



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:50 AM UTC

PDB ID : 1G0N / pdb_00001g0n
Title : STRUCTURE OF TRIHYDROXYNAPHTHALENE REDUCTASE IN
COMPLEX WITH NADPH AND 4,5,6,7-TETRACHLORO-PHTHALID
E
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Deposited on : 2000-10-06
Resolution : 2.00 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

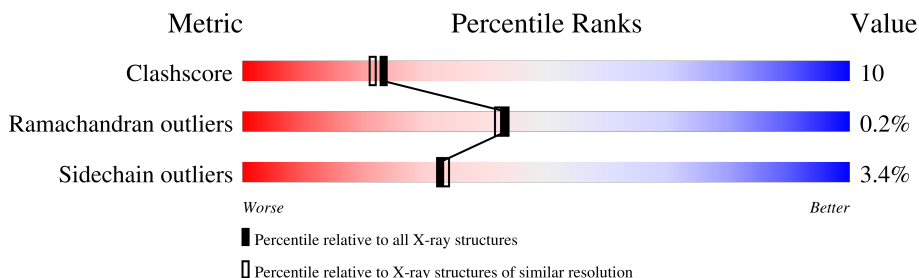
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	283	
1	B	283	

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	PHH	A	402	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4417 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.

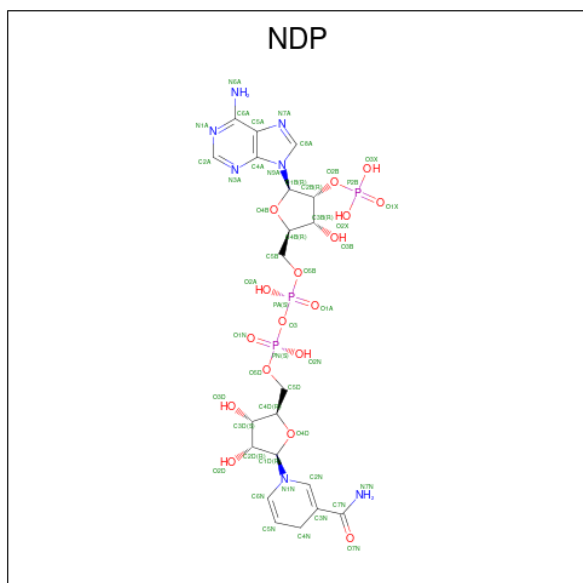
- Molecule 1 is a protein called TRIHYDROXYNAPHTHALENE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2036	1279	357	385	15			
1	B	254	Total	C	N	O	S	0	0	0
			1885	1188	331	352	14			

There are 6 discrepancies between the modelled and reference sequences:

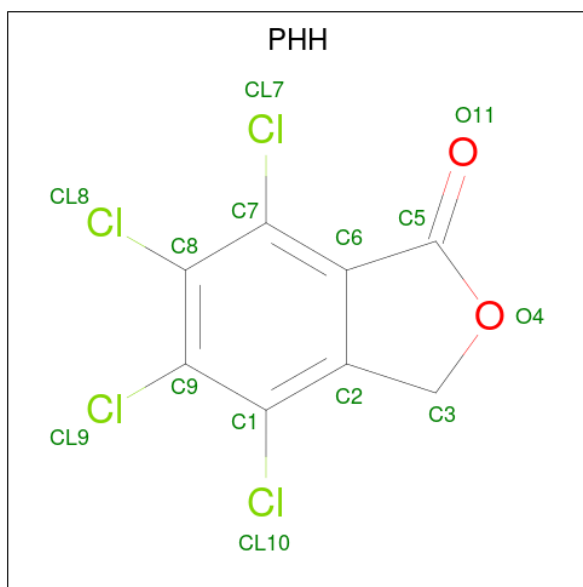
Chain	Residue	Modelled	Actual	Comment	Reference
A	241	VAL	SER	engineered mutation	UNP Q12634
A	242	GLN	ALA	engineered mutation	UNP Q12634
A	247	ARG	HIS	engineered mutation	UNP Q12634
B	241	VAL	SER	engineered mutation	UNP Q12634
B	242	GLN	ALA	engineered mutation	UNP Q12634
B	247	ARG	HIS	engineered mutation	UNP Q12634

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



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

- Molecule 3 is 4,5,6,7-TETRACHLORO-PHTHALIDE (CCD ID: PHH) (formula: $C_8H_2Cl_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	O	0	0
			14	8	4	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	202	Total	O	0	0
			202	202		
4	B	184	Total	O	0	0
			184	184		

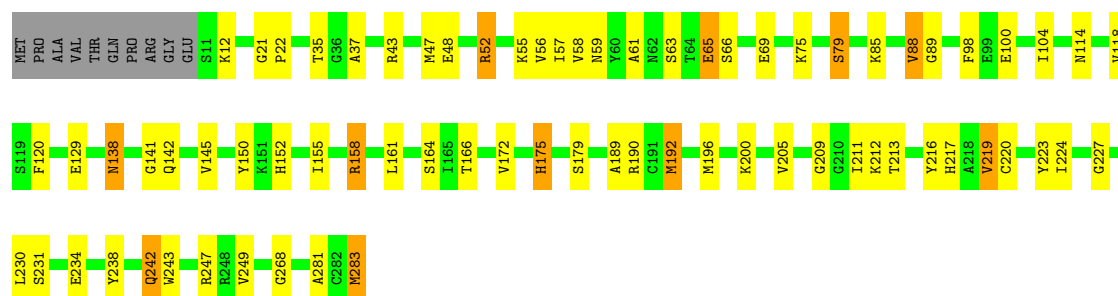
3 Residue-property plots

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. 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.

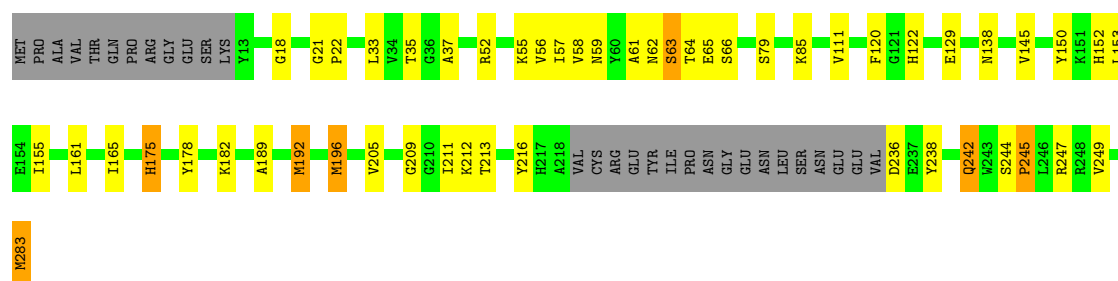
• Molecule 1: TRIHYDROXYNAPHTHALENE REDUCTASE

Chain A: 



• Molecule 1: TRIHYDROXYNAPHTHALENE REDUCTASE

Chain B: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.50 Å 140.50 Å 72.70 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.00)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.195 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4417	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PHH, 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	0.67	2/2068 (0.1%)	1.08	14/2790 (0.5%)
1	B	0.69	3/1914 (0.2%)	1.06	14/2580 (0.5%)
All	All	0.68	5/3982 (0.1%)	1.07	28/5370 (0.5%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	196	MET	SD-CE	-11.84	1.50	1.79
1	A	192	MET	SD-CE	-10.63	1.52	1.79
1	B	283	MET	SD-CE	-6.88	1.62	1.79
1	A	283	MET	SD-CE	-6.82	1.62	1.79
1	B	192	MET	SD-CE	-6.15	1.64	1.79

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	21	GLY	CA-C-N	9.54	130.54	119.47
1	B	21	GLY	C-N-CA	9.54	130.54	119.47
1	B	209	GLY	N-CA-C	-8.15	98.58	112.77
1	A	21	GLY	CA-C-N	8.10	128.87	119.47
1	A	21	GLY	C-N-CA	8.10	128.87	119.47
1	A	209	GLY	N-CA-C	-7.41	99.88	112.77
1	B	55	LYS	N-CA-C	-7.29	97.72	109.96
1	A	145	VAL	N-CA-C	-7.12	103.53	110.72
1	A	56	VAL	N-CA-C	7.00	117.91	108.11
1	A	55	LYS	N-CA-C	-6.79	98.20	109.46
1	B	145	VAL	N-CA-C	-6.72	103.93	110.72
1	B	155	ILE	N-CA-C	-6.24	100.95	109.37
1	B	247	ARG	N-CA-C	6.13	120.10	111.52
1	A	12	LYS	N-CA-C	-6.10	105.32	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	242	GLN	N-CA-C	6.03	117.93	111.36
1	B	242	GLN	N-CA-C	5.95	119.00	111.69
1	A	219	VAL	N-CA-C	5.80	118.57	112.83
1	A	268	GLY	N-CA-C	5.70	119.78	112.83
1	A	155	ILE	N-CA-C	-5.45	102.01	109.37
1	A	247	ARG	N-CA-C	5.44	118.90	111.39
1	A	138	ASN	N-CA-C	5.42	118.11	111.82
1	B	56	VAL	N-CA-C	5.39	115.66	108.11
1	B	122	HIS	N-CA-C	-5.39	102.41	110.23
1	B	245	PRO	N-CA-C	-5.36	107.16	113.86
1	B	18	GLY	CA-C-N	5.32	125.08	119.76
1	B	18	GLY	C-N-CA	5.32	125.08	119.76
1	B	138	ASN	N-CA-C	5.27	117.93	111.82
1	A	231	SER	N-CA-C	-5.27	103.73	110.53

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	2036	0	2051	51	0
1	B	1885	0	1908	32	0
2	A	48	0	26	3	0
2	B	48	0	26	0	0
3	A	14	0	2	4	0
4	A	202	0	0	8	0
4	B	184	0	0	3	0
All	All	4417	0	4013	83	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 10.

All (83) 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:283:MET:HE1	4:A:440:HOH:O	1.70	0.91
1:A:120:PHE:HE1	1:A:219:VAL:HB	1.37	0.87
1:A:120:PHE:CE1	1:A:219:VAL:HB	2.14	0.82
1:B:165:ILE:HD13	1:B:283:MET:HE3	1.62	0.81
1:B:283:MET:HE1	4:B:525:HOH:O	1.79	0.81
1:A:216:TYR:HA	3:A:402:PHH:CL10	2.18	0.80
1:B:189:ALA:HB2	1:B:205:VAL:HG23	1.70	0.72
1:B:238:TYR:O	1:B:242:GLN:HB2	1.90	0.70
1:A:238:TYR:O	1:A:242:GLN:HB2	1.91	0.70
1:B:161:LEU:HD11	1:B:192:MET:HE1	1.75	0.69
1:A:48:GLU:O	1:A:52:ARG:HG2	1.95	0.67
1:B:213:THR:H	1:B:216:TYR:HB3	1.60	0.66
1:B:211:ILE:O	1:B:216:TYR:HB2	1.96	0.65
1:A:211:ILE:O	1:A:216:TYR:HB2	1.98	0.64
1:A:213:THR:O	1:A:217:HIS:HD2	1.82	0.62
1:B:61:ALA:O	1:B:85:LYS:HE3	2.00	0.61
1:A:118:VAL:HG21	1:A:219:VAL:HG21	1.84	0.59
1:A:213:THR:H	1:A:216:TYR:HB3	1.68	0.58
1:B:161:LEU:HD11	1:B:192:MET:CE	2.34	0.57
1:A:65:GLU:O	1:A:69:GLU:HG3	2.05	0.57
1:A:100:GLU:O	1:A:104:ILE:HG13	2.05	0.57
1:B:212:LYS:HG3	1:B:249:VAL:HG11	1.86	0.56
1:B:150:TYR:HB2	1:B:196:MET:HE3	1.88	0.56
1:B:238:TYR:CZ	1:B:242:GLN:HG3	2.41	0.56
1:A:158:ARG:NH2	4:A:463:HOH:O	2.39	0.56
1:B:189:ALA:CB	1:B:205:VAL:HG23	2.35	0.55
1:B:175:HIS:CE1	1:B:178:TYR:CD2	2.95	0.55
1:B:189:ALA:HB2	1:B:205:VAL:CG2	2.36	0.55
1:A:189:ALA:HB2	1:A:205:VAL:HG23	1.88	0.55
2:A:400:NDP:H41N	3:A:402:PHH:C5	2.37	0.54
1:B:238:TYR:CE2	1:B:242:GLN:HG3	2.42	0.54
1:A:161:LEU:HD11	1:A:192:MET:HE1	1.90	0.54
1:A:150:TYR:HB2	1:A:196:MET:HE2	1.91	0.52
2:A:400:NDP:H52A	4:A:498:HOH:O	2.09	0.52
1:A:243:TRP:O	1:A:281:ALA:HB2	2.10	0.52
1:A:238:TYR:CZ	1:A:242:GLN:HG3	2.44	0.52
1:B:63:SER:HB3	1:B:66:SER:HB2	1.93	0.51
1:A:190:ARG:NH2	4:A:447:HOH:O	2.44	0.51
1:A:227:GLY:HA2	1:A:230:LEU:HG	1.93	0.50
1:A:164:SER:OG	3:A:402:PHH:O11	2.28	0.50
1:A:172:VAL:HB	4:A:554:HOH:O	2.11	0.50
1:B:213:THR:O	1:B:216:TYR:HB3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:ARG:NH1	4:A:587:HOH:O	2.41	0.50
1:A:63:SER:HB3	1:A:66:SER:HB2	1.93	0.49
1:A:52:ARG:HA	1:B:22:PRO:HG3	1.95	0.49
1:B:120:PHE:CE2	1:B:175:HIS:CD2	3.00	0.48
1:A:161:LEU:HD11	1:A:192:MET:CE	2.43	0.48
1:A:89:GLY:HA3	4:A:595:HOH:O	2.13	0.48
1:B:182:LYS:NZ	4:B:583:HOH:O	2.45	0.48
1:A:35:THR:HA	1:A:59:ASN:HB3	1.96	0.48
1:B:35:THR:HA	1:B:59:ASN:HB3	1.95	0.48
1:A:43:ARG:O	1:A:47:MET:HG2	2.14	0.48
1:B:152:HIS:HE1	4:B:444:HOH:O	1.97	0.48
1:A:220:CYS:HA	3:A:402:PHH:CL9	2.51	0.47
1:A:219:VAL:HG23	1:A:223:TYR:HE1	1.79	0.47
1:A:166:THR:HB	1:A:179:SER:OG	2.14	0.47
1:B:33:LEU:C	1:B:33:LEU:HD23	2.39	0.47
1:A:189:ALA:CB	1:A:205:VAL:HG23	2.45	0.47
1:A:114:ASN:O	2:A:400:NDP:H4D	2.16	0.46
1:B:111:VAL:HG23	1:B:153:LEU:HD21	1.98	0.46
1:A:213:THR:O	1:A:216:TYR:HB3	2.16	0.45
1:A:224:ILE:HG12	1:A:238:TYR:CE1	2.51	0.45
1:A:98:PHE:HB3	1:A:152:HIS:CD2	2.51	0.45
1:B:165:ILE:CD1	1:B:283:MET:HE3	2.38	0.45
1:A:230:LEU:HD22	1:A:234:GLU:HB3	1.98	0.45
1:A:216:TYR:O	1:A:220:CYS:HB3	2.16	0.45
1:A:75:LYS:HA	1:A:79:SER:O	2.17	0.44
1:B:244:SER:HA	1:B:245:PRO:HD3	1.88	0.44
1:A:37:ALA:HB3	1:A:58:VAL:HG13	1.98	0.44
1:A:200:LYS:NZ	4:A:490:HOH:O	2.49	0.43
1:A:138:ASN:O	1:A:142:GLN:HG3	2.19	0.43
1:B:213:THR:H	1:B:216:TYR:CB	2.30	0.42
1:A:224:ILE:HG12	1:A:238:TYR:CD1	2.55	0.42
1:A:212:LYS:HG3	1:A:249:VAL:HG11	2.02	0.42
1:B:62:ASN:O	1:B:64:THR:N	2.53	0.42
1:A:224:ILE:HG21	1:A:230:LEU:HD11	2.01	0.41
1:A:88:VAL:CG1	1:A:141:GLY:HA3	2.50	0.41
1:A:61:ALA:O	1:A:85:LYS:HE3	2.21	0.41
1:B:37:ALA:HB3	1:B:58:VAL:HG13	2.03	0.41
1:A:22:PRO:HG3	1:B:52:ARG:HA	2.02	0.41
1:A:238:TYR:CE2	1:A:242:GLN:HG3	2.56	0.41
1:B:150:TYR:CB	1:B:196:MET:HE3	2.49	0.41
1:A:172:VAL:HG12	1:A:175:HIS:HB2	2.02	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	271/283 (96%)	266 (98%)	5 (2%)	0	100	100
1	B	250/283 (88%)	239 (96%)	10 (4%)	1 (0%)	30	27
All	All	521/566 (92%)	505 (97%)	15 (3%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	63	SER

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	213/221 (96%)	205 (96%)	8 (4%)	29	29
1	B	195/221 (88%)	189 (97%)	6 (3%)	35	37
All	All	408/442 (92%)	394 (97%)	14 (3%)	32	33

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	ARG
1	A	57	ILE

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Mol	Chain	Res	Type
1	A	65	GLU
1	A	79	SER
1	A	88	VAL
1	A	129	GLU
1	A	158	ARG
1	A	175	HIS
1	B	57	ILE
1	B	65	GLU
1	B	79	SER
1	B	129	GLU
1	B	175	HIS
1	B	236	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	HIS
1	B	62	ASN
1	B	152	HIS
1	B	175	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 oligosaccharides 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
2	NDP	A	400	-	51,52,52	1.43	7 (13%)	71,80,80	1.93	14 (19%)
2	NDP	B	401	-	51,52,52	1.40	7 (13%)	71,80,80	1.90	14 (19%)
3	PHH	A	402	-	15,15,15	3.63	3 (20%)	20,23,23	2.44	5 (25%)

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	NDP	A	400	-	-	9/34/77/77	0/5/5/5
2	NDP	B	401	-	-	12/34/77/77	0/5/5/5
3	PHH	A	402	-	-	-	0/2/2/2

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	PHH	O4-C5	9.86	1.48	1.36
3	A	402	PHH	O4-C3	-8.93	1.34	1.45
2	A	400	NDP	C4N-C3N	-5.18	1.40	1.50
2	B	401	NDP	C4N-C3N	-4.76	1.41	1.50
2	B	401	NDP	C5A-N7A	-3.81	1.32	1.39
3	A	402	PHH	O11-C5	3.42	1.32	1.22
2	A	400	NDP	C5A-N7A	-3.33	1.33	1.39
2	A	400	NDP	C4N-C5N	-3.14	1.40	1.49
2	B	401	NDP	C4N-C5N	-3.09	1.41	1.49
2	A	400	NDP	C6N-C5N	2.71	1.41	1.33
2	B	401	NDP	PA-O3	2.66	1.62	1.59
2	A	400	NDP	C7N-C3N	2.61	1.54	1.48
2	B	401	NDP	C4A-N9A	-2.57	1.32	1.37
2	B	401	NDP	C6N-C5N	2.36	1.40	1.33
2	B	401	NDP	C7N-C3N	2.35	1.53	1.48
2	A	400	NDP	C2N-C3N	2.23	1.41	1.35
2	A	400	NDP	P2B-O3X	-2.12	1.46	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	NDP	C5A-C4A-N3A	-6.47	117.81	126.72
3	A	402	PHH	O4-C3-C2	6.34	109.59	104.57
2	B	401	NDP	N3A-C4A-N9A	6.22	137.75	127.17
2	A	400	NDP	N3A-C4A-N9A	6.21	137.73	127.17
2	B	401	NDP	C5A-C4A-N3A	-6.15	118.25	126.72
2	A	400	NDP	O4B-C1B-N9A	5.47	118.59	108.09
3	A	402	PHH	C3-O4-C5	-5.33	107.16	110.64
3	A	402	PHH	O4-C5-O11	5.11	126.48	121.09
2	B	401	NDP	O4B-C1B-N9A	4.91	117.53	108.09
2	B	401	NDP	N3A-C2A-N1A	-4.75	121.40	128.58
2	A	400	NDP	N3A-C2A-N1A	-4.50	121.76	128.58
2	B	401	NDP	C6A-C5A-C4A	3.82	122.40	117.18
2	A	400	NDP	C6A-C5A-C4A	3.78	122.34	117.18
2	A	400	NDP	C4B-O4B-C1B	-3.55	101.64	109.47
2	A	400	NDP	C2A-N3A-C4A	3.47	120.32	111.83
2	B	401	NDP	C2A-N3A-C4A	3.47	120.30	111.83
2	B	401	NDP	C6A-C5A-N7A	-3.45	125.44	132.09
3	A	402	PHH	O11-C5-C6	-3.21	125.27	130.99
2	B	401	NDP	C4A-N9A-C8A	3.20	109.10	105.74
2	B	401	NDP	C4B-O4B-C1B	-3.11	102.61	109.47
2	A	400	NDP	C6A-C5A-N7A	-2.87	126.55	132.09
2	A	400	NDP	C4A-N9A-C8A	2.86	108.74	105.74
2	A	400	NDP	N9A-C8A-N7A	-2.82	109.94	113.94
2	A	400	NDP	C2B-C1B-N9A	-2.78	109.17	113.75
2	B	401	NDP	N9A-C8A-N7A	-2.64	110.18	113.94
2	B	401	NDP	O3B-C3B-C2B	-2.54	104.07	111.19
2	A	400	NDP	O5B-C5B-C4B	2.37	117.06	108.99
3	A	402	PHH	C2-C6-C5	-2.29	105.84	108.39
2	A	400	NDP	O3B-C3B-C2B	-2.21	105.00	111.19
2	B	401	NDP	O5B-C5B-C4B	2.18	116.42	108.99
2	B	401	NDP	O2N-PN-O1N	2.06	122.02	112.44
2	A	400	NDP	C2A-N1A-C6A	2.05	122.10	118.73
2	B	401	NDP	O2A-PA-O1A	2.04	121.91	112.44

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	NDP	C5B-O5B-PA-O1A
2	A	400	NDP	C5B-O5B-PA-O2A
2	B	401	NDP	O4B-C4B-C5B-O5B

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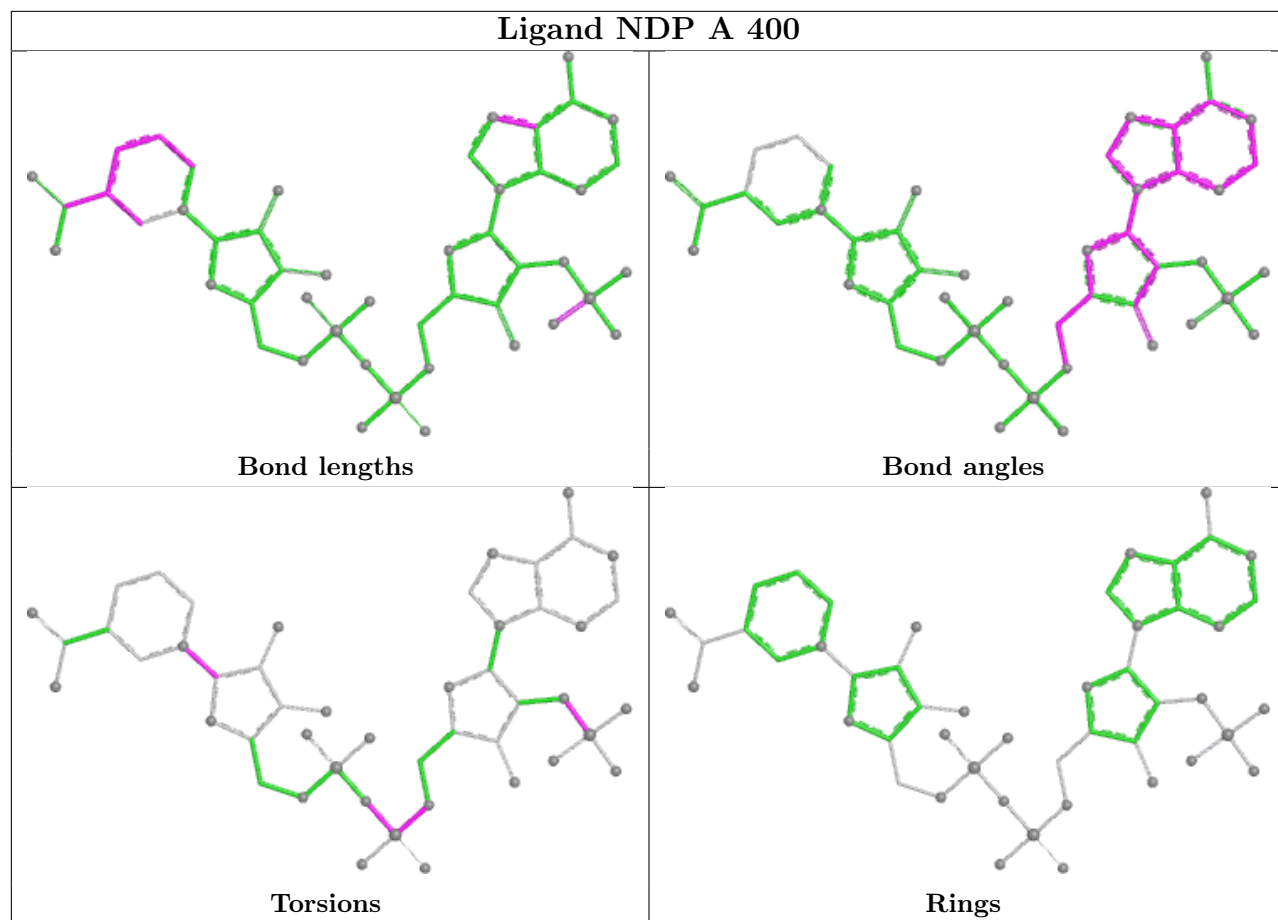
Mol	Chain	Res	Type	Atoms
2	B	401	NDP	C5D-O5D-PN-O1N
2	B	401	NDP	C3B-C4B-C5B-O5B
2	A	400	NDP	PN-O3-PA-O2A
2	B	401	NDP	PN-O3-PA-O2A
2	A	400	NDP	O4D-C1D-N1N-C6N
2	A	400	NDP	C5B-O5B-PA-O3
2	B	401	NDP	C5B-O5B-PA-O1A
2	B	401	NDP	C5B-O5B-PA-O2A
2	B	401	NDP	C5B-O5B-PA-O3
2	B	401	NDP	C5D-O5D-PN-O3
2	B	401	NDP	C5D-O5D-PN-O2N
2	B	401	NDP	PN-O3-PA-O1A
2	B	401	NDP	O4D-C1D-N1N-C6N
2	A	400	NDP	C2B-O2B-P2B-O1X
2	B	401	NDP	C2B-O2B-P2B-O1X
2	A	400	NDP	C2B-O2B-P2B-O2X
2	A	400	NDP	C2B-O2B-P2B-O3X
2	A	400	NDP	PN-O3-PA-O1A

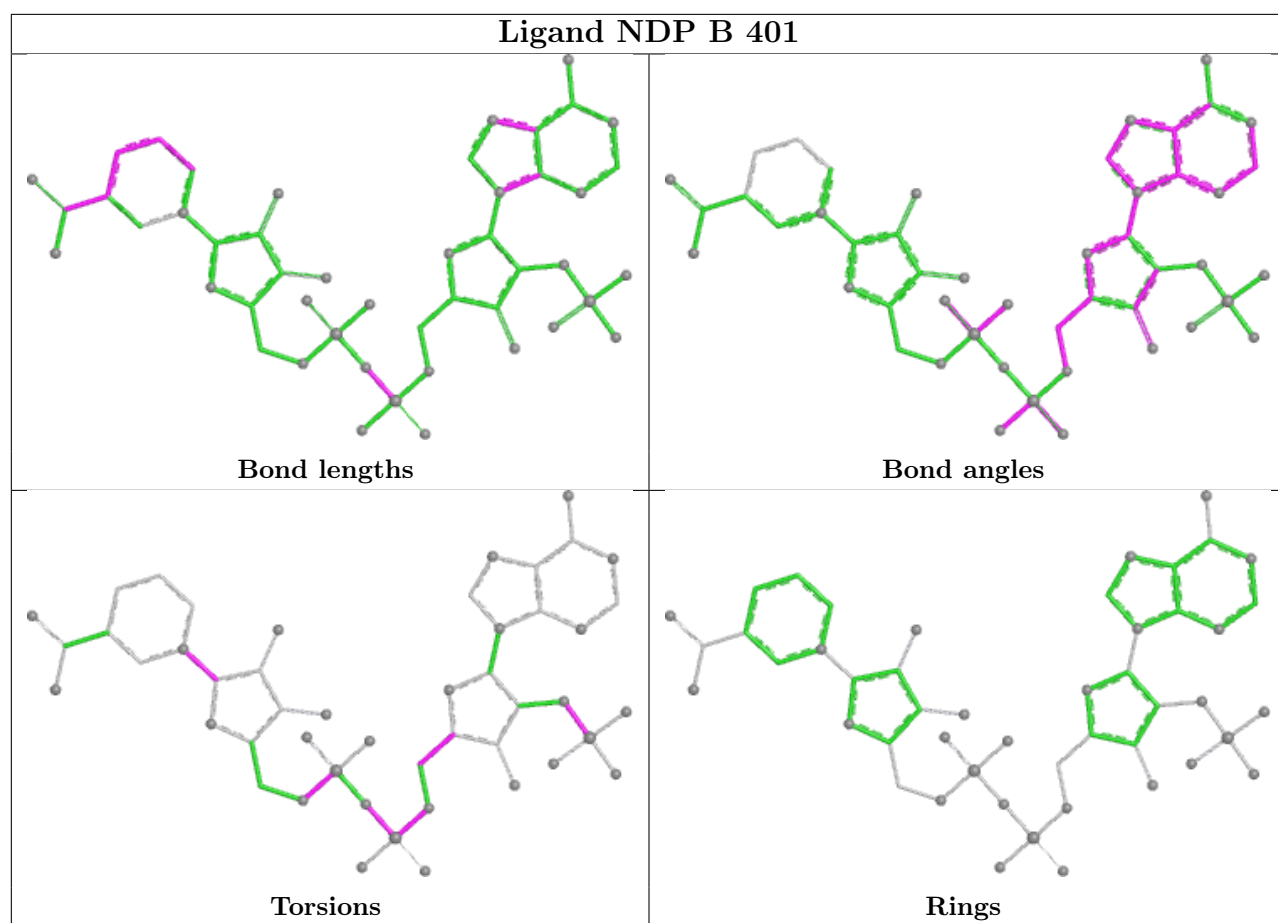
There are no ring outliers.

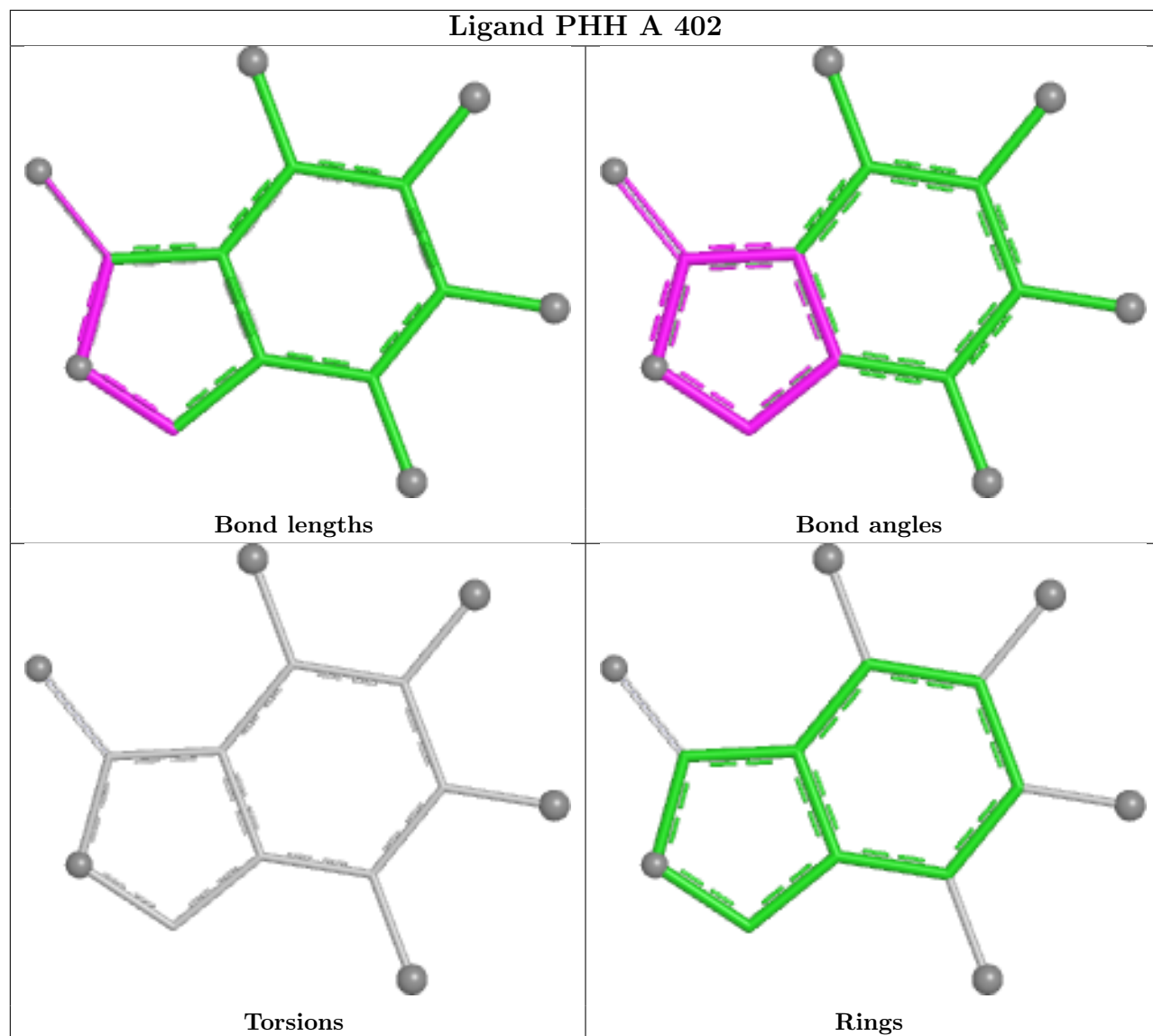
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	NDP	3	0
3	A	402	PHH	4	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.