Full wwPDB X-ray Structure Validation Report (i)

Aug 26, 2023 – 12:27 PM EDT

PDB ID	:	3F6X
Title	:	c-Src kinase domain in complex with small molecule inhibitor
Authors	:	Seeliger, M.A.; Statsuk, A.V.; Maly, D.J.; Patrick, P.Z.; Kuriyan, J.; Shokat, K.M.
Deposited on	:	2008-11-06
Resolution	:	2.35 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:

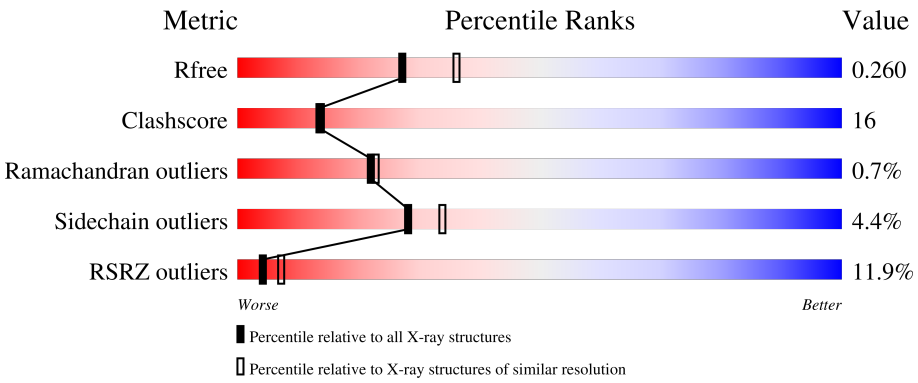
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	286	13% 66% 24% • 8%
1	B	286	9% 67% 25% • 7%
1	C	286	12% 59% 30% • 7%
1	D	286	10% 66% 26% • 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
2	IHH	B	1	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8872 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.

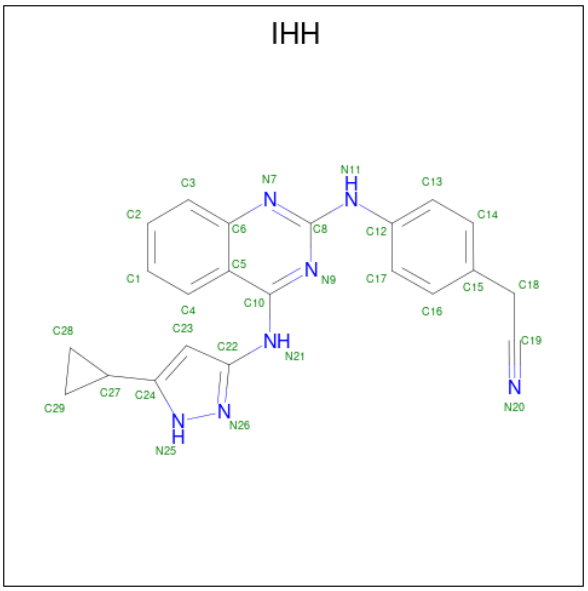
- Molecule 1 is a protein called Proto-oncogene tyrosine-protein kinase Src.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	264	2124	1364	355	389	16	0	0	0
1	B	267	2137	1372	358	391	16	0	0	0
1	C	265	2133	1373	357	387	16	0	0	0
1	D	267	2142	1375	359	391	17	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	248	GLY	-	expression tag	UNP P00523
A	249	HIS	-	expression tag	UNP P00523
A	250	MET	-	expression tag	UNP P00523
B	248	GLY	-	expression tag	UNP P00523
B	249	HIS	-	expression tag	UNP P00523
B	250	MET	-	expression tag	UNP P00523
C	248	GLY	-	expression tag	UNP P00523
C	249	HIS	-	expression tag	UNP P00523
C	250	MET	-	expression tag	UNP P00523
D	248	GLY	-	expression tag	UNP P00523
D	249	HIS	-	expression tag	UNP P00523
D	250	MET	-	expression tag	UNP P00523

- Molecule 2 is [4-({4-[(5-cyclopropyl-1H-pyrazol-3-yl)amino]quinazolin-2-yl}amino)phenyl]acetonitrile (three-letter code: IHH) (formula: $C_{22}H_{19}N_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N 29 22 7	0	0
2	B	1	Total C N 29 22 7	0	0
2	C	1	Total C N 29 22 7	0	0
2	D	1	Total C N 29 22 7	0	0

- Molecule 3 is water.

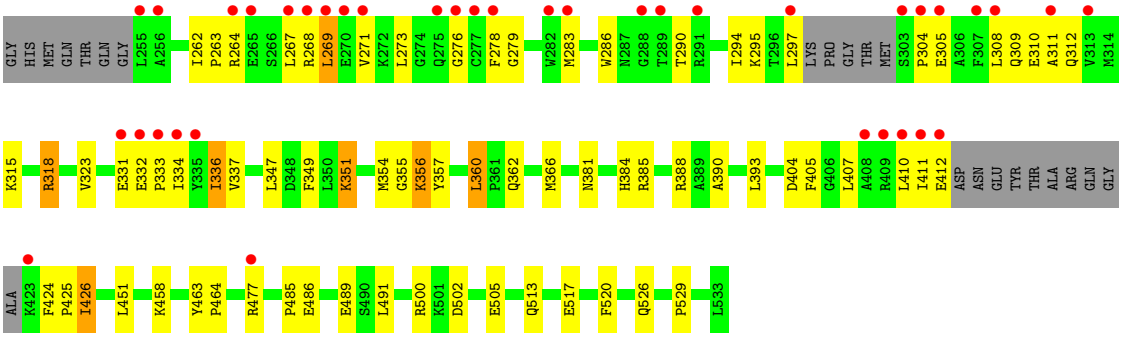
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	51	Total O 51 51	0	0
3	B	57	Total O 57 57	0	0
3	C	47	Total O 47 47	0	0
3	D	65	Total O 65 65	0	0

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.

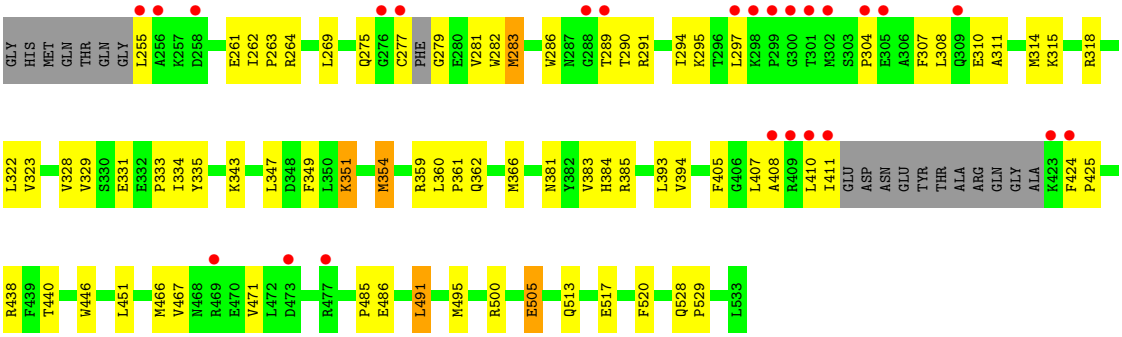
• Molecule 1: Proto-oncogene tyrosine-protein kinase Src

Chain A:



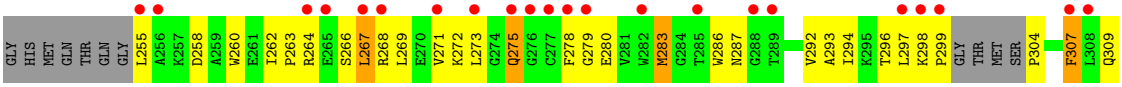
• Molecule 1: Proto-oncogene tyrosine-protein kinase Src

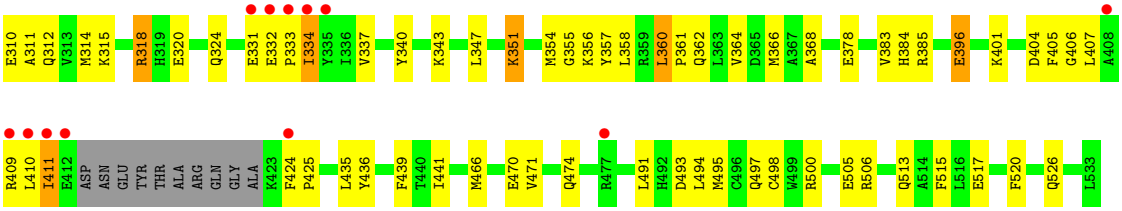
Chain B:



• Molecule 1: Proto-oncogene tyrosine-protein kinase Src

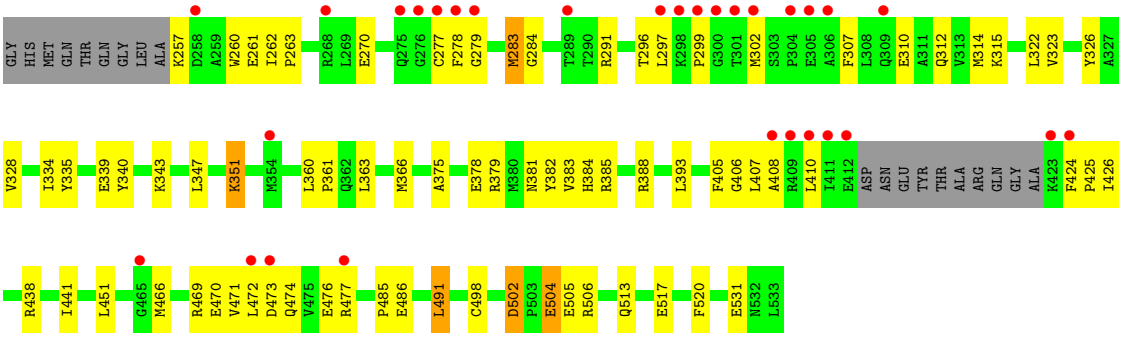
Chain C:





- Molecule 1: Proto-oncogene tyrosine-protein kinase Src

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	63.70Å 74.49Å 84.65Å 89.06° 89.97° 78.43°	Depositor
Resolution (Å)	34.99 – 2.35 34.98 – 2.35	Depositor EDS
% Data completeness (in resolution range)	79.8 (34.99-2.35) 80.2 (34.98-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	0.00 (at 2.34Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.234 , 0.266 0.227 , 0.260	Depositor DCC
R_{free} test set	2984 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 55.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8872	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 91.92 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1880e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: IHH

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	0.36	0/2174	0.58	0/2942
1	B	0.35	0/2187	0.61	0/2961
1	C	0.36	0/2184	0.61	0/2955
1	D	0.36	0/2194	0.60	0/2970
All	All	0.36	0/8739	0.60	0/11828

There are no bond length outliers.

There are no bond angle outliers.

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	2124	0	2104	68	0
1	B	2137	0	2123	57	0
1	C	2133	0	2125	75	0
1	D	2142	0	2124	54	0
2	A	29	0	19	6	0
2	B	29	0	19	9	0
2	C	29	0	19	5	0
2	D	29	0	19	8	0
3	A	51	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	57	0	0	1	0
3	C	47	0	0	1	0
3	D	65	0	0	1	0
All	All	8872	0	8552	275	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 16.

All (275) 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:C:334:ILE:HD13	1:C:334:ILE:H	1.29	0.96
2:B:1:IHH:H17	2:B:1:IHH:H23	1.52	0.91
2:D:1:IHH:H17	2:D:1:IHH:H23	1.54	0.90
1:C:262:ILE:HD13	1:C:286:TRP:HE1	1.44	0.83
2:B:1:IHH:H17	2:B:1:IHH:N9	1.93	0.82
1:B:289:THR:HG23	1:B:290:THR:HG23	1.61	0.80
1:D:504:GLU:CD	1:D:504:GLU:H	1.85	0.80
1:C:513:GLN:O	1:C:517:GLU:HG3	1.84	0.77
2:C:1:IHH:H16	2:C:1:IHH:H28	1.66	0.77
2:A:1:IHH:H16	2:A:1:IHH:H28	1.64	0.76
2:D:1:IHH:H17	2:D:1:IHH:N9	1.99	0.75
2:A:1:IHH:N9	2:A:1:IHH:H17	2.02	0.75
1:C:334:ILE:H	1:C:334:ILE:CD1	1.99	0.75
1:C:310:GLU:HG3	1:C:405:PHE:C	2.08	0.74
1:B:269:LEU:HD23	1:B:294:ILE:HD12	1.68	0.74
1:C:262:ILE:CD1	1:C:286:TRP:HE1	2.02	0.73
1:D:283:MET:HG3	1:D:340:TYR:CE1	2.24	0.72
1:C:309:GLN:NE2	1:C:312:GLN:HG3	2.05	0.72
1:B:347:LEU:O	1:B:351:LYS:HG2	1.88	0.71
1:D:363:LEU:HA	1:D:366:MET:HE3	1.71	0.71
1:C:410:LEU:H	1:C:410:LEU:HD23	1.55	0.71
1:C:297:LEU:O	1:C:297:LEU:HD23	1.90	0.71
1:C:298:LYS:HB3	1:C:299:PRO:HD3	1.73	0.70
1:B:297:LEU:HD22	1:B:307:PHE:HB2	1.73	0.70
1:C:334:ILE:HD13	1:C:334:ILE:N	2.07	0.69
1:D:297:LEU:HD11	1:D:302:MET:HB2	1.73	0.69
1:D:283:MET:HE1	1:D:291:ARG:HH21	1.58	0.69
1:A:347:LEU:O	1:A:351:LYS:HG2	1.93	0.69
1:B:466:MET:CE	1:B:471:VAL:HA	2.23	0.69
1:C:356:LYS:HE3	1:C:357:TYR:CZ	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:473:ASP:HB2	1:D:477:ARG:NH2	2.07	0.69
1:B:311:ALA:O	1:B:315:LYS:HG3	1.93	0.68
1:B:393:LEU:HD11	2:B:1:IHH:H28	1.75	0.68
1:A:268:ARG:O	1:A:268:ARG:HG3	1.93	0.68
1:D:278:PHE:O	1:D:297:LEU:HD12	1.93	0.68
2:C:1:IHH:H17	2:C:1:IHH:N9	2.08	0.68
1:C:378:GLU:HG3	1:C:441:ILE:HG12	1.76	0.66
1:A:310:GLU:HG3	1:A:405:PHE:C	2.16	0.66
1:A:407:LEU:HD13	1:A:424:PHE:CE1	2.31	0.65
1:A:295:LYS:HB2	1:A:336:ILE:CG2	2.27	0.65
1:B:381:ASN:O	1:B:410:LEU:HD12	1.97	0.64
1:A:349:PHE:CZ	1:A:354:MET:HG2	2.32	0.64
2:A:1:IHH:H17	2:A:1:IHH:H23	1.79	0.64
1:A:513:GLN:O	1:A:517:GLU:HG3	1.98	0.64
1:A:334:ILE:HD12	1:A:334:ILE:H	1.62	0.63
2:C:1:IHH:H17	2:C:1:IHH:H23	1.80	0.63
1:D:257:LYS:HE3	1:D:261:GLU:HG2	1.80	0.63
1:A:407:LEU:HD13	1:A:424:PHE:HE1	1.63	0.63
1:C:262:ILE:CG1	1:C:263:PRO:HD2	2.29	0.62
1:C:280:GLU:HG3	1:C:299:PRO:HG2	1.79	0.62
1:A:331:GLU:O	1:A:334:ILE:HA	2.00	0.62
1:B:297:LEU:HD22	1:B:307:PHE:CB	2.29	0.62
1:B:323:VAL:HG21	1:B:393:LEU:HD12	1.82	0.62
1:A:410:LEU:O	1:A:412:GLU:HG3	2.01	0.61
1:D:278:PHE:HA	1:D:302:MET:HG3	1.82	0.61
1:A:500:ARG:HD3	1:A:505:GLU:HG3	1.82	0.61
1:D:381:ASN:O	1:D:410:LEU:HD12	2.00	0.61
1:D:426:ILE:HD11	1:D:472:LEU:HB2	1.83	0.61
1:B:383:VAL:HG23	1:B:411:ILE:HB	1.84	0.60
1:C:356:LYS:HE3	1:C:357:TYR:OH	2.02	0.60
1:A:279:GLY:HA3	1:A:297:LEU:HA	1.84	0.59
1:C:262:ILE:HG13	1:C:263:PRO:HD2	1.84	0.59
1:C:267:LEU:N	1:C:267:LEU:HD23	2.16	0.59
1:A:354:MET:HE1	1:C:435:LEU:HB3	1.84	0.59
1:D:312:GLN:HA	1:D:315:LYS:HD3	1.85	0.58
1:A:273:LEU:HD11	1:A:283:MET:HB2	1.85	0.58
1:D:466:MET:CE	1:D:471:VAL:HA	2.34	0.57
1:B:362:GLN:O	1:B:366:MET:HG3	2.03	0.57
1:C:318:ARG:HD3	1:C:324:GLN:OE1	2.03	0.57
1:C:355:GLY:HA2	1:C:358:LEU:HD12	1.86	0.57
1:B:354:MET:HB2	3:B:72:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:513:GLN:O	1:B:517:GLU:HG3	2.04	0.57
1:D:513:GLN:O	1:D:517:GLU:HG3	2.05	0.56
2:B:1:IHH:N9	2:B:1:IHH:C17	2.63	0.56
1:A:269:LEU:HD23	1:A:269:LEU:H	1.70	0.56
1:A:334:ILE:HD12	1:A:334:ILE:N	2.20	0.56
2:D:1:IHH:H16	2:D:1:IHH:C28	2.34	0.56
1:A:295:LYS:HB2	1:A:336:ILE:HG23	1.86	0.56
1:D:322:LEU:HG	1:D:405:PHE:HZ	1.70	0.56
1:A:311:ALA:O	1:A:315:LYS:HG3	2.05	0.56
1:D:378:GLU:HG3	1:D:441:ILE:HG12	1.86	0.56
1:B:467:VAL:O	1:B:471:VAL:HG23	2.06	0.55
1:A:451:LEU:O	1:A:451:LEU:HD23	2.07	0.55
1:A:485:PRO:O	1:A:486:GLU:HB2	2.06	0.55
2:B:1:IHH:H16	2:B:1:IHH:C28	2.35	0.55
1:C:310:GLU:HG2	1:C:314:MET:HE2	1.89	0.55
1:C:294:ILE:N	1:C:294:ILE:HD12	2.22	0.54
1:C:383:VAL:HG23	1:C:411:ILE:HG23	1.90	0.54
1:B:451:LEU:HD23	1:B:451:LEU:O	2.07	0.54
1:C:466:MET:CE	1:C:471:VAL:HA	2.38	0.54
1:B:261:GLU:HA	1:B:328:VAL:O	2.06	0.54
1:B:318:ARG:NH1	1:B:318:ARG:HB2	2.23	0.54
1:B:264:ARG:NH2	1:B:331:GLU:O	2.40	0.54
1:C:384:HIS:O	1:C:385:ARG:HB2	2.08	0.54
1:B:411:ILE:HD12	1:B:440:THR:HA	1.89	0.54
1:C:273:LEU:HD11	1:C:283:MET:HB2	1.88	0.54
1:D:283:MET:CE	1:D:291:ARG:HH21	2.21	0.54
1:D:375:ALA:O	1:D:379:ARG:HG3	2.08	0.53
2:A:1:IHH:N9	2:A:1:IHH:C17	2.70	0.53
1:D:283:MET:HE1	1:D:291:ARG:NH2	2.23	0.53
1:A:331:GLU:OE2	1:A:331:GLU:HA	2.07	0.53
1:C:354:MET:HB2	3:C:129:HOH:O	2.09	0.53
2:A:1:IHH:H28	2:A:1:IHH:C16	2.34	0.53
1:A:502:ASP:HB3	1:A:505:GLU:HG2	1.90	0.52
1:C:258:ASP:OD1	1:C:260:TRP:HB2	2.09	0.52
1:C:383:VAL:CG2	1:C:411:ILE:HG23	2.39	0.52
2:D:1:IHH:H16	2:D:1:IHH:H28	1.91	0.52
1:D:393:LEU:HD11	2:D:1:IHH:H28	1.92	0.52
1:A:318:ARG:HG2	1:A:318:ARG:HH11	1.75	0.52
1:B:318:ARG:HB2	1:B:318:ARG:HH11	1.74	0.52
2:B:1:IHH:H28	2:B:1:IHH:H16	1.91	0.52
1:D:310:GLU:O	1:D:314:MET:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:270:GLU:HG3	1:D:284:GLY:HA2	1.92	0.52
1:C:268:ARG:O	1:C:268:ARG:HG3	2.10	0.51
1:C:260:TRP:NE1	1:C:315:LYS:HB2	2.26	0.51
1:A:332:GLU:HA	1:A:333:PRO:C	2.29	0.51
2:B:1:IHH:H23	2:B:1:IHH:C17	2.35	0.51
1:C:362:GLN:O	1:C:366:MET:HG3	2.10	0.51
1:C:409:ARG:NH2	1:C:439:PHE:HB2	2.26	0.51
1:D:261:GLU:HA	1:D:328:VAL:O	2.09	0.51
1:C:292:VAL:HG23	1:C:294:ILE:HD11	1.93	0.51
1:A:354:MET:CE	1:C:435:LEU:HB3	2.41	0.50
1:C:283:MET:HG3	1:C:340:TYR:CE1	2.45	0.50
1:D:424:PHE:HB3	1:D:425:PRO:HD2	1.92	0.50
1:B:310:GLU:HG2	1:B:314:MET:CE	2.41	0.50
1:B:269:LEU:HD23	1:B:294:ILE:CD1	2.40	0.50
1:C:493:ASP:O	1:C:497:GLN:HG3	2.10	0.50
1:B:269:LEU:HD22	1:B:282:TRP:HB2	1.92	0.50
1:B:304:PRO:O	1:B:308:LEU:HG	2.11	0.50
1:A:269:LEU:HD23	1:A:269:LEU:N	2.26	0.50
1:C:385:ARG:HG2	1:C:439:PHE:CD2	2.47	0.50
1:A:323:VAL:HG21	1:A:393:LEU:HD12	1.94	0.50
1:B:261:GLU:OE2	1:B:329:VAL:HA	2.12	0.50
1:D:279:GLY:HA3	1:D:296:THR:O	2.12	0.50
2:D:1:IHH:N9	2:D:1:IHH:C17	2.68	0.50
1:A:308:LEU:O	1:A:312:GLN:HG2	2.12	0.50
1:B:485:PRO:O	1:B:486:GLU:HB2	2.13	0.49
1:B:383:VAL:CG2	1:B:411:ILE:HB	2.43	0.49
1:D:323:VAL:HG21	1:D:393:LEU:HD12	1.94	0.49
1:A:276:GLY:HA3	1:A:279:GLY:O	2.13	0.49
1:A:309:GLN:HA	1:A:312:GLN:HG2	1.95	0.49
1:C:500:ARG:HD3	1:C:505:GLU:HG3	1.95	0.49
1:D:382:TYR:CE1	1:D:410:LEU:HD13	2.48	0.48
1:D:485:PRO:O	1:D:486:GLU:HB2	2.13	0.48
1:A:304:PRO:HG2	1:A:305:GLU:H	1.77	0.48
1:C:360:LEU:HD22	1:C:364:VAL:HG23	1.96	0.48
1:C:470:GLU:O	1:C:474:GLN:HG2	2.14	0.48
1:A:384:HIS:O	1:A:385:ARG:HB2	2.13	0.48
1:C:310:GLU:O	1:C:314:MET:HG3	2.13	0.48
1:D:466:MET:HE2	1:D:471:VAL:HG22	1.95	0.48
1:A:267:LEU:HD21	1:A:337:VAL:HG21	1.95	0.48
1:D:361:PRO:HA	1:D:520:PHE:CE2	2.48	0.48
1:C:271:VAL:HG12	1:C:272:LYS:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:310:GLU:HB3	1:D:406:GLY:HA2	1.95	0.47
1:A:297:LEU:HD23	1:A:297:LEU:C	2.35	0.47
1:A:362:GLN:O	1:A:366:MET:HG3	2.14	0.47
1:B:329:VAL:HB	1:B:335:TYR:HB2	1.96	0.47
1:D:470:GLU:O	1:D:474:GLN:HB2	2.14	0.47
1:B:500:ARG:HD3	1:B:505:GLU:HB3	1.97	0.47
1:C:310:GLU:HG3	1:C:406:GLY:N	2.30	0.47
1:D:473:ASP:O	1:D:477:ARG:HG3	2.14	0.47
1:C:304:PRO:O	1:C:307:PHE:HB3	2.15	0.47
1:C:407:LEU:HD13	1:C:424:PHE:CE2	2.50	0.47
1:A:269:LEU:HD22	1:A:294:ILE:HD12	1.97	0.46
1:B:405:PHE:HB3	1:B:408:ALA:HB3	1.97	0.46
1:C:278:PHE:O	1:C:297:LEU:HG	2.16	0.46
1:C:495:MET:O	1:C:498:CYS:HB2	2.16	0.46
1:C:269:LEU:N	1:C:269:LEU:HD23	2.29	0.46
1:C:272:LYS:O	1:C:273:LEU:HD23	2.15	0.46
1:C:424:PHE:HD1	1:C:424:PHE:H	1.61	0.46
1:A:354:MET:HE1	1:C:435:LEU:CB	2.44	0.46
1:A:463:TYR:N	1:A:464:PRO:HD3	2.30	0.45
1:B:407:LEU:HD13	1:B:424:PHE:CE1	2.51	0.45
1:B:491:LEU:O	1:B:495:MET:HG3	2.16	0.45
1:C:309:GLN:HA	1:C:312:GLN:HG2	1.98	0.45
1:A:336:ILE:HG23	1:A:336:ILE:O	2.15	0.45
1:C:293:ALA:C	1:C:294:ILE:HD12	2.37	0.45
1:C:320:GLU:O	1:C:401:LYS:HE2	2.16	0.45
1:A:489:GLU:HG2	3:A:181:HOH:O	2.16	0.45
1:A:356:LYS:HE3	1:A:357:TYR:CZ	2.52	0.45
1:D:451:LEU:O	1:D:451:LEU:HD23	2.17	0.45
1:A:424:PHE:O	1:A:426:ILE:N	2.49	0.45
1:C:361:PRO:HA	1:C:520:PHE:CE2	2.52	0.45
1:A:388:ARG:NH1	1:A:390:ALA:HB3	2.32	0.45
1:B:359:ARG:HA	1:B:359:ARG:HD3	1.82	0.45
1:D:283:MET:HG3	1:D:340:TYR:CZ	2.52	0.45
1:A:381:ASN:O	1:A:411:ILE:HG22	2.17	0.45
1:B:329:VAL:O	1:B:334:ILE:HG23	2.17	0.45
1:C:494:LEU:HD22	1:C:515:PHE:CE1	2.52	0.44
1:A:262:ILE:HB	1:A:263:PRO:HD2	1.99	0.44
1:C:262:ILE:HG12	1:C:263:PRO:HD2	1.98	0.44
1:B:393:LEU:HD11	2:B:1:IHH:C28	2.45	0.44
1:C:267:LEU:HD13	1:C:337:VAL:HG21	1.98	0.44
1:C:383:VAL:HG12	1:C:385:ARG:HG3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:297:LEU:O	1:D:299:PRO:HD3	2.18	0.44
1:A:286:TRP:HB3	1:A:290:THR:OG1	2.18	0.44
2:B:1:IHH:H16	2:B:1:IHH:C29	2.48	0.44
2:C:1:IHH:N9	2:C:1:IHH:H23	2.33	0.44
1:D:297:LEU:HD22	1:D:307:PHE:CB	2.48	0.44
1:B:286:TRP:HB3	1:B:290:THR:HG1	1.82	0.44
1:D:476:GLU:OE2	1:D:476:GLU:HA	2.16	0.44
1:A:278:PHE:CD1	1:A:278:PHE:N	2.86	0.44
1:A:310:GLU:HG3	1:A:405:PHE:CA	2.47	0.44
1:B:262:ILE:HB	1:B:263:PRO:HD2	2.00	0.43
1:B:297:LEU:HD22	1:B:307:PHE:CG	2.53	0.43
1:A:354:MET:HE2	1:C:436:TYR:CE1	2.53	0.43
1:C:332:GLU:OE1	1:C:334:ILE:HD12	2.17	0.43
1:C:347:LEU:O	1:C:351:LYS:HG2	2.17	0.43
1:A:477:ARG:HG3	1:A:477:ARG:HH11	1.83	0.43
1:A:529:PRO:HB3	3:A:82:HOH:O	2.19	0.43
1:B:264:ARG:HH12	1:B:333:PRO:HG2	1.83	0.43
1:C:498:CYS:O	1:C:506:ARG:HG2	2.18	0.43
1:A:500:ARG:HB3	1:A:505:GLU:HG3	2.00	0.43
1:D:262:ILE:HB	1:D:263:PRO:HD2	2.00	0.43
1:A:311:ALA:HB2	1:A:336:ILE:HD12	2.01	0.43
1:C:311:ALA:O	1:C:315:LYS:HG2	2.18	0.43
1:C:466:MET:CE	1:C:474:GLN:HG3	2.49	0.43
1:D:326:TYR:CE2	1:D:339:GLU:HA	2.54	0.43
2:D:1:IHH:H23	2:D:1:IHH:N9	2.34	0.43
1:A:502:ASP:O	1:A:505:GLU:HG2	2.19	0.43
1:A:360:LEU:HD13	1:A:520:PHE:HZ	1.84	0.43
1:B:310:GLU:HG2	1:B:314:MET:HE2	2.00	0.43
1:B:466:MET:HE3	1:B:471:VAL:HA	1.96	0.43
1:C:264:ARG:HH11	1:C:333:PRO:HD2	1.84	0.42
1:D:407:LEU:HD13	1:D:424:PHE:CE2	2.54	0.42
1:D:347:LEU:O	1:D:351:LYS:HG2	2.19	0.42
2:D:1:IHH:H16	2:D:1:IHH:C29	2.50	0.42
1:B:277:CYS:O	1:B:279:GLY:N	2.52	0.42
1:B:322:LEU:HG	1:B:405:PHE:HZ	1.85	0.42
1:B:384:HIS:O	1:B:385:ARG:HB2	2.19	0.42
1:A:310:GLU:HG3	1:A:405:PHE:HB2	2.01	0.42
1:A:332:GLU:HA	1:A:334:ILE:N	2.35	0.42
1:A:500:ARG:CB	1:A:505:GLU:HG3	2.50	0.42
1:D:502:ASP:HB3	1:D:505:GLU:HG2	2.02	0.42
1:A:477:ARG:O	1:A:477:ARG:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:GLN:NE2	1:A:312:GLN:HG3	2.35	0.42
1:D:334:ILE:HG22	1:D:335:TYR:N	2.35	0.42
1:D:498:CYS:O	1:D:506:ARG:HG2	2.20	0.42
1:A:273:LEU:CD1	1:A:283:MET:HB2	2.49	0.42
1:B:361:PRO:HA	1:B:520:PHE:CE2	2.55	0.42
1:C:279:GLY:HA3	1:C:296:THR:O	2.19	0.42
1:D:383:VAL:O	1:D:408:ALA:HA	2.20	0.42
1:A:308:LEU:HD21	1:A:334:ILE:CG2	2.50	0.41
2:A:1:IHH:N9	2:A:1:IHH:H23	2.34	0.41
1:B:446:TRP:CD1	1:B:446:TRP:C	2.94	0.41
1:B:275:GLN:NE2	1:B:275:GLN:HA	2.35	0.41
1:B:407:LEU:HD13	1:B:424:PHE:HE1	1.85	0.41
1:B:343:LYS:HB2	1:B:394:VAL:HB	2.03	0.41
1:B:349:PHE:CZ	1:B:354:MET:HG3	2.56	0.41
1:C:266:SER:HB2	1:C:287:ASN:OD1	2.21	0.41
2:C:1:IHH:N9	2:C:1:IHH:C23	2.83	0.41
1:A:318:ARG:HH11	1:A:318:ARG:CG	2.33	0.41
1:B:281:VAL:HG22	1:B:295:LYS:HG2	2.02	0.41
1:D:260:TRP:NE1	1:D:315:LYS:HG2	2.35	0.41
1:B:411:ILE:HD11	1:B:440:THR:HG22	2.01	0.41
1:D:260:TRP:CD1	1:D:315:LYS:HE3	2.55	0.41
1:D:384:HIS:O	1:D:385:ARG:HB2	2.20	0.41
1:D:491:LEU:HD23	1:D:491:LEU:HA	1.94	0.41
1:D:466:MET:HE3	1:D:471:VAL:HA	2.03	0.41
1:A:332:GLU:HA	1:A:334:ILE:HD12	2.03	0.41
1:A:411:ILE:HG23	1:A:411:ILE:O	2.21	0.41
1:C:255:LEU:HD11	1:C:331:GLU:OE2	2.20	0.41
1:A:355:GLY:O	1:A:458:LYS:HE2	2.20	0.41
1:C:272:LYS:HE2	1:C:275:GLN:OE1	2.20	0.41
1:C:343:LYS:HE3	1:C:396:GLU:OE1	2.21	0.41
1:D:351:LYS:HE3	3:D:176:HOH:O	2.20	0.41
1:C:309:GLN:HA	1:C:312:GLN:CG	2.51	0.40
1:D:531:GLU:HA	1:D:531:GLU:OE2	2.22	0.40
1:B:275:GLN:CA	1:B:275:GLN:HE21	2.35	0.40
1:C:368:ALA:HB1	1:C:517:GLU:HG2	2.03	0.40
1:D:363:LEU:HD23	1:D:366:MET:CE	2.51	0.40
1:B:283:MET:HE2	1:B:291:ARG:HD3	2.04	0.40
1:A:334:ILE:H	1:A:334:ILE:CD1	2.32	0.40
1:B:528:GLN:HA	1:B:529:PRO:HD3	1.98	0.40
1:B:269:LEU:CD2	1:B:294:ILE:HD12	2.45	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	258/286 (90%)	240 (93%)	15 (6%)	3 (1%)	13	11
1	B	261/286 (91%)	246 (94%)	14 (5%)	1 (0%)	34	38
1	C	259/286 (91%)	245 (95%)	11 (4%)	3 (1%)	13	11
1	D	263/286 (92%)	254 (97%)	9 (3%)	0	100	100
All	All	1041/1144 (91%)	985 (95%)	49 (5%)	7 (1%)	22	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	404	ASP
1	A	336	ILE
1	A	425	PRO
1	A	404	ASP
1	C	411	ILE
1	C	425	PRO
1	B	425	PRO

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	227/245 (93%)	217 (96%)	10 (4%)	28	34
1	B	228/245 (93%)	220 (96%)	8 (4%)	36	44
1	C	228/245 (93%)	217 (95%)	11 (5%)	25	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	229/245 (94%)	218 (95%)	11 (5%)	25	30
All	All	912/980 (93%)	872 (96%)	40 (4%)	28	34

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	264	ARG
1	A	269	LEU
1	A	271	VAL
1	A	318	ARG
1	A	351	LYS
1	A	356	LYS
1	A	360	LEU
1	A	426	ILE
1	A	491	LEU
1	A	526	GLN
1	B	255	LEU
1	B	283	MET
1	B	351	LYS
1	B	354	MET
1	B	360	LEU
1	B	438	ARG
1	B	491	LEU
1	B	505	GLU
1	C	267	LEU
1	C	275	GLN
1	C	283	MET
1	C	307	PHE
1	C	318	ARG
1	C	334	ILE
1	C	351	LYS
1	C	360	LEU
1	C	396	GLU
1	C	491	LEU
1	C	526	GLN
1	D	277	CYS
1	D	283	MET
1	D	343	LYS
1	D	351	LYS
1	D	360	LEU
1	D	388	ARG
1	D	438	ARG

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Mol	Chain	Res	Type
1	D	469	ARG
1	D	491	LEU
1	D	502	ASP
1	D	504	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	309	GLN
1	C	309	GLN
1	C	381	ASN

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)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	IHH	B	1	-	32,33,33	2.03	6 (18%)	42,46,46	2.04	11 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	IHH	D	1	-	32,33,33	2.03	6 (18%)	42,46,46	2.06	12 (28%)
2	IHH	C	1	-	32,33,33	2.00	6 (18%)	42,46,46	2.10	12 (28%)
2	IHH	A	1	-	32,33,33	2.00	6 (18%)	42,46,46	2.09	13 (30%)

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	IHH	B	1	-	-	0/12/17/17	0/5/5/5
2	IHH	D	1	-	-	0/12/17/17	0/5/5/5
2	IHH	C	1	-	-	0/12/17/17	0/5/5/5
2	IHH	A	1	-	-	0/12/17/17	0/5/5/5

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	IHH	N26-N25	-7.25	1.23	1.37
2	D	1	IHH	N26-N25	-7.24	1.23	1.37
2	A	1	IHH	N26-N25	-7.18	1.23	1.37
2	C	1	IHH	N26-N25	-7.12	1.23	1.37
2	B	1	IHH	C10-C5	-4.48	1.39	1.44
2	D	1	IHH	C10-C5	-4.46	1.39	1.44
2	C	1	IHH	C10-C5	-4.22	1.39	1.44
2	A	1	IHH	C10-C5	-4.20	1.39	1.44
2	B	1	IHH	C12-N11	-3.31	1.33	1.40
2	C	1	IHH	C12-N11	-3.31	1.33	1.40
2	A	1	IHH	C12-N11	-3.29	1.33	1.40
2	D	1	IHH	C23-C24	-3.29	1.33	1.39
2	A	1	IHH	C23-C24	-3.28	1.33	1.39
2	C	1	IHH	C23-C24	-3.28	1.33	1.39
2	B	1	IHH	C23-C24	-3.27	1.33	1.39
2	D	1	IHH	C12-N11	-3.22	1.33	1.40
2	D	1	IHH	C22-N21	-2.98	1.33	1.38
2	B	1	IHH	C22-N21	-2.88	1.33	1.38
2	A	1	IHH	C22-N21	-2.74	1.33	1.38
2	C	1	IHH	C22-N21	-2.73	1.33	1.38
2	C	1	IHH	C6-N7	-2.25	1.33	1.37
2	D	1	IHH	C6-N7	-2.14	1.34	1.37
2	A	1	IHH	C6-N7	-2.14	1.34	1.37
2	B	1	IHH	C6-N7	-2.08	1.34	1.37

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1	IHH	C10-C5-C6	5.79	119.53	115.88
2	B	1	IHH	C10-C5-C6	5.78	119.52	115.88
2	D	1	IHH	N21-C10-N9	5.65	123.96	118.66
2	C	1	IHH	C10-C5-C6	5.60	119.40	115.88
2	D	1	IHH	C10-C5-C6	5.47	119.32	115.88
2	B	1	IHH	N21-C10-N9	5.40	123.73	118.66
2	A	1	IHH	N21-C10-N9	5.03	123.38	118.66
2	C	1	IHH	N21-C10-N9	4.76	123.12	118.66
2	C	1	IHH	C28-C27-C24	-4.62	114.91	120.33
2	A	1	IHH	C28-C27-C24	-4.46	115.09	120.33
2	B	1	IHH	C28-C27-C24	-3.89	115.76	120.33
2	C	1	IHH	C29-C27-C24	-3.80	115.87	120.33
2	D	1	IHH	C28-C27-C24	-3.76	115.92	120.33
2	B	1	IHH	C5-C6-N7	-3.71	118.88	122.81
2	A	1	IHH	C29-C27-C24	-3.69	116.00	120.33
2	C	1	IHH	C5-C6-N7	-3.61	118.98	122.81
2	D	1	IHH	C5-C6-N7	-3.57	119.03	122.81
2	D	1	IHH	C29-C27-C24	-3.52	116.20	120.33
2	A	1	IHH	C5-C6-N7	-3.50	119.10	122.81
2	D	1	IHH	N7-C8-N9	-3.43	120.81	126.23
2	B	1	IHH	C29-C27-C24	-3.40	116.33	120.33
2	D	1	IHH	C8-N7-C6	3.36	121.07	115.60
2	C	1	IHH	C8-N7-C6	3.35	121.07	115.60
2	C	1	IHH	C24-C23-C22	-3.30	103.16	106.07
2	A	1	IHH	N7-C8-N9	-3.29	121.02	126.23
2	C	1	IHH	N7-C8-N9	-3.27	121.06	126.23
2	B	1	IHH	C8-N7-C6	3.27	120.93	115.60
2	A	1	IHH	C8-N7-C6	3.22	120.85	115.60
2	D	1	IHH	C24-C23-C22	-3.21	103.23	106.07
2	B	1	IHH	N7-C8-N9	-3.20	121.17	126.23
2	B	1	IHH	C24-C23-C22	-3.18	103.26	106.07
2	A	1	IHH	C24-C23-C22	-2.97	103.45	106.07
2	A	1	IHH	C27-C24-N25	2.86	124.45	120.23
2	C	1	IHH	C27-C24-N25	2.74	124.27	120.23
2	D	1	IHH	C5-C10-N21	-2.49	116.99	119.73
2	C	1	IHH	N11-C8-N9	2.43	125.17	116.92
2	D	1	IHH	C27-C24-N25	2.38	123.74	120.23
2	C	1	IHH	C23-C24-C27	-2.32	125.69	129.25
2	A	1	IHH	C23-C24-C27	-2.28	125.75	129.25
2	A	1	IHH	N11-C8-N9	2.27	124.62	116.92
2	B	1	IHH	C5-C10-N21	-2.23	117.28	119.73
2	D	1	IHH	C4-C5-C10	-2.20	121.04	124.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1	IHH	C4-C5-C10	-2.11	121.17	124.06
2	B	1	IHH	C27-C24-N25	2.09	123.31	120.23
2	A	1	IHH	C5-C10-N9	-2.08	118.57	121.98
2	C	1	IHH	C3-C6-C5	2.06	121.49	119.13
2	D	1	IHH	N11-C8-N9	2.01	123.74	116.92
2	A	1	IHH	C4-C5-C10	-2.01	121.31	124.06

There are no chirality outliers.

There are no torsion outliers.

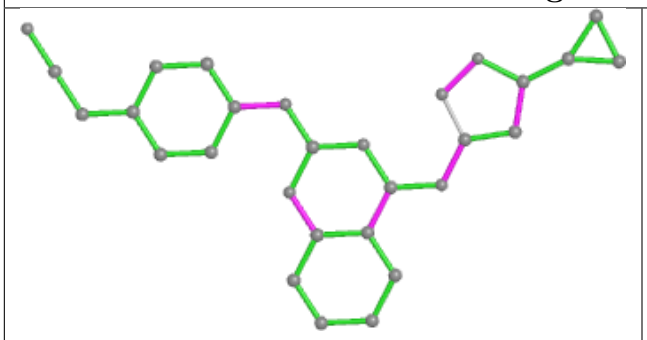
There are no ring outliers.

4 monomers are involved in 28 short contacts:

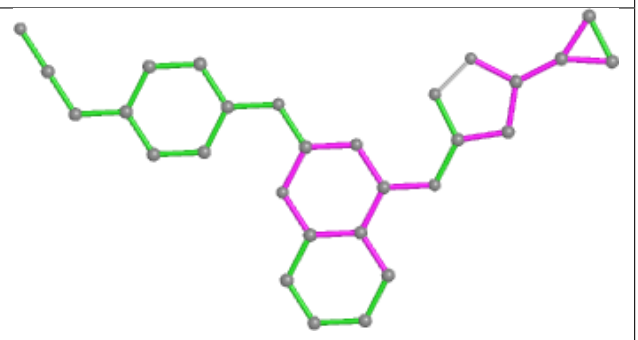
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	IHH	9	0
2	D	1	IHH	8	0
2	C	1	IHH	5	0
2	A	1	IHH	6	0

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.

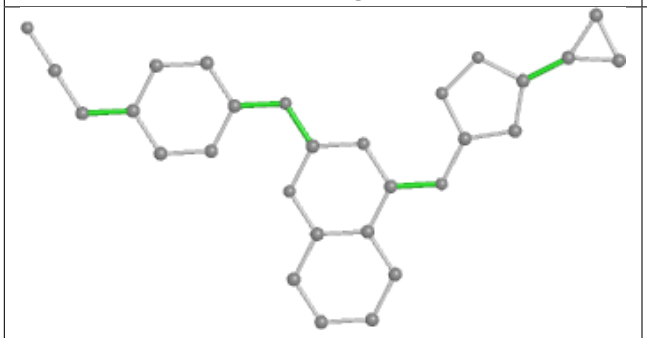
Ligand IHH B 1



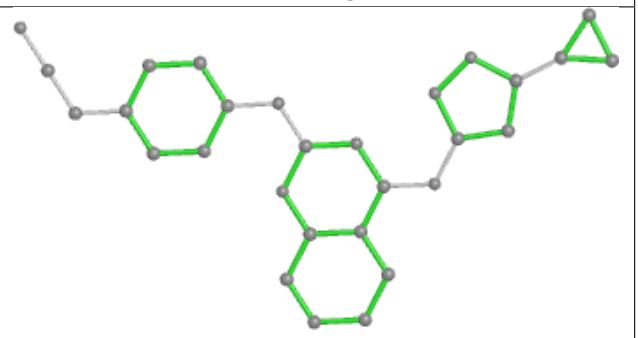
Bond lengths



Bond angles

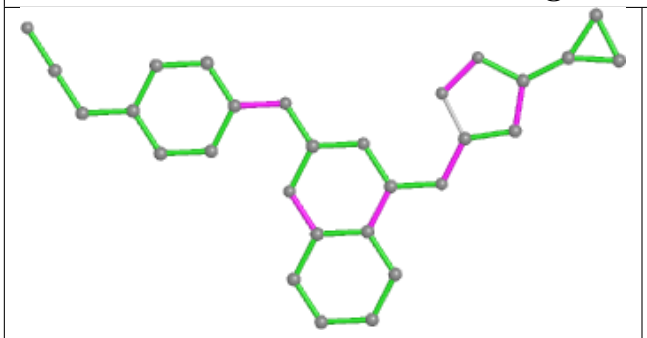


Torsions

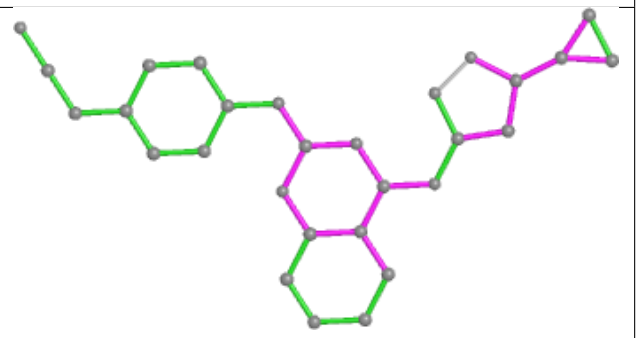


Rings

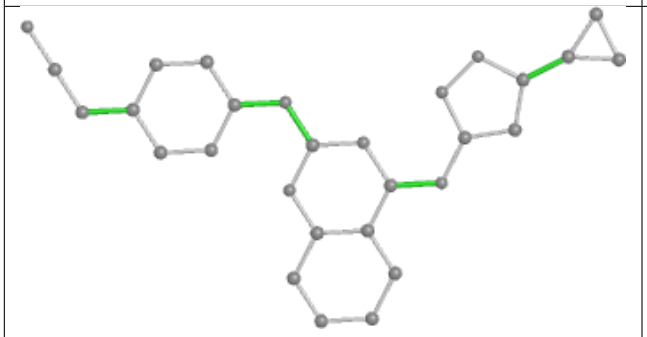
Ligand IHH D 1



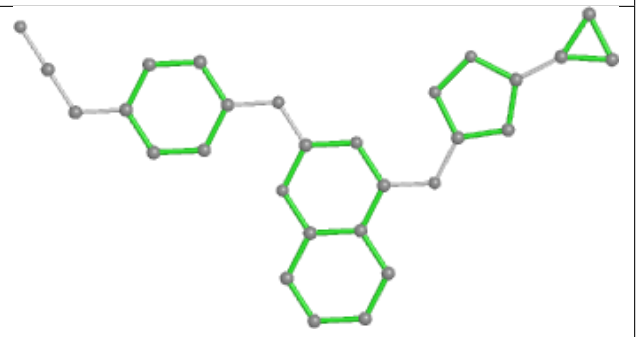
Bond lengths



Bond angles

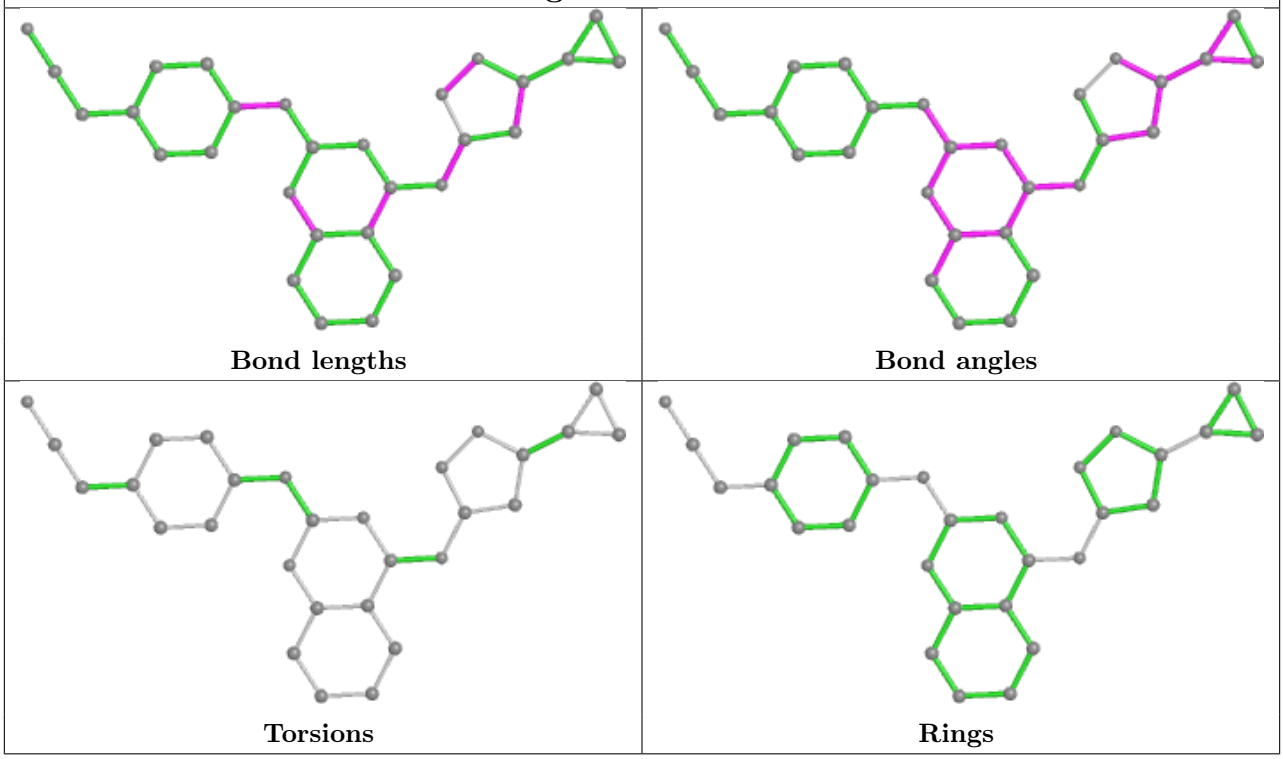


Torsions

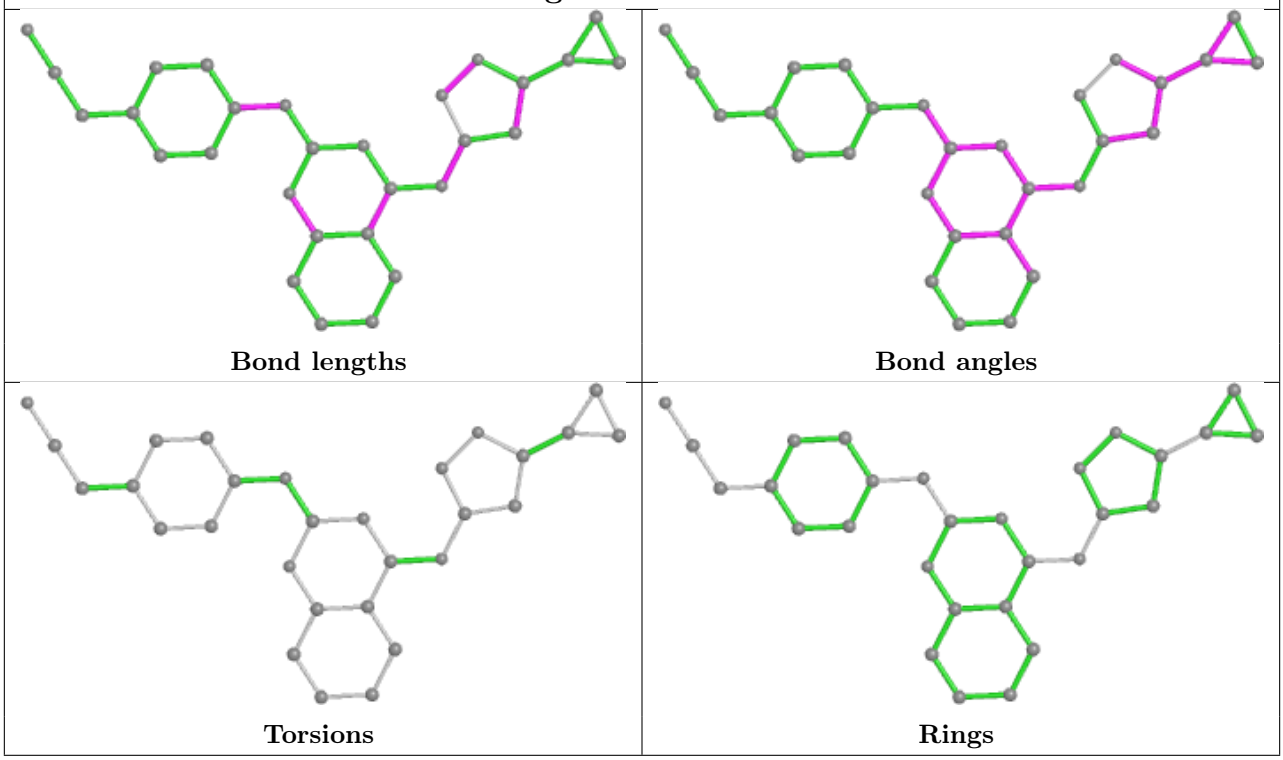


Rings

Ligand IHH C 1



Ligand IHH A 1



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	264/286 (92%)	0.67	38 (14%)	2	4	12, 30, 76, 98	0
1	B	267/286 (93%)	0.53	25 (9%)	8	13	11, 31, 75, 86	0
1	C	265/286 (92%)	0.70	34 (12%)	3	6	12, 31, 78, 90	0
1	D	267/286 (93%)	0.71	30 (11%)	5	8	12, 32, 75, 84	0
All	All	1063/1144 (92%)	0.65	127 (11%)	4	7	11, 31, 77, 98	0

All (127) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	277	CYS	11.6
1	C	299	PRO	10.9
1	D	412	GLU	10.8
1	D	278	PHE	9.2
1	B	300	GLY	8.3
1	D	300	GLY	8.0
1	B	255	LEU	7.8
1	C	298	LYS	7.2
1	B	301	THR	6.5
1	D	301	THR	6.3
1	B	256	ALA	6.1
1	C	411	ILE	5.9
1	D	299	PRO	5.8
1	A	304	PRO	5.4
1	A	271	VAL	5.3
1	A	333	PRO	5.2
1	A	411	ILE	5.2
1	B	299	PRO	5.0
1	D	354	MET	4.9
1	D	411	ILE	4.8
1	A	412	GLU	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	269	LEU	4.5
1	C	267	LEU	4.4
1	C	271	VAL	4.4
1	B	277	CYS	4.4
1	A	278	PHE	4.3
1	D	472	LEU	4.3
1	B	276	GLY	4.2
1	D	279	GLY	4.2
1	D	408	ALA	4.0
1	C	477	ARG	4.0
1	C	331	GLU	3.9
1	C	424	PHE	3.9
1	C	277	CYS	3.9
1	D	424	PHE	3.9
1	A	477	ARG	3.8
1	C	332	GLU	3.8
1	A	332	GLU	3.7
1	C	297	LEU	3.6
1	B	469	ARG	3.6
1	C	273	LEU	3.6
1	C	278	PHE	3.5
1	B	423	LYS	3.5
1	C	333	PRO	3.4
1	B	409	ARG	3.4
1	B	424	PHE	3.4
1	D	423	LYS	3.4
1	A	276	GLY	3.3
1	C	412	GLU	3.3
1	A	268	ARG	3.3
1	D	473	ASP	3.3
1	A	288	GLY	3.3
1	A	334	ILE	3.3
1	C	409	ARG	3.3
1	C	268	ARG	3.3
1	A	277	CYS	3.3
1	C	307	PHE	3.2
1	A	270	GLU	3.1
1	A	267	LEU	3.1
1	A	410	LEU	3.1
1	C	255	LEU	3.1
1	C	308	LEU	3.1
1	B	411	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	289	THR	3.1
1	B	298	LYS	3.1
1	A	305	GLU	3.1
1	C	408	ALA	3.1
1	D	302	MET	3.0
1	D	309	GLN	3.0
1	B	297	LEU	3.0
1	A	307	PHE	3.0
1	D	276	GLY	3.0
1	B	473	ASP	3.0
1	D	258	ASP	3.0
1	C	289	THR	3.0
1	A	423	LYS	3.0
1	C	275	GLN	2.9
1	C	276	GLY	2.9
1	C	334	ILE	2.9
1	B	309	GLN	2.8
1	B	302	MET	2.8
1	D	268	ARG	2.8
1	D	297	LEU	2.7
1	A	408	ALA	2.7
1	A	264	ARG	2.7
1	A	308	LEU	2.7
1	D	289	THR	2.7
1	A	331	GLU	2.6
1	A	335	TYR	2.6
1	D	306	ALA	2.6
1	A	409	ARG	2.6
1	B	477	ARG	2.6
1	A	289	THR	2.5
1	D	304	PRO	2.5
1	C	265	GLU	2.5
1	C	410	LEU	2.5
1	C	288	GLY	2.5
1	D	477	ARG	2.4
1	A	265	GLU	2.4
1	A	255	LEU	2.4
1	A	275	GLN	2.4
1	D	465	GLY	2.4
1	C	335	TYR	2.4
1	A	283	MET	2.4
1	A	311	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	288	GLY	2.4
1	A	297	LEU	2.4
1	C	264	ARG	2.3
1	B	410	LEU	2.3
1	C	282	TRP	2.3
1	A	303	SER	2.3
1	A	291	ARG	2.3
1	C	279	GLY	2.2
1	A	282	TRP	2.2
1	B	304	PRO	2.2
1	B	408	ALA	2.2
1	D	409	ARG	2.2
1	A	313	VAL	2.2
1	A	256	ALA	2.2
1	D	275	GLN	2.2
1	C	256	ALA	2.1
1	B	258	ASP	2.1
1	D	410	LEU	2.1
1	D	305	GLU	2.1
1	B	305	GLU	2.1
1	D	298	LYS	2.1
1	C	285	THR	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
2	IHH	A	1	29/29	0.90	0.21	30,33,54,55	0

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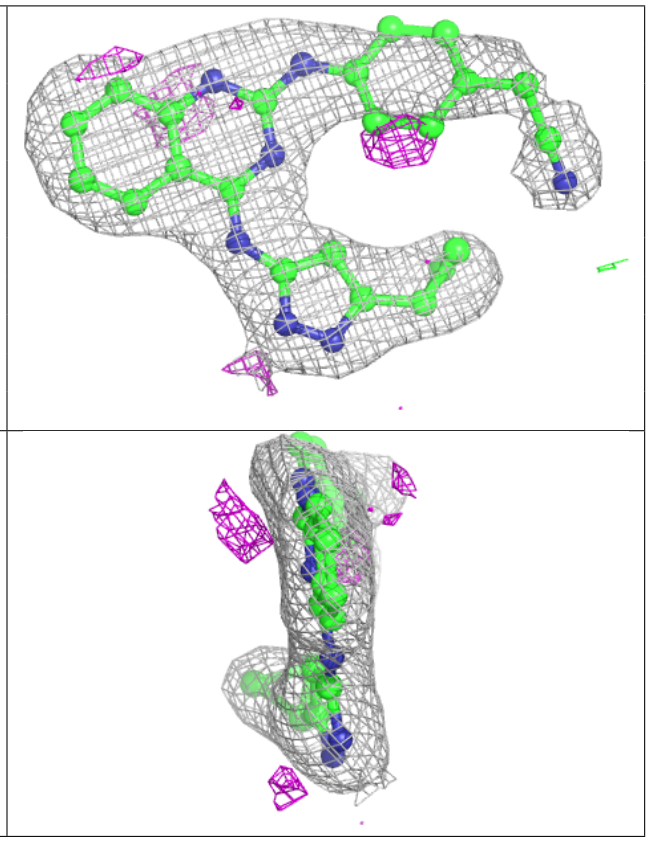
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	IHH	C	1	29/29	0.91	0.18	32,34,54,56	0
2	IHH	B	1	29/29	0.92	0.20	31,34,59,61	0
2	IHH	D	1	29/29	0.93	0.18	28,30,51,53	0

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

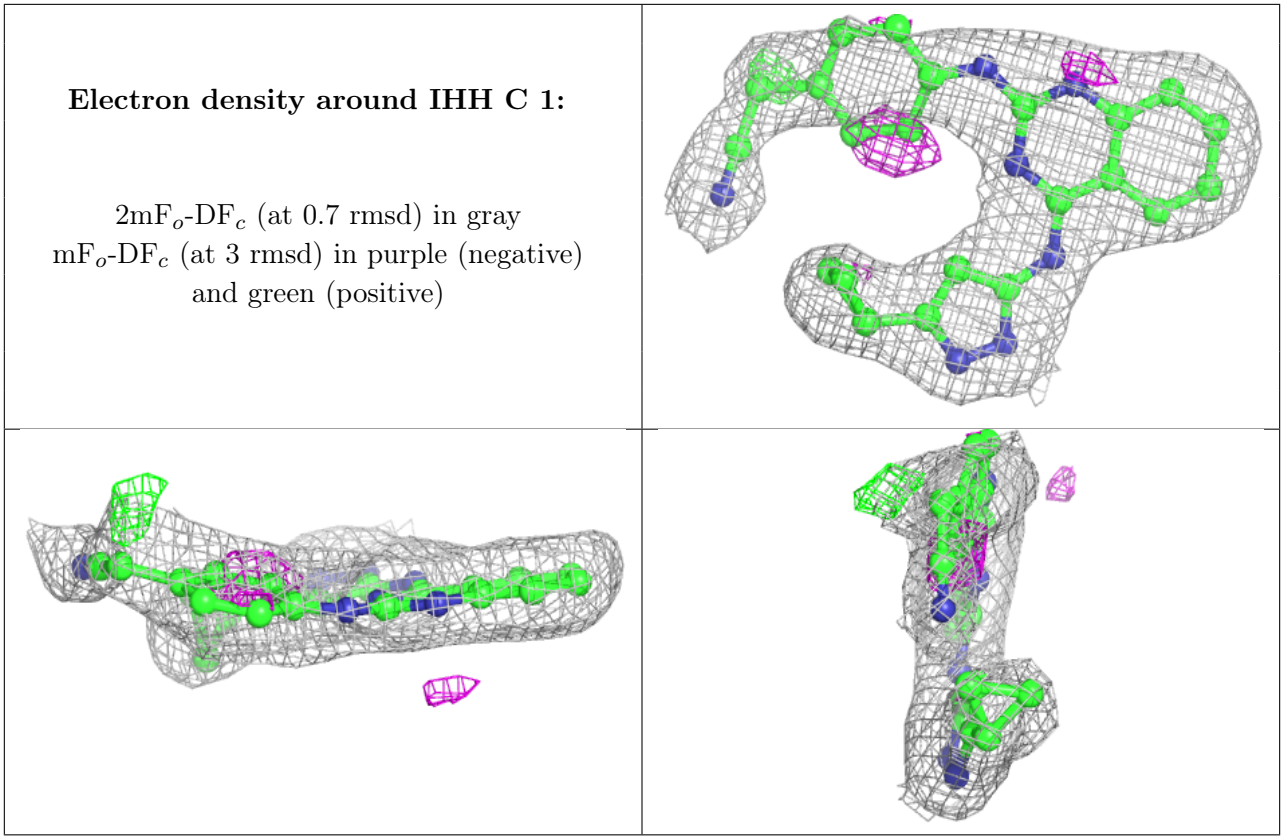
Electron density around IHH A 1:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



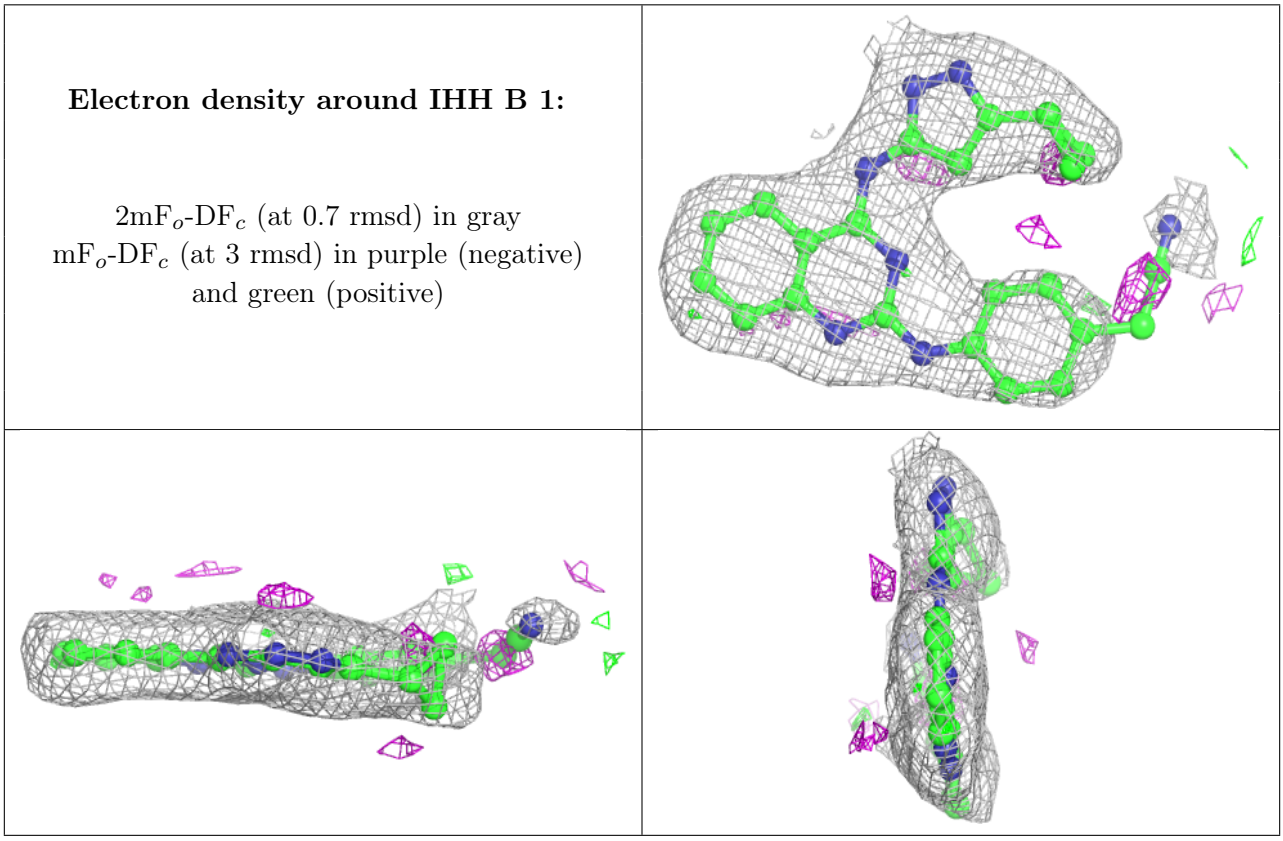
Electron density around IHH C 1:

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



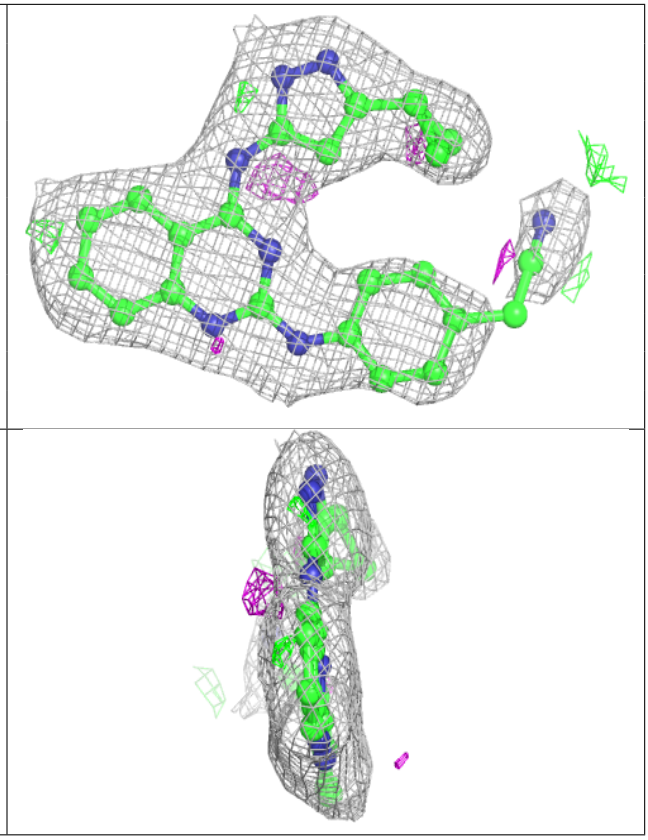
Electron density around IHH B 1:

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



Electron density around IHH D 1:

$2mF_o$-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.