



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 04:50 pm BST

PDB ID : 1B12
Title : CRYSTAL STRUCTURE OF TYPE 1 SIGNAL PEPTIDASE FROM ES-
CHERICHIA COLI IN COMPLEX WITH A BETA-LACTAM INHIBITOR
Authors : Paetzel, M.; Dalbey, R.; Strynadka, N.C.J.
Deposited on : 1999-11-24
Resolution : 1.95 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

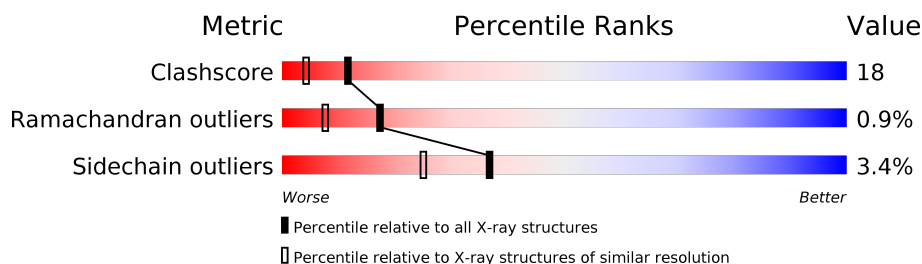
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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))
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)

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 on 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	248	78% 15% . . .
1	B	248	69% 15% . 15%
1	C	248	68% 19% . . 9%
1	D	248	69% 17% . 10%

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	1PN	A	1001	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	1PN	C	1001	-	-	X	-
2	1PN	D	1001	-	-	X	-

2 Entry composition [i](#)

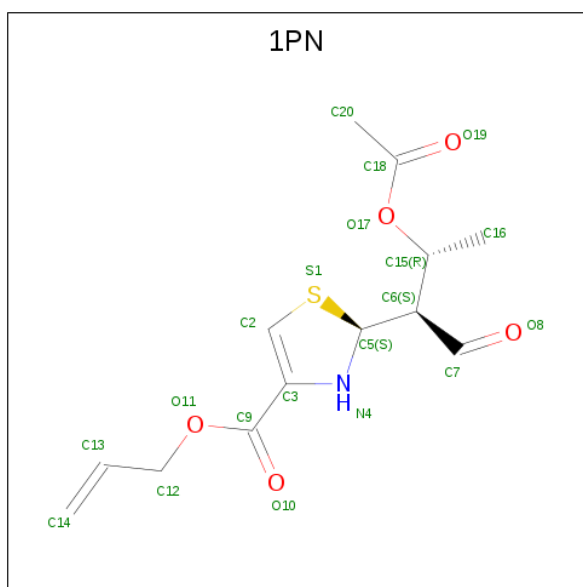
There are 4 unique types of molecules in this entry. The entry contains 7646 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SIGNAL PEPTIDASE I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1887	1204	319	356	8			
1	B	211	Total	C	N	O	S	0	0	0
			1679	1077	280	314	8			
1	C	226	Total	C	N	O	S	0	0	0
			1775	1135	295	337	8			
1	D	222	Total	C	N	O	S	0	0	0
			1772	1143	293	330	6			

- Molecule 2 is prop-2-en-1-yl (2S)-2-[(2S,3R)-3-(acetyloxy)-1-oxobutan-2-yl]-2,3-dihydro-1,3-thiazole-4-carboxylate (three-letter code: 1PN) (formula: C₁₃H₁₇NO₅S).



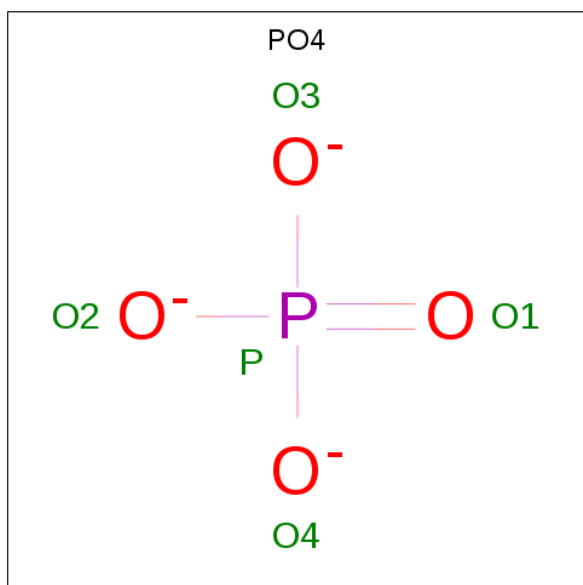
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	13	1	5	1		
2	B	1	Total	C	N	O	S	0	0
			20	13	1	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			20	13	1	5	1		
2	D	1	Total	C	N	O	S	0	0
			20	13	1	5	1		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	129	Total	O	0	0
			129	129		
4	B	111	Total	O	0	0
			111	111		
4	C	122	Total	O	0	0
			122	122		
4	D	86	Total	O	0	0
			86	86		

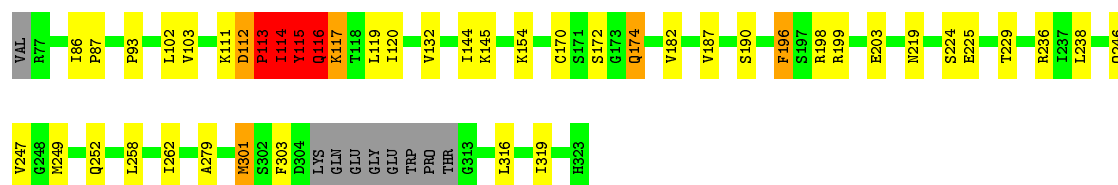
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

Note EDS was not executed.

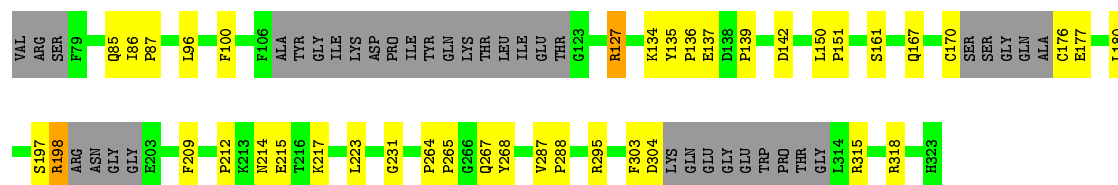
• Molecule 1: SIGNAL PEPTIDASE I

Chain A: 



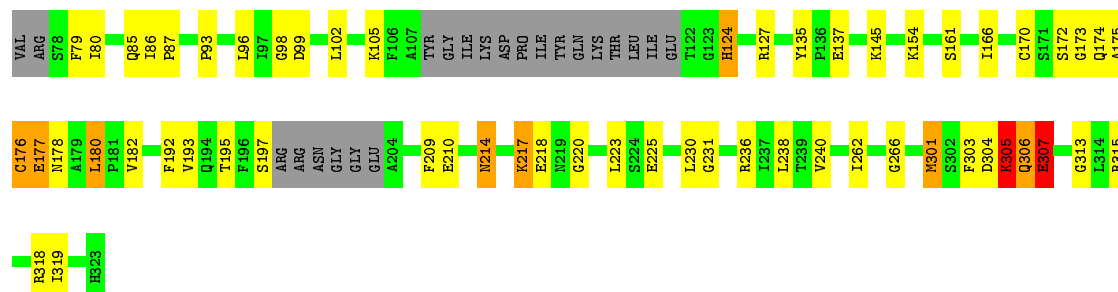
• Molecule 1: SIGNAL PEPTIDASE I

Chain B: 



• Molecule 1: SIGNAL PEPTIDASE I

Chain C: 



• Molecule 1: SIGNAL PEPTIDASE I

Chain D: 

VAL	GLN
R77	ALA
S78	CYS
F79	GLU
I80	ASN
Y81	A179
E82	L180
I86	P181
P87	V187
P93	D191
F100	S197
K105	ARG
F106	ASN
I110	GLY
K111	GLY
D112	GLU
P113	ALA
I114	T205
Y115	S206
Q116	K213
K117	L223
I118	S224
L119	G231
I120	L238
E121	T239
P125	V240
R126	Q246
R127	Q267
I130	P288
V131	R295
V132	F303
Y135	D304
P136	LYS
E137	GLN
D138	GLU
P139	GLY
K140	GLU
L141	GLY
I144	GLU
K145	GLY
D153	TRP
I166	PRO
Q167	THR
P168	G313
GLY	I322
CYS	H323
SER	
SER	
GLY	

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.70 Å 113.20 Å 99.20 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.50 – 1.95	Depositor
% Data completeness (in resolution range)	96.8 (19.50-1.95)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.11	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.220 , 0.246	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7646	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 1PN

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	0.81	5/1935 (0.3%)	1.06	13/2622 (0.5%)
1	B	0.48	0/1721	0.77	0/2330
1	C	0.65	1/1822 (0.1%)	0.90	8/2472 (0.3%)
1	D	0.54	1/1818 (0.1%)	0.76	1/2464 (0.0%)
All	All	0.63	7/7296 (0.1%)	0.88	22/9888 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	D	0	2
All	All	0	5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114	ILE	C-N	-17.71	0.93	1.34
1	C	307	GLU	N-CA	16.00	1.78	1.46
1	A	115	TYR	C-N	-13.86	1.02	1.34
1	D	116	GLN	N-CA	10.98	1.68	1.46
1	A	114	ILE	CA-C	-9.01	1.29	1.52
1	A	117	LYS	N-CA	6.03	1.58	1.46
1	A	115	TYR	CA-C	-5.81	1.37	1.52

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	TYR	N-CA-C	16.08	154.41	111.00
1	A	114	ILE	O-C-N	-15.77	97.46	122.70
1	A	115	TYR	N-CA-CB	-12.23	88.58	110.60
1	A	117	LYS	N-CA-CB	11.77	131.79	110.60
1	A	114	ILE	CA-C-N	10.58	140.47	117.20
1	C	306	GLN	C-N-CA	-9.01	99.17	121.70
1	A	117	LYS	N-CA-C	-8.62	87.73	111.00
1	A	113	PRO	O-C-N	8.60	136.45	122.70
1	D	116	GLN	N-CA-CB	8.55	125.99	110.60
1	A	116	GLN	N-CA-C	8.01	132.63	111.00
1	A	116	GLN	C-N-CA	7.71	140.98	121.70
1	C	307	GLU	N-CA-C	-7.23	91.47	111.00
1	C	175	ALA	N-CA-C	-6.58	93.23	111.00
1	C	306	GLN	O-C-N	-6.49	112.31	122.70
1	A	113	PRO	CB-CA-C	6.46	128.14	112.00
1	A	116	GLN	N-CA-CB	-6.19	99.45	110.60
1	C	180	LEU	N-CA-C	-5.95	94.95	111.00
1	A	115	TYR	CA-CB-CG	5.78	124.39	113.40
1	C	176	CYS	C-N-CA	5.59	135.67	121.70
1	C	176	CYS	CA-C-N	-5.32	105.50	117.20
1	C	176	CYS	N-CA-C	5.21	125.07	111.00
1	A	113	PRO	CA-C-N	-5.21	105.74	117.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	PRO	Peptide
1	A	114	ILE	Mainchain
1	A	116	GLN	Peptide
1	D	115	TYR	Sidechain,Peptide

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	1887	0	1843	62	0
1	B	1679	0	1635	47	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1775	0	1718	82	6
1	D	1772	0	1745	54	8
2	A	20	0	16	8	0
2	B	20	0	16	6	0
2	C	20	0	16	8	0
2	D	20	0	16	8	0
3	B	5	0	0	0	0
4	A	129	0	0	2	1
4	B	111	0	0	2	0
4	C	122	0	0	4	2
4	D	86	0	0	3	0
All	All	7646	0	7005	249	9

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 18.

All (249) 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:D:116:GLN:N	1:D:116:GLN:CA	1.68	1.53
1:C:307:GLU:N	1:C:307:GLU:CA	1.78	1.44
1:A:114:ILE:HD12	1:A:115:TYR:CB	1.63	1.27
1:C:177:GLU:HG3	1:C:178:ASN:H	1.11	1.11
1:A:87:PRO:HD2	2:A:1001:1PN:H20	1.37	1.05
1:A:114:ILE:HD12	1:A:115:TYR:HB2	1.04	1.03
1:A:112:ASP:O	1:A:114:ILE:HD13	1.62	0.98
1:B:87:PRO:HD2	2:B:1001:1PN:H20	1.42	0.98
1:D:87:PRO:HD2	2:D:1001:1PN:H20	1.41	0.98
1:C:96:LEU:HG	1:C:318:ARG:NH1	1.79	0.96
1:A:114:ILE:CD1	1:A:115:TYR:HB2	1.94	0.96
1:C:304:ASP:HB3	1:C:307:GLU:HB2	1.47	0.96
1:C:98:GLY:HA3	1:C:305:LYS:NZ	1.81	0.96
1:C:306:GLN:C	1:C:307:GLU:CA	2.36	0.94
1:C:99:ASP:OD2	1:C:318:ARG:NH1	2.01	0.93
1:A:301:MET:HE3	1:A:319:ILE:HD11	1.46	0.93
1:A:301:MET:CE	1:A:319:ILE:HD11	2.00	0.90
1:C:305:LYS:HD2	1:C:305:LYS:H	1.37	0.89
1:C:306:GLN:O	1:C:307:GLU:CA	2.21	0.88
1:C:182:VAL:HG22	1:C:230:LEU:HD23	1.54	0.87
1:B:198:ARG:CZ	1:B:198:ARG:H	1.91	0.84
1:C:98:GLY:HA3	1:C:305:LYS:HZ1	1.37	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ILE:HD12	1:A:115:TYR:HB3	1.59	0.83
1:C:98:GLY:CA	1:C:305:LYS:HZ1	1.92	0.83
1:C:174:GLN:HE21	1:C:176:CYS:CB	1.92	0.82
1:D:87:PRO:HD2	2:D:1001:1PN:C20	2.10	0.81
2:D:1001:1PN:H16	2:D:1001:1PN:H20A	1.62	0.81
1:A:87:PRO:HD2	2:A:1001:1PN:C20	2.11	0.80
1:B:198:ARG:HG3	1:B:198:ARG:HH11	1.45	0.80
1:D:115:TYR:C	1:D:116:GLN:CA	2.50	0.79
1:C:305:LYS:CD	1:C:305:LYS:H	1.96	0.78
1:A:112:ASP:HB2	1:A:117:LYS:O	1.83	0.78
1:D:112:ASP:OD2	1:D:119:LEU:HG	1.82	0.78
1:B:197:SER:HB2	1:B:198:ARG:NE	1.99	0.78
1:C:86:ILE:HG23	2:C:1001:1PN:H20A	1.66	0.78
1:C:307:GLU:N	1:C:307:GLU:C	2.38	0.77
1:C:305:LYS:HD2	1:C:305:LYS:N	2.00	0.77
1:C:177:GLU:HG3	1:C:178:ASN:N	1.94	0.76
1:A:111:LYS:HB3	1:A:114:ILE:HD11	1.67	0.76
1:C:301:MET:CE	1:C:319:ILE:HD11	2.15	0.76
1:D:132:VAL:HG22	1:D:144:ILE:HG12	1.68	0.76
1:C:180:LEU:HD23	4:C:1035:HOH:O	1.85	0.75
1:B:198:ARG:NH1	1:B:198:ARG:HA	2.02	0.74
1:A:87:PRO:CD	2:A:1001:1PN:H20	2.17	0.74
1:A:111:LYS:CB	1:A:114:ILE:HD11	2.17	0.74
1:B:86:ILE:HG23	2:B:1001:1PN:H20A	1.70	0.74
1:C:87:PRO:HD2	2:C:1001:1PN:H20	1.70	0.74
1:C:177:GLU:CG	1:C:178:ASN:H	1.85	0.73
1:A:114:ILE:CD1	1:A:115:TYR:CB	2.55	0.73
1:B:87:PRO:HD2	2:B:1001:1PN:C20	2.16	0.72
1:B:198:ARG:HH11	1:B:198:ARG:HA	1.55	0.72
1:C:87:PRO:HD2	2:C:1001:1PN:C20	2.19	0.72
1:C:127:ARG:HD2	1:C:230:LEU:HD11	1.72	0.72
1:C:180:LEU:HD22	1:C:230:LEU:HD22	1.69	0.72
2:A:1001:1PN:H16	2:A:1001:1PN:H20A	1.72	0.71
1:C:304:ASP:HB2	1:C:315:ARG:HG3	1.71	0.71
1:B:134:LYS:HE3	1:B:142:ASP:OD1	1.89	0.71
2:C:1001:1PN:H16	2:C:1001:1PN:H20A	1.71	0.71
1:D:295:ARG:HB3	1:D:295:ARG:HH11	1.55	0.71
1:A:301:MET:HE2	1:A:301:MET:HA	1.71	0.71
1:A:279:ALA:HB3	2:A:1001:1PN:H12	1.73	0.70
1:A:301:MET:SD	1:A:303:PHE:CE1	2.84	0.70
1:B:198:ARG:HH11	1:B:198:ARG:CG	2.05	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ILE:HA	1:A:115:TYR:CB	2.22	0.70
1:A:252:GLN:HE22	1:C:172:SER:HB3	1.58	0.68
1:D:113:PRO:HG2	1:D:114:ILE:H	1.59	0.68
1:C:193:VAL:HG21	1:C:217:LYS:HB2	1.76	0.68
1:C:214:ASN:ND2	1:C:214:ASN:H	1.93	0.67
1:B:85:GLN:NE2	1:B:303:PHE:HB3	2.10	0.67
1:D:267:GLN:HG3	1:D:288:PRO:HA	1.77	0.67
1:A:225:GLU:OE2	1:A:236:ARG:HD2	1.95	0.66
1:C:135:TYR:CE2	1:C:137:GLU:HB2	2.30	0.66
1:A:93:PRO:HD3	1:A:238:LEU:HG	1.77	0.66
2:B:1001:1PN:H16	2:B:1001:1PN:H20A	1.76	0.66
1:B:198:ARG:NH1	1:B:198:ARG:HG3	2.10	0.66
1:D:322:ILE:O	1:D:323:HIS:HB3	1.95	0.66
1:A:114:ILE:CA	1:A:115:TYR:HB3	2.21	0.66
1:B:198:ARG:CA	1:B:198:ARG:HH11	2.09	0.65
1:C:301:MET:HE1	1:C:319:ILE:HD11	1.78	0.65
1:D:93:PRO:HD3	1:D:238:LEU:HG	1.80	0.64
1:D:187:VAL:HG13	1:D:224:SER:HB3	1.80	0.64
1:B:212:PRO:HB2	1:B:215:GLU:HG2	1.80	0.63
1:C:93:PRO:HD3	1:C:238:LEU:HG	1.80	0.63
1:C:230:LEU:HD13	1:C:230:LEU:C	2.18	0.63
1:D:295:ARG:HB3	1:D:295:ARG:NH1	2.14	0.63
1:A:247:VAL:CG1	1:A:258:LEU:HB3	2.30	0.62
1:C:214:ASN:HD22	1:C:214:ASN:H	1.47	0.62
1:C:304:ASP:C	1:C:306:GLN:H	2.03	0.61
1:D:86:ILE:HG23	2:D:1001:1PN:H20A	1.82	0.61
1:C:192:PHE:CE1	1:C:210:GLU:HG2	2.35	0.61
1:A:112:ASP:O	1:A:113:PRO:C	2.39	0.61
1:D:80:ILE:N	1:D:80:ILE:HD12	2.15	0.60
1:C:193:VAL:HG11	1:C:217:LYS:HG3	1.83	0.60
1:B:198:ARG:NE	1:B:198:ARG:H	1.99	0.60
1:A:301:MET:HE2	1:A:319:ILE:HD11	1.84	0.60
1:B:198:ARG:CA	1:B:198:ARG:NH1	2.64	0.60
1:A:145:LYS:NZ	2:A:1001:1PN:HN4	2.00	0.60
1:C:217:LYS:HD2	1:C:220:GLY:O	2.01	0.60
1:C:195:THR:CG2	1:C:209:PHE:HE2	2.14	0.59
1:C:124:HIS:N	1:C:124:HIS:CD2	2.70	0.59
1:D:187:VAL:CG1	1:D:224:SER:HB3	2.32	0.59
1:D:110:ILE:CD1	1:D:121:GLU:CD	2.71	0.58
1:D:82:GLU:OE1	1:D:105:LYS:HE2	2.03	0.58
1:D:140:LYS:HG3	1:D:141:LEU:HD13	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:LYS:NZ	2:D:1001:1PN:HN4	2.01	0.58
1:D:116:GLN:N	1:D:116:GLN:C	2.54	0.58
1:C:306:GLN:O	1:C:307:GLU:HA	2.03	0.58
1:A:114:ILE:HA	1:A:115:TYR:HB3	1.79	0.57
1:C:174:GLN:HE21	1:C:176:CYS:HB3	1.66	0.57
1:A:103:VAL:HG21	1:A:132:VAL:HG21	1.86	0.57
1:C:180:LEU:HD22	1:C:230:LEU:CD2	2.34	0.57
1:A:114:ILE:CD1	1:A:115:TYR:HB3	2.32	0.57
1:D:140:LYS:CG	1:D:141:LEU:HD13	2.35	0.56
1:A:120:ILE:HD12	1:A:120:ILE:N	2.20	0.56
1:C:174:GLN:HE21	1:C:176:CYS:HB2	1.69	0.56
1:C:96:LEU:CG	1:C:318:ARG:NH1	2.61	0.56
1:A:113:PRO:O	1:A:116:GLN:NE2	2.26	0.56
1:D:79:PHE:C	1:D:80:ILE:HD12	2.25	0.56
1:A:236:ARG:HH11	1:A:236:ARG:HG3	1.71	0.55
1:C:195:THR:HG21	1:C:209:PHE:HE2	1.70	0.55
1:D:127:ARG:HH11	1:D:166:ILE:HG21	1.71	0.55
1:A:112:ASP:O	1:A:114:ILE:CD1	2.46	0.55
1:C:304:ASP:O	1:C:306:GLN:N	2.37	0.55
1:A:174:GLN:H	1:A:174:GLN:NE2	2.04	0.55
1:B:198:ARG:NH1	1:B:198:ARG:H	2.05	0.55
1:B:198:ARG:N	1:B:198:ARG:CD	2.70	0.54
1:D:323:HIS:CB	4:D:1011:HOH:O	2.54	0.54
1:C:172:SER:O	1:C:174:GLN:N	2.41	0.54
1:D:180:LEU:HD12	1:D:181:PRO:HD2	1.89	0.54
2:C:1001:1PN:H14	4:C:1048:HOH:O	2.07	0.54
1:B:304:ASP:HB2	1:B:315:ARG:CZ	2.38	0.53
1:D:205:THR:O	1:D:205:THR:HG22	2.07	0.53
1:A:115:TYR:HA	1:A:116:GLN:HG3	1.89	0.53
1:A:187:VAL:HG13	1:A:224:SER:HB3	1.90	0.53
1:A:170:CYS:SG	1:A:174:GLN:NE2	2.82	0.53
1:B:127:ARG:HG3	1:B:150:LEU:HG	1.90	0.53
1:B:209:PHE:CD2	1:B:217:LYS:HE3	2.43	0.53
1:C:195:THR:HG23	4:C:1087:HOH:O	2.08	0.53
1:C:166:ILE:HD12	1:C:230:LEU:CD2	2.38	0.53
1:D:138:ASP:OD2	1:D:140:LYS:NZ	2.41	0.53
1:A:86:ILE:HA	2:A:1001:1PN:H20	1.91	0.52
1:B:177:GLU:HG3	4:B:1031:HOH:O	2.09	0.52
1:D:323:HIS:HB2	4:D:1011:HOH:O	2.08	0.52
1:A:174:GLN:H	1:A:174:GLN:CD	2.13	0.52
1:B:170:CYS:HA	1:B:176:CYS:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:SER:HB2	1:B:198:ARG:HE	1.71	0.52
1:C:124:HIS:H	1:C:124:HIS:CD2	2.28	0.52
1:C:154:LYS:HD2	1:C:176:CYS:SG	2.49	0.52
1:C:170:CYS:O	1:C:262:ILE:HG23	2.10	0.52
1:B:180:LEU:HD11	1:B:231:GLY:N	2.25	0.51
1:D:127:ARG:NH1	1:D:166:ILE:HG21	2.25	0.51
1:D:322:ILE:O	1:D:323:HIS:CB	2.59	0.51
1:C:124:HIS:H	1:C:124:HIS:HD2	1.57	0.51
1:C:301:MET:CE	1:C:301:MET:HA	2.41	0.51
1:D:86:ILE:HA	2:D:1001:1PN:H20	1.92	0.51
1:D:87:PRO:CD	2:D:1001:1PN:H20	2.28	0.50
1:C:225:GLU:OE2	1:C:236:ARG:NH1	2.44	0.50
1:C:304:ASP:HB2	1:C:315:ARG:CG	2.39	0.50
1:A:172:SER:HB2	1:A:174:GLN:OE1	2.12	0.49
1:A:112:ASP:OD1	1:A:113:PRO:HD2	2.13	0.49
1:D:295:ARG:CB	1:D:295:ARG:HH11	2.22	0.49
1:A:187:VAL:CG1	1:A:224:SER:HB3	2.43	0.49
1:C:307:GLU:N	1:C:307:GLU:CB	2.67	0.48
1:A:279:ALA:CB	2:A:1001:1PN:H12	2.43	0.48
1:D:115:TYR:O	1:D:117:LYS:N	2.45	0.48
1:B:96:LEU:HG	1:B:318:ARG:CZ	2.43	0.48
1:D:112:ASP:OD2	1:D:117:LYS:O	2.30	0.48
1:C:96:LEU:HG	1:C:318:ARG:HH12	1.71	0.48
1:C:301:MET:HE2	1:C:319:ILE:HD11	1.95	0.48
1:D:223:LEU:HD22	1:D:240:VAL:HG22	1.95	0.48
1:B:295:ARG:O	1:B:295:ARG:HG3	2.14	0.47
1:C:180:LEU:CD2	4:C:1035:HOH:O	2.53	0.47
1:A:111:LYS:HB2	1:A:114:ILE:HD11	1.91	0.47
1:C:145:LYS:NZ	2:C:1001:1PN:HN4	2.13	0.47
1:B:151:PRO:HB3	1:B:265:PRO:HA	1.96	0.47
1:A:246:GLN:OE1	1:A:249:MET:HE1	2.15	0.47
1:B:135:TYR:CZ	1:B:137:GLU:HB2	2.49	0.47
1:C:127:ARG:HB3	1:C:230:LEU:HD12	1.96	0.47
1:C:86:ILE:HG23	2:C:1001:1PN:C20	2.40	0.47
1:A:132:VAL:HG22	1:A:144:ILE:HG12	1.96	0.47
1:C:87:PRO:HD2	2:C:1001:1PN:H20B	1.94	0.47
1:A:182:VAL:HA	1:A:229:THR:O	2.14	0.46
1:B:315:ARG:HD3	1:B:318:ARG:HD3	1.96	0.46
1:D:135:TYR:CE2	1:D:137:GLU:HB2	2.50	0.46
1:A:114:ILE:HA	1:A:114:ILE:HD12	1.65	0.46
1:B:268:TYR:HB2	1:B:287:VAL:HG13	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:GLN:HE21	1:B:303:PHE:HB3	1.80	0.46
1:C:127:ARG:CD	1:C:230:LEU:HD11	2.44	0.46
1:C:301:MET:SD	1:C:303:PHE:HE1	2.38	0.46
1:B:197:SER:HB2	1:B:198:ARG:CD	2.45	0.46
1:B:161:SER:CB	1:C:161:SER:HB3	2.46	0.46
1:B:315:ARG:NH2	4:B:1022:HOH:O	2.17	0.46
1:D:112:ASP:CG	1:D:119:LEU:HG	2.35	0.46
1:B:267:GLN:HG2	1:B:288:PRO:HA	1.98	0.46
1:D:110:ILE:HD12	1:D:121:GLU:CD	2.35	0.46
1:D:116:GLN:N	1:D:117:LYS:N	2.63	0.46
1:A:246:GLN:OE1	1:A:249:MET:CE	2.63	0.45
2:B:1001:1PN:C3	2:B:1001:1PN:C7	2.93	0.45
1:D:130:ILE:HD12	1:D:322:ILE:HG21	1.98	0.45
1:B:136:PRO:HG2	1:B:137:GLU:OE1	2.16	0.45
1:B:198:ARG:NH1	1:B:198:ARG:N	2.64	0.45
1:C:307:GLU:HB3	1:C:313:GLY:H	1.80	0.45
1:B:264:PRO:HG2	1:B:267:GLN:HB2	1.97	0.45
1:A:301:MET:SD	1:A:303:PHE:CZ	3.10	0.45
1:C:85:GLN:NE2	1:C:98:GLY:HA2	2.32	0.45
1:A:252:GLN:HE22	1:C:172:SER:CB	2.28	0.45
1:A:114:ILE:CG1	1:A:115:TYR:HB3	2.47	0.44
1:D:106:PHE:CZ	1:D:125:PRO:HG3	2.52	0.44
1:A:198:ARG:HD3	1:A:219:ASN:OD1	2.17	0.44
1:B:198:ARG:N	1:B:198:ARG:HD2	2.31	0.44
1:D:127:ARG:NH1	1:D:153:ASP:OD2	2.50	0.44
1:C:223:LEU:HD22	1:C:240:VAL:HG22	1.99	0.44
1:C:193:VAL:CG1	1:C:217:LYS:HG3	2.47	0.44
1:A:112:ASP:OD1	1:A:119:LEU:HD11	2.18	0.44
1:A:301:MET:HE1	1:A:316:LEU:HD23	2.00	0.44
1:C:96:LEU:H	1:C:318:ARG:HH12	1.65	0.44
1:D:110:ILE:HD11	1:D:121:GLU:OE1	2.17	0.43
1:A:236:ARG:NH1	1:A:236:ARG:HG3	2.33	0.43
1:D:267:GLN:HB2	1:D:267:GLN:HE21	1.63	0.43
1:D:323:HIS:OXT	1:D:323:HIS:CG	2.72	0.43
1:C:79:PHE:HB3	1:C:102:LEU:HD11	1.99	0.43
1:A:196:PHE:CD2	1:A:203:GLU:HG2	2.53	0.43
1:C:223:LEU:CD2	1:C:240:VAL:HG22	2.49	0.43
1:B:127:ARG:CG	1:B:150:LEU:HG	2.49	0.43
1:C:124:HIS:N	1:C:124:HIS:HD2	2.16	0.43
1:C:177:GLU:CG	1:C:178:ASN:N	2.57	0.43
1:D:80:ILE:CD1	1:D:80:ILE:N	2.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:ASP:HB2	1:B:315:ARG:NE	2.34	0.42
1:C:230:LEU:HD13	1:C:231:GLY:N	2.35	0.42
1:C:170:CYS:O	1:C:262:ILE:CG2	2.67	0.42
1:B:198:ARG:NH1	1:B:198:ARG:CG	2.72	0.41
1:D:191:ASP:OD1	1:D:213:LYS:NZ	2.52	0.41
1:A:154:LYS:HG3	1:A:262:ILE:CD1	2.51	0.41
1:D:140:LYS:HG2	1:D:141:LEU:HD13	2.02	0.41
1:B:176:CYS:O	1:B:176:CYS:SG	2.79	0.41
1:D:180:LEU:HD11	1:D:231:GLY:HA3	2.02	0.41
1:A:301:MET:CE	1:A:316:LEU:HD23	2.50	0.41
1:B:85:GLN:HG2	1:B:100:PHE:CE1	2.55	0.41
1:C:96:LEU:N	1:C:318:ARG:HH12	2.19	0.41
2:D:1001:1PN:H16	2:D:1001:1PN:C20	2.43	0.41
1:D:323:HIS:HA	4:D:1011:HOH:O	2.21	0.41
1:A:190:SER:HB2	4:A:1078:HOH:O	2.20	0.41
1:A:102:LEU:HA	1:A:102:LEU:HD12	1.94	0.41
1:B:86:ILE:HG23	2:B:1001:1PN:C20	2.43	0.41
1:C:80:ILE:HG13	1:C:105:LYS:HE3	2.03	0.41
1:A:154:LYS:HE3	4:A:1116:HOH:O	2.21	0.40
1:D:100:PHE:HE1	1:D:303:PHE:HD2	1.68	0.40
1:D:110:ILE:HD11	1:D:121:GLU:CD	2.42	0.40
1:B:127:ARG:HG3	1:B:150:LEU:CD1	2.51	0.40

All (9) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:GLY:CA	1:D:115:TYR:OH[4_455]	1.34	0.86
1:D:115:TYR:CD1	4:C:1045:HOH:O[4_555]	1.52	0.68
1:C:266:GLY:N	1:D:115:TYR:CE1[4_455]	1.74	0.46
1:D:115:TYR:CE1	4:C:1045:HOH:O[4_555]	1.76	0.44
1:C:266:GLY:CA	1:D:115:TYR:CZ[4_455]	1.97	0.23
1:C:266:GLY:N	1:D:115:TYR:OH[4_455]	1.99	0.21
1:C:266:GLY:N	1:D:115:TYR:CZ[4_455]	2.07	0.13
1:C:266:GLY:C	1:D:115:TYR:OH[4_455]	2.15	0.05
1:B:198:ARG:NH2	4:A:1095:HOH:O[4_455]	2.16	0.04

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	235/248 (95%)	221 (94%)	13 (6%)	1 (0%)	34	22
1	B	201/248 (81%)	197 (98%)	3 (2%)	1 (0%)	29	17
1	C	220/248 (89%)	209 (95%)	7 (3%)	4 (2%)	8	2
1	D	214/248 (86%)	207 (97%)	5 (2%)	2 (1%)	17	8
All	All	870/992 (88%)	834 (96%)	28 (3%)	8 (1%)	17	8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	173	GLY
1	C	177	GLU
1	C	305	LYS
1	D	113	PRO
1	C	218	GLU
1	A	199	ARG
1	B	139	PRO
1	D	139	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	205/213 (96%)	198 (97%)	7 (3%)	37	25
1	B	184/213 (86%)	179 (97%)	5 (3%)	44	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	193/213 (91%)	186 (96%)	7 (4%)	35	23
1	D	194/213 (91%)	187 (96%)	7 (4%)	35	23
All	All	776/852 (91%)	750 (97%)	26 (3%)	37	25

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	ASP
1	A	114	ILE
1	A	115	TYR
1	A	116	GLN
1	A	174	GLN
1	A	196	PHE
1	A	301	MET
1	B	127	ARG
1	B	167	GLN
1	B	198	ARG
1	B	214	ASN
1	B	223	LEU
1	C	124	HIS
1	C	197	SER
1	C	214	ASN
1	C	217	LYS
1	C	301	MET
1	C	305	LYS
1	C	307	GLU
1	D	87	PRO
1	D	105	LYS
1	D	114	ILE
1	D	141	LEU
1	D	206	SER
1	D	246	GLN
1	D	267	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	252	GLN
1	A	257	GLN

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Mol	Chain	Res	Type
1	B	85	GLN
1	B	167	GLN
1	B	214	ASN
1	C	85	GLN
1	C	124	HIS
1	C	174	GLN
1	C	214	ASN
1	D	167	GLN
1	D	186	ASN
1	D	267	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 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	1PN	C	1001	1	17,20,20	2.17	2 (11%)	17,26,26	2.15	6 (35%)
3	PO4	B	1002	-	4,4,4	1.39	0	6,6,6	0.46	0
2	1PN	A	1001	1	17,20,20	2.29	3 (17%)	17,26,26	2.09	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	1PN	D	1001	1	17,20,20	2.11	2 (11%)	17,26,26	2.04	5 (29%)
2	1PN	B	1001	1	17,20,20	2.32	2 (11%)	17,26,26	1.91	5 (29%)

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	1PN	C	1001	1	-	7/18/31/31	0/1/1/1
2	1PN	A	1001	1	-	7/18/31/31	0/1/1/1
2	1PN	D	1001	1	-	8/18/31/31	0/1/1/1
2	1PN	B	1001	1	-	8/18/31/31	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	1PN	C6-C15	-7.94	1.45	1.54
2	B	1001	1PN	C6-C15	-7.79	1.45	1.54
2	C	1001	1PN	C6-C15	-7.22	1.46	1.54
2	D	1001	1PN	C6-C15	-6.94	1.46	1.54
2	B	1001	1PN	C2-S1	4.30	1.80	1.74
2	C	1001	1PN	C2-S1	4.02	1.80	1.74
2	D	1001	1PN	C2-S1	3.78	1.80	1.74
2	A	1001	1PN	C2-S1	3.06	1.78	1.74
2	A	1001	1PN	C6-C7	2.58	1.54	1.50

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	1PN	O11-C9-C3	-5.12	106.23	112.00
2	C	1001	1PN	O11-C9-C3	-4.88	106.51	112.00
2	D	1001	1PN	O11-C9-C3	-4.88	106.51	112.00
2	B	1001	1PN	O11-C9-C3	-4.51	106.92	112.00
2	C	1001	1PN	O17-C15-C16	3.49	115.15	108.21
2	B	1001	1PN	O17-C15-C16	3.37	114.91	108.21
2	C	1001	1PN	C15-O17-C18	3.35	124.15	117.81
2	A	1001	1PN	O17-C15-C16	3.18	114.55	108.21
2	D	1001	1PN	O17-C15-C16	3.17	114.52	108.21
2	A	1001	1PN	C2-C3-C9	-2.60	124.41	131.89
2	D	1001	1PN	C2-C3-C9	-2.52	124.65	131.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	1PN	C12-O11-C9	2.50	119.58	116.45
2	B	1001	1PN	C15-O17-C18	2.46	122.47	117.81
2	C	1001	1PN	C2-C3-C9	-2.43	124.90	131.89
2	B	1001	1PN	C2-C3-C9	-2.16	125.69	131.89
2	C	1001	1PN	C2-C3-N4	2.08	114.23	111.58
2	A	1001	1PN	C2-C3-N4	2.07	114.22	111.58
2	D	1001	1PN	C15-O17-C18	2.06	121.72	117.81
2	B	1001	1PN	O8-C7-C6	-2.02	120.13	125.23
2	D	1001	1PN	O8-C7-C6	-2.01	120.14	125.23

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1001	1PN	C3-C9-O11-C12
2	A	1001	1PN	C3-C9-O11-C12
2	D	1001	1PN	C3-C9-O11-C12
2	B	1001	1PN	C3-C9-O11-C12
2	C	1001	1PN	O10-C9-O11-C12
2	A	1001	1PN	O10-C9-O11-C12
2	D	1001	1PN	O10-C9-O11-C12
2	B	1001	1PN	O10-C9-O11-C12
2	A	1001	1PN	C20-C18-O17-C15
2	B	1001	1PN	C20-C18-O17-C15
2	A	1001	1PN	O19-C18-O17-C15
2	B	1001	1PN	O19-C18-O17-C15
2	C	1001	1PN	C20-C18-O17-C15
2	D	1001	1PN	C20-C18-O17-C15
2	C	1001	1PN	O19-C18-O17-C15
2	D	1001	1PN	O19-C18-O17-C15
2	C	1001	1PN	O11-C12-C13-C14
2	A	1001	1PN	O11-C12-C13-C14
2	D	1001	1PN	O11-C12-C13-C14
2	B	1001	1PN	O11-C12-C13-C14
2	D	1001	1PN	C16-C15-O17-C18
2	D	1001	1PN	C6-C15-O17-C18
2	B	1001	1PN	C6-C15-O17-C18
2	C	1001	1PN	C16-C15-O17-C18
2	A	1001	1PN	C16-C15-O17-C18
2	B	1001	1PN	C16-C15-O17-C18
2	C	1001	1PN	C5-C6-C7-O8
2	A	1001	1PN	C5-C6-C7-O8

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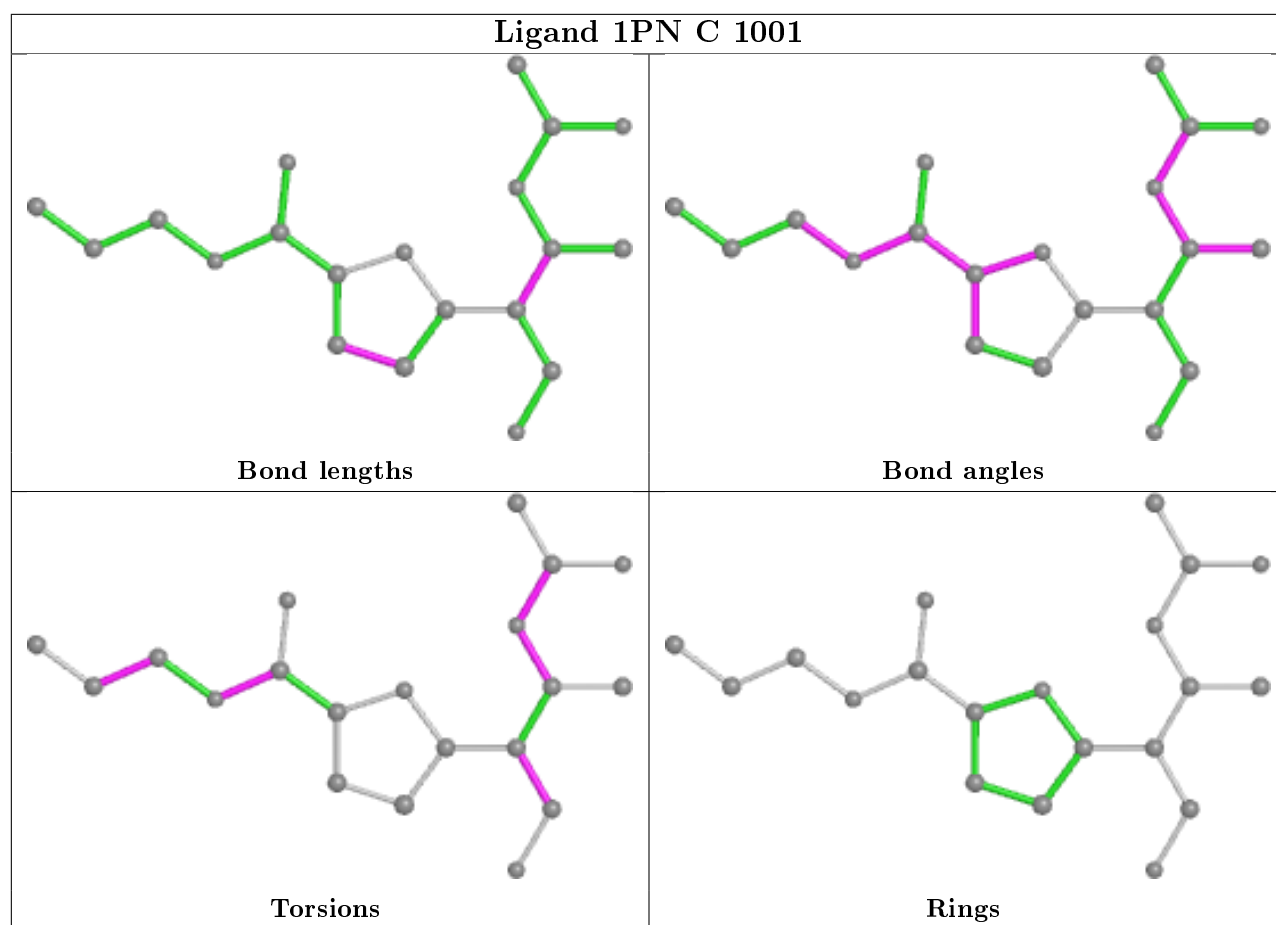
Mol	Chain	Res	Type	Atoms
2	D	1001	1PN	C5-C6-C7-O8
2	B	1001	1PN	C5-C6-C7-O8

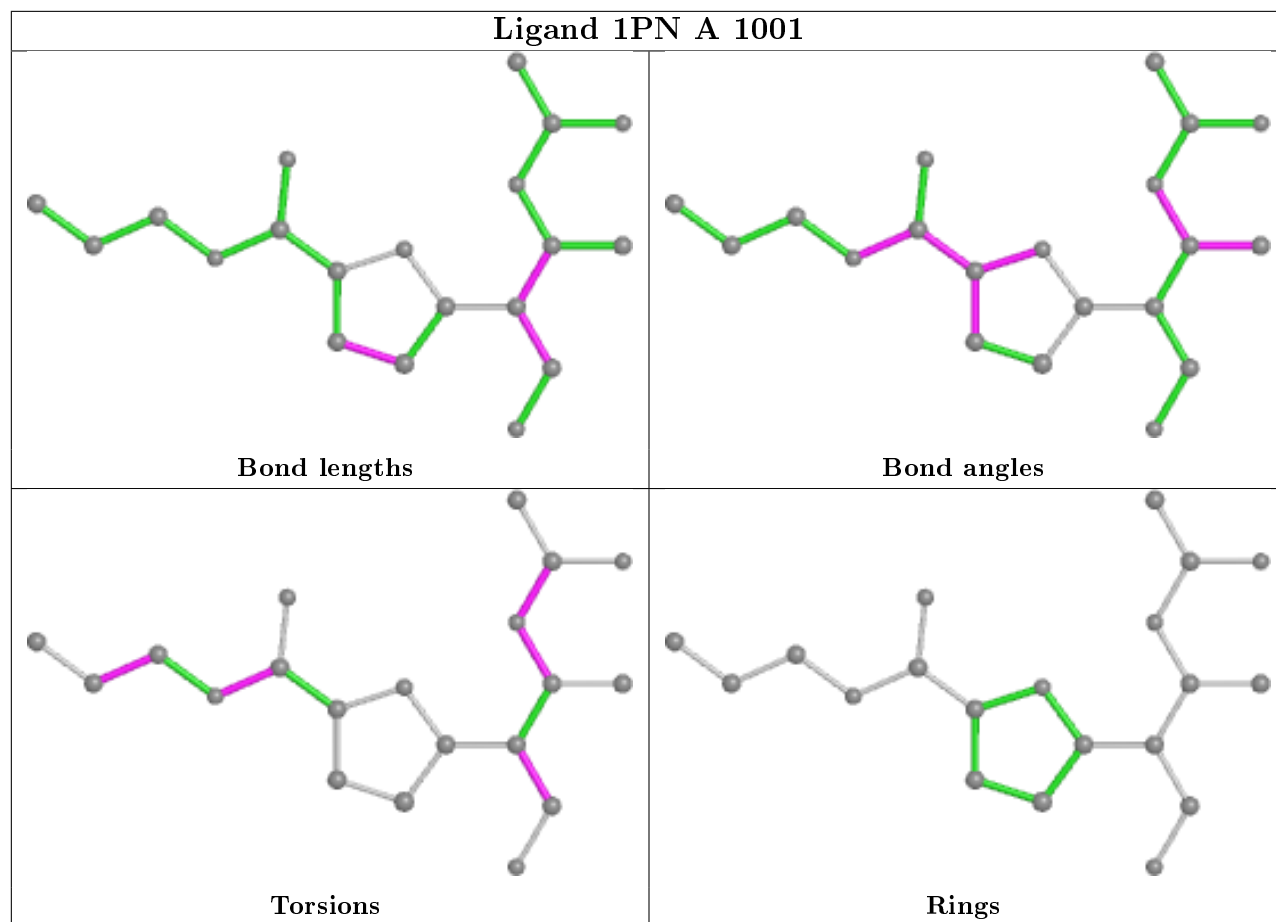
There are no ring outliers.

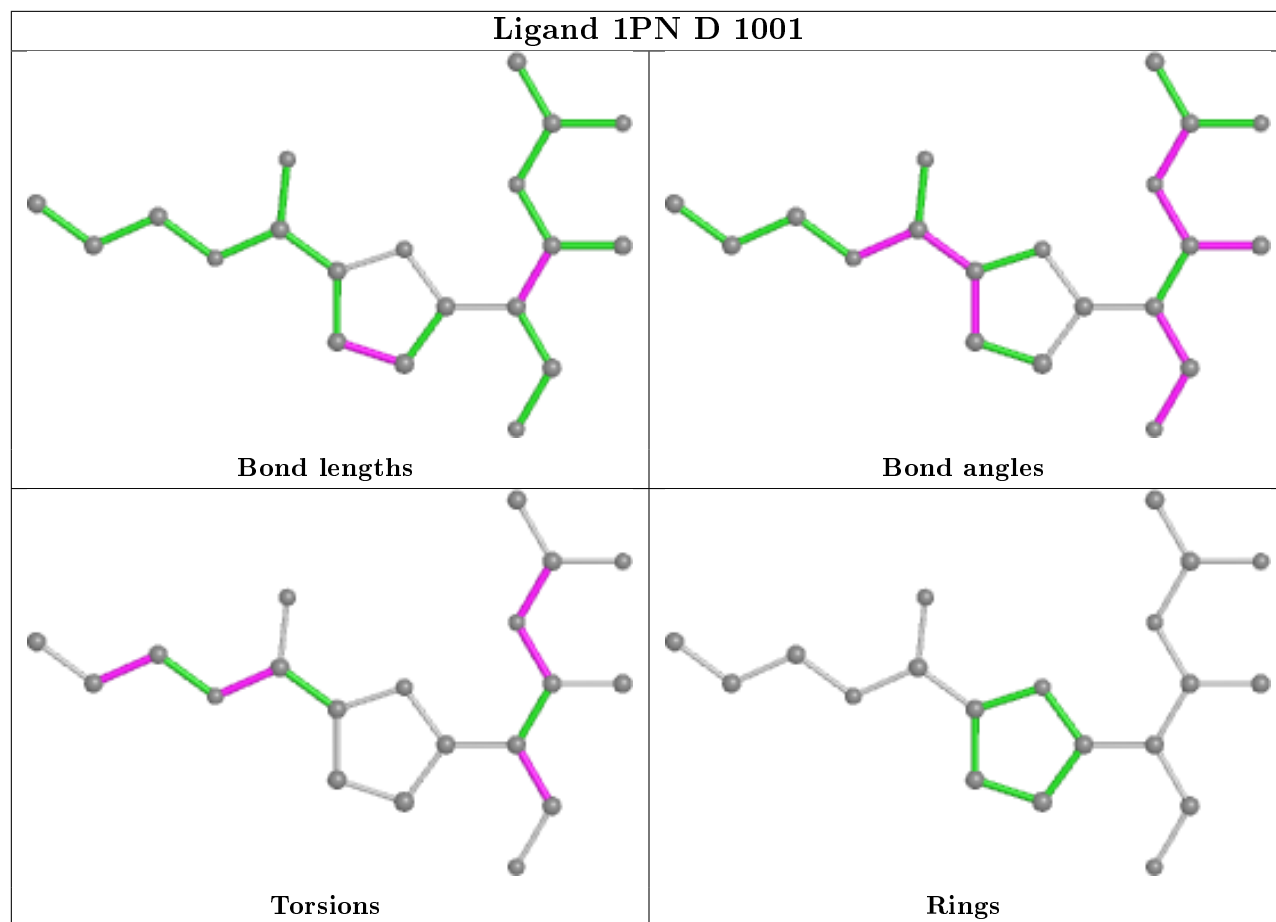
4 monomers are involved in 30 short contacts:

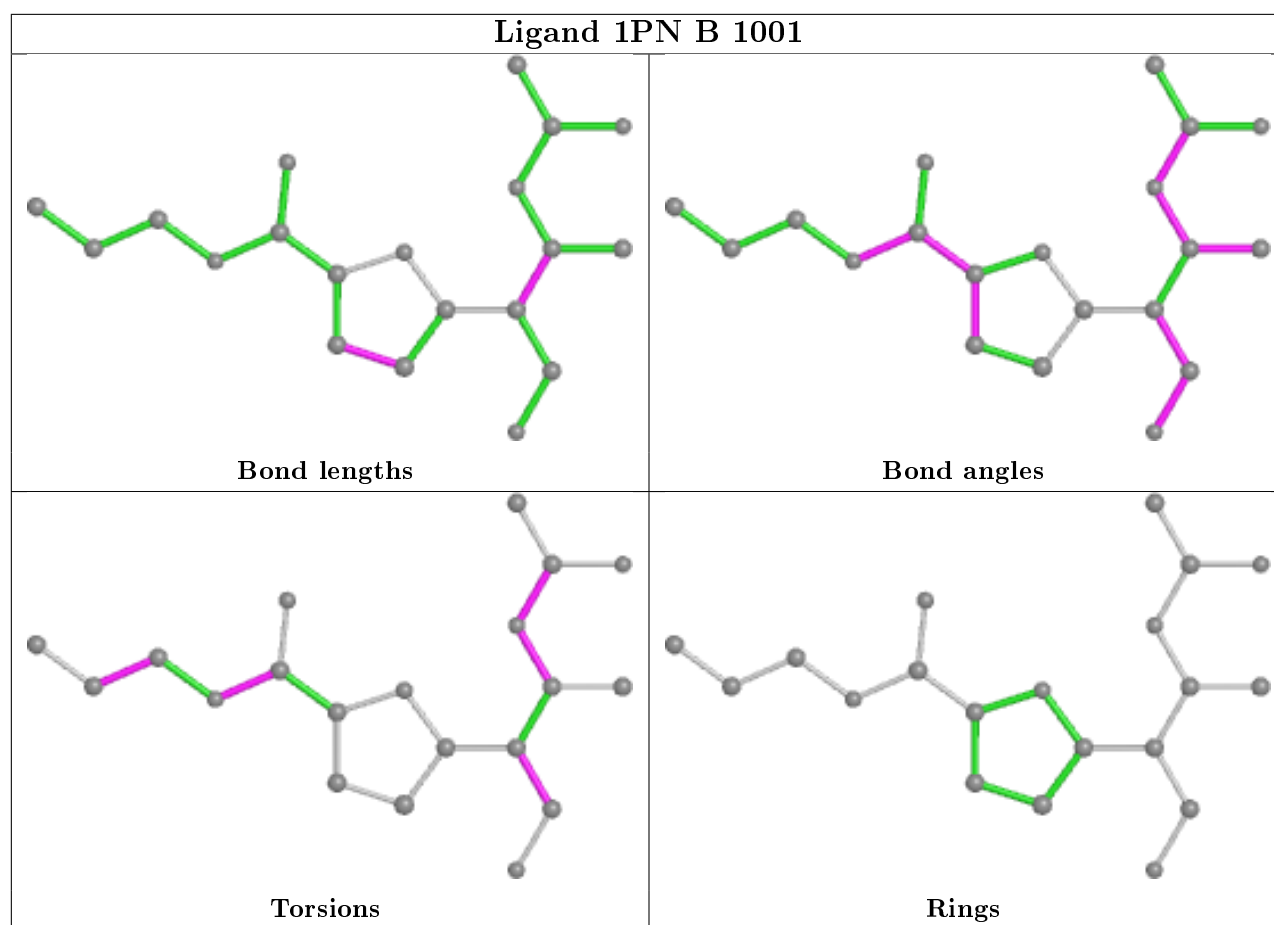
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1001	1PN	8	0
2	A	1001	1PN	8	0
2	D	1001	1PN	8	0
2	B	1001	1PN	6	0

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 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	115:TYR	C	116:GLN	N	1.02
1	A	114:ILE	C	115:TYR	N	0.93

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.