



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2026 – 07:19 AM UTC

PDB ID : 1W1U / pdb_00001w1u
Title : Inactive Urocanase-SA cocrystallized with urocanate
Authors : Kessler, D.; Retey, J.; Schulz, G.E.
Deposited on : 2004-06-24
Resolution : 2.23 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

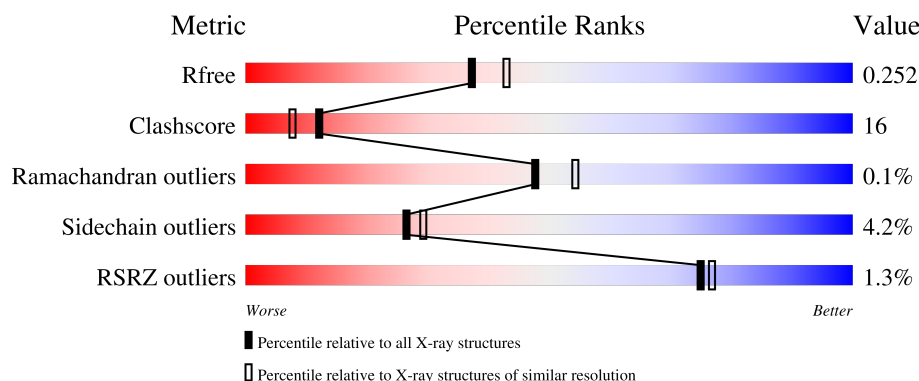
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

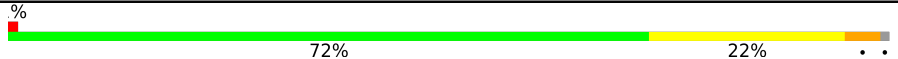
The reported resolution of this entry is 2.23 Å.

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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	557	
1	B	557	

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	URO	A	1556	-	X	X	-
3	URO	B	1556	-	X	X	-

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 9000 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.

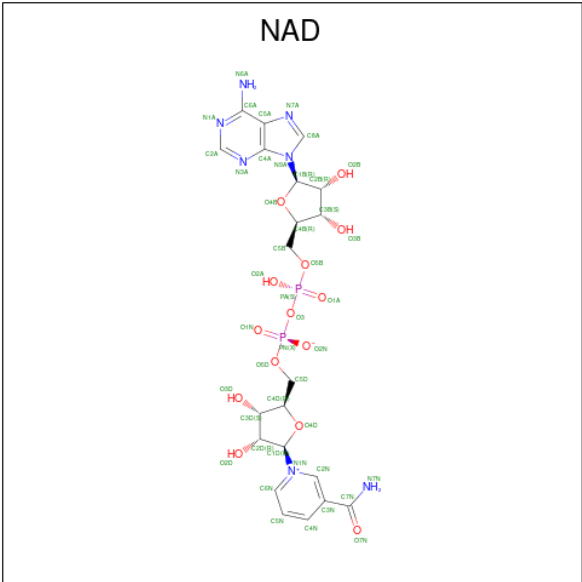
- Molecule 1 is a protein called UROCANATE HYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	1
			4217	2645	753	797	22			
1	B	551	Total	C	N	O	S	0	0	1
			4217	2645	753	797	22			

There are 6 discrepancies between the modelled and reference sequences:

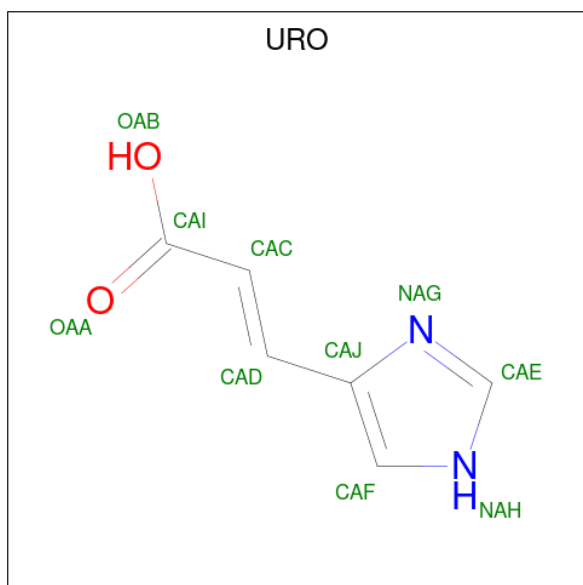
Chain	Residue	Modelled	Actual	Comment	Reference
A	6	ASN	LYS	conflict	UNP P25080
A	198	SER	CYS	engineered mutation	UNP P25080
A	455	ALA	ARG	engineered mutation	UNP P25080
B	6	ASN	LYS	conflict	UNP P25080
B	198	SER	CYS	engineered mutation	UNP P25080
B	455	ALA	ARG	engineered mutation	UNP P25080

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is (2E)-3-(1H-IMIDAZOL-4-YL)ACRYLIC ACID (CCD ID: URO) (formula: $C_6H_6N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	6	2	2		
3	B	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	241	Total	O	0	0
			241	241		
4	B	217	Total	O	0	0
			217	217		

• Molecule 1: UROCANATE HYDRATASE



V507	C508	D509	G510	T511	D512	E513	A514	A515	E516	R517	I518	+	V521	L522	T523	H524	D525	P526	+	A534	+	I540	A541	I542	+	Q548	+	M554	I555	+	THR	GLY
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.43Å 71.62Å 129.16Å 90.00° 98.75° 90.00°	Depositor
Resolution (Å)	19.18 – 2.23 19.18 – 2.23	Depositor EDS
% Data completeness (in resolution range)	98.4 (19.18-2.23) 98.0 (19.18-2.23)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 2.23Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.179 , 0.231 0.206 , 0.252	Depositor DCC
R_{free} test set	2414 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	9.8	Xtriage
Anisotropy	0.410	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9000	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: URO, NAD

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	0.49	1/4310 (0.0%)	1.01	26/5852 (0.4%)
1	B	0.47	1/4310 (0.0%)	1.01	22/5852 (0.4%)
All	All	0.48	2/8620 (0.0%)	1.01	48/11704 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	554	MET	C-N	-5.80	1.25	1.33
1	B	554	MET	C-N	-5.53	1.25	1.33

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	164	SER	N-CA-C	8.87	124.23	113.23
1	B	447	SER	N-CA-C	7.71	120.36	111.11
1	B	343	PRO	N-CA-C	7.52	121.74	111.22
1	B	146	VAL	N-CA-C	7.38	118.11	110.36
1	A	523	THR	N-CA-C	-7.30	102.98	112.68
1	A	247	ILE	N-CA-C	7.06	117.17	110.53
1	A	21	THR	N-CA-C	-6.99	103.64	112.93
1	A	447	SER	N-CA-C	6.64	119.08	111.11
1	A	103	ILE	N-CA-C	6.61	117.36	108.11
1	A	192	CYS	N-CA-C	-6.48	101.20	110.59
1	A	265	THR	N-CA-C	-6.44	102.22	110.53
1	B	443	ASP	N-CA-C	-6.44	102.22	110.53
1	B	21	THR	N-CA-C	-6.43	105.08	113.12
1	A	443	ASP	N-CA-C	-6.42	102.24	110.53
1	B	534	ALA	N-CA-C	-6.26	104.54	111.36
1	A	456	GLU	N-CA-C	6.18	119.67	111.75
1	B	265	THR	N-CA-C	-6.18	102.56	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	460	MET	N-CA-C	-6.14	101.19	110.28
1	A	343	PRO	N-CA-C	6.04	121.77	111.03
1	A	42	VAL	N-CA-C	5.90	116.63	111.56
1	A	137	TRP	N-CA-C	5.90	119.53	111.39
1	B	456	GLU	N-CA-C	5.78	119.15	111.75
1	B	342	PHE	CA-C-N	5.78	125.76	120.21
1	B	342	PHE	C-N-CA	5.78	125.76	120.21
1	B	523	THR	N-CA-C	-5.75	105.03	112.68
1	A	127	ALA	N-CA-C	5.66	117.79	109.24
1	B	399	GLU	N-CA-C	5.59	117.18	111.14
1	A	450	VAL	N-CA-C	5.55	115.68	108.35
1	B	466	ALA	N-CA-C	5.45	119.67	113.19
1	A	280	THR	N-CA-C	-5.39	102.07	110.42
1	B	345	PHE	N-CA-C	5.34	117.85	111.71
1	B	280	THR	N-CA-C	-5.33	101.84	110.32
1	A	483	ALA	N-CA-C	-5.31	103.53	110.53
1	A	105	ASN	N-CA-C	5.29	117.53	108.90
1	B	103	ILE	N-CA-C	5.29	115.52	108.11
1	A	146	VAL	N-CA-C	5.23	116.58	110.62
1	A	131	GLN	CB-CA-C	-5.21	110.54	116.54
1	A	444	HIS	N-CA-C	-5.21	106.88	113.18
1	B	213	ASP	N-CA-C	5.19	117.73	111.40
1	A	213	ASP	N-CA-C	5.17	117.70	111.40
1	A	344	GLY	N-CA-C	-5.14	105.72	112.82
1	B	45	ASN	CA-C-N	5.12	124.91	119.28
1	B	45	ASN	C-N-CA	5.12	124.91	119.28
1	A	530	VAL	N-CA-C	-5.11	105.51	110.42
1	B	247	ILE	N-CA-C	5.10	115.31	110.42
1	B	131	GLN	CB-CA-C	-5.08	110.69	116.54
1	A	34	LEU	N-CA-C	-5.04	105.68	111.07
1	A	324	GLY	N-CA-C	5.02	123.29	114.76

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	4217	0	4116	133	1
1	B	4217	0	4116	143	0
2	A	44	0	26	4	0
2	B	44	0	26	7	0
3	A	10	0	5	9	0
3	B	10	0	5	8	0
4	A	241	0	0	7	0
4	B	217	0	0	7	0
All	All	9000	0	8294	262	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 (262) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:ILE:HD13	1:B:540:ILE:H	1.22	1.01
1:B:346:VAL:HB	4:B:2125:HOH:O	1.61	0.99
1:A:132:MET:SD	3:A:1556:URO:NAH	2.38	0.96
1:B:346:VAL:HG11	1:B:456:GLU:HA	1.53	0.90
1:B:132:MET:SD	3:B:1556:URO:NAH	2.48	0.87
1:A:196:ILE:HD13	2:A:1555:NAD:H2A	1.58	0.85
1:B:253:LYS:NZ	1:B:253:LYS:HB3	1.92	0.83
1:B:346:VAL:HG13	1:B:347:PRO:HD3	1.61	0.83
1:B:514:ALA:O	1:B:518:ILE:HG12	1.79	0.82
1:A:554:MET:HE2	1:B:64:CYS:SG	2.22	0.80
1:A:392:ARG:CD	1:A:396:MET:HE3	2.13	0.79
1:B:517:ARG:O	1:B:521:VAL:HG13	1.83	0.79
1:B:307:VAL:HA	1:B:310:MET:HE3	1.64	0.78
1:B:55:ILE:HD11	1:B:57:ARG:CZ	2.13	0.78
1:A:397:ALA:HA	1:A:401:ILE:HD13	1.66	0.78
1:A:392:ARG:HH21	1:A:396:MET:HE1	1.51	0.74
1:A:13:ARG:NH1	1:A:13:ARG:HB3	2.01	0.74
1:A:415:LEU:HD12	1:A:530:VAL:HG21	1.67	0.74
1:A:178:MET:SD	3:A:1556:URO:HAD	2.29	0.73
1:B:540:ILE:H	1:B:540:ILE:CD1	1.96	0.73
1:B:511:THR:HG22	1:B:513:GLU:H	1.54	0.72
1:A:6:ASN:O	1:A:48:GLU:HA	1.87	0.72
1:A:13:ARG:HB3	1:A:13:ARG:HH11	1.54	0.71
1:B:534:ALA:HB1	1:B:542:ILE:HD13	1.73	0.71
1:B:253:LYS:HB3	1:B:253:LYS:HZ3	1.56	0.70
1:A:440:ILE:HD13	1:A:508:CYS:SG	2.32	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:PHE:HB3	1:A:318:VAL:HG21	1.74	0.70
1:A:196:ILE:HG13	1:A:247:ILE:HD12	1.74	0.70
1:A:196:ILE:CD1	2:A:1555:NAD:H2A	2.22	0.70
1:A:64:CYS:SG	1:B:554:MET:HE2	2.32	0.69
1:A:163:GLY:O	4:A:2081:HOH:O	2.11	0.69
1:A:291:GLU:HB3	4:A:2047:HOH:O	1.92	0.69
1:A:392:ARG:HD2	1:A:396:MET:HE3	1.73	0.69
1:B:415:LEU:HD13	1:B:548:GLN:HG3	1.75	0.68
1:A:392:ARG:HD3	1:A:396:MET:HE3	1.74	0.68
1:A:52:TYR:CE2	3:A:1556:URO:NAG	2.63	0.67
1:B:214:GLU:CD	1:B:236:ILE:HD11	2.18	0.67
1:B:443:ASP:OD2	3:B:1556:URO:HAF	1.94	0.67
1:A:415:LEU:HD13	1:A:548:GLN:HG3	1.77	0.67
1:A:132:MET:SD	3:A:1556:URO:CAF	2.82	0.67
1:B:511:THR:HG22	1:B:513:GLU:N	2.10	0.67
1:A:392:ARG:HH21	1:A:396:MET:CE	2.08	0.66
1:B:55:ILE:HD11	1:B:57:ARG:NH2	2.10	0.66
1:A:93:LYS:NZ	1:A:93:LYS:H	1.92	0.66
1:B:132:MET:SD	3:B:1556:URO:CAF	2.84	0.66
1:B:440:ILE:HD13	1:B:508:CYS:SG	2.35	0.66
1:B:50:VAL:HG22	1:B:55:ILE:HD12	1.77	0.66
1:B:248:LEU:HB2	1:B:249:PRO:HD3	1.79	0.65
1:B:346:VAL:HG12	4:B:2124:HOH:O	1.96	0.65
1:B:346:VAL:HG13	1:B:347:PRO:CD	2.27	0.65
1:B:293:ALA:O	1:B:297:LYS:HG3	1.96	0.64
1:B:196:ILE:CD1	2:B:1555:NAD:H2A	2.28	0.64
1:A:306:HIS:CD2	1:A:310:MET:HE2	2.33	0.63
1:B:178:MET:SD	3:B:1556:URO:HAD	2.38	0.63
1:B:68:ILE:HG22	1:B:88:PRO:HB3	1.81	0.63
1:A:307:VAL:HA	1:A:310:MET:HE3	1.81	0.63
1:A:443:ASP:OD2	3:A:1556:URO:HAF	1.98	0.63
1:B:147:GLN:HB2	1:B:393:TRP:CZ2	2.34	0.63
1:B:184:LEU:HD21	1:B:401:ILE:HD13	1.81	0.62
1:B:307:VAL:HA	1:B:310:MET:CE	2.29	0.62
1:B:511:THR:HG21	4:B:2190:HOH:O	1.98	0.62
1:B:511:THR:CG2	1:B:513:GLU:OE2	2.47	0.62
1:A:80:THR:HG21	1:A:102:LEU:HD22	1.81	0.62
1:A:272:ASN:HB3	1:A:285:ARG:HH22	1.65	0.62
1:B:214:GLU:CG	1:B:236:ILE:HD11	2.29	0.62
1:A:551:ASP:O	1:A:553:PRO:HD3	1.99	0.62
1:A:20:LEU:HD11	1:A:25:TRP:CE2	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:THR:OG1	1:B:283:GLN:HG3	2.01	0.61
1:A:170:VAL:HG22	1:A:260:MET:HG2	1.81	0.61
1:B:96:SER:HA	1:B:100:ARG:NH2	2.15	0.61
1:B:65:TYR:O	1:B:68:ILE:HG12	2.00	0.60
1:A:447:SER:OG	1:A:529:GLY:HA3	2.01	0.60
1:A:291:GLU:HB2	1:A:294:ALA:HB3	1.83	0.59
1:B:555:ILE:N	4:B:2213:HOH:O	2.36	0.59
1:A:64:CYS:SG	1:B:554:MET:CE	2.90	0.59
1:A:517:ARG:O	1:A:521:VAL:HG13	2.02	0.59
1:A:329:GLN:O	1:A:333:GLU:HG3	2.02	0.58
1:B:20:LEU:HD21	1:B:25:TRP:CE2	2.38	0.58
1:A:52:TYR:HE2	3:A:1556:URO:NAG	2.00	0.58
1:A:170:VAL:HG22	1:A:260:MET:CG	2.33	0.58
1:A:131:GLN:OE1	3:A:1556:URO:HAE	2.03	0.58
1:A:367:SER:HB3	1:A:436:ALA:HB3	1.84	0.58
1:A:102:LEU:HG	1:B:476:LEU:HB3	1.85	0.57
1:A:257:ARG:HD2	1:A:318:VAL:CG1	2.34	0.57
1:B:392:ARG:O	1:B:396:MET:HG3	2.04	0.57
1:B:397:ALA:HA	1:B:401:ILE:HG12	1.86	0.57
1:A:313:PHE:HB3	1:A:318:VAL:CG2	2.34	0.57
1:B:133:THR:HA	1:B:136:SER:OG	2.05	0.57
1:B:178:MET:HE2	1:B:178:MET:HA	1.88	0.56
1:B:229:TYR:CE2	1:B:236:ILE:HD13	2.39	0.56
1:B:196:ILE:HD13	2:B:1555:NAD:H2A	1.85	0.56
1:A:287:ARG:NH1	1:A:291:GLU:OE2	2.39	0.56
1:A:489:HIS:HB2	1:A:501:HIS:CE1	2.40	0.55
1:A:109:VAL:HA	1:B:500:GLN:HE21	1.72	0.55
1:A:383:LEU:HD22	1:A:420:LYS:HE2	1.89	0.55
1:B:196:ILE:HD12	1:B:244:ALA:HB2	1.88	0.55
1:A:196:ILE:HD13	1:A:196:ILE:C	2.31	0.55
1:A:297:LYS:O	1:A:301:GLN:HG3	2.07	0.55
1:B:55:ILE:HG13	1:B:55:ILE:O	2.07	0.54
1:A:257:ARG:HG3	1:A:257:ARG:HH11	1.73	0.54
1:A:384:ILE:CD1	1:A:417:LEU:HD22	2.36	0.54
1:A:453:PRO:HB2	1:B:112:TRP:CH2	2.42	0.54
1:B:196:ILE:HD11	1:B:244:ALA:N	2.22	0.54
1:B:52:TYR:CE2	3:B:1556:URO:NAG	2.76	0.54
1:A:12:ILE:O	1:A:39:ASP:HA	2.07	0.54
1:A:338:ASN:HB2	1:A:341:ASP:OD1	2.08	0.54
1:B:94:THR:HG22	1:B:95:HIS:N	2.23	0.54
1:A:475:ALA:HA	4:A:2204:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:GLU:HG2	1:B:236:ILE:HD11	1.90	0.53
1:A:164:SER:HA	4:A:2082:HOH:O	2.07	0.53
1:A:254:ARG:HD3	4:A:2110:HOH:O	2.08	0.53
1:B:196:ILE:HD13	1:B:196:ILE:C	2.34	0.52
1:B:489:HIS:HB2	1:B:501:HIS:CE1	2.44	0.52
1:A:497:GLY:HA2	1:B:110:PRO:HB2	1.91	0.52
1:B:52:TYR:HE2	3:B:1556:URO:NAG	2.07	0.52
1:A:13:ARG:HH11	1:A:13:ARG:CB	2.19	0.52
1:A:383:LEU:O	1:A:384:ILE:HD13	2.09	0.52
1:A:112:TRP:CH2	1:B:453:PRO:HB2	2.44	0.51
1:A:234:LYS:HG3	4:A:2103:HOH:O	2.11	0.51
1:A:440:ILE:CD1	1:A:508:CYS:SG	2.98	0.51
1:B:94:THR:HG22	1:B:95:HIS:H	1.75	0.51
1:A:9:ASP:OD2	1:A:47:LYS:HE3	2.10	0.51
1:B:50:VAL:HG22	1:B:55:ILE:CD1	2.40	0.51
1:B:196:ILE:HD11	2:B:1555:NAD:H2A	1.93	0.51
1:A:37:ASN:O	1:A:43:ALA:HB2	2.11	0.51
1:B:253:LYS:HB3	1:B:253:LYS:HZ2	1.73	0.51
1:A:173:ALA:HB2	1:A:196:ILE:HG23	1.92	0.51
1:B:291:GLU:HB2	1:B:294:ALA:HB3	1.93	0.50
1:A:257:ARG:HH11	1:A:257:ARG:CG	2.24	0.50
1:B:338:ASN:HB2	1:B:341:ASP:OD1	2.11	0.50
1:B:324:GLY:HA3	2:B:1555:NAD:H71N	1.76	0.50
1:A:461:ARG:HH21	1:A:461:ARG:HG3	1.77	0.50
1:A:554:MET:CE	1:B:64:CYS:SG	2.97	0.50
1:B:12:ILE:O	1:B:39:ASP:HA	2.11	0.50
1:A:324:GLY:HA3	2:A:1555:NAD:H71N	1.77	0.50
1:A:290:THR:HG22	1:A:291:GLU:HG2	1.92	0.49
1:A:312:ASP:O	1:A:316:GLN:HG2	2.12	0.49
1:A:525:ASP:HB3	1:A:526:PRO:CD	2.42	0.49
1:B:12:ILE:HD12	4:B:2026:HOH:O	2.11	0.49
1:A:307:VAL:HG22	1:A:310:MET:HE3	1.94	0.49
1:B:45:ASN:HB3	1:B:50:VAL:HB	1.94	0.49
1:A:93:LYS:H	1:A:93:LYS:HZ2	1.61	0.49
1:A:313:PHE:C	1:A:318:VAL:HG22	2.37	0.49
1:A:504:MET:SD	1:A:504:MET:C	2.96	0.49
1:B:266:SER:HB2	1:B:273:GLY:HA3	1.94	0.49
1:A:427:GLU:HG2	1:A:430:ARG:NH2	2.29	0.48
1:B:138:ILE:HG13	1:B:138:ILE:O	2.13	0.48
1:A:19:LYS:HD3	1:A:19:LYS:C	2.39	0.48
1:A:480:ALA:HB1	1:B:94:THR:HG21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:THR:HG21	1:B:102:LEU:HD22	1.95	0.48
1:B:307:VAL:HG21	1:B:331:ALA:HB2	1.96	0.48
1:B:263:ASP:OD2	1:B:265:THR:HG23	2.14	0.48
1:B:511:THR:HG23	1:B:513:GLU:OE2	2.13	0.48
1:A:93:LYS:H	1:A:93:LYS:CE	2.26	0.48
1:B:214:GLU:OE1	1:B:225:ARG:HD3	2.14	0.48
1:A:443:ASP:OD1	1:A:444:HIS:N	2.46	0.48
1:A:20:LEU:HD11	1:A:25:TRP:NE1	2.29	0.47
1:A:214:GLU:OE1	1:A:225:ARG:HD3	2.13	0.47
1:A:93:LYS:N	1:A:93:LYS:HD3	2.29	0.47
1:B:362:ARG:HH12	3:B:1556:URO:CAI	2.27	0.47
1:B:94:THR:HG21	1:B:98:ALA:CB	2.45	0.47
1:B:506:ILE:CD1	1:B:518:ILE:HD13	2.45	0.47
1:A:288:ALA:HA	1:A:295:VAL:HG21	1.96	0.47
1:B:513:GLU:O	1:B:516:GLU:HG3	2.15	0.47
1:B:297:LYS:O	1:B:301:GLN:HG3	2.15	0.47
1:B:439:VAL:HG12	1:B:507:VAL:HG22	1.97	0.47
1:A:130:GLY:O	1:A:131:GLN:C	2.57	0.46
1:A:554:MET:SD	1:B:67:LYS:HG2	2.55	0.46
1:A:196:ILE:HD11	1:A:242:GLY:O	2.15	0.46
1:B:220:ASP:O	1:B:224:VAL:HG23	2.15	0.46
1:B:204:ASP:O	1:B:208:GLU:HG3	2.15	0.46
1:B:230:THR:HG23	1:B:231:ALA:N	2.30	0.46
1:A:64:CYS:O	1:A:68:ILE:HG13	2.16	0.45
1:B:227:ALA:O	1:B:230:THR:HG22	2.17	0.45
1:B:376:THR:O	1:B:380:VAL:HG23	2.15	0.45
1:A:109:VAL:HG21	1:B:453:PRO:HD3	1.98	0.45
1:A:249:PRO:O	1:A:253:LYS:HG3	2.16	0.45
1:B:439:VAL:C	1:B:440:ILE:HD12	2.42	0.45
1:A:313:PHE:O	1:A:318:VAL:HG22	2.16	0.45
1:A:68:ILE:O	1:A:72:LEU:HB2	2.15	0.45
1:B:387:ASP:CG	1:B:390:LEU:HB2	2.42	0.45
1:A:525:ASP:HB3	1:A:526:PRO:HD3	1.99	0.45
2:A:1555:NAD:C4N	3:A:1556:URO:CAF	2.94	0.45
1:B:540:ILE:HD13	1:B:540:ILE:N	2.07	0.45
1:A:257:ARG:HD2	1:A:318:VAL:HG12	1.98	0.45
1:A:93:LYS:N	1:A:93:LYS:CD	2.80	0.45
1:A:290:THR:HG22	1:A:291:GLU:N	2.30	0.45
1:B:362:ARG:HD3	1:B:442:ARG:HA	1.98	0.45
1:B:525:ASP:HB3	1:B:526:PRO:CD	2.47	0.45
1:B:397:ALA:HA	1:B:401:ILE:CG1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:MET:HE3	4:A:2069:HOH:O	2.17	0.44
1:B:144:GLY:HA3	4:B:2217:HOH:O	2.16	0.44
1:B:59:ALA:C	1:B:128:MET:SD	3.01	0.44
1:A:16:ARG:HD2	1:A:368:GLY:HA3	2.00	0.44
1:A:111:HIS:HB3	1:A:112:TRP:CE3	2.53	0.44
1:B:85:SER:HA	1:B:105:ASN:OD1	2.17	0.44
1:B:15:PRO:HD3	1:B:35:MET:SD	2.57	0.44
1:B:257:ARG:HG2	1:B:257:ARG:HH11	1.82	0.44
1:A:470:TRP:CZ3	1:A:531:MET:HE3	2.52	0.44
1:A:445:LEU:C	1:A:445:LEU:HD23	2.43	0.43
1:A:390:LEU:C	1:A:390:LEU:HD23	2.43	0.43
1:B:131:GLN:HB3	1:B:132:MET:H	1.65	0.43
1:B:169:TRP:CE2	1:B:258:PRO:HA	2.54	0.43
1:A:118:PHE:CD1	1:A:118:PHE:C	2.97	0.43
1:B:443:ASP:OD1	1:B:444:HIS:N	2.50	0.43
1:A:246:GLU:OE1	1:A:278:GLY:N	2.52	0.43
1:A:138:ILE:HG13	1:A:138:ILE:O	2.18	0.43
1:A:400:ARG:C	1:A:401:ILE:HD12	2.44	0.43
1:B:230:THR:CG2	1:B:231:ALA:N	2.82	0.43
1:B:251:LEU:HD22	1:B:256:VAL:HG11	2.01	0.43
1:B:540:ILE:CD1	1:B:540:ILE:N	2.75	0.43
1:B:506:ILE:HD13	1:B:518:ILE:HD13	2.01	0.43
1:A:105:ASN:HB2	1:A:489:HIS:CD2	2.54	0.42
1:A:553:PRO:HB2	1:B:88:PRO:O	2.19	0.42
1:B:103:ILE:O	1:B:487:SER:HA	2.18	0.42
1:B:445:LEU:HD23	1:B:445:LEU:C	2.44	0.42
2:B:1555:NAD:C4N	3:B:1556:URO:CAF	2.97	0.42
1:A:19:LYS:HD3	1:A:20:LEU:O	2.19	0.42
1:A:285:ARG:HA	1:A:285:ARG:HD3	1.69	0.42
1:B:9:ASP:OD2	1:B:47:LYS:HE3	2.19	0.42
1:B:184:LEU:CD2	1:B:401:ILE:HD13	2.49	0.42
1:A:150:TYR:O	1:A:154:VAL:HG23	2.19	0.42
1:B:67:LYS:HG3	1:B:88:PRO:HG3	2.01	0.42
1:B:362:ARG:HD2	1:B:410:ILE:HD13	2.02	0.42
1:A:110:PRO:HB2	1:B:497:GLY:HA2	2.00	0.42
1:A:453:PRO:HD3	1:B:109:VAL:HG21	2.01	0.42
1:B:445:LEU:HD12	2:B:1555:NAD:O7N	2.20	0.42
1:A:361:PHE:CD1	1:A:418:ARG:HD3	2.54	0.42
1:B:281:TRP:HE1	1:B:285:ARG:HD2	1.85	0.42
1:B:475:ALA:HA	4:B:2184:HOH:O	2.19	0.42
1:A:310:MET:HE1	1:A:327:ILE:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:ARG:HG3	1:A:461:ARG:NH2	2.34	0.42
1:B:509:ASP:OD1	1:B:511:THR:HB	2.20	0.41
1:A:52:TYR:CZ	3:A:1556:URO:NAG	2.88	0.41
1:A:109:VAL:HA	1:B:500:GLN:NE2	2.33	0.41
1:B:266:SER:CB	1:B:273:GLY:HA3	2.49	0.41
1:A:93:LYS:H	1:A:93:LYS:CD	2.33	0.41
1:B:5:ASN:O	1:B:61:ASN:HB2	2.19	0.41
1:B:324:GLY:O	2:B:1555:NAD:H2N	2.20	0.41
1:A:234:LYS:HD3	1:A:234:LYS:HA	1.81	0.41
1:B:196:ILE:CD1	1:B:244:ALA:N	2.83	0.41
1:B:346:VAL:HG11	1:B:456:GLU:CA	2.38	0.41
1:B:130:GLY:O	1:B:131:GLN:C	2.61	0.41
1:B:158:ARG:HA	1:B:163:GLY:HA2	2.03	0.41
1:A:131:GLN:HB3	1:A:132:MET:H	1.60	0.41
1:A:220:ASP:O	1:A:224:VAL:HG23	2.21	0.41
1:A:383:LEU:C	1:A:384:ILE:HD13	2.45	0.41
1:B:221:ASP:O	1:B:225:ARG:HG3	2.21	0.41
1:B:245:ALA:O	1:B:249:PRO:HG2	2.20	0.41
1:B:111:HIS:HB3	1:B:112:TRP:CE3	2.56	0.41
1:B:445:LEU:HD23	1:B:445:LEU:O	2.21	0.41
1:A:184:LEU:HD22	1:A:188:LEU:HG	2.03	0.40
1:A:276:PRO:HB2	1:A:279:TRP:CD1	2.56	0.40
1:A:384:ILE:HD12	1:A:417:LEU:HD22	2.04	0.40
1:B:68:ILE:HG23	1:B:88:PRO:HD3	2.02	0.40
1:B:316:GLN:NE2	1:B:316:GLN:HA	2.36	0.40
1:A:485:TRP:CH2	1:A:505:VAL:HG11	2.57	0.40
1:A:93:LYS:HB2	1:A:93:LYS:HE2	1.70	0.40
1:B:425:PHE:O	1:B:429:VAL:HG23	2.21	0.40
1:A:14:ALA:O	1:A:16:ARG:NH1	2.52	0.40
1:B:67:LYS:HA	1:B:67:LYS:HE3	2.04	0.40
1:B:362:ARG:HB3	1:B:410:ILE:HD12	2.04	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:280:THR:OG1	1:A:399:GLU:O[4_544]	2.06	0.14

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	549/557 (99%)	522 (95%)	27 (5%)	0	100	100
1	B	549/557 (99%)	520 (95%)	28 (5%)	1 (0%)	43	48
All	All	1098/1114 (99%)	1042 (95%)	55 (5%)	1 (0%)	48	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	138	ILE

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	432/438 (99%)	413 (96%)	19 (4%)	25	27
1	B	432/438 (99%)	415 (96%)	17 (4%)	28	33
All	All	864/876 (99%)	828 (96%)	36 (4%)	26	29

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

Mol	Chain	Res	Type
1	A	6	ASN
1	A	13	ARG
1	A	72	LEU
1	A	93	LYS
1	A	102	LEU

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Mol	Chain	Res	Type
1	A	126	LEU
1	A	184	LEU
1	A	196	ILE
1	A	206	ARG
1	A	248	LEU
1	A	257	ARG
1	A	318	VAL
1	A	343	PRO
1	A	400	ARG
1	A	401	ILE
1	A	415	LEU
1	A	476	LEU
1	A	521	VAL
1	A	553	PRO
1	B	19	LYS
1	B	72	LEU
1	B	100	ARG
1	B	102	LEU
1	B	110	PRO
1	B	126	LEU
1	B	196	ILE
1	B	206	ARG
1	B	236	ILE
1	B	248	LEU
1	B	291	GLU
1	B	346	VAL
1	B	362	ARG
1	B	415	LEU
1	B	476	LEU
1	B	521	VAL
1	B	540	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	119	ASN
1	A	160	HIS
1	A	199	GLN
1	A	241	HIS
1	A	548	GLN
1	B	18	ASN

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Mol	Chain	Res	Type
1	B	199	GLN
1	B	264	GLN
1	B	272	ASN
1	B	301	GLN
1	B	314	GLN
1	B	316	GLN
1	B	500	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	NAD	B	1555	-	46,48,48	1.58	9 (19%)	64,73,73	1.58	12 (18%)
3	URO	B	1556	-	10,10,10	2.37	6 (60%)	11,12,12	1.96	3 (27%)
2	NAD	A	1555	-	46,48,48	1.57	9 (19%)	64,73,73	1.57	12 (18%)
3	URO	A	1556	-	10,10,10	2.36	6 (60%)	11,12,12	1.96	2 (18%)

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	NAD	B	1555	-	-	2/30/62/62	0/5/5/5
3	URO	B	1556	-	-	4/5/5/5	0/1/1/1
2	NAD	A	1555	-	-	2/30/62/62	0/5/5/5
3	URO	A	1556	-	-	4/5/5/5	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1555	NAD	C2N-N1N	5.14	1.40	1.35
2	B	1555	NAD	C2N-N1N	5.11	1.40	1.35
3	B	1556	URO	CAF-CAJ	-4.56	1.33	1.37
3	A	1556	URO	CAF-CAJ	-4.53	1.33	1.37
2	B	1555	NAD	C2A-N1A	3.79	1.40	1.33
2	A	1555	NAD	C2A-N1A	3.76	1.40	1.33
3	A	1556	URO	CAJ-NAG	-3.23	1.33	1.39
3	B	1556	URO	CAJ-NAG	-3.21	1.33	1.39
2	B	1555	NAD	C5A-N7A	-3.09	1.33	1.39
2	B	1555	NAD	PN-O1N	3.08	1.61	1.50
2	A	1555	NAD	C5A-N7A	-3.08	1.33	1.39
2	B	1555	NAD	PA-O1A	3.08	1.61	1.50
2	A	1555	NAD	PN-O1N	3.05	1.61	1.50
2	A	1555	NAD	PA-O1A	3.03	1.61	1.50
3	A	1556	URO	OAB-CAI	-2.72	1.23	1.30
3	B	1556	URO	OAB-CAI	-2.71	1.23	1.30
3	B	1556	URO	CAD-CAJ	2.68	1.53	1.45
3	A	1556	URO	CAD-CAJ	2.63	1.53	1.45
2	A	1555	NAD	O4D-C1D	2.52	1.44	1.40
2	B	1555	NAD	O4D-C1D	2.49	1.44	1.40
2	B	1555	NAD	C6N-N1N	2.35	1.40	1.35
2	A	1555	NAD	C6N-N1N	2.31	1.40	1.35
2	B	1555	NAD	C8A-N9A	-2.28	1.33	1.37
2	A	1555	NAD	C8A-N9A	-2.28	1.33	1.37
3	B	1556	URO	CAD-CAC	2.15	1.39	1.33
3	A	1556	URO	CAD-CAC	2.12	1.39	1.33
2	A	1555	NAD	C4A-N9A	-2.10	1.33	1.37
2	B	1555	NAD	C4A-N9A	-2.08	1.33	1.37
3	B	1556	URO	CAC-CAI	2.07	1.53	1.48
3	A	1556	URO	CAC-CAI	2.04	1.53	1.48

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1555	NAD	N3A-C2A-N1A	-5.47	120.30	128.58
2	B	1555	NAD	N3A-C2A-N1A	-5.44	120.35	128.58
2	B	1555	NAD	C5A-C4A-N3A	-4.29	120.81	126.72
2	B	1555	NAD	C2A-N3A-C4A	4.11	121.86	111.83
3	A	1556	URO	CAC-CAD-CAJ	-4.03	111.94	129.21
2	A	1555	NAD	C5A-C4A-N3A	-3.95	121.27	126.72
2	A	1555	NAD	C2A-N3A-C4A	3.93	121.44	111.83
3	B	1556	URO	CAC-CAD-CAJ	-3.87	112.63	129.21
2	A	1555	NAD	C4D-O4D-C1D	-3.28	106.92	109.92
3	A	1556	URO	CAD-CAC-CAI	-3.12	111.17	127.05
2	B	1555	NAD	N9A-C8A-N7A	-3.11	109.53	113.94
3	B	1556	URO	CAD-CAC-CAI	-3.09	111.34	127.05
2	A	1555	NAD	N9A-C8A-N7A	-3.02	109.66	113.94
2	B	1555	NAD	N3A-C4A-N9A	2.98	132.24	127.17
2	B	1555	NAD	C5A-N7A-C8A	2.83	107.90	103.45
2	A	1555	NAD	N3A-C4A-N9A	2.80	131.92	127.17
2	B	1555	NAD	C4A-C5A-N7A	-2.76	107.43	110.58
2	A	1555	NAD	C5A-N7A-C8A	2.72	107.73	103.45
2	B	1555	NAD	C4D-O4D-C1D	-2.68	107.47	109.92
2	A	1555	NAD	C4A-C5A-N7A	-2.63	107.58	110.58
2	B	1555	NAD	O2N-PN-O3	2.47	113.95	107.27
2	A	1555	NAD	O2N-PN-O3	2.46	113.91	107.27
2	A	1555	NAD	C4A-N9A-C8A	2.38	108.23	105.74
2	B	1555	NAD	C4A-N9A-C8A	2.34	108.19	105.74
2	B	1555	NAD	O4B-C1B-C2B	-2.29	101.72	106.62
2	A	1555	NAD	C4B-O4B-C1B	-2.27	104.44	109.47
2	A	1555	NAD	C3N-C7N-N7N	-2.13	115.12	117.74
2	B	1555	NAD	C3N-C7N-N7N	-2.10	115.16	117.74
3	B	1556	URO	CAE-NAG-CAJ	2.01	108.32	104.97

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1555	NAD	O4D-C1D-N1N-C6N
2	B	1555	NAD	O4D-C1D-N1N-C6N
3	A	1556	URO	CAC-CAD-CAJ-CAF
3	A	1556	URO	CAC-CAD-CAJ-NAG
3	B	1556	URO	CAC-CAD-CAJ-CAF
3	B	1556	URO	CAC-CAD-CAJ-NAG
3	B	1556	URO	CAD-CAC-CAI-OAA

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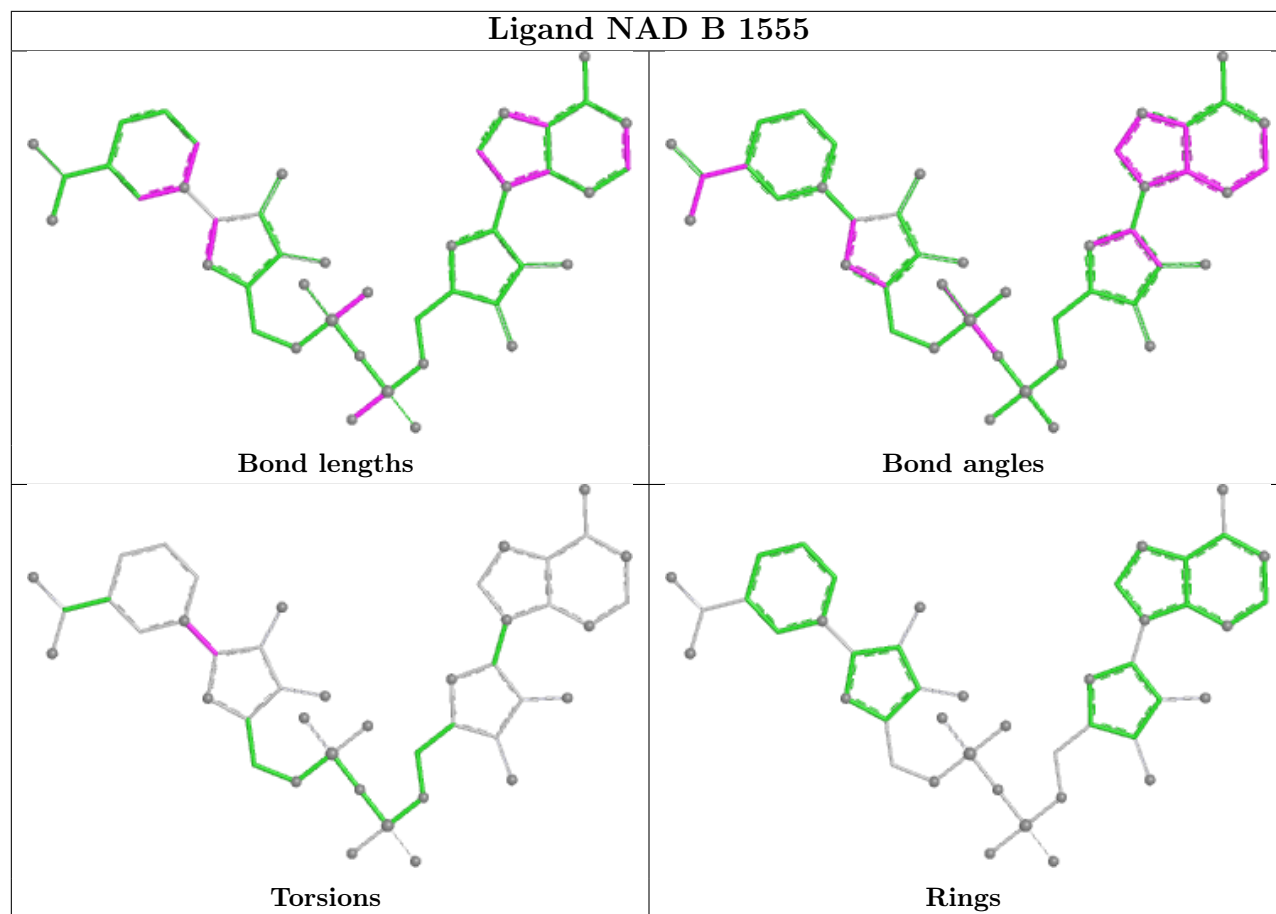
Mol	Chain	Res	Type	Atoms
3	B	1556	URO	CAD-CAC-CAI-OAB
3	A	1556	URO	CAD-CAC-CAI-OAB
3	A	1556	URO	CAD-CAC-CAI-OAA
2	A	1555	NAD	O4D-C1D-N1N-C2N
2	B	1555	NAD	O4D-C1D-N1N-C2N

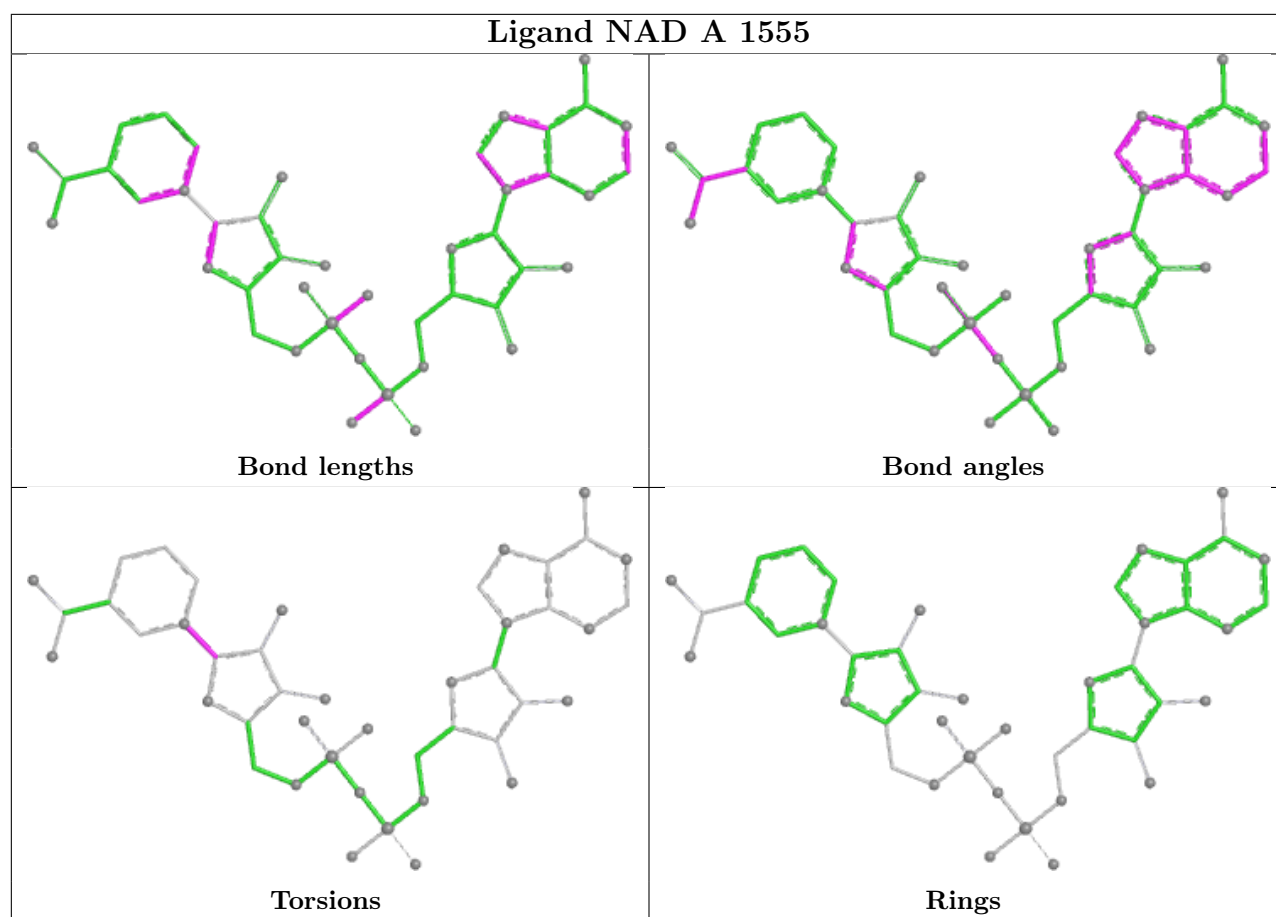
There are no ring outliers.

4 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1555	NAD	7	0
3	B	1556	URO	8	0
2	A	1555	NAD	4	0
3	A	1556	URO	9	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 [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	551/557 (98%)	0.34	7 (1%) 75 77	2, 8, 21, 30	0
1	B	551/557 (98%)	-0.03	7 (1%) 75 77	3, 11, 24, 36	0
All	All	1102/1114 (98%)	0.16	14 (1%) 75 77	2, 10, 23, 36	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	SER	4.2
1	B	163	GLY	3.3
1	B	5	ASN	2.9
1	B	555	ILE	2.7
1	B	6	ASN	2.6
1	A	163	GLY	2.6
1	A	164	SER	2.6
1	A	231	ALA	2.5
1	A	167	GLY	2.5
1	A	291	GLU	2.2
1	A	13	ARG	2.1
1	A	290	THR	2.1
1	B	55	ILE	2.1
1	B	285	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

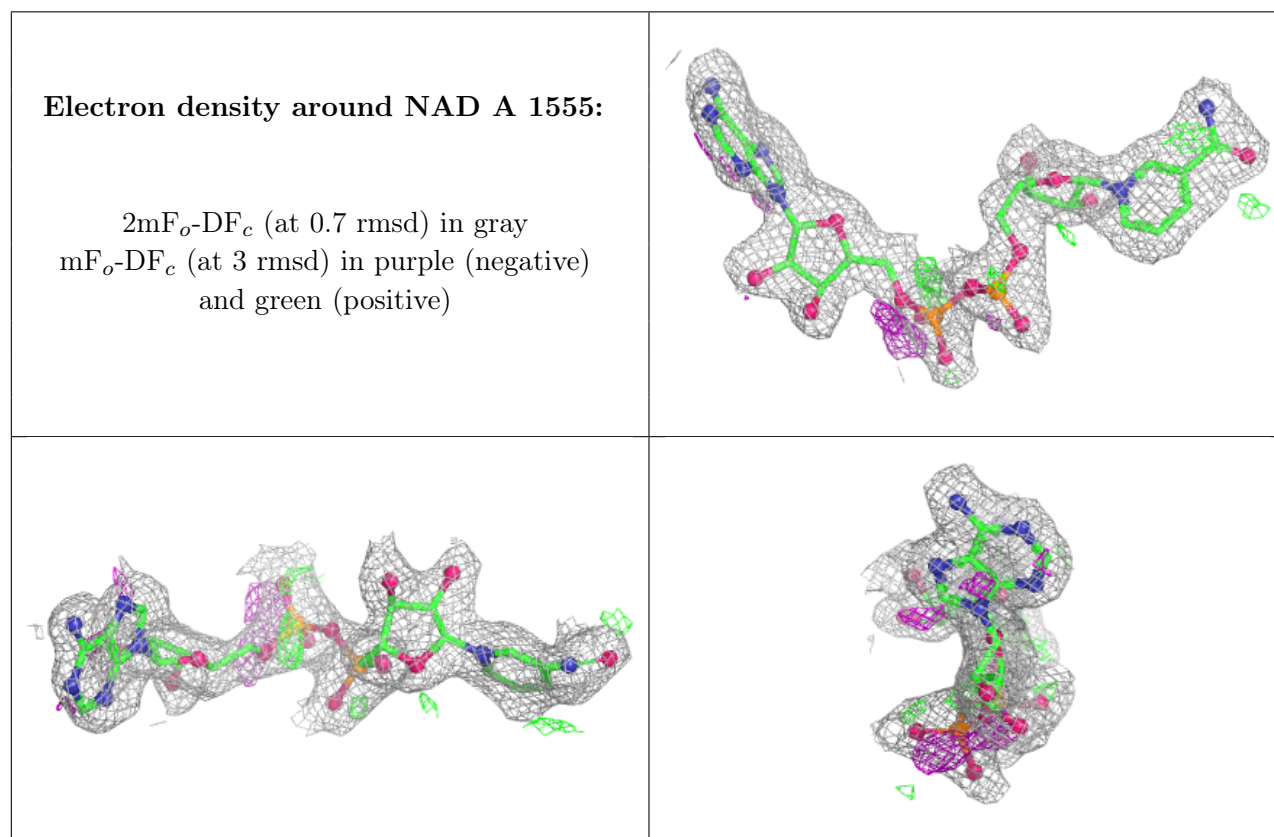
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

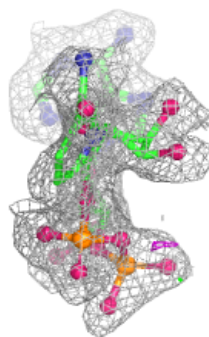
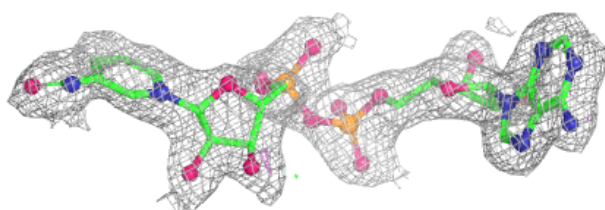
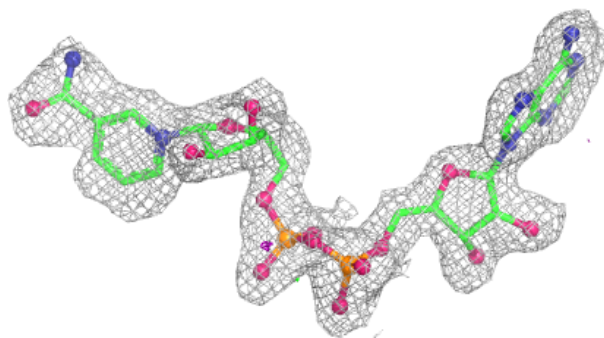
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	URO	A	1556	10/10	0.84	0.26	26,39,44,46	0
3	URO	B	1556	10/10	0.86	0.23	25,39,43,43	0
2	NAD	A	1555	44/44	0.91	0.10	2,6,22,27	0
2	NAD	B	1555	44/44	0.96	0.07	7,11,26,30	0

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



Electron density around NAD B 1555:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.