Full wwPDB X-ray Structure Validation Report (i)

Feb 15, 2024 – 02:18 PM EST

PDB ID	:	3QJ9
Title	:	Crystal structure of fatty acid amide hydrolase with small molecule inhibitor
Authors	:	Min, X.; Walker, N.P.C.; Wang, Z.
Deposited on	:	2011-01-28
Resolution	:	2.30 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 (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:

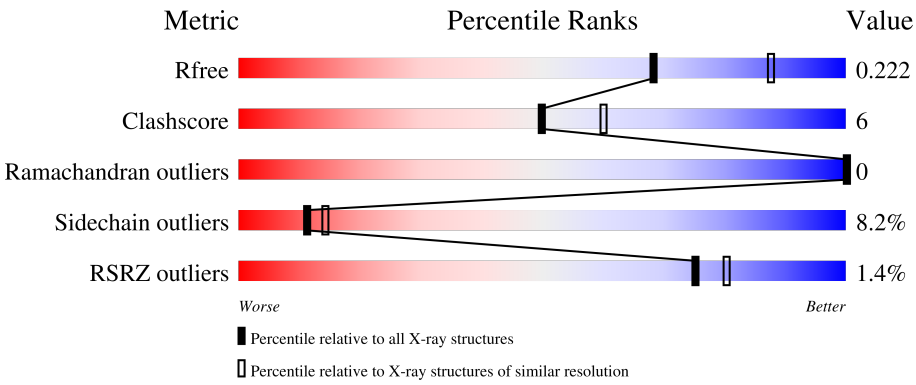
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	587	.% 80% 12% • 6%
1	B	587	2% 77% 13% • 7%

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	EDO	A	584	-	-	-	X
3	EDO	A	586	-	-	X	X
3	EDO	A	588	-	-	-	X
3	EDO	B	581	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 9341 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 Fatty-acid amide hydrolase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	549	4244	2708	722	784	30	0	4	0
1	B	544	4229	2701	719	778	31	0	9	0

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP P97612
A	-6	GLY	-	expression tag	UNP P97612
A	-5	GLY	-	expression tag	UNP P97612
A	-4	SER	-	expression tag	UNP P97612
A	-3	HIS	-	expression tag	UNP P97612
A	-2	HIS	-	expression tag	UNP P97612
A	-1	HIS	-	expression tag	UNP P97612
A	0	HIS	-	expression tag	UNP P97612
A	1	HIS	-	expression tag	UNP P97612
A	2	HIS	-	expression tag	UNP P97612
A	3	GLY	-	expression tag	UNP P97612
A	4	MET	-	expression tag	UNP P97612
A	5	ALA	-	expression tag	UNP P97612
A	6	SER	-	expression tag	UNP P97612
A	7	MET	-	expression tag	UNP P97612
A	8	THR	-	expression tag	UNP P97612
A	9	GLY	-	expression tag	UNP P97612
A	10	GLY	-	expression tag	UNP P97612
A	11	GLN	-	expression tag	UNP P97612
A	12	GLN	-	expression tag	UNP P97612
A	13	MET	-	expression tag	UNP P97612
A	14	GLY	-	expression tag	UNP P97612
A	15	ARG	-	expression tag	UNP P97612
A	16	ASP	-	expression tag	UNP P97612
A	17	LEU	-	expression tag	UNP P97612

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Chain	Residue	Modelled	Actual	Comment	Reference
A	18	TYR	-	expression tag	UNP P97612
A	19	ASP	-	expression tag	UNP P97612
A	20	ASP	-	expression tag	UNP P97612
A	21	ASP	-	expression tag	UNP P97612
A	22	ASP	-	expression tag	UNP P97612
A	23	LYS	-	expression tag	UNP P97612
A	24	ASP	-	expression tag	UNP P97612
A	25	ARG	-	expression tag	UNP P97612
A	26	TRP	-	expression tag	UNP P97612
A	27	GLY	-	expression tag	UNP P97612
A	28	SER	-	expression tag	UNP P97612
A	29	GLU	-	expression tag	UNP P97612
A	30	LEU	-	expression tag	UNP P97612
A	31	GLU	-	expression tag	UNP P97612
B	-7	MET	-	expression tag	UNP P97612
B	-6	GLY	-	expression tag	UNP P97612
B	-5	GLY	-	expression tag	UNP P97612
B	-4	SER	-	expression tag	UNP P97612
B	-3	HIS	-	expression tag	UNP P97612
B	-2	HIS	-	expression tag	UNP P97612
B	-1	HIS	-	expression tag	UNP P97612
B	0	HIS	-	expression tag	UNP P97612
B	1	HIS	-	expression tag	UNP P97612
B	2	HIS	-	expression tag	UNP P97612
B	3	GLY	-	expression tag	UNP P97612
B	4	MET	-	expression tag	UNP P97612
B	5	ALA	-	expression tag	UNP P97612
B	6	SER	-	expression tag	UNP P97612
B	7	MET	-	expression tag	UNP P97612
B	8	THR	-	expression tag	UNP P97612
B	9	GLY	-	expression tag	UNP P97612
B	10	GLY	-	expression tag	UNP P97612
B	11	GLN	-	expression tag	UNP P97612
B	12	GLN	-	expression tag	UNP P97612
B	13	MET	-	expression tag	UNP P97612
B	14	GLY	-	expression tag	UNP P97612
B	15	ARG	-	expression tag	UNP P97612
B	16	ASP	-	expression tag	UNP P97612
B	17	LEU	-	expression tag	UNP P97612
B	18	TYR	-	expression tag	UNP P97612
B	19	ASP	-	expression tag	UNP P97612
B	20	ASP	-	expression tag	UNP P97612

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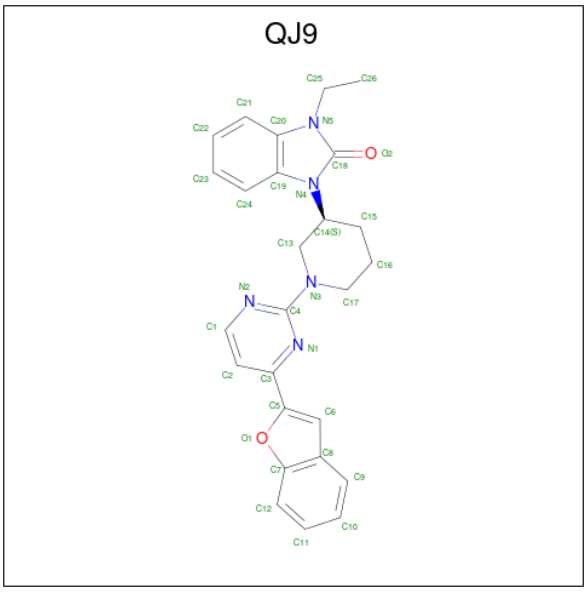
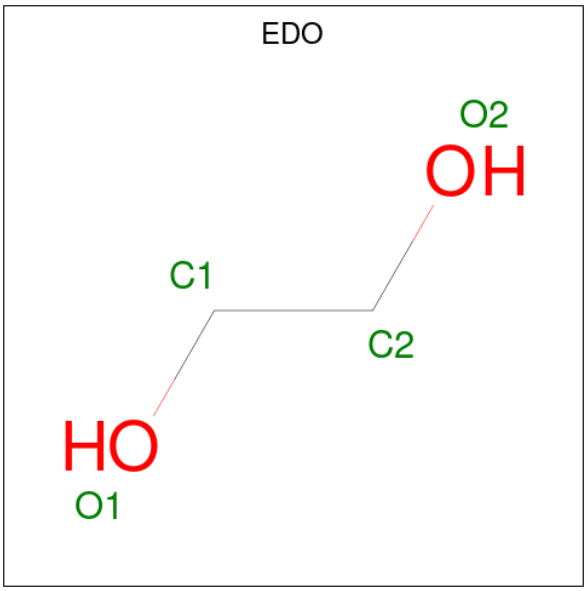
Chain	Residue	Modelled	Actual	Comment	Reference
B	21	ASP	-	expression tag	UNP P97612
B	22	ASP	-	expression tag	UNP P97612
B	23	LYS	-	expression tag	UNP P97612
B	24	ASP	-	expression tag	UNP P97612
B	25	ARG	-	expression tag	UNP P97612
B	26	TRP	-	expression tag	UNP P97612
B	27	GLY	-	expression tag	UNP P97612
B	28	SER	-	expression tag	UNP P97612
B	29	GLU	-	expression tag	UNP P97612
B	30	LEU	-	expression tag	UNP P97612
B	31	GLU	-	expression tag	UNP P97612

- Molecule 2 is 1-{(3S)-1-[4-(1-benzofuran-2-yl)pyrimidin-2-yl]piperidin-3-yl}-3-ethyl-1,3-dihydro-2H-benzimidazol-2-one (three-letter code: QJ9) (formula: $C_{26}H_{25}N_5O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	33	26	5	2	0	0
			Total	C	N	O		
2	B	1	33	26	5	2	0	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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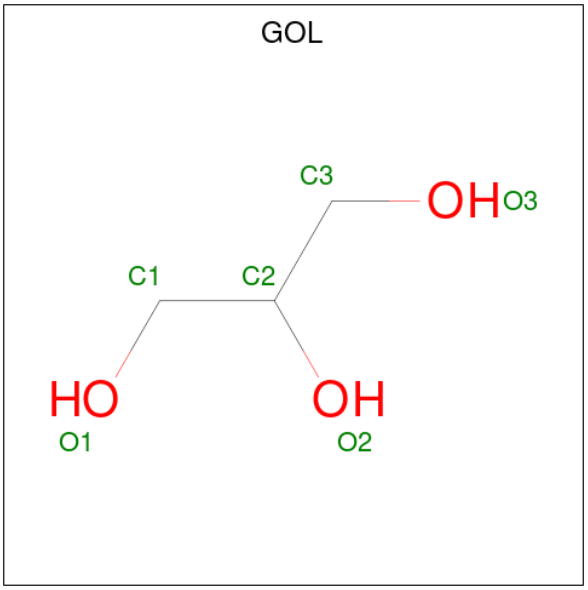
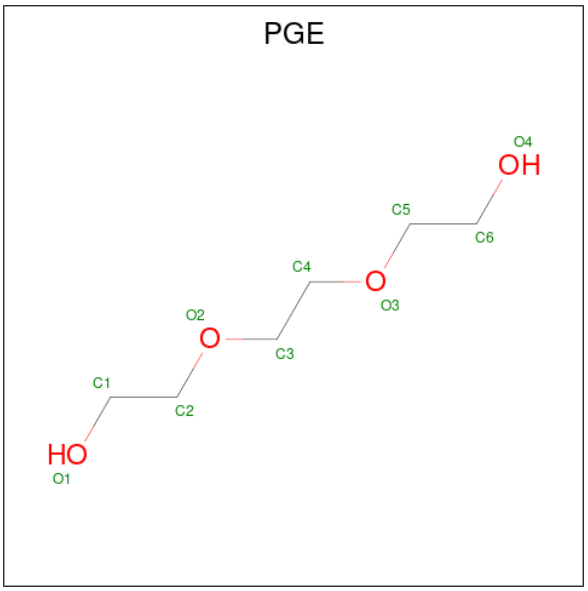
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0
			Total	C	O			
3	B	1	4	2	2		0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	10	6	4	0	0

- Molecule 6 is water.

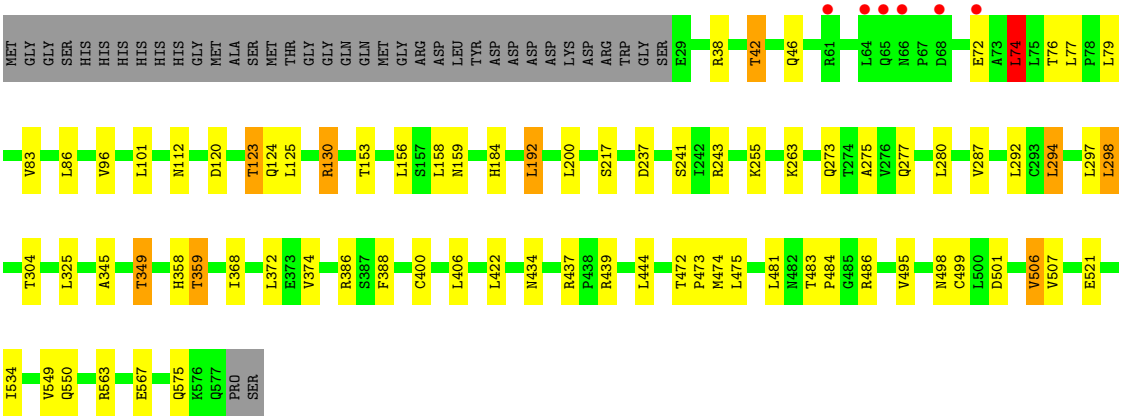
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	331	331	331	0	0
			Total	O		
6	B	365	365	365	0	0

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.

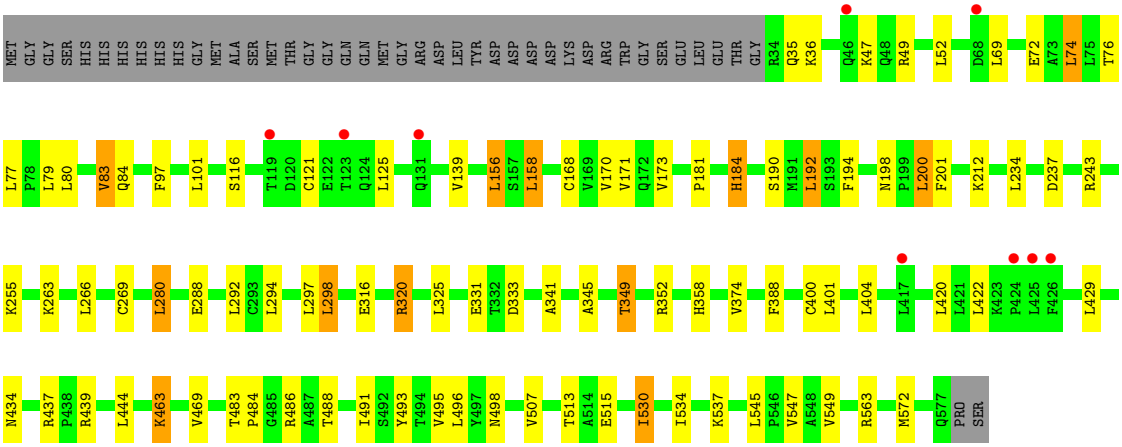
• Molecule 1: Fatty-acid amide hydrolase 1

Chain A:



• Molecule 1: Fatty-acid amide hydrolase 1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	91.56Å 104.72Å 148.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 44.77 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.1 (30.00-2.30) 94.1 (44.77-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$< I/\sigma(I) >$ [1]	3.54 (at 2.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.178 , 0.221 0.179 , 0.222	Depositor DCC
R_{free} test set	3067 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	16.8	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 46.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9341	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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, QJ9, EDO, PGE

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	0.40	0/4349	0.58	1/5899 (0.0%)
1	B	0.40	0/4349	0.60	1/5898 (0.0%)
All	All	0.40	0/8698	0.59	2/11797 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	74	LEU	CA-CB-CG	5.24	127.35	115.30
1	B	74	LEU	CA-CB-CG	5.21	127.27	115.30

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	4244	0	4313	49	0
1	B	4229	0	4313	51	0
2	A	33	0	25	3	0
2	B	33	0	25	2	0
3	A	36	0	54	9	0
3	B	48	0	72	11	0
4	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	6	0	8	0	0
5	A	10	0	14	0	0
6	A	331	0	0	2	0
6	B	365	0	0	3	0
All	All	9341	0	8832	102	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 6.

All (102) 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:192:LEU:HD11	1:B:388:PHE:CE1	1.91	1.04
1:A:325:LEU:H	1:A:358:HIS:HD2	1.08	0.94
1:A:42:THR:HG21	3:A:587:EDO:H21	1.49	0.91
1:B:325:LEU:H	1:B:358:HIS:HD2	1.22	0.85
1:A:495:VAL:HA	1:A:498:ASN:HD22	1.40	0.84
1:B:192:LEU:HD11	1:B:388:PHE:CZ	2.12	0.84
1:B:198:ASN:HD22	1:B:200:LEU:H	1.34	0.75
1:A:192:LEU:HG	1:A:388:PHE:CE1	2.22	0.74
1:A:325:LEU:H	1:A:358:HIS:CD2	1.99	0.73
1:B:194:PHE:H	3:B:583:EDO:H11	1.56	0.71
1:A:472:THR:HG22	1:A:473:PRO:O	1.90	0.71
1:A:325:LEU:N	1:A:358:HIS:HD2	1.86	0.70
1:B:513:THR:HG21	3:B:589:EDO:H21	1.72	0.69
1:A:275:ALA:O	3:A:586:EDO:H22	1.95	0.67
1:B:198:ASN:HD21	1:B:201:PHE:H	1.40	0.67
3:B:585:EDO:H22	6:B:694:HOH:O	1.97	0.65
1:A:192:LEU:HD13	2:A:600:QJ9:C23	2.27	0.64
1:A:501:ASP:HB2	3:A:586:EDO:H12	1.80	0.62
1:B:316:GLU:OE1	1:B:320:ARG:HD2	2.00	0.61
1:B:325:LEU:H	1:B:358:HIS:CD2	2.13	0.60
1:A:474:MET:HG2	1:A:506:VAL:HG21	1.85	0.59
1:A:130:ARG:HH21	1:A:130:ARG:HG3	1.66	0.59
1:B:192:LEU:HB3	2:B:600:QJ9:C22	2.33	0.59
1:A:243:ARG:HH22	1:A:498:ASN:ND2	2.01	0.58
1:A:294:LEU:HD22	1:A:298:LEU:HD22	1.85	0.58
1:B:198:ASN:ND2	1:B:201:PHE:H	2.02	0.58
1:B:79:LEU:O	1:B:83:VAL:HG13	2.04	0.57
1:A:499:CYS:O	3:A:586:EDO:H21	2.05	0.56
1:B:97:PHE:HZ	1:B:121:CYS:HG	1.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:ARG:NH2	1:A:498:ASN:HD21	2.04	0.56
1:B:496:LEU:HD23	3:B:581:EDO:H12	1.87	0.55
1:A:486:ARG:HB3	1:A:534:ILE:HD13	1.88	0.55
1:B:345:ALA:O	1:B:349:THR:HG23	2.07	0.55
1:A:304:THR:O	1:B:463:LYS:NZ	2.34	0.55
1:B:495:VAL:HA	1:B:498:ASN:HD22	1.72	0.54
1:B:434:ASN:ND2	1:B:437:ARG:HH11	2.06	0.54
1:A:74:LEU:HD11	1:A:96:VAL:HB	1.89	0.54
1:B:72:GLU:O	1:B:76:THR:HG23	2.08	0.54
1:B:434:ASN:HD21	1:B:437:ARG:HH11	1.55	0.54
1:A:243:ARG:HD2	1:A:550:GLN:HE21	1.73	0.53
1:B:192:LEU:HD11	1:B:388:PHE:HE1	1.64	0.53
1:B:190[A]:SER:HB2	1:B:269[A]:CYS:SG	2.48	0.53
1:A:38:ARG:O	1:A:42:THR:HG23	2.10	0.52
1:A:120:ASP:O	1:A:123:THR:HG23	2.10	0.52
1:B:237:ASP:HB3	1:B:255:LYS:HG3	1.92	0.51
1:A:277:GLN:HG2	3:A:586:EDO:H11	1.93	0.51
1:B:80:LEU:HG	1:B:288:GLU:HG2	1.93	0.50
2:A:600:QJ9:H15	2:A:600:QJ9:H24	1.92	0.50
1:A:217:SER:H	1:A:241:SER:CB	2.25	0.50
1:B:170:VAL:HG11	1:B:280:LEU:HD11	1.94	0.50
1:B:493:TYR:HA	3:B:581:EDO:C1	2.42	0.50
1:B:345:ALA:O	1:B:349:THR:CG2	2.60	0.50
2:B:600:QJ9:H24	2:B:600:QJ9:H15	1.94	0.50
1:A:345:ALA:O	1:A:349:THR:CG2	2.60	0.49
1:A:192:LEU:HG	1:A:388:PHE:HE1	1.71	0.49
1:A:130:ARG:HH21	1:A:130:ARG:CG	2.24	0.49
1:B:493:TYR:HA	3:B:581:EDO:H11	1.94	0.49
1:B:325:LEU:N	1:B:358:HIS:HD2	2.00	0.49
1:A:243:ARG:HD2	1:A:550:GLN:NE2	2.28	0.49
1:A:243:ARG:HH22	1:A:498:ASN:HD21	1.57	0.48
1:A:72:GLU:O	1:A:76:THR:HG23	2.14	0.47
1:A:112:ASN:HB2	3:A:583:EDO:H11	1.97	0.47
1:A:42:THR:HG21	3:A:587:EDO:C2	2.32	0.47
1:A:243:ARG:HH11	1:A:550:GLN:NE2	2.11	0.47
1:A:358:HIS:HE1	1:A:567:GLU:OE1	1.98	0.47
1:A:79:LEU:O	1:A:83:VAL:HG13	2.15	0.47
1:B:294:LEU:HG	1:B:298:LEU:HD22	1.97	0.46
1:A:46:GLN:HG3	3:A:588:EDO:H12	1.98	0.46
1:B:486:ARG:HB3	1:B:534:ILE:HD13	1.97	0.46
1:B:35:GLN:HB3	6:B:606:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ARG:O	1:A:130:ARG:HD3	2.15	0.46
1:A:192:LEU:HD13	2:A:600:QJ9:H23	1.98	0.46
1:A:474:MET:HG2	1:A:506:VAL:CG2	2.46	0.46
1:B:194:PHE:CD2	3:B:583:EDO:H11	2.51	0.45
1:B:168:CYS:SG	1:B:171:VAL:HG23	2.56	0.45
1:B:243:ARG:NH2	1:B:498:ASN:HD21	2.14	0.45
1:B:352:ARG:HD3	1:B:572:MET:SD	2.56	0.45
1:B:212:LYS:HG3	1:B:545:LEU:HD21	2.00	0.44
1:A:483:THR:N	1:A:484:PRO:CD	2.81	0.44
1:B:496:LEU:CD2	3:B:581:EDO:H12	2.48	0.44
1:A:368:ILE:O	1:A:372:LEU:HG	2.18	0.43
1:B:139:VAL:O	1:B:181:PRO:HA	2.18	0.43
1:B:263:LYS:HG2	1:B:266:LEU:HD22	2.00	0.43
1:A:345:ALA:O	1:A:349:THR:HG22	2.19	0.43
1:B:116:SER:HB3	1:B:184:HIS:HB2	2.01	0.43
1:B:515:GLU:OE2	3:B:589:EDO:H11	2.18	0.43
1:B:537:LYS:HB3	6:B:815:HOH:O	2.18	0.42
1:B:156:LEU:HB3	1:B:158:LEU:HD23	2.01	0.42
1:A:277:GLN:CG	3:A:586:EDO:H11	2.48	0.42
1:B:333:ASP:HB3	3:B:581:EDO:H22	2.01	0.42
1:B:47:LYS:O	3:B:580:EDO:H12	2.20	0.42
1:A:153:THR:HA	1:A:159:ASN:HB2	2.02	0.42
1:A:237:ASP:HB3	1:A:255:LYS:HG3	2.01	0.41
1:A:243:ARG:HH11	1:A:550:GLN:HE22	1.68	0.41
1:A:434:ASN:HD21	1:A:437:ARG:HH11	1.67	0.41
1:B:243:ARG:HH22	1:B:498:ASN:ND2	2.18	0.41
1:A:124:GLN:NE2	6:A:891:HOH:O	2.50	0.41
1:B:341:ALA:HB1	1:B:547:VAL:HG21	2.02	0.41
1:B:488:THR:O	1:B:491:ILE:HG12	2.21	0.41
1:B:483:THR:N	1:B:484:PRO:CD	2.84	0.40
1:A:359:THR:HG21	6:A:725:HOH:O	2.22	0.40
1:B:429:LEU:HD21	1:B:530: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	551/587 (94%)	536 (97%)	15 (3%)	0	100	100
1	B	551/587 (94%)	538 (98%)	13 (2%)	0	100	100
All	All	1102/1174 (94%)	1074 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	470/496 (95%)	432 (92%)	38 (8%)	11	15
1	B	471/496 (95%)	433 (92%)	38 (8%)	11	15
All	All	941/992 (95%)	865 (92%)	76 (8%)	11	15

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	A	74	LEU
1	A	77	LEU
1	A	86	LEU
1	A	101	LEU
1	A	123	THR
1	A	125	LEU
1	A	130	ARG
1	A	156	LEU
1	A	158	LEU
1	A	184	HIS
1	A	192	LEU
1	A	200	LEU
1	A	263	LYS

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Mol	Chain	Res	Type
1	A	273	GLN
1	A	280	LEU
1	A	287	VAL
1	A	292	LEU
1	A	294	LEU
1	A	297	LEU
1	A	298	LEU
1	A	349	THR
1	A	359	THR
1	A	374	VAL
1	A	386	ARG
1	A	400	CYS
1	A	406	LEU
1	A	422	LEU
1	A	439	ARG
1	A	444	LEU
1	A	475	LEU
1	A	481	LEU
1	A	506	VAL
1	A	507	VAL
1	A	521	GLU
1	A	549	VAL
1	A	563	ARG
1	A	575	GLN
1	B	36	LYS
1	B	49	ARG
1	B	52	LEU
1	B	69	LEU
1	B	74	LEU
1	B	77	LEU
1	B	83	VAL
1	B	84	GLN
1	B	101	LEU
1	B	125	LEU
1	B	156	LEU
1	B	158	LEU
1	B	173	VAL
1	B	184	HIS
1	B	192	LEU
1	B	200	LEU
1	B	234	LEU
1	B	280	LEU

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Mol	Chain	Res	Type
1	B	292	LEU
1	B	297	LEU
1	B	298	LEU
1	B	320	ARG
1	B	331	GLU
1	B	349	THR
1	B	374	VAL
1	B	400	CYS
1	B	401	LEU
1	B	404	LEU
1	B	420	LEU
1	B	422	LEU
1	B	439	ARG
1	B	444	LEU
1	B	463	LYS
1	B	469	VAL
1	B	507	VAL
1	B	530	ILE
1	B	549	VAL
1	B	563	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	124	GLN
1	A	127	GLN
1	A	150	HIS
1	A	159	ASN
1	A	203	GLN
1	A	334	ASN
1	A	351	GLN
1	A	358	HIS
1	A	367	ASN
1	A	391	ASN
1	A	434	ASN
1	A	456	GLN
1	A	498	ASN
1	A	550	GLN
1	B	46	GLN
1	B	81	GLN
1	B	87	GLN

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Mol	Chain	Res	Type
1	B	159	ASN
1	B	198	ASN
1	B	259	ASN
1	B	334	ASN
1	B	358	HIS
1	B	391	ASN
1	B	434	ASN
1	B	456	GLN
1	B	498	ASN
1	B	577	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)

26 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	$\#\|Z\| > 2$
3	EDO	B	580	-	3,3,3	0.40	0	2,2,2	0.30	0
3	EDO	B	588	-	3,3,3	0.45	0	2,2,2	0.30	0
3	EDO	A	587	-	3,3,3	0.49	0	2,2,2	0.17	0
3	EDO	B	586	-	3,3,3	0.45	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	EDO	A	583	-	3,3,3	0.46	0	2,2,2	0.26	0
3	EDO	B	592	-	3,3,3	0.46	0	2,2,2	0.37	0
3	EDO	B	584	-	3,3,3	0.43	0	2,2,2	0.42	0
3	EDO	B	583	-	3,3,3	0.45	0	2,2,2	0.28	0
2	QJ9	A	600	-	34,38,38	1.70	8 (23%)	42,55,55	1.85	8 (19%)
3	EDO	B	587	-	3,3,3	0.41	0	2,2,2	0.48	0
3	EDO	A	585	-	3,3,3	0.48	0	2,2,2	0.34	0
5	PGE	A	589	-	9,9,9	0.47	0	8,8,8	0.20	0
3	EDO	A	580	-	3,3,3	0.45	0	2,2,2	0.31	0
3	EDO	A	590	-	3,3,3	0.42	0	2,2,2	0.38	0
3	EDO	B	585	-	3,3,3	0.46	0	2,2,2	0.30	0
3	EDO	B	591	-	3,3,3	0.50	0	2,2,2	0.21	0
3	EDO	A	588	-	3,3,3	0.47	0	2,2,2	0.35	0
3	EDO	A	581	-	3,3,3	0.48	0	2,2,2	0.29	0
3	EDO	B	589	-	3,3,3	0.43	0	2,2,2	0.34	0
4	GOL	A	582	-	5,5,5	0.39	0	5,5,5	0.45	0
3	EDO	A	584	-	3,3,3	0.56	0	2,2,2	0.36	0
3	EDO	B	581	-	3,3,3	0.35	0	2,2,2	0.39	0
4	GOL	B	582	-	5,5,5	0.45	0	5,5,5	0.37	0
3	EDO	B	590	-	3,3,3	0.45	0	2,2,2	0.36	0
2	QJ9	B	600	-	34,38,38	1.68	7 (20%)	42,55,55	1.80	8 (19%)
3	EDO	A	586	-	3,3,3	0.41	0	2,2,2	0.50	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
3	EDO	B	580	-	-	0/1/1/1	-
3	EDO	B	588	-	-	1/1/1/1	-
3	EDO	A	587	-	-	1/1/1/1	-
3	EDO	B	586	-	-	1/1/1/1	-
3	EDO	A	583	-	-	0/1/1/1	-
3	EDO	B	592	-	-	0/1/1/1	-
3	EDO	B	584	-	-	1/1/1/1	-
3	EDO	B	583	-	-	1/1/1/1	-
2	QJ9	A	600	-	-	0/10/24/24	0/6/6/6
3	EDO	B	587	-	-	0/1/1/1	-
3	EDO	A	585	-	-	0/1/1/1	-
5	PGE	A	589	-	-	2/7/7/7	-
3	EDO	A	580	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	590	-	-	1/1/1/1	-
3	EDO	B	585	-	-	1/1/1/1	-
3	EDO	B	591	-	-	1/1/1/1	-
3	EDO	A	588	-	-	1/1/1/1	-
3	EDO	A	581	-	-	1/1/1/1	-
3	EDO	B	589	-	-	1/1/1/1	-
4	GOL	A	582	-	-	4/4/4/4	-
3	EDO	A	584	-	-	1/1/1/1	-
3	EDO	B	581	-	-	1/1/1/1	-
4	GOL	B	582	-	-	4/4/4/4	-
3	EDO	B	590	-	-	1/1/1/1	-
2	QJ9	B	600	-	-	0/10/24/24	0/6/6/6
3	EDO	A	586	-	-	1/1/1/1	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	QJ9	C18-N5	-4.60	1.33	1.38
2	A	600	QJ9	C18-N5	-4.34	1.33	1.38
2	B	600	QJ9	C19-N4	-3.43	1.34	1.40
2	A	600	QJ9	C19-N4	-3.39	1.34	1.40
2	A	600	QJ9	C4-N3	3.29	1.41	1.35
2	A	600	QJ9	C20-N5	-3.06	1.33	1.39
2	B	600	QJ9	C20-N5	-2.99	1.33	1.39
2	B	600	QJ9	C18-N4	-2.92	1.33	1.38
2	A	600	QJ9	C18-N4	-2.90	1.33	1.38
2	B	600	QJ9	C4-N3	2.80	1.40	1.35
2	B	600	QJ9	C2-C1	2.40	1.39	1.36
2	A	600	QJ9	C2-C1	2.40	1.39	1.36
2	A	600	QJ9	C6-C5	-2.06	1.34	1.42
2	B	600	QJ9	C6-C5	-2.05	1.34	1.42
2	A	600	QJ9	C6-C8	-2.00	1.33	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	600	QJ9	N2-C4-N1	-5.57	119.83	126.07
2	A	600	QJ9	C1-N2-C4	5.41	121.05	115.08
2	B	600	QJ9	C1-N2-C4	5.27	120.91	115.08
2	B	600	QJ9	N2-C4-N1	-4.94	120.53	126.07
2	B	600	QJ9	C2-C1-N2	-4.19	119.68	123.81
2	A	600	QJ9	N2-C4-N3	3.94	121.16	116.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	600	QJ9	C2-C1-N2	-3.79	120.08	123.81
2	B	600	QJ9	N2-C4-N3	3.50	120.69	116.90
2	B	600	QJ9	N4-C18-N5	3.14	109.17	106.89
2	A	600	QJ9	N4-C18-N5	3.14	109.17	106.89
2	B	600	QJ9	C13-C14-N4	-3.00	110.36	116.59
2	A	600	QJ9	C13-C14-N4	-2.89	110.61	116.59
2	B	600	QJ9	C17-N3-C13	2.50	119.82	112.55
2	A	600	QJ9	C17-N3-C13	2.37	119.45	112.55
2	A	600	QJ9	C25-N5-C18	2.29	125.95	123.55
2	B	600	QJ9	C6-C8-C7	2.01	108.02	106.27

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	582	GOL	O1-C1-C2-C3
4	A	582	GOL	C1-C2-C3-O3
4	B	582	GOL	C1-C2-C3-O3
4	B	582	GOL	O2-C2-C3-O3
4	B	582	GOL	O1-C1-C2-C3
4	A	582	GOL	O1-C1-C2-O2
4	A	582	GOL	O2-C2-C3-O3
4	B	582	GOL	O1-C1-C2-O2
3	A	580	EDO	O1-C1-C2-O2
3	A	584	EDO	O1-C1-C2-O2
3	A	588	EDO	O1-C1-C2-O2
3	B	581	EDO	O1-C1-C2-O2
3	B	588	EDO	O1-C1-C2-O2
3	B	590	EDO	O1-C1-C2-O2
5	A	589	PGE	O2-C3-C4-O3
3	B	586	EDO	O1-C1-C2-O2
3	B	589	EDO	O1-C1-C2-O2
3	B	591	EDO	O1-C1-C2-O2
3	A	586	EDO	O1-C1-C2-O2
3	A	587	EDO	O1-C1-C2-O2
5	A	589	PGE	C4-C3-O2-C2
3	A	581	EDO	O1-C1-C2-O2
3	A	590	EDO	O1-C1-C2-O2
3	B	584	EDO	O1-C1-C2-O2
3	B	585	EDO	O1-C1-C2-O2
3	B	583	EDO	O1-C1-C2-O2

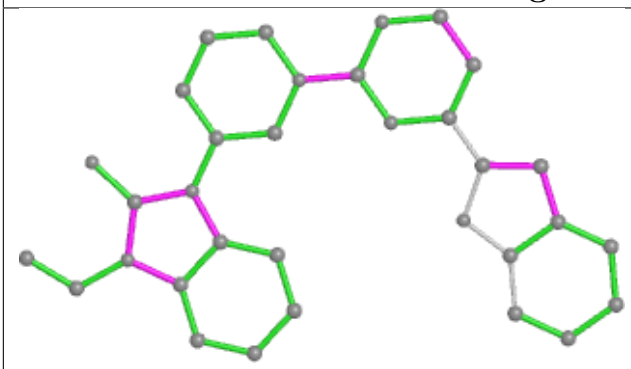
There are no ring outliers.

11 monomers are involved in 25 short contacts:

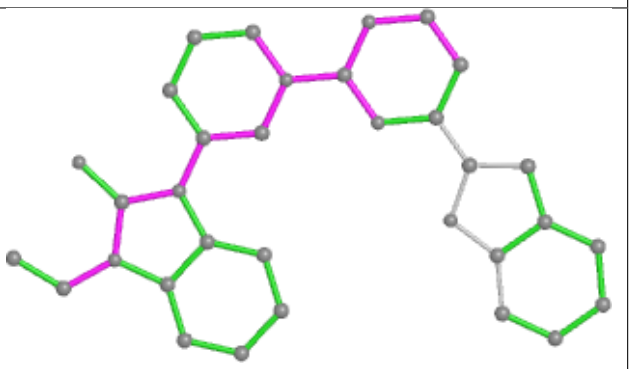
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	580	EDO	1	0
3	A	587	EDO	2	0
3	A	583	EDO	1	0
3	B	583	EDO	2	0
2	A	600	QJ9	3	0
3	B	585	EDO	1	0
3	A	588	EDO	1	0
3	B	589	EDO	2	0
3	B	581	EDO	5	0
2	B	600	QJ9	2	0
3	A	586	EDO	5	0

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.

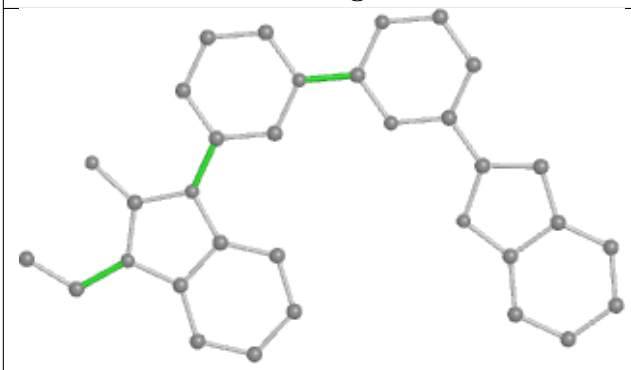
Ligand QJ9 A 600



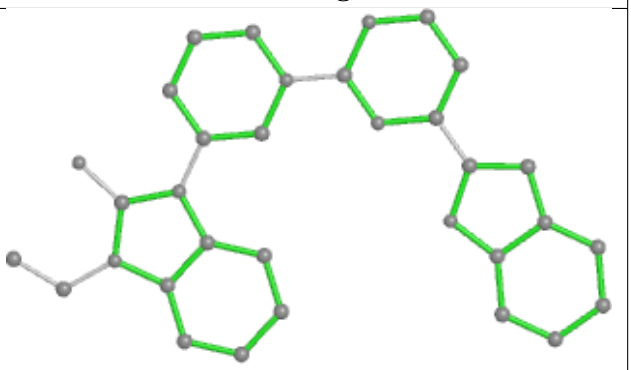
Bond lengths



Bond angles

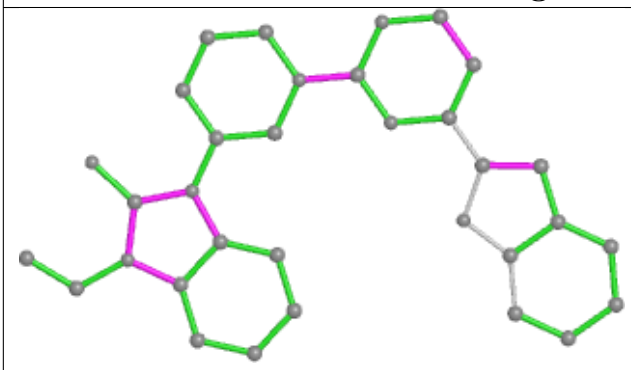


Torsions

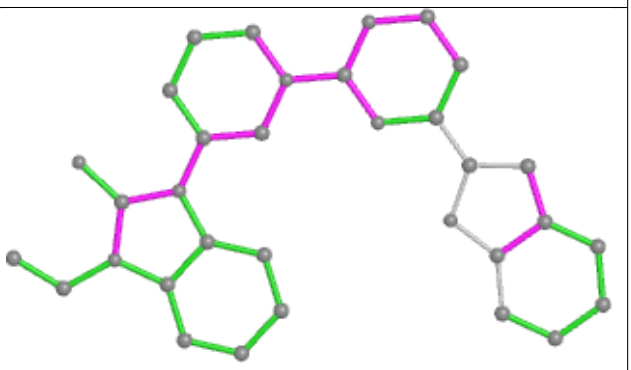


Rings

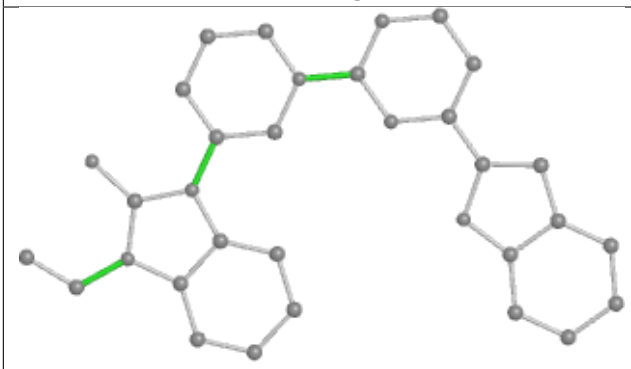
Ligand QJ9 B 600



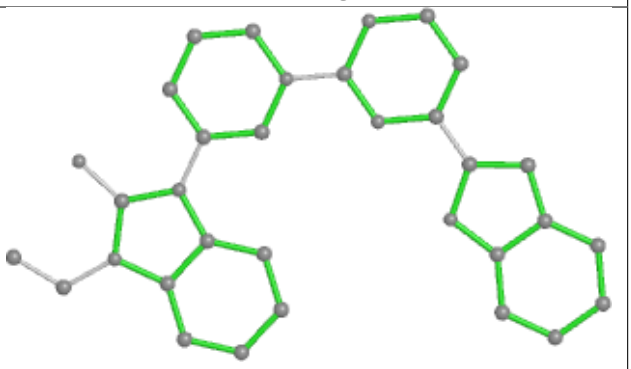
Bond lengths



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	549/587 (93%)	-0.36	6 (1%)	80	85	8, 15, 28, 44	0
1	B	544/587 (92%)	-0.43	9 (1%)	70	76	7, 13, 29, 37	0
All	All	1093/1174 (93%)	-0.40	15 (1%)	75	80	7, 14, 28, 44	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	425	LEU	3.1
1	A	64	LEU	3.0
1	A	72	GLU	2.5
1	B	424	PRO	2.5
1	B	123	THR	2.4
1	B	46	GLN	2.3
1	B	131	GLN	2.3
1	B	68	ASP	2.3
1	A	65	GLN	2.3
1	B	119	THR	2.2
1	A	61	ARG	2.2
1	B	426	PHE	2.1
1	A	66	ASN	2.1
1	B	417	LEU	2.1
1	A	68	ASP	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)

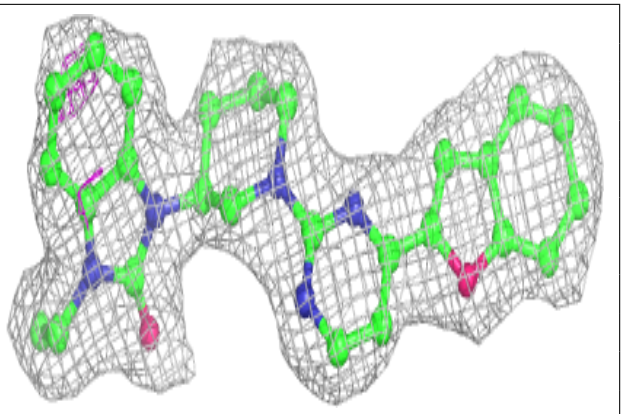
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

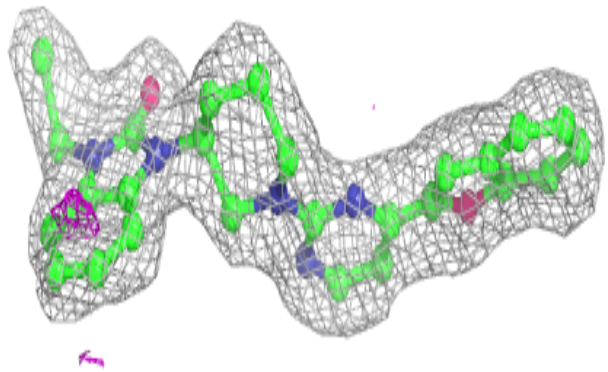
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	A	584	4/4	0.35	0.75	103,103,103,104	0
3	EDO	A	585	4/4	0.61	0.36	49,49,50,50	0
3	EDO	A	587	4/4	0.61	0.32	43,43,43,44	0
3	EDO	A	586	4/4	0.66	0.52	65,66,66,66	0
3	EDO	B	591	4/4	0.68	0.28	41,41,41,41	0
4	GOL	A	582	6/6	0.74	0.21	39,39,40,40	0
3	EDO	A	588	4/4	0.77	0.46	49,49,49,50	0
3	EDO	A	581	4/4	0.79	0.22	42,43,43,44	0
3	EDO	B	588	4/4	0.79	0.30	44,44,45,45	0
4	GOL	B	582	6/6	0.80	0.21	28,30,31,31	0
3	EDO	B	583	4/4	0.81	0.29	48,48,48,48	0
3	EDO	A	580	4/4	0.86	0.18	34,35,36,37	0
3	EDO	B	584	4/4	0.86	0.30	45,45,46,46	0
3	EDO	B	585	4/4	0.87	0.19	28,29,29,30	0
3	EDO	B	586	4/4	0.87	0.18	38,38,38,39	0
3	EDO	B	587	4/4	0.88	0.30	48,48,48,49	0
3	EDO	B	592	4/4	0.88	0.22	37,37,38,38	0
5	PGE	A	589	10/10	0.88	0.21	44,45,45,45	0
3	EDO	B	589	4/4	0.89	0.22	25,26,26,26	0
3	EDO	B	590	4/4	0.90	0.12	37,37,37,37	0
2	QJ9	B	600	33/33	0.90	0.13	24,26,28,28	0
3	EDO	A	590	4/4	0.91	0.15	19,20,20,21	0
2	QJ9	A	600	33/33	0.91	0.12	18,21,27,27	0
3	EDO	A	583	4/4	0.92	0.17	37,37,37,38	0
3	EDO	B	581	4/4	0.93	0.27	31,32,33,34	0
3	EDO	B	580	4/4	0.94	0.33	26,27,28,29	0

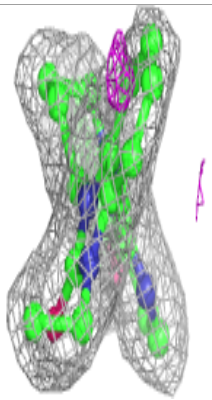
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around QJ9 B 600:

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

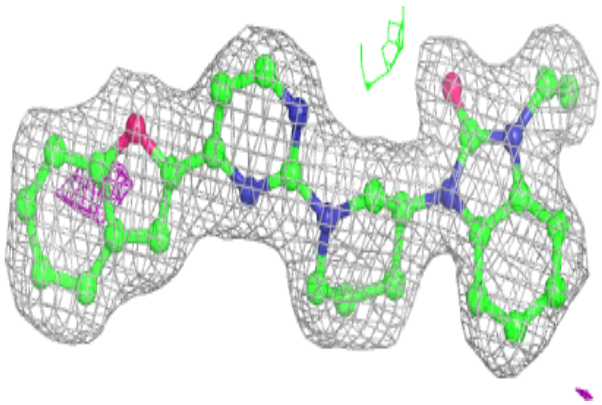


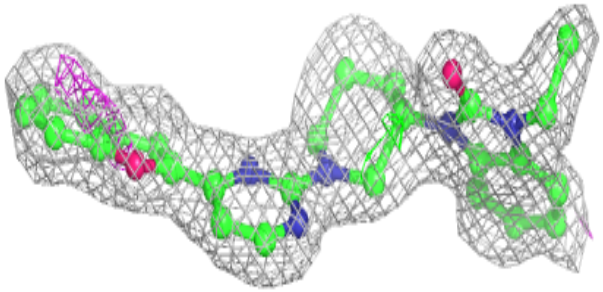


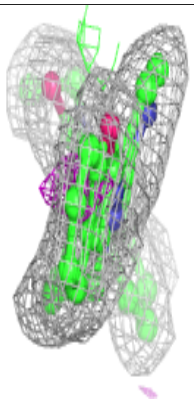


Electron density around QJ9 A 600:

$2mF_o$-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.