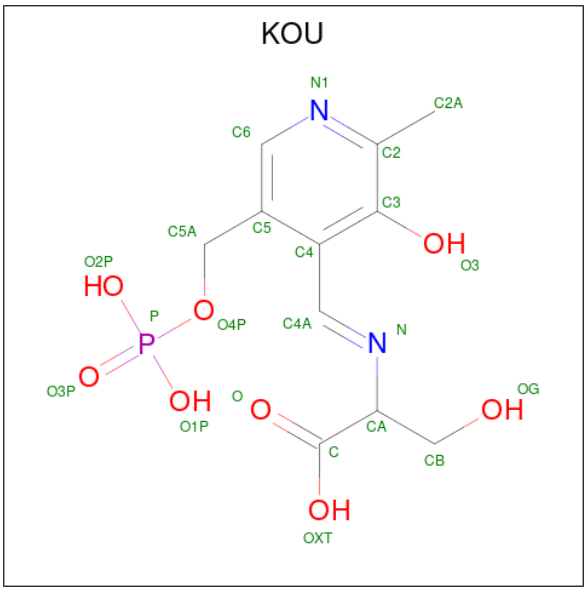
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
5	A	1	21	11	2	7	1	0	0

- Molecule 6 is PYRUVIC ACID (three-letter code: PYR) (formula: $C_3H_4O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
6	B	1	6	3	3	0	0
			Total	C	O		
6	D	1	6	3	3	0	0

- Molecule 7 is (E)-N-({3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridin-4-yl}methylidene)-L-serine (three-letter code: KOU) (formula: $C_{11}H_{15}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
7	B	1	22	11	2	8	1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
7	C	1	22	11	2	8	1	0	0
			Total	C	N	O	P		
7	D	1	22	11	2	8	1	0	0

- Molecule 8 is water.

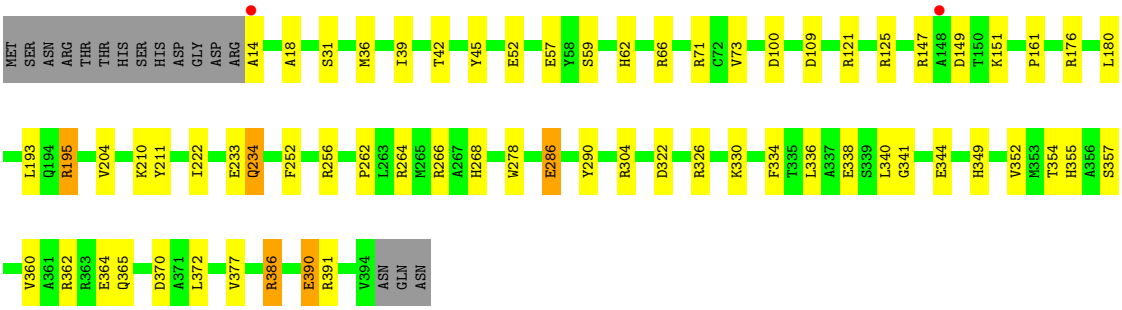
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	346	346	346	0	0
			Total	O		
8	B	389	389	389	0	0
			Total	O		
8	C	332	332	332	0	0
			Total	O		
8	D	361	361	361	0	0

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.

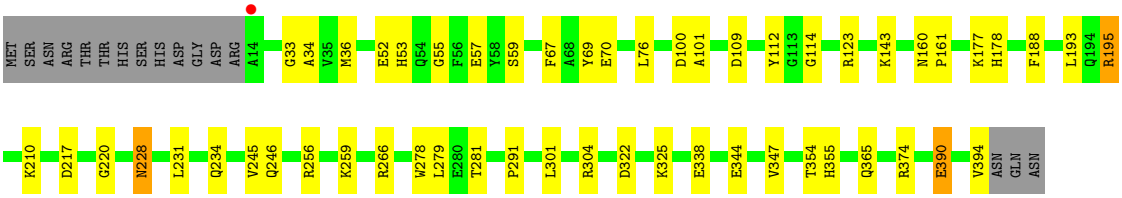
• Molecule 1: Cystathionine gamma-lyase-like protein

Chain A:



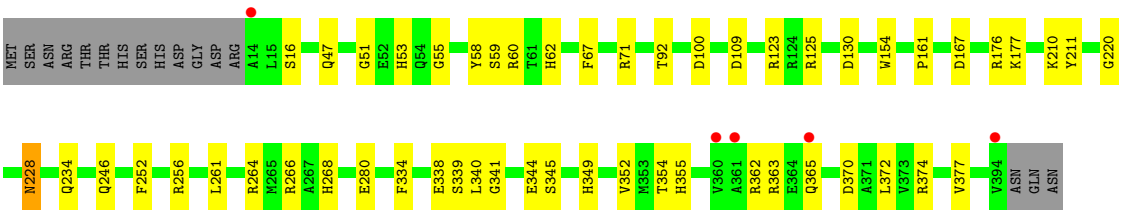
• Molecule 1: Cystathionine gamma-lyase-like protein

Chain B:



• Molecule 1: Cystathionine gamma-lyase-like protein

Chain C:



• Molecule 1: Cystathionine gamma-lyase-like protein

Chain D:

MET
SER
ASN
ARG
THR
THR
HIS
SER
HIS
ASP
GLY
ASP
ARG
A14
L15
S16
P28
S31
M36
I39
Q47
G51
E52
H53
Q54
G55
R60
R71
R81
D100
D109
F116
R117
R121
V122
R123
R147
A148
D149
P161
D185
S190
R195
P196
V204
K210
N228
A229
E230
E233
Q234
K259
N271
P283
R304
L317
K318
R326
E338
S339
L340
E344
H349
V352
M353
T354
H355
A356
S357
E364
D370
A371
L372
R386
Q396
ASN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	76.22Å 86.07Å 226.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.23 – 1.70 30.23 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.23-1.70) 100.0 (30.23-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.71 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.134 , 0.169 0.134 , 0.169	Depositor DCC
R_{free} test set	8207 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	13.6	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 45.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	13157	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, SO4, PYR, KOU, 0JO

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	1.41	16/2965 (0.5%)	1.26	26/4022 (0.6%)
1	B	1.41	13/2971 (0.4%)	1.21	19/4030 (0.5%)
1	C	1.39	5/2959 (0.2%)	1.17	19/4014 (0.5%)
1	D	1.39	9/3000 (0.3%)	1.20	16/4067 (0.4%)
All	All	1.40	43/11895 (0.4%)	1.21	80/16133 (0.5%)

The worst 5 of 43 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	338	GLU	CD-OE2	-11.75	1.12	1.25
1	A	233	GLU	CG-CD	9.78	1.66	1.51
1	D	338	GLU	CD-OE1	-8.94	1.15	1.25
1	A	57	GLU	CD-OE1	8.39	1.34	1.25
1	A	344	GLU	CG-CD	7.68	1.63	1.51

The worst 5 of 80 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	304	ARG	NE-CZ-NH2	-10.25	115.18	120.30
1	A	66	ARG	NE-CZ-NH2	-10.17	115.22	120.30
1	A	109	ASP	CB-CG-OD1	9.90	127.21	118.30
1	A	121	ARG	NE-CZ-NH1	9.58	125.09	120.30
1	C	109	ASP	CB-CG-OD1	9.54	126.89	118.30

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	2889	0	2919	40	0
1	B	2892	0	2924	48	0
1	C	2886	0	2914	33	0
1	D	2918	0	2959	38	0
2	A	7	0	4	0	0
2	B	7	0	4	0	0
2	C	7	0	4	0	0
2	D	7	0	4	0	0
3	A	12	0	16	0	0
4	A	5	0	0	0	0
5	A	21	0	9	3	0
6	B	6	0	0	1	0
6	D	6	0	0	2	0
7	B	22	0	11	11	0
7	C	22	0	11	12	0
7	D	22	0	12	13	0
8	A	346	0	0	14	0
8	B	389	0	0	13	0
8	C	332	0	0	8	0
8	D	361	0	0	11	0
All	All	13157	0	11791	157	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 7.

The worst 5 of 157 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:B:210[A]:LYS:NZ	7:B:403:KOU:C4A	2.20	1.05
1:C:210[A]:LYS:HZ1	7:C:402:KOU:C4A	1.75	0.98
1:C:210[B]:LYS:HZ1	7:C:402:KOU:HA	1.29	0.97
1:D:304:ARG:NH1	8:D:794:HOH:O	1.98	0.95
1:B:53:HIS:HD2	1:B:55:GLY:H	1.11	0.95

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	385/397 (97%)	376 (98%)	9 (2%)	0	100	100
1	B	386/397 (97%)	376 (97%)	10 (3%)	0	100	100
1	C	384/397 (97%)	373 (97%)	10 (3%)	1 (0%)	41	24
1	D	389/397 (98%)	380 (98%)	9 (2%)	0	100	100
All	All	1544/1588 (97%)	1505 (98%)	38 (2%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	339	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	304/313 (97%)	298 (98%)	6 (2%)	55	38
1	B	305/313 (97%)	302 (99%)	3 (1%)	76	67
1	C	303/313 (97%)	300 (99%)	3 (1%)	76	67
1	D	308/313 (98%)	302 (98%)	6 (2%)	57	41
All	All	1220/1252 (97%)	1202 (98%)	18 (2%)	67	51

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	161	PRO
1	D	340	LEU
1	D	317	LEU
1	B	365	GLN
1	D	16[B]	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	234	GLN
1	D	240	ASN
1	D	396	GLN
1	B	271	ASN
1	B	246	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)

13 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	SER	C	401	-	5,6,6	1.04	0	5,7,7	1.69	2 (40%)
6	PYR	B	402	-	5,5,5	3.99	4 (80%)	3,6,6	1.84	1 (33%)
2	SER	A	401	-	5,6,6	0.94	0	5,7,7	1.57	2 (40%)
3	GOL	A	403	-	5,5,5	0.91	0	5,5,5	1.85	1 (20%)
6	PYR	D	402	-	5,5,5	2.09	2 (40%)	3,6,6	3.92	2 (66%)
3	GOL	A	402	-	5,5,5	0.78	0	5,5,5	0.40	0
5	0JO	A	405	-	20,21,21	2.66	6 (30%)	23,30,30	1.62	4 (17%)
4	SO4	A	404	-	4,4,4	0.25	0	6,6,6	0.50	0
7	KOU	C	402	-	22,22,22	2.55	9 (40%)	27,31,31	2.18	6 (22%)
7	KOU	B	403	-	22,22,22	2.34	5 (22%)	27,31,31	1.72	6 (22%)
2	SER	B	401	-	5,6,6	1.09	0	5,7,7	1.53	1 (20%)
2	SER	D	401	-	5,6,6	0.73	0	5,7,7	1.73	1 (20%)
7	KOU	D	403	-	22,22,22	2.61	8 (36%)	27,31,31	2.25	6 (22%)

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	SER	C	401	-	-	0/6/6/6	-
6	PYR	B	402	-	-	0/4/4/4	-
2	SER	A	401	-	-	2/6/6/6	-
3	GOL	A	403	-	-	3/4/4/4	-
6	PYR	D	402	-	-	2/4/4/4	-
3	GOL	A	402	-	-	0/4/4/4	-
5	0JO	A	405	-	-	0/10/15/15	0/1/1/1
7	KOU	C	402	-	-	4/17/17/17	0/1/1/1
7	KOU	B	403	-	-	5/17/17/17	0/1/1/1
2	SER	B	401	-	-	0/6/6/6	-
2	SER	D	401	-	-	0/6/6/6	-
7	KOU	D	403	-	-	4/17/17/17	0/1/1/1

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	403	KOU	C3-C2	6.70	1.47	1.40
5	A	405	0JO	C3-C2	5.78	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	405	0JO	C4-C5	5.75	1.49	1.42
7	C	402	KOU	C3-C2	5.74	1.46	1.40
6	B	402	PYR	O3-CA	5.72	1.35	1.23

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	402	KOU	CA-N-C4A	7.65	128.36	117.31
7	D	403	KOU	C3-C4-C5	-6.29	113.43	118.26
7	B	403	KOU	C3-C4-C5	-5.44	114.09	118.26
6	D	402	PYR	OXT-C-O	-5.23	111.64	123.61
7	D	403	KOU	C2A-C2-C3	-5.11	114.58	120.89

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	GOL	O1-C1-C2-C3
6	D	402	PYR	OXT-C-CA-CB
7	B	403	KOU	CB-CA-N-C4A
7	B	403	KOU	N-CA-CB-OG
7	C	402	KOU	C-CA-N-C4A

There are no ring outliers.

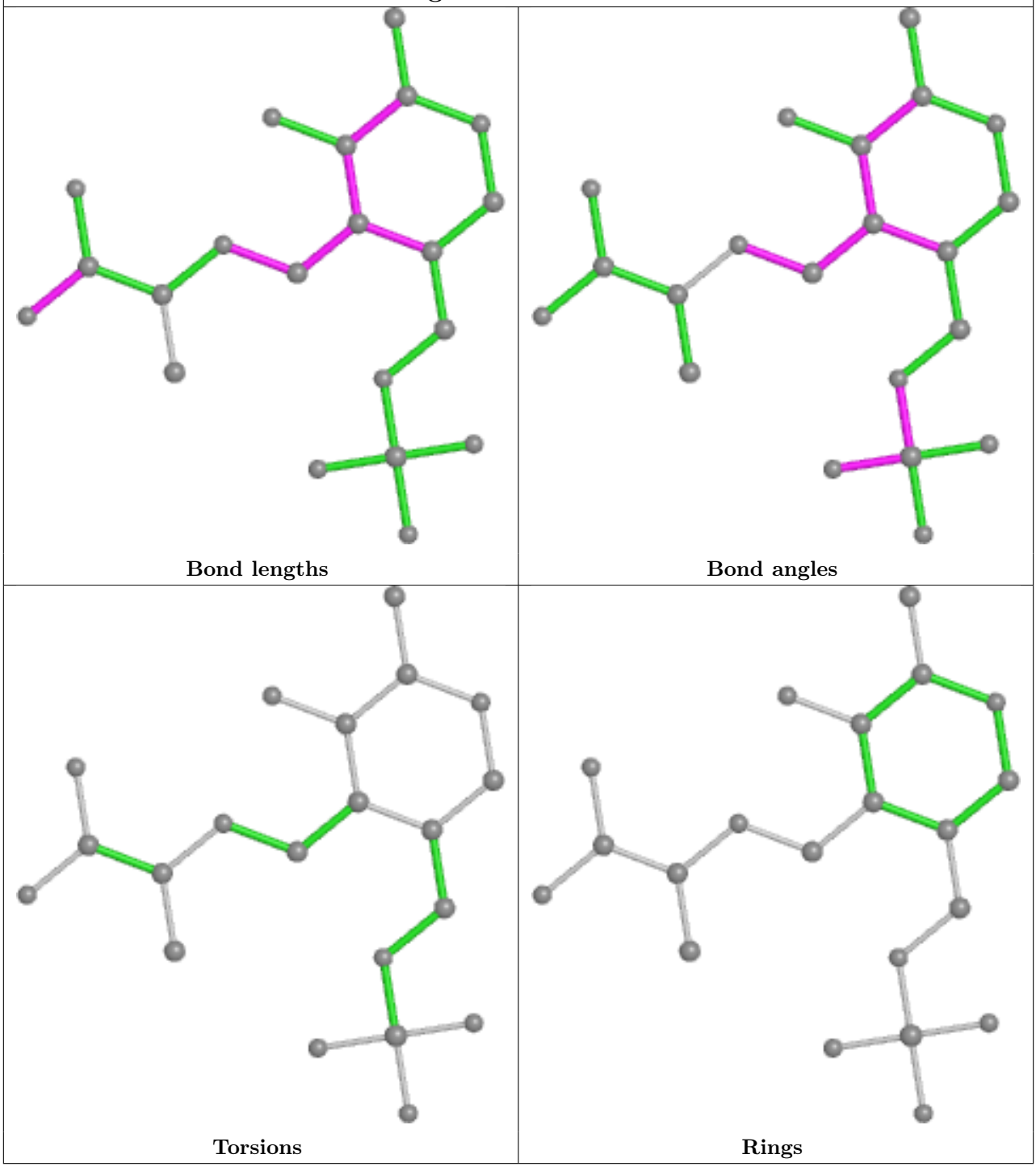
6 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	402	PYR	1	0
6	D	402	PYR	2	0
5	A	405	0JO	3	0
7	C	402	KOU	12	0
7	B	403	KOU	11	0
7	D	403	KOU	13	0

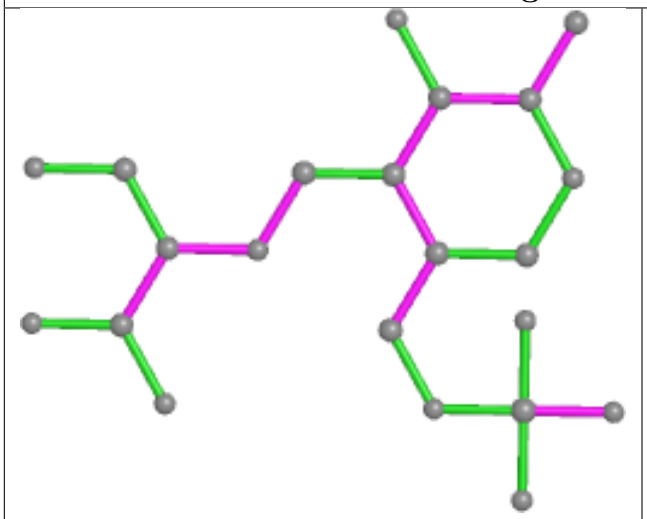
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

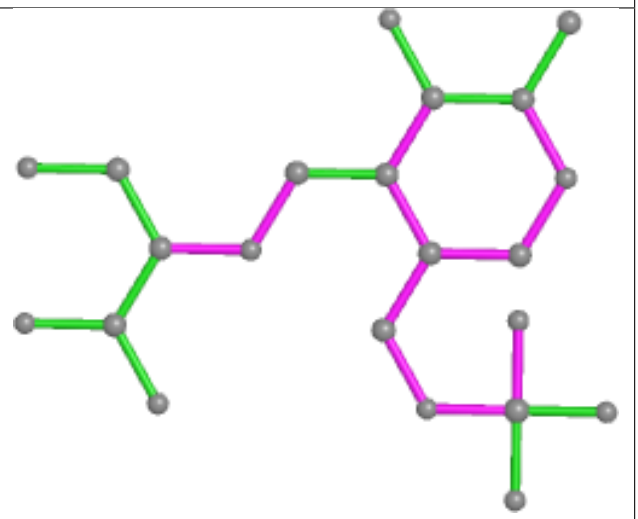
Ligand 0JO A 405



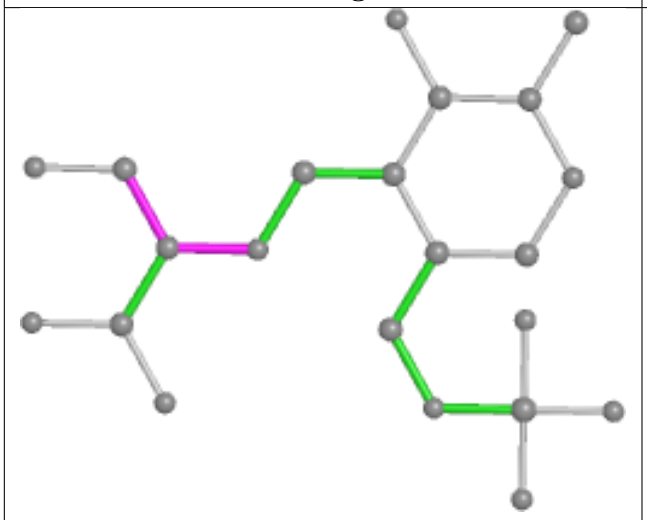
Ligand KOU C 402



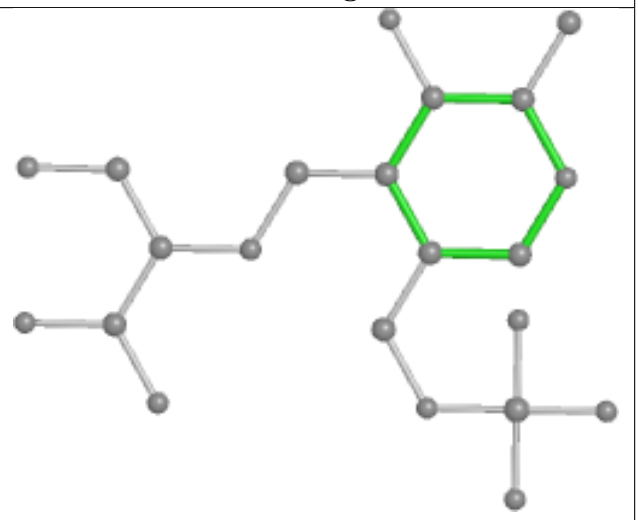
Bond lengths



Bond angles

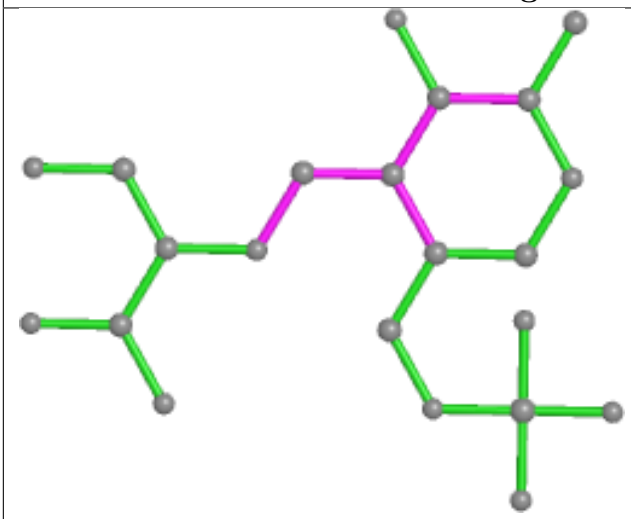


Torsions

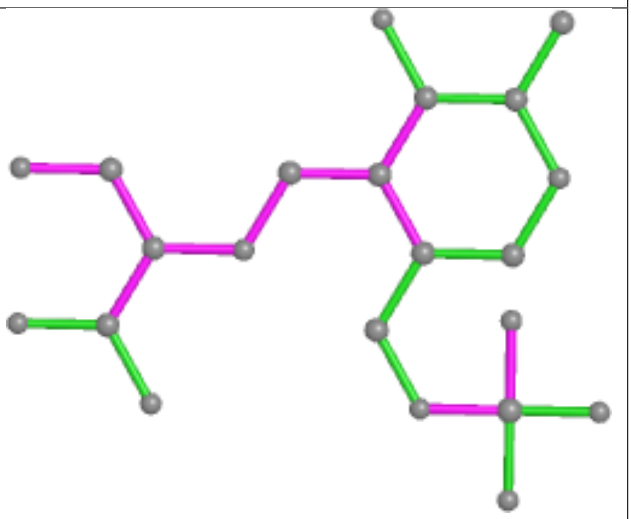


Rings

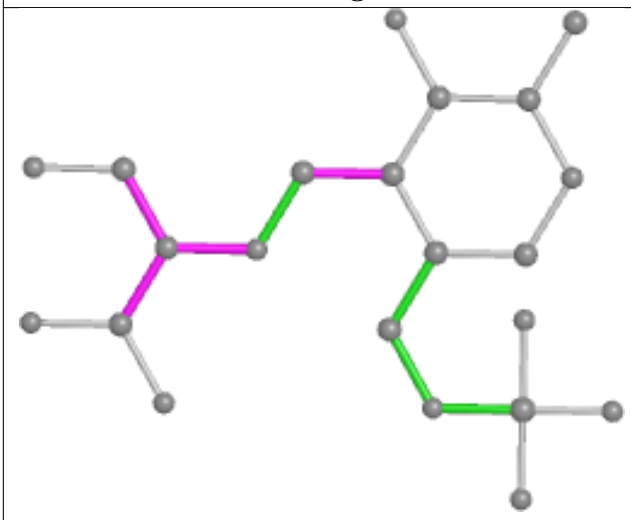
Ligand KOU B 403



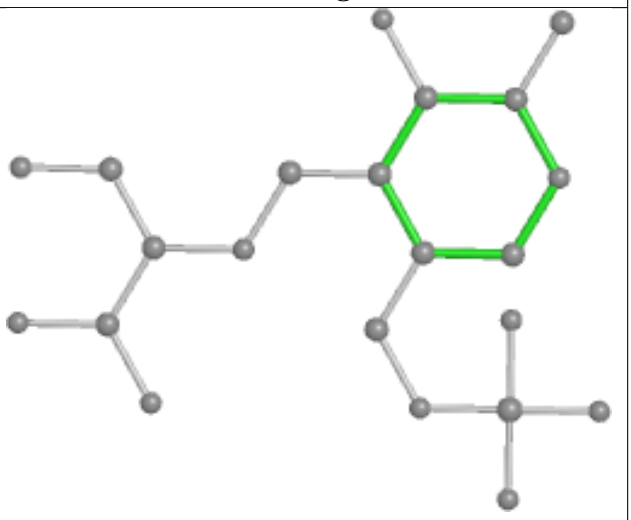
Bond lengths



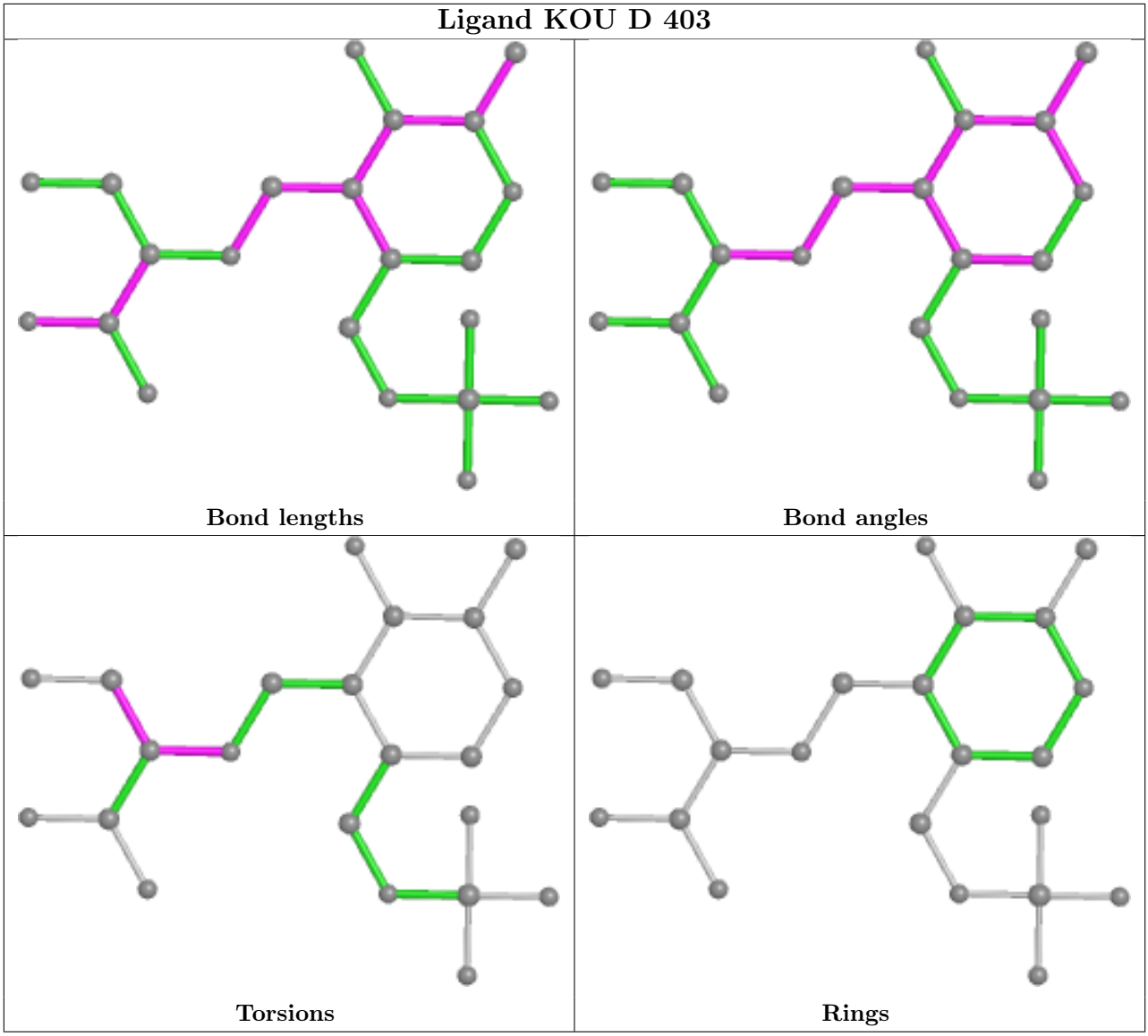
Bond angles



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(Å^2)	Q<0.9
1	A	381/397 (95%)	-0.52	2 (0%)	91	92	7, 12, 24, 35	0
1	B	381/397 (95%)	-0.58	1 (0%)	94	94	7, 11, 22, 33	0
1	C	381/397 (95%)	-0.32	5 (1%)	77	81	7, 14, 28, 36	0
1	D	383/397 (96%)	-0.50	2 (0%)	91	92	8, 12, 23, 39	0
All	All	1526/1588 (96%)	-0.48	10 (0%)	87	90	7, 12, 25, 39	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	ALA	5.3
1	A	14	ALA	4.4
1	C	394	VAL	4.2
1	D	14	ALA	4.0
1	C	14	ALA	3.4

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)

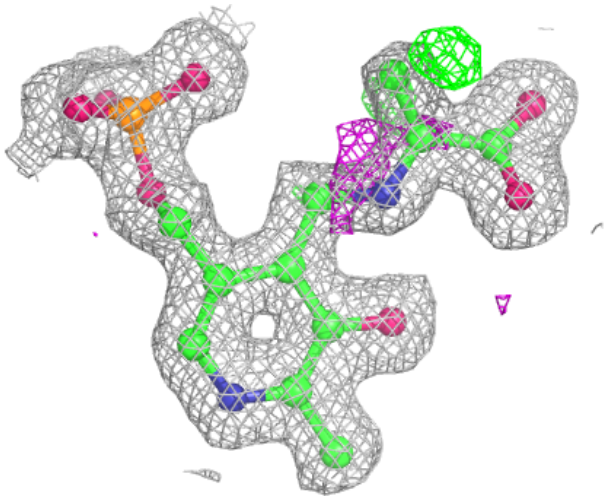
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

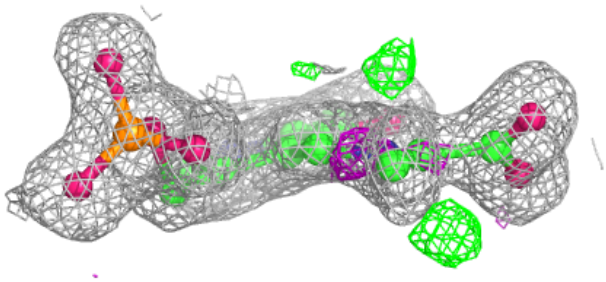
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	A	403	6/6	0.85	0.11	28,33,35,39	0
6	PYR	D	402	6/6	0.85	0.21	25,28,32,34	0
6	PYR	B	402	6/6	0.90	0.15	13,25,31,35	0
2	SER	C	401	7/7	0.95	0.07	18,20,21,24	0
2	SER	A	401	7/7	0.95	0.09	19,22,23,26	0
2	SER	B	401	7/7	0.97	0.05	14,16,17,21	0
2	SER	D	401	7/7	0.97	0.06	17,19,21,21	0
3	GOL	A	402	6/6	0.97	0.06	13,14,17,17	0
4	SO4	A	404	5/5	0.98	0.17	30,38,42,46	0
5	0JO	A	405	21/21	0.98	0.07	7,12,18,20	0
7	KOU	B	403	22/22	0.98	0.08	9,13,20,29	0
7	KOU	C	402	22/22	0.98	0.10	8,17,25,33	0
7	KOU	D	403	22/22	0.98	0.09	9,14,19,29	0

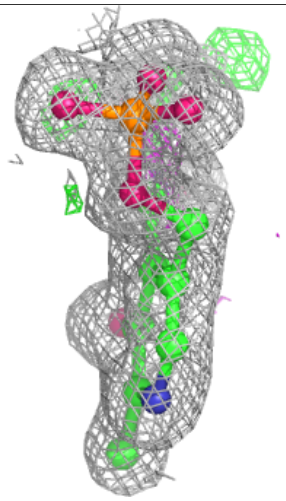
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 0JO A 405:

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

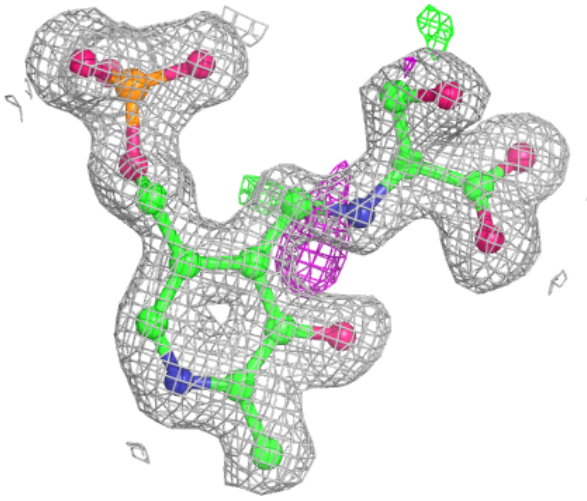


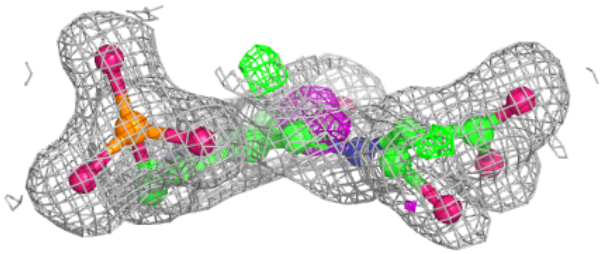




Electron density around KOU B 403:

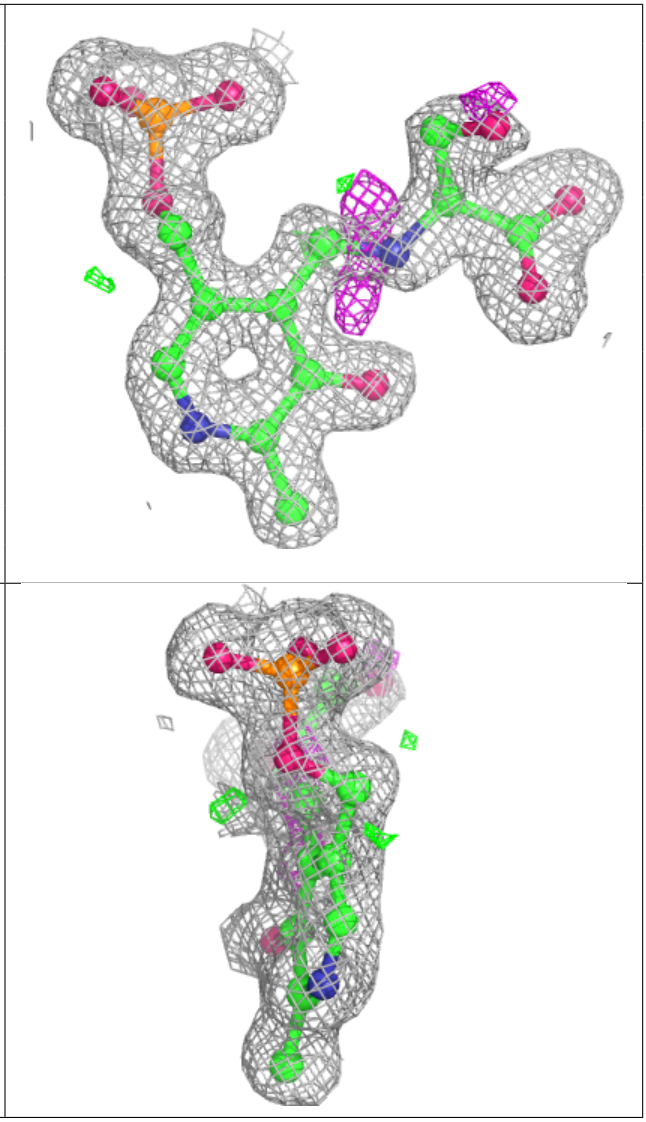
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





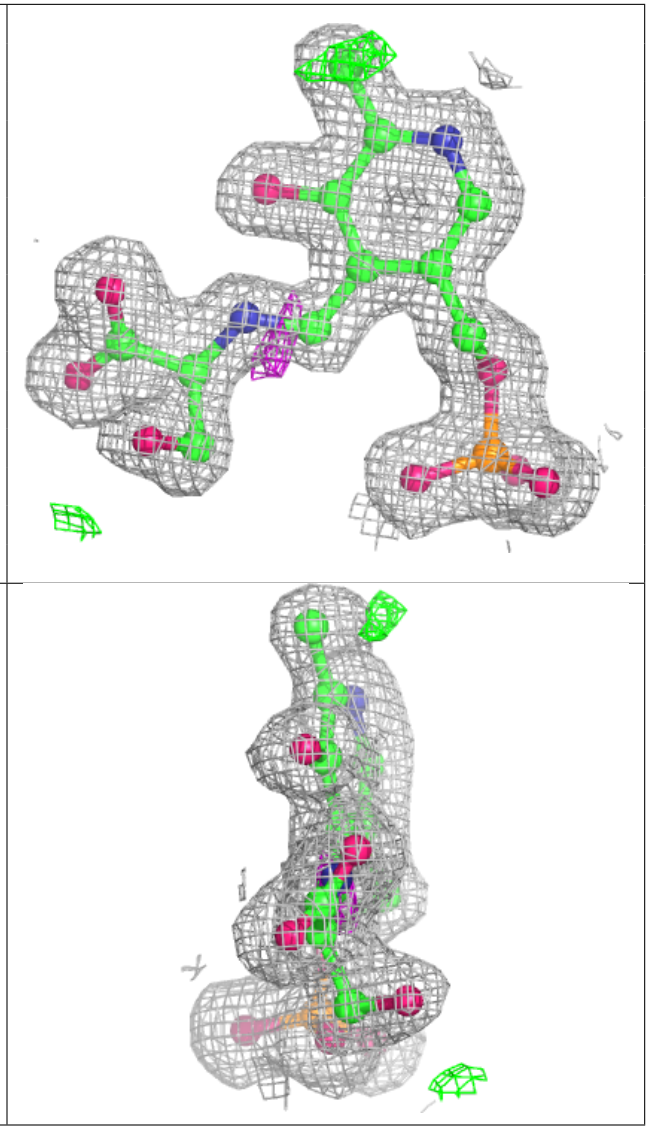
Electron density around KOU C 402:

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



Electron density around KOU D 403:

$2mF_o$-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.