Full wwPDB X-ray Structure Validation Report (i)

Jan 8, 2024 – 04:33 am GMT

PDB ID	:	6FM4
Title	:	The crystal structure of S. aureus Gyrase complex with ID-130 and DNA
Authors	:	Ombrato, R.; Garofalo, B.; Mangano, G.; Mancini, F.
Deposited on	:	2018-01-30
Resolution	:	2.70 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

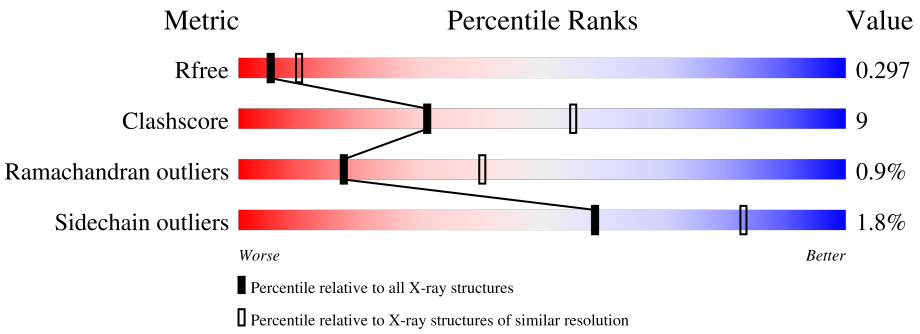
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Mol	Chain	Length	Quality of chain
1	B	692	74% 22% • •
1	D	692	76% 19% • •
2	E	19	58% 42%
3	F	18	67% 33%

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 11536 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 DNA gyrase subunit B,DNA gyrase subunit B,DNA gyrase subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	672	5324	3318	958	1023	25	0	1	0
1	D	670	5291	3298	952	1016	25	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	409	MET	LEU	conflict	UNP P66937
B	544	THR	-	linker	UNP P66937
B	545	GLY	-	linker	UNP P66937
B	1123	PHE	TYR	engineered mutation	UNP Q99XG5
D	409	MET	LEU	conflict	UNP P66937
D	544	THR	-	linker	UNP P66937
D	545	GLY	-	linker	UNP P66937
D	1123	PHE	TYR	engineered mutation	UNP Q99XG5

- Molecule 2 is a DNA chain called DNA (5'-5UA*D(P*GP*CP*CP*GP*TP*AP*GP*GP*GP*CP*CP*CP*TP*AP*CP*GP*GP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	19	389	183	72	115	19	0	0	0

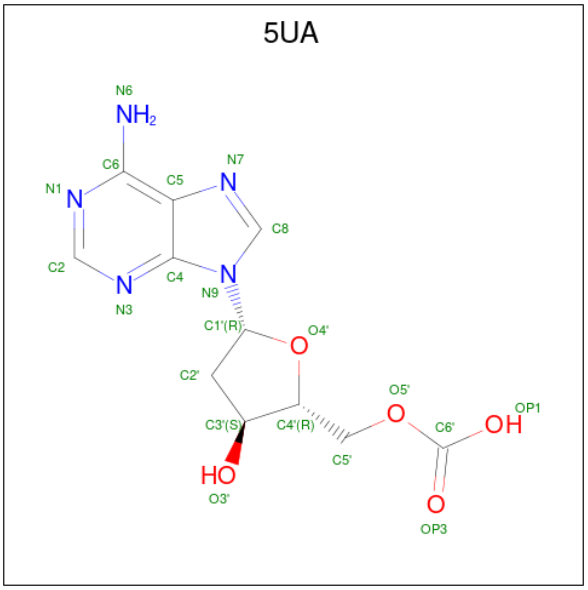
- Molecule 3 is a DNA chain called DNA (5'-5UA*D(P*GP*CP*CP*GP*TP*AP*GP*GP*GP*CP*CP*CP*TP*AP*CP*GP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	F	18	369	173	70	108	18	0	0	0

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

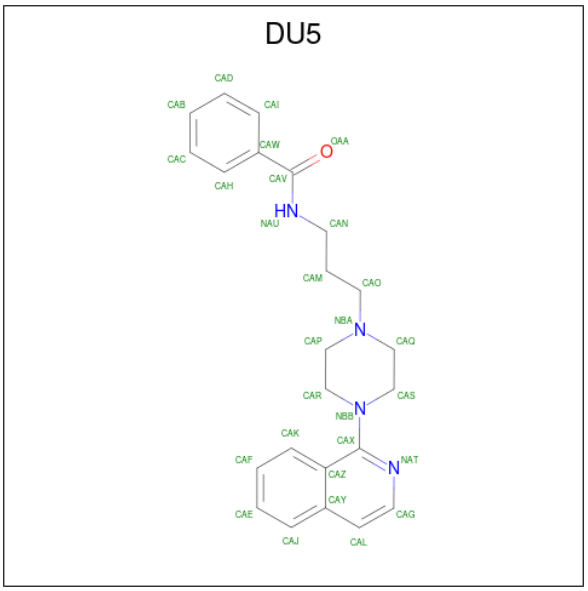
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mn 1 1	1	0
4	E	1	Total Mn 1 1	1	0
4	F	1	Total Mn 1 1	1	0

- Molecule 5 is 5'-O-CARBOXY-2'-DEOXYADENOSINE (three-letter code: 5UA) (formula: $C_{11}H_{13}N_5O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	E	1	Total C N O 21 11 5 5	21	0
5	F	1	Total C N O 21 11 5 5	21	0

- Molecule 6 is {N}-[3-(4-isoquinolin-1-ylpiperazin-1-yl)propyl]benzamide (three-letter code: DU5) (formula: $C_{23}H_{26}N_4O$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	E	1	56	46	8	2	0	1

- Molecule 7 is water.

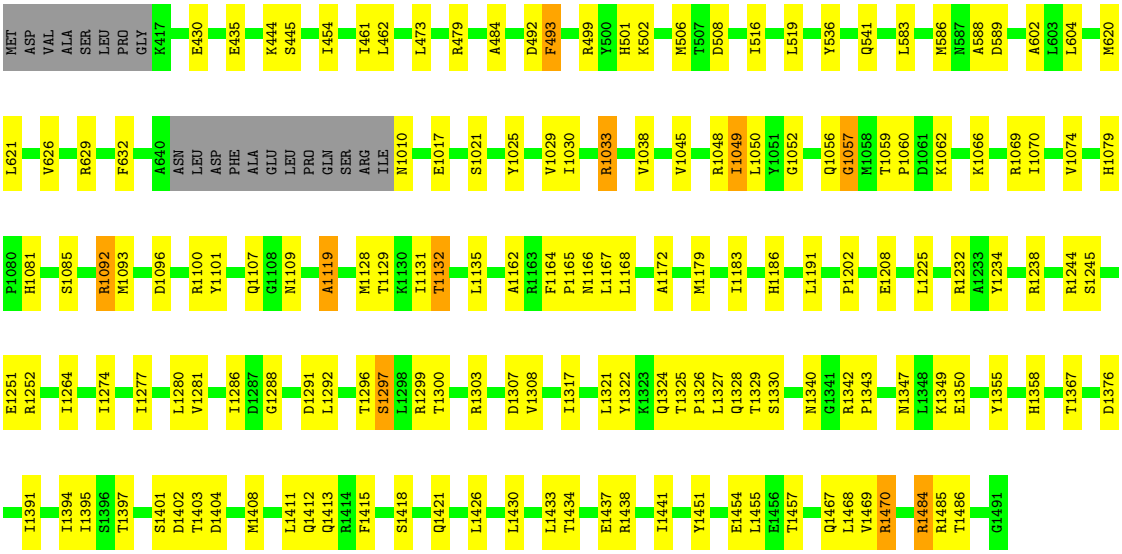
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	B	21	21	21	0	0
			Total	O		
7	D	37	37	37	0	0
			Total	O		
7	E	3	3	3	0	0
			Total	O		
7	F	1	1	1	0	0

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. 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.

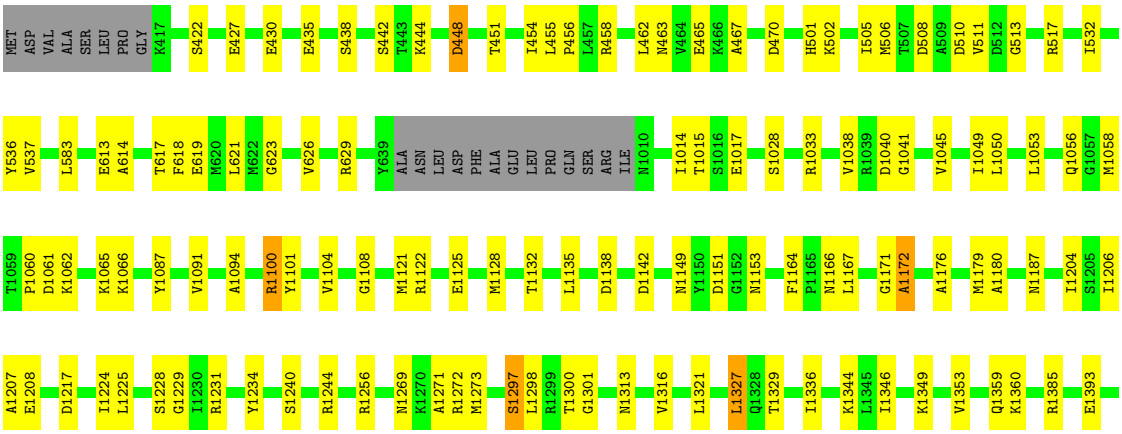
- Molecule 1: DNA gyrase subunit B,DNA gyrase subunit B,DNA gyrase subunit A

Chain B:



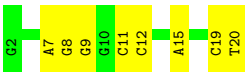
- Molecule 1: DNA gyrase subunit B,DNA gyrase subunit B,DNA gyrase subunit A

Chain D:



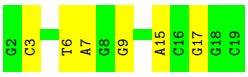
• Molecule 2: DNA (5'-5UA*D(P*GP*CP*CP*GP*TP*AP*GP*GP*GP*CP*CP*CP*TP*AP*CP*GP*GP*CP*T)-3')

Chain E: 58% 42%



• Molecule 3: DNA (5'-5UA*D(P*GP*CP*CP*GP*TP*AP*GP*GP*GP*CP*CP*CP*TP*AP*CP*GP*GP*C)-3')

Chain F: 67% 33%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	92.95Å 92.95Å 407.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.05 – 2.70 39.43 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.05-2.70) 99.8 (39.43-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	8.23 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.225 , 0.300 0.226 , 0.297	Depositor DCC
R_{free} test set	2750 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	39.7	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 3.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.31$, $< L^2 > = 0.14$	Xtriage
Estimated twinning fraction	0.379 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	11536	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.49% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MN, 5UA, DU5

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	B	0.69	0/5395	0.89	5/7267 (0.1%)
1	D	0.69	0/5362	0.89	5/7227 (0.1%)
2	E	0.60	0/435	0.93	0/669
3	F	0.58	0/413	0.90	1/635 (0.2%)
All	All	0.68	0/11605	0.89	11/15798 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	1484	ARG	NE-CZ-NH2	-6.05	117.28	120.30
1	B	1470	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	B	1485	ARG	NE-CZ-NH1	5.95	123.27	120.30
3	F	3	DC	C1'-O4'-C4'	-5.66	104.44	110.10
1	D	1122	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	D	1231	ARG	NE-CZ-NH2	-5.39	117.60	120.30
1	B	1470	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	B	1092	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	D	1385	ARG	NE-CZ-NH2	-5.17	117.71	120.30
1	D	1033	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	D	1100	ARG	NE-CZ-NH1	5.05	122.83	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	5324	0	5356	110	0
1	D	5291	0	5308	93	1
2	E	389	0	213	7	0
3	F	369	0	201	6	0
4	B	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
5	E	21	0	11	0	1
5	F	21	0	11	0	0
6	E	56	0	0	6	0
7	B	21	0	0	8	0
7	D	37	0	0	9	0
7	E	3	0	0	1	0
7	F	1	0	0	2	0
All	All	11536	0	11100	204	1

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 9.

All (204) 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:435:GLU:OE2	1:B:508:ASP:OD2	1.81	0.96
6:E:103[A]:DU5:CAR	6:E:103[A]:DU5:CAK	2.50	0.89
1:D:427:GLU:HA	7:D:1619:HOH:O	1.77	0.83
6:E:103[B]:DU5:CAK	6:E:103[B]:DU5:CAR	2.56	0.82
1:D:1297:SER:HG	1:D:1300:THR:HG1	1.22	0.80
6:E:103[B]:DU5:CAN	6:E:103[B]:DU5:CAI	2.59	0.80
1:D:435:GLU:OE2	1:D:508:ASP:OD2	1.99	0.79
1:D:1176:ALA:HB3	1:D:1179:MET:O	1.83	0.77
1:D:1313:ASN:HD22	1:D:1316:VAL:HG23	1.52	0.73
1:D:462:LEU:O	7:D:1601:HOH:O	2.08	0.72
1:B:430:GLU:HB3	1:B:502:LYS:HB2	1.74	0.70
1:D:614:ALA:O	1:D:618:PHE:HD1	1.76	0.69
1:D:1321:LEU:HB3	1:D:1327:LEU:HD22	1.76	0.68
1:B:1397:THR:HG21	1:B:1411:LEU:HD23	1.75	0.67
1:D:1297:SER:O	1:D:1301:GLY:N	2.26	0.67
1:B:1340:ASN:HB2	7:B:1602:HOH:O	1.95	0.65
1:B:435:GLU:HG2	1:B:516:ILE:HD13	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:621:LEU:O	1:D:629:ARG:NH1	2.28	0.64
1:B:519:LEU:O	7:B:1601:HOH:O	2.15	0.63
1:B:1070:ILE:O	1:B:1074:VAL:HG23	1.99	0.63
1:D:619:GLU:HG3	7:D:1602:HOH:O	1.99	0.63
1:D:455:LEU:HD12	1:D:456:PRO:HD2	1.80	0.62
1:D:462:LEU:HD11	1:D:467:ALA:HB2	1.80	0.61
1:D:1056:GLN:HB2	1:D:1058:MET:HG3	1.82	0.61
1:B:1129:THR:OG1	1:B:1132:THR:OG1	2.17	0.61
1:D:1403:THR:OG1	1:D:1406:VAL:HG23	2.01	0.61
1:D:427:GLU:HG2	7:D:1617:HOH:O	2.00	0.60
1:D:1050:LEU:HD22	1:D:1132:THR:HG21	1.82	0.60
1:D:1349:LYS:O	1:D:1353:VAL:HG23	2.01	0.60
1:B:604:LEU:HD23	1:B:1010:ASN:HB3	1.83	0.60
1:B:541:GLN:NE2	1:B:604:LEU:HD11	2.17	0.60
1:B:1030:ILE:HG21	1:B:1343:PRO:HG3	1.84	0.60
1:B:1277:ILE:HG22	1:B:1292:LEU:HD22	1.83	0.59
1:B:1376:ASP:HB3	7:B:1618:HOH:O	2.02	0.59
1:D:1058:MET:HG2	1:D:1065:LYS:HG3	1.84	0.59
1:D:501:HIS:CD2	7:D:1619:HOH:O	2.55	0.58
1:D:1321:LEU:HB3	1:D:1327:LEU:CD2	2.33	0.58
1:D:1427:ASP:O	1:D:1428:MET:C	2.42	0.58
1:D:1445:TYR:CE2	1:D:1449:LEU:HD11	2.39	0.58
1:D:1045:VAL:O	1:D:1049:ILE:HG13	2.03	0.58
1:D:1151:ASP:OD1	1:D:1153:ASN:HB2	2.05	0.57
1:D:1204:ILE:HA	1:D:1208:GLU:OE2	2.05	0.57
1:B:1033:ARG:CZ	2:E:7:DA:H4'	2.34	0.57
1:B:435:GLU:OE2	1:B:508:ASP:CG	2.43	0.56
1:B:462:LEU:HD12	3:F:15:DA:OP1	2.05	0.56
1:D:506:MET:HG2	1:D:583:LEU:HD11	1.88	0.55
3:F:9:DG:N2	7:F:1601:HOH:O	2.30	0.55
1:B:1107:GLN:CG	7:B:1609:HOH:O	2.54	0.55
1:B:1300:THR:HG21	1:B:1303:ARG:HD3	1.88	0.55
1:B:1355:TYR:O	1:B:1358:HIS:HB3	2.07	0.55
1:B:1434:THR:O	1:B:1437:GLU:HB2	2.07	0.54
1:B:1062:LYS:NZ	7:B:1604:HOH:O	2.41	0.54
1:B:1202:PRO:O	1:B:1349:LYS:NZ	2.28	0.54
1:B:1280:LEU:HD11	1:B:1324:GLN:HB3	1.89	0.53
1:D:1344:LYS:HD3	1:D:1346:ILE:HG22	1.89	0.53
1:B:435:GLU:OE2	2:E:9:DG:OP1	2.26	0.53
1:D:1171:GLY:O	1:D:1172:ALA:HB2	2.07	0.53
1:D:1187:ASN:HB2	1:D:1217:ASP:OD2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1346:ILE:HA	7:D:1603:HOH:O	2.08	0.53
1:B:541:GLN:OE1	1:B:602:ALA:O	2.27	0.53
1:D:448:ASP:O	1:D:448:ASP:CG	2.46	0.52
1:D:1430:LEU:O	1:D:1432:ARG:N	2.43	0.52
1:B:1291:ASP:OD2	1:B:1292:LEU:N	2.43	0.52
1:B:1038:VAL:HA	1:B:1167:LEU:HD22	1.91	0.51
1:B:1252:ARG:NH1	1:B:1308:VAL:O	2.43	0.51
1:B:1438:ARG:O	1:B:1441:ILE:HB	2.09	0.51
1:B:588:ALA:CB	1:D:1298:LEU:HD21	2.41	0.51
1:B:1025:TYR:O	1:B:1029:VAL:HG23	2.10	0.51
1:D:532:ILE:CG1	1:D:537:VAL:HG21	2.41	0.51
1:B:1096:ASP:C	1:B:1096:ASP:OD2	2.49	0.51
1:B:1059:THR:HA	1:B:1128:MET:HE3	1.93	0.50
1:B:516:ILE:HA	1:B:519:LEU:HD12	1.94	0.50
1:B:1296:THR:HG22	1:B:1297:SER:N	2.27	0.50
1:B:1167:LEU:HD11	1:B:1183:ILE:HD12	1.94	0.50
1:D:532:ILE:HG13	1:D:537:VAL:HG21	1.95	0.49
1:B:1288:GLY:HA3	1:B:1317:ILE:HD13	1.93	0.49
1:B:1402:ASP:OD1	1:B:1403:THR:HG23	2.12	0.49
1:D:1225:LEU:HD11	1:D:1244:ARG:NE	2.27	0.49
1:D:451:THR:HG22	7:D:1616:HOH:O	2.12	0.49
1:D:458:ARG:HB3	7:F:1601:HOH:O	2.12	0.49
1:B:1329:THR:HG22	1:B:1330:SER:N	2.28	0.48
1:B:1048:ARG:O	1:B:1052:GLY:N	2.32	0.48
1:B:1092:ARG:HD3	7:E:203:HOH:O	2.13	0.48
1:D:1040:ASP:O	1:D:1166:ASN:HB3	2.13	0.48
1:B:589:ASP:N	1:B:589:ASP:OD1	2.47	0.48
1:B:1277:ILE:CG2	1:B:1292:LEU:HD22	2.43	0.48
1:D:430:GLU:HB3	1:D:502:LYS:HB2	1.96	0.48
1:B:1234:TYR:O	1:B:1347:ASN:HB2	2.14	0.48
1:D:1060:PRO:HD3	1:D:1128:MET:HB2	1.95	0.48
1:D:1014:ILE:HG23	1:D:1015:THR:N	2.28	0.48
1:B:1225:LEU:HD11	1:B:1244:ARG:NE	2.29	0.48
1:B:1433:LEU:N	1:B:1433:LEU:HD12	2.28	0.48
1:D:1180:ALA:O	1:D:1336:ILE:HD12	2.14	0.47
3:F:6:DT:C2	3:F:7:DA:C8	3.02	0.47
1:B:588:ALA:HB3	1:D:1298:LEU:HD21	1.97	0.47
1:B:506:MET:HG2	1:B:583:LEU:HD11	1.96	0.47
1:B:1109:ASN:HB3	1:B:1119:ALA:HB2	1.97	0.47
1:B:1119:ALA:HA	1:D:438:SER:HA	1.95	0.47
1:D:1273:MET:SD	1:D:1327:LEU:HD12	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:586:MET:O	1:D:1108:GLY:HA2	2.15	0.47
1:B:1467:GLN:HA	1:B:1470:ARG:HD2	1.97	0.47
1:B:445:SER:O	1:D:1298:LEU:HD22	2.15	0.47
1:B:461:ILE:HD12	1:B:462:LEU:H	1.80	0.47
1:D:1053:LEU:HD23	1:D:1058:MET:SD	2.55	0.47
1:D:613:GLU:O	1:D:617:THR:OG1	2.20	0.46
1:D:1359:GLN:NE2	1:D:1359:GLN:HA	2.31	0.46
1:D:510:ASP:O	1:D:513:GLY:N	2.48	0.46
1:D:1038:VAL:HA	1:D:1167:LEU:HD22	1.96	0.46
1:B:1135:LEU:HA	1:B:1162:ALA:HA	1.98	0.46
1:D:1094:ALA:HA	1:D:1104:VAL:O	2.16	0.46
1:B:492:ASP:O	1:B:493:PHE:C	2.54	0.46
1:B:1172:ALA:O	1:B:1183:ILE:N	2.47	0.46
2:E:19:DC:H2'	2:E:20:DT:C6	2.51	0.46
1:B:620:MET:SD	1:B:632:PHE:HB2	2.57	0.46
1:D:462:LEU:HD12	2:E:15:DA:OP1	2.16	0.46
1:B:1165:PRO:HG2	1:B:1191:LEU:HD13	1.98	0.45
1:D:444:LYS:HA	1:D:454:ILE:CD1	2.46	0.45
1:D:1206:ILE:HD12	1:D:1234:TYR:O	2.16	0.45
1:B:473:LEU:O	1:B:479:ARG:HD2	2.15	0.45
1:B:1179:MET:O	3:F:17:DG:H4'	2.16	0.45
1:B:1056:GLN:O	1:B:1057:GLY:C	2.55	0.45
1:B:1408:MET:SD	1:B:1426:LEU:HD12	2.56	0.45
1:B:1225:LEU:HD21	1:B:1244:ARG:HD2	1.98	0.45
1:B:1404:ASP:OD2	1:D:1431:ARG:NH2	2.50	0.45
1:D:501:HIS:CE1	1:D:536:TYR:HH	2.34	0.45
1:B:1281:VAL:HG22	1:B:1286:ILE:CG2	2.47	0.45
1:D:1229:GLY:HA3	1:D:1240:SER:O	2.16	0.45
1:B:1391:ILE:O	1:B:1395:ILE:HG12	2.17	0.44
1:B:621:LEU:O	1:B:629:ARG:HD3	2.18	0.44
1:D:619:GLU:O	1:D:623:GLY:HA3	2.18	0.44
1:B:1325:THR:HB	1:B:1326:PRO:CD	2.47	0.44
1:B:1166:ASN:O	1:B:1167:LEU:C	2.56	0.44
1:B:1050:LEU:HD22	1:B:1132:THR:HG21	2.00	0.44
1:B:1397:THR:HG21	1:B:1411:LEU:CD2	2.47	0.44
1:D:463:ASN:OD1	1:D:463:ASN:C	2.55	0.44
1:D:1066:LYS:HA	1:D:1125:GLU:HG2	1.99	0.44
1:B:1107:GLN:HG3	7:B:1609:HOH:O	2.16	0.44
1:D:1430:LEU:C	1:D:1432:ARG:N	2.72	0.44
1:B:1252:ARG:NH2	1:B:1307:ASP:OD2	2.52	0.43
1:D:1061:ASP:OD1	1:D:1061:ASP:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1475:GLU:O	1:D:1478:ASP:HB2	2.18	0.43
1:B:1232:ARG:HG2	1:B:1238:ARG:O	2.19	0.43
1:D:1100:ARG:HG3	1:D:1101:TYR:CE2	2.52	0.43
1:B:1418:SER:OG	1:B:1421:GLN:HG3	2.19	0.43
1:D:1017:GLU:HA	1:D:1017:GLU:OE1	2.18	0.43
1:B:1342:ARG:NH2	7:B:1602:HOH:O	2.37	0.43
1:D:1121:MET:HE3	6:E:103[A]:DU5:CAW	2.49	0.43
1:B:1413:GLN:HG3	7:B:1605:HOH:O	2.18	0.43
1:D:1041:GLY:HA2	1:D:1167:LEU:HD13	2.00	0.43
1:D:1121:MET:HE3	6:E:103[B]:DU5:CAW	2.49	0.43
1:B:1049:ILE:HG22	1:B:1050:LEU:N	2.33	0.42
1:D:501:HIS:NE2	1:D:536:TYR:OH	2.51	0.42
1:B:1168:LEU:O	1:B:1186:HIS:ND1	2.37	0.42
1:D:1269:ASN:HB3	1:D:1272:ARG:HB3	2.02	0.42
1:D:1135:LEU:HD21	1:D:1164:PHE:CZ	2.55	0.42
1:D:1058:MET:CG	1:D:1065:LYS:HG3	2.50	0.42
1:D:1224:ILE:HG13	1:D:1486:THR:HG21	2.02	0.42
1:B:1274:ILE:HG23	1:B:1292:LEU:HD21	2.01	0.42
1:D:505:ILE:HG21	1:D:517:ARG:HG3	2.01	0.42
1:D:1062:LYS:N	1:D:1062:LYS:HD2	2.34	0.42
1:D:1414:ARG:HG2	1:D:1415:PHE:CE1	2.55	0.42
1:B:501:HIS:CE1	1:B:536:TYR:OH	2.73	0.42
1:B:1245:SER:OG	1:B:1264:ILE:HA	2.20	0.42
1:B:1347:ASN:OD1	1:B:1350[A]:GLU:HG3	2.19	0.42
1:B:1468:LEU:O	1:B:1469:VAL:C	2.58	0.42
1:D:1207:ALA:O	1:D:1208:GLU:C	2.57	0.42
1:D:1271:ALA:O	1:D:1272:ARG:C	2.57	0.42
1:B:1045:VAL:HB	1:B:1079:HIS:CE1	2.55	0.42
1:B:1402:ASP:HA	1:D:1435:GLY:H	1.85	0.42
1:B:1451:TYR:CZ	1:B:1455:LEU:HD11	2.55	0.42
1:D:626:VAL:O	1:D:629:ARG:N	2.53	0.42
1:D:1359:GLN:O	1:D:1360:LYS:C	2.58	0.42
1:B:1060:PRO:HD3	1:B:1128:MET:HB2	2.01	0.42
1:D:465:GLU:OE2	7:D:1602:HOH:O	2.22	0.42
1:B:1179:MET:C	3:F:17:DG:H4'	2.39	0.41
1:D:1430:LEU:C	1:D:1432:ARG:H	2.24	0.41
2:E:11:DC:C2	2:E:12:DC:C5	3.07	0.41
1:B:1367:THR:HG23	1:B:1455:LEU:HD22	2.03	0.41
1:B:1164:PHE:HB2	1:B:1165:PRO:HD2	2.03	0.41
1:B:1347:ASN:OD1	1:B:1347:ASN:C	2.59	0.41
1:B:1430:LEU:HA	1:B:1433:LEU:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1050:LEU:HD22	1:D:1132:THR:CG2	2.50	0.41
1:D:1121:MET:CE	6:E:103[A]:DU5:CAI	2.98	0.41
1:B:1017:GLU:OE2	1:B:1021:SER:OG	2.32	0.41
1:B:1066:LYS:O	1:B:1069:ARG:HG2	2.21	0.41
1:B:1286:ILE:O	1:B:1286:ILE:HG23	2.19	0.41
1:B:1322:TYR:O	1:B:1328:GLN:HB3	2.21	0.41
1:B:1394:ILE:HD13	1:B:1415:PHE:CE1	2.56	0.41
1:B:1454:GLU:O	1:B:1457:THR:HB	2.21	0.41
1:D:501:HIS:CG	7:D:1619:HOH:O	2.73	0.41
1:D:511:VAL:HG11	1:D:1028:SER:HB2	2.02	0.41
1:B:1411:LEU:O	1:B:1412:GLN:C	2.59	0.41
2:E:11:DC:C2	2:E:12:DC:C6	3.08	0.41
1:B:1484:ARG:HD3	1:B:1486:THR:O	2.21	0.41
1:D:1313:ASN:ND2	1:D:1316:VAL:HG23	2.27	0.41
1:B:1049:ILE:HG21	1:B:1093:MET:CE	2.51	0.40
1:D:1087:TYR:O	1:D:1091:VAL:HG23	2.21	0.40
1:B:1081:HIS:HA	2:E:8:DG:OP1	2.21	0.40
1:B:626:VAL:HG11	3:F:17:DG:H3'	2.03	0.40
1:B:1100:ARG:HG3	1:B:1101:TYR:CE2	2.57	0.40
1:B:1101:TYR:HB3	1:B:1131:ILE:HD13	2.02	0.40
1:B:444:LYS:HA	1:B:454:ILE:CD1	2.52	0.40
1:B:484:ALA:O	1:B:499:ARG:HD3	2.21	0.40
1:B:1321:LEU:O	1:B:1325:THR:OG1	2.18	0.40

All (1) 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:D:1393:GLU:OE2	5:E:102:5UA:N1[1_655]	2.09	0.11

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	B	667/692 (96%)	610 (92%)	51 (8%)	6 (1%)	17	40
1	D	664/692 (96%)	612 (92%)	46 (7%)	6 (1%)	17	40
All	All	1331/1384 (96%)	1222 (92%)	97 (7%)	12 (1%)	17	40

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1033	ARG
1	B	1057	GLY
1	B	1119	ALA
1	B	1327	LEU
1	D	1142	ASP
1	D	1431	ARG
1	D	1172	ALA
1	B	493	PHE
1	B	1049	ILE
1	D	1327	LEU
1	D	1149	ASN
1	D	1256	ARG

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	B	573/591 (97%)	566 (99%)	7 (1%)	71	88
1	D	568/591 (96%)	555 (98%)	13 (2%)	50	78
All	All	1141/1182 (96%)	1121 (98%)	20 (2%)	59	83

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1085	SER
1	B	1132	THR
1	B	1208	GLU
1	B	1251	GLU

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Mol	Chain	Res	Type
1	B	1297	SER
1	B	1299	ARG
1	B	1401	SER
1	D	422	SER
1	D	442	SER
1	D	448	ASP
1	D	470	ASP
1	D	1138	ASP
1	D	1228	SER
1	D	1297	SER
1	D	1329	THR
1	D	1396	SER
1	D	1401	SER
1	D	1414	ARG
1	D	1450	ASN
1	D	1465	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	501	HIS
1	B	541	GLN
1	B	605	GLN
1	B	1368	GLN
1	D	476	ASN
1	D	480	GLN
1	D	1313	ASN
1	D	1334	ASN
1	D	1390	HIS

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)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

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	$\#\|Z\| > 2$
6	DU5	E	103[A]	-	31,31,31	1.37	3 (9%)	40,41,41	2.03	10 (25%)
5	5UA	F	1502	3	21,23,23	2.76	3 (14%)	20,33,33	2.52	6 (30%)
5	5UA	E	102	-	21,23,23	2.70	3 (14%)	20,33,33	2.79	6 (30%)
6	DU5	E	103[B]	-	31,31,31	1.36	3 (9%)	40,41,41	2.00	11 (27%)

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
6	DU5	E	103[A]	-	-	5/15/25/25	0/4/4/4
5	5UA	F	1502	3	-	1/4/21/21	0/3/3/3
5	5UA	E	102	-	-	3/4/21/21	0/3/3/3
6	DU5	E	103[B]	-	-	6/15/25/25	0/4/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	1502	5UA	C2-N3	8.79	1.46	1.32
5	E	102	5UA	C2-N3	8.26	1.45	1.32
5	F	1502	5UA	C2-N1	6.61	1.46	1.33
5	E	102	5UA	C2-N1	6.49	1.46	1.33
5	E	102	5UA	O5'-C6'	5.40	1.48	1.33
5	F	1502	5UA	O5'-C6'	5.21	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	103[A]	DU5	CAW-CAV	-4.74	1.40	1.50
6	E	103[B]	DU5	CAW-CAV	-4.50	1.40	1.50
6	E	103[A]	DU5	CAG-NAT	2.82	1.40	1.34
6	E	103[B]	DU5	CAG-NAT	2.77	1.40	1.34
6	E	103[B]	DU5	CAS-NBB	2.26	1.50	1.46
6	E	103[A]	DU5	CAS-NBB	2.16	1.50	1.46

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	E	102	5UA	N3-C2-N1	-8.76	114.98	128.68
5	F	1502	5UA	N3-C2-N1	-8.68	115.10	128.68
6	E	103[A]	DU5	CAO-NBA-CAQ	-5.90	96.14	111.23
5	E	102	5UA	O5'-C5'-C4'	5.84	120.80	108.43
6	E	103[B]	DU5	CAO-NBA-CAQ	-5.61	96.89	111.23
6	E	103[B]	DU5	CAR-NBB-CAS	5.60	123.87	111.52
6	E	103[A]	DU5	CAR-NBB-CAS	5.58	123.83	111.52
6	E	103[A]	DU5	CAO-NBA-CAP	-5.02	98.40	111.23
6	E	103[B]	DU5	CAO-NBA-CAP	-4.53	99.64	111.23
5	E	102	5UA	C4-C5-N7	-3.64	105.61	109.40
5	F	1502	5UA	C4-C5-N7	-3.58	105.67	109.40
6	E	103[A]	DU5	CAP-NBA-CAQ	3.37	116.42	108.83
6	E	103[A]	DU5	CAO-CAM-CAN	-3.35	100.83	113.84
5	E	102	5UA	C2-N1-C6	3.27	124.34	118.75
6	E	103[B]	DU5	CAL-CAG-NAT	-3.26	120.60	123.81
6	E	103[A]	DU5	CAL-CAG-NAT	-3.22	120.64	123.81
5	E	102	5UA	C5'-O5'-C6'	3.10	125.08	115.89
5	F	1502	5UA	C2-N1-C6	3.05	123.98	118.75
6	E	103[B]	DU5	CAP-NBA-CAQ	3.01	115.61	108.83
5	F	1502	5UA	O5'-C5'-C4'	2.83	114.43	108.43
6	E	103[B]	DU5	CAQ-CAS-NBB	2.58	115.71	110.70
5	F	1502	5UA	C5'-O5'-C6'	2.57	123.52	115.89
6	E	103[B]	DU5	CAO-CAM-CAN	-2.57	103.88	113.84
6	E	103[B]	DU5	CAS-CAQ-NBA	-2.53	105.45	110.64
6	E	103[A]	DU5	CAQ-CAS-NBB	2.49	115.54	110.70
6	E	103[B]	DU5	CAG-NAT-CAX	2.45	120.83	115.15
6	E	103[A]	DU5	CAG-NAT-CAX	2.43	120.78	115.15
6	E	103[A]	DU5	CAS-CAQ-NBA	-2.34	105.85	110.64
5	F	1502	5UA	O5'-C6'-OP3	-2.32	121.23	126.62
5	E	102	5UA	O5'-C6'-OP3	-2.30	121.29	126.62
6	E	103[B]	DU5	CAM-CAO-NBA	-2.26	108.14	113.84
6	E	103[A]	DU5	CAK-CAZ-CAX	-2.20	120.47	124.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	E	103[B]	DU5	CAK-CAZ-CAX	-2.04	120.80	124.83

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	102	5UA	O4'-C4'-C5'-O5'
5	E	102	5UA	C3'-C4'-C5'-O5'
6	E	103[A]	DU5	CAZ-CAX-NBB-CAR
6	E	103[A]	DU5	NAT-CAX-NBB-CAR
6	E	103[B]	DU5	CAZ-CAX-NBB-CAR
6	E	103[B]	DU5	NAT-CAX-NBB-CAR
6	E	103[B]	DU5	OAA-CAV-NAU-CAN
6	E	103[B]	DU5	CAW-CAV-NAU-CAN
6	E	103[A]	DU5	CAO-CAM-CAN-NAU
5	E	102	5UA	OP3-C6'-O5'-C5'
5	F	1502	5UA	OP3-C6'-O5'-C5'
6	E	103[A]	DU5	CAM-CAO-NBA-CAQ
6	E	103[B]	DU5	CAM-CAO-NBA-CAQ
6	E	103[A]	DU5	CAM-CAO-NBA-CAP
6	E	103[B]	DU5	CAM-CAO-NBA-CAP

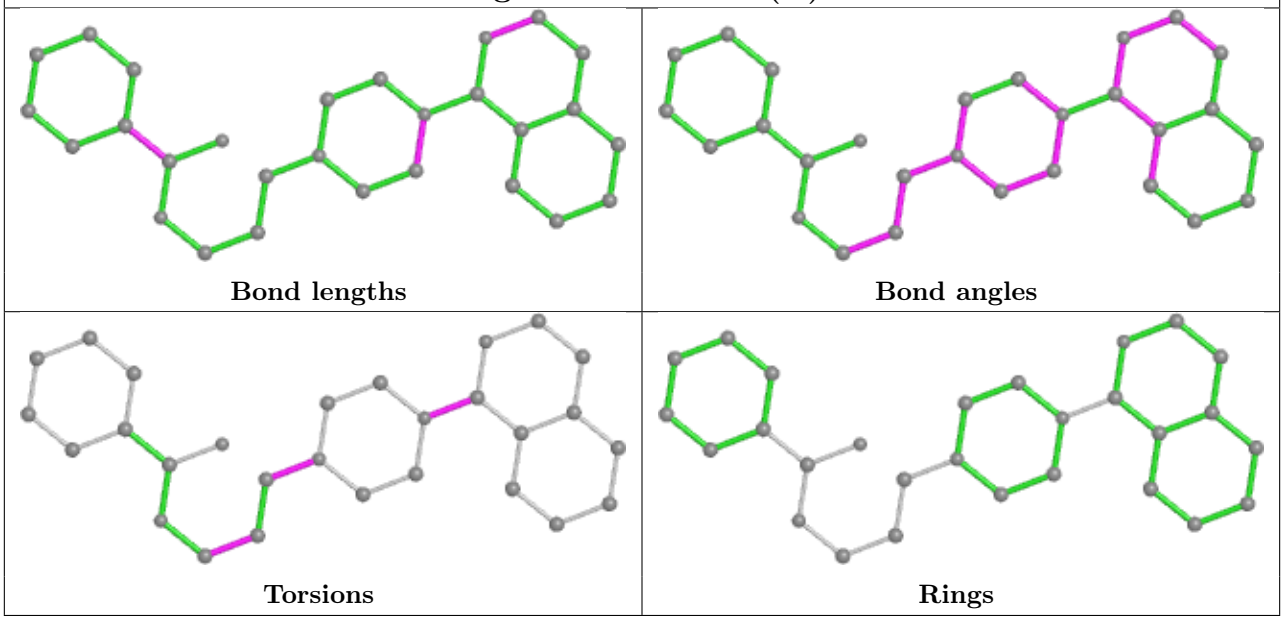
There are no ring outliers.

3 monomers are involved in 7 short contacts:

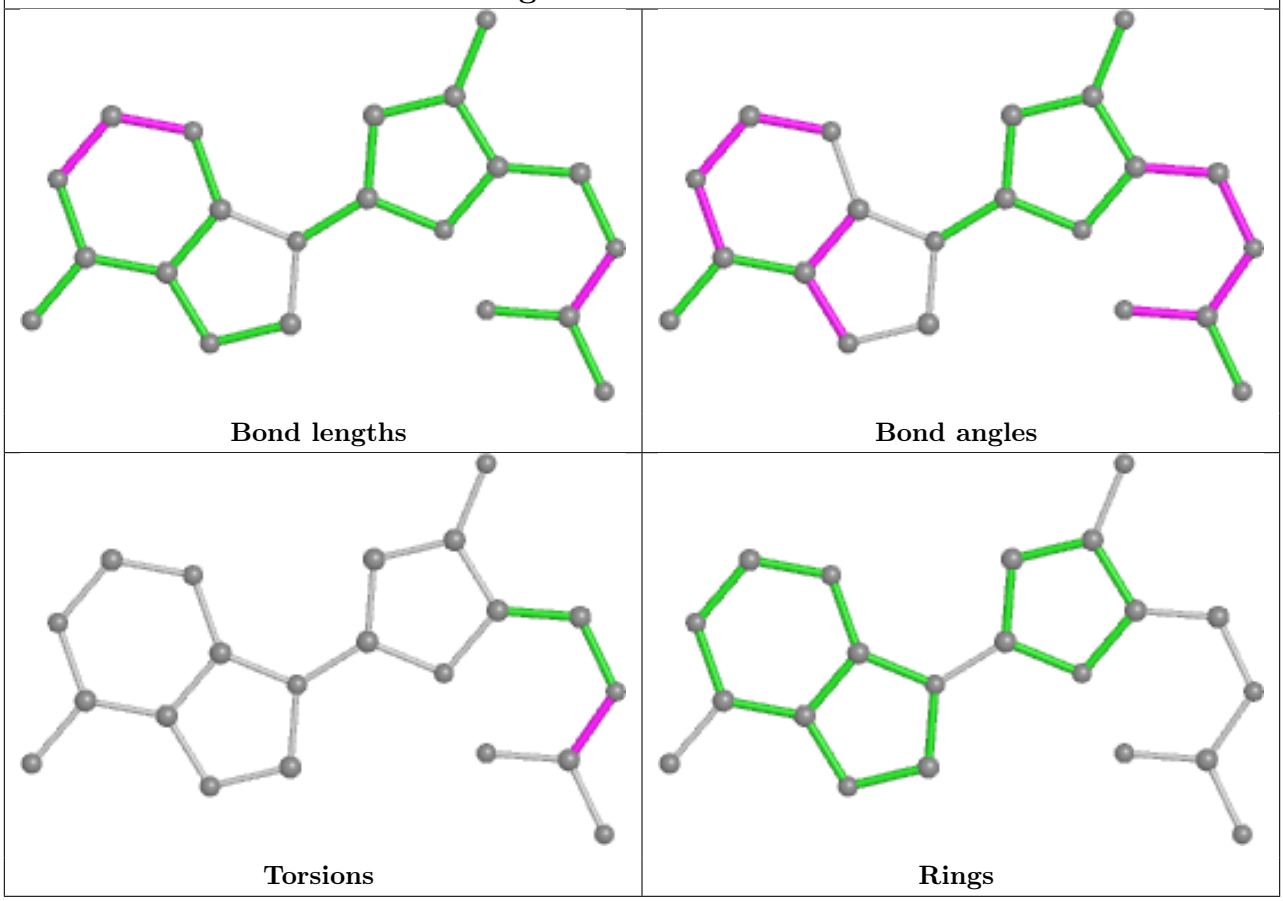
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	103[A]	DU5	3	0
5	E	102	5UA	0	1
6	E	103[B]	DU5	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

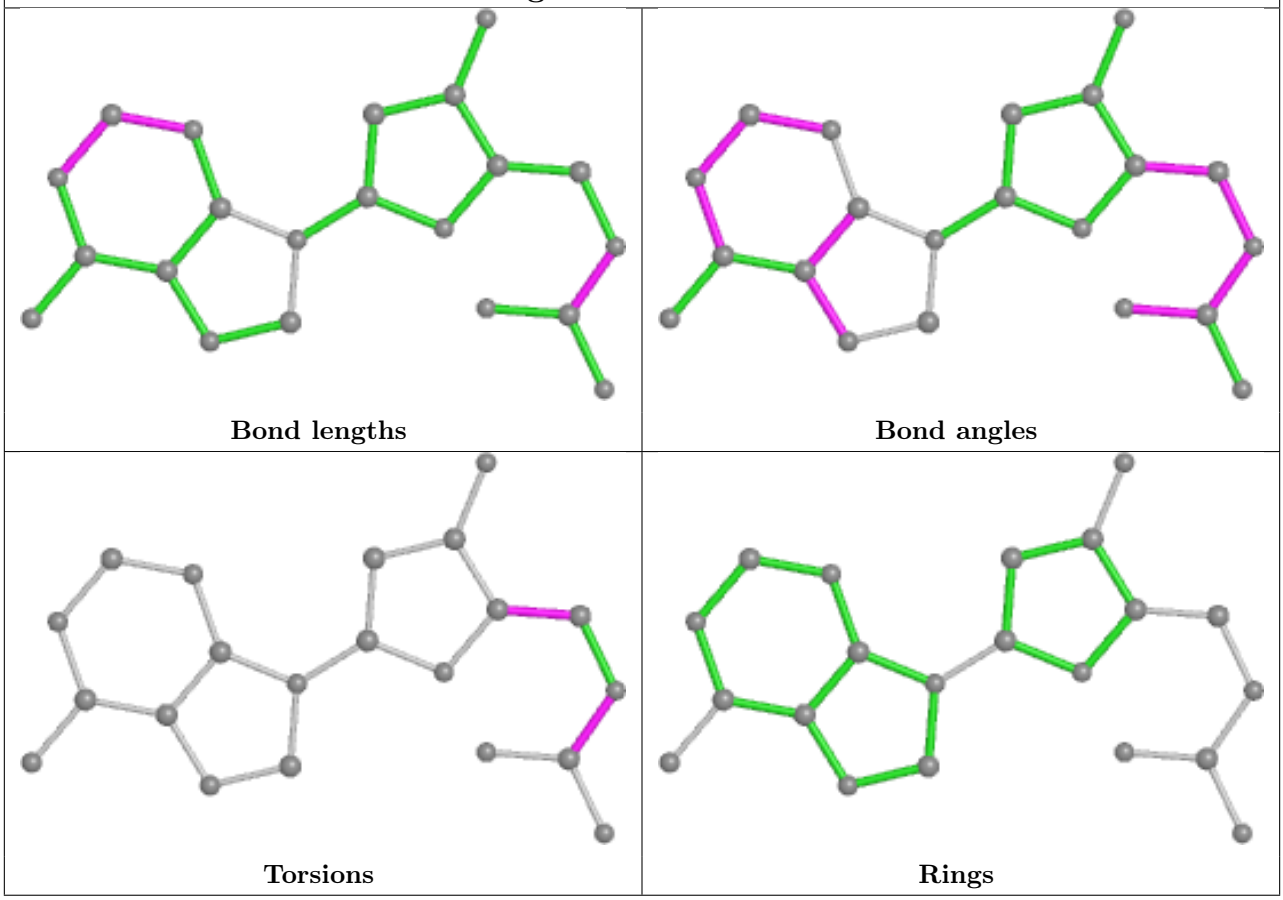
Ligand DU5 E 103 (A)



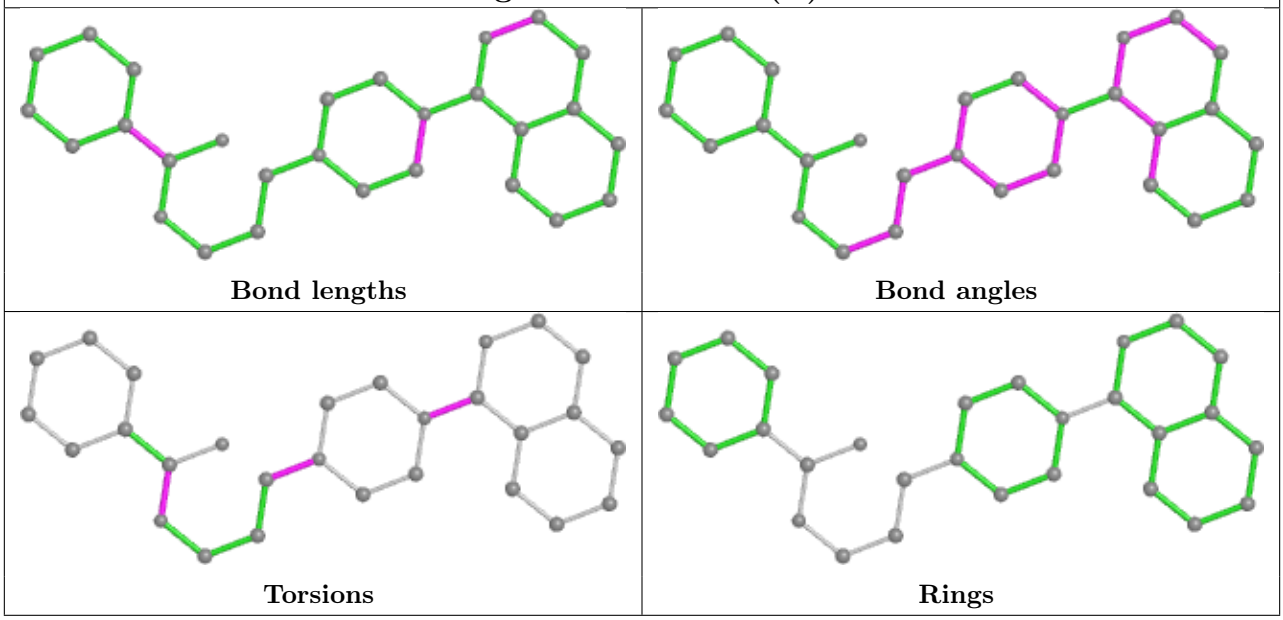
Ligand 5UA F 1502



Ligand 5UA E 102



Ligand DU5 E 103 (B)



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	D	1
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	545:GLY	C	580:TYR	N	3.21
1	B	545:GLY	C	580:TYR	N	3.09

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

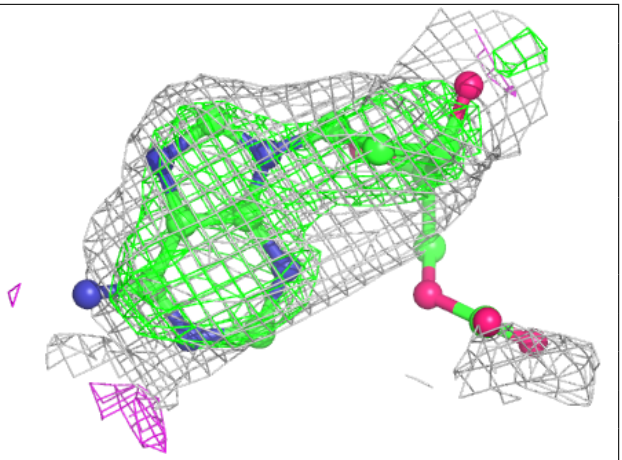
6.4 Ligands (i)

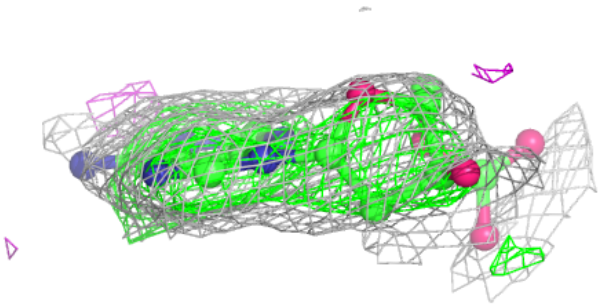
Unable to reproduce the depositors R factor - this section is therefore empty.

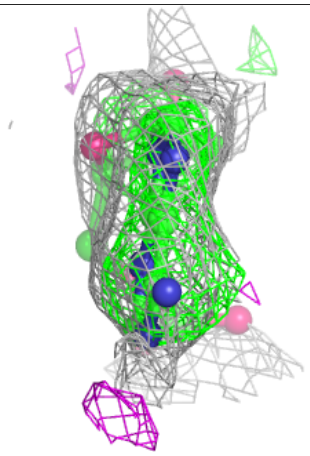
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 5UA E 102:

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

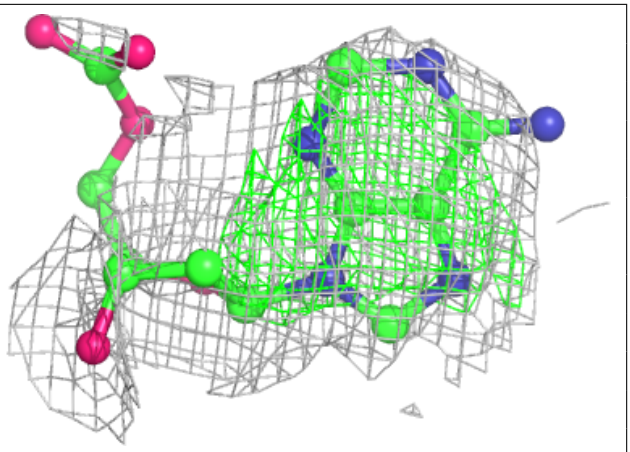


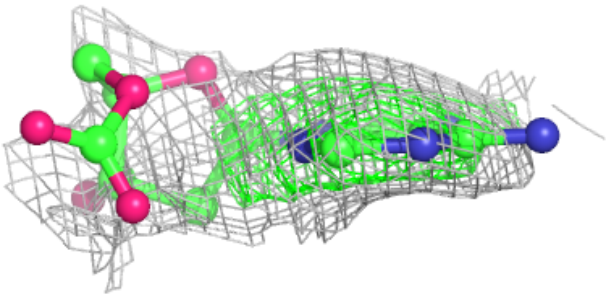


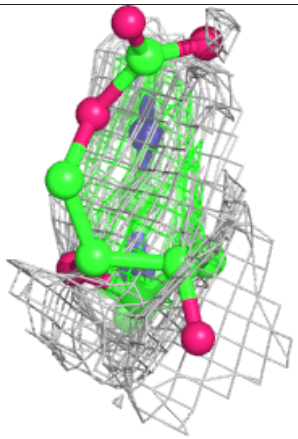


Electron density around 5UA F 1502:

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

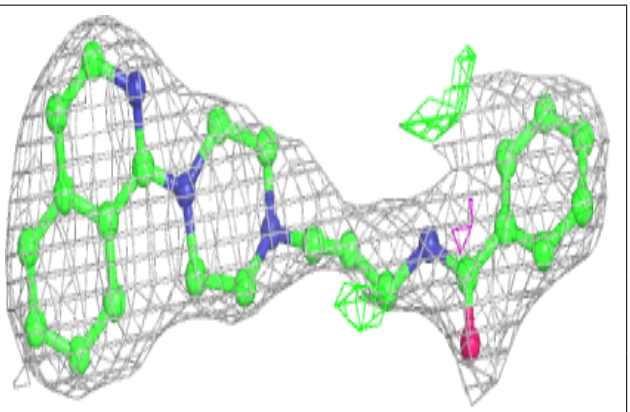


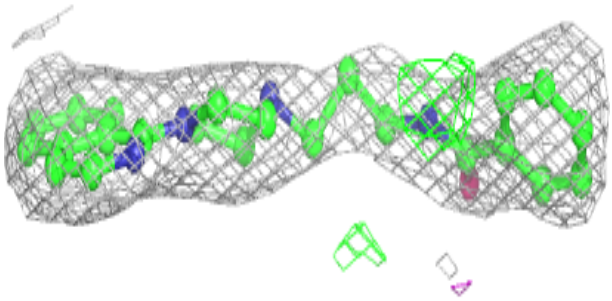


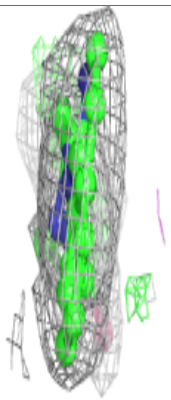


Electron density around DU5 E 103 (A):

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

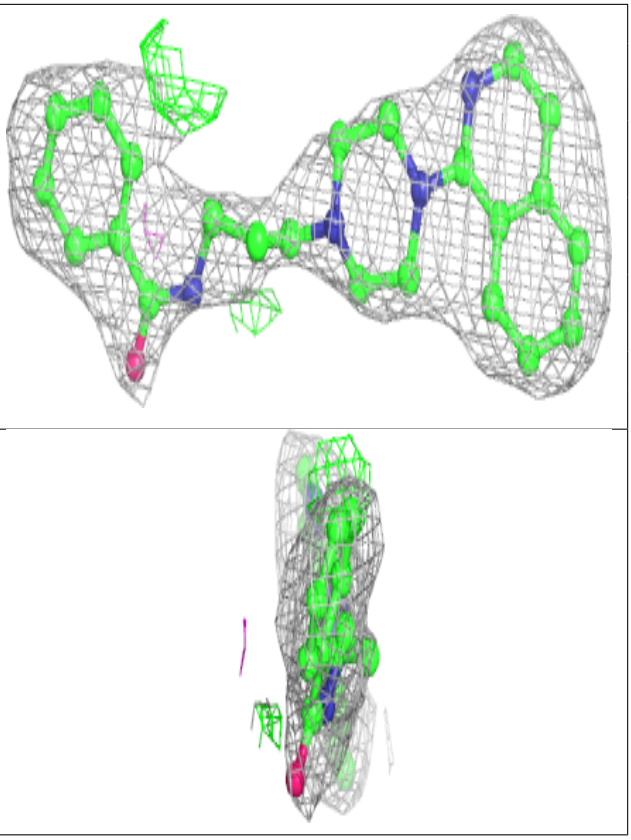






Electron density around DU5 E 103 (B):

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.