



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:29 PM UTC

PDB ID : 1IDE / pdb_00001ide
Title : ISOCITRATE DEHYDROGENASE Y160F MUTANT STEADY-STATE INTERMEDIATE COMPLEX (LAUE DETERMINATION)
Authors : Bolduc, J.M.; Dyer, D.H.; Scott, W.G.; Singer, P.; Sweet, R.M.; Koshland Junior, D.E.; Stoddard, B.L.
Deposited on : 1995-01-18
Resolution : 2.50 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

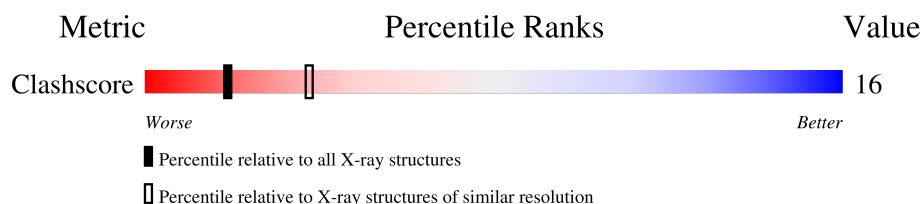
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)

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	416	 65% 30% .

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
4	NAP	A	419	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3954 atoms, of which 695 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 ISOCITRATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	414	3879	2035	684	538	604	18	0	0	0

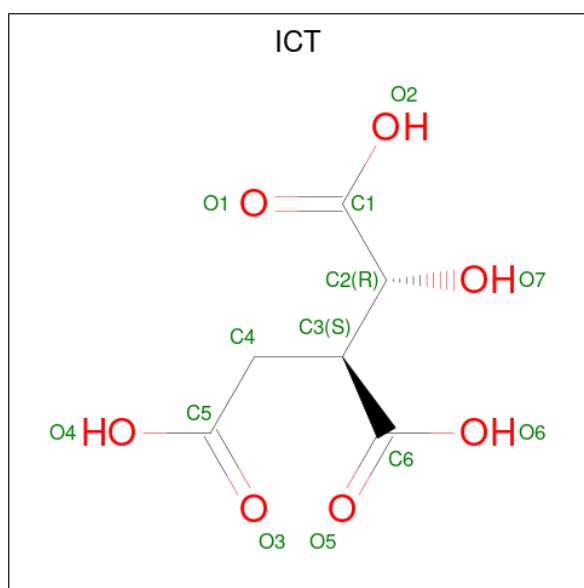
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	160	PHE	TYR	engineered mutation	UNP P08200

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

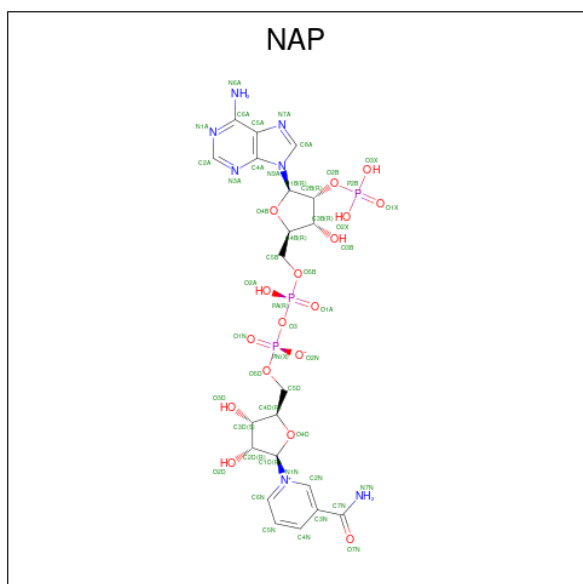
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ISOCITRIC ACID (CCD ID: ICT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $\text{C}_{21}\text{H}_{28}\text{N}_7\text{O}_{17}\text{P}_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	0	0
			55	21	7	7	17	3		

- Molecule 5 is water.

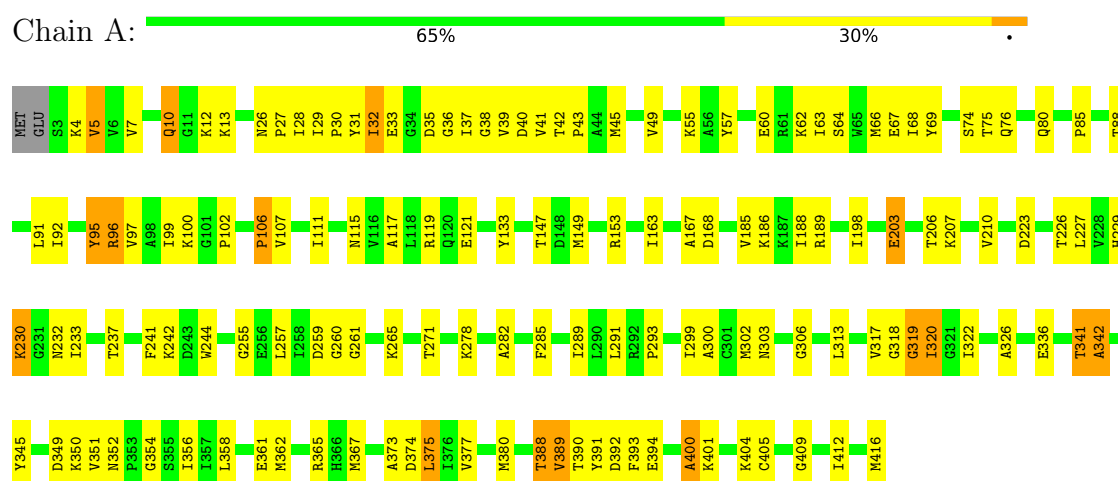
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	2	Total	H	O	0	0
			6	4	2		

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: ISOCITRATE DEHYDROGENASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.10Å 105.10Å 150.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.192 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3954	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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: MG, NAP, ICT

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	0.52	1/3256 (0.0%)	1.27	39/4403 (0.9%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	341	THR	CA-C	-5.42	1.45	1.52

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	230	LYS	N-CA-C	-8.91	96.26	109.62
1	A	4	LYS	N-CA-C	8.23	121.89	108.55
1	A	38	GLY	N-CA-C	-7.87	103.23	112.83
1	A	189	ARG	N-CA-C	7.66	119.26	111.07
1	A	133	TYR	N-CA-C	7.32	120.20	108.41
1	A	342	ALA	CA-C-N	7.31	128.98	119.84
1	A	342	ALA	C-N-CA	7.31	128.98	119.84
1	A	63	ILE	N-CA-C	7.26	118.33	108.17
1	A	282	ALA	N-CA-C	7.01	119.52	111.11
1	A	188	ILE	N-CA-C	-6.94	98.27	108.54
1	A	400	ALA	N-CA-C	-6.79	99.90	110.42
1	A	203	GLU	N-CA-C	-6.70	103.07	111.11

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	TYR	N-CA-C	6.68	120.05	108.76
1	A	107	VAL	N-CA-C	-6.63	100.98	109.80
1	A	233	ILE	N-CA-C	-6.42	106.04	111.56
1	A	241	PHE	N-CA-C	-6.37	104.26	111.07
1	A	375	LEU	N-CA-C	-6.31	104.33	111.14
1	A	401	LYS	N-CA-C	-6.24	101.25	110.48
1	A	388	THR	N-CA-C	-6.24	96.50	107.73
1	A	320	ILE	N-CA-C	-5.95	103.65	111.09
1	A	7	VAL	N-CA-C	-5.95	103.49	109.02
1	A	291	LEU	N-CA-C	5.76	118.68	111.24
1	A	336	GLU	N-CA-C	5.76	117.55	108.96
1	A	223	ASP	N-CA-C	5.57	119.25	112.23
1	A	186	LYS	N-CA-C	-5.56	105.64	113.20
1	A	32	ILE	N-CA-C	-5.48	100.00	107.99
1	A	185	VAL	N-CA-C	5.43	112.03	106.21
1	A	100	LYS	N-CA-C	5.39	117.59	109.23
1	A	271	THR	N-CA-C	5.38	122.81	114.64
1	A	5	VAL	N-CA-C	-5.37	99.91	107.75
1	A	303	ASN	N-CA-C	5.34	116.80	110.97
1	A	33	GLU	N-CA-C	5.27	116.71	110.97
1	A	237	THR	N-CA-C	-5.13	103.70	111.34
1	A	389	VAL	N-CA-C	5.10	115.04	108.82
1	A	106	PRO	N-CA-C	5.07	119.16	110.95
1	A	302	MET	N-CA-C	-5.06	103.37	110.35
1	A	319	GLY	N-CA-C	5.04	125.13	113.18
1	A	10	GLN	N-CA-C	5.04	119.14	112.89
1	A	96	ARG	N-CA-C	5.00	121.46	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	95	TYR	Sidechain

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	3195	684	3222	103	1
2	A	1	0	0	0	0
3	A	13	0	4	0	1
4	A	48	7	24	0	0
5	A	2	4	0	0	0
All	All	3259	695	3250	103	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 16.

All (103) 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:342:ALA:HB1	1:A:345:TYR:HD2	1.30	0.95
1:A:37:ILE:HG12	1:A:341:THR:HA	1.50	0.89
1:A:37:ILE:HD12	1:A:351:VAL:HG21	1.55	0.88
1:A:5:VAL:HG11	1:A:91:LEU:HD21	1.55	0.87
1:A:37:ILE:H	1:A:341:THR:HB	1.40	0.85
1:A:35:ASP:O	1:A:341:THR:HG21	1.80	0.81
1:A:37:ILE:N	1:A:341:THR:HB	1.97	0.79
1:A:37:ILE:HG23	1:A:341:THR:O	1.83	0.78
1:A:41:VAL:HG13	1:A:356:ILE:HG21	1.66	0.77
1:A:106:PRO:HG2	1:A:111:ILE:HD13	1.66	0.76
1:A:102:PRO:HB2	1:A:341:THR:HG23	1.68	0.76
1:A:75:THR:HB	1:A:80:GLN:HA	1.68	0.75
1:A:37:ILE:CD1	1:A:351:VAL:HG21	2.17	0.74
1:A:389:VAL:HG22	1:A:390:THR:H	1.54	0.71
1:A:342:ALA:HB1	1:A:345:TYR:CD2	2.21	0.71
1:A:257:LEU:HD11	1:A:261:GLY:HA2	1.74	0.70
1:A:358:LEU:HD21	1:A:380:MET:HE2	1.74	0.70
1:A:361:GLU:HG3	1:A:377:VAL:HG12	1.74	0.67
1:A:389:VAL:HG22	1:A:390:THR:N	2.13	0.64
1:A:394:GLU:HG2	1:A:400:ALA:HB3	1.80	0.64
1:A:404:LYS:HZ2	1:A:404:LYS:HB2	1.61	0.64
1:A:342:ALA:CB	1:A:345:TYR:HD2	2.08	0.63
1:A:349:ASP:OD1	1:A:405:CYS:HB3	2.00	0.62
1:A:389:VAL:HG21	1:A:393:PHE:CD2	2.34	0.62
1:A:26:ASN:HA	1:A:62:LYS:O	2.01	0.60
1:A:37:ILE:HG12	1:A:341:THR:CA	2.25	0.60
1:A:255:GLY:HA2	1:A:265:LYS:O	2.01	0.60
1:A:37:ILE:HG21	1:A:351:VAL:HG11	1.83	0.60
1:A:68:ILE:HD12	1:A:88:THR:HG23	1.83	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ASP:O	1:A:43:PRO:HD2	2.02	0.60
1:A:147:THR:HG23	1:A:149:MET:HE3	1.84	0.59
1:A:36:GLY:O	1:A:39:VAL:HG12	2.01	0.59
1:A:319:GLY:HA3	1:A:322:ILE:HG12	1.85	0.59
1:A:76:GLN:HE21	1:A:76:GLN:HA	1.68	0.58
1:A:388:THR:HG23	1:A:388:THR:O	2.04	0.58
1:A:41:VAL:HG13	1:A:356:ILE:CG2	2.34	0.58
1:A:45:MET:HE2	1:A:99:ILE:HD13	1.85	0.58
1:A:289:ILE:O	1:A:293:PRO:HB3	2.04	0.58
1:A:374:ASP:HA	1:A:377:VAL:HG13	1.86	0.58
1:A:57:TYR:O	1:A:60:GLU:HG2	2.05	0.57
1:A:365:ARG:HH21	1:A:373:ALA:HB1	1.71	0.56
1:A:163:ILE:HB	1:A:198:ILE:HB	1.88	0.55
1:A:203:GLU:HG2	1:A:244:TRP:CE2	2.41	0.55
1:A:37:ILE:HD12	1:A:351:VAL:CG2	2.34	0.55
1:A:30:PRO:HD2	1:A:97:VAL:O	2.06	0.55
1:A:102:PRO:HB2	1:A:341:THR:CG2	2.36	0.54
1:A:29:ILE:HD11	1:A:367:MET:HE3	1.89	0.53
1:A:32:ILE:HG12	1:A:68:ILE:HG13	1.91	0.53
1:A:37:ILE:HG12	1:A:341:THR:HG22	1.91	0.53
1:A:358:LEU:O	1:A:361:GLU:HB3	2.09	0.52
1:A:389:VAL:CG2	1:A:390:THR:H	2.21	0.52
1:A:227:LEU:HD12	1:A:300:ALA:HB3	1.93	0.50
1:A:115:ASN:O	1:A:119:ARG:HD3	2.11	0.50
1:A:12:LYS:O	1:A:27:PRO:HA	2.12	0.49
1:A:37:ILE:CG1	1:A:341:THR:HA	2.32	0.49
1:A:102:PRO:CB	1:A:341:THR:HG23	2.38	0.49
1:A:55:LYS:HE2	1:A:416:MET:O	2.13	0.49
1:A:391:TYR:CG	1:A:392:ASP:N	2.80	0.49
1:A:285:PHE:CZ	1:A:299:ILE:HD13	2.48	0.49
1:A:285:PHE:HZ	1:A:299:ILE:CD1	2.26	0.48
1:A:74:SER:OG	1:A:85:PRO:HD3	2.13	0.48
1:A:10:GLN:O	1:A:26:ASN:ND2	2.47	0.48
1:A:37:ILE:HG21	1:A:351:VAL:CG1	2.43	0.48
1:A:358:LEU:HD21	1:A:380:MET:CE	2.43	0.48
1:A:5:VAL:HG11	1:A:91:LEU:CD2	2.36	0.47
1:A:257:LEU:HD11	1:A:261:GLY:CA	2.44	0.47
1:A:45:MET:O	1:A:49:VAL:HG23	2.15	0.46
1:A:326:ALA:HB2	1:A:362:MET:CE	2.46	0.46
1:A:389:VAL:CG2	1:A:390:THR:N	2.79	0.46
1:A:88:THR:O	1:A:92:ILE:HG13	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:THR:O	1:A:210:VAL:HG23	2.15	0.45
1:A:149:MET:SD	1:A:313:LEU:HD21	2.57	0.45
1:A:117:ALA:O	1:A:121:GLU:HB2	2.18	0.45
1:A:203:GLU:HG2	1:A:244:TRP:NE1	2.32	0.44
1:A:153:ARG:HG3	1:A:306:GLY:HA3	1.99	0.44
1:A:167:ALA:O	1:A:168:ASP:HB2	2.17	0.44
1:A:389:VAL:HG21	1:A:393:PHE:HD2	1.82	0.44
1:A:352:ASN:HA	1:A:390:THR:HG21	1.99	0.44
1:A:28:ILE:HA	1:A:64:SER:O	2.19	0.43
1:A:326:ALA:HB2	1:A:362:MET:HE3	2.00	0.43
1:A:30:PRO:HA	1:A:66:MET:O	2.18	0.43
1:A:354:GLY:CA	1:A:380:MET:HE1	2.49	0.43
1:A:31:TYR:CE1	1:A:67:GLU:HG3	2.54	0.43
1:A:229:HIS:CE1	1:A:242:LYS:HD3	2.54	0.42
1:A:313:LEU:O	1:A:317:VAL:HG22	2.19	0.42
1:A:230:LYS:HD2	1:A:230:LYS:HA	1.65	0.42
1:A:75:THR:HB	1:A:80:GLN:CA	2.43	0.42
1:A:350:LYS:O	1:A:350:LYS:HG2	2.19	0.42
1:A:111:ILE:HD12	1:A:111:ILE:H	1.84	0.42
1:A:341:THR:O	1:A:342:ALA:C	2.60	0.42
1:A:226:THR:HA	1:A:278:LYS:O	2.21	0.41
1:A:259:ASP:HB3	1:A:260:GLY:H	1.71	0.41
1:A:285:PHE:HZ	1:A:299:ILE:HD13	1.84	0.41
1:A:409:GLY:HA2	1:A:412:ILE:HD12	2.02	0.41
1:A:203:GLU:O	1:A:207:LYS:HB2	2.21	0.41
1:A:37:ILE:HG12	1:A:341:THR:CB	2.50	0.41
1:A:41:VAL:HG12	1:A:42:THR:N	2.36	0.41
1:A:318:GLY:O	1:A:320:ILE:N	2.54	0.41
1:A:95:TYR:O	1:A:96:ARG:HB2	2.20	0.41
1:A:12:LYS:HD2	1:A:13:LYS:N	2.37	0.41
1:A:341:THR:O	1:A:342:ALA:O	2.39	0.40
1:A:375:LEU:HD12	1:A:375:LEU:HA	1.67	0.40
1:A:62:LYS:HB2	1:A:62:LYS:HE3	1.81	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:A:232:ASN:HD21	3:A:418:ICT:O4[7_555]	1.38	0.22

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z > 2$
4	NAP	A	419	-	50,52,52	1.96	10 (20%)	71,80,80	1.96	18 (25%)
3	ICT	A	418	2	12,12,12	2.24	2 (16%)	13,16,16	2.45	6 (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
4	NAP	A	419	-	1/1/12/12	12/35/67/67	0/5/5/5
3	ICT	A	418	2	-	2/16/16/16	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	419	NAP	C3N-C7N	-6.22	1.41	1.50
4	A	419	NAP	PA-O3	5.38	1.65	1.59
4	A	419	NAP	PN-O3	5.21	1.65	1.59
3	A	418	ICT	C3-C6	4.95	1.59	1.51
3	A	418	ICT	C3-C2	-4.64	1.48	1.54
4	A	419	NAP	C2N-C3N	4.14	1.45	1.39
4	A	419	NAP	C7N-N7N	3.84	1.40	1.33
4	A	419	NAP	C4N-C3N	2.86	1.43	1.39
4	A	419	NAP	O4D-C4D	-2.70	1.39	1.45
4	A	419	NAP	C5N-C4N	-2.43	1.34	1.38
4	A	419	NAP	P2B-O3X	-2.30	1.46	1.54
4	A	419	NAP	O4B-C4B	-2.24	1.40	1.45

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	419	NAP	O4B-C1B-N9A	6.53	120.64	108.09
3	A	418	ICT	O6-C6-O5	-5.37	111.90	124.08
4	A	419	NAP	O2A-PA-O3	5.13	121.13	107.27
4	A	419	NAP	C3B-C2B-C1B	-4.63	93.94	102.81
3	A	418	ICT	O6-C6-C3	4.22	124.88	113.93
4	A	419	NAP	C4D-O4D-C1D	3.80	113.40	109.92
4	A	419	NAP	N3A-C2A-N1A	-3.57	123.17	128.58
4	A	419	NAP	C3N-C7N-N7N	3.30	121.80	117.74
4	A	419	NAP	O7N-C7N-N7N	-3.28	117.87	122.62
4	A	419	NAP	C4A-N9A-C8A	-3.09	102.50	105.74
4	A	419	NAP	C6N-N1N-C2N	-3.06	119.27	121.88
4	A	419	NAP	O4B-C1B-C2B	-3.01	101.41	106.59
4	A	419	NAP	O3-PA-O1A	-2.65	102.72	110.70
4	A	419	NAP	O4D-C4D-C5D	2.64	117.78	109.33
3	A	418	ICT	O7-C2-C1	-2.54	105.26	110.69
4	A	419	NAP	O3B-C3B-C4B	2.51	118.29	111.08
3	A	418	ICT	O7-C2-C3	2.46	116.19	110.63
4	A	419	NAP	O4B-C4B-C3B	-2.43	100.33	105.15
4	A	419	NAP	O5D-C5D-C4D	2.30	116.82	108.99

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	419	NAP	O3B-C3B-C2B	2.25	117.48	111.19
3	A	418	ICT	O2-C1-C2	2.23	119.50	113.31
3	A	418	ICT	O2-C1-O1	-2.11	119.29	124.08
4	A	419	NAP	O3-PN-O1N	2.11	117.05	110.70
4	A	419	NAP	PN-O5D-C5D	2.01	132.88	121.35

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	419	NAP	C3D

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	418	ICT	C1-C2-C3-C6
4	A	419	NAP	C5D-O5D-PN-O2N
4	A	419	NAP	O4D-C4D-C5D-O5D
4	A	419	NAP	O4D-C1D-N1N-C2N
4	A	419	NAP	O4D-C1D-N1N-C6N
4	A	419	NAP	C2D-C1D-N1N-C2N
4	A	419	NAP	C2D-C1D-N1N-C6N
4	A	419	NAP	C3D-C4D-C5D-O5D
4	A	419	NAP	C3B-C2B-O2B-P2B
4	A	419	NAP	C1B-C2B-O2B-P2B
3	A	418	ICT	O7-C2-C3-C6
4	A	419	NAP	C5D-O5D-PN-O3
4	A	419	NAP	C5D-O5D-PN-O1N
4	A	419	NAP	C4D-C5D-O5D-PN

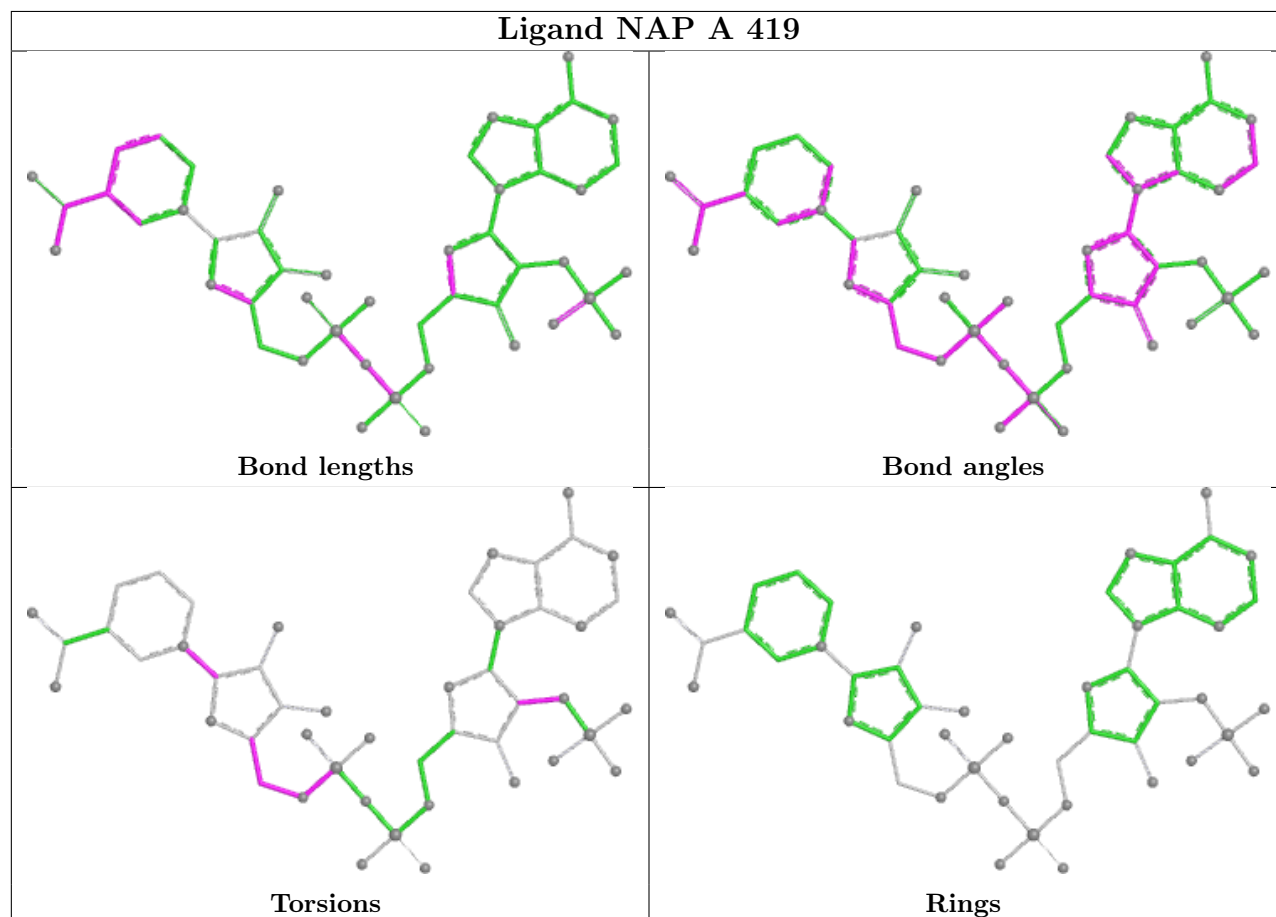
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	418	ICT	0	1

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.