



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:02 PM UTC

PDB ID : 8PW1 / pdb_00008pw1
Title : Structure of human UCHL1 in complex with CG341 inhibitor
Authors : Grethe, C.; Gersch, M.
Deposited on : 2023-07-19
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

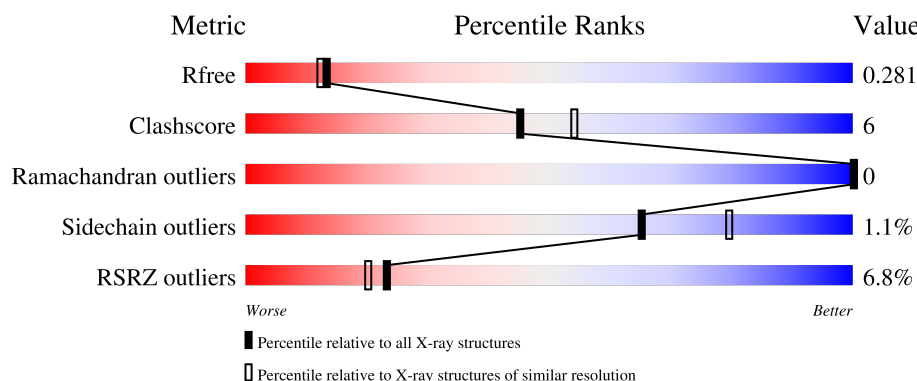
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	227	3% 92% 6% .
1	B	227	4% 86% 10% .
1	C	227	5% 85% 11% ..
1	D	227	4% 80% 15% 5%
1	E	227	2% 83% 13% .

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Mol	Chain	Length	Quality of chain
1	F	227	 2% 83% 12% .
1	G	227	 8% 82% 14% .
1	H	227	 4% 85% 11% .
1	I	227	 7% 78% 17% .
1	J	227	 26% 75% 21% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16561 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 Ubiquitin carboxyl-terminal hydrolase isozyme L1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1665	1051	290	313	11			
1	B	219	Total	C	N	O	S	0	0	0
			1609	1020	280	298	11			
1	C	220	Total	C	N	O	S	0	0	0
			1598	1013	279	295	11			
1	D	216	Total	C	N	O	S	0	0	0
			1602	1013	273	305	11			
1	E	219	Total	C	N	O	S	0	0	0
			1623	1025	284	302	12			
1	F	217	Total	C	N	O	S	0	0	0
			1614	1022	282	298	12			
1	G	219	Total	C	N	O	S	0	0	0
			1593	1009	279	294	11			
1	H	217	Total	C	N	O	S	0	0	0
			1578	997	274	296	11			
1	I	217	Total	C	N	O	S	0	0	0
			1590	1006	278	295	11			
1	J	219	Total	C	N	O	S	0	0	0
			1552	977	270	293	12			

There are 40 discrepancies between the modelled and reference sequences:

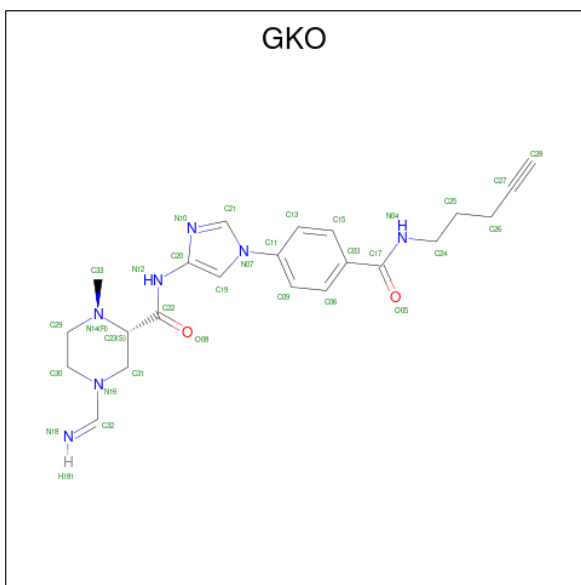
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P09936
A	-2	PRO	-	expression tag	UNP P09936
A	-1	GLY	-	expression tag	UNP P09936
A	0	SER	-	expression tag	UNP P09936
B	-3	GLY	-	expression tag	UNP P09936
B	-2	PRO	-	expression tag	UNP P09936
B	-1	GLY	-	expression tag	UNP P09936
B	0	SER	-	expression tag	UNP P09936
C	-3	GLY	-	expression tag	UNP P09936

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	PRO	-	expression tag	UNP P09936
C	-1	GLY	-	expression tag	UNP P09936
C	0	SER	-	expression tag	UNP P09936
D	-3	GLY	-	expression tag	UNP P09936
D	-2	PRO	-	expression tag	UNP P09936
D	-1	GLY	-	expression tag	UNP P09936
D	0	SER	-	expression tag	UNP P09936
E	-3	GLY	-	expression tag	UNP P09936
E	-2	PRO	-	expression tag	UNP P09936
E	-1	GLY	-	expression tag	UNP P09936
E	0	SER	-	expression tag	UNP P09936
F	-3	GLY	-	expression tag	UNP P09936
F	-2	PRO	-	expression tag	UNP P09936
F	-1	GLY	-	expression tag	UNP P09936
F	0	SER	-	expression tag	UNP P09936
G	-3	GLY	-	expression tag	UNP P09936
G	-2	PRO	-	expression tag	UNP P09936
G	-1	GLY	-	expression tag	UNP P09936
G	0	SER	-	expression tag	UNP P09936
H	-3	GLY	-	expression tag	UNP P09936
H	-2	PRO	-	expression tag	UNP P09936
H	-1	GLY	-	expression tag	UNP P09936
H	0	SER	-	expression tag	UNP P09936
I	-3	GLY	-	expression tag	UNP P09936
I	-2	PRO	-	expression tag	UNP P09936
I	-1	GLY	-	expression tag	UNP P09936
I	0	SER	-	expression tag	UNP P09936
J	-3	GLY	-	expression tag	UNP P09936
J	-2	PRO	-	expression tag	UNP P09936
J	-1	GLY	-	expression tag	UNP P09936
J	0	SER	-	expression tag	UNP P09936

- Molecule 2 is (2 {S})-4-(iminomethyl)-1-methyl- {N}-[1-[4-(pent-4-ynylcarbamoyl)phenyl]imidazol-4-yl]piperazine-2-carboxamide (CCD ID: GKO) (formula: C₂₂H₂₇N₇O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 31	C 22	N 7	O 2	0	0
2	B	1	Total 31	C 22	N 7	O 2	0	0
2	C	1	Total 31	C 22	N 7	O 2	0	0
2	D	1	Total 31	C 22	N 7	O 2	0	0
2	E	1	Total 31	C 22	N 7	O 2	0	0
2	F	1	Total 31	C 22	N 7	O 2	0	0
2	G	1	Total 31	C 22	N 7	O 2	0	0
2	H	1	Total 31	C 22	N 7	O 2	0	0
2	I	1	Total 31	C 22	N 7	O 2	0	0
2	J	1	Total 31	C 22	N 7	O 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	46	Total O 46 46	0	0
3	B	34	Total O 34 34	0	0

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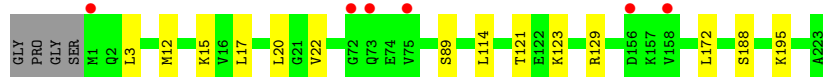
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	18	Total 18	O 18	0	0
3	D	16	Total 16	O 16	0	0
3	E	34	Total 34	O 34	0	0
3	F	21	Total 21	O 21	0	0
3	G	20	Total 20	O 20	0	0
3	H	14	Total 14	O 14	0	0
3	I	4	Total 4	O 4	0	0
3	J	20	Total 20	O 20	0	0

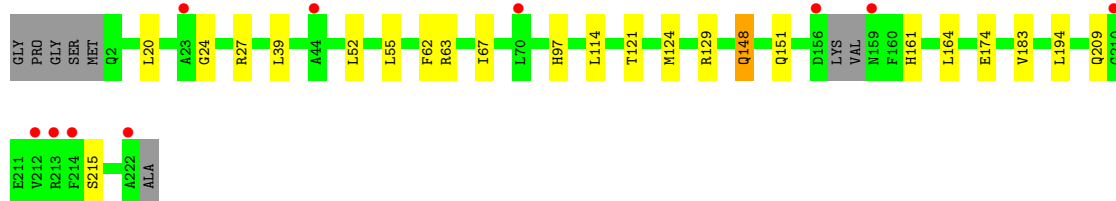
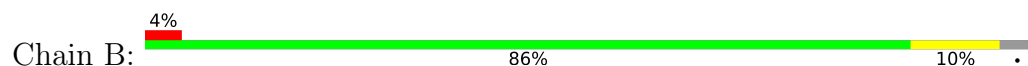
3 Residue-property plots [i](#)

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.

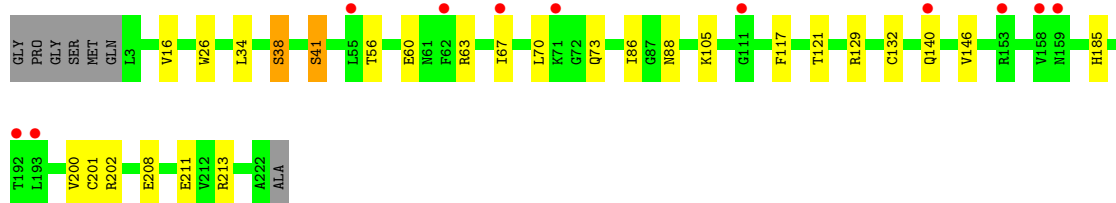
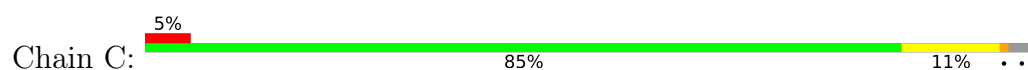
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1



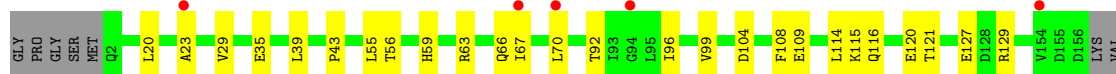
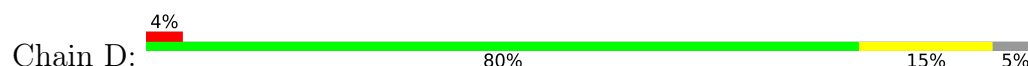
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1

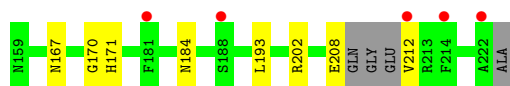


- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1

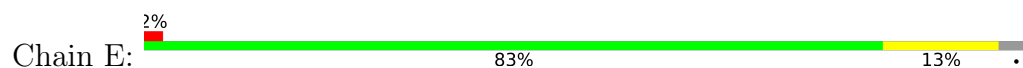


- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1

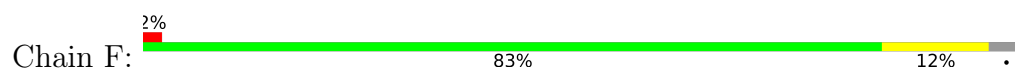




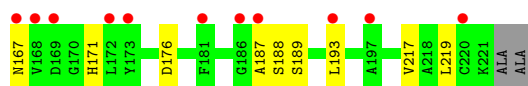
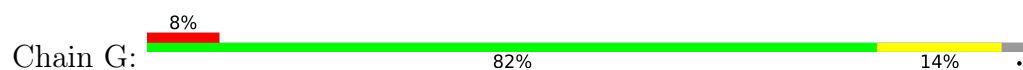
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1



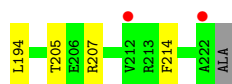
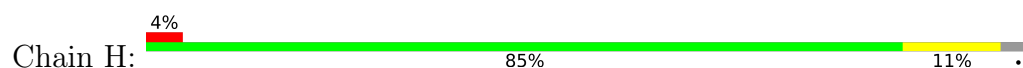
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1



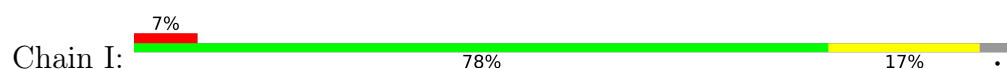
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1

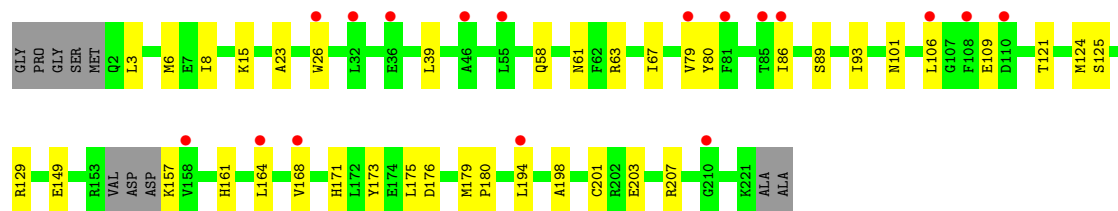


- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1

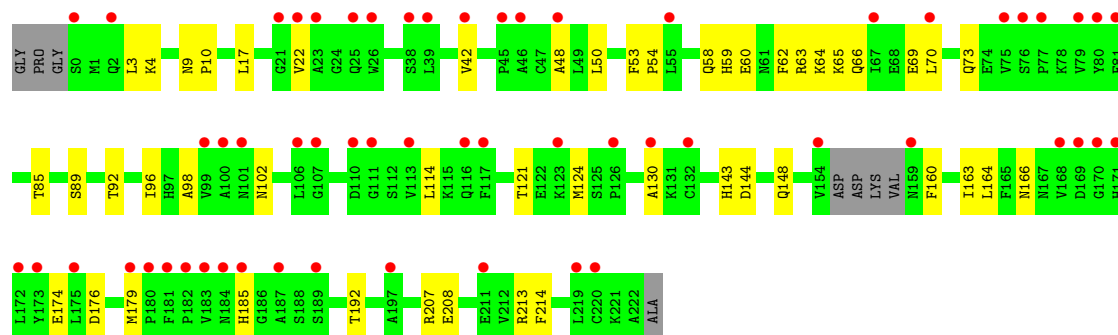
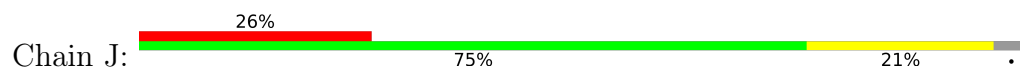


- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1





- Molecule 1: Ubiquitin carboxyl-terminal hydrolase isozyme L1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	69.63Å 77.75Å 126.47Å 94.09° 104.06° 113.80°	Depositor
Resolution (Å)	51.61 – 2.20 51.61 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.0 (51.61-2.20) 98.1 (51.61-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.227 , 0.271 (Not available) , 0.281	Depositor DCC
R_{free} test set	5770 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 76.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16561	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.12	0/1696	0.26	0/2296
1	B	0.10	0/1639	0.25	0/2222
1	C	0.20	0/1629	0.33	0/2212
1	D	0.09	0/1631	0.22	0/2211
1	E	0.10	0/1653	0.23	0/2241
1	F	0.10	0/1644	0.25	0/2226
1	G	0.10	0/1622	0.25	0/2201
1	H	0.11	0/1607	0.25	0/2184
1	I	0.15	0/1619	0.27	0/2197
1	J	0.14	0/1582	0.26	0/2154
All	All	0.12	0/16322	0.26	0/22144

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

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	1665	0	1588	8	0
1	B	1609	0	1518	12	0
1	C	1598	0	1483	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1602	0	1498	22	0
1	E	1623	0	1532	20	0
1	F	1614	0	1531	18	0
1	G	1593	0	1484	18	0
1	H	1578	0	1465	16	0
1	I	1590	0	1491	28	0
1	J	1552	0	1371	27	0
2	A	31	0	0	0	0
2	B	31	0	0	0	0
2	C	31	0	0	0	0
2	D	31	0	0	0	0
2	E	31	0	0	0	0
2	F	31	0	0	1	0
2	G	31	0	0	1	0
2	H	31	0	0	0	0
2	I	31	0	0	0	0
2	J	31	0	0	0	0
3	A	46	0	0	0	0
3	B	34	0	0	0	0
3	C	18	0	0	0	0
3	D	16	0	0	1	0
3	E	34	0	0	0	0
3	F	21	0	0	0	0
3	G	20	0	0	0	0
3	H	14	0	0	0	0
3	I	4	0	0	0	0
3	J	20	0	0	0	0
All	All	16561	0	14961	184	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 (184) 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:G:99:VAL:HG11	1:G:114:LEU:HD21	1.61	0.82
1:I:121:THR:HA	1:I:124:MET:HG3	1.66	0.76
1:D:99:VAL:HG21	1:D:114:LEU:HD21	1.67	0.75
1:H:121:THR:HB	1:H:129:ARG:HG2	1.70	0.73
1:J:4:LYS:HD2	1:J:148:GLN:HE22	1.54	0.71
1:D:202:ARG:NH2	3:D:401:HOH:O	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:80:TYR:H	1:I:101:ASN:HD21	1.42	0.68
1:I:23:ALA:HB2	1:I:109:GLU:HG2	1.74	0.67
1:A:121:THR:HB	1:A:129:ARG:HG2	1.76	0.67
1:D:23:ALA:HB2	1:D:109:GLU:HG3	1.77	0.67
1:D:171:HIS:HD2	1:D:184:ASN:HD21	1.43	0.67
1:D:56:THR:HG23	1:D:59:HIS:H	1.62	0.65
1:J:50:LEU:HG	1:J:164:LEU:HD13	1.77	0.64
1:G:85:THR:HG23	1:G:86:ILE:HG13	1.81	0.63
1:E:1:MET:SD	1:E:1:MET:N	2.65	0.61
1:F:34:LEU:O	1:F:202:ARG:NH1	2.33	0.61
1:E:172:LEU:HD12	1:E:185:HIS:HB2	1.81	0.61
1:J:85:THR:OG1	1:J:130:ALA:HB1	2.01	0.61
1:I:121:THR:HG22	1:I:124:MET:HE3	1.83	0.61
1:E:121:THR:HB	1:E:129:ARG:HG2	1.83	0.61
1:I:80:TYR:H	1:I:101:ASN:ND2	2.00	0.59
1:J:48:ALA:HB1	1:J:164:LEU:HD11	1.83	0.59
1:E:3:LEU:HG	1:E:86:ILE:HG21	1.85	0.59
1:C:34:LEU:O	1:C:202:ARG:NH1	2.36	0.58
1:H:139:ILE:HD12	1:H:139:ILE:H	1.68	0.58
1:F:5:PRO:HB2	1:F:152:CYS:HB3	1.84	0.58
1:F:3:LEU:HD21	1:F:140:GLN:HG2	1.86	0.57
1:J:144:ASP:O	1:J:148:GLN:HG2	2.04	0.57
1:I:3:LEU:HG	1:I:86:ILE:HG21	1.87	0.57
1:B:63:ARG:O	1:B:67:ILE:HD12	2.06	0.56
1:E:7:GLU:OE1	1:E:151:GLN:NE2	2.39	0.56
1:G:171:HIS:HA	1:G:187:ALA:HA	1.86	0.56
1:J:66:GLN:O	1:J:70:LEU:N	2.39	0.55
1:E:1:MET:HE1	1:E:134:GLU:HB3	1.88	0.55
1:I:179:MET:HG3	1:I:180:PRO:HD2	1.87	0.55
1:H:20:LEU:HD13	1:H:114:LEU:HD13	1.89	0.54
1:I:26:TRP:HB2	1:I:106:LEU:HD21	1.90	0.54
1:E:164:LEU:HD21	1:E:166:ASN:HD22	1.73	0.54
1:F:64:LYS:N	1:F:64:LYS:HD3	2.23	0.53
1:G:144:ASP:O	1:G:148:GLN:HG3	2.08	0.53
1:C:63:ARG:O	1:C:67:ILE:HG13	2.08	0.53
1:E:23:ALA:HB2	1:E:109:GLU:HG2	1.90	0.53
1:G:106:LEU:HD11	1:G:219:LEU:HD21	1.91	0.53
1:F:3:LEU:HD13	1:F:143:HIS:CD2	2.44	0.53
1:C:70:LEU:HD12	1:C:73:GLN:HB2	1.92	0.52
1:B:174:GLU:HB3	1:B:183:VAL:HB	1.92	0.52
1:F:164:LEU:HD21	1:F:166:ASN:ND2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:176:ASP:HB3	1:I:179:MET:HB2	1.91	0.52
1:J:174:GLU:OE1	1:J:185:HIS:NE2	2.39	0.52
1:B:55:LEU:HD13	1:B:161:HIS:HB2	1.92	0.51
1:F:96:ILE:HG23	1:F:114:LEU:HD11	1.93	0.51
1:E:29:VAL:HG21	1:E:43:PRO:HD2	1.92	0.51
1:I:6:MET:HG2	1:I:8:ILE:HD12	1.92	0.50
1:G:9:ASN:HB2	1:G:10:PRO:HD2	1.94	0.50
1:C:117:PHE:O	1:C:121:THR:OG1	2.23	0.50
1:E:73:GLN:NE2	1:E:182:PRO:O	2.44	0.50
1:H:3:LEU:HD11	1:H:140:GLN:HG2	1.94	0.50
1:F:98:ALA:HA	1:F:166:ASN:HD21	1.77	0.50
1:F:63:ARG:O	1:F:67:ILE:HD12	2.12	0.50
1:J:17:LEU:HD22	1:J:22:VAL:HG21	1.94	0.50
1:J:102:ASN:OD1	1:J:166:ASN:ND2	2.45	0.50
1:H:49:LEU:HD12	1:H:194:LEU:HD23	1.93	0.49
1:I:79:VAL:HG22	1:I:168:VAL:HG21	1.93	0.49
1:E:124:MET:HE2	1:E:128:ASP:HB3	1.94	0.49
1:E:73:GLN:HE22	1:E:183:VAL:HA	1.77	0.49
1:D:63:ARG:O	1:D:67:ILE:HD12	2.12	0.49
1:B:97:HIS:HB3	1:B:164:LEU:HD22	1.95	0.49
1:B:24:GLY:O	1:B:27:ARG:NH1	2.44	0.48
1:E:6:MET:HE2	1:E:143:HIS:CE1	2.48	0.48
1:H:92:THR:O	1:H:96:ILE:HG13	2.13	0.48
1:A:20:LEU:HB3	1:A:114:LEU:HB2	1.95	0.48
1:J:58:GLN:OE1	1:J:207:ARG:HG2	2.13	0.48
1:F:6:MET:HE1	1:F:16:VAL:HG21	1.96	0.48
1:F:121:THR:HB	1:F:129:ARG:HG2	1.96	0.48
1:B:52:LEU:HB3	1:B:215:SER:HB3	1.94	0.48
1:J:208:GLU:HG3	1:J:214:PHE:HZ	1.79	0.48
1:I:86:ILE:HB	1:I:89:SER:HB3	1.94	0.48
1:G:188:SER:OG	1:G:189:SER:N	2.46	0.47
1:I:15:LYS:NZ	1:I:149:GLU:OE1	2.35	0.47
1:I:161:HIS:NE2	1:I:176:ASP:OD1	2.38	0.47
1:D:92:THR:O	1:D:96:ILE:HG13	2.15	0.47
1:J:160:PHE:HE2	1:J:213:ARG:HB3	1.80	0.47
1:C:56:THR:OG1	1:C:208:GLU:OE2	2.25	0.47
1:I:63:ARG:O	1:I:67:ILE:HG13	2.14	0.47
1:E:164:LEU:HD21	1:E:166:ASN:ND2	2.29	0.47
1:H:121:THR:HA	1:H:124:MET:HG3	1.97	0.47
1:C:121:THR:HB	1:C:129:ARG:HG2	1.97	0.46
1:J:192:THR:O	1:J:192:THR:OG1	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:58:GLN:O	1:I:61:ASN:N	2.48	0.46
1:J:9:ASN:HB2	1:J:10:PRO:HD2	1.98	0.46
1:F:120:GLU:HG3	1:F:121:THR:HG23	1.97	0.46
1:J:176:ASP:HB3	1:J:179:MET:HG3	1.98	0.46
1:D:167:ASN:HB2	1:D:193:LEU:HD12	1.97	0.46
1:C:185:HIS:CE1	1:C:200:VAL:HG11	2.50	0.46
1:E:211:GLU:OE2	1:E:213:ARG:NH2	2.49	0.46
1:J:60:GLU:O	1:J:64:LYS:N	2.38	0.46
1:D:96:ILE:HG23	1:D:114:LEU:HD11	1.98	0.46
1:D:121:THR:HB	1:D:129:ARG:HG2	1.98	0.46
1:J:92:THR:HG21	1:J:143:HIS:HB2	1.98	0.46
1:D:63:ARG:O	1:D:66:GLN:N	2.42	0.46
1:J:98:ALA:O	1:J:102:ASN:HB2	2.16	0.46
1:H:53:PHE:HE1	1:H:163:ILE:HG12	1.81	0.45
1:J:65:LYS:O	1:J:69:GLU:HG3	2.17	0.45
1:E:121:THR:HA	1:E:124:MET:SD	2.56	0.45
1:F:52:LEU:HD21	2:F:301:GKO:N07	2.32	0.45
1:I:80:TYR:N	1:I:101:ASN:HD21	2.12	0.45
1:B:121:THR:HB	1:B:129:ARG:HG2	1.99	0.45
1:H:176:ASP:HB3	1:H:179:MET:HB2	1.99	0.45
1:G:160:PHE:HB2	2:G:301:GKO:O08	2.17	0.45
1:B:121:THR:HA	1:B:124:MET:HE3	1.99	0.45
1:J:96:ILE:HG12	1:J:114:LEU:HD13	1.99	0.44
1:A:172:LEU:HB2	1:A:188:SER:HB2	1.99	0.44
1:D:171:HIS:CD2	1:D:184:ASN:HD21	2.30	0.44
1:H:114:LEU:HD12	1:H:114:LEU:HA	1.88	0.44
1:G:56:THR:OG1	1:G:58:GLN:HG3	2.18	0.44
1:E:121:THR:HA	1:E:124:MET:HG3	2.00	0.44
1:H:205:THR:HA	1:H:214:PHE:CE2	2.52	0.44
1:J:121:THR:HA	1:J:124:MET:HE3	2.00	0.44
1:F:23:ALA:HB2	1:F:109:GLU:HG2	1.98	0.44
1:I:171:HIS:HB2	1:I:173:TYR:HE1	1.83	0.44
1:E:39:LEU:HD22	1:E:194:LEU:HG	1.99	0.44
1:J:53:PHE:HE1	1:J:163:ILE:HG23	1.82	0.44
1:C:60:GLU:HA	1:C:63:ARG:HB2	2.00	0.43
1:H:53:PHE:CE1	1:H:163:ILE:HG12	2.53	0.43
1:G:66:GLN:O	1:G:69:GLU:N	2.50	0.43
1:D:104:ASP:OD1	1:D:104:ASP:N	2.51	0.43
1:H:89:SER:O	1:H:93:ILE:HG12	2.19	0.43
1:A:17:LEU:HG	1:A:22:VAL:HB	2.01	0.43
1:D:116:GLN:O	1:D:120:GLU:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:20:LEU:HB3	1:E:114:LEU:HB2	1.99	0.43
1:I:125:SER:O	1:I:129:ARG:HG3	2.18	0.43
1:G:50:LEU:HD13	1:G:217:VAL:HG23	2.00	0.43
1:C:38:SER:O	1:C:41:SER:OG	2.36	0.43
1:B:39:LEU:HB3	1:B:194:LEU:HD23	2.01	0.43
1:B:62:PHE:C	1:B:62:PHE:CD1	2.95	0.43
1:F:64:LYS:HD3	1:F:64:LYS:H	1.82	0.43
1:I:164:LEU:HD23	1:I:175:LEU:HD22	2.01	0.43
1:B:148:GLN:HE21	1:B:148:GLN:HB2	1.67	0.43
1:D:35:GLU:O	1:D:39:LEU:HG	2.18	0.43
1:H:179:MET:HA	1:H:179:MET:HE2	1.99	0.43
1:I:106:LEU:HD23	1:I:106:LEU:HA	1.80	0.43
1:J:3:LEU:HD12	1:J:89:SER:HB3	2.01	0.43
1:G:132:CYS:HA	1:G:135:LYS:HE2	2.01	0.42
1:I:198:ALA:HA	1:I:201:CYS:HB2	2.00	0.42
1:D:29:VAL:HG11	1:D:43:PRO:HD2	2.01	0.42
1:I:39:LEU:HB3	1:I:194:LEU:HD23	2.02	0.42
1:I:63:ARG:HD2	1:I:179:MET:HE1	2.00	0.42
1:J:54:PRO:HD3	1:J:213:ARG:O	2.19	0.42
1:D:108:PHE:HB2	1:D:115:LYS:HE2	2.01	0.42
1:D:170:GLY:C	1:D:171:HIS:HD1	2.27	0.42
1:F:66:GLN:HG2	1:F:70:LEU:CD1	2.48	0.42
1:F:66:GLN:HG2	1:F:70:LEU:HD12	2.02	0.42
1:G:62:PHE:O	1:G:66:GLN:HB2	2.19	0.42
1:J:62:PHE:O	1:J:66:GLN:HB2	2.20	0.42
1:J:70:LEU:C	1:J:73:GLN:HE22	2.26	0.42
1:D:56:THR:HG21	1:D:208:GLU:OE2	2.19	0.42
1:E:124:MET:CE	1:E:128:ASP:HB3	2.48	0.42
1:A:123:LYS:HB3	1:A:123:LYS:HE3	1.93	0.42
1:B:20:LEU:HB3	1:B:114:LEU:HB2	2.01	0.42
1:C:16:VAL:HG22	1:C:146:VAL:HG21	2.01	0.42
1:C:86:ILE:HG22	1:C:88:ASN:H	1.84	0.42
1:I:80:TYR:HA	1:I:129:ARG:HH12	1.83	0.42
1:F:144:ASP:O	1:F:148:GLN:HG3	2.20	0.42
1:G:167:ASN:H	1:G:193:LEU:HD11	1.84	0.42
1:C:26:TRP:CZ2	1:C:105:LYS:HD2	2.54	0.42
1:I:93:ILE:HD13	1:I:93:ILE:HA	1.88	0.42
1:D:55:LEU:HD23	1:D:55:LEU:HA	1.79	0.41
1:I:203:GLU:O	1:I:207:ARG:HG3	2.20	0.41
1:G:135:LYS:HE3	1:G:135:LYS:HB2	1.74	0.41
1:A:195:LYS:HB3	1:A:195:LYS:HE2	1.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:58:GLN:OE1	1:H:207:ARG:HB3	2.20	0.41
1:A:3:LEU:HD12	1:A:89:SER:HB3	2.02	0.41
1:G:161:HIS:HE1	1:G:176:ASP:OD1	2.04	0.41
1:I:124:MET:HB2	1:I:129:ARG:HG2	2.02	0.41
1:C:63:ARG:HD2	1:C:63:ARG:HA	1.81	0.41
1:C:211:GLU:OE2	1:C:213:ARG:NH2	2.54	0.41
1:J:59:HIS:O	1:J:63:ARG:N	2.51	0.41
1:G:103:GLN:HE21	1:G:103:GLN:HB3	1.70	0.41
1:A:12:MET:HE2	1:A:15:LYS:HD2	2.02	0.41
1:C:34:LEU:HD13	1:C:201:CYS:HB3	2.02	0.41
1:D:20:LEU:HB3	1:D:114:LEU:HB2	2.02	0.41
1:D:127:GLU:H	1:D:127:GLU:CD	2.29	0.41
1:H:194:LEU:HD23	1:H:194:LEU:HA	1.97	0.40
1:G:15:LYS:NZ	1:G:149:GLU:OE1	2.42	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	221/227 (97%)	215 (97%)	6 (3%)	0	100	100
1	B	215/227 (95%)	212 (99%)	3 (1%)	0	100	100
1	C	218/227 (96%)	214 (98%)	4 (2%)	0	100	100
1	D	210/227 (92%)	207 (99%)	3 (1%)	0	100	100
1	E	215/227 (95%)	212 (99%)	3 (1%)	0	100	100
1	F	213/227 (94%)	210 (99%)	3 (1%)	0	100	100
1	G	215/227 (95%)	210 (98%)	5 (2%)	0	100	100
1	H	213/227 (94%)	210 (99%)	3 (1%)	0	100	100
1	I	213/227 (94%)	211 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	J	215/227 (95%)	209 (97%)	6 (3%)	0	100	100
All	All	2148/2270 (95%)	2110 (98%)	38 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	168/190 (88%)	168 (100%)	0	100	100
1	B	159/190 (84%)	156 (98%)	3 (2%)	50	66
1	C	153/190 (80%)	149 (97%)	4 (3%)	40	55
1	D	161/190 (85%)	159 (99%)	2 (1%)	63	78
1	E	162/190 (85%)	160 (99%)	2 (1%)	63	78
1	F	160/190 (84%)	159 (99%)	1 (1%)	78	89
1	G	154/190 (81%)	153 (99%)	1 (1%)	78	89
1	H	154/190 (81%)	151 (98%)	3 (2%)	50	66
1	I	156/190 (82%)	155 (99%)	1 (1%)	78	89
1	J	141/190 (74%)	140 (99%)	1 (1%)	76	87
All	All	1568/1900 (82%)	1550 (99%)	18 (1%)	65	79

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	148	GLN
1	B	151	GLN
1	B	209	GLN
1	C	38	SER
1	C	41	SER
1	C	132	CYS
1	C	140	GLN
1	D	70	LEU

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Mol	Chain	Res	Type
1	D	212	VAL
1	E	75	VAL
1	E	164	LEU
1	F	38	SER
1	G	78	LYS
1	H	34	LEU
1	H	39	LEU
1	H	113	VAL
1	I	157	LYS
1	J	42	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	HIS
1	A	116	GLN
1	B	103	GLN
1	B	166	ASN
1	B	167	ASN
1	C	102	ASN
1	C	140	GLN
1	C	185	HIS
1	D	84	GLN
1	D	140	GLN
1	D	171	HIS
1	E	61	ASN
1	E	116	GLN
1	F	116	GLN
1	F	140	GLN
1	F	148	GLN
1	F	151	GLN
1	G	84	GLN
1	G	103	GLN
1	H	59	HIS
1	I	101	ASN
1	I	102	ASN
1	I	185	HIS
1	J	73	GLN
1	J	148	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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	GKO	H	301	1	31,33,33	0.86	1 (3%)	35,44,44	0.89	2 (5%)
2	GKO	A	301	1	31,33,33	0.93	2 (6%)	35,44,44	1.32	4 (11%)
2	GKO	E	301	1	31,33,33	0.95	1 (3%)	35,44,44	0.98	2 (5%)
2	GKO	B	301	1	31,33,33	0.88	1 (3%)	35,44,44	0.98	2 (5%)
2	GKO	D	301	1	31,33,33	0.87	1 (3%)	35,44,44	0.82	0
2	GKO	J	301	1	31,33,33	0.86	1 (3%)	35,44,44	0.80	1 (2%)
2	GKO	I	301	1	31,33,33	0.90	1 (3%)	35,44,44	0.78	1 (2%)
2	GKO	F	301	1	31,33,33	0.90	2 (6%)	35,44,44	0.96	2 (5%)
2	GKO	C	301	1	31,33,33	0.92	1 (3%)	35,44,44	0.86	1 (2%)
2	GKO	G	301	1	31,33,33	0.88	2 (6%)	35,44,44	0.85	1 (2%)

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	GKO	H	301	1	-	1/21/37/37	0/3/3/3
2	GKO	A	301	1	-	2/21/37/37	0/3/3/3
2	GKO	E	301	1	-	0/21/37/37	0/3/3/3
2	GKO	B	301	1	-	3/21/37/37	0/3/3/3
2	GKO	D	301	1	-	6/21/37/37	0/3/3/3
2	GKO	J	301	1	-	2/21/37/37	0/3/3/3
2	GKO	I	301	1	-	0/21/37/37	0/3/3/3
2	GKO	F	301	1	-	1/21/37/37	0/3/3/3
2	GKO	C	301	1	-	4/21/37/37	0/3/3/3
2	GKO	G	301	1	-	1/21/37/37	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	GKO	C31-C23	2.51	1.55	1.53
2	C	301	GKO	C31-C23	2.51	1.55	1.53
2	E	301	GKO	C31-C23	2.48	1.55	1.53
2	A	301	GKO	C31-C23	2.48	1.55	1.53
2	I	301	GKO	C31-C23	2.43	1.55	1.53
2	B	301	GKO	C31-C23	2.35	1.55	1.53
2	H	301	GKO	C31-C23	2.30	1.55	1.53
2	D	301	GKO	C31-C23	2.28	1.55	1.53
2	J	301	GKO	C31-C23	2.19	1.55	1.53
2	G	301	GKO	C31-C23	2.11	1.55	1.53
2	F	301	GKO	C23-N14	-2.05	1.45	1.48
2	A	301	GKO	C17-N04	2.04	1.38	1.33
2	G	301	GKO	C23-N14	-2.01	1.45	1.48

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	GKO	C24-N04-C17	3.85	130.81	122.11
2	E	301	GKO	C23-C31-N16	3.75	114.93	109.33
2	F	301	GKO	C23-C31-N16	3.50	114.56	109.33
2	B	301	GKO	C23-C31-N16	3.19	114.09	109.33
2	C	301	GKO	C23-C31-N16	2.82	113.54	109.33
2	A	301	GKO	C23-C31-N16	2.64	113.27	109.33
2	J	301	GKO	C23-C31-N16	2.50	113.06	109.33
2	A	301	GKO	N07-C21-N10	-2.47	111.39	112.45
2	G	301	GKO	C23-C31-N16	2.46	113.00	109.33
2	A	301	GKO	C25-C24-N04	2.36	118.82	112.20
2	I	301	GKO	C23-C31-N16	2.33	112.80	109.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	GKO	C33-N14-C23	2.07	113.83	111.64
2	E	301	GKO	C33-N14-C23	2.05	113.81	111.64
2	H	301	GKO	C33-N14-C23	2.05	113.81	111.64
2	H	301	GKO	C23-C31-N16	2.01	112.33	109.33
2	B	301	GKO	C21-N10-C20	2.00	105.79	103.52

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	GKO	C25-C24-N04-C17
2	D	301	GKO	N04-C24-C25-C26
2	G	301	GKO	N04-C24-C25-C26
2	J	301	GKO	N04-C24-C25-C26
2	C	301	GKO	C15-C03-C17-O05
2	C	301	GKO	C15-C03-C17-N04
2	B	301	GKO	N04-C24-C25-C26
2	A	301	GKO	C24-C25-C26-C27
2	B	301	GKO	C24-C25-C26-C27
2	F	301	GKO	C24-C25-C26-C27
2	J	301	GKO	C24-C25-C26-C27
2	C	301	GKO	C06-C03-C17-O05
2	C	301	GKO	C06-C03-C17-N04
2	D	301	GKO	C06-C03-C17-O05
2	D	301	GKO	C06-C03-C17-N04
2	B	301	GKO	N12-C22-C23-N14
2	D	301	GKO	C15-C03-C17-O05
2	D	301	GKO	C15-C03-C17-N04
2	D	301	GKO	C24-C25-C26-C27
2	H	301	GKO	C19-C20-N12-C22

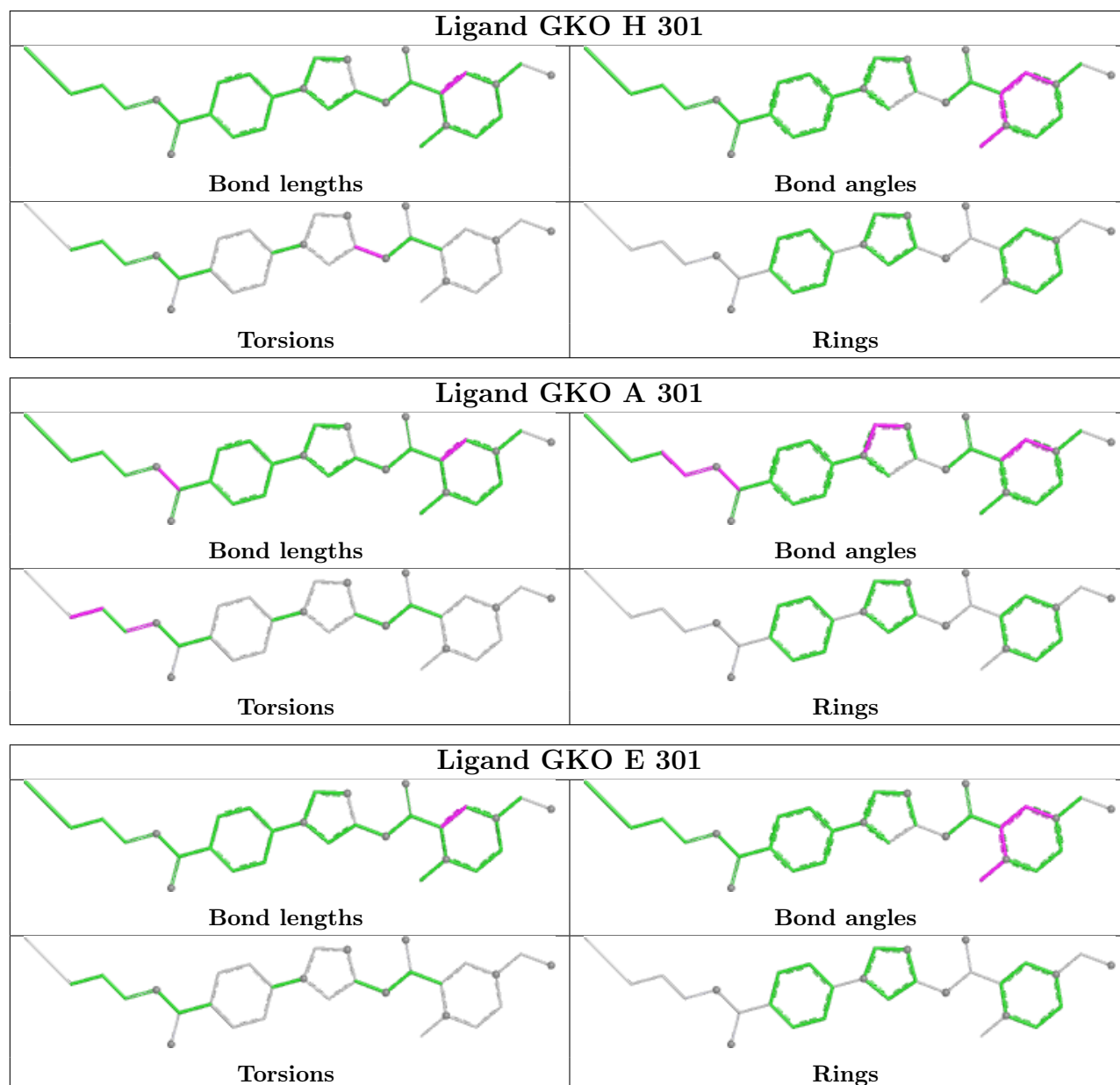
There are no ring outliers.

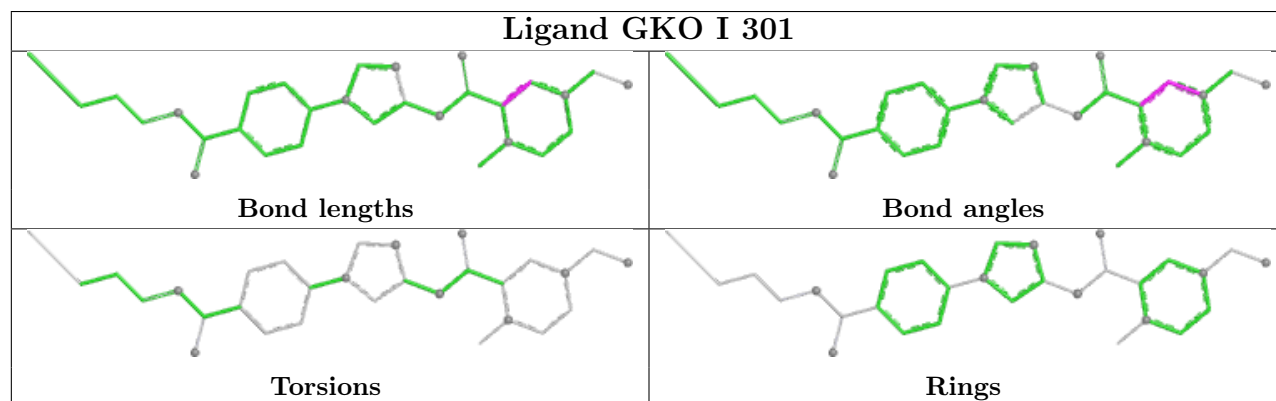
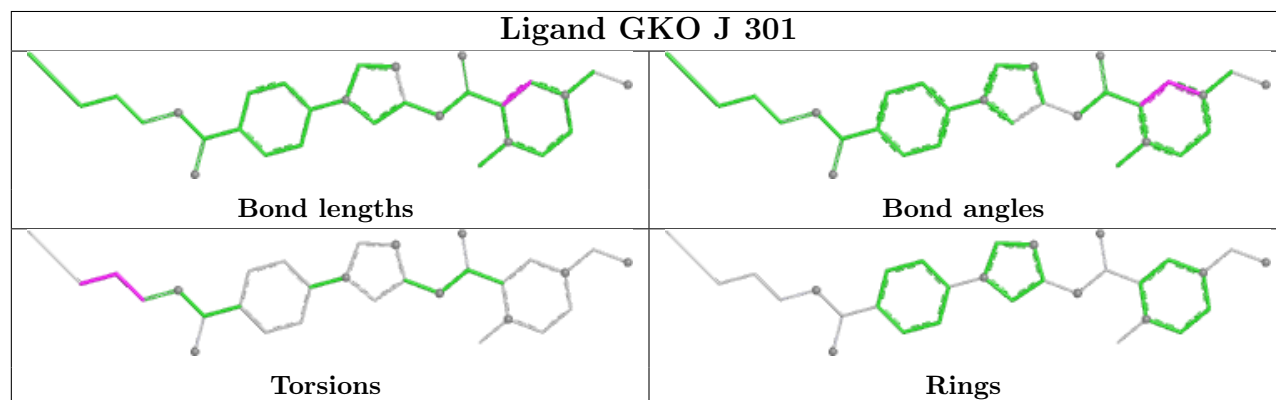
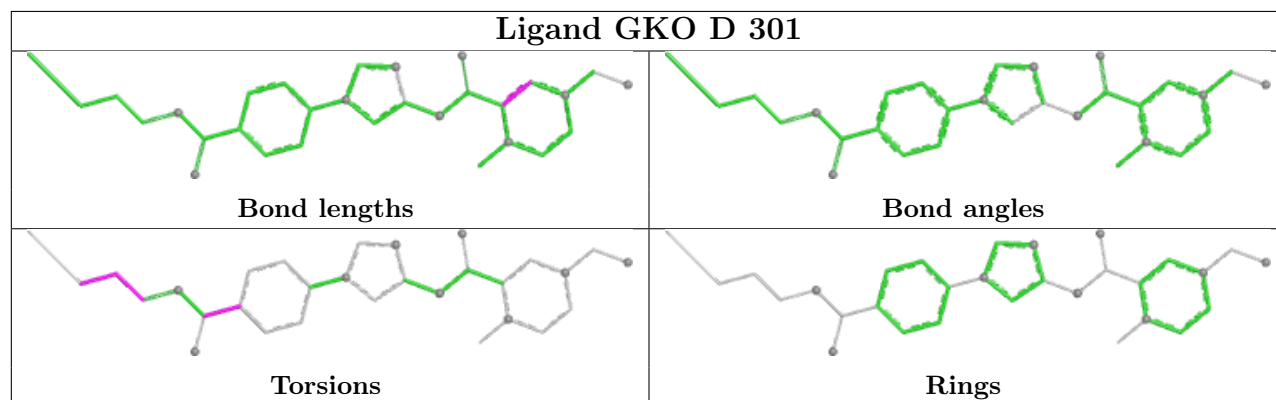
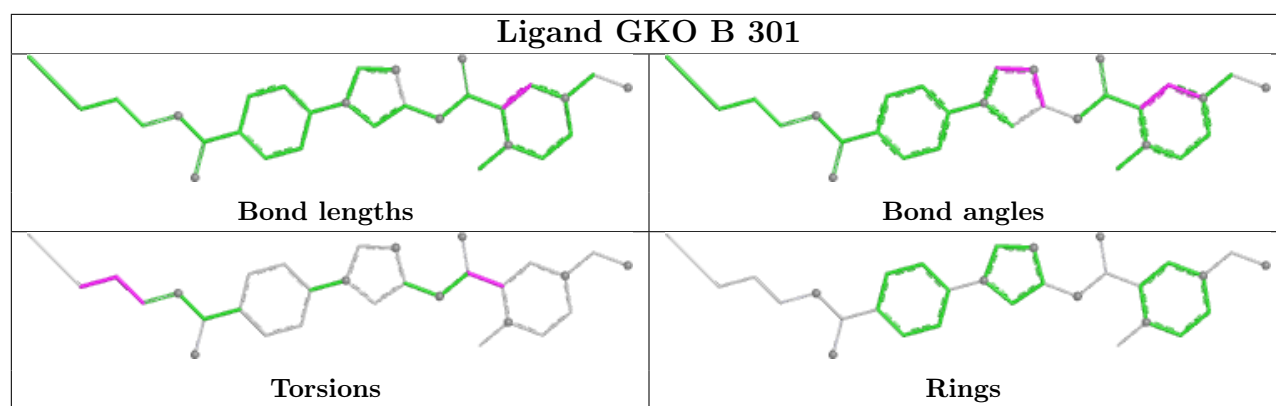
2 monomers are involved in 2 short contacts:

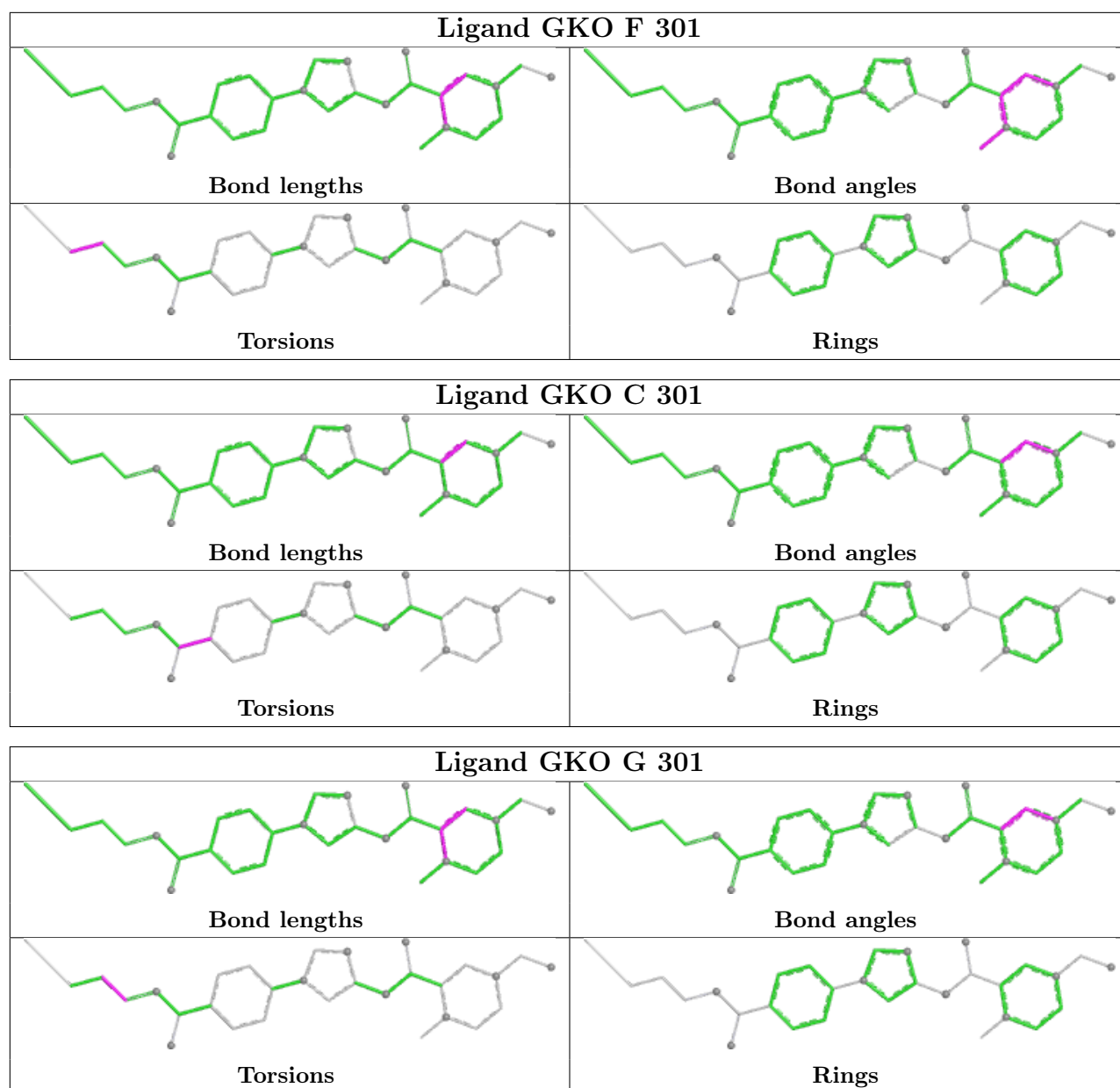
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	301	GKO	1	0
2	G	301	GKO	1	0

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

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	223/227 (98%)	0.24	6 (2%) 56 53	39, 57, 107, 158	0
1	B	219/227 (96%)	0.39	10 (4%) 37 34	43, 66, 110, 154	0
1	C	220/227 (96%)	0.73	11 (5%) 34 31	48, 79, 122, 144	0
1	D	216/227 (95%)	0.65	10 (4%) 37 34	51, 79, 111, 138	0
1	E	219/227 (96%)	0.36	4 (1%) 67 64	45, 71, 109, 139	0
1	F	217/227 (95%)	0.54	5 (2%) 61 58	53, 74, 106, 129	0
1	G	219/227 (96%)	0.91	19 (8%) 16 13	51, 91, 127, 151	0
1	H	217/227 (95%)	0.75	8 (3%) 45 42	58, 91, 120, 153	0
1	I	217/227 (95%)	0.93	17 (7%) 19 16	61, 89, 123, 161	0
1	J	219/227 (96%)	1.36	58 (26%) 1 1	50, 104, 150, 182	0
All	All	2186/2270 (96%)	0.68	148 (6%) 23 20	39, 79, 124, 182	0

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	187	ALA	4.3
1	B	212	VAL	3.9
1	J	189	SER	3.9
1	G	67	ILE	3.8
1	J	46	ALA	3.6
1	J	172	LEU	3.6
1	J	79	VAL	3.6
1	J	168	VAL	3.5
1	J	80	TYR	3.4
1	J	100	ALA	3.4
1	J	113	VAL	3.4
1	G	168	VAL	3.4
1	B	222	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	I	55	LEU	3.3
1	G	158	VAL	3.3
1	G	108	PHE	3.2
1	J	181	PHE	3.2
1	J	77	PRO	3.1
1	C	71	LYS	3.1
1	J	70	LEU	3.0
1	G	26	TRP	3.0
1	J	169	ASP	3.0
1	B	214	PHE	2.9
1	H	212	VAL	2.9
1	I	158	VAL	2.9
1	J	220	CYS	2.9
1	I	81	PHE	2.9
1	D	154	VAL	2.9
1	I	26	TRP	2.8
1	J	26	TRP	2.8
1	J	219	LEU	2.8
1	C	62	PHE	2.8
1	J	107	GLY	2.8
1	J	106	LEU	2.8
1	J	110	ASP	2.8
1	J	197	ALA	2.7
1	G	169	ASP	2.7
1	A	158	VAL	2.7
1	J	99	VAL	2.7
1	C	192	THR	2.7
1	I	164	LEU	2.7
1	J	175	LEU	2.7
1	G	100	ALA	2.7
1	J	211	GLU	2.7
1	J	22	VAL	2.6
1	G	79	VAL	2.6
1	I	106	LEU	2.6
1	I	46	ALA	2.6
1	J	117	PHE	2.6
1	A	72	GLY	2.6
1	I	86	ILE	2.6
1	F	32	LEU	2.6
1	G	193	LEU	2.5
1	J	0	SER	2.5
1	C	153	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	J	101	ASN	2.5
1	J	185	HIS	2.5
1	I	108	PHE	2.5
1	A	75	VAL	2.5
1	J	170	GLY	2.5
1	J	184	ASN	2.5
1	J	126	PRO	2.5
1	I	36	GLU	2.4
1	H	70	LEU	2.4
1	E	87	GLY	2.4
1	D	212	VAL	2.4
1	F	108	PHE	2.4
1	J	75	VAL	2.4
1	G	80	TYR	2.4
1	H	191	ASP	2.4
1	C	140	GLN	2.4
1	I	110	ASP	2.3
1	J	173	TYR	2.3
1	J	123	LYS	2.3
1	C	67	ILE	2.3
1	J	154	VAL	2.3
1	I	32	LEU	2.3
1	I	79	VAL	2.3
1	I	168	VAL	2.3
1	G	197	ALA	2.3
1	J	187	ALA	2.3
1	H	3	LEU	2.3
1	I	194	LEU	2.3
1	F	159	ASN	2.3
1	J	180	PRO	2.3
1	H	177	GLY	2.3
1	J	111	GLY	2.3
1	D	188	SER	2.3
1	J	76	SER	2.3
1	C	159	ASN	2.2
1	G	220	CYS	2.2
1	D	94	GLY	2.2
1	G	186	GLY	2.2
1	H	147	ALA	2.2
1	J	23	ALA	2.2
1	I	85	THR	2.2
1	J	67	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	55	LEU	2.2
1	D	70	LEU	2.2
1	J	81	PHE	2.2
1	G	167	ASN	2.2
1	J	132	CYS	2.2
1	F	46	ALA	2.2
1	F	153	ARG	2.2
1	J	21	GLY	2.2
1	J	2	GLN	2.2
1	J	25	GLN	2.2
1	J	130	ALA	2.2
1	J	39	LEU	2.2
1	A	1	MET	2.2
1	J	179	MET	2.2
1	J	182	PRO	2.2
1	C	111	GLY	2.2
1	E	166	ASN	2.1
1	J	159	ASN	2.1
1	J	171	HIS	2.1
1	B	210	GLY	2.1
1	I	210	GLY	2.1
1	D	222	ALA	2.1
1	H	222	ALA	2.1
1	B	156	ASP	2.1
1	C	158	VAL	2.1
1	D	181	PHE	2.1
1	J	45	PRO	2.1
1	B	213	ARG	2.1
1	J	42	VAL	2.1
1	D	214	PHE	2.1
1	G	181	PHE	2.1
1	J	48	ALA	2.1
1	A	73	GLN	2.1
1	A	156	ASP	2.1
1	B	159	ASN	2.1
1	G	77	PRO	2.1
1	J	183	VAL	2.1
1	H	87	GLY	2.1
1	J	55	LEU	2.0
1	D	67	ILE	2.0
1	G	173	TYR	2.0
1	E	38	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	190	GLU	2.0
1	J	38	SER	2.0
1	B	23	ALA	2.0
1	B	44	ALA	2.0
1	D	23	ALA	2.0
1	J	116	GLN	2.0
1	B	70	LEU	2.0
1	C	193	LEU	2.0
1	G	172	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

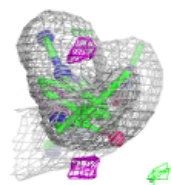
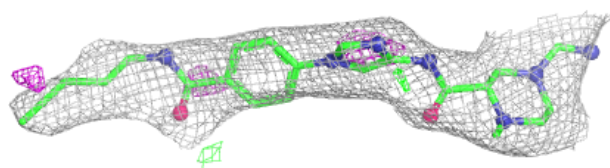
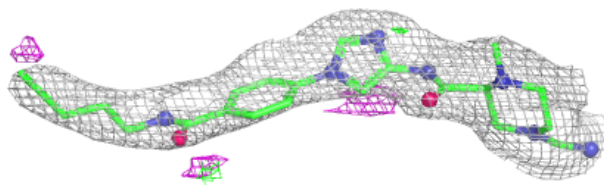
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GKO	H	301	31/31	0.88	0.16	73,80,90,93	0
2	GKO	B	301	31/31	0.89	0.13	41,52,92,99	0
2	GKO	C	301	31/31	0.90	0.13	57,70,77,78	0
2	GKO	F	301	31/31	0.91	0.13	55,64,74,79	0
2	GKO	D	301	31/31	0.91	0.12	61,70,83,88	0
2	GKO	I	301	31/31	0.91	0.12	59,65,74,78	0
2	GKO	A	301	31/31	0.92	0.12	41,54,78,80	0
2	GKO	E	301	31/31	0.93	0.11	47,61,68,70	0
2	GKO	G	301	31/31	0.93	0.11	51,62,74,79	0
2	GKO	J	301	31/31	0.93	0.11	43,60,71,80	0

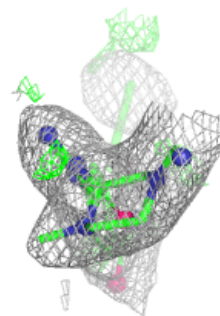
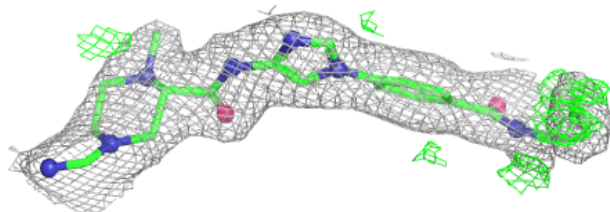
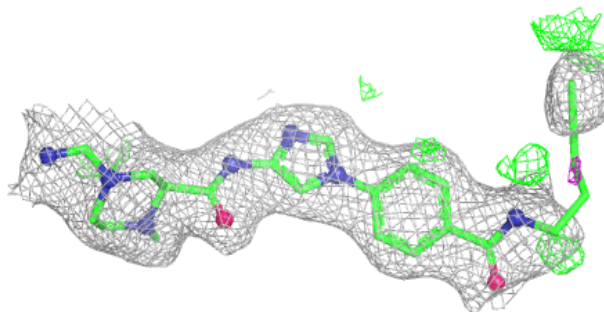
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GKO H 301:

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

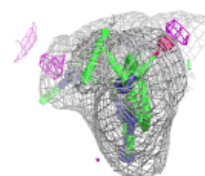
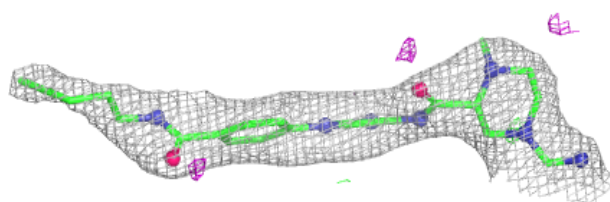
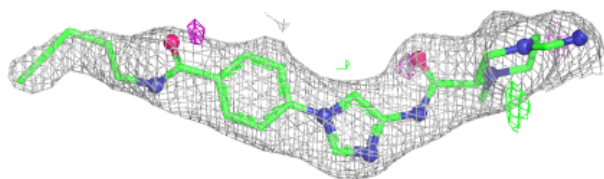
**Electron density around GKO B 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

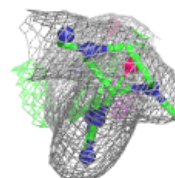
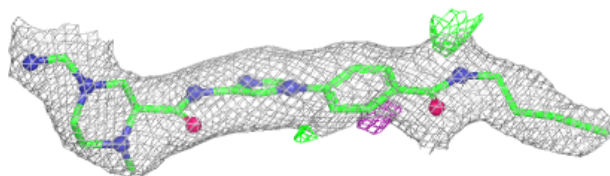
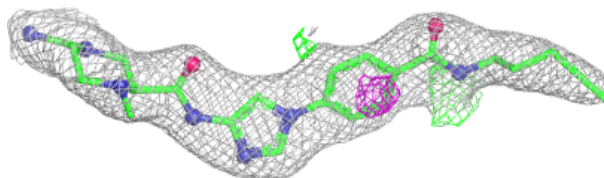


Electron density around GKO C 301:

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

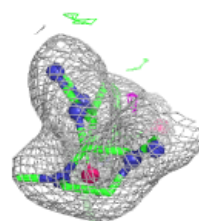
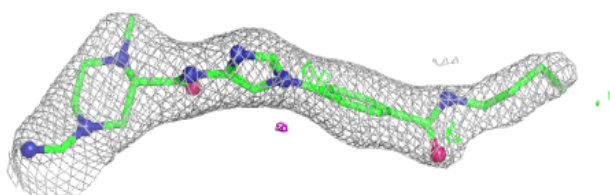
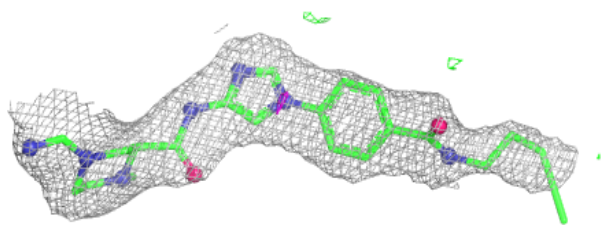
**Electron density around GKO F 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

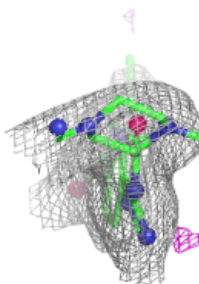
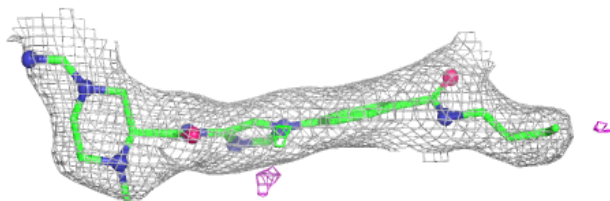
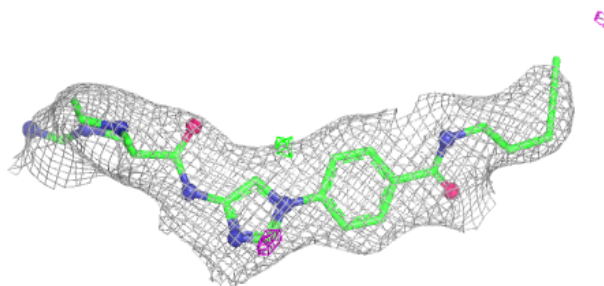


Electron density around GKO D 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

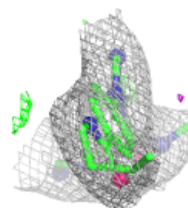
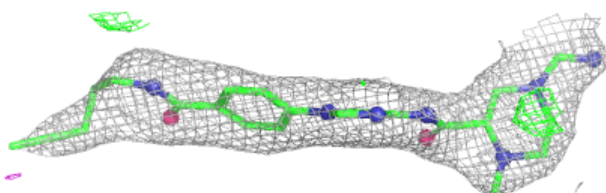
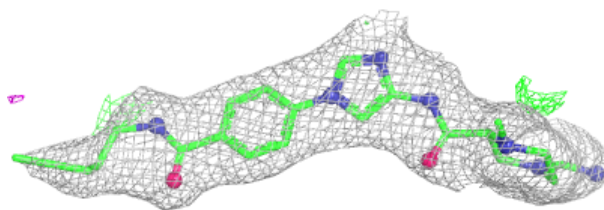
**Electron density around GKO I 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

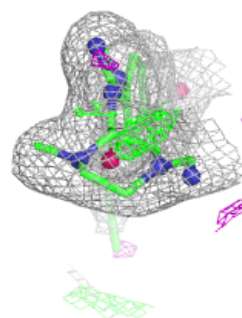
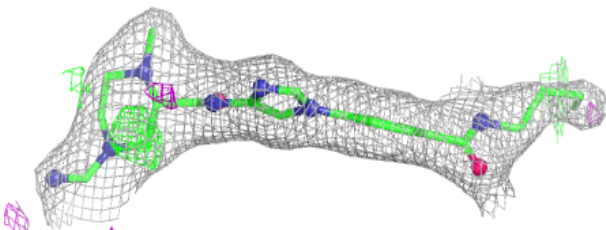
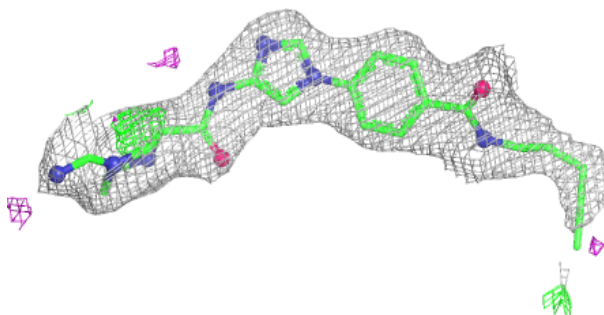


Electron density around GKO A 301:

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and green (positive)

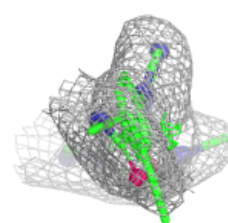
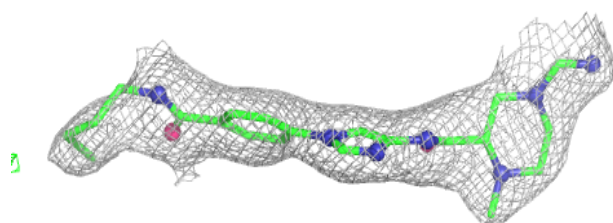
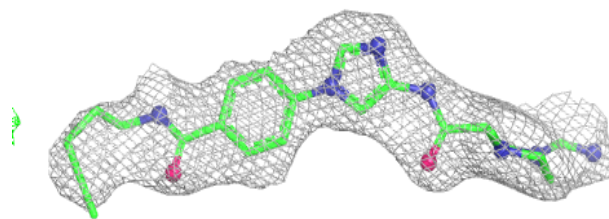
**Electron density around GKO E 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

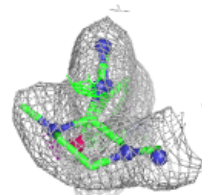
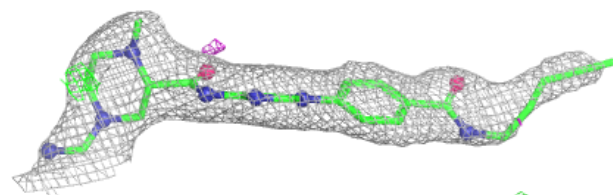
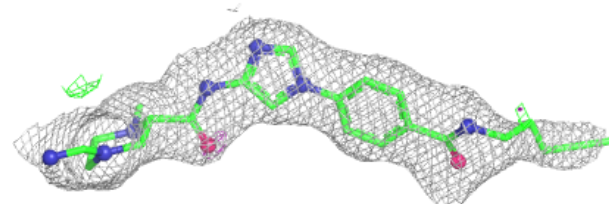


Electron density around GKO G 301:

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

**Electron density around GKO J 301:**

$2mF_o-DF_c$ (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.