



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:22 AM UTC

PDB ID : 3UM8 / pdb_00003um8
Title : Wild-type Plasmodium falciparum DHFR-TS complexed with cycloguanil and NADPH
Authors : Vanichtanankul, J.; Chitnumsub, P.; Kamchonwongpaisan, S.; Yuthavong, Y.
Deposited on : 2011-11-12
Resolution : 2.60 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

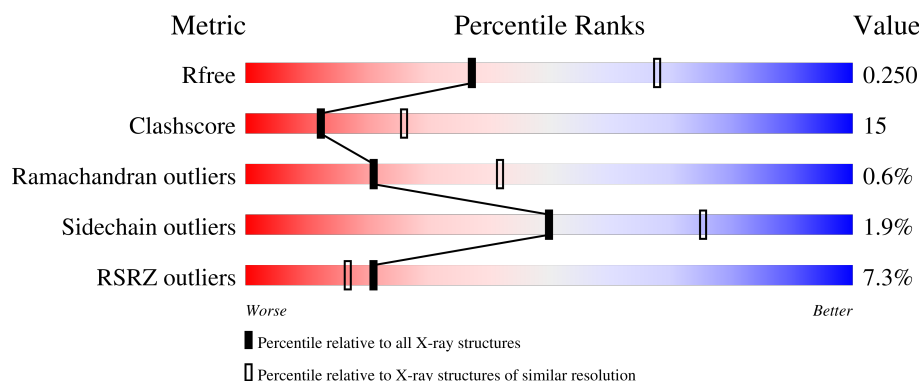
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	608	5% 61% 27% 10%
1	B	608	8% 59% 27% 11%

2 Entry composition [i](#)

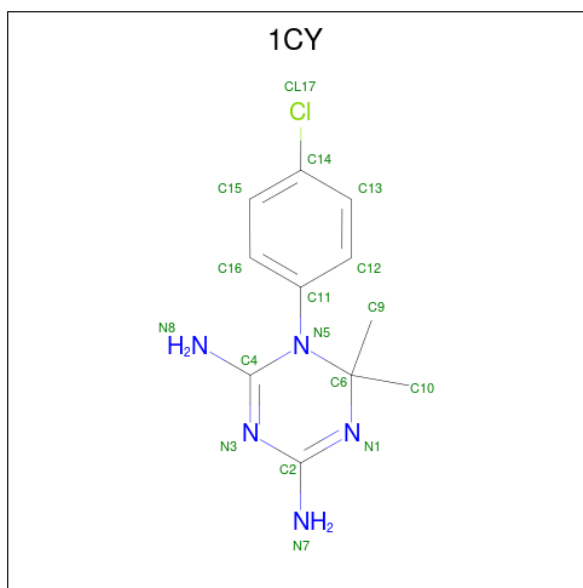
There are 5 unique types of molecules in this entry. The entry contains 9640 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional dihydrofolate reductase-thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	547	Total	C	N	O	S	0	0	0
			4543	2931	750	834	28			
1	B	542	Total	C	N	O	S	0	0	0
			4501	2906	742	825	28			

- Molecule 2 is 1-(4-chlorophenyl)-6,6-dimethyl-1,6-dihydro-1,3,5-triazine-2,4-diamine (CCD ID: 1CY) (formula: C₁₁H₁₄ClN₅).



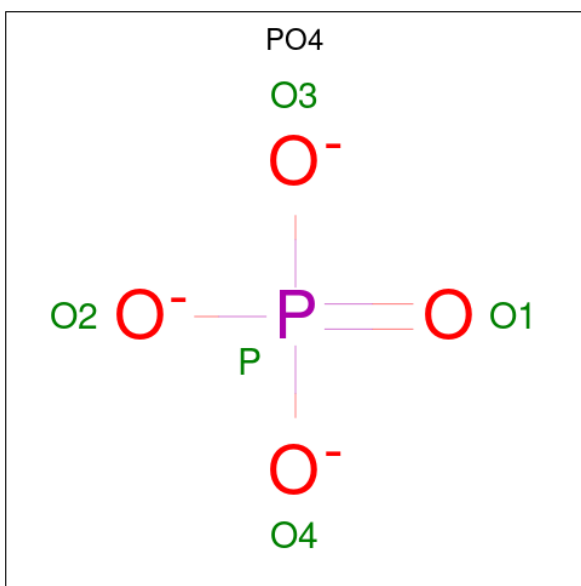
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	0	0
			17	11	1	5		
2	B	1	Total	C	Cl	N	0	0
			17	11	1	5		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	B	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

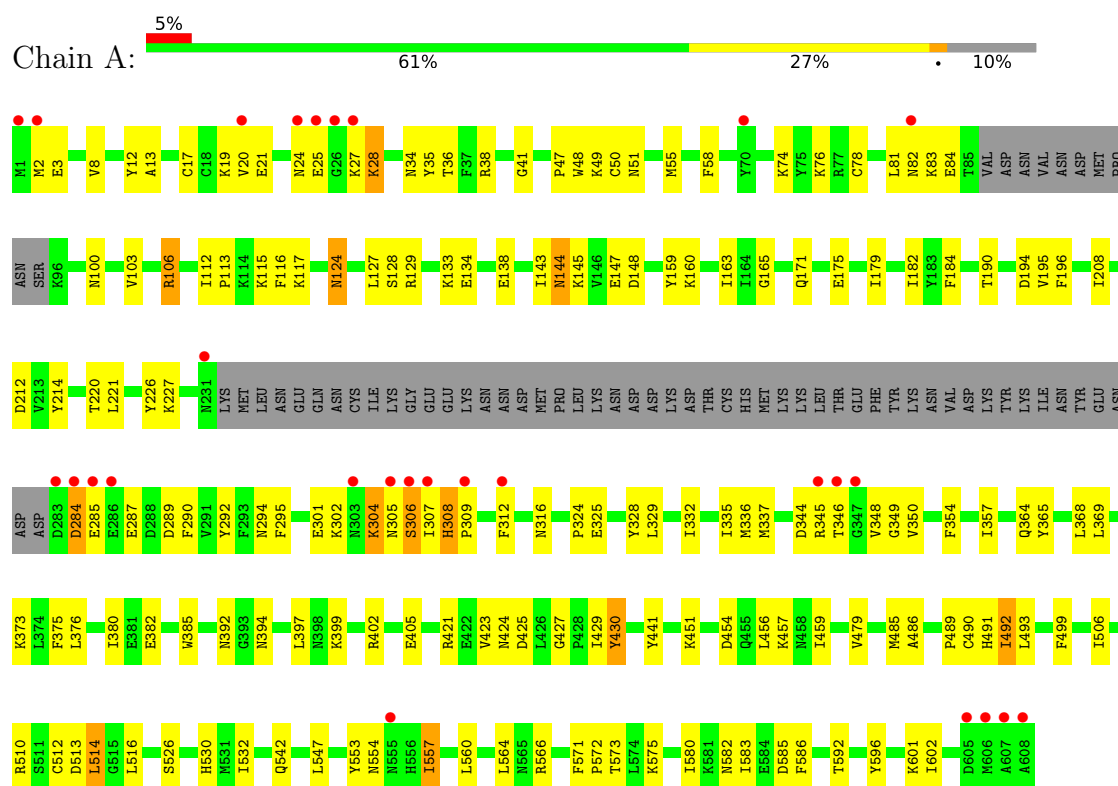
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	250	Total 250	O 250	0	0
5	B	206	Total 206	O 206	0	0

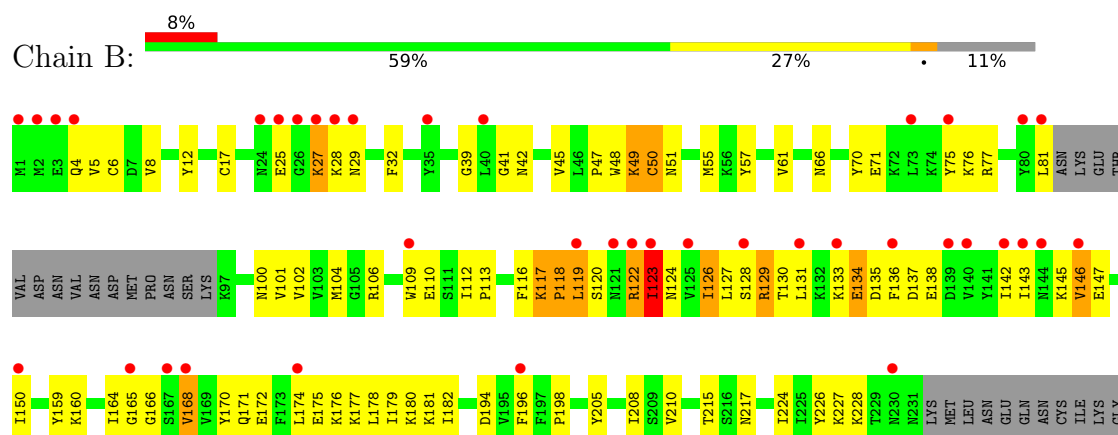
3 Residue-property plots

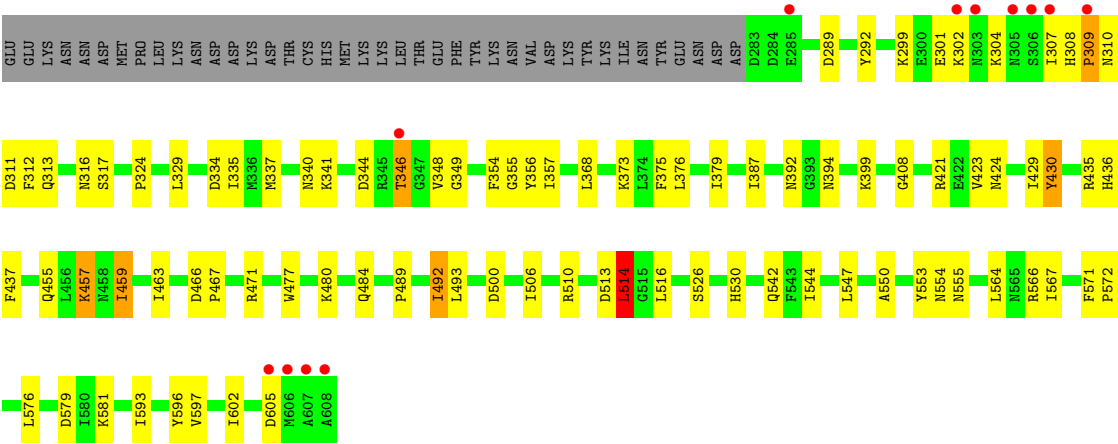
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Bifunctional dihydrofolate reductase-thymidylate synthase



- Molecule 1: Bifunctional dihydrofolate reductase-thymidylate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.76Å 157.54Å 164.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.16 – 2.60 39.16 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.4 (39.16-2.60) 98.3 (39.16-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 2.61Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.212 , 0.258 0.208 , 0.250	Depositor DCC
R_{free} test set	2369 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9640	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.46	0/4648	0.93	15/6272 (0.2%)
1	B	0.49	2/4606 (0.0%)	0.93	16/6217 (0.3%)
All	All	0.47	2/9254 (0.0%)	0.93	31/12489 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	117	LYS	C-N	-6.50	1.21	1.34
1	B	118	PRO	CG-CD	6.31	1.68	1.51

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	123	ILE	N-CA-C	-7.41	97.77	108.36
1	A	492	ILE	N-CA-C	7.24	117.34	110.53
1	B	492	ILE	N-CA-C	7.24	117.34	110.53
1	A	557	ILE	N-CA-C	6.86	117.57	110.36
1	B	130	THR	N-CA-C	-6.37	105.53	113.55
1	A	302	LYS	N-CA-C	6.34	120.58	112.34
1	B	70	TYR	N-CA-C	6.28	117.79	111.07
1	A	138	GLU	N-CA-C	6.18	118.69	110.53
1	B	25	GLU	N-CA-C	6.00	120.79	112.45
1	A	212	ASP	N-CA-C	-5.72	102.97	110.53
1	A	284	ASP	N-CA-C	5.69	119.94	112.89
1	A	308	HIS	CA-C-N	5.68	125.36	119.05
1	A	308	HIS	C-N-CA	5.68	125.36	119.05
1	A	542	GLN	N-CA-C	5.56	118.30	109.24
1	A	399	LYS	N-CA-C	-5.50	105.19	112.94
1	A	306	SER	N-CA-C	-5.47	106.66	113.38
1	B	118	PRO	N-CA-CB	5.33	108.46	102.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	486	ALA	N-CA-C	-5.28	104.94	111.33
1	B	119	LEU	N-CA-C	5.25	118.17	110.30
1	B	514	LEU	N-CA-C	5.23	117.72	111.71
1	B	368	LEU	N-CA-C	-5.22	100.79	109.46
1	B	117	LYS	CA-C-O	-5.22	115.81	120.19
1	A	129	ARG	N-CA-C	-5.21	106.25	113.18
1	B	27	LYS	N-CA-C	5.21	116.76	111.14
1	B	567	ILE	CA-C-N	5.17	125.16	119.89
1	B	567	ILE	C-N-CA	5.17	125.16	119.89
1	B	542	GLN	N-CA-C	5.14	118.19	109.76
1	B	526	SER	N-CA-C	-5.11	105.62	111.14
1	A	526	SER	N-CA-C	-5.08	105.82	111.36
1	A	514	LEU	N-CA-C	5.08	116.82	111.28
1	B	399	LYS	N-CA-C	-5.03	106.39	112.88

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4543	0	4497	132	0
1	B	4501	0	4452	154	0
2	A	17	0	14	1	0
2	B	17	0	14	0	0
3	A	48	0	26	3	0
3	B	48	0	26	6	0
4	A	10	0	0	1	0
5	A	250	0	0	5	0
5	B	206	0	0	8	0
All	All	9640	0	9029	280	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 15.

All (280) 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:376:LEU:HD22	1:B:379:ILE:HD11	1.41	1.00
1:B:102:VAL:HG22	1:B:123:ILE:O	1.64	0.97
1:B:145:LYS:HG2	1:B:146:VAL:H	1.28	0.97
1:B:119:LEU:HB3	1:B:122:ARG:CZ	1.95	0.96
1:A:127:LEU:HD23	1:A:143:ILE:HG13	1.53	0.90
1:A:165:GLY:HA3	3:A:610:NDP:H5N	1.57	0.86
1:A:451:LYS:NZ	5:A:1450:HOH:O	2.11	0.83
1:B:101:VAL:HA	1:B:123:ILE:HB	1.61	0.82
1:A:51:ASN:H	1:A:55:MET:HE3	1.46	0.81
1:B:165:GLY:HA3	3:B:710:NDP:H5N	1.59	0.81
1:B:51:ASN:O	1:B:55:MET:HG2	1.83	0.79
1:A:50:CYS:HA	1:A:55:MET:HE1	1.63	0.79
1:A:8:VAL:HA	1:A:76:LYS:HD3	1.65	0.78
1:A:51:ASN:H	1:A:55:MET:CE	1.99	0.76
1:A:346:THR:HG22	1:A:346:THR:O	1.83	0.76
1:A:28:LYS:NZ	1:A:28:LYS:HA	2.01	0.75
1:B:329:LEU:HD22	1:B:564:LEU:HD23	1.69	0.75
1:A:20:VAL:HG12	1:A:21:GLU:N	2.01	0.75
1:B:301:GLU:HB2	1:B:304:LYS:HB2	1.67	0.75
1:B:29:ASN:HB3	1:B:32:PHE:CZ	2.21	0.75
1:A:28:LYS:HA	1:A:28:LYS:HZ3	1.51	0.73
1:B:102:VAL:CG2	1:B:123:ILE:O	2.35	0.73
1:A:301:GLU:O	1:A:304:LYS:HD2	1.87	0.73
1:A:41:GLY:HA2	1:A:47:PRO:HD3	1.72	0.72
1:B:100:ASN:O	1:B:123:ILE:HB	1.90	0.72
1:B:145:LYS:HG2	1:B:146:VAL:N	2.03	0.72
1:A:78:CYS:HB3	1:A:83:LYS:O	1.90	0.71
1:B:210:VAL:HG12	1:B:224:ILE:HG22	1.72	0.70
1:A:106:ARG:HG3	3:A:610:NDP:O3X	1.91	0.70
1:B:120:SER:N	1:B:122:ARG:HH21	1.89	0.70
1:A:24:ASN:O	1:A:27:LYS:HG2	1.92	0.69
1:A:124:ASN:N	1:A:124:ASN:HD22	1.89	0.69
1:B:5:VAL:HG11	1:B:150:ILE:HD12	1.72	0.69
1:B:335:ILE:HD12	1:B:514:LEU:HD13	1.74	0.69
1:A:304:LYS:HD3	1:A:305:ASN:N	2.08	0.68
1:B:119:LEU:HB3	1:B:122:ARG:NH2	2.09	0.68
1:B:344:ASP:HB3	1:B:348:VAL:O	1.94	0.68
1:A:112:ILE:O	1:A:117:LYS:HE3	1.93	0.68
1:B:104:MET:HA	1:B:165:GLY:O	1.93	0.67
1:A:491:HIS:CE1	5:A:1413:HOH:O	2.46	0.67
1:A:106:ARG:HG3	1:A:106:ARG:HH11	1.60	0.67
1:A:20:VAL:CG1	1:A:21:GLU:N	2.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ASP:OD1	1:A:195:VAL:HG23	1.95	0.67
1:B:126:ILE:HD12	1:B:126:ILE:N	2.10	0.66
1:A:147:GLU:H	1:A:147:GLU:CD	2.04	0.66
1:B:500:ASP:O	5:B:1433:HOH:O	2.13	0.66
1:B:471:ARG:NH1	5:B:1334:HOH:O	2.26	0.66
1:A:165:GLY:HA3	3:A:610:NDP:C5N	2.27	0.65
1:B:182:ILE:HB	1:B:226:TYR:HB2	1.79	0.65
1:B:136:PHE:CD1	1:B:142:ILE:HD11	2.32	0.65
1:A:21:GLU:O	1:A:21:GLU:HG2	1.97	0.64
1:A:304:LYS:HD3	1:A:305:ASN:H	1.61	0.64
1:B:51:ASN:H	1:B:55:MET:HE3	1.61	0.64
1:A:145:LYS:HG2	1:A:148:ASP:OD2	1.98	0.64
1:A:336:MET:HE2	1:A:557:ILE:HG23	1.80	0.63
1:A:304:LYS:HD2	1:A:304:LYS:H	1.64	0.63
1:B:492:ILE:HD11	1:B:510:ARG:HD3	1.81	0.63
1:B:106:ARG:O	1:B:110:GLU:HG3	1.99	0.62
1:B:346:THR:HG22	1:B:348:VAL:HG22	1.82	0.60
1:A:328:TYR:CZ	1:A:332:ILE:HD11	2.35	0.60
1:B:29:ASN:HB3	1:B:32:PHE:CE1	2.36	0.60
1:B:131:LEU:HD22	1:B:136:PHE:CE1	2.37	0.60
1:B:27:LYS:HG3	1:B:28:LYS:N	2.17	0.59
1:B:346:THR:CG2	1:B:348:VAL:HG22	2.33	0.59
1:B:179:ILE:HB	1:B:205:TYR:OH	2.03	0.59
1:A:285:GLU:OE2	1:B:160:LYS:NZ	2.30	0.59
1:B:166:GLY:HA2	3:B:710:NDP:H5N	1.85	0.59
1:B:50:CYS:O	1:B:217:ASN:ND2	2.36	0.58
1:A:12:TYR:CE2	1:A:160:LYS:HD3	2.39	0.58
1:A:27:LYS:HG3	1:A:27:LYS:O	2.02	0.58
1:B:133:LYS:H	1:B:133:LYS:HD2	1.67	0.58
1:B:127:LEU:HA	1:B:143:ILE:HG13	1.84	0.58
1:B:42:ASN:HA	1:B:194:ASP:OD2	2.04	0.57
1:B:579:ASP:O	1:B:581:LYS:HG2	2.05	0.57
1:A:100:ASN:OD1	1:A:159:TYR:HB3	2.05	0.56
1:B:373:LYS:HE2	1:B:375:PHE:CE1	2.40	0.56
1:A:306:SER:O	1:A:307:ILE:HD13	2.06	0.56
1:B:109:TRP:O	1:B:117:LYS:HE2	2.05	0.56
1:A:289:ASP:HA	1:A:292:TYR:CD2	2.40	0.56
1:A:214:TYR:O	1:A:220:THR:HA	2.06	0.55
1:A:506:ILE:HG13	1:B:354:PHE:CE2	2.41	0.55
1:A:345:ARG:O	1:A:346:THR:HB	2.06	0.55
1:B:118:PRO:HB2	1:B:124:ASN:ND2	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:302:LYS:HZ3	1:B:341:LYS:H	1.53	0.55
1:B:349:GLY:C	1:B:554:ASN:ND2	2.64	0.55
1:B:166:GLY:HA2	3:B:710:NDP:C5N	2.37	0.55
1:A:13:ALA:HB2	1:A:179:ILE:HD12	1.89	0.55
1:A:485:MET:SD	1:A:489:PRO:HD3	2.47	0.55
1:B:168:VAL:O	1:B:172:GLU:HG2	2.06	0.55
1:B:553:TYR:HB3	1:B:555:ASN:OD1	2.07	0.55
1:A:159:TYR:CD2	1:A:160:LYS:HG3	2.42	0.55
1:A:329:LEU:HD22	1:A:564:LEU:HD23	1.88	0.54
1:B:457:LYS:NZ	1:B:457:LYS:HB3	2.23	0.54
1:A:350:VAL:HG12	1:A:553:TYR:CD1	2.42	0.54
1:A:376:LEU:O	1:A:380:ILE:HG13	2.07	0.54
1:B:493:LEU:HD12	1:B:493:LEU:C	2.32	0.54
1:B:48:TRP:O	1:B:49:LYS:HB3	2.07	0.54
1:B:109:TRP:CE2	1:B:117:LYS:HD3	2.42	0.54
1:A:81:LEU:O	1:A:82:ASN:HB2	2.08	0.54
1:A:301:GLU:OE1	1:A:337:MET:HB3	2.08	0.53
1:A:324:PRO:HB2	1:A:571:PHE:HE2	1.73	0.53
1:B:208:ILE:HD13	1:B:227:LYS:HB2	1.89	0.53
1:A:421:ARG:HD2	1:A:425:ASP:CG	2.34	0.53
1:B:168:VAL:HG23	3:B:710:NDP:O1N	2.07	0.53
1:B:308:HIS:HB2	1:B:311:ASP:OD1	2.08	0.53
1:A:499:PHE:CE1	1:B:340:ASN:HB3	2.44	0.53
1:A:513:ASP:OD2	1:A:516:LEU:HB2	2.08	0.53
1:A:368:LEU:HD13	1:A:376:LEU:HD11	1.90	0.53
1:B:506:ILE:HG12	1:B:544:ILE:HB	1.91	0.52
1:B:100:ASN:OD1	1:B:159:TYR:HB3	2.09	0.52
1:B:120:SER:O	1:B:122:ARG:NH2	2.42	0.52
1:B:547:LEU:HD12	1:B:547:LEU:N	2.25	0.52
1:B:210:VAL:HG12	1:B:224:ILE:CG2	2.39	0.52
1:B:109:TRP:CZ2	1:B:117:LYS:HD3	2.45	0.52
1:A:490:CYS:SG	5:A:1303:HOH:O	2.59	0.52
1:B:172:GLU:O	1:B:176:LYS:HG2	2.09	0.52
1:B:459:ILE:O	1:B:463:ILE:HG13	2.10	0.52
1:B:133:LYS:H	1:B:133:LYS:CD	2.22	0.52
1:B:102:VAL:HG21	1:B:122:ARG:HG2	1.93	0.51
1:B:477:TRP:HE3	1:B:489:PRO:HG2	1.75	0.51
1:A:2:MET:HG2	1:A:3:GLU:N	2.26	0.51
1:A:332:ILE:HD13	1:A:560:LEU:HD22	1.93	0.51
1:A:335:ILE:HD12	1:A:514:LEU:HD13	1.93	0.51
1:B:376:LEU:CD2	1:B:379:ILE:HD11	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:GLN:O	1:B:8:VAL:HG23	2.10	0.51
1:B:572:PRO:HB3	1:B:596:TYR:HA	1.92	0.51
1:A:171:GLN:O	1:A:175:GLU:HB2	2.10	0.50
1:A:382:GLU:O	1:A:385:TRP:HB3	2.11	0.50
1:B:324:PRO:HB2	1:B:571:PHE:HE2	1.76	0.50
1:B:480:LYS:HE2	5:B:1233:HOH:O	2.11	0.50
1:A:304:LYS:HD2	1:A:304:LYS:N	2.27	0.50
1:A:344:ASP:HB3	1:A:348:VAL:HB	1.94	0.50
1:A:373:LYS:HB3	1:A:601:LYS:HB3	1.94	0.50
1:B:48:TRP:O	1:B:49:LYS:CB	2.59	0.50
1:B:104:MET:HE2	1:B:164:ILE:HD11	1.92	0.50
1:A:423:VAL:O	1:A:424:ASN:HB2	2.11	0.50
1:A:115:LYS:HE3	1:A:116:PHE:CE2	2.48	0.49
1:A:582:ASN:HB3	1:A:585:ASP:OD2	2.11	0.49
1:B:41:GLY:HA2	1:B:47:PRO:HD3	1.94	0.49
1:A:392:ASN:OD1	1:A:394:ASN:HB2	2.13	0.49
1:A:530:HIS:CE1	5:A:1434:HOH:O	2.66	0.49
1:B:133:LYS:HD2	1:B:133:LYS:N	2.28	0.49
1:A:456:LEU:O	1:A:459:ILE:HG13	2.13	0.49
1:A:124:ASN:N	1:A:124:ASN:ND2	2.60	0.49
1:A:19:LYS:O	1:A:190:THR:HA	2.13	0.49
1:B:530:HIS:HB3	1:B:576:LEU:HD11	1.94	0.49
1:A:346:THR:O	1:A:346:THR:CG2	2.55	0.48
1:A:421:ARG:HG2	1:A:421:ARG:HH11	1.79	0.48
1:B:47:PRO:HB2	1:B:48:TRP:CE3	2.49	0.48
1:B:376:LEU:HD12	1:B:593:ILE:HG13	1.96	0.48
1:B:138:GLU:H	1:B:138:GLU:CD	2.22	0.48
1:A:58:PHE:CZ	2:A:609:1CY:H16	2.49	0.47
1:B:307:ILE:CG2	1:B:312:PHE:HE2	2.27	0.47
1:A:290:PHE:CE1	1:A:294:ASN:ND2	2.82	0.47
1:B:171:GLN:HA	1:B:198:PRO:HG2	1.96	0.47
1:B:423:VAL:O	1:B:424:ASN:HB2	2.14	0.47
1:B:109:TRP:CH2	1:B:118:PRO:HB3	2.49	0.47
1:A:144:ASN:C	1:A:144:ASN:HD22	2.21	0.47
1:B:113:PRO:HG2	1:B:116:PHE:HD2	1.79	0.47
1:A:128:SER:O	1:A:144:ASN:HA	2.14	0.47
1:A:290:PHE:O	1:A:294:ASN:ND2	2.46	0.47
1:B:12:TYR:CE1	1:B:180:LYS:HD3	2.49	0.47
1:B:355:GLY:HA2	1:B:547:LEU:O	2.15	0.47
1:A:159:TYR:CE2	1:A:160:LYS:HE3	2.49	0.47
1:A:402:ARG:HG2	1:A:402:ARG:HH11	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:532:ILE:HD13	1:A:583:ILE:HD13	1.96	0.47
1:A:575:LYS:HB2	1:A:592:THR:HB	1.96	0.47
1:B:32:PHE:CD1	1:B:597:VAL:HG13	2.50	0.47
1:B:309:PRO:C	1:B:311:ASP:H	2.22	0.47
1:B:312:PHE:HB2	1:B:316:ASN:HD22	1.80	0.47
1:A:20:VAL:HG21	1:A:196:PHE:HE1	1.80	0.46
1:B:17:CYS:HA	1:B:39:GLY:O	2.15	0.46
1:A:144:ASN:HD21	1:A:145:LYS:NZ	2.14	0.46
1:B:129:ARG:HG3	3:B:710:NDP:O1X	2.16	0.46
1:B:12:TYR:CD1	1:B:181:LYS:HB2	2.51	0.46
1:B:75:TYR:HD1	1:B:76:LYS:HG3	1.80	0.46
1:A:17:CYS:HB2	1:A:184:PHE:CE1	2.49	0.46
1:A:103:VAL:HB	1:A:163:ILE:HD13	1.96	0.46
1:B:313:GLN:O	1:B:317:SER:HB3	2.16	0.46
1:B:566:ARG:NH1	1:B:602:ILE:HD11	2.30	0.46
1:A:402:ARG:HG2	1:A:402:ARG:NH1	2.31	0.46
1:A:429:ILE:O	1:A:430:TYR:C	2.58	0.46
1:A:20:VAL:HG21	1:A:196:PHE:CE1	2.51	0.46
1:B:102:VAL:CG2	1:B:122:ARG:HG2	2.46	0.46
1:B:134:GLU:O	1:B:135:ASP:C	2.58	0.46
1:B:457:LYS:HB3	1:B:457:LYS:HZ2	1.81	0.46
1:B:120:SER:N	1:B:122:ARG:NH2	2.60	0.46
1:B:455:GLN:HA	5:B:1420:HOH:O	2.16	0.45
1:B:45:VAL:HB	5:B:1362:HOH:O	2.16	0.45
1:A:285:GLU:C	1:A:287:GLU:N	2.73	0.45
1:A:20:VAL:CG1	1:A:21:GLU:H	2.27	0.45
1:A:83:LYS:O	1:A:84:GLU:HG3	2.17	0.45
1:B:137:ASP:O	1:B:138:GLU:C	2.60	0.45
1:A:492:ILE:HD11	1:A:510:ARG:HD3	1.99	0.45
1:B:28:LYS:HD3	1:B:28:LYS:C	2.42	0.45
1:B:112:ILE:HB	1:B:117:LYS:HE2	1.98	0.45
1:B:307:ILE:O	1:B:307:ILE:HG22	2.16	0.44
1:B:484:GLN:HG3	5:B:1161:HOH:O	2.16	0.44
1:A:493:LEU:C	1:A:493:LEU:HD12	2.42	0.44
1:A:354:PHE:CE2	1:B:506:ILE:HG13	2.51	0.44
1:A:301:GLU:O	1:A:304:LYS:CD	2.62	0.44
1:A:304:LYS:N	1:A:304:LYS:CD	2.80	0.44
1:B:134:GLU:OE1	1:B:134:GLU:N	2.51	0.44
1:A:48:TRP:O	1:A:49:LYS:HB2	2.16	0.44
1:B:5:VAL:O	1:B:8:VAL:N	2.46	0.44
1:A:106:ARG:HH11	1:A:106:ARG:CG	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:GLU:HG3	1:A:369:LEU:HD22	2.00	0.44
1:A:479:VAL:HB	1:B:437:PHE:CD1	2.52	0.44
1:B:127:LEU:O	3:B:710:NDP:H1B	2.18	0.44
1:B:310:ASN:ND2	1:B:310:ASN:O	2.51	0.44
1:A:144:ASN:HD21	1:A:145:LYS:HZ2	1.66	0.44
1:B:119:LEU:C	1:B:122:ARG:HH21	2.25	0.44
1:B:131:LEU:HD22	1:B:136:PHE:CZ	2.53	0.44
1:B:334:ASP:OD2	1:B:356:TYR:OH	2.35	0.44
1:A:208:ILE:HD13	1:A:227:LYS:HB2	2.00	0.44
1:A:285:GLU:C	1:A:287:GLU:H	2.25	0.44
1:B:101:VAL:HG22	1:B:123:ILE:HG21	2.00	0.44
1:A:113:PRO:HB2	1:A:116:PHE:HD2	1.83	0.43
1:B:101:VAL:HG22	1:B:123:ILE:HD13	1.99	0.43
1:B:308:HIS:HB2	1:B:311:ASP:CG	2.44	0.43
1:A:349:GLY:C	1:A:554:ASN:OD1	2.62	0.43
1:A:357:ILE:HB	1:B:357:ILE:HD11	2.01	0.43
1:A:572:PRO:HB3	1:A:596:TYR:HA	2.01	0.43
1:A:221:LEU:N	1:A:221:LEU:HD23	2.33	0.43
1:B:71:GLU:OE1	1:B:71:GLU:HA	2.18	0.43
1:B:126:ILE:N	1:B:126:ILE:CD1	2.81	0.43
1:A:427:GLY:HA2	1:A:441:TYR:CE2	2.54	0.43
1:B:205:TYR:CD1	1:B:228:LYS:HA	2.53	0.43
1:B:387:ILE:O	1:B:435:ARG:NH1	2.52	0.43
1:B:196:PHE:CD1	1:B:196:PHE:N	2.86	0.42
1:A:182:ILE:HB	1:A:226:TYR:HB2	2.00	0.42
1:A:373:LYS:HE3	1:A:375:PHE:CZ	2.53	0.42
1:A:510:ARG:NH1	4:A:801:PO4:O2	2.51	0.42
1:B:75:TYR:HD1	1:B:76:LYS:CG	2.32	0.42
1:A:335:ILE:HD12	1:A:514:LEU:CD1	2.48	0.42
1:A:573:THR:HA	5:A:1159:HOH:O	2.18	0.42
1:B:57:TYR:O	1:B:61:VAL:HG23	2.19	0.42
1:B:75:TYR:CD1	1:B:75:TYR:C	2.97	0.42
1:B:429:ILE:O	1:B:430:TYR:C	2.62	0.42
1:B:170:TYR:O	1:B:174:LEU:HG	2.20	0.42
1:B:421:ARG:HH11	1:B:421:ARG:HG2	1.85	0.42
1:B:466:ASP:N	1:B:467:PRO:CD	2.83	0.42
1:A:454:ASP:OD2	1:A:457:LYS:HD2	2.20	0.42
1:B:175:GLU:O	1:B:177:LYS:HG3	2.20	0.42
1:B:301:GLU:OE1	1:B:337:MET:HE3	2.20	0.42
1:B:112:ILE:O	1:B:117:LYS:HE3	2.20	0.42
1:A:51:ASN:N	1:A:55:MET:HE3	2.24	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:HIS:HB3	1:B:309:PRO:HD2	2.01	0.42
1:A:35:TYR:O	1:A:38:ARG:NH1	2.48	0.41
1:B:77:ARG:O	1:B:81:LEU:HG	2.19	0.41
1:B:513:ASP:OD2	1:B:516:LEU:HB3	2.19	0.41
1:A:34:ASN:C	1:A:36:THR:H	2.29	0.41
1:B:514:LEU:HD11	1:B:550:ALA:HB1	2.02	0.41
1:A:397:LEU:HD22	1:A:402:ARG:NH1	2.35	0.41
1:B:392:ASN:OD1	1:B:394:ASN:HB2	2.20	0.41
1:A:34:ASN:C	1:A:36:THR:N	2.78	0.41
1:A:312:PHE:HB2	1:A:316:ASN:ND2	2.36	0.41
1:A:512:CYS:SG	1:A:547:LEU:HD22	2.60	0.41
1:B:6:CYS:HB3	1:B:178:LEU:C	2.45	0.41
1:B:66:ASN:C	1:B:66:ASN:HD22	2.28	0.41
1:B:309:PRO:C	1:B:311:ASP:N	2.77	0.41
1:A:74:LYS:O	1:A:78:CYS:HB2	2.21	0.41
1:B:572:PRO:HD3	5:B:1042:HOH:O	2.20	0.41
1:A:292:TYR:O	1:A:295:PHE:HB3	2.21	0.41
1:A:566:ARG:CZ	1:A:602:ILE:HD11	2.51	0.41
1:B:32:PHE:CG	1:B:597:VAL:HG13	2.56	0.41
1:B:106:ARG:HB2	1:B:128:SER:HB2	2.03	0.41
1:B:289:ASP:HA	1:B:292:TYR:CD2	2.55	0.41
1:B:436:HIS:O	1:B:437:PHE:C	2.63	0.41
1:B:605:ASP:HB3	5:B:1169:HOH:O	2.20	0.41
1:A:50:CYS:CA	1:A:55:MET:HE1	2.42	0.41
1:A:308:HIS:HB3	1:A:309:PRO:HD2	2.02	0.41
1:B:408:GLY:O	1:B:423:VAL:HG13	2.21	0.40
1:A:133:LYS:HE2	1:A:134:GLU:OE2	2.22	0.40
1:B:335:ILE:CD1	1:B:514:LEU:HD13	2.47	0.40
1:A:402:ARG:NH1	1:A:405:GLU:OE2	2.54	0.40
1:A:580:ILE:HG21	1:A:586:PHE:CG	2.56	0.40
1:A:144:ASN:ND2	1:A:144:ASN:H	2.20	0.40
1:A:364:GLN:O	1:A:365:TYR:HB3	2.20	0.40
1:B:299:LYS:HA	1:B:299:LYS:HD2	1.88	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	541/608 (89%)	503 (93%)	36 (7%)	2 (0%)	30	51
1	B	536/608 (88%)	490 (91%)	42 (8%)	4 (1%)	18	38
All	All	1077/1216 (89%)	993 (92%)	78 (7%)	6 (1%)	21	42

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	TYR
1	B	49	LYS
1	B	430	TYR
1	A	25	GLU
1	B	309	PRO
1	B	146	VAL

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	510/570 (90%)	504 (99%)	6 (1%)	63	83
1	B	505/570 (89%)	492 (97%)	13 (3%)	40	68
All	All	1015/1140 (89%)	996 (98%)	19 (2%)	50	75

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	106	ARG
1	A	124	ASN
1	A	144	ASN
1	A	284	ASP
1	A	304	LYS
1	B	50	CYS
1	B	122	ARG
1	B	123	ILE
1	B	126	ILE
1	B	129	ARG
1	B	134	GLU
1	B	147	GLU
1	B	168	VAL
1	B	215	THR
1	B	346	THR
1	B	457	LYS
1	B	459	ILE
1	B	514	LEU

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

Mol	Chain	Res	Type
1	A	24	ASN
1	A	99	GLN
1	A	144	ASN
1	A	157	ASN
1	A	201	ASN
1	A	294	ASN
1	A	394	ASN
1	A	424	ASN
1	B	4	GLN
1	B	24	ASN
1	B	66	ASN
1	B	99	GLN
1	B	124	ASN
1	B	310	ASN
1	B	313	GLN
1	B	316	ASN
1	B	340	ASN
1	B	394	ASN
1	B	407	ASN
1	B	424	ASN

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Mol	Chain	Res	Type
1	B	509	GLN
1	B	521	ASN
1	B	530	HIS
1	B	554	ASN
1	B	582	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 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	$\# Z > 2$
4	PO4	A	802	-	4,4,4	1.81	1 (25%)	6,6,6	0.46	0
4	PO4	A	801	-	4,4,4	1.82	3 (75%)	6,6,6	0.46	0
2	1CY	A	609	-	16,18,18	3.58	12 (75%)	18,27,27	1.30	2 (11%)
3	NDP	B	710	-	51,52,52	1.48	8 (15%)	71,80,80	1.68	12 (16%)
3	NDP	A	610	-	51,52,52	1.43	8 (15%)	71,80,80	1.73	15 (21%)
2	1CY	B	709	-	16,18,18	3.70	11 (68%)	18,27,27	1.62	3 (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	1CY	B	709	-	-	0/4/23/23	0/2/2/2
3	NDP	A	610	-	-	2/34/77/77	0/5/5/5
3	NDP	B	710	-	-	5/34/77/77	0/5/5/5
2	1CY	A	609	-	-	0/4/23/23	0/2/2/2

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	609	1CY	C16-C11	6.84	1.52	1.39
2	B	709	1CY	C16-C11	6.23	1.51	1.39
2	A	609	1CY	C15-C14	5.71	1.48	1.38
2	B	709	1CY	C15-C14	5.52	1.48	1.38
2	B	709	1CY	C4-N5	5.42	1.47	1.37
3	B	710	NDP	C6N-N1N	5.08	1.49	1.37
2	B	709	1CY	C16-C15	5.01	1.46	1.38
3	A	610	NDP	C6N-N1N	4.74	1.48	1.37
2	A	609	1CY	C12-C11	4.49	1.47	1.39
2	A	609	1CY	C14-CL17	-4.46	1.64	1.74
2	B	709	1CY	C12-C11	4.22	1.47	1.39
2	B	709	1CY	C13-C14	4.08	1.45	1.38
2	A	609	1CY	C2-N3	4.03	1.44	1.36
2	B	709	1CY	C2-N3	4.02	1.44	1.36
2	A	609	1CY	C4-N5	3.89	1.44	1.37
3	B	710	NDP	C8A-N7A	3.73	1.38	1.31
2	B	709	1CY	C4-N8	3.73	1.40	1.34
2	A	609	1CY	C13-C14	3.36	1.44	1.38
3	A	610	NDP	C8A-N7A	3.30	1.38	1.31
2	B	709	1CY	C13-C12	3.22	1.44	1.38
2	A	609	1CY	C16-C15	3.21	1.44	1.38
3	A	610	NDP	C7N-C3N	3.08	1.55	1.48
2	A	609	1CY	C13-C12	2.93	1.43	1.38
3	A	610	NDP	C4A-N3A	2.93	1.39	1.34
3	B	710	NDP	C7N-C3N	2.90	1.54	1.48
2	A	609	1CY	C2-N7	2.83	1.42	1.34
3	B	710	NDP	C4A-N3A	2.77	1.39	1.34
2	B	709	1CY	C2-N7	2.72	1.41	1.34
2	A	609	1CY	C4-N8	2.58	1.38	1.34
3	B	710	NDP	C5D-C4D	2.51	1.59	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	710	NDP	C5A-N7A	-2.43	1.34	1.39
2	B	709	1CY	C14-CL17	-2.43	1.68	1.74
3	A	610	NDP	C5A-N7A	-2.35	1.34	1.39
3	A	610	NDP	O4D-C1D	2.30	1.47	1.42
3	B	710	NDP	O4D-C1D	2.21	1.47	1.42
4	A	802	PO4	P-O4	-2.19	1.48	1.54
4	A	801	PO4	P-O2	-2.16	1.48	1.54
3	A	610	NDP	C5D-C4D	2.12	1.57	1.51
3	A	610	NDP	PA-O2A	-2.11	1.45	1.55
3	B	710	NDP	C1D-N1N	2.08	1.52	1.46
4	A	801	PO4	P-O4	-2.01	1.48	1.54
4	A	801	PO4	P-O3	-2.01	1.48	1.54
2	A	609	1CY	C10-C6	2.00	1.58	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	710	NDP	N3A-C2A-N1A	-6.37	118.93	128.58
3	A	610	NDP	N3A-C2A-N1A	-5.84	119.74	128.58
3	A	610	NDP	C1D-N1N-C2N	-4.68	113.43	121.14
3	B	710	NDP	C5A-C4A-N3A	-4.26	120.85	126.72
3	A	610	NDP	C5A-C4A-N3A	-4.16	120.98	126.72
2	B	709	1CY	C10-C6-N5	3.87	113.13	110.24
3	B	710	NDP	O4B-C1B-C2B	-3.68	100.26	106.59
2	B	709	1CY	C10-C6-C9	-3.58	104.75	110.68
3	A	610	NDP	O2B-C2B-C3B	-3.52	99.06	111.68
3	B	710	NDP	C1B-N9A-C8A	3.45	134.76	127.09
3	B	710	NDP	C2A-N3A-C4A	3.27	119.81	111.83
3	B	710	NDP	O2B-C2B-C3B	-3.25	100.00	111.68
2	A	609	1CY	C10-C6-C9	-3.18	105.41	110.68
3	B	710	NDP	C2A-N1A-C6A	3.10	123.82	118.73
3	A	610	NDP	C2A-N1A-C6A	3.06	123.76	118.73
3	A	610	NDP	O4B-C1B-C2B	-3.04	101.35	106.59
3	A	610	NDP	C2A-N3A-C4A	2.90	118.92	111.83
2	B	709	1CY	C16-C11-N5	2.85	123.26	120.08
3	B	710	NDP	C4A-N9A-C1B	-2.79	120.12	126.63
3	B	710	NDP	P2B-O2B-C2B	-2.62	116.43	123.43
2	A	609	1CY	C10-C6-N5	2.58	112.17	110.24
3	A	610	NDP	C6N-N1N-C2N	2.54	122.04	119.32
3	B	710	NDP	C5A-C4A-N9A	2.46	108.49	105.81
3	A	610	NDP	C1B-N9A-C8A	2.34	132.29	127.09
3	A	610	NDP	C2D-C1D-N1N	2.33	119.03	113.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	610	NDP	N3A-C4A-N9A	2.30	131.08	127.17
3	A	610	NDP	O3X-P2B-O2B	2.29	114.79	105.85
3	A	610	NDP	C4A-N9A-C1B	-2.22	121.45	126.63
3	A	610	NDP	C3N-C2N-N1N	-2.18	120.01	123.20
3	B	710	NDP	C3N-C2N-N1N	-2.13	120.08	123.20
3	B	710	NDP	N3A-C4A-N9A	2.04	130.64	127.17
3	A	610	NDP	N9A-C8A-N7A	-2.04	111.04	113.94

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	710	NDP	O4D-C1D-N1N-C2N
3	A	610	NDP	O4D-C1D-N1N-C2N
3	A	610	NDP	C2N-C3N-C7N-N7N
3	B	710	NDP	C2N-C3N-C7N-N7N
3	B	710	NDP	O4D-C4D-C5D-O5D
3	B	710	NDP	C3D-C4D-C5D-O5D
3	B	710	NDP	C2B-O2B-P2B-O1X

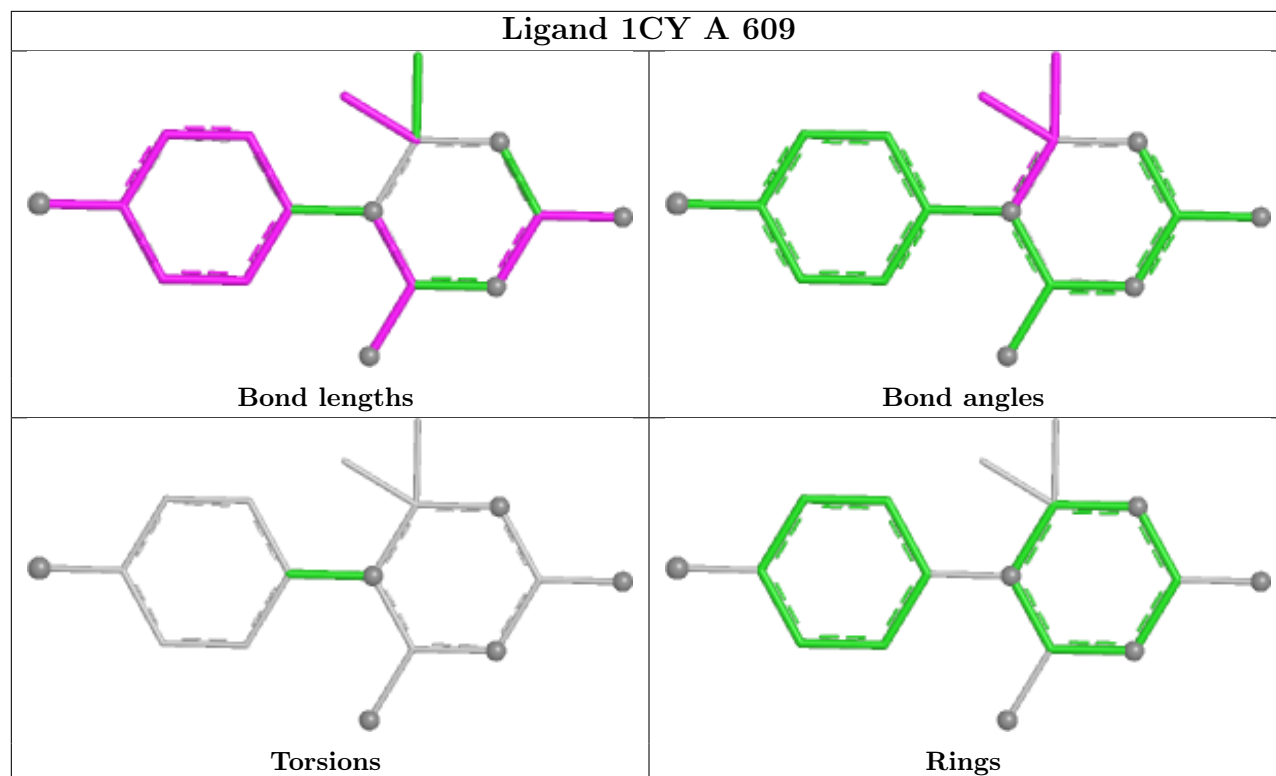
There are no ring outliers.

4 monomers are involved in 11 short contacts:

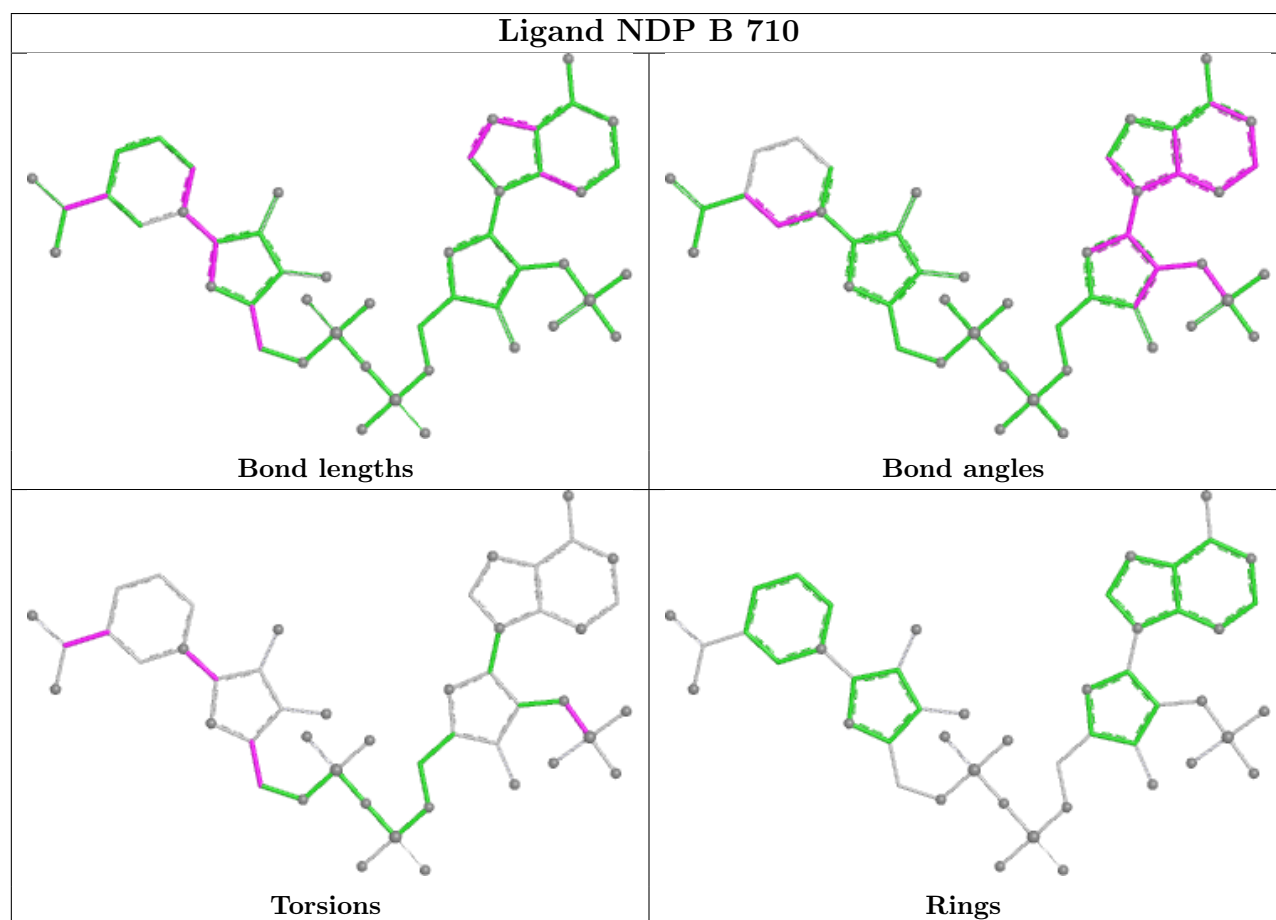
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	PO4	1	0
2	A	609	1CY	1	0
3	B	710	NDP	6	0
3	A	610	NDP	3	0

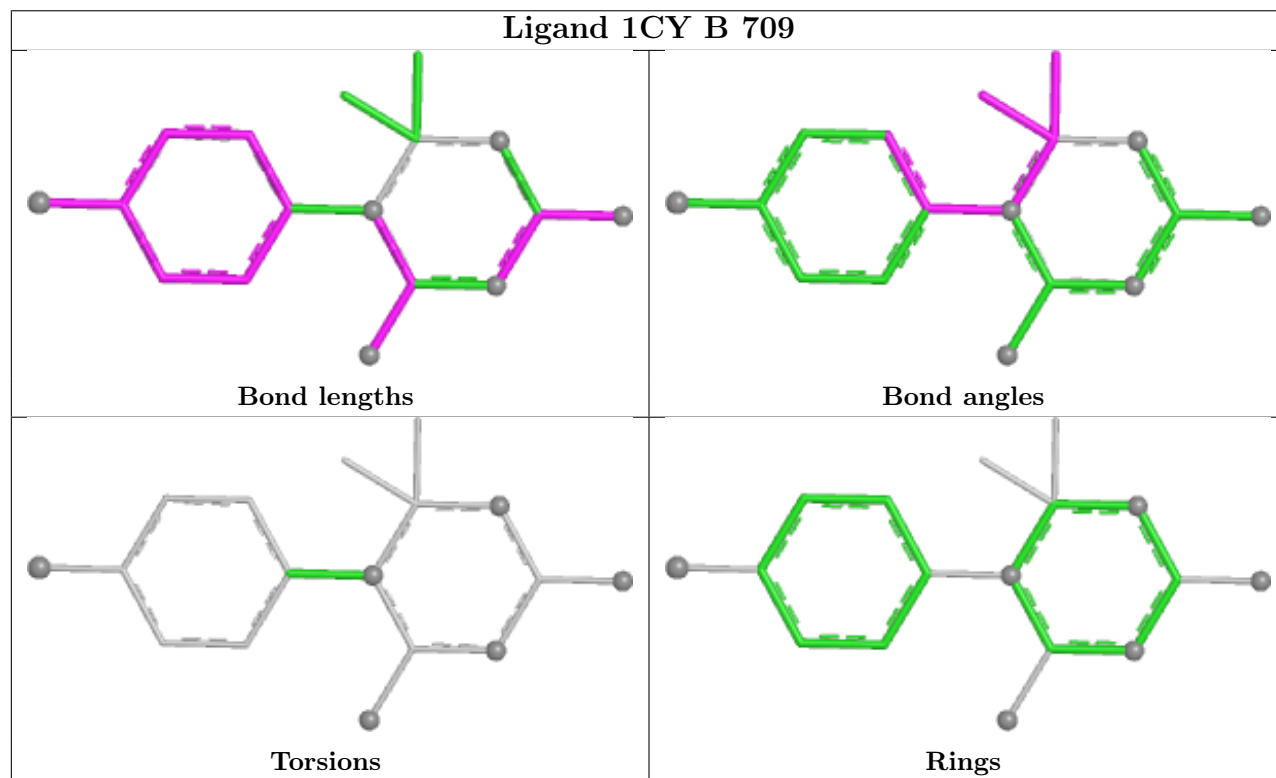
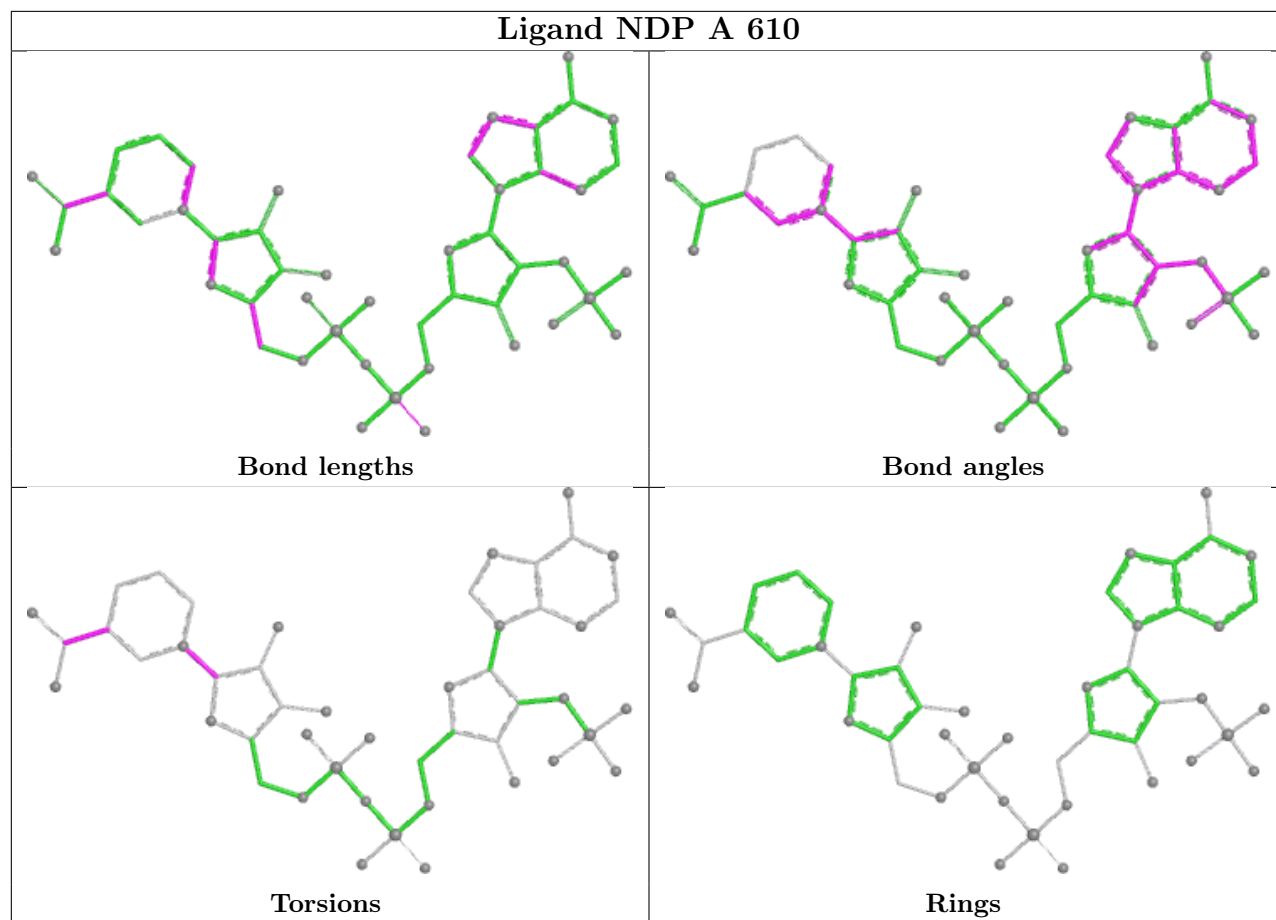
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand 1CY A 609



Ligand NDP B 710





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	547/608 (89%)	0.09	28 (5%)	33 28	23, 42, 90, 91	0
1	B	542/608 (89%)	0.35	51 (9%)	14 11	22, 48, 91, 91	0
All	All	1089/1216 (89%)	0.22	79 (7%)	21 17	22, 44, 91, 91	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	607	ALA	6.0
1	A	607	ALA	5.9
1	A	307	ILE	5.5
1	B	123	ILE	5.4
1	B	307	ILE	5.2
1	B	26	GLY	4.9
1	B	27	LYS	4.7
1	A	346	THR	4.3
1	A	608	ALA	4.2
1	A	286	GLU	4.0
1	A	305	ASN	4.0
1	A	283	ASP	3.8
1	B	608	ALA	3.7
1	B	28	LYS	3.7
1	B	2	MET	3.6
1	B	165	GLY	3.5
1	A	347	GLY	3.5
1	B	121	ASN	3.4
1	A	285	GLU	3.3
1	B	29	ASN	3.3
1	A	2	MET	3.2
1	A	309	PRO	3.2
1	B	306	SER	3.2
1	A	24	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	167	SER	3.2
1	B	230	ASN	3.1
1	B	346	THR	3.1
1	B	122	ARG	3.1
1	B	606	MET	3.1
1	B	128	SER	3.0
1	A	284	ASP	2.9
1	A	26	GLY	2.9
1	B	75	TYR	2.8
1	B	309	PRO	2.8
1	A	605	ASP	2.8
1	B	146	VAL	2.8
1	B	142	ILE	2.7
1	B	139	ASP	2.6
1	B	1	MET	2.6
1	B	80	TYR	2.6
1	B	302	LYS	2.6
1	A	82	ASN	2.5
1	B	303	ASN	2.5
1	A	27	LYS	2.5
1	B	605	ASP	2.5
1	B	125	VAL	2.5
1	B	143	ILE	2.4
1	A	606	MET	2.4
1	A	70	TYR	2.3
1	B	40	LEU	2.3
1	B	140	VAL	2.3
1	B	305	ASN	2.3
1	A	312	PHE	2.3
1	B	131	LEU	2.3
1	B	24	ASN	2.3
1	B	4	GLN	2.3
1	A	1	MET	2.2
1	B	133	LYS	2.2
1	B	3	GLU	2.2
1	B	25	GLU	2.2
1	A	303	ASN	2.2
1	B	35	TYR	2.2
1	A	345	ARG	2.2
1	B	136	PHE	2.2
1	B	144	ASN	2.1
1	B	168	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	25	GLU	2.1
1	B	285	GLU	2.1
1	B	174	LEU	2.1
1	A	231	ASN	2.1
1	B	196	PHE	2.1
1	B	109	TRP	2.1
1	B	73	LEU	2.1
1	A	20	VAL	2.1
1	A	306	SER	2.0
1	B	81	LEU	2.0
1	A	555	ASN	2.0
1	B	150	ILE	2.0
1	B	119	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

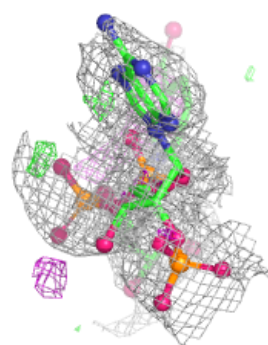
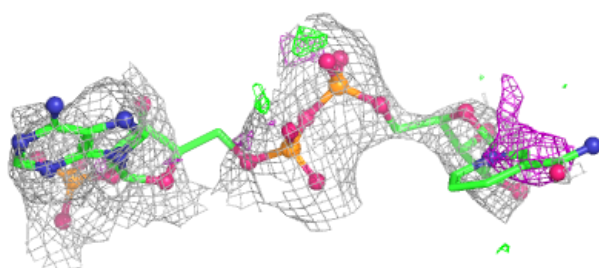
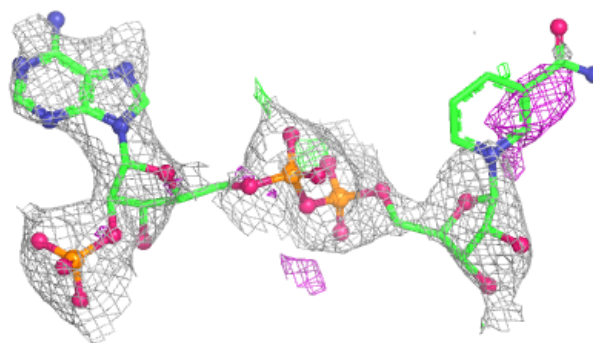
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NDP	B	710	48/48	0.77	0.16	88,90,90,90	0
2	1CY	B	709	17/17	0.82	0.20	83,84,89,90	0
4	PO4	A	801	5/5	0.84	0.17	77,78,78,79	0
4	PO4	A	802	5/5	0.87	0.19	63,65,66,67	0
3	NDP	A	610	48/48	0.93	0.10	47,53,58,60	0
2	1CY	A	609	17/17	0.95	0.08	27,31,38,44	0

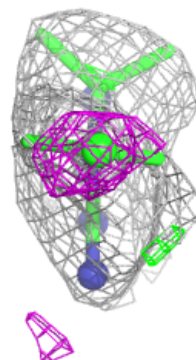
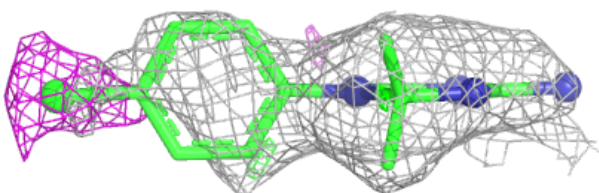
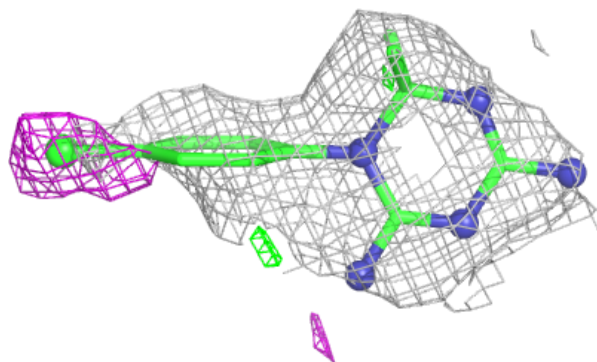
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NDP B 710:

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

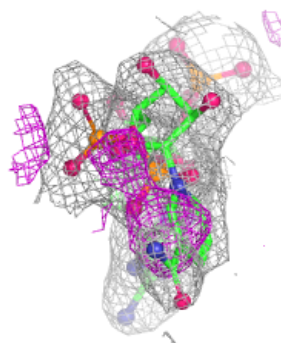
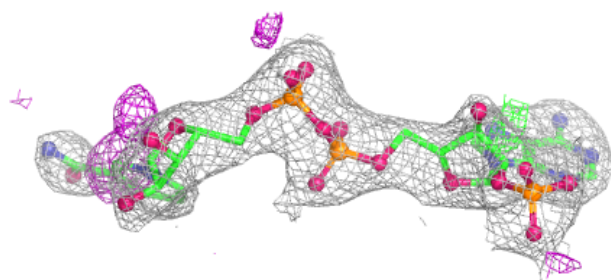
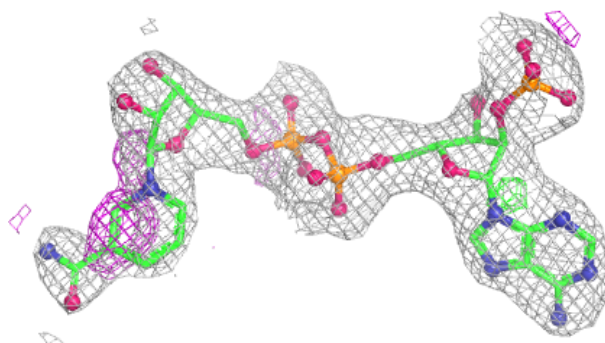
**Electron density around 1CY B 709:**

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

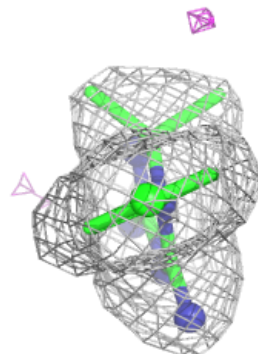
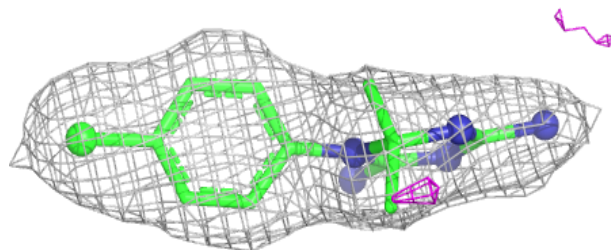
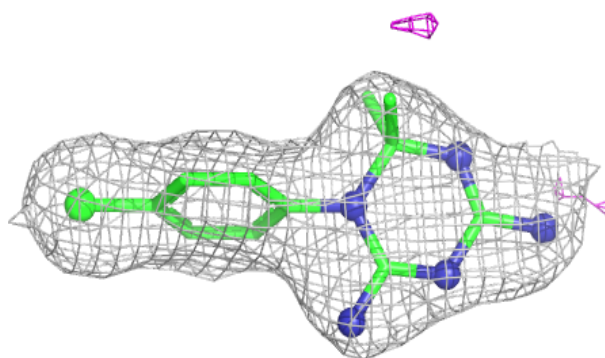


Electron density around NDP A 610:

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

**Electron density around 1CY A 609:**

$2mF_o-DF_c$ (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.