



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:39 PM UTC

PDB ID : 4AOU / pdb_00004aou
Title : CtIDH bound to NADP. The complex structures of Isocitrate dehydrogenase from *Clostridium thermocellum* and *Desulfotalea psychrophila*, support a new active site locking mechanism
Authors : Leiros, H.-K.S.; Fedoy, A.-E.; Leiros, I.; Steen, I.H.
Deposited on : 2012-03-30
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)	:	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

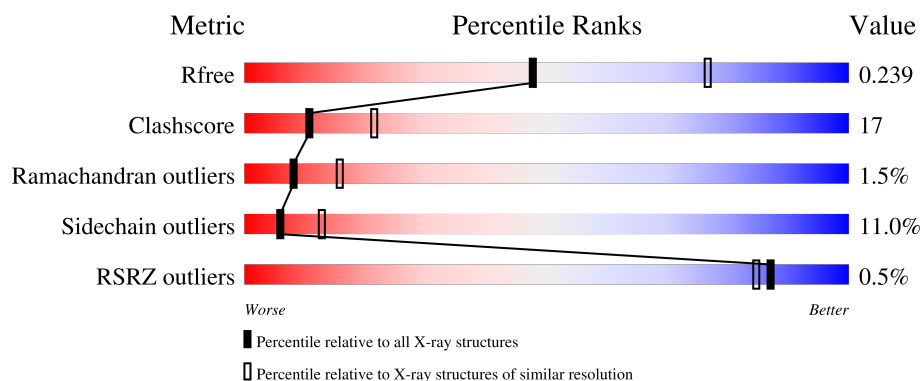
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 $\leq 5\%$. 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	402	63% 29% 7%

2 Entry composition [i](#)

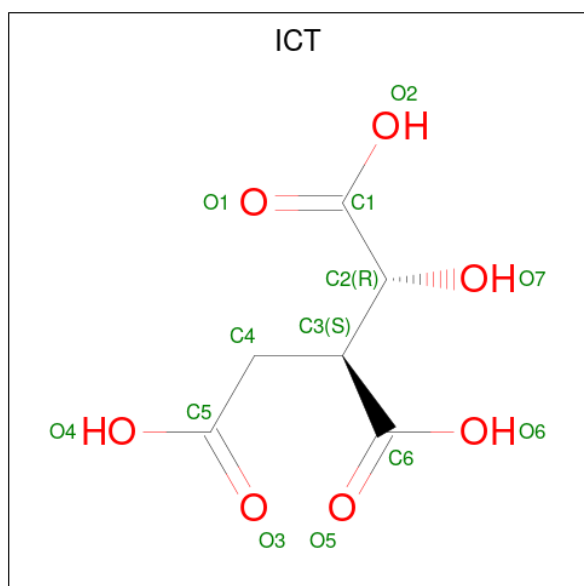
There are 5 unique types of molecules in this entry. The entry contains 3313 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ISOCITRATE DEHYDROGENASE [NADP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	29	0	0
			3210	2042	535	615	18			

- Molecule 2 is ISOCITRIC ACID (CCD ID: ICT) (formula: $C_6H_8O_7$).



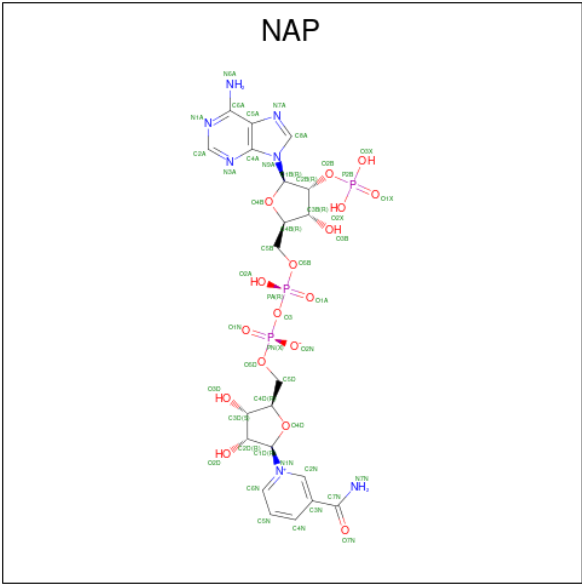
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD

ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	8	0
			48	21	7	17		

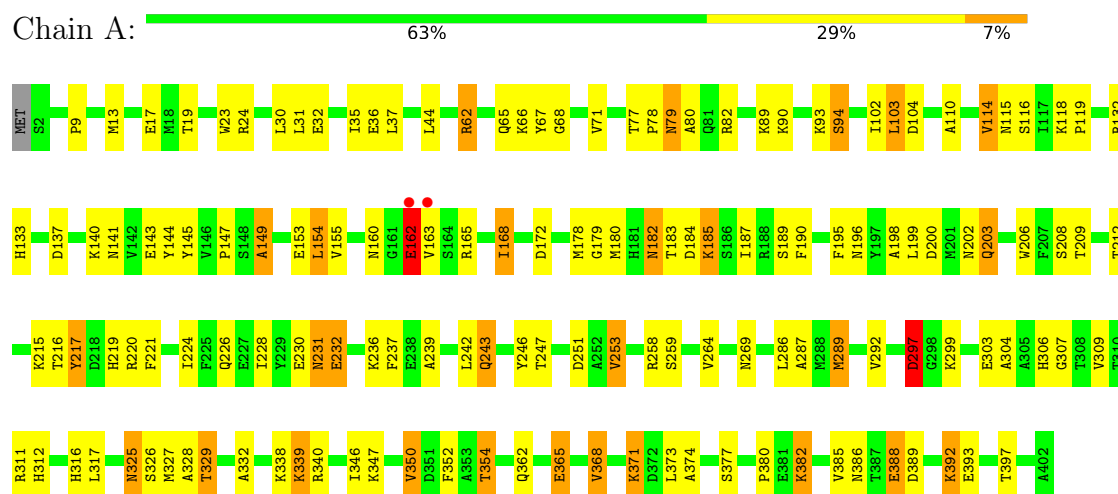
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	41	Total	0	0
			41 O		

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 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: ISOCITRATE DEHYDROGENASE [NADP]



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.20Å 129.20Å 60.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.00 – 2.50 45.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.00-2.50) 99.8 (45.00-2.50)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, R_{free}	0.189 , 0.244 0.189 , 0.239	Depositor DCC
R_{free} test set	1043 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	48.2	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 27.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.044 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3313	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: NAP, MG, 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	1.20	9/3275 (0.3%)	1.36	25/4416 (0.6%)

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 (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	371	LYS	CE-NZ	-21.84	0.83	1.49
1	A	392	LYS	CE-NZ	-16.75	0.99	1.49
1	A	162	GLU	CG-CD	-15.21	1.14	1.52
1	A	392	LYS	CG-CD	-14.52	1.08	1.52
1	A	243	GLN	CD-NE2	10.17	1.54	1.33
1	A	185	LYS	CD-CE	-7.24	1.30	1.52
1	A	371	LYS	CG-CD	-5.59	1.35	1.52
1	A	354	THR	CA-CB	5.48	1.61	1.53
1	A	350	VAL	CA-CB	5.35	1.61	1.54

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	GLN	OE1-CD-NE2	-24.19	98.41	122.60
1	A	371	LYS	CB-CG-CD	20.54	158.53	111.30
1	A	392	LYS	CB-CG-CD	13.43	142.18	111.30
1	A	32	GLU	CA-C-N	-9.71	109.56	120.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	GLU	C-N-CA	-9.71	109.56	120.04
1	A	253	VAL	CB-CA-C	-9.41	99.44	112.14
1	A	339	LYS	CG-CD-CE	7.73	129.07	111.30
1	A	36	GLU	CB-CG-CD	7.49	125.33	112.60
1	A	371	LYS	CG-CD-CE	6.95	127.30	111.30
1	A	347	LYS	CA-CB-CG	6.58	127.26	114.10
1	A	346	ILE	N-CA-C	6.35	113.00	106.21
1	A	94	SER	CA-C-N	-5.90	113.60	119.56
1	A	94	SER	C-N-CA	-5.90	113.60	119.56
1	A	162	GLU	CB-CG-CD	5.81	122.47	112.60
1	A	346	ILE	CB-CA-C	-5.78	106.75	113.22
1	A	68	GLY	N-CA-C	5.62	121.38	114.69
1	A	332	ALA	N-CA-C	-5.58	105.28	111.36
1	A	340	ARG	N-CA-C	-5.54	105.14	111.07
1	A	297	ASP	CB-CA-C	5.49	120.31	110.81
1	A	212	THR	CA-C-N	-5.35	117.47	122.66
1	A	212	THR	C-N-CA	-5.35	117.47	122.66
1	A	312	HIS	N-CA-C	-5.33	105.55	111.36
1	A	259	SER	N-CA-C	5.27	116.73	110.19
1	A	149	ALA	N-CA-C	5.17	116.97	110.24
1	A	17	GLU	CB-CA-C	-5.02	110.77	116.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	GLN	Sidechain

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	3210	0	3180	107	0
2	A	13	0	4	0	0
3	A	1	0	0	0	0
4	A	48	0	25	0	0
5	A	41	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3313	0	3209	107	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 17.

All (107) 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:368:VAL:HG21	1:A:393:GLU:HG2	1.44	1.00
1:A:303:GLU:HG3	1:A:304:ALA:O	1.72	0.89
1:A:226:GLN:O	1:A:230:GLU:HG2	1.73	0.88
1:A:297:ASP:O	1:A:297:ASP:OD1	1.92	0.87
1:A:325:ASN:ND2	1:A:327:MET:H	1.76	0.83
1:A:242:LEU:HD12	1:A:242:LEU:O	1.81	0.80
1:A:180:MET:HE3	1:A:217:TYR:OH	1.81	0.79
1:A:231:ASN:HD22	1:A:231:ASN:C	1.90	0.79
1:A:237:PHE:CD2	1:A:242:LEU:CD1	2.67	0.78
1:A:66:LYS:HE2	1:A:67:TYR:CZ	2.21	0.76
1:A:297:ASP:OD1	1:A:297:ASP:C	2.31	0.73
1:A:203:GLN:HE21	1:A:264:VAL:HG23	1.53	0.73
1:A:180:MET:HE3	1:A:217:TYR:CZ	2.24	0.73
1:A:237:PHE:CD2	1:A:242:LEU:HD12	2.25	0.72
1:A:325:ASN:C	1:A:325:ASN:HD22	1.98	0.71
1:A:306:HIS:HD2	1:A:307:GLY:O	1.73	0.71
1:A:325:ASN:HD21	1:A:327:MET:HG2	1.58	0.69
1:A:237:PHE:HD2	1:A:242:LEU:O	1.76	0.68
1:A:154:LEU:HD22	1:A:155:VAL:N	2.14	0.63
1:A:153:GLU:C	1:A:168:ILE:HD12	2.24	0.62
1:A:325:ASN:HD22	1:A:327:MET:H	1.45	0.62
1:A:23:TRP:HH2	1:A:71:VAL:HG13	1.63	0.62
1:A:103:LEU:N	1:A:103:LEU:HD23	2.15	0.61
1:A:180:MET:CE	1:A:217:TYR:CZ	2.83	0.61
1:A:23:TRP:CH2	1:A:71:VAL:HG13	2.38	0.59
1:A:110:ALA:C	1:A:289:MET:HE2	2.28	0.59
1:A:325:ASN:HD22	1:A:326:SER:N	2.00	0.59
1:A:65:GLN:O	1:A:299:LYS:HE3	2.01	0.59
1:A:102:ILE:HG22	1:A:103:LEU:HD23	1.84	0.58
1:A:236:LYS:O	1:A:239:ALA:HB3	2.03	0.58
1:A:377:SER:HB3	1:A:382:LYS:HD2	1.86	0.58
1:A:79:ASN:ND2	1:A:82:ARG:H	2.02	0.58
1:A:311:ARG:HG3	5:A:2032:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:SER:OG	1:A:365:GLU:OE1	2.14	0.57
1:A:89:LYS:O	1:A:90:LYS:HG2	2.04	0.57
1:A:306:HIS:HE1	5:A:2031:HOH:O	1.88	0.57
1:A:182:ASN:HD22	1:A:183:THR:H	1.54	0.55
1:A:287:ALA:O	1:A:328:ALA:HB2	2.07	0.55
1:A:154:LEU:N	1:A:168:ILE:HD12	2.22	0.54
1:A:71:VAL:O	1:A:71:VAL:CG1	2.54	0.53
1:A:153:GLU:OE1	1:A:165:ARG:NH2	2.42	0.53
1:A:368:VAL:HG21	1:A:393:GLU:CG	2.30	0.53
1:A:231:ASN:C	1:A:231:ASN:ND2	2.62	0.53
1:A:237:PHE:CD2	1:A:242:LEU:O	2.60	0.52
1:A:237:PHE:CE2	1:A:242:LEU:HD13	2.44	0.52
1:A:199:LEU:O	1:A:200:ASP:C	2.51	0.52
1:A:228:ILE:O	1:A:232:GLU:HB2	2.09	0.52
1:A:149:ALA:HB2	1:A:172:ASP:HA	1.92	0.51
1:A:114:VAL:O	1:A:115:ASN:C	2.54	0.51
1:A:137:ASP:HB3	5:A:2018:HOH:O	2.11	0.51
1:A:71:VAL:HG13	1:A:71:VAL:O	2.11	0.51
1:A:386:ASN:HD21	1:A:388:GLU:HB2	1.76	0.51
1:A:325:ASN:HD21	1:A:327:MET:CG	2.23	0.50
1:A:178:MET:HE3	1:A:180:MET:HG2	1.93	0.50
1:A:13:MET:HB3	1:A:44:LEU:HD22	1.94	0.50
1:A:114:VAL:HG23	1:A:365:GLU:HG2	1.93	0.50
1:A:187:ILE:HG12	1:A:221:PHE:CE2	2.47	0.49
1:A:79:ASN:C	1:A:79:ASN:HD22	2.19	0.49
1:A:154:LEU:HD22	1:A:154:LEU:C	2.37	0.49
1:A:220:ARG:HG3	1:A:224:ILE:HD12	1.95	0.49
1:A:231:ASN:O	1:A:232:GLU:HG3	2.12	0.48
1:A:325:ASN:ND2	1:A:325:ASN:C	2.69	0.48
1:A:35:ILE:HD11	1:A:352:PHE:CD2	2.48	0.48
1:A:247:THR:OG1	1:A:251:ASP:OD2	2.21	0.48
1:A:62:ARG:NH2	5:A:2012:HOH:O	2.47	0.48
1:A:237:PHE:CD2	1:A:242:LEU:HD13	2.47	0.47
1:A:137:ASP:OD1	1:A:217:TYR:OH	2.29	0.47
1:A:246:TYR:C	1:A:246:TYR:CD1	2.92	0.47
1:A:132:ARG:HE	1:A:269:ASN:HD22	1.62	0.47
1:A:118:LYS:HD3	1:A:118:LYS:HA	1.31	0.47
1:A:385:VAL:HB	1:A:389:ASP:HB2	1.97	0.47
1:A:195:PHE:O	1:A:196:ASN:C	2.56	0.47
1:A:141:ASN:OD1	1:A:143:GLU:HG3	2.15	0.47
1:A:79:ASN:HD22	1:A:82:ARG:H	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:HIS:CD2	1:A:190:PHE:HB2	2.51	0.46
1:A:144:TYR:CE2	1:A:145:TYR:O	2.69	0.46
1:A:30:LEU:C	1:A:31:LEU:HD23	2.41	0.45
1:A:208:SER:C	1:A:209:THR:CG2	2.89	0.45
1:A:231:ASN:O	1:A:232:GLU:CG	2.65	0.45
1:A:329:THR:HG23	5:A:2031:HOH:O	2.16	0.45
1:A:37:LEU:HA	1:A:37:LEU:HD12	1.68	0.45
1:A:89:LYS:C	1:A:90:LYS:HG2	2.42	0.45
1:A:203:GLN:HE21	1:A:264:VAL:CG2	2.27	0.45
1:A:178:MET:HG2	1:A:179:GLY:N	2.32	0.45
1:A:230:GLU:HA	1:A:230:GLU:OE1	2.17	0.44
1:A:178:MET:CE	1:A:180:MET:HG2	2.48	0.44
1:A:208:SER:O	1:A:209:THR:HG22	2.16	0.44
1:A:326:SER:O	1:A:327:MET:C	2.57	0.44
1:A:242:LEU:HD12	1:A:242:LEU:C	2.43	0.44
1:A:373:LEU:O	1:A:374:ALA:C	2.62	0.43
1:A:118:LYS:HA	1:A:119:PRO:HD3	1.87	0.42
1:A:386:ASN:ND2	1:A:388:GLU:N	2.68	0.42
1:A:154:LEU:C	1:A:154:LEU:CD2	2.92	0.42
1:A:154:LEU:CD2	1:A:155:VAL:N	2.80	0.42
1:A:339:LYS:HA	1:A:339:LYS:HD3	1.92	0.42
1:A:215:LYS:O	1:A:219:HIS:HB2	2.20	0.42
1:A:198:ALA:O	1:A:202:ASN:N	2.53	0.42
1:A:206:TRP:CD1	1:A:206:TRP:N	2.88	0.42
1:A:215:LYS:O	1:A:219:HIS:CB	2.68	0.41
1:A:79:ASN:O	1:A:80:ALA:C	2.61	0.41
1:A:162:GLU:H	1:A:162:GLU:HG3	1.59	0.41
1:A:184:ASP:O	1:A:185:LYS:C	2.61	0.41
1:A:386:ASN:ND2	1:A:388:GLU:H	2.19	0.41
1:A:66:LYS:HE2	1:A:67:TYR:CE1	2.55	0.41
1:A:9:PRO:HG2	1:A:67:TYR:CD1	2.57	0.40
1:A:77:THR:HA	1:A:78:PRO:HD3	1.85	0.40
1:A:93:LYS:O	1:A:94:SER:C	2.65	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	399/402 (99%)	361 (90%)	32 (8%)	6 (2%)	8	16

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	ASN
1	A	104	ASP
1	A	232	GLU
1	A	217	TYR
1	A	147	PRO
1	A	380	PRO

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	346/347 (100%)	308 (89%)	38 (11%)	6	13

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	THR
1	A	24	ARG
1	A	62	ARG
1	A	79	ASN
1	A	103	LEU

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Mol	Chain	Res	Type
1	A	114	VAL
1	A	140	LYS
1	A	154	LEU
1	A	162	GLU
1	A	163	VAL
1	A	168	ILE
1	A	182	ASN
1	A	189	SER
1	A	203	GLN
1	A	216	THR
1	A	231	ASN
1	A	253	VAL
1	A	258	ARG
1	A	286	LEU
1	A	289	MET
1	A	292	VAL
1	A	297	ASP
1	A	309	VAL
1	A	316	HIS
1	A	317	LEU
1	A	325	ASN
1	A	329	THR
1	A	338	LYS
1	A	350	VAL
1	A	354	THR
1	A	362	GLN
1	A	365	GLU
1	A	368	VAL
1	A	371	LYS
1	A	382	LYS
1	A	388	GLU
1	A	392	LYS
1	A	397	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	87	ASN
1	A	115	ASN
1	A	160	ASN
1	A	169	HIS

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Mol	Chain	Res	Type
1	A	182	ASN
1	A	203	GLN
1	A	219	HIS
1	A	231	ASN
1	A	269	ASN
1	A	306	HIS
1	A	325	ASN
1	A	386	ASN

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	1520	-	50,52,52	2.06	8 (16%)	71,80,80	1.84	17 (23%)
2	ICT	A	1403	3	12,12,12	1.56	2 (16%)	13,16,16	2.17	5 (38%)

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	1520	-	-	8/35/67/67	0/5/5/5
2	ICT	A	1403	3	-	0/16/16/16	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1520	NAP	O7N-C7N	9.61	1.42	1.24
4	A	1520	NAP	C2N-N1N	-6.55	1.27	1.35
4	A	1520	NAP	C6N-N1N	5.40	1.47	1.35
2	A	1403	ICT	C3-C2	-2.93	1.50	1.54
4	A	1520	NAP	C2A-N1A	2.16	1.37	1.33
4	A	1520	NAP	C8A-N7A	2.08	1.35	1.31
2	A	1403	ICT	O2-C1	-2.07	1.24	1.30
4	A	1520	NAP	C2A-N3A	2.07	1.37	1.33
4	A	1520	NAP	PA-O3	2.02	1.61	1.59
4	A	1520	NAP	O4D-C1D	2.00	1.43	1.40

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1520	NAP	N3A-C2A-N1A	-5.76	119.87	128.58
4	A	1520	NAP	C5A-C4A-N3A	-4.38	120.69	126.72
2	A	1403	ICT	O7-C2-C3	-4.34	100.83	110.63
4	A	1520	NAP	C2B-C1B-N9A	-4.15	106.92	113.75
2	A	1403	ICT	O2-C1-O1	-3.91	115.21	124.08
4	A	1520	NAP	C2A-N3A-C4A	3.60	120.62	111.83
4	A	1520	NAP	O2B-P2B-O1X	-3.48	96.92	109.33
4	A	1520	NAP	C5N-C6N-N1N	-3.37	115.79	120.38
4	A	1520	NAP	O7N-C7N-C3N	-3.31	115.55	119.60
4	A	1520	NAP	N9A-C8A-N7A	-3.28	109.29	113.94
4	A	1520	NAP	C5A-N7A-C8A	3.25	108.56	103.45
4	A	1520	NAP	C6N-N1N-C1D	-3.25	113.35	119.73
4	A	1520	NAP	O2A-PA-O3	2.83	114.93	107.27
4	A	1520	NAP	N3A-C4A-N9A	2.77	131.88	127.17
4	A	1520	NAP	O3-PN-O1N	-2.64	102.78	110.70
2	A	1403	ICT	C3-C4-C5	-2.46	109.34	113.95
2	A	1403	ICT	O4-C5-O3	-2.33	117.33	123.33
4	A	1520	NAP	O2X-P2B-O2B	2.21	114.46	105.85
4	A	1520	NAP	C4A-C5A-N7A	-2.19	108.07	110.58
4	A	1520	NAP	C6N-N1N-C2N	2.15	123.70	121.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1403	ICT	O2-C1-C2	2.03	118.95	113.31
4	A	1520	NAP	O7N-C7N-N7N	2.02	125.53	122.62

There are no chirality outliers.

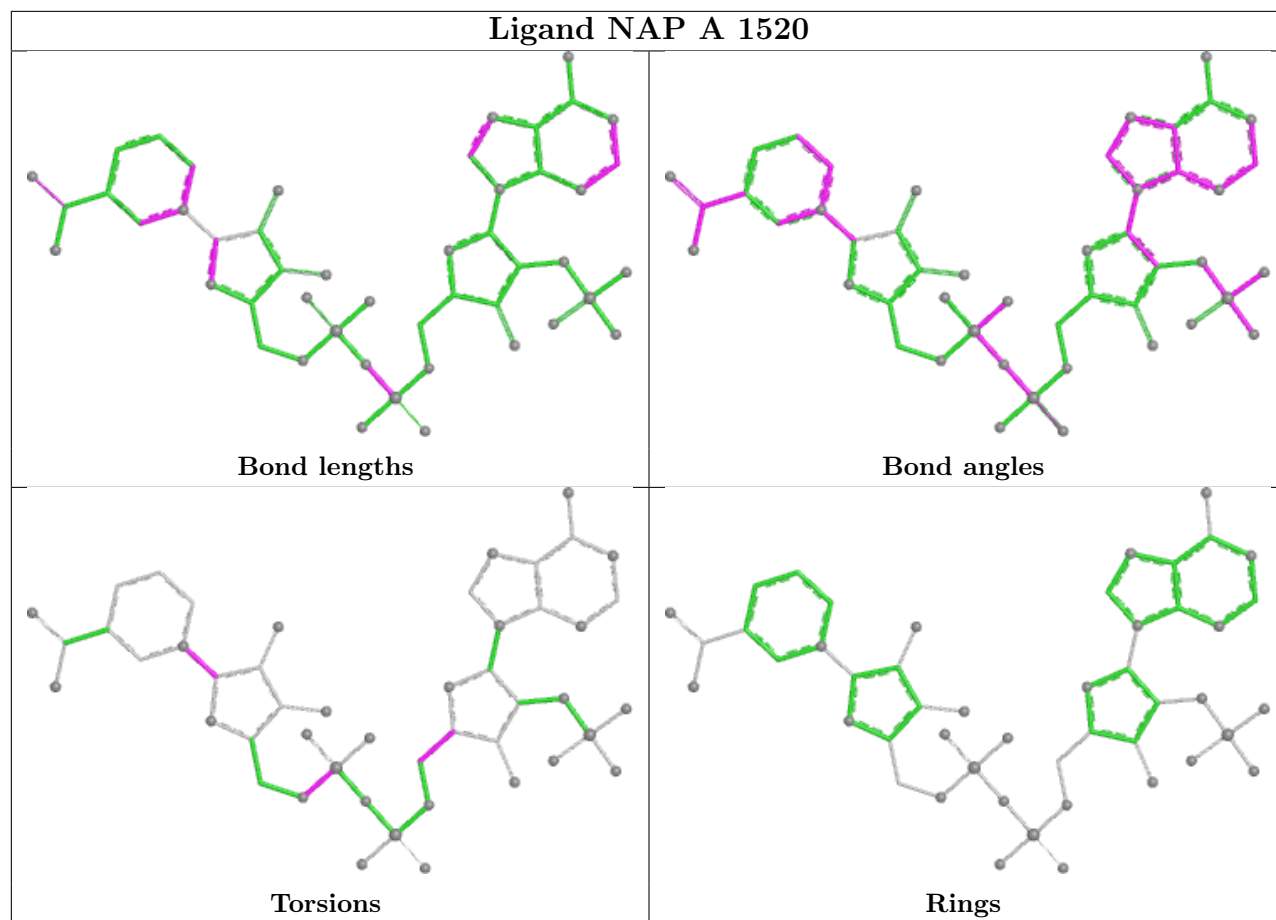
All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1520	NAP	C5D-O5D-PN-O3
4	A	1520	NAP	C5D-O5D-PN-O2N
4	A	1520	NAP	O4D-C1D-N1N-C2N
4	A	1520	NAP	O4D-C1D-N1N-C6N
4	A	1520	NAP	C2D-C1D-N1N-C2N
4	A	1520	NAP	O4B-C4B-C5B-O5B
4	A	1520	NAP	C3B-C4B-C5B-O5B
4	A	1520	NAP	C5D-O5D-PN-O1N

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	401/402 (99%)	-0.22	2 (0%) 87 85	24, 42, 65, 85	13 (3%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	GLU	4.6
1	A	163	VAL	2.4

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.

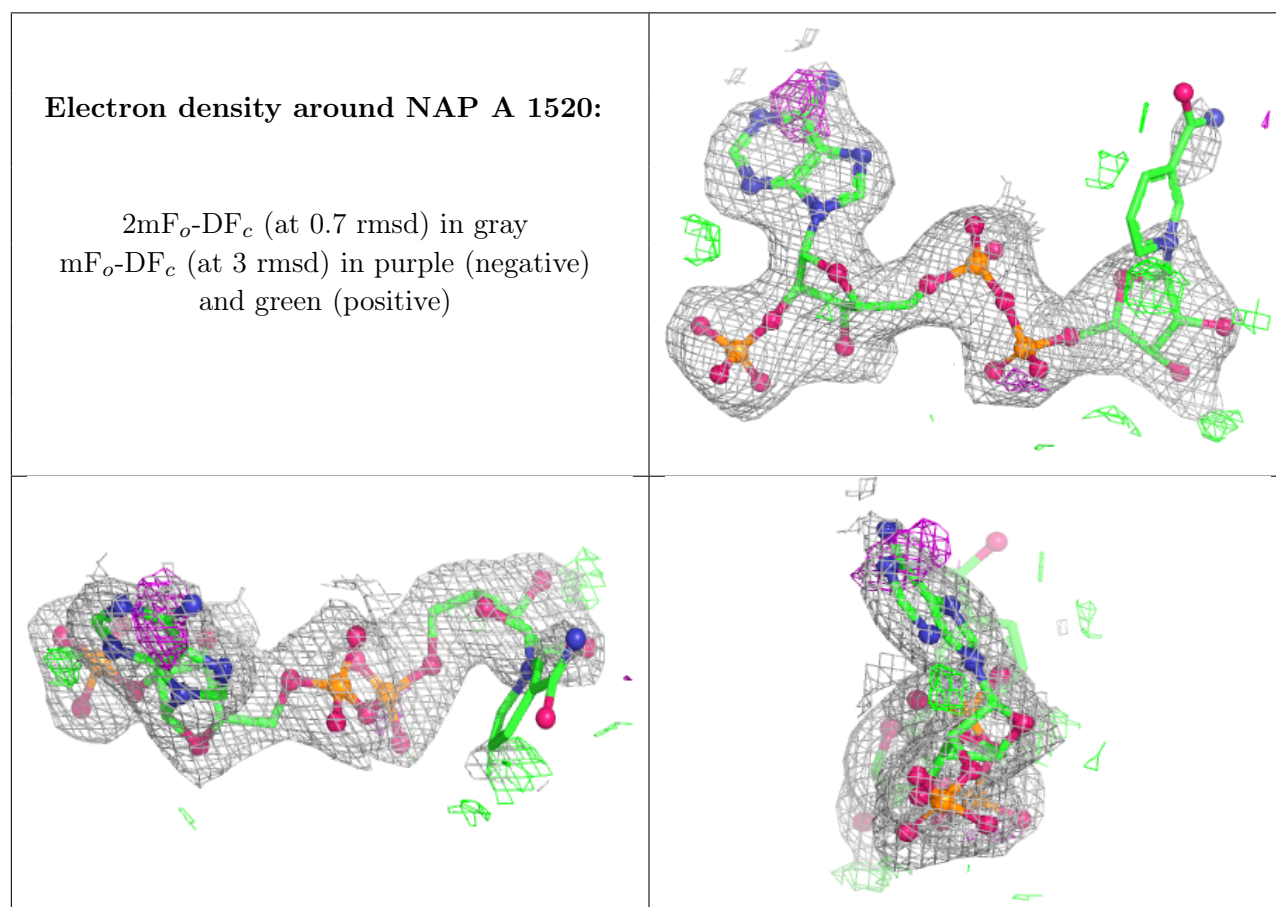
6.4 Ligands [i](#)

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	MG	A	1404	1/1	0.93	0.07	43,43,43,43	0
4	NAP	A	1520	48/48	0.96	0.08	38,46,79,80	8
2	ICT	A	1403	13/13	0.97	0.05	34,37,40,41	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.