Full wwPDB X-ray Structure Validation Report (i)

Jun 11, 2024 – 07:49 PM EDT

PDB ID	:	1VJ3
Title	:	STRUCTURAL STUDIES ON BIO-ACTIVE COMPOUNDS. CRYSTAL STRUCTURE AND MOLECULAR MODELING STUDIES ON THE PNEUMOCYSTIS CARINII DIHYDROFOLATE REDUCTASE COFACTOR COMPLEX WITH TAB, A HIGHLY SELECTIVE ANTIFOLATE.
Authors	:	Cody, V.; Galitsky, N.; Rak, D.; Luft, J.R.; Queener, S.F.; Laughton, C.A.; Malcolm, F.G.
Deposited on	:	2004-01-13
Resolution	:	2.10 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:

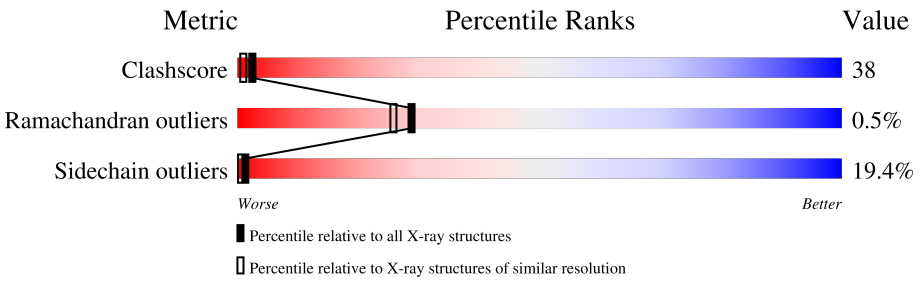
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.10 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	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	205	39% 40% 15% 6%

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
3	TAB	A	300[A]	-	-	X	-

2 Entry composition (i)

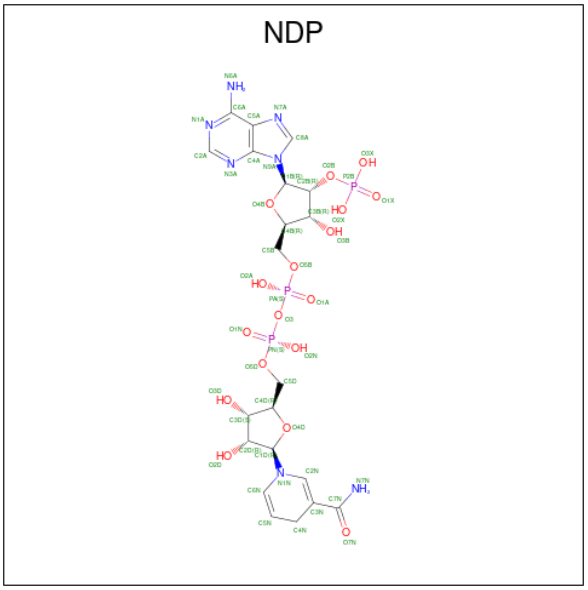
There are 4 unique types of molecules in this entry. The entry contains 1847 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DIHYDROFOLATE REDUCTASE.

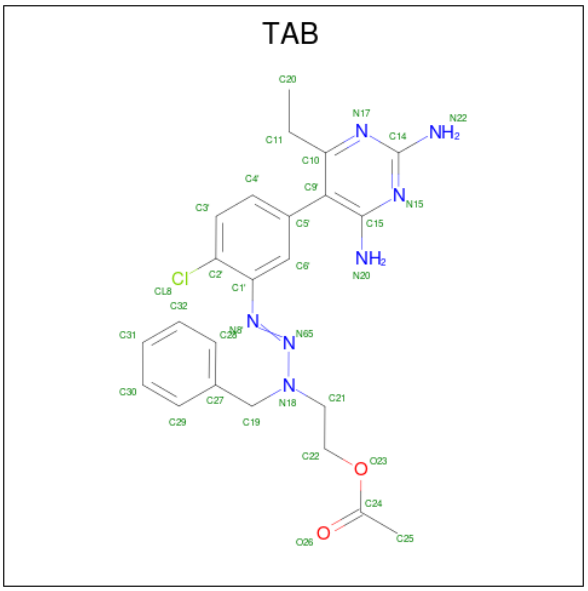
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	205	1678	1081	287	304	6	0	0	0

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



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

- Molecule 3 is ACETIC ACID N-[2-CHLORO-5-[6-ETHYL-2,4-DIAMINO-PYRIMID-5-YL]-PHENYL]-[BENZYL-TRIAZEN-3-YL]ETHYL ESTER (three-letter code: TAB) (formula: $C_{23}H_{26}ClN_7O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
3	A	1	46	34	1	7	4	0	1

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	75	75	75	0	0

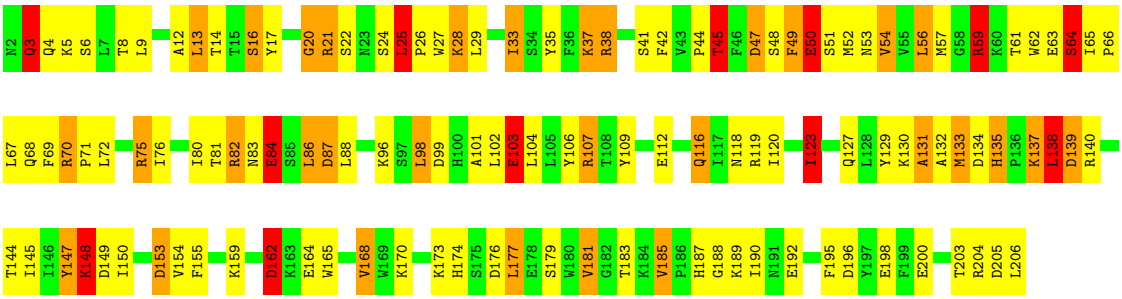
3 Residue-property plots (i)

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.

Note EDS was not executed.

• Molecule 1: DIHYDROFOLATE REDUCTASE

Chain A: 39% 40% 15% 6%



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	37.48Å 43.14Å 61.29Å 90.00° 94.74° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	97.8 (8.00-2.10)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.190 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1847	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

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: TAB, NDP

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	1.03	0/1720	2.68	106/2320 (4.6%)

There are no bond length outliers.

All (106) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	204	ARG	CD-NE-CZ	40.01	179.61	123.60
1	A	38	ARG	CD-NE-CZ	21.66	153.92	123.60
1	A	38	ARG	NE-CZ-NH2	21.20	130.90	120.30
1	A	99	ASP	CB-CG-OD2	-18.14	101.98	118.30
1	A	21	ARG	NE-CZ-NH2	-15.59	112.50	120.30
1	A	119	ARG	NE-CZ-NH2	-15.51	112.55	120.30
1	A	87	ASP	CB-CG-OD1	14.09	130.98	118.30
1	A	204	ARG	NE-CZ-NH2	14.00	127.30	120.30
1	A	140	ARG	NE-CZ-NH1	13.44	127.02	120.30
1	A	13	LEU	CB-CG-CD1	13.32	133.64	111.00
1	A	75	ARG	NE-CZ-NH2	-12.97	113.81	120.30
1	A	82	ARG	NE-CZ-NH1	12.57	126.59	120.30
1	A	107	ARG	NE-CZ-NH1	-12.56	114.02	120.30
1	A	75	ARG	NE-CZ-NH1	12.30	126.45	120.30
1	A	205	ASP	CB-CG-OD1	11.90	129.01	118.30
1	A	21	ARG	NE-CZ-NH1	11.67	126.14	120.30
1	A	63	GLU	OE1-CD-OE2	-11.44	109.58	123.30
1	A	38	ARG	CA-CB-CG	11.28	138.22	113.40
1	A	38	ARG	NH1-CZ-NH2	-10.68	107.65	119.40
1	A	99	ASP	OD1-CG-OD2	10.38	143.02	123.30
1	A	176	ASP	CA-CB-CG	9.72	134.79	113.40
1	A	70	ARG	NE-CZ-NH1	9.71	125.15	120.30
1	A	25	LEU	CB-CA-C	9.63	128.51	110.20
1	A	134	ASP	CB-CG-OD1	9.62	126.96	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	139	ASP	CB-CG-OD2	-9.62	109.64	118.30
1	A	63	GLU	CA-CB-CG	9.04	133.30	113.40
1	A	82	ARG	CD-NE-CZ	9.03	136.24	123.60
1	A	204	ARG	NH1-CZ-NH2	-8.95	109.55	119.40
1	A	35	TYR	CB-CG-CD1	8.87	126.32	121.00
1	A	3	GLN	CA-CB-CG	8.54	132.19	113.40
1	A	138	LEU	CB-CA-C	8.46	126.27	110.20
1	A	168	VAL	CA-CB-CG2	7.97	122.86	110.90
1	A	153	ASP	CB-CG-OD2	7.92	125.43	118.30
1	A	63	GLU	CB-CG-CD	7.89	135.49	114.20
1	A	139	ASP	CB-CG-OD1	7.70	125.23	118.30
1	A	205	ASP	CB-CG-OD2	-7.50	111.55	118.30
1	A	35	TYR	CB-CG-CD2	-7.34	116.60	121.00
1	A	200	GLU	CG-CD-OE1	7.25	132.79	118.30
1	A	132	ALA	N-CA-CB	-7.09	100.17	110.10
1	A	82	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	37	LYS	CA-CB-CG	7.01	128.81	113.40
1	A	107	ARG	CD-NE-CZ	-6.97	113.84	123.60
1	A	176	ASP	CB-CA-C	6.86	124.12	110.40
1	A	101	ALA	CB-CA-C	6.83	120.35	110.10
1	A	195	PHE	CA-CB-CG	6.71	130.01	113.90
1	A	123	ILE	CB-CG1-CD1	6.58	132.31	113.90
1	A	16	SER	N-CA-CB	-6.53	100.71	110.50
1	A	107	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	A	134	ASP	CB-CG-OD2	-6.37	112.57	118.30
1	A	177	LEU	CA-CB-CG	6.33	129.85	115.30
1	A	103	GLU	CB-CG-CD	6.28	131.16	114.20
1	A	192	GLU	CG-CD-OE2	-6.27	105.76	118.30
1	A	59	ARG	CA-CB-CG	6.23	127.11	113.40
1	A	47	ASP	CB-CG-OD2	-6.14	112.78	118.30
1	A	144	THR	N-CA-CB	6.10	121.89	110.30
1	A	54	VAL	CG1-CB-CG2	6.09	120.64	110.90
1	A	27	TRP	C-N-CA	5.88	136.40	121.70
1	A	103	GLU	OE1-CD-OE2	-5.87	116.25	123.30
1	A	168	VAL	CA-CB-CG1	5.86	119.69	110.90
1	A	50	GLU	C-N-CA	5.86	136.35	121.70
1	A	133	MET	CG-SD-CE	5.83	109.53	100.20
1	A	21	ARG	CD-NE-CZ	-5.79	115.50	123.60
1	A	147	TYR	CB-CG-CD2	-5.72	117.57	121.00
1	A	181	VAL	CA-CB-CG2	5.65	119.37	110.90
1	A	147	TYR	CB-CG-CD1	5.63	124.38	121.00
1	A	162	ASP	CB-CG-OD2	-5.62	113.24	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	101	ALA	N-CA-CB	-5.61	102.25	110.10
1	A	192	GLU	OE1-CD-OE2	5.60	130.03	123.30
1	A	96	LYS	CA-CB-CG	5.56	125.63	113.40
1	A	54	VAL	CB-CA-C	-5.54	100.86	111.40
1	A	14	THR	CA-CB-CG2	5.54	120.15	112.40
1	A	140	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	A	56	LEU	CA-CB-CG	5.52	127.99	115.30
1	A	13	LEU	CA-CB-CG	5.49	127.92	115.30
1	A	59	ARG	CD-NE-CZ	5.42	131.19	123.60
1	A	25	LEU	N-CA-C	-5.42	96.37	111.00
1	A	135	HIS	CB-CA-C	5.41	121.23	110.40
1	A	154	VAL	CA-CB-CG2	5.40	119.00	110.90
1	A	84	GLU	CB-CG-CD	5.39	128.75	114.20
1	A	130	LYS	O-C-N	-5.39	114.08	122.70
1	A	45	THR	N-CA-CB	-5.37	100.10	110.30
1	A	64	SER	O-C-N	5.37	131.28	122.70
1	A	127	GLN	CA-C-O	-5.34	108.88	120.10
1	A	28	LYS	CB-CA-C	-5.31	99.78	110.40
1	A	107	ARG	CA-CB-CG	5.30	125.07	113.40
1	A	53	ASN	N-CA-CB	5.29	120.13	110.60
1	A	109	TYR	CB-CG-CD2	-5.29	117.83	121.00
1	A	168	VAL	CG1-CB-CG2	-5.28	102.46	110.90
1	A	3	GLN	CA-C-N	-5.27	105.61	117.20
1	A	132	ALA	CA-C-O	5.27	131.16	120.10
1	A	59	ARG	CA-C-O	-5.26	109.06	120.10
1	A	131	ALA	N-CA-CB	-5.22	102.80	110.10
1	A	119	ARG	NH1-CZ-NH2	5.21	125.13	119.40
1	A	8	THR	N-CA-CB	5.16	120.09	110.30
1	A	168	VAL	CB-CA-C	5.15	121.19	111.40
1	A	83	ASN	CB-CA-C	5.14	120.69	110.40
1	A	155	PHE	N-CA-CB	-5.14	101.35	110.60
1	A	20	GLY	O-C-N	5.11	130.88	122.70
1	A	203	THR	CA-CB-CG2	5.11	119.55	112.40
1	A	204	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	12	ALA	N-CA-CB	5.10	117.24	110.10
1	A	64	SER	N-CA-CB	5.09	118.13	110.50
1	A	148	LYS	CA-CB-CG	5.08	124.58	113.40
1	A	75	ARG	CD-NE-CZ	5.07	130.69	123.60
1	A	148	LYS	CB-CA-C	5.04	120.48	110.40
1	A	99	ASP	O-C-N	5.03	130.75	122.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1678	0	1681	119	0
2	A	48	0	26	13	0
3	A	46	0	28	16	0
4	A	75	0	0	6	0
All	All	1847	0	1735	131	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 38.

All (131) 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:A:42:PHE:HB3	4:A:270:HOH:O	1.45	1.15
1:A:107:ARG:HH11	1:A:107:ARG:HG2	1.09	1.14
1:A:50:GLU:CD	1:A:116:GLN:HG3	1.82	1.00
1:A:170:LYS:HG2	4:A:222:HOH:O	1.67	0.94
1:A:107:ARG:HG2	1:A:107:ARG:NH1	1.82	0.94
1:A:98:LEU:HD12	1:A:98:LEU:H	1.30	0.92
1:A:50:GLU:HG3	1:A:116:GLN:HG2	1.54	0.89
1:A:20:GLY:HA2	1:A:26:PRO:HD3	1.56	0.87
1:A:50:GLU:HG3	1:A:116:GLN:CG	2.05	0.86
1:A:86:LEU:H	1:A:86:LEU:HD12	1.41	0.85
1:A:112:GLU:HG2	1:A:112:GLU:O	1.77	0.84
1:A:50:GLU:HG2	1:A:51:SER:N	1.96	0.81
1:A:3:GLN:HA	1:A:106:TYR:HE2	1.48	0.78
1:A:187:HIS:ND1	1:A:188:GLY:N	2.33	0.77
1:A:50:GLU:OE2	1:A:116:GLN:HG3	1.86	0.76
1:A:107:ARG:HH11	1:A:107:ARG:CG	1.79	0.74
1:A:54:VAL:HG12	1:A:56:LEU:HD12	1.68	0.73
1:A:33:ILE:HG13	3:A:300[A]:TAB:C32	2.19	0.73
1:A:50:GLU:CG	1:A:116:GLN:HG3	2.19	0.73
1:A:33:ILE:HG13	3:A:300[A]:TAB:H321	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:300[A]:TAB:C29	3:A:300[A]:TAB:H211	2.20	0.70
1:A:54:VAL:CG1	1:A:56:LEU:HD12	2.21	0.70
1:A:80:ILE:O	2:A:207:NDP:H1B	1.92	0.70
1:A:38:ARG:NE	1:A:183:THR:HG21	2.07	0.69
1:A:135:HIS:HD2	1:A:137:LYS:H	1.38	0.69
1:A:44:PRO:O	1:A:48:SER:N	2.26	0.69
1:A:66:PRO:HG2	3:A:300[A]:TAB:H291	1.73	0.69
1:A:104:LEU:HD23	1:A:104:LEU:C	2.14	0.68
1:A:66:PRO:HG2	3:A:300[A]:TAB:C29	2.23	0.68
1:A:145:ILE:H	1:A:145:ILE:HD12	1.58	0.68
1:A:86:LEU:HD12	1:A:86:LEU:N	2.10	0.67
1:A:70:ARG:HA	1:A:71:PRO:C	2.14	0.67
1:A:61:THR:O	1:A:65:ILE:HG13	1.95	0.67
1:A:98:LEU:HD12	4:A:236:HOH:O	1.95	0.67
1:A:37:LYS:HG3	3:A:300[B]:TAB:H311	1.77	0.66
1:A:21:ARG:HA	1:A:153:ASP:OD2	1.97	0.65
1:A:38:ARG:HD2	1:A:183:THR:HG21	1.77	0.65
1:A:107:ARG:NH1	1:A:107:ARG:CG	2.48	0.65
1:A:5:LYS:NZ	1:A:118:ASN:O	2.23	0.65
1:A:98:LEU:CD1	4:A:236:HOH:O	2.44	0.65
1:A:50:GLU:CG	1:A:116:GLN:CG	2.73	0.64
1:A:148:LYS:HG3	1:A:150:ILE:HG13	1.79	0.63
1:A:145:ILE:HD12	1:A:145:ILE:N	2.13	0.63
1:A:68:GLN:HG2	1:A:69:PHE:CE1	2.33	0.63
1:A:82:ARG:HG3	2:A:207:NDP:H2A	1.81	0.63
1:A:25:LEU:HD12	1:A:25:LEU:O	2.00	0.61
1:A:56:LEU:CD1	1:A:120:ILE:CG2	2.78	0.61
1:A:147:TYR:CE2	1:A:189:LYS:HE3	2.36	0.61
1:A:148:LYS:HG3	1:A:150:ILE:CG1	2.31	0.60
1:A:38:ARG:CZ	1:A:183:THR:CG2	2.79	0.60
1:A:38:ARG:CD	1:A:183:THR:HG21	2.31	0.59
1:A:162:ASP:OD2	4:A:212:HOH:O	2.17	0.59
1:A:9:LEU:HB2	1:A:138:LEU:HD21	1.85	0.59
1:A:69:PHE:CD1	3:A:300[A]:TAB:H301	2.38	0.59
1:A:66:PRO:CG	3:A:300[A]:TAB:H291	2.34	0.57
1:A:174:HIS:CE1	1:A:185:VAL:HG22	2.39	0.57
1:A:51:SER:HB3	1:A:118:ASN:HB2	1.87	0.56
1:A:135:HIS:CD2	1:A:137:LYS:H	2.21	0.56
1:A:129:TYR:OH	2:A:207:NDP:H41N	2.06	0.55
1:A:66:PRO:HD3	3:A:300[A]:TAB:H192	1.88	0.55
1:A:84:GLU:OE2	1:A:84:GLU:O	2.25	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:ARG:CZ	1:A:183:THR:HG22	2.36	0.54
1:A:56:LEU:HD13	1:A:120:ILE:HG22	1.89	0.54
1:A:38:ARG:CZ	1:A:183:THR:HG21	2.38	0.54
1:A:66:PRO:CG	3:A:300[A]:TAB:C29	2.86	0.54
1:A:38:ARG:NE	1:A:183:THR:CG2	2.71	0.53
1:A:98:LEU:HD11	1:A:131:ALA:HB1	1.90	0.53
1:A:50:GLU:HG3	1:A:116:GLN:HG3	1.80	0.53
1:A:47:ASP:O	1:A:51:SER:HB2	2.09	0.53
1:A:3:GLN:HG2	1:A:137:LYS:HG3	1.91	0.51
1:A:69:PHE:CE1	3:A:300[A]:TAB:H301	2.47	0.50
1:A:164:GLU:HG3	1:A:165:TRP:CD1	2.46	0.50
1:A:68:GLN:HG2	1:A:69:PHE:CZ	2.47	0.50
1:A:82:ARG:HG3	2:A:207:NDP:C2A	2.41	0.50
1:A:174:HIS:CE1	1:A:185:VAL:CG2	2.94	0.50
3:A:300[A]:TAB:H211	3:A:300[A]:TAB:H291	1.93	0.49
1:A:57:MET:HB3	1:A:123:ILE:HD11	1.96	0.47
1:A:65:ILE:CG2	1:A:66:PRO:HD2	2.43	0.47
1:A:66:PRO:CD	3:A:300[A]:TAB:H291	2.44	0.47
1:A:98:LEU:HD12	1:A:98:LEU:N	2.10	0.47
1:A:103:GLU:O	1:A:107:ARG:HB2	2.15	0.47
1:A:159:LYS:CE	1:A:162:ASP:OD2	2.64	0.46
1:A:56:LEU:HD11	1:A:120:ILE:CG2	2.44	0.46
1:A:45:THR:O	1:A:49:PHE:HD1	1.98	0.46
1:A:72:LEU:HB3	1:A:75:ARG:NH2	2.31	0.46
1:A:3:GLN:HG2	1:A:137:LYS:CG	2.46	0.45
1:A:86:LEU:N	1:A:86:LEU:CD1	2.71	0.45
1:A:159:LYS:HB3	1:A:159:LYS:HE3	1.70	0.45
1:A:70:ARG:HB2	1:A:71:PRO:HA	1.99	0.45
1:A:81:THR:O	1:A:81:THR:HG23	2.17	0.45
1:A:52:MET:HE2	1:A:76:ILE:HD11	1.99	0.44
1:A:145:ILE:HD12	1:A:198:GLU:O	2.18	0.44
1:A:37:LYS:HG3	3:A:300[A]:TAB:H251	2.00	0.43
1:A:16:SER:O	1:A:17:TYR:HB2	2.17	0.43
1:A:56:LEU:CD1	1:A:120:ILE:HG23	2.49	0.43
1:A:159:LYS:HE3	1:A:162:ASP:OD2	2.18	0.43
1:A:66:PRO:CD	3:A:300[A]:TAB:H192	2.49	0.43
1:A:137:LYS:HB3	1:A:137:LYS:HE2	1.76	0.43
1:A:33:ILE:HD13	1:A:33:ILE:HA	1.89	0.43
1:A:25:LEU:O	1:A:25:LEU:CD1	2.67	0.43
1:A:59:ARG:O	1:A:62:TRP:HB3	2.19	0.43
1:A:3:GLN:HA	1:A:106:TYR:CE2	2.40	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:VAL:HG11	1:A:56:LEU:HD12	1.98	0.42
1:A:86:LEU:HD12	1:A:86:LEU:O	2.19	0.42
1:A:87:ASP:HA	4:A:265:HOH:O	2.19	0.42
1:A:54:VAL:HG11	1:A:56:LEU:CD1	2.50	0.42
1:A:3:GLN:HB3	1:A:102:LEU:HD13	2.02	0.42
1:A:47:ASP:C	1:A:49:PHE:H	2.19	0.42
1:A:65:ILE:HG23	1:A:66:PRO:HD2	2.01	0.42
1:A:56:LEU:HD13	1:A:120:ILE:CG2	2.46	0.42
1:A:38:ARG:O	1:A:42:PHE:HB2	2.20	0.41
1:A:139:ASP:HB3	1:A:206:LEU:HD11	2.03	0.41
1:A:129:TYR:O	1:A:133:MET:HG2	2.20	0.41
1:A:188:GLY:O	1:A:190:ILE:HD12	2.21	0.41
3:A:300[A]:TAB:C29	3:A:300[A]:TAB:C21	2.93	0.41
1:A:59:ARG:O	1:A:62:TRP:N	2.53	0.40
1:A:187:HIS:CG	1:A:188:GLY:N	2.90	0.40
1:A:21:ARG:HH21	1:A:21:ARG:HD3	1.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	203/205 (99%)	190 (94%)	12 (6%)	1 (0%)	29 26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	GLN

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	186/186 (100%)	150 (81%)	36 (19%)	1 0

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	4	GLN
1	A	6	SER
1	A	13	LEU
1	A	22	SER
1	A	24	SER
1	A	25	LEU
1	A	28	LYS
1	A	29	LEU
1	A	33	ILE
1	A	41	SER
1	A	45	THR
1	A	49	PHE
1	A	50	GLU
1	A	59	ARG
1	A	64	SER
1	A	67	LEU
1	A	84	GLU
1	A	86	LEU
1	A	88	LEU
1	A	98	LEU
1	A	103	GLU
1	A	116	GLN
1	A	123	ILE
1	A	137	LYS
1	A	138	LEU
1	A	148	LYS
1	A	149	ASP
1	A	162	ASP
1	A	168	VAL

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Mol	Chain	Res	Type
1	A	173	LYS
1	A	177	LEU
1	A	179	SER
1	A	181	VAL
1	A	185	VAL
1	A	196	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	116	GLN
1	A	135	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)

3 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	$\#\|Z\| > 2$
3	TAB	A	300[B]	-	34,35,35	1.78	8 (23%)	39,47,47	3.20	16 (41%)
2	NDP	A	207	-	47,52,52	2.51	16 (34%)	61,80,80	2.34	19 (31%)
3	TAB	A	300[A]	-	34,35,35	1.84	8 (23%)	39,47,47	3.32	18 (46%)

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
3	TAB	A	300[B]	-	-	9/17/21/21	0/3/3/3
2	NDP	A	207	-	-	3/30/77/77	0/5/5/5
3	TAB	A	300[A]	-	-	2/17/21/21	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	207	NDP	O4B-C4B	-9.70	1.23	1.45
2	A	207	NDP	P2B-O2B	7.26	1.72	1.59
3	A	300[A]	TAB	C21-N18	-5.07	1.41	1.46
3	A	300[A]	TAB	C1'-N8'	-4.66	1.32	1.42
3	A	300[B]	TAB	C1'-N8'	-4.66	1.32	1.42
3	A	300[A]	TAB	N8'-N65	4.41	1.35	1.27
3	A	300[B]	TAB	N8'-N65	4.41	1.35	1.27
2	A	207	NDP	C3B-C4B	4.40	1.64	1.53
3	A	300[B]	TAB	C21-N18	-4.01	1.42	1.46
2	A	207	NDP	C6A-C5A	3.58	1.56	1.43
2	A	207	NDP	O4B-C1B	3.31	1.45	1.40
3	A	300[A]	TAB	N18-N65	2.82	1.44	1.34
3	A	300[B]	TAB	N18-N65	2.82	1.44	1.34
2	A	207	NDP	PN-O1N	-2.77	1.41	1.50
2	A	207	NDP	C1D-N1N	2.61	1.53	1.46
2	A	207	NDP	C4N-C3N	2.58	1.54	1.50
2	A	207	NDP	O3B-C3B	2.51	1.49	1.43
2	A	207	NDP	O4D-C1D	2.46	1.47	1.42
2	A	207	NDP	C6A-N1A	-2.29	1.27	1.36
2	A	207	NDP	O4D-C4D	2.27	1.50	1.45
2	A	207	NDP	O2D-C2D	-2.26	1.37	1.43
3	A	300[A]	TAB	C6'-C1'	2.22	1.41	1.37
3	A	300[B]	TAB	C6'-C1'	2.22	1.41	1.37
3	A	300[A]	TAB	C10-N17	2.21	1.35	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	300[B]	TAB	C10-N17	2.21	1.35	1.32
2	A	207	NDP	PA-O2A	-2.19	1.45	1.55
2	A	207	NDP	PN-O3	2.15	1.61	1.59
3	A	300[A]	TAB	C4'-C5'	-2.10	1.37	1.42
3	A	300[B]	TAB	C4'-C5'	-2.10	1.37	1.42
2	A	207	NDP	C1B-N9A	-2.07	1.44	1.49
3	A	300[A]	TAB	C4'-C3'	2.00	1.41	1.36
3	A	300[B]	TAB	C4'-C3'	2.00	1.41	1.36

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	300[A]	TAB	C1'-N8'-N65	9.33	123.89	112.17
3	A	300[B]	TAB	C1'-N8'-N65	9.33	123.89	112.17
3	A	300[A]	TAB	C3'-C2'-C1'	-8.51	114.16	121.48
3	A	300[B]	TAB	C3'-C2'-C1'	-8.51	114.16	121.48
2	A	207	NDP	O3B-C3B-C4B	-7.32	90.05	111.08
3	A	300[A]	TAB	C10-C9'-C15	-6.58	117.45	123.34
3	A	300[B]	TAB	C10-C9'-C15	-6.58	117.45	123.34
3	A	300[A]	TAB	C6'-C1'-C2'	5.74	124.05	117.99
3	A	300[B]	TAB	C6'-C1'-C2'	5.74	124.05	117.99
2	A	207	NDP	C3N-C2N-N1N	-5.34	115.36	123.20
3	A	300[A]	TAB	N22-C14-N15	4.91	124.58	117.22
3	A	300[B]	TAB	N22-C14-N15	4.91	124.58	117.22
2	A	207	NDP	C4B-O4B-C1B	4.89	114.40	109.92
2	A	207	NDP	P2B-O2B-C2B	4.72	136.05	123.43
2	A	207	NDP	C6N-N1N-C2N	4.34	123.96	119.32
2	A	207	NDP	O3X-P2B-O1X	4.28	127.52	110.83
3	A	300[A]	TAB	O26-C24-C25	-4.25	109.73	124.77
3	A	300[B]	TAB	O26-C24-C25	-4.23	109.78	124.77
3	A	300[A]	TAB	C19-C27-C28	4.15	128.39	120.75
3	A	300[A]	TAB	N15-C14-N17	-3.93	119.46	125.48
3	A	300[B]	TAB	N15-C14-N17	-3.93	119.46	125.48
2	A	207	NDP	C1D-N1N-C6N	-3.80	112.74	120.77
3	A	300[A]	TAB	C19-C27-C29	-3.68	113.97	120.75
2	A	207	NDP	C3N-C7N-N7N	-3.67	111.16	117.67
2	A	207	NDP	C5B-C4B-C3B	-3.60	102.24	115.21
3	A	300[A]	TAB	C1'-C2'-CL8	3.48	123.61	119.70
3	A	300[B]	TAB	C1'-C2'-CL8	3.48	123.61	119.70
2	A	207	NDP	N6A-C6A-N1A	3.29	125.37	118.33
3	A	300[A]	TAB	C14-N15-C15	3.25	126.01	116.72
3	A	300[B]	TAB	C14-N15-C15	3.25	126.01	116.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	207	NDP	C4A-C5A-N7A	3.16	112.68	109.34
2	A	207	NDP	C2B-C3B-C4B	-3.15	95.21	101.99
2	A	207	NDP	C3B-C2B-C1B	-3.06	96.94	102.81
3	A	300[A]	TAB	C4'-C3'-C2'	3.02	122.35	119.68
3	A	300[B]	TAB	C4'-C3'-C2'	3.02	122.35	119.68
2	A	207	NDP	O7N-C7N-N7N	2.84	129.26	122.89
2	A	207	NDP	O2A-PA-O3	2.77	114.77	107.27
3	A	300[A]	TAB	N20-C15-N15	2.68	124.31	117.11
3	A	300[B]	TAB	N20-C15-N15	2.68	124.31	117.11
3	A	300[A]	TAB	C6'-C1'-N8'	-2.38	116.31	122.42
3	A	300[B]	TAB	C6'-C1'-N8'	-2.38	116.31	122.42
2	A	207	NDP	N3A-C2A-N1A	2.36	131.88	128.67
3	A	300[A]	TAB	C9'-C15-N15	-2.32	116.53	122.61
3	A	300[B]	TAB	C9'-C15-N15	-2.32	116.53	122.61
2	A	207	NDP	O5D-PN-O1N	2.24	117.80	108.94
2	A	207	NDP	PN-O5D-C5D	2.13	133.56	121.35
3	A	300[A]	TAB	C20-C11-C10	-2.12	108.91	114.59
3	A	300[B]	TAB	C20-C11-C10	-2.12	108.91	114.59
3	A	300[A]	TAB	C9'-C15-N20	-2.09	118.99	122.70
3	A	300[B]	TAB	C9'-C15-N20	-2.09	118.99	122.70
3	A	300[A]	TAB	C30-C29-C27	2.08	123.54	120.61
3	A	300[B]	TAB	C30-C29-C27	2.05	123.50	120.61
2	A	207	NDP	C5A-C6A-N1A	-2.04	115.37	120.23

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	300[A]	TAB	N18-N65-N8'-C1'
3	A	300[B]	TAB	N18-N65-N8'-C1'
3	A	300[B]	TAB	C22-C21-N18-C19
3	A	300[B]	TAB	C19-N18-N65-N8'
3	A	300[B]	TAB	C21-N18-N65-N8'
3	A	300[B]	TAB	N18-C21-C22-O23
3	A	300[B]	TAB	C25-C24-O23-C22
3	A	300[B]	TAB	O26-C24-O23-C22
2	A	207	NDP	C3B-C4B-C5B-O5B
2	A	207	NDP	O4D-C1D-N1N-C2N
3	A	300[A]	TAB	C9'-C10-C11-C20
3	A	300[B]	TAB	C9'-C10-C11-C20
3	A	300[B]	TAB	C27-C19-N18-C21
2	A	207	NDP	C2N-C3N-C7N-N7N

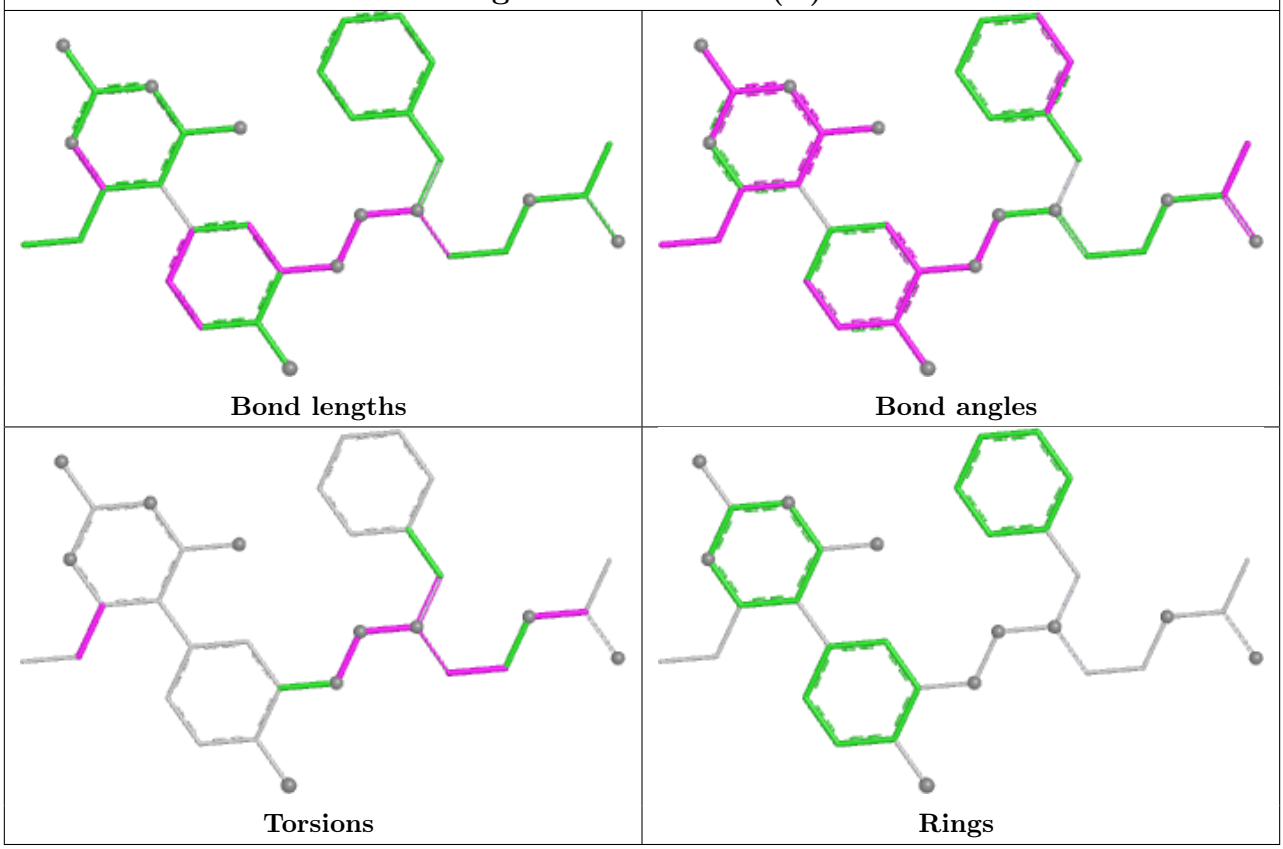
There are no ring outliers.

3 monomers are involved in 29 short contacts:

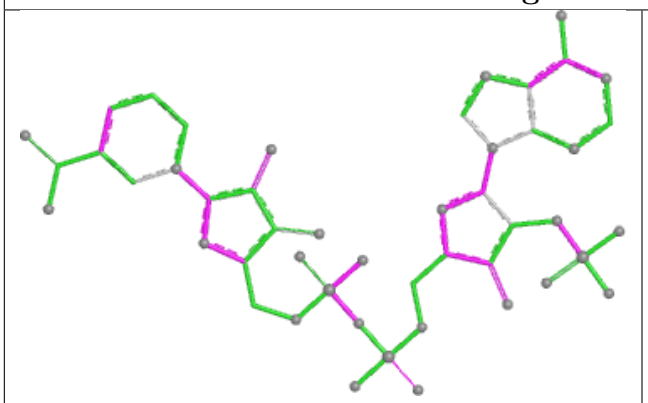
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	300[B]	TAB	1	0
2	A	207	NDP	13	0
3	A	300[A]	TAB	15	0

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.

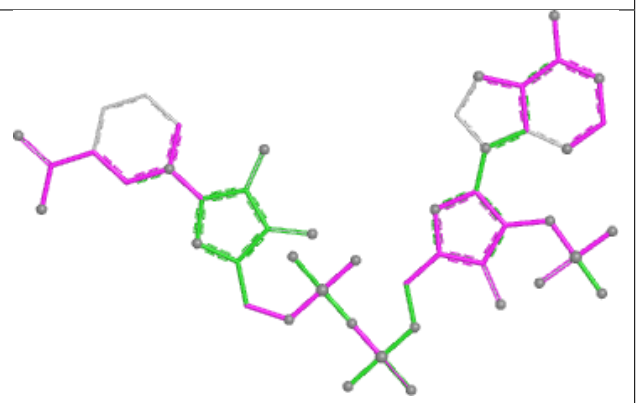
Ligand TAB A 300 (B)



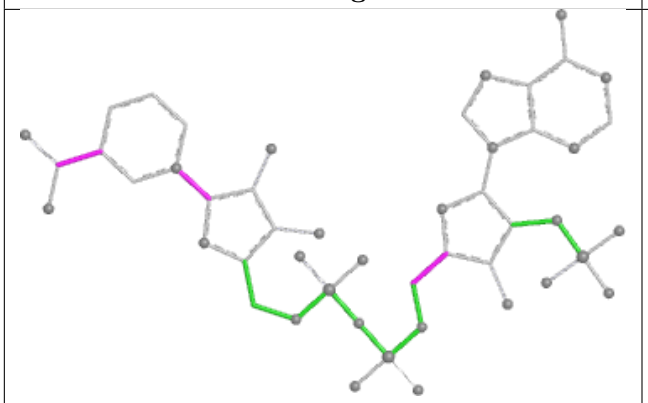
Ligand NDP A 207



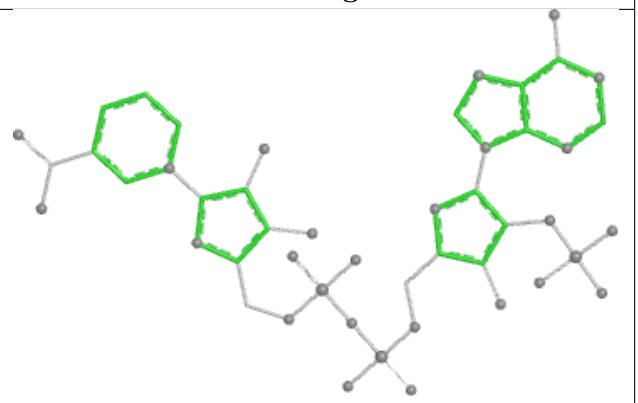
Bond lengths



Bond angles

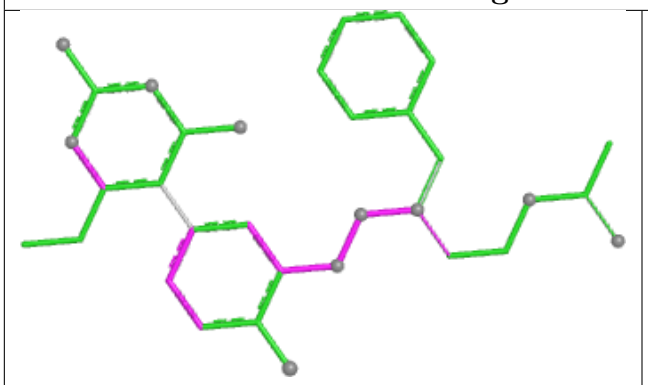


Torsions

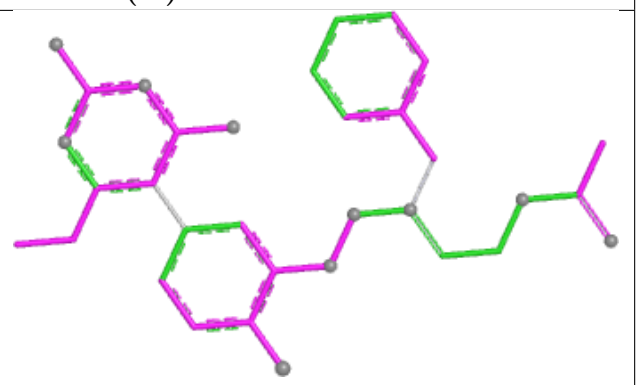


Rings

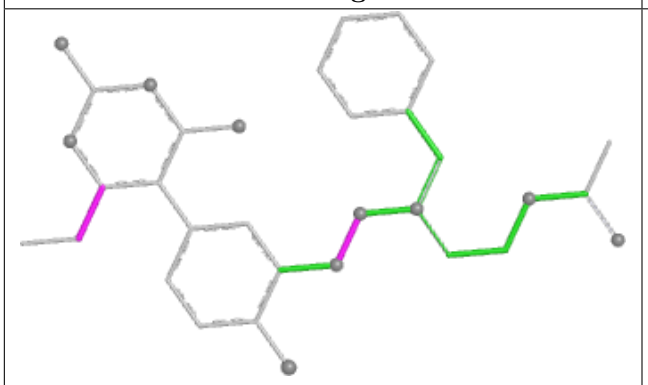
Ligand TAB A 300 (A)



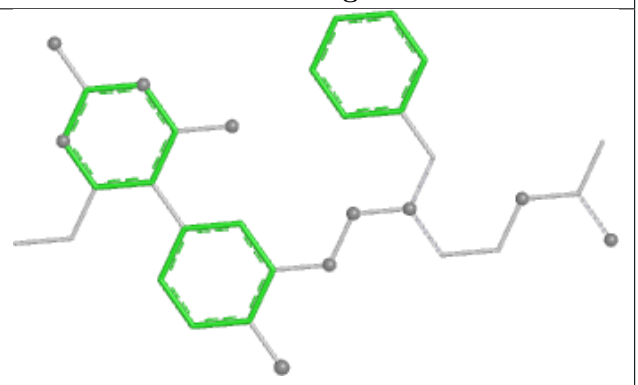
Bond lengths



Bond angles



Torsions



Rings

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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.