



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2021 – 04:23 PM EST

PDB ID : 5Q04
Title : Human liver fructose-1,6-bisphosphatase 1 (fructose 1,6-bisphosphate 1-phosphatase, E.C.3.1.3.11) complexed with the allosteric inhibitor 1-(4-bromo-5-chlorothiophen-2-yl)sulfonyl-3-(5-bromo-1,3-thiazol-2-yl)urea
Authors : Ruf, A.; Joseph, C.; Alker, A.; Banner, D.; Tetaz, T.; Benz, J.; Kuhn, B.; Rudolph, M.G.; Yang, H.; Shao, C.; Burley, S.K.
Deposited on : 2017-04-18
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.16
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) : 2.16

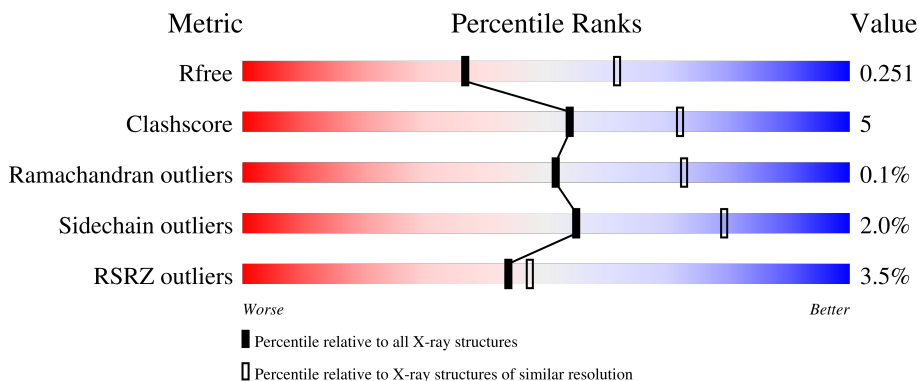
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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.



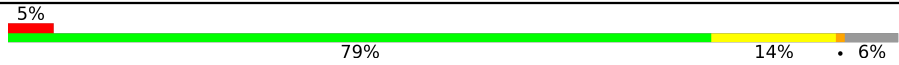


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	338	
1	B	338	
1	C	338	
1	D	338	
1	E	338	

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Mol	Chain	Length	Quality of chain
1	F	338	 5% 79% 14% • 6%
1	G	338	 % 80% 13% • 6%
1	H	338	 3% 81% 12% • 6%

2 Entry composition [i](#)

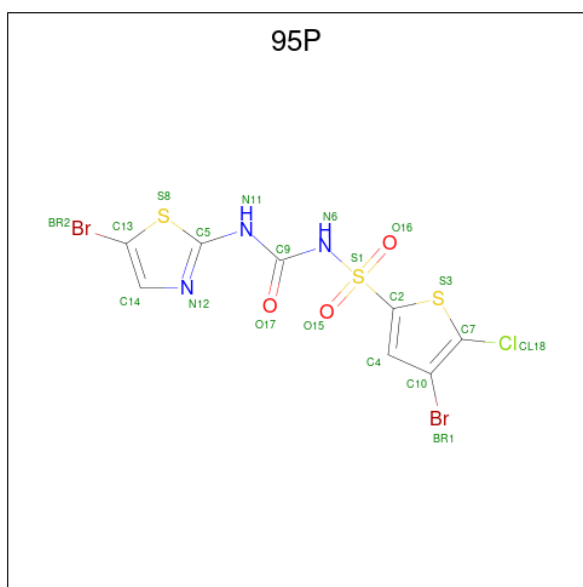
There are 3 unique types of molecules in this entry. The entry contains 19992 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	319	2446	1558	412	459	17	0	1	0
1	B	317	2431	1548	409	457	17	0	1	0
1	C	317	2425	1545	406	457	17	0	0	0
1	D	317	2431	1548	409	457	17	0	1	0
1	E	319	2447	1559	412	459	17	0	1	0
1	F	317	2430	1547	409	457	17	0	1	0
1	G	317	2430	1547	409	457	17	0	1	0
1	H	317	2425	1545	406	457	17	0	0	0

- Molecule 2 is 4-bromo-N-[(5-bromo-1,3-thiazol-2-yl)carbamoyl]-5-chlorothiophene-2-sulfonamide (three-letter code: 95P) (formula: $C_8H_4Br_2ClN_3O_3S_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	
			Total	Br	C	Cl	N	O			S
2	A	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	B	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	C	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	D	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	E	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	F	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	G	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		
2	H	1	Total	Br	C	Cl	N	O	S	0	0
			20	2	8	1	3	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	45	Total	O	0	0
			45	45		
3	B	38	Total	O	0	0
			38	38		
3	C	53	Total	O	0	0
			53	53		
3	D	60	Total	O	0	0
			60	60		

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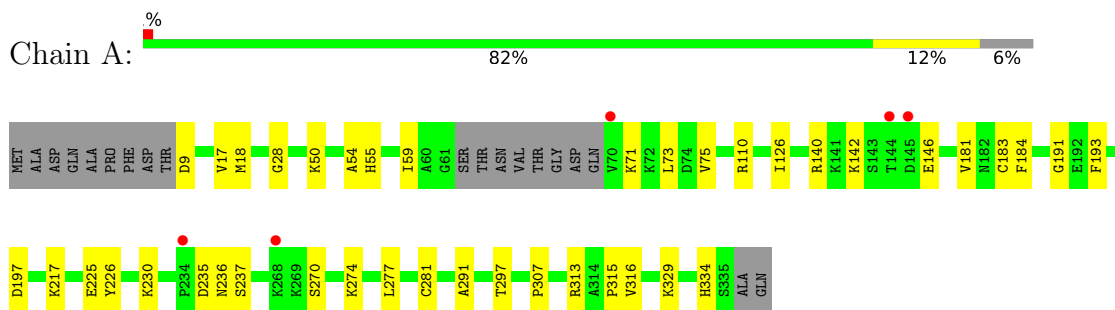
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	35	Total 35	O 35	0	0
3	F	19	Total 19	O 19	0	0
3	G	67	Total 67	O 67	0	0
3	H	50	Total 50	O 50	0	0

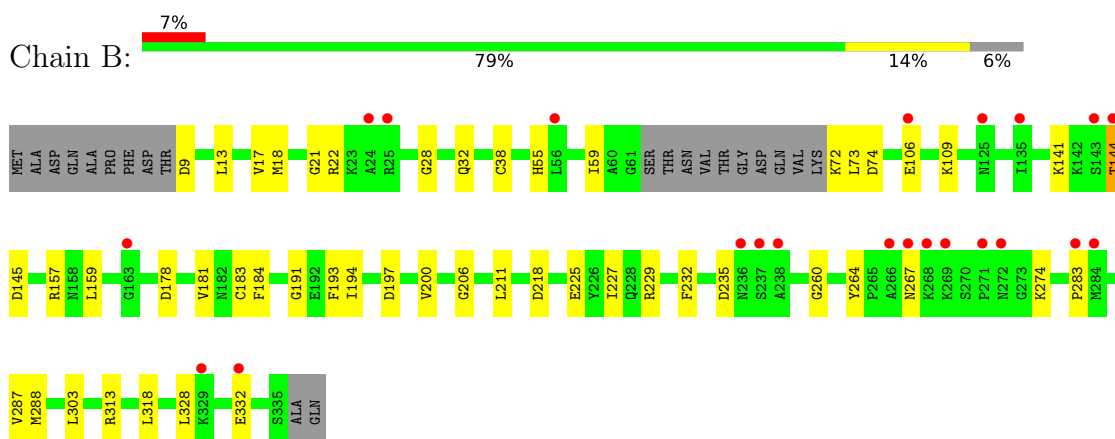
3 Residue-property plots [i](#)

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.

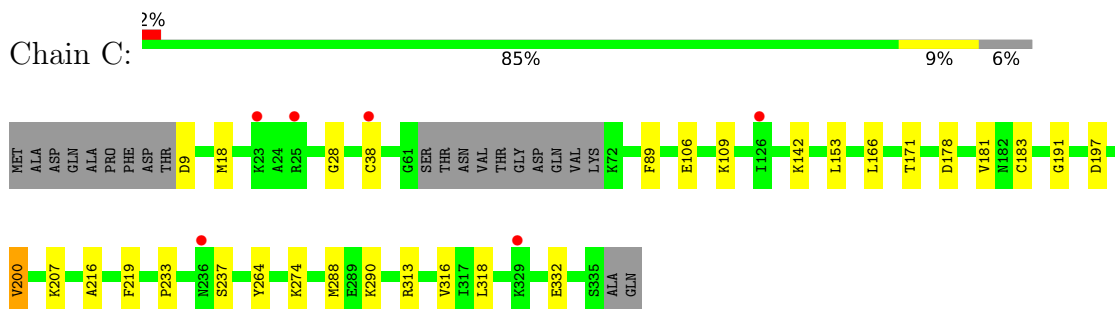
- Molecule 1: Fructose-1,6-bisphosphatase 1




- Molecule 1: Fructose-1,6-bisphosphatase 1

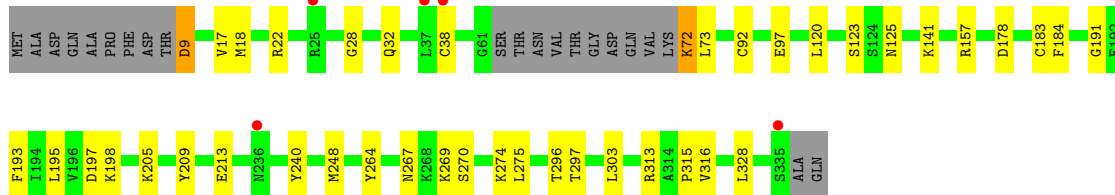


- Molecule 1: Fructose-1,6-bisphosphatase 1




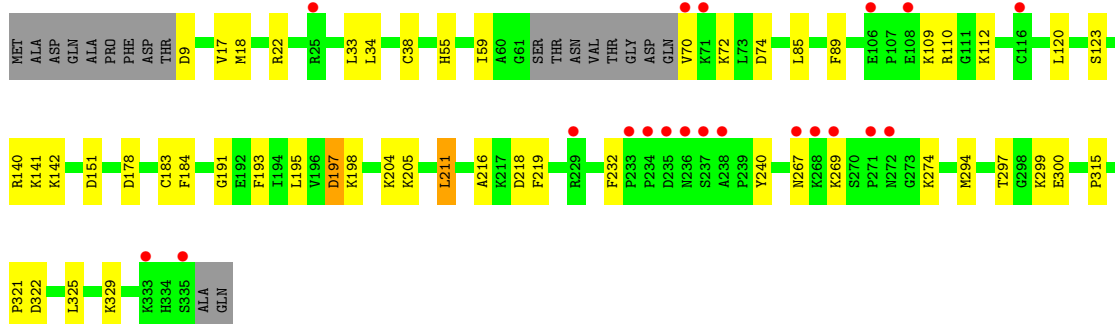
- Molecule 1: Fructose-1,6-bisphosphatase 1

Chain D: 




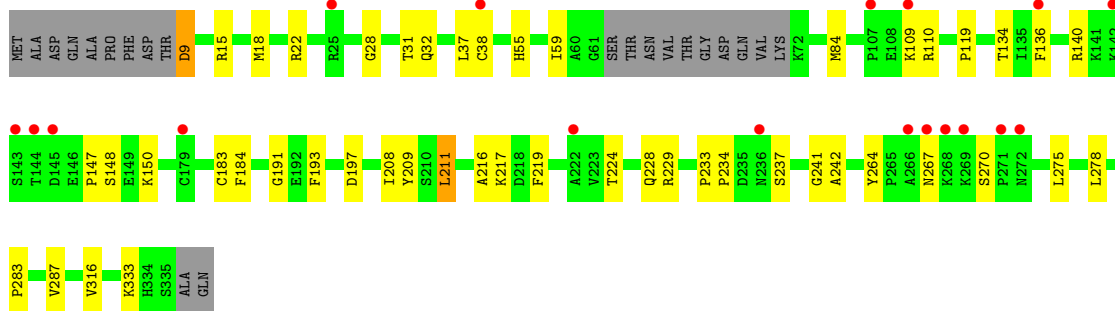
• Molecule 1: Fructose-1,6-bisphosphatase 1

Chain E: 




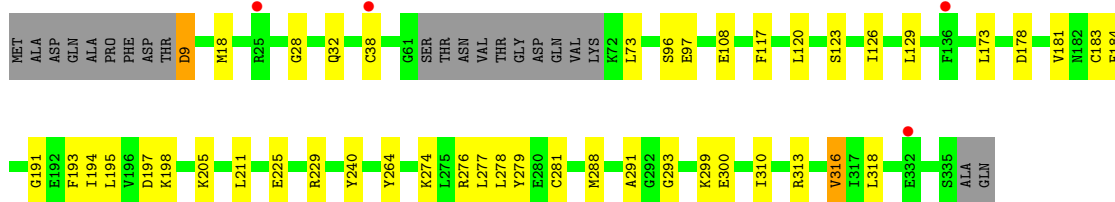
• Molecule 1: Fructose-1,6-bisphosphatase 1

Chain F: 

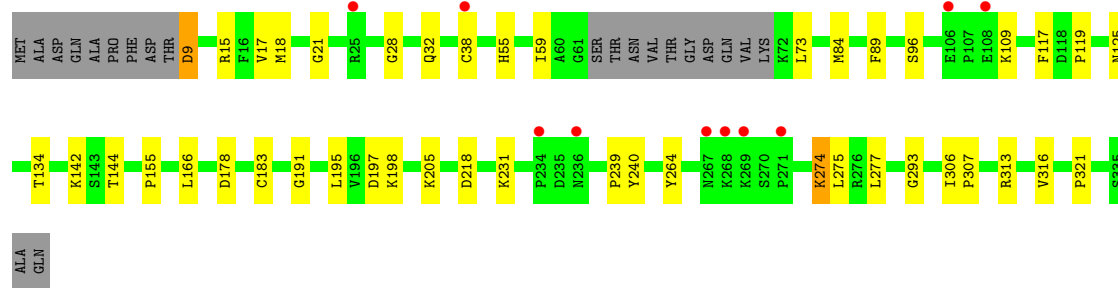
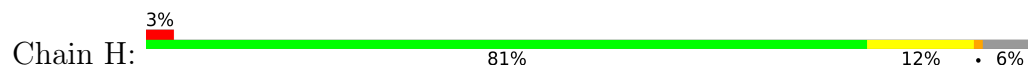


• Molecule 1: Fructose-1,6-bisphosphatase 1

Chain G: 



• Molecule 1: Fructose-1,6-bisphosphatase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.28Å 286.68Å 83.27Å 90.00° 97.68° 90.00°	Depositor
Resolution (Å)	29.23 – 2.50 29.24 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.9 (29.23-2.50) 94.9 (29.24-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.51Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.186 , 0.251 0.186 , 0.251	Depositor DCC
R_{free} test set	5006 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtrriage
Anisotropy	0.416	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 65.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19992	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 95P

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	0.43	0/2495	0.63	0/3368
1	B	0.42	0/2479	0.61	0/3347
1	C	0.48	1/2468 (0.0%)	0.63	0/3333
1	D	0.50	1/2479 (0.0%)	0.68	0/3347
1	E	0.40	0/2495	0.62	1/3368 (0.0%)
1	F	0.42	0/2479	0.62	1/3347 (0.0%)
1	G	0.44	0/2479	0.63	1/3347 (0.0%)
1	H	0.45	0/2468	0.65	0/3333
All	All	0.44	2/19842 (0.0%)	0.63	3/26790 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	97	GLU	CG-CD	5.61	1.60	1.51
1	C	153	LEU	C-N	-5.55	1.21	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	211	LEU	CA-CB-CG	6.26	129.70	115.30
1	F	211	LEU	CA-CB-CG	6.01	129.13	115.30
1	G	173	LEU	CB-CG-CD2	-5.58	101.51	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2446	0	2496	28	0
1	B	2431	0	2478	34	0
1	C	2425	0	2469	16	0
1	D	2431	0	2478	24	0
1	E	2447	0	2500	32	0
1	F	2430	0	2474	35	0
1	G	2430	0	2474	28	0
1	H	2425	0	2469	36	0
2	A	20	0	0	3	0
2	B	20	0	0	4	0
2	C	20	0	0	3	0
2	D	20	0	0	3	0
2	E	20	0	0	2	0
2	F	20	0	0	3	0
2	G	20	0	0	2	0
2	H	20	0	0	5	0
3	A	45	0	0	4	0
3	B	38	0	0	0	0
3	C	53	0	0	1	0
3	D	60	0	0	1	0
3	E	35	0	0	1	0
3	F	19	0	0	3	0
3	G	67	0	0	2	0
3	H	50	0	0	2	0
All	All	19992	0	19838	211	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 5.

All (211) 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:22:ARG:NH1	1:D:32:GLN:OE1	2.08	0.86
1:H:183:CYS:HB2	1:H:197:ASP:HB2	1.61	0.83
1:A:274:LYS:O	1:A:313:ARG:NH1	2.12	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:LEU:HD11	1:B:227:ILE:HD11	1.65	0.79
1:H:17:VAL:HG13	2:H:401:95P:BR1	2.41	0.75
1:A:235:ASP:O	1:A:237:SER:N	2.20	0.74
1:G:274:LYS:O	1:G:313:ARG:NH1	2.21	0.73
1:A:146:GLU:O	3:A:501:HOH:O	2.05	0.73
1:B:191:GLY:HA3	1:D:191:GLY:HA3	1.71	0.72
1:H:274:LYS:O	1:H:313:ARG:NH1	2.21	0.72
1:B:109:LYS:NZ	1:D:9:ASP:OD2	2.24	0.71
1:E:294:MET:HE1	1:E:325:LEU:HG	1.73	0.70
1:F:31:THR:HG22	1:H:18:MET:HE3	1.73	0.70
1:G:108:GLU:H	1:G:108:GLU:CD	1.98	0.67
1:C:183:CYS:HB2	1:C:197:ASP:HB2	1.78	0.66
1:A:274:LYS:HA	1:A:313:ARG:HD3	1.77	0.65
1:B:183:CYS:HB2	1:B:197:ASP:HB2	1.77	0.65
1:H:119:PRO:HA	1:H:134:THR:HG23	1.79	0.65
1:E:183:CYS:HB2	1:E:197:ASP:HB2	1.80	0.64
1:F:233:PRO:HG2	1:F:237:SER:O	1.97	0.64
1:F:9:ASP:N	3:F:501:HOH:O	2.32	0.63
1:F:191:GLY:HA3	1:H:191:GLY:HA3	1.80	0.63
1:H:18:MET:HG2	2:H:401:95P:BR2	2.55	0.62
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.82	0.61
1:G:195:LEU:HD21	1:G:198:LYS:HG2	1.82	0.61
1:A:28:GLY:HA2	2:A:401:95P:O17	2.01	0.60
1:E:22:ARG:NH1	1:G:32:GLN:OE1	2.35	0.59
1:A:183:CYS:HB2	1:A:197:ASP:HB2	1.86	0.58
1:F:333:LYS:NZ	3:F:502:HOH:O	2.36	0.58
1:D:92:CYS:SG	1:D:141:LYS:HD2	2.43	0.58
1:F:109:LYS:NZ	1:H:9:ASP:OD2	2.32	0.58
1:B:18:MET:HG2	2:B:401:95P:BR2	2.59	0.58
1:D:183:CYS:HB2	1:D:197:ASP:HB2	1.86	0.58
1:B:106:GLU:OE2	1:B:109:LYS:HE2	2.03	0.58
1:F:148:SER:OG	1:F:150:LYS:HG2	2.02	0.57
1:G:18:MET:HG2	2:G:401:95P:BR2	2.59	0.57
1:A:297:THR:HB	1:A:315:PRO:HG2	1.86	0.57
1:B:32:GLN:OE1	1:D:22:ARG:NH1	2.37	0.57
1:F:183:CYS:HB2	1:F:197:ASP:HB2	1.86	0.56
1:G:205:LYS:HE2	1:G:240:TYR:OH	2.06	0.56
1:G:316:VAL:HG21	1:G:318:LEU:HD23	1.88	0.56
1:D:205:LYS:HE2	1:D:240:TYR:OH	2.06	0.56
1:F:84:MET:O	1:H:15:ARG:NH2	2.39	0.56
1:H:89:PHE:HE1	1:H:109:LYS:HG2	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:17:VAL:HG13	2:D:401:95P:BR1	2.61	0.55
1:G:97:GLU:OE1	3:G:501:HOH:O	2.18	0.55
1:C:264:TYR:CZ	1:C:274:LYS:HD3	2.42	0.55
1:H:89:PHE:CE1	1:H:109:LYS:HG2	2.43	0.54
1:H:28:GLY:HA2	2:H:401:95P:C9	2.38	0.54
1:B:17:VAL:HG13	2:B:401:95P:BR1	2.63	0.54
1:H:9:ASP:N	3:H:503:HOH:O	2.41	0.54
1:E:269:LYS:NZ	1:E:274:LYS:HE2	2.23	0.53
1:B:328:LEU:O	1:B:332:GLU:HG3	2.09	0.53
1:F:9:ASP:N	3:F:504:HOH:O	2.42	0.52
1:D:18:MET:HG2	2:D:401:95P:BR2	2.64	0.52
1:F:229:ARG:HH21	1:F:234:PRO:HG3	1.73	0.52
1:E:18:MET:HG2	2:E:401:95P:BR2	2.64	0.52
1:C:216:ALA:HA	1:C:219:PHE:CD2	2.45	0.52
1:C:332:GLU:HB3	3:C:538:HOH:O	2.09	0.52
1:G:73:LEU:HD23	1:G:126:ILE:HD13	1.90	0.52
1:G:181:VAL:HG21	1:G:291:ALA:HB2	1.91	0.52
1:E:204:LYS:HE3	1:E:322:ASP:OD2	2.10	0.51
1:E:191:GLY:HA3	1:G:191:GLY:HA3	1.91	0.51
1:D:296:THR:HG21	1:D:328:LEU:HD21	1.93	0.51
1:E:33:LEU:HD11	1:E:85:LEU:HD22	1.92	0.51
1:F:32:GLN:HG2	1:H:18:MET:HE2	1.93	0.51
1:G:28:GLY:HA2	2:G:401:95P:O17	2.11	0.50
1:D:120:LEU:HG	1:D:123:SER:HB3	1.94	0.50
1:B:235:ASP:HA	1:E:70:VAL:HG21	1.93	0.50
1:B:218:ASP:HB3	1:B:267:ASN:HB2	1.93	0.49
1:D:195:LEU:HD21	1:D:198:LYS:HG2	1.94	0.49
1:G:129:LEU:HD12	1:H:166:LEU:HD23	1.94	0.49
1:A:140[B]:ARG:NH1	3:A:505:HOH:O	2.33	0.49
1:A:50:LYS:NZ	3:A:511:HOH:O	2.44	0.49
1:C:181:VAL:HB	1:C:200:VAL:HG22	1.94	0.49
1:H:125:ASN:ND2	3:H:502:HOH:O	2.35	0.49
1:G:108:GLU:N	1:G:108:GLU:OE1	2.43	0.49
1:E:195:LEU:HD21	1:E:198:LYS:HG2	1.94	0.49
1:G:276:ARG:HG3	1:G:279:TYR:CE1	2.47	0.49
1:A:270:SER:OG	1:A:274:LYS:HE3	2.13	0.49
1:A:277:LEU:HD23	1:A:307:PRO:HB3	1.95	0.49
1:F:18:MET:HG2	2:F:401:95P:BR2	2.68	0.49
1:A:329:LYS:NZ	3:A:509:HOH:O	2.42	0.49
1:B:274:LYS:HA	1:B:313:ARG:HD3	1.94	0.49
1:C:274:LYS:O	1:C:313:ARG:NH1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:GLY:HA2	2:B:401:95P:S3	2.53	0.48
1:D:157:ARG:HG2	1:D:303:LEU:O	2.12	0.48
1:B:197:ASP:HB3	1:B:200:VAL:CG1	2.43	0.48
1:H:195:LEU:HD21	1:H:198:LYS:HG2	1.95	0.48
1:F:31:THR:HG22	1:H:18:MET:CE	2.41	0.48
1:H:55:HIS:HA	1:H:59:ILE:HG22	1.95	0.48
1:E:112:LYS:HD2	1:E:140[B]:ARG:NH2	2.29	0.48
1:A:181:VAL:HG21	1:A:291:ALA:HB2	1.95	0.48
1:G:277:LEU:HA	1:G:281:CYS:HB2	1.95	0.48
1:C:18:MET:HG2	2:C:401:95P:BR2	2.68	0.48
1:G:299:LYS:HG3	1:G:300:GLU:HG2	1.94	0.48
1:E:232:PHE:HE1	1:F:217:LYS:HA	1.78	0.47
1:F:216:ALA:HA	1:F:219:PHE:CD2	2.49	0.47
1:B:157:ARG:HG2	1:B:303:LEU:O	2.15	0.47
1:E:184:PHE:HB3	1:E:193:PHE:HB3	1.96	0.47
1:H:73:LEU:HA	1:H:73:LEU:HD12	1.62	0.47
1:B:55:HIS:HA	1:B:59:ILE:HG22	1.96	0.47
1:D:297:THR:HB	1:D:315:PRO:HG2	1.96	0.47
1:F:110:ARG:HD3	1:F:147:PRO:HG3	1.96	0.47
1:F:275:LEU:HD12	1:F:316:VAL:HG21	1.97	0.47
1:F:55:HIS:HA	1:F:59:ILE:HG22	1.97	0.47
1:H:205:LYS:HE2	1:H:240:TYR:OH	2.14	0.47
1:F:32:GLN:CG	1:H:18:MET:HE2	2.45	0.47
1:B:197:ASP:HB3	1:B:200:VAL:HG11	1.95	0.47
1:H:264:TYR:CE1	1:H:274:LYS:HB2	2.50	0.47
1:B:73:LEU:HD12	1:B:73:LEU:HA	1.49	0.47
1:D:125:ASN:ND2	3:D:511:HOH:O	2.47	0.47
1:A:17:VAL:HG13	2:A:401:95P:BR1	2.70	0.46
1:A:55:HIS:HA	1:A:59:ILE:HG22	1.97	0.46
1:E:216:ALA:HA	1:E:219:PHE:CD2	2.51	0.46
1:F:28:GLY:HA2	2:F:401:95P:O17	2.14	0.46
1:C:142:LYS:HA	1:C:142:LYS:HD2	1.71	0.46
1:F:119:PRO:HA	1:F:134:THR:HG23	1.98	0.46
1:H:89:PHE:CD1	1:H:109:LYS:HA	2.51	0.46
1:B:225:GLU:O	1:B:229:ARG:HG3	2.16	0.46
1:E:17:VAL:HG13	2:E:401:95P:BR1	2.71	0.46
1:F:224:THR:O	1:F:228:GLN:HG2	2.15	0.46
1:C:166:LEU:O	1:C:171:THR:HA	2.16	0.46
1:F:9:ASP:OD2	1:H:109:LYS:NZ	2.45	0.46
1:H:21:GLY:HA2	2:H:401:95P:S3	2.56	0.46
1:G:184:PHE:HA	1:G:194:ILE:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:297:THR:HB	1:E:315:PRO:HG2	1.97	0.45
1:C:233:PRO:HG2	1:C:237:SER:O	2.16	0.45
1:A:142:LYS:HD3	1:A:142:LYS:HA	1.71	0.45
1:D:270:SER:OG	1:D:274:LYS:HE3	2.16	0.45
1:F:184:PHE:HB3	1:F:193:PHE:HB3	1.99	0.45
1:E:218:ASP:O	1:E:267:ASN:HB2	2.17	0.45
1:F:209:TYR:CZ	1:F:242:ALA:HB2	2.52	0.45
1:F:229:ARG:HE	1:F:234:PRO:HG3	1.82	0.45
1:H:155:PRO:HB3	1:H:306:ILE:HD13	1.97	0.45
1:G:120:LEU:HG	1:G:123:SER:HB3	1.99	0.45
1:H:277:LEU:HD23	1:H:307:PRO:HB3	1.99	0.45
1:A:235:ASP:C	1:A:237:SER:H	2.19	0.44
1:G:96:SER:HB2	1:G:117:PHE:CZ	2.52	0.44
1:B:28:GLY:HA2	2:B:401:95P:C9	2.46	0.44
1:D:274:LYS:O	1:D:313:ARG:NH1	2.51	0.44
1:A:226:TYR:CE1	1:A:230:LYS:HE3	2.52	0.44
1:B:181:VAL:HB	1:B:200:VAL:HG22	1.99	0.44
1:B:288:MET:HG3	1:B:318:LEU:HD13	1.99	0.44
1:F:22:ARG:NH1	1:H:32:GLN:OE1	2.50	0.44
1:D:28:GLY:HA2	2:D:401:95P:C9	2.48	0.44
1:A:225:GLU:OE1	1:A:334:HIS:NE2	2.37	0.44
1:B:141:LYS:HE2	1:B:141:LYS:HB3	1.70	0.44
1:H:142:LYS:HA	1:H:142:LYS:HD2	1.79	0.44
1:E:329:LYS:HE3	1:E:329:LYS:HB2	1.77	0.43
1:F:15:ARG:NH2	1:H:84:MET:O	2.51	0.43
1:E:299:LYS:HG3	1:E:300:GLU:HG2	2.00	0.43
1:B:144:THR:O	1:B:144:THR:OG1	2.35	0.43
1:B:283:PRO:O	1:B:287:VAL:HG23	2.19	0.43
1:B:264:TYR:OH	1:B:274:LYS:HD3	2.18	0.43
1:C:28:GLY:HA2	2:C:401:95P:C9	2.49	0.43
1:C:89:PHE:CD2	1:C:109:LYS:HA	2.53	0.43
1:D:267:ASN:OD1	1:D:269:LYS:HB2	2.19	0.43
1:B:206:GLY:HA3	1:B:260:GLY:N	2.34	0.43
1:A:217:LYS:HA	1:B:232:PHE:CE1	2.53	0.43
1:E:142:LYS:HD3	1:E:142:LYS:HA	1.70	0.43
1:F:28:GLY:HA2	2:F:401:95P:C9	2.48	0.43
2:C:401:95P:O17	2:C:401:95P:S8	2.77	0.43
1:E:72:LYS:HB3	1:E:74:ASP:OD1	2.18	0.43
1:E:109:LYS:NZ	1:G:9:ASP:OD2	2.30	0.43
1:D:72:LYS:HE3	1:D:72:LYS:HB2	1.64	0.43
1:E:120:LEU:HG	1:E:123:SER:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:37:LEU:HD21	1:F:136:PHE:CD2	2.53	0.42
1:G:274:LYS:NZ	3:G:509:HOH:O	2.52	0.42
1:E:34:LEU:HD23	1:E:34:LEU:HA	1.65	0.42
1:A:277:LEU:HA	1:A:281:CYS:HB2	2.01	0.42
1:E:294:MET:HE2	1:E:321:PRO:HA	2.02	0.42
1:D:73:LEU:HA	1:D:73:LEU:HD12	1.72	0.42
1:H:231:LYS:O	1:H:239:PRO:HB3	2.19	0.42
1:B:184:PHE:HB3	1:B:193:PHE:HB3	2.00	0.42
1:H:293:GLY:HA2	1:H:321:PRO:HD3	2.01	0.42
1:G:288:MET:HG3	1:G:318:LEU:HD13	2.01	0.42
1:E:89:PHE:HA	1:E:109:LYS:O	2.20	0.42
1:A:54:ALA:HB2	1:B:194:ILE:HD11	2.01	0.41
1:C:290:LYS:HA	1:C:290:LYS:HD3	1.86	0.41
1:H:144:THR:HG22	1:H:144:THR:O	2.20	0.41
1:A:73:LEU:HD23	1:A:126:ILE:HD13	2.03	0.41
1:A:71:LYS:HD3	1:A:75:VAL:HG12	2.02	0.41
1:B:13:LEU:HA	1:B:184:PHE:CE1	2.55	0.41
1:E:110:ARG:NH2	3:E:506:HOH:O	2.48	0.41
1:E:33:LEU:HD12	1:E:33:LEU:HA	1.81	0.41
1:A:18:MET:HG2	2:A:401:95P:BR2	2.74	0.41
1:C:106:GLU:OE2	1:C:109:LYS:HE2	2.20	0.41
1:D:184:PHE:HB3	1:D:193:PHE:HB3	2.02	0.41
1:E:74:ASP:OD1	1:E:74:ASP:N	2.52	0.41
1:F:267:ASN:OD1	1:F:270:SER:N	2.53	0.41
1:G:184:PHE:HB3	1:G:193:PHE:HB3	2.01	0.41
1:D:248:MET:HE3	1:D:275:LEU:HD13	2.02	0.41
1:E:205:LYS:HE2	1:E:240:TYR:OH	2.21	0.41
1:E:55:HIS:HA	1:E:59:ILE:HG22	2.03	0.41
1:H:96:SER:HB2	1:H:117:PHE:CZ	2.56	0.41
1:G:183:CYS:HB2	1:G:197:ASP:HB2	2.03	0.41
1:G:225:GLU:O	1:G:229:ARG:HG3	2.20	0.41
1:B:74:ASP:OD1	1:B:74:ASP:N	2.47	0.41
1:H:28:GLY:HA2	2:H:401:95P:O17	2.21	0.41
1:F:208:ILE:HA	1:F:241:GLY:O	2.21	0.41
1:F:278:LEU:HD23	1:F:278:LEU:HA	1.84	0.41
1:D:209:TYR:OH	1:D:213:GLU:OE2	2.21	0.41
1:F:283:PRO:O	1:F:287:VAL:HG23	2.21	0.41
1:C:288:MET:HG3	1:C:318:LEU:HD13	2.03	0.40
1:G:278:LEU:HD12	1:G:310:ILE:HA	2.03	0.40
1:A:184:PHE:HB3	1:A:193:PHE:HB3	2.04	0.40
1:B:159:LEU:HD23	1:B:159:LEU:HA	1.94	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LEU:HD12	1:A:73:LEU:HA	1.89	0.40
1:E:141:LYS:HE2	1:E:151:ASP:OD2	2.22	0.40
1:G:288:MET:HE3	1:G:293:GLY:HA3	2.03	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	316/338 (94%)	309 (98%)	6 (2%)	1 (0%)	41	61
1	B	314/338 (93%)	304 (97%)	9 (3%)	1 (0%)	41	61
1	C	313/338 (93%)	305 (97%)	8 (3%)	0	100	100
1	D	314/338 (93%)	308 (98%)	6 (2%)	0	100	100
1	E	316/338 (94%)	309 (98%)	7 (2%)	0	100	100
1	F	314/338 (93%)	305 (97%)	9 (3%)	0	100	100
1	G	314/338 (93%)	306 (98%)	8 (2%)	0	100	100
1	H	313/338 (93%)	306 (98%)	7 (2%)	0	100	100
All	All	2514/2704 (93%)	2452 (98%)	60 (2%)	2 (0%)	51	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	ASN
1	B	145	ASP

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	267/281 (95%)	264 (99%)	3 (1%)	73	89
1	B	265/281 (94%)	260 (98%)	5 (2%)	57	80
1	C	264/281 (94%)	258 (98%)	6 (2%)	50	76
1	D	265/281 (94%)	259 (98%)	6 (2%)	50	76
1	E	267/281 (95%)	262 (98%)	5 (2%)	57	80
1	F	265/281 (94%)	259 (98%)	6 (2%)	50	76
1	G	265/281 (94%)	259 (98%)	6 (2%)	50	76
1	H	264/281 (94%)	257 (97%)	7 (3%)	44	71
All	All	2122/2248 (94%)	2078 (98%)	44 (2%)	55	78

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASP
1	A	110	ARG
1	A	316	VAL
1	B	9	ASP
1	B	38	CYS
1	B	72	LYS
1	B	144	THR
1	B	178	ASP
1	C	9	ASP
1	C	38	CYS
1	C	178	ASP
1	C	200	VAL
1	C	207	LYS
1	C	316	VAL
1	D	9	ASP
1	D	38	CYS
1	D	72	LYS
1	D	178	ASP
1	D	264	TYR

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Mol	Chain	Res	Type
1	D	316	VAL
1	E	9	ASP
1	E	38	CYS
1	E	178	ASP
1	E	197	ASP
1	E	211	LEU
1	F	9	ASP
1	F	38	CYS
1	F	140[A]	ARG
1	F	140[B]	ARG
1	F	211	LEU
1	F	264	TYR
1	G	9	ASP
1	G	38	CYS
1	G	178	ASP
1	G	211	LEU
1	G	264	TYR
1	G	316	VAL
1	H	9	ASP
1	H	38	CYS
1	H	178	ASP
1	H	218	ASP
1	H	274	LYS
1	H	275	LEU
1	H	316	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

8 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	95P	B	401	-	14,21,21	1.18	1 (7%)	12,31,31	1.86	2 (16%)
2	95P	D	401	-	14,21,21	1.22	1 (7%)	12,31,31	1.36	2 (16%)
2	95P	F	401	-	14,21,21	1.23	1 (7%)	12,31,31	1.15	2 (16%)
2	95P	H	401	-	14,21,21	1.21	1 (7%)	12,31,31	1.40	2 (16%)
2	95P	G	401	-	14,21,21	1.50	3 (21%)	12,31,31	1.19	1 (8%)
2	95P	A	401	-	14,21,21	1.39	2 (14%)	12,31,31	1.62	3 (25%)
2	95P	C	401	-	14,21,21	1.40	1 (7%)	12,31,31	1.98	2 (16%)
2	95P	E	401	-	14,21,21	1.52	1 (7%)	12,31,31	1.64	2 (16%)

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	95P	B	401	-	-	1/7/15/15	0/2/2/2
2	95P	D	401	-	-	1/7/15/15	0/2/2/2
2	95P	F	401	-	-	1/7/15/15	0/2/2/2
2	95P	H	401	-	-	1/7/15/15	0/2/2/2
2	95P	G	401	-	-	1/7/15/15	0/2/2/2
2	95P	A	401	-	-	2/7/15/15	0/2/2/2
2	95P	C	401	-	-	1/7/15/15	0/2/2/2
2	95P	E	401	-	-	1/7/15/15	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	95P	C5-N11	4.39	1.44	1.36
2	C	401	95P	C2-S1	-3.58	1.71	1.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	95P	BR1-C10	-3.29	1.82	1.89
2	G	401	95P	C2-S1	-2.94	1.72	1.76
2	B	401	95P	C2-S1	-2.86	1.72	1.76
2	D	401	95P	C5-N11	2.85	1.41	1.36
2	G	401	95P	BR1-C10	-2.79	1.83	1.89
2	F	401	95P	C5-N11	2.79	1.41	1.36
2	H	401	95P	BR1-C10	-2.61	1.83	1.89
2	G	401	95P	C5-N11	2.46	1.41	1.36
2	A	401	95P	C2-S1	-2.28	1.73	1.76

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	95P	C2-S1-N6	5.62	113.84	107.27
2	B	401	95P	C2-S1-N6	5.46	113.65	107.27
2	E	401	95P	C2-S1-N6	4.34	112.34	107.27
2	H	401	95P	C2-S1-N6	3.63	111.51	107.27
2	D	401	95P	C2-S1-N6	3.24	111.06	107.27
2	A	401	95P	O16-S1-N6	3.23	115.83	106.74
2	G	401	95P	BR2-C13-C14	2.43	128.34	124.51
2	E	401	95P	BR2-C13-C14	2.37	128.24	124.51
2	C	401	95P	BR2-C13-C14	2.31	128.14	124.51
2	H	401	95P	BR2-C13-C14	2.26	128.07	124.51
2	A	401	95P	O15-S1-N6	-2.23	100.47	106.74
2	B	401	95P	BR2-C13-C14	2.19	127.96	124.51
2	F	401	95P	BR2-C13-C14	2.19	127.95	124.51
2	F	401	95P	C2-S1-N6	2.17	109.80	107.27
2	D	401	95P	BR2-C13-C14	2.14	127.88	124.51
2	A	401	95P	BR2-C13-C14	2.13	127.87	124.51

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	95P	N6-C9-N11-C5
2	D	401	95P	N6-C9-N11-C5
2	F	401	95P	N6-C9-N11-C5
2	H	401	95P	N6-C9-N11-C5
2	G	401	95P	N6-C9-N11-C5
2	A	401	95P	N6-C9-N11-C5
2	C	401	95P	N6-C9-N11-C5
2	E	401	95P	N6-C9-N11-C5

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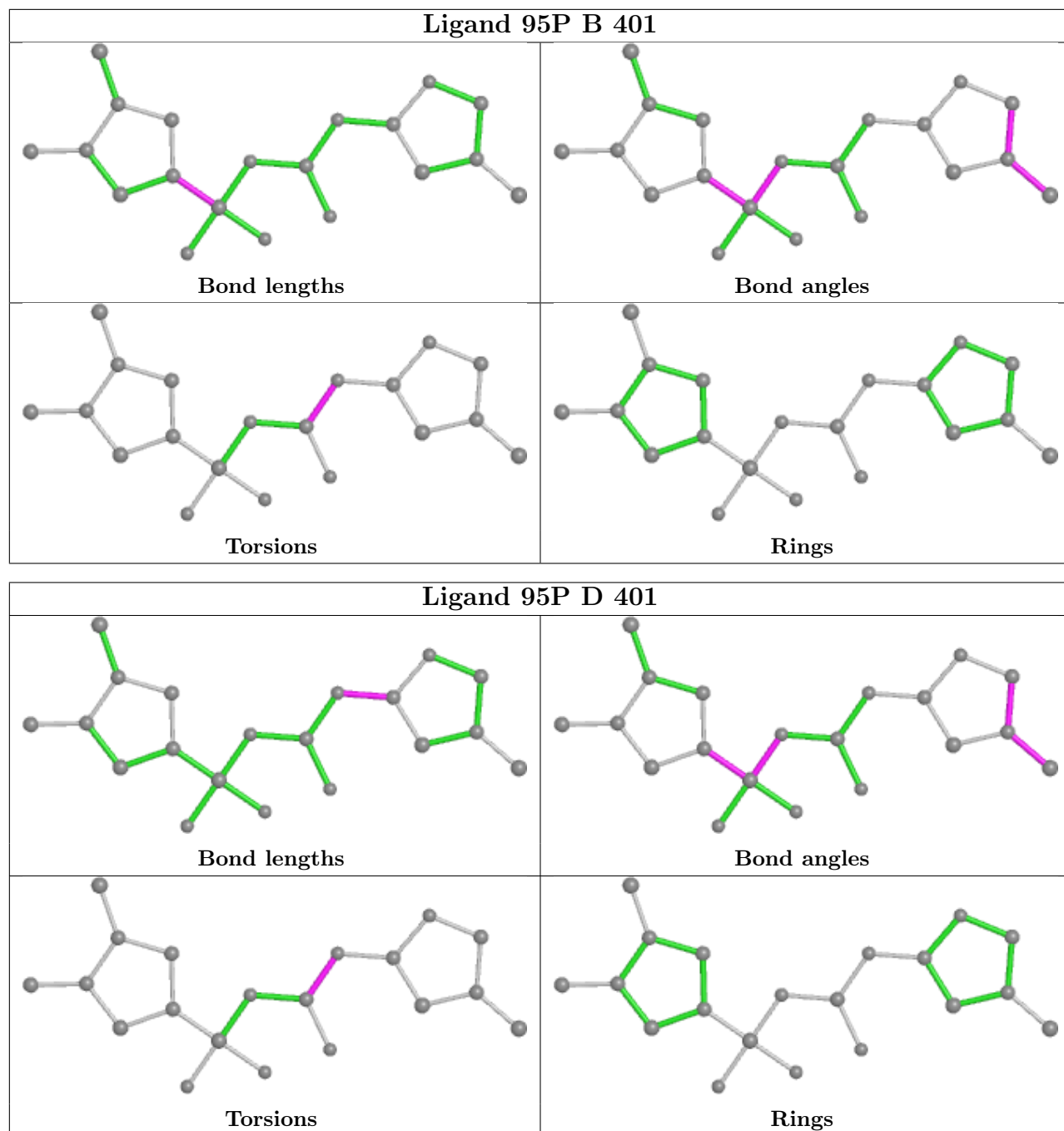
Mol	Chain	Res	Type	Atoms
2	A	401	95P	C9-N6-S1-C2

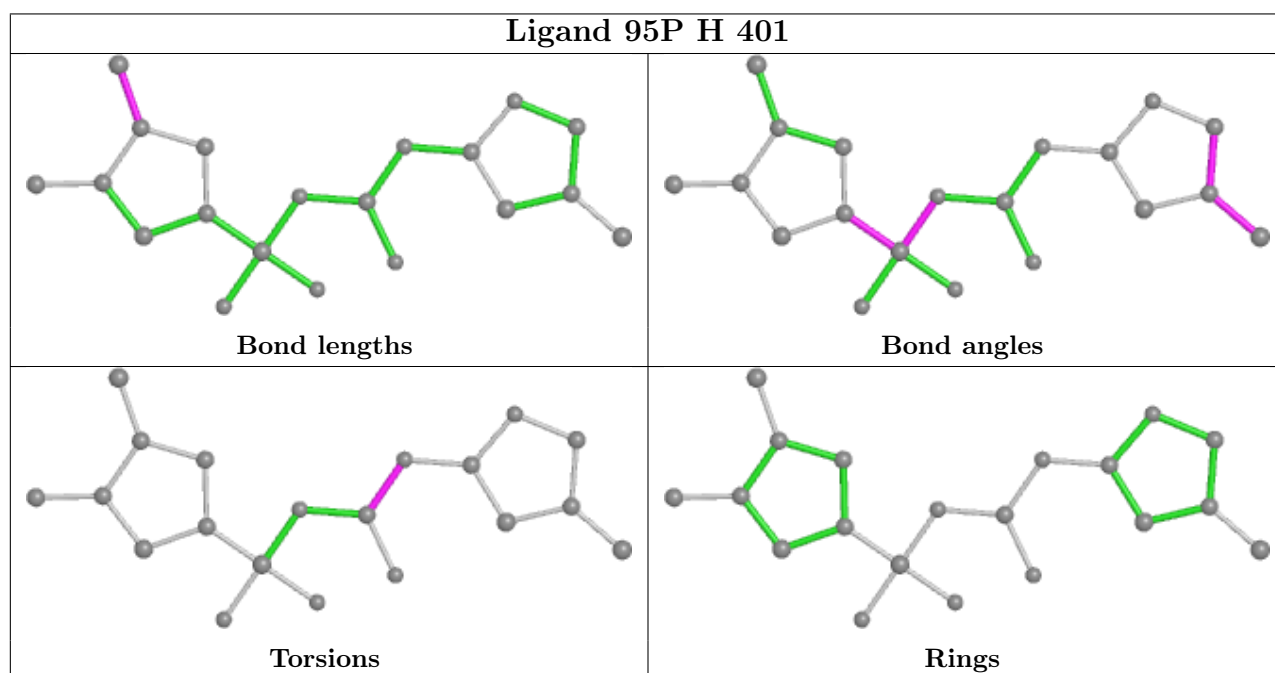
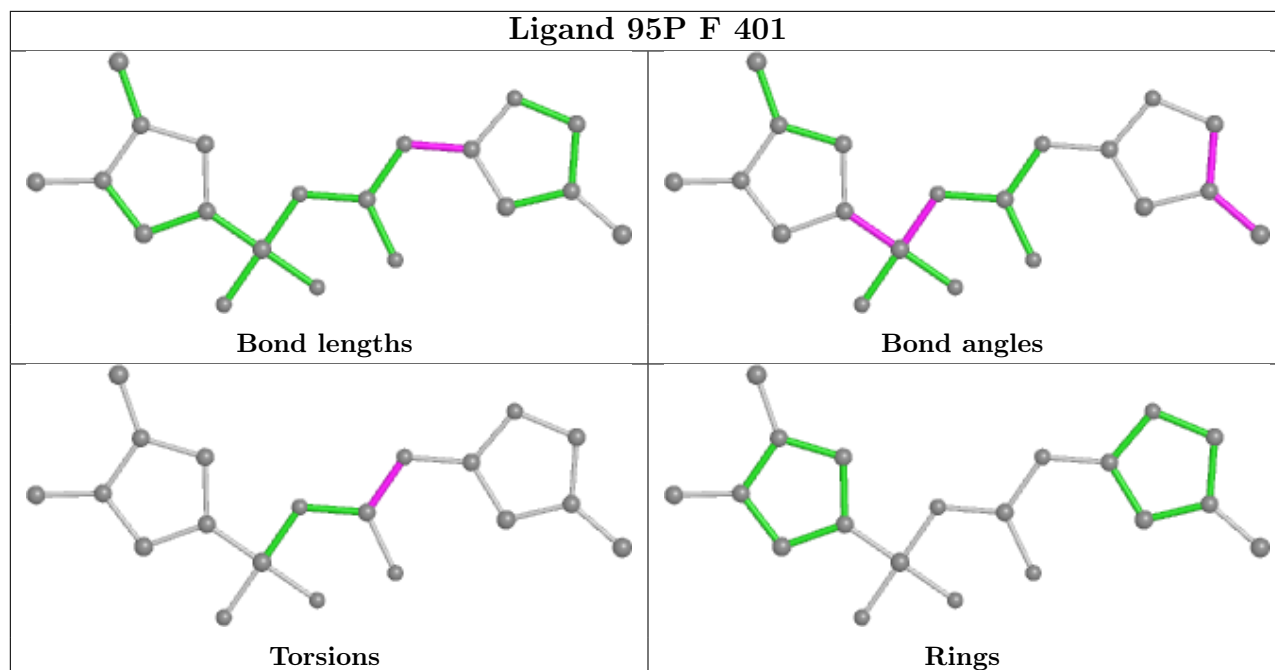
There are no ring outliers.

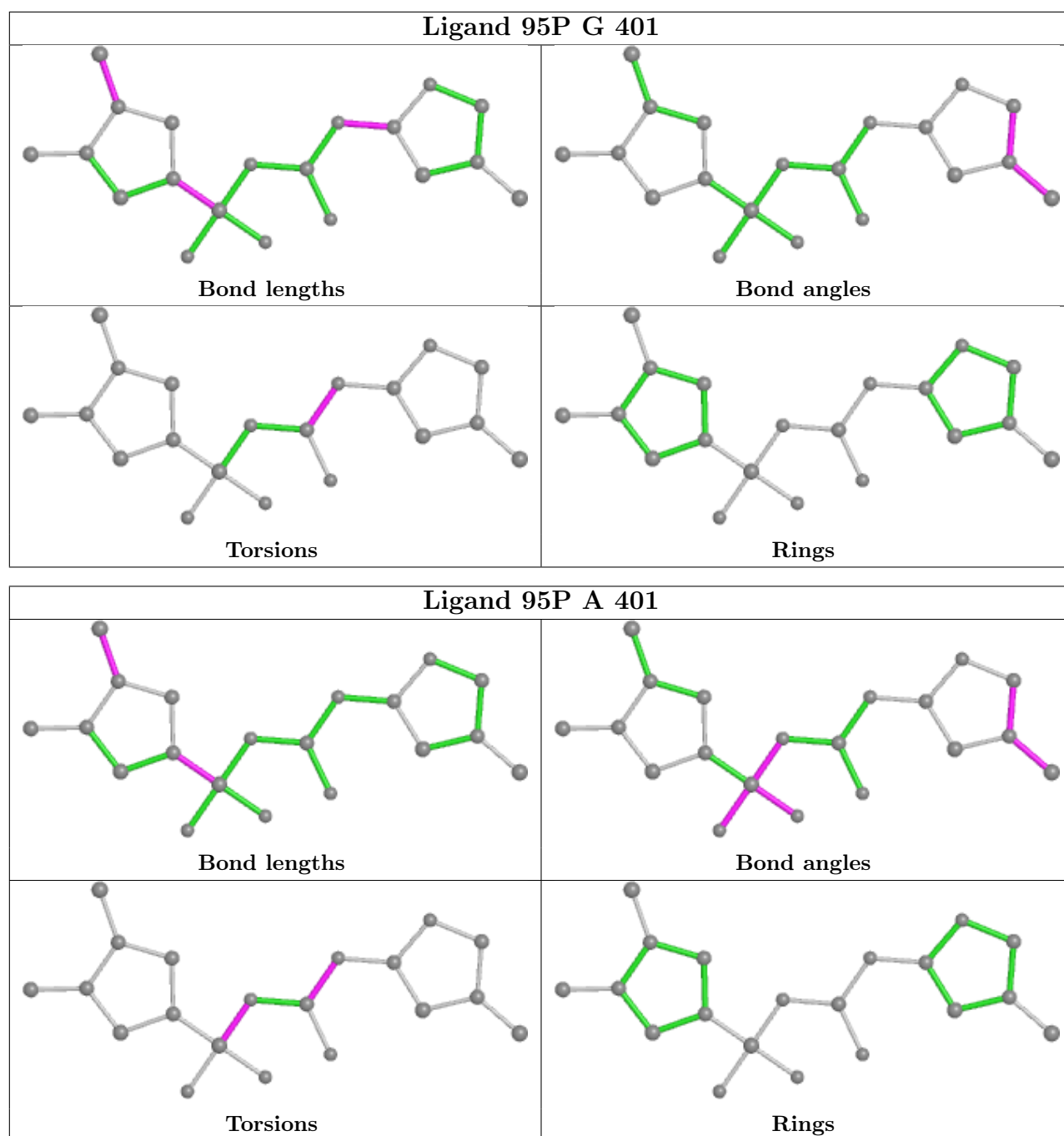
8 monomers are involved in 25 short contacts:

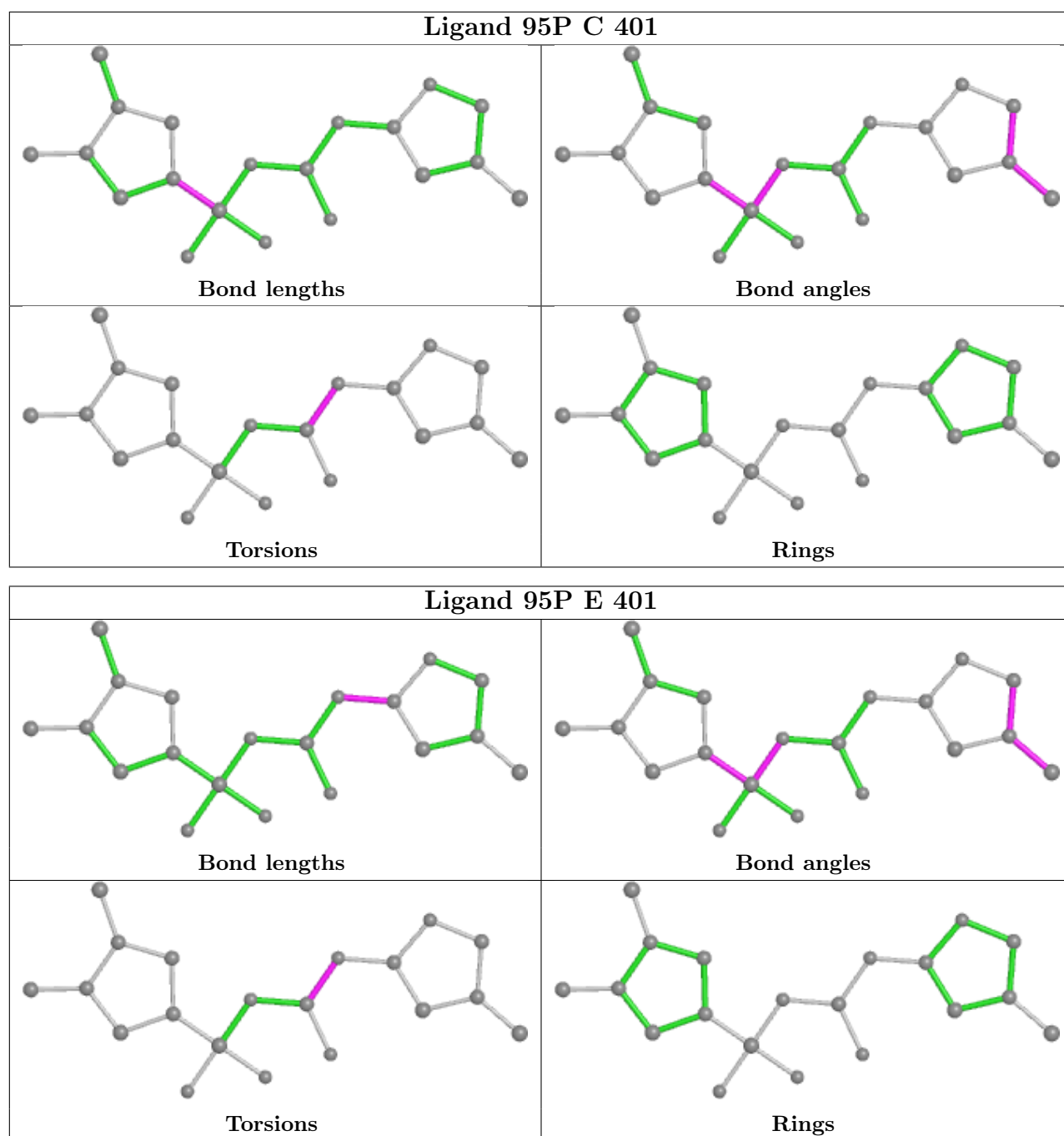
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	95P	4	0
2	D	401	95P	3	0
2	F	401	95P	3	0
2	H	401	95P	5	0
2	G	401	95P	2	0
2	A	401	95P	3	0
2	C	401	95P	3	0
2	E	401	95P	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	319/338 (94%)	-0.08	5 (1%) 72 74	26, 48, 83, 102	0
1	B	317/338 (93%)	0.18	22 (6%) 16 17	28, 49, 89, 124	0
1	C	317/338 (93%)	-0.12	6 (1%) 66 69	27, 42, 70, 106	0
1	D	317/338 (93%)	-0.21	5 (1%) 72 74	25, 41, 69, 112	0
1	E	319/338 (94%)	0.13	20 (6%) 20 21	28, 53, 88, 126	0
1	F	317/338 (93%)	0.19	18 (5%) 23 25	32, 53, 90, 138	0
1	G	317/338 (93%)	-0.13	4 (1%) 77 79	25, 44, 70, 105	0
1	H	317/338 (93%)	-0.07	10 (3%) 47 51	26, 45, 77, 123	0
All	All	2540/2704 (93%)	-0.01	90 (3%) 44 47	25, 47, 81, 138	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	70	VAL	13.5
1	F	268	LYS	6.9
1	H	269	LYS	6.2
1	E	236	ASN	4.9
1	E	234	PRO	4.8
1	D	236	ASN	4.5
1	F	271	PRO	4.5
1	F	236	ASN	4.4
1	E	235	ASP	4.2
1	H	236	ASN	4.1
1	B	268	LYS	4.0
1	G	25	ARG	3.9
1	A	144	THR	3.9
1	F	143	SER	3.8
1	H	268	LYS	3.8
1	B	269	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	106	GLU	3.6
1	E	335	SER	3.6
1	A	268	LYS	3.6
1	B	332	GLU	3.5
1	F	272	ASN	3.4
1	C	126	ILE	3.3
1	E	268	LYS	3.3
1	E	269	LYS	3.3
1	A	70	VAL	3.2
1	F	107	PRO	3.2
1	C	236	ASN	3.2
1	A	234	PRO	3.2
1	B	144	THR	3.2
1	D	38	CYS	3.2
1	E	237	SER	3.2
1	C	25	ARG	3.1
1	B	267	ASN	3.1
1	D	335	SER	3.1
1	E	238	ALA	3.1
1	H	234	PRO	3.0
1	D	25	ARG	3.0
1	C	38	CYS	3.0
1	B	25	ARG	2.9
1	F	267	ASN	2.9
1	B	236	ASN	2.8
1	F	144	THR	2.8
1	F	25	ARG	2.7
1	B	135	ILE	2.7
1	H	271	PRO	2.6
1	E	108	GLU	2.6
1	G	332	GLU	2.6
1	B	163	GLY	2.6
1	B	56	LEU	2.6
1	F	142	LYS	2.6
1	E	333	LYS	2.5
1	F	222	ALA	2.5
1	B	283	PRO	2.5
1	B	143	SER	2.4
1	A	145	ASP	2.4
1	B	125	ASN	2.4
1	E	267	ASN	2.4
1	B	24	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	229	ARG	2.4
1	H	25	ARG	2.4
1	G	38	CYS	2.4
1	H	267	ASN	2.3
1	F	179	CYS	2.3
1	H	106	GLU	2.3
1	B	271	PRO	2.3
1	F	145	ASP	2.3
1	C	23	LYS	2.3
1	C	329	LYS	2.3
1	F	38	CYS	2.3
1	B	237	SER	2.3
1	H	108	GLU	2.3
1	F	109	LYS	2.2
1	B	329	LYS	2.2
1	B	272	ASN	2.2
1	B	284	MET	2.1
1	B	266	ALA	2.1
1	E	233	PRO	2.1
1	E	25	ARG	2.1
1	F	266	ALA	2.1
1	E	71	LYS	2.1
1	D	37	LEU	2.1
1	F	269	LYS	2.1
1	E	116	CYS	2.1
1	F	136	PHE	2.1
1	E	272	ASN	2.1
1	E	271	PRO	2.1
1	H	38	CYS	2.1
1	E	106	GLU	2.0
1	B	238	ALA	2.0
1	G	136	PHE	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](#)

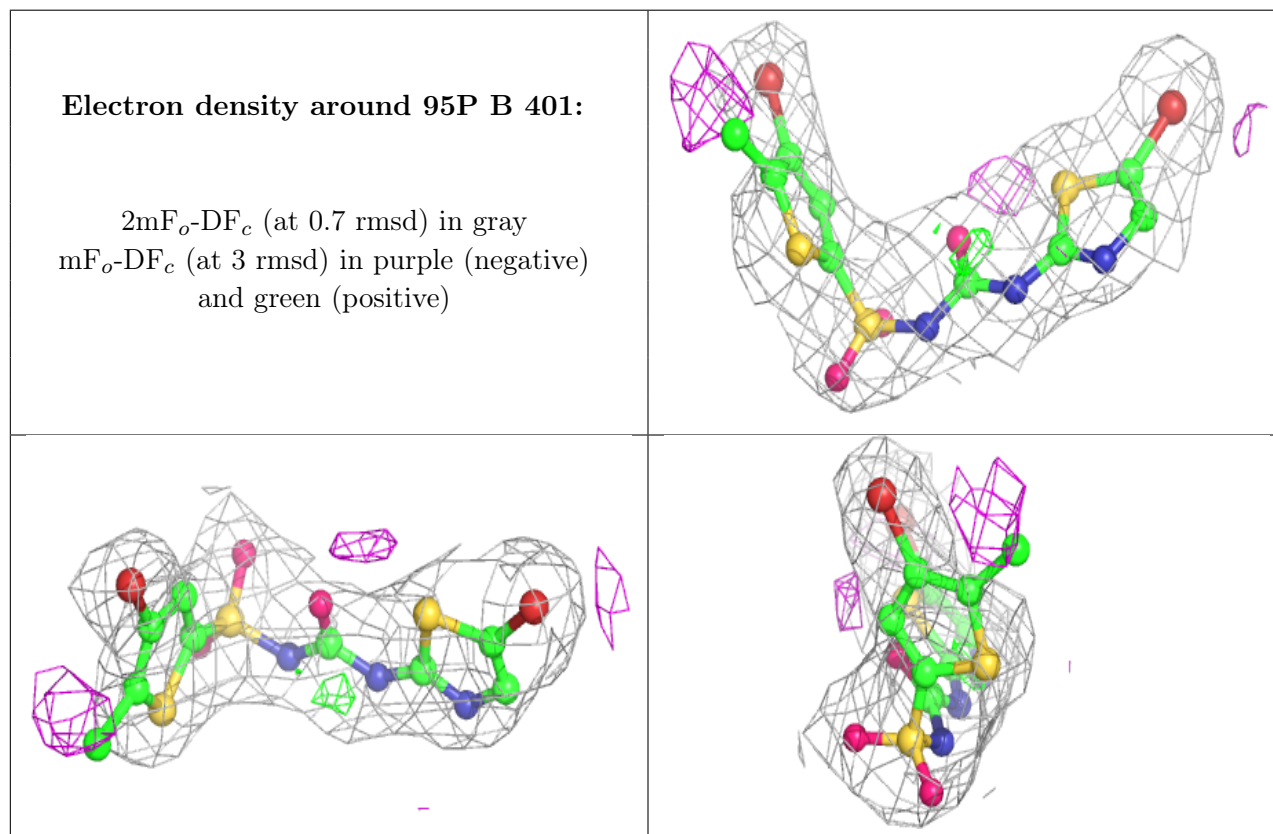
There are no monosaccharides in this entry.

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.

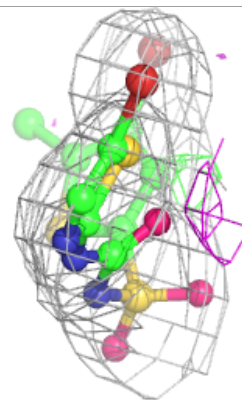
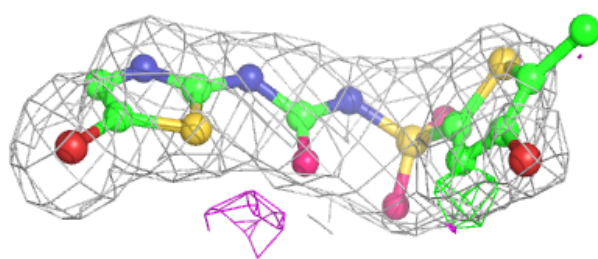
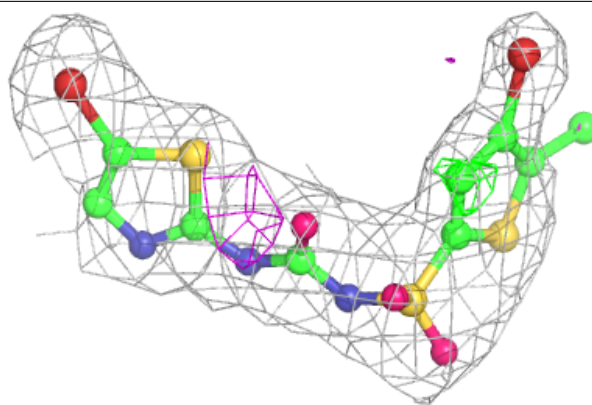
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	95P	B	401	20/20	0.93	0.13	54,70,98,124	2
2	95P	D	401	20/20	0.93	0.15	57,68,103,123	2
2	95P	F	401	20/20	0.93	0.14	57,69,111,131	2
2	95P	H	401	20/20	0.93	0.15	49,75,101,119	2
2	95P	E	401	20/20	0.93	0.13	58,70,80,110	2
2	95P	C	401	20/20	0.94	0.12	56,66,79,100	2
2	95P	A	401	20/20	0.95	0.14	51,65,81,86	2
2	95P	G	401	20/20	0.96	0.11	42,62,89,109	2

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.

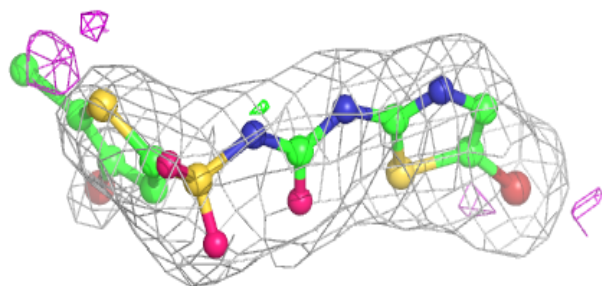
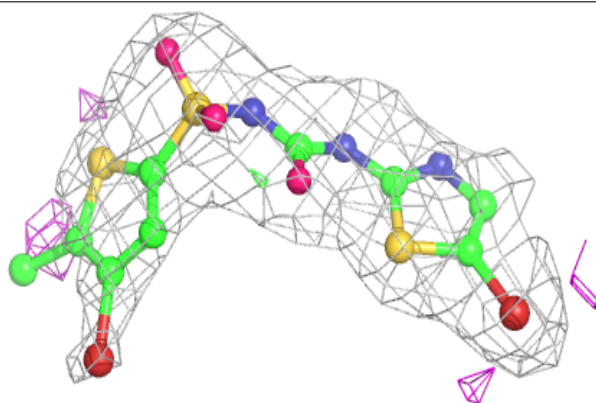


Electron density around 95P D 401:

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

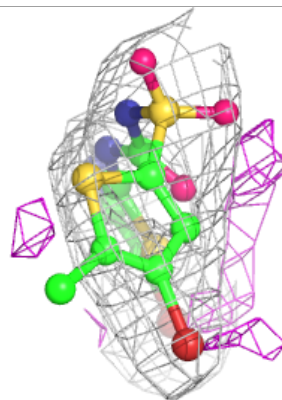
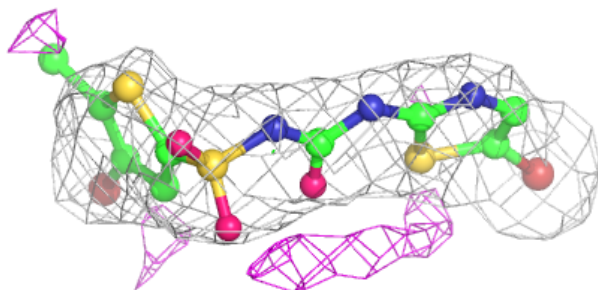
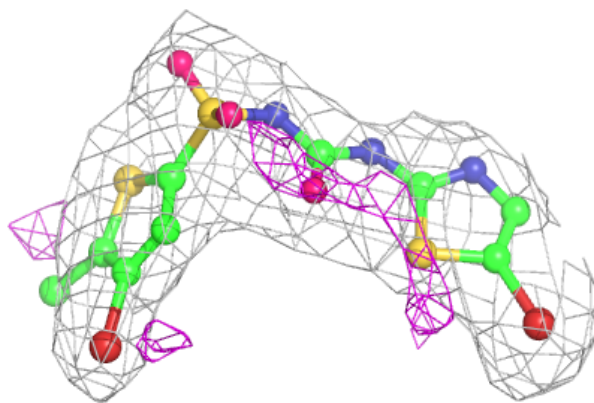
**Electron density around 95P F 401:**

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

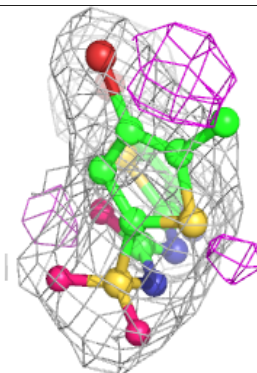
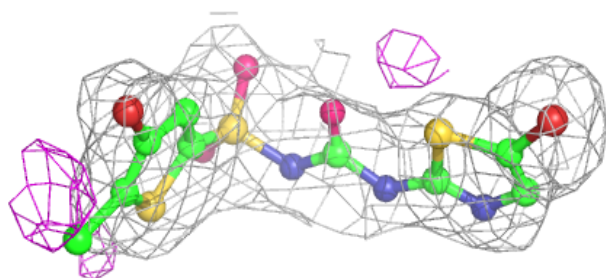
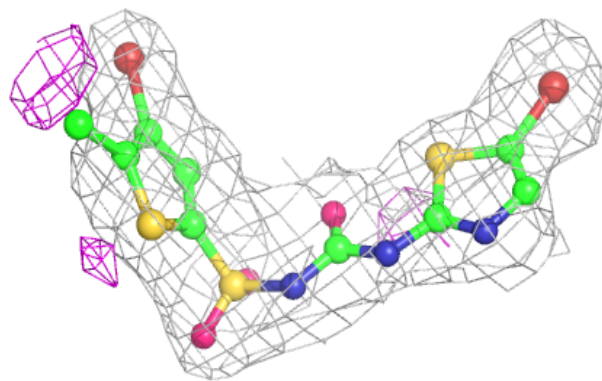


Electron density around 95P H 401:

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

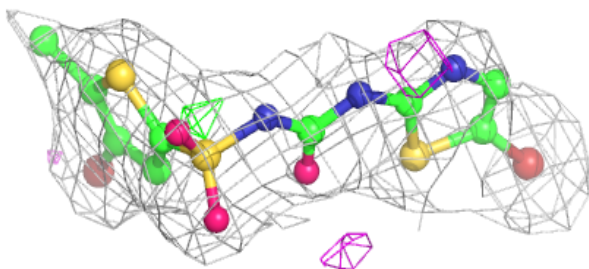
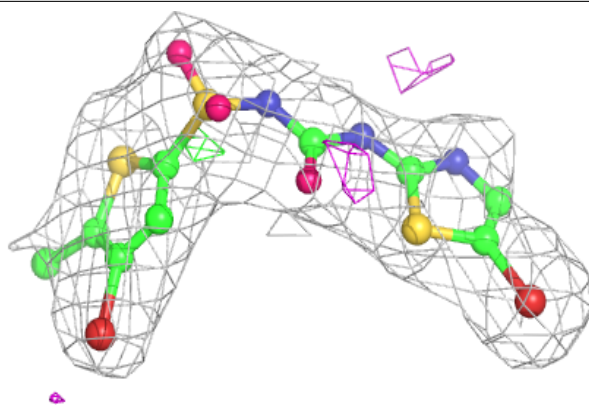
**Electron density around 95P E 401:**

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

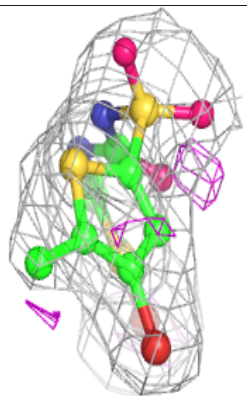
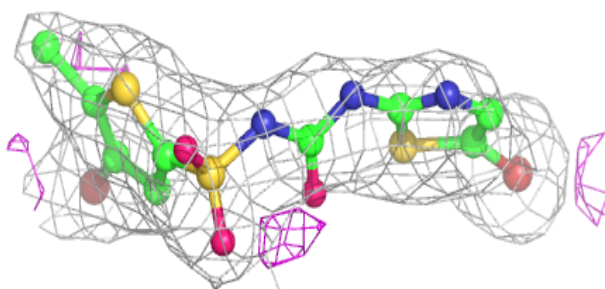
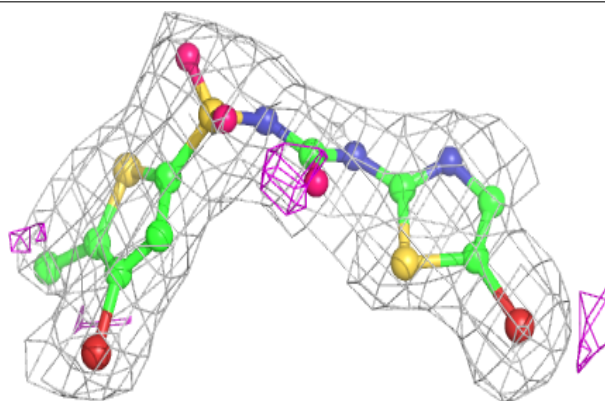


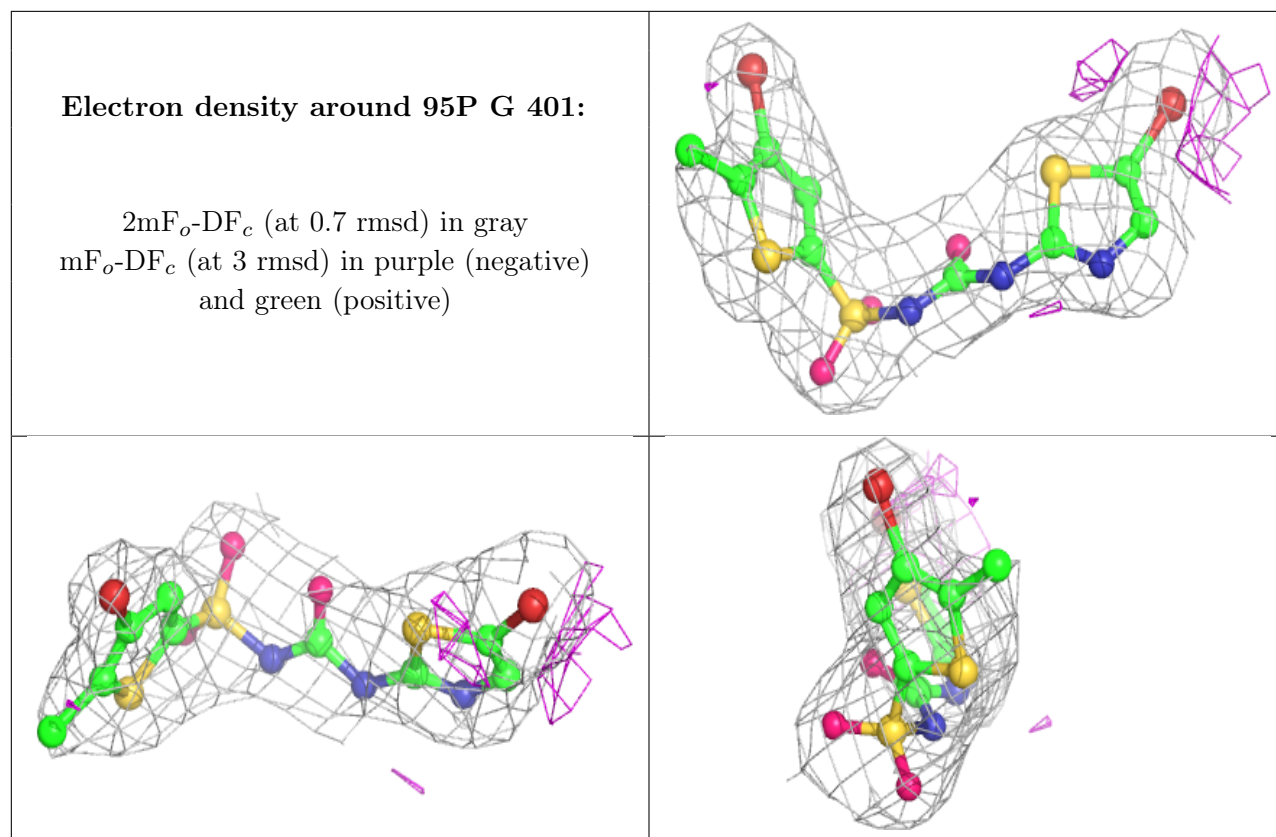
Electron density around 95P C 401:

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

**Electron density around 95P A 401:**

$2mF_o-DF_c$ (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.