



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:05 AM UTC

PDB ID : 3OM9 / pdb_00003om9
Title : T. Gondii bradyzoite-specific LDH (LDH1) in complex with NAD and OXQ
Authors : Susko, M.S.; Wilson, D.K.
Deposited on : 2010-08-26
Resolution : 1.98 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

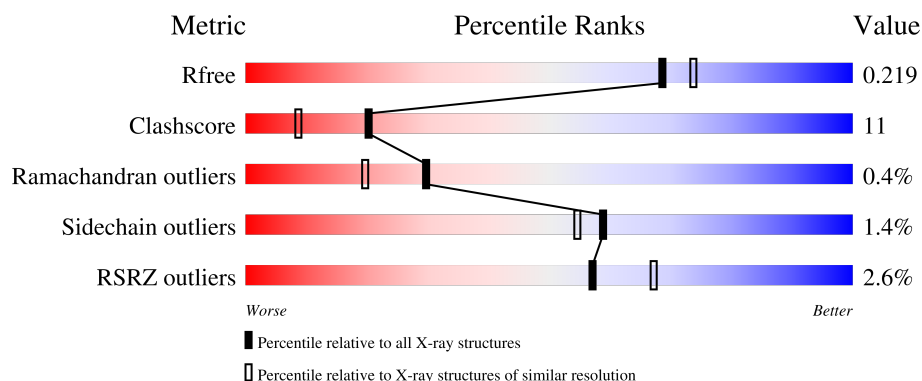
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	329	2% 75% 23% .
1	B	329	2% 78% 21% .
1	C	329	2% 77% 20% ..
1	D	329	3% 79% 20% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10521 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.

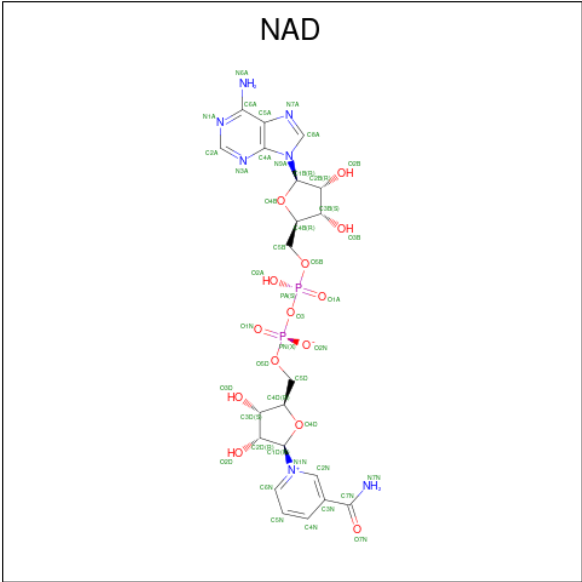
- Molecule 1 is a protein called Lactate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2479	1568	419	469	23			
1	B	329	Total	C	N	O	S	0	0	0
			2479	1568	419	469	23			
1	C	326	Total	C	N	O	S	0	0	0
			2462	1558	416	465	23			
1	D	326	Total	C	N	O	S	0	0	0
			2462	1558	416	465	23			

There are 8 discrepancies between the modelled and reference sequences:

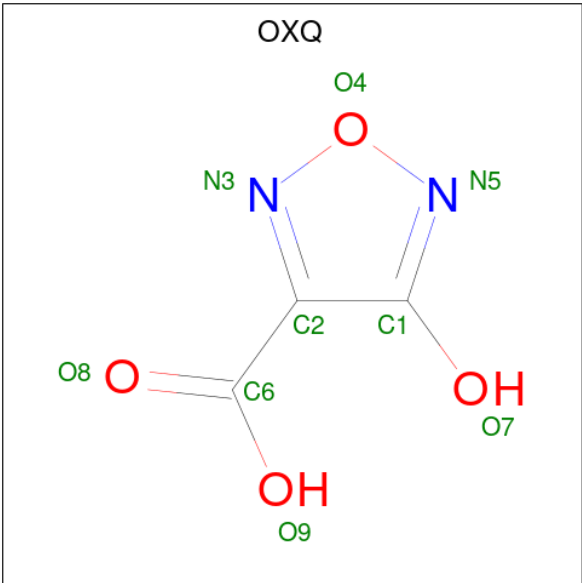
Chain	Residue	Modelled	Actual	Comment	Reference
A	334	PRO	-	expression tag	UNP P90613
A	335	GLY	-	expression tag	UNP P90613
B	334	PRO	-	expression tag	UNP P90613
B	335	GLY	-	expression tag	UNP P90613
C	334	PRO	-	expression tag	UNP P90613
C	335	GLY	-	expression tag	UNP P90613
D	334	PRO	-	expression tag	UNP P90613
D	335	GLY	-	expression tag	UNP P90613

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is 4-HYDROXY-1,2,5-OXADIAZOLE-3-CARBOXYLIC ACID (CCD ID: OXQ) (formula: C₃H₂N₂O₄).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 9 3 2 4	0	0
3	B	1	Total C N O 9 3 2 4	0	0
3	C	1	Total C N O 9 3 2 4	0	0
3	D	1	Total C N O 9 3 2 4	0	0

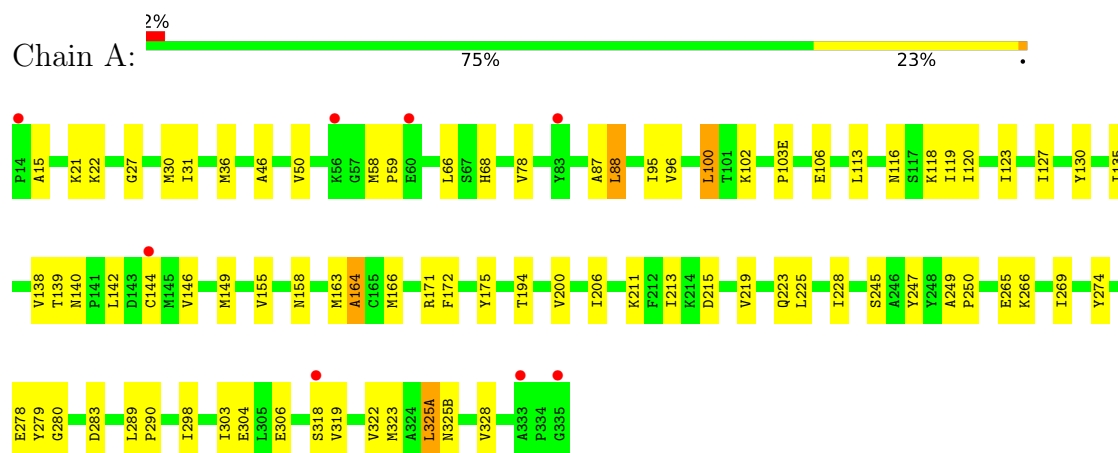
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	131	Total O 131 131	0	0
4	B	104	Total O 104 104	0	0
4	C	98	Total O 98 98	0	0
4	D	94	Total O 94 94	0	0

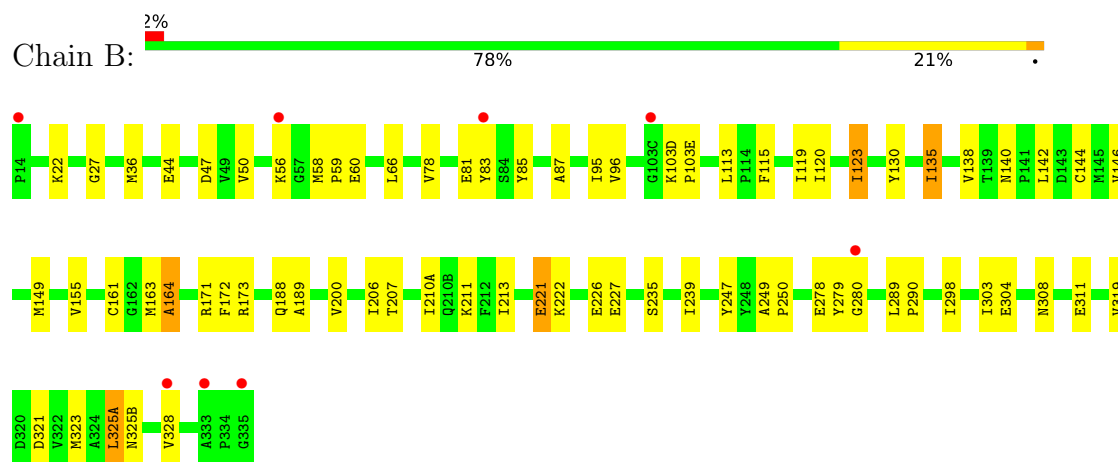
3 Residue-property plots

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 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.

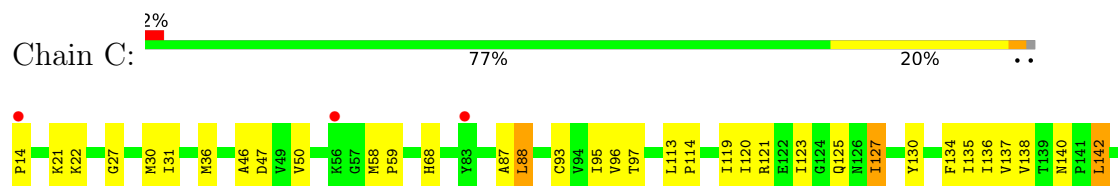
• Molecule 1: Lactate dehydrogenase

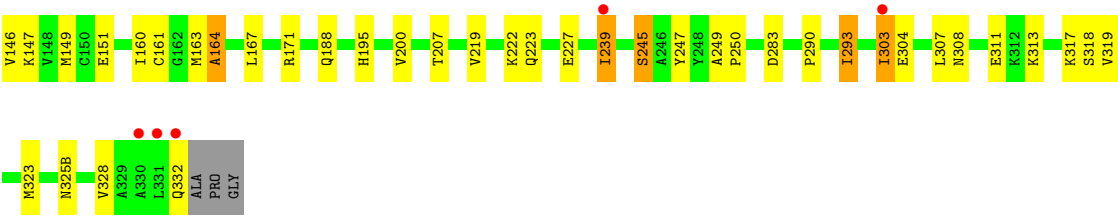


• Molecule 1: Lactate dehydrogenase

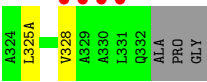
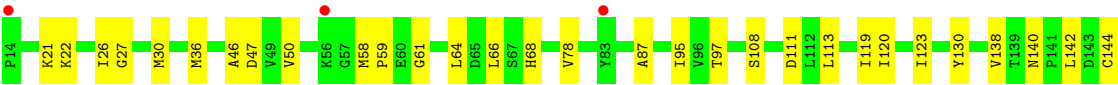
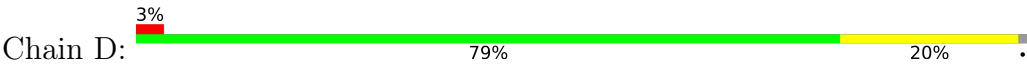


• Molecule 1: Lactate dehydrogenase





• Molecule 1: Lactate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.31Å 124.46Å 86.72Å 90.00° 105.61° 90.00°	Depositor
Resolution (Å)	30.00 – 1.98 30.00 – 1.99	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.98) 98.4 (30.00-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	26.73 (at 1.98Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.196 , 0.223 0.196 , 0.219	Depositor DCC
R_{free} test set	2822 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	15.7	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 31.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10521	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: OXQ, 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	0.36	0/2518	0.86	5/3405 (0.1%)
1	B	0.37	0/2518	0.90	6/3405 (0.2%)
1	C	0.36	0/2500	0.88	7/3381 (0.2%)
1	D	0.35	0/2500	0.87	4/3381 (0.1%)
All	All	0.36	0/10036	0.88	22/13572 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	200	VAL	N-CA-C	9.32	117.38	108.15
1	D	200	VAL	N-CA-C	8.79	116.85	108.15
1	B	200	VAL	N-CA-C	8.66	116.73	108.15
1	A	200	VAL	N-CA-C	8.56	116.62	108.15
1	C	88	LEU	N-CA-C	6.18	119.29	111.69
1	A	96	VAL	N-CA-C	5.83	116.20	107.51
1	B	96	VAL	N-CA-C	5.66	115.94	107.51
1	C	47	ASP	N-CA-C	-5.66	102.53	110.50
1	C	27	GLY	N-CA-C	-5.64	99.82	113.18
1	A	130	TYR	N-CA-C	5.49	120.09	112.90
1	D	130	TYR	N-CA-C	5.49	120.25	113.50
1	C	130	TYR	N-CA-C	5.46	120.08	113.41
1	C	96	VAL	N-CA-C	5.46	115.72	107.80
1	D	47	ASP	N-CA-C	-5.41	102.87	110.50
1	A	27	GLY	N-CA-C	-5.38	100.44	113.18
1	D	27	GLY	N-CA-C	-5.35	100.50	113.18
1	B	47	ASP	N-CA-C	-5.34	102.97	110.50
1	B	27	GLY	N-CA-C	-5.28	100.66	113.18
1	B	213	ILE	N-CA-C	-5.20	105.78	110.82
1	B	130	TYR	N-CA-C	5.11	120.17	113.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	127	ILE	N-CA-C	-5.04	105.63	110.72
1	A	88	LEU	N-CA-C	5.00	119.45	112.90

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	2479	0	2533	68	0
1	B	2479	0	2533	60	0
1	C	2462	0	2518	66	0
1	D	2462	0	2518	47	0
2	A	44	0	26	1	0
2	B	44	0	26	1	0
2	C	44	0	26	4	0
2	D	44	0	26	1	0
3	A	9	0	0	1	0
3	B	9	0	0	0	0
3	C	9	0	0	2	0
3	D	9	0	0	1	0
4	A	131	0	0	1	0
4	B	104	0	0	1	0
4	C	98	0	0	4	0
4	D	94	0	0	2	0
All	All	10521	0	10206	230	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 11.

All (230) 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:D:142:LEU:HD22	1:D:164:ALA:HB2	1.42	0.99
1:A:213:ILE:HD13	1:A:219:VAL:HG23	1.46	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325(A):LEU:HG	4:B:422:HOH:O	1.77	0.85
1:B:119:ILE:O	1:B:123:ILE:HD13	1.75	0.84
1:C:113:LEU:HD11	1:C:328:VAL:CG1	2.06	0.83
1:C:123:ILE:O	1:C:127:ILE:HD13	1.77	0.83
1:A:175:TYR:HB3	1:A:228:ILE:HD12	1.57	0.83
1:A:325(A):LEU:O	1:A:328:VAL:HG23	1.83	0.79
1:C:119:ILE:HD12	2:C:705:NAD:H61A	1.48	0.78
1:A:120:ILE:HD11	1:A:139:THR:HG21	1.66	0.78
1:D:220:THR:OG1	1:D:223:GLN:HG3	1.85	0.75
1:C:135:ILE:HB	1:C:160:ILE:HD13	1.69	0.75
1:C:134:PHE:CD2	1:C:136:ILE:HD11	2.21	0.75
1:D:26:ILE:HG21	1:D:123:ILE:HD12	1.71	0.73
1:D:142:LEU:HD22	1:D:164:ALA:CB	2.18	0.72
1:A:103(E):PRO:HG2	1:A:106:GLU:HG3	1.71	0.72
1:C:93:CYS:SG	1:C:136:ILE:HD13	2.30	0.72
1:C:119:ILE:HD12	2:C:705:NAD:N6A	2.05	0.71
1:D:113:LEU:HD21	1:D:328:VAL:CG2	2.21	0.70
1:C:14:PRO:HB3	4:D:371:HOH:O	1.92	0.69
1:D:222:LYS:O	1:D:227:GLU:HG3	1.91	0.69
1:B:239:ILE:HD11	1:D:64:LEU:HB2	1.75	0.69
1:B:325(A):LEU:O	1:B:328:VAL:HG23	1.94	0.68
1:B:113:LEU:HD22	1:B:325(A):LEU:HD13	1.76	0.68
1:A:123:ILE:O	1:A:127:ILE:HD13	1.93	0.68
1:A:116:ASN:O	1:A:120:ILE:HD13	1.94	0.68
1:A:135:ILE:HD13	1:A:155:VAL:HG21	1.74	0.68
1:B:161:CYS:HB3	1:B:298:ILE:HD11	1.77	0.67
1:A:15:ALA:HB1	4:A:5:HOH:O	1.95	0.66
1:C:303:ILE:HD12	1:C:303:ILE:N	2.11	0.65
1:B:135:ILE:N	1:B:135:ILE:HD12	2.10	0.65
1:C:113:LEU:HD11	1:C:328:VAL:HG13	1.77	0.65
1:A:30:MET:HE1	1:C:30:MET:HE1	1.78	0.65
1:A:158:ASN:O	1:A:298:ILE:HD12	1.97	0.64
1:A:144:CYS:SG	1:A:325(A):LEU:HB2	2.38	0.64
1:B:113:LEU:HD11	1:B:328:VAL:HG22	1.79	0.64
1:C:319:VAL:HG12	1:C:323:MET:HE2	1.80	0.63
1:A:50:VAL:HG11	1:A:87:ALA:HA	1.80	0.63
1:B:115:PHE:O	1:B:119:ILE:HD13	1.99	0.63
1:A:120:ILE:HG23	1:A:149:MET:HE2	1.81	0.62
1:D:50:VAL:HG11	1:D:87:ALA:HA	1.81	0.62
1:A:135:ILE:HD12	1:A:135:ILE:N	2.15	0.62
1:C:120:ILE:HG23	1:C:149:MET:HE2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LEU:HD21	1:A:328:VAL:HG22	1.81	0.61
1:C:121:ARG:NH1	1:C:332:GLN:HG3	2.15	0.61
1:D:113:LEU:HD21	1:D:328:VAL:HG23	1.83	0.61
1:A:113:LEU:HD11	1:A:328:VAL:HG22	1.82	0.61
1:D:144:CYS:SG	1:D:328:VAL:HG21	2.42	0.60
1:A:135:ILE:HD11	1:A:155:VAL:HG11	1.84	0.60
1:C:293:ILE:HD12	1:C:293:ILE:N	2.16	0.60
1:D:142:LEU:O	1:D:146:VAL:HG23	2.03	0.59
1:D:290:PRO:HB2	1:D:303:ILE:HB	1.84	0.59
1:B:113:LEU:HD22	1:B:325(A):LEU:CD1	2.33	0.59
1:B:144:CYS:SG	1:B:328:VAL:HG21	2.43	0.58
1:B:221:GLU:OE2	1:B:226:GLU:HG3	2.02	0.58
1:C:121:ARG:O	1:C:125:GLN:HG3	2.03	0.58
1:A:306:GLU:HG3	1:D:211:LYS:HD3	1.86	0.58
1:B:142:LEU:HG	1:B:164:ALA:HB2	1.85	0.58
1:C:135:ILE:C	1:C:136:ILE:HD12	2.28	0.58
1:D:22:LYS:HE3	1:D:50:VAL:HG21	1.84	0.58
1:C:142:LEU:O	1:C:146:VAL:HG23	2.04	0.57
1:D:322:VAL:HA	1:D:325(A):LEU:HD12	1.86	0.57
1:C:136:ILE:HD12	1:C:136:ILE:N	2.19	0.57
1:A:68:HIS:CE1	1:C:171:ARG:HG2	2.39	0.57
1:A:144:CYS:SG	1:A:325(A):LEU:CB	2.92	0.57
1:B:140:ASN:HA	1:B:142:LEU:N	2.18	0.57
1:B:239:ILE:HD12	1:D:61:GLY:O	2.04	0.57
1:C:313:LYS:HG2	1:C:317:LYS:HE3	1.87	0.57
1:A:319:VAL:O	1:A:323:MET:HG3	2.05	0.57
1:B:239:ILE:HD11	1:D:64:LEU:CB	2.34	0.57
1:B:119:ILE:HD12	1:B:119:ILE:H	1.69	0.57
1:A:166:MET:HE3	1:A:269:ILE:HD12	1.87	0.56
1:B:135:ILE:HD13	1:B:155:VAL:HG21	1.87	0.56
1:C:140:ASN:HA	1:C:142:LEU:N	2.21	0.56
1:A:194:THR:H	1:A:318:SER:HB2	1.69	0.56
1:B:119:ILE:HD12	1:B:119:ILE:N	2.20	0.56
1:C:142:LEU:HG	1:C:164:ALA:HB2	1.88	0.56
1:D:166:MET:HE3	1:D:269:ILE:HD12	1.87	0.56
1:C:50:VAL:HG11	1:C:87:ALA:HA	1.87	0.56
1:C:308:ASN:OD1	1:C:311:GLU:HG3	2.06	0.56
1:B:319:VAL:HG12	1:B:323:MET:HE2	1.87	0.56
1:D:140:ASN:HA	1:D:142:LEU:N	2.20	0.56
1:B:142:LEU:O	1:B:146:VAL:HG23	2.06	0.55
1:C:222:LYS:O	1:C:227:GLU:HG3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:PHE:CZ	1:A:206:ILE:HD11	2.41	0.55
1:B:290:PRO:HG2	1:B:303:ILE:HB	1.88	0.55
1:D:301:ARG:HD3	1:D:303:ILE:HD11	1.89	0.55
1:A:36:MET:SD	1:A:95:ILE:HG21	2.47	0.55
1:B:172:PHE:CZ	1:B:206:ILE:HD11	2.41	0.55
1:B:222:LYS:O	1:B:227:GLU:HG3	2.06	0.54
1:B:113:LEU:HD11	1:B:328:VAL:CG2	2.37	0.54
1:C:22:LYS:HE3	1:C:50:VAL:HG21	1.88	0.54
1:D:119:ILE:O	1:D:123:ILE:HG12	2.07	0.54
1:D:138:VAL:O	2:D:707:NAD:H2N	2.06	0.54
1:B:138:VAL:O	2:B:703:NAD:H2N	2.07	0.53
1:A:171:ARG:HG2	1:C:68:HIS:CE1	2.44	0.53
1:A:113:LEU:HD11	1:A:328:VAL:CG2	2.38	0.52
1:A:213:ILE:CD1	1:A:219:VAL:HG23	2.30	0.52
1:A:135:ILE:HD12	1:A:135:ILE:H	1.75	0.52
1:A:142:LEU:HG	1:A:164:ALA:HB2	1.92	0.52
1:A:223:GLN:O	1:A:228:ILE:HG12	2.10	0.52
1:A:140:ASN:HA	1:A:142:LEU:N	2.24	0.52
1:D:66:LEU:HB2	1:D:78:VAL:HG11	1.92	0.52
1:B:22:LYS:HE2	1:B:50:VAL:HG21	1.91	0.52
1:B:36:MET:SD	1:B:95:ILE:HG21	2.50	0.52
1:B:60:GLU:HG3	1:B:83:TYR:OH	2.10	0.52
1:C:147:LYS:O	1:C:151:GLU:HG3	2.10	0.52
1:C:21:LYS:O	1:C:46:ALA:HB1	2.10	0.52
1:C:313:LYS:HE2	1:C:317:LYS:NZ	2.25	0.51
1:C:219:VAL:HG23	1:C:223:GLN:CD	2.35	0.51
1:C:249:ALA:HB3	1:C:250:PRO:HD3	1.93	0.51
1:A:213:ILE:HD11	1:A:225:LEU:CD1	2.41	0.51
1:D:22:LYS:HE3	1:D:50:VAL:CG2	2.40	0.51
1:B:279:TYR:CE2	1:B:289:LEU:HD22	2.45	0.51
1:D:308:ASN:OD1	1:D:311:GLU:HG3	2.10	0.51
1:A:194:THR:HG21	1:A:322:VAL:HB	1.92	0.50
1:B:50:VAL:HG11	1:B:87:ALA:HA	1.93	0.50
1:B:172:PHE:CE2	1:B:206:ILE:HD11	2.46	0.50
1:C:138:VAL:O	2:C:705:NAD:H2N	2.11	0.50
1:D:188:GLN:HB2	1:D:207:THR:OG1	2.10	0.49
1:A:290:PRO:HG2	1:A:303:ILE:HB	1.94	0.49
1:B:85:TYR:OH	1:B:123:ILE:HD12	2.12	0.49
1:A:279:TYR:CE2	1:A:289:LEU:HD22	2.46	0.49
1:B:247:TYR:C	1:B:250:PRO:HD2	2.38	0.49
1:D:163:MET:SD	1:D:163:MET:C	2.96	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:MET:SD	1:C:163:MET:C	2.96	0.48
1:A:303:ILE:HD12	1:A:303:ILE:N	2.28	0.48
1:A:211:LYS:HE2	4:D:349:HOH:O	2.13	0.48
1:A:172:PHE:CE2	1:A:206:ILE:HD11	2.48	0.48
1:A:138:VAL:O	2:A:701:NAD:H2N	2.13	0.48
1:C:222:LYS:HG3	4:C:345:HOH:O	2.14	0.48
1:A:118:LYS:HE3	1:A:119:ILE:HG13	1.95	0.48
1:C:58:MET:HB3	1:C:59:PRO:HD3	1.95	0.48
1:A:21:LYS:O	1:A:46:ALA:HB1	2.14	0.47
1:B:135:ILE:HD11	1:B:155:VAL:HG11	1.95	0.47
1:A:213:ILE:HD11	1:A:225:LEU:HD11	1.97	0.47
1:C:97:THR:HB	2:C:705:NAD:H51N	1.96	0.47
1:A:22:LYS:HE3	1:A:50:VAL:HG21	1.96	0.47
1:C:195:HIS:NE2	3:C:706:OXQ:O7	2.44	0.47
1:B:119:ILE:H	1:B:119:ILE:CD1	2.28	0.46
1:B:120:ILE:HG23	1:B:149:MET:HE2	1.97	0.46
1:C:239:ILE:HD13	4:C:394:HOH:O	2.14	0.46
1:D:120:ILE:HG23	1:D:149:MET:HE2	1.97	0.46
1:B:163:MET:SD	1:B:163:MET:C	2.98	0.46
1:C:160:ILE:HD12	1:C:161:CYS:H	1.80	0.46
1:D:26:ILE:CG2	1:D:123:ILE:HD12	2.43	0.46
1:A:163:MET:C	1:A:163:MET:SD	2.98	0.46
1:C:245:SER:HB2	3:C:706:OXQ:N3	2.31	0.46
1:A:66:LEU:HB2	1:A:78:VAL:HG11	1.96	0.46
1:B:142:LEU:HD13	1:B:142:LEU:C	2.41	0.46
1:A:58:MET:HB3	1:A:59:PRO:HD3	1.96	0.46
1:A:249:ALA:N	1:A:250:PRO:CD	2.79	0.46
1:A:142:LEU:O	1:A:146:VAL:HG23	2.16	0.45
1:D:249:ALA:HB3	1:D:250:PRO:HD3	1.99	0.45
1:D:319:VAL:O	1:D:323:MET:HG3	2.17	0.45
1:A:245:SER:HB2	3:A:702:OXQ:N3	2.32	0.45
1:D:97:THR:HG22	1:D:138:VAL:HB	1.97	0.45
1:A:88:LEU:HD13	1:A:127:ILE:HD12	1.99	0.44
1:A:211:LYS:HE3	1:A:215:ASP:OD2	2.16	0.44
1:C:303:ILE:N	1:C:303:ILE:CD1	2.80	0.44
1:C:134:PHE:CE2	1:C:136:ILE:HD11	2.52	0.44
1:D:58:MET:HB3	1:D:59:PRO:HD3	1.99	0.44
1:A:247:TYR:C	1:A:250:PRO:HD2	2.42	0.44
1:C:142:LEU:HD13	1:C:142:LEU:C	2.42	0.44
1:C:167:LEU:HD22	1:C:250:PRO:HG3	1.99	0.44
1:C:188:GLN:HB2	1:C:207:THR:OG1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:ARG:HG2	1:C:68:HIS:CG	2.53	0.44
1:C:88:LEU:CD1	1:C:127:ILE:HD12	2.48	0.44
1:B:171:ARG:HG2	1:D:68:HIS:CE1	2.53	0.44
1:D:36:MET:SD	1:D:95:ILE:HG21	2.58	0.44
1:C:22:LYS:HE3	1:C:50:VAL:CG2	2.47	0.43
1:C:319:VAL:O	1:C:323:MET:HG3	2.17	0.43
1:C:160:ILE:HD12	1:C:161:CYS:N	2.32	0.43
1:C:239:ILE:CD1	4:C:394:HOH:O	2.66	0.43
1:C:290:PRO:HB2	1:C:303:ILE:HB	2.00	0.43
1:D:21:LYS:O	1:D:46:ALA:HB1	2.18	0.43
1:C:31:ILE:O	1:C:36:MET:HG3	2.17	0.43
1:B:66:LEU:HB2	1:B:78:VAL:HG11	2.00	0.43
1:D:30:MET:HE3	1:D:247:TYR:CZ	2.53	0.43
1:A:142:LEU:C	1:A:142:LEU:HD13	2.43	0.43
1:C:325(B):ASN:HA	1:C:328:VAL:HG22	2.01	0.43
1:A:194:THR:CG2	1:A:322:VAL:HB	2.49	0.43
1:B:211:LYS:HE2	1:C:304:GLU:O	2.19	0.43
1:B:278:GLU:C	1:B:280:GLY:N	2.76	0.43
1:D:108:SER:HB3	1:D:111:ASP:OD2	2.19	0.43
1:A:100:LEU:HD13	1:A:102:LYS:O	2.19	0.43
1:D:171:ARG:HD3	1:D:235:SER:HB2	2.01	0.43
1:D:319:VAL:HG12	1:D:323:MET:HE2	2.00	0.43
1:B:249:ALA:N	1:B:250:PRO:CD	2.82	0.42
1:B:173:ARG:HG3	1:B:189:ALA:HB3	2.01	0.42
1:A:22:LYS:HE3	1:A:50:VAL:CG2	2.50	0.42
1:B:188:GLN:HB2	1:B:207:THR:OG1	2.19	0.42
1:B:279:TYR:CZ	1:B:304:GLU:HA	2.55	0.42
1:B:321:ASP:O	1:B:325(A):LEU:HD23	2.19	0.42
1:C:135:ILE:CB	1:C:160:ILE:HD13	2.45	0.42
1:D:249:ALA:N	1:D:250:PRO:CD	2.83	0.42
1:A:265:GLU:OE2	1:B:44:GLU:OE2	2.37	0.42
1:A:279:TYR:CZ	1:A:304:GLU:HA	2.54	0.42
1:A:274:TYR:CE1	1:A:283:ASP:HA	2.54	0.42
1:A:278:GLU:C	1:A:280:GLY:N	2.78	0.42
1:B:278:GLU:C	1:B:280:GLY:H	2.28	0.42
1:B:58:MET:HB3	1:B:59:PRO:HD3	2.02	0.42
1:B:172:PHE:CD2	1:B:189:ALA:HB1	2.55	0.42
1:B:308:ASN:OD1	1:B:311:GLU:HG3	2.20	0.41
1:D:172:PHE:CD2	1:D:189:ALA:HB1	2.55	0.41
1:B:103(D):LYS:HA	1:B:103(E):PRO:HD3	1.94	0.41
1:D:278:GLU:O	1:D:279:TYR:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:113:LEU:HB3	1:C:114:PRO:CD	2.51	0.41
1:D:140:ASN:ND2	3:D:708:OXQ:N5	2.68	0.41
1:A:278:GLU:C	1:A:280:GLY:H	2.29	0.41
1:C:164:ALA:HB3	4:C:419:HOH:O	2.21	0.41
1:A:265:GLU:O	1:A:266:LYS:HB2	2.21	0.41
1:B:56:LYS:HG2	1:B:83:TYR:CE2	2.55	0.41
1:C:247:TYR:C	1:C:250:PRO:HD2	2.45	0.41
1:B:22:LYS:HE2	1:B:50:VAL:CG2	2.50	0.41
1:B:171:ARG:HD3	1:B:235:SER:HB2	2.03	0.41
1:B:206:ILE:HB	1:B:210(A):ILE:HB	2.02	0.41
1:C:36:MET:SD	1:C:95:ILE:HG21	2.61	0.41
1:C:307:LEU:HB3	1:C:311:GLU:HB2	2.02	0.41
1:C:137:VAL:HG23	1:C:160:ILE:HD11	2.03	0.41
1:B:211:LYS:HA	1:B:211:LYS:HD2	1.88	0.40
1:D:147:LYS:O	1:D:151:GLU:HG3	2.20	0.40
1:D:301:ARG:HD3	1:D:303:ILE:CD1	2.50	0.40
1:A:144:CYS:SG	1:A:325(A):LEU:HB3	2.61	0.40
1:C:134:PHE:HD2	1:C:136:ILE:HD11	1.82	0.40
1:A:31:ILE:O	1:A:36:MET:HG3	2.21	0.40
1:A:120:ILE:HD12	1:A:120:ILE:N	2.37	0.40
1:B:50:VAL:HG13	1:B:81:GLU:HG2	2.03	0.40
1:C:88:LEU:HD13	1:C:127:ILE:HD12	2.03	0.40
1:D:247:TYR:C	1:D:250:PRO:HD2	2.46	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	327/329 (99%)	318 (97%)	8 (2%)	1 (0%)	36 27
1	B	327/329 (99%)	319 (98%)	7 (2%)	1 (0%)	36 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	324/329 (98%)	316 (98%)	6 (2%)	2 (1%)	21	12
1	D	324/329 (98%)	316 (98%)	7 (2%)	1 (0%)	36	27
All	All	1302/1316 (99%)	1269 (98%)	28 (2%)	5 (0%)	30	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	ALA
1	D	164	ALA
1	B	164	ALA
1	C	164	ALA
1	C	142	LEU

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	270/270 (100%)	267 (99%)	3 (1%)	65	61
1	B	270/270 (100%)	265 (98%)	5 (2%)	50	44
1	C	269/270 (100%)	263 (98%)	6 (2%)	45	38
1	D	269/270 (100%)	268 (100%)	1 (0%)	84	82
All	All	1078/1080 (100%)	1063 (99%)	15 (1%)	59	54

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	LEU
1	A	325(A)	LEU
1	A	325(B)	ASN
1	B	123	ILE
1	B	135	ILE
1	B	221	GLU
1	B	325(A)	LEU

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Mol	Chain	Res	Type
1	B	325(B)	ASN
1	C	239	ILE
1	C	245	SER
1	C	283	ASP
1	C	293	ILE
1	C	303	ILE
1	C	318	SER
1	D	318	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	276	ASN
1	A	314	GLN
1	B	276	ASN
1	B	314	GLN
1	C	314	GLN
1	D	332	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 ⓘ

8 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	A	701	-	46,48,48	1.92	11 (23%)	64,73,73	1.37	11 (17%)
3	OXQ	C	706	-	9,9,9	1.50	2 (22%)	12,12,12	1.45	3 (25%)
2	NAD	B	703	-	46,48,48	1.90	11 (23%)	64,73,73	1.34	10 (15%)
2	NAD	D	707	-	46,48,48	1.95	11 (23%)	64,73,73	1.32	11 (17%)
3	OXQ	D	708	-	9,9,9	1.51	2 (22%)	12,12,12	1.45	3 (25%)
3	OXQ	A	702	-	9,9,9	1.52	2 (22%)	12,12,12	1.45	3 (25%)
3	OXQ	B	704	-	9,9,9	1.54	2 (22%)	12,12,12	1.45	3 (25%)
2	NAD	C	705	-	46,48,48	1.87	11 (23%)	64,73,73	1.36	12 (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	A	701	-	-	4/30/62/62	0/5/5/5
3	OXQ	C	706	-	-	0/4/4/4	0/1/1/1
2	NAD	B	703	-	-	5/30/62/62	0/5/5/5
2	NAD	D	707	-	-	4/30/62/62	0/5/5/5
3	OXQ	D	708	-	-	0/4/4/4	0/1/1/1
3	OXQ	A	702	-	-	0/4/4/4	0/1/1/1
3	OXQ	B	704	-	-	0/4/4/4	0/1/1/1
2	NAD	C	705	-	-	4/30/62/62	0/5/5/5

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	707	NAD	C2N-N1N	5.56	1.41	1.35
2	A	701	NAD	C2N-N1N	5.40	1.40	1.35
2	B	703	NAD	C2N-N1N	5.21	1.40	1.35
2	C	705	NAD	C2N-N1N	5.14	1.40	1.35
2	D	707	NAD	C4N-C3N	5.12	1.47	1.39
2	A	701	NAD	C4N-C3N	5.12	1.47	1.39
2	B	703	NAD	C4N-C3N	5.08	1.47	1.39
2	C	705	NAD	C4N-C3N	5.02	1.47	1.39
2	A	701	NAD	C2A-N1A	4.20	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	707	NAD	C2A-N1A	4.19	1.41	1.33
2	B	703	NAD	C2A-N1A	4.01	1.41	1.33
2	C	705	NAD	C2A-N1A	3.88	1.40	1.33
2	D	707	NAD	C6N-N1N	3.84	1.44	1.35
2	A	701	NAD	C6N-N1N	3.83	1.44	1.35
2	B	703	NAD	O4D-C1D	3.81	1.45	1.40
2	C	705	NAD	C6N-N1N	3.80	1.44	1.35
2	B	703	NAD	C6N-N1N	3.69	1.43	1.35
2	C	705	NAD	O4D-C1D	3.57	1.45	1.40
2	D	707	NAD	O4D-C1D	3.57	1.45	1.40
2	A	701	NAD	O4D-C1D	3.45	1.45	1.40
2	B	703	NAD	C2A-N3A	3.19	1.39	1.33
2	D	707	NAD	C2A-N3A	3.12	1.39	1.33
2	C	705	NAD	C2A-N3A	3.02	1.39	1.33
2	A	701	NAD	C2A-N3A	3.00	1.39	1.33
2	B	703	NAD	C4A-N3A	2.82	1.39	1.34
2	C	705	NAD	PA-O1A	-2.80	1.41	1.50
2	A	701	NAD	PA-O1A	-2.77	1.41	1.50
2	D	707	NAD	PA-O1A	-2.76	1.41	1.50
2	D	707	NAD	C4A-N3A	2.66	1.39	1.34
2	A	701	NAD	C4A-N3A	2.64	1.39	1.34
2	B	703	NAD	C5N-C4N	2.62	1.43	1.38
2	B	703	NAD	PA-O1A	-2.61	1.41	1.50
2	D	707	NAD	PA-O3	-2.50	1.56	1.59
2	D	707	NAD	C5N-C4N	2.47	1.43	1.38
2	A	701	NAD	C5N-C4N	2.42	1.43	1.38
2	C	705	NAD	C5N-C4N	2.42	1.43	1.38
2	C	705	NAD	C6N-C5N	2.39	1.43	1.38
2	C	705	NAD	C4A-N3A	2.37	1.38	1.34
2	B	703	NAD	C6N-C5N	2.36	1.43	1.38
3	B	704	OXQ	C2-C6	2.34	1.53	1.48
3	A	702	OXQ	C2-C6	2.33	1.53	1.48
3	C	706	OXQ	C2-C6	2.26	1.53	1.48
3	D	708	OXQ	C2-C6	2.26	1.53	1.48
3	B	704	OXQ	C1-N5	2.24	1.33	1.31
3	D	708	OXQ	C1-N5	2.23	1.33	1.31
3	C	706	OXQ	C1-N5	2.18	1.33	1.31
3	A	702	OXQ	C1-N5	2.17	1.33	1.31
2	D	707	NAD	C6N-C5N	2.17	1.43	1.38
2	B	703	NAD	C6A-N1A	2.05	1.41	1.35
2	C	705	NAD	PA-O3	-2.05	1.57	1.59
2	A	701	NAD	C6N-C5N	2.04	1.42	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	NAD	PA-O3	-2.04	1.57	1.59

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	NAD	N3A-C2A-N1A	-3.87	122.73	128.58
2	B	703	NAD	N3A-C2A-N1A	-3.78	122.86	128.58
2	D	707	NAD	N3A-C2A-N1A	-3.78	122.86	128.58
2	C	705	NAD	N3A-C2A-N1A	-3.65	123.06	128.58
2	C	705	NAD	C2N-C3N-C4N	3.00	121.75	118.26
3	A	702	OXQ	N5-O4-N3	-3.00	107.77	111.49
3	B	704	OXQ	N5-O4-N3	-2.97	107.80	111.49
2	A	701	NAD	C2N-C3N-C4N	2.97	121.72	118.26
3	D	708	OXQ	N5-O4-N3	-2.94	107.84	111.49
2	B	703	NAD	C2N-C3N-C4N	2.94	121.67	118.26
3	C	706	OXQ	N5-O4-N3	-2.92	107.86	111.49
2	D	707	NAD	C2N-C3N-C4N	2.83	121.55	118.26
2	C	705	NAD	C5N-C4N-C3N	-2.77	117.64	120.36
2	A	701	NAD	O4B-C1B-C2B	-2.73	100.77	106.62
2	C	705	NAD	O4B-C1B-C2B	-2.68	100.87	106.62
2	B	703	NAD	C5N-C4N-C3N	-2.68	117.73	120.36
2	C	705	NAD	O5B-C5B-C4B	-2.67	99.91	108.99
2	D	707	NAD	C5N-C4N-C3N	-2.58	117.83	120.36
2	D	707	NAD	O4B-C1B-C2B	-2.58	101.10	106.62
2	A	701	NAD	C5N-C4N-C3N	-2.57	117.84	120.36
2	D	707	NAD	O5B-C5B-C4B	-2.56	100.28	108.99
2	B	703	NAD	O4B-C1B-C2B	-2.55	101.16	106.62
2	B	703	NAD	O5B-C5B-C4B	-2.55	100.32	108.99
2	A	701	NAD	O5B-C5B-C4B	-2.53	100.38	108.99
2	C	705	NAD	O4B-C1B-N9A	2.50	112.89	108.09
2	A	701	NAD	O7N-C7N-C3N	-2.47	116.58	119.60
2	A	701	NAD	C3N-C7N-N7N	2.43	120.73	117.74
2	B	703	NAD	O4B-C1B-N9A	2.42	112.73	108.09
3	D	708	OXQ	C6-C2-N3	2.38	124.88	119.71
2	C	705	NAD	O7N-C7N-C3N	-2.33	116.74	119.60
2	A	701	NAD	O4B-C1B-N9A	2.32	112.54	108.09
2	C	705	NAD	C4D-O4D-C1D	-2.31	107.81	109.92
3	B	704	OXQ	C6-C2-N3	2.28	124.67	119.71
3	B	704	OXQ	O4-N5-C1	2.26	106.53	104.28
3	C	706	OXQ	C6-C2-N3	2.26	124.61	119.71
2	D	707	NAD	C5N-C6N-N1N	-2.26	117.30	120.38
3	A	702	OXQ	O4-N5-C1	2.25	106.52	104.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	706	OXQ	O4-N5-C1	2.25	106.52	104.28
2	D	707	NAD	O7N-C7N-C3N	-2.24	116.86	119.60
2	C	705	NAD	C3N-C7N-N7N	2.23	120.49	117.74
2	A	701	NAD	C4D-O4D-C1D	-2.21	107.90	109.92
3	A	702	OXQ	C6-C2-N3	2.20	124.47	119.71
3	D	708	OXQ	O4-N5-C1	2.19	106.46	104.28
2	D	707	NAD	O4B-C1B-N9A	2.18	112.28	108.09
2	C	705	NAD	O3B-C3B-C4B	-2.17	104.86	111.08
2	C	705	NAD	C1B-N9A-C8A	2.14	131.85	127.09
2	B	703	NAD	O7N-C7N-C3N	-2.14	116.99	119.60
2	D	707	NAD	C1B-N9A-C8A	2.13	131.83	127.09
2	B	703	NAD	C3N-C7N-N7N	2.12	120.35	117.74
2	B	703	NAD	C5N-C6N-N1N	-2.12	117.49	120.38
2	D	707	NAD	C3N-C7N-N7N	2.10	120.32	117.74
2	A	701	NAD	C1B-N9A-C8A	2.09	131.74	127.09
2	D	707	NAD	O3B-C3B-C4B	-2.05	105.18	111.08
2	B	703	NAD	C1B-N9A-C8A	2.05	131.65	127.09
2	C	705	NAD	C5N-C6N-N1N	-2.05	117.59	120.38
2	A	701	NAD	C5N-C6N-N1N	-2.03	117.62	120.38

There are no chirality outliers.

All (17) torsion outliers are listed below:

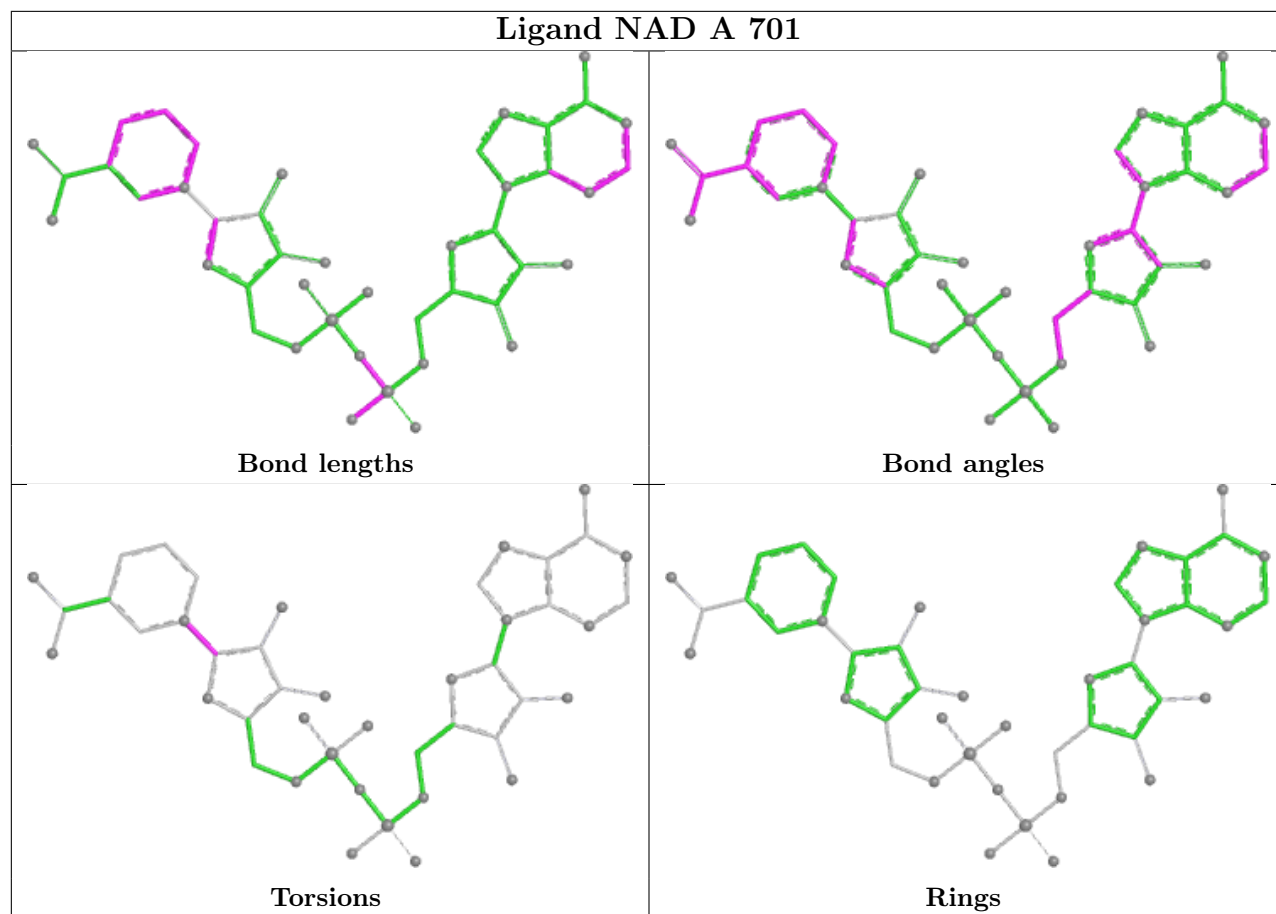
Mol	Chain	Res	Type	Atoms
2	A	701	NAD	O4D-C1D-N1N-C2N
2	A	701	NAD	O4D-C1D-N1N-C6N
2	A	701	NAD	C2D-C1D-N1N-C2N
2	A	701	NAD	C2D-C1D-N1N-C6N
2	B	703	NAD	O4D-C1D-N1N-C2N
2	B	703	NAD	O4D-C1D-N1N-C6N
2	B	703	NAD	C2D-C1D-N1N-C2N
2	B	703	NAD	C2D-C1D-N1N-C6N
2	C	705	NAD	O4D-C1D-N1N-C2N
2	C	705	NAD	O4D-C1D-N1N-C6N
2	C	705	NAD	C2D-C1D-N1N-C2N
2	C	705	NAD	C2D-C1D-N1N-C6N
2	D	707	NAD	O4D-C1D-N1N-C2N
2	D	707	NAD	O4D-C1D-N1N-C6N
2	D	707	NAD	C2D-C1D-N1N-C2N
2	D	707	NAD	C2D-C1D-N1N-C6N
2	B	703	NAD	C2B-C1B-N9A-C8A

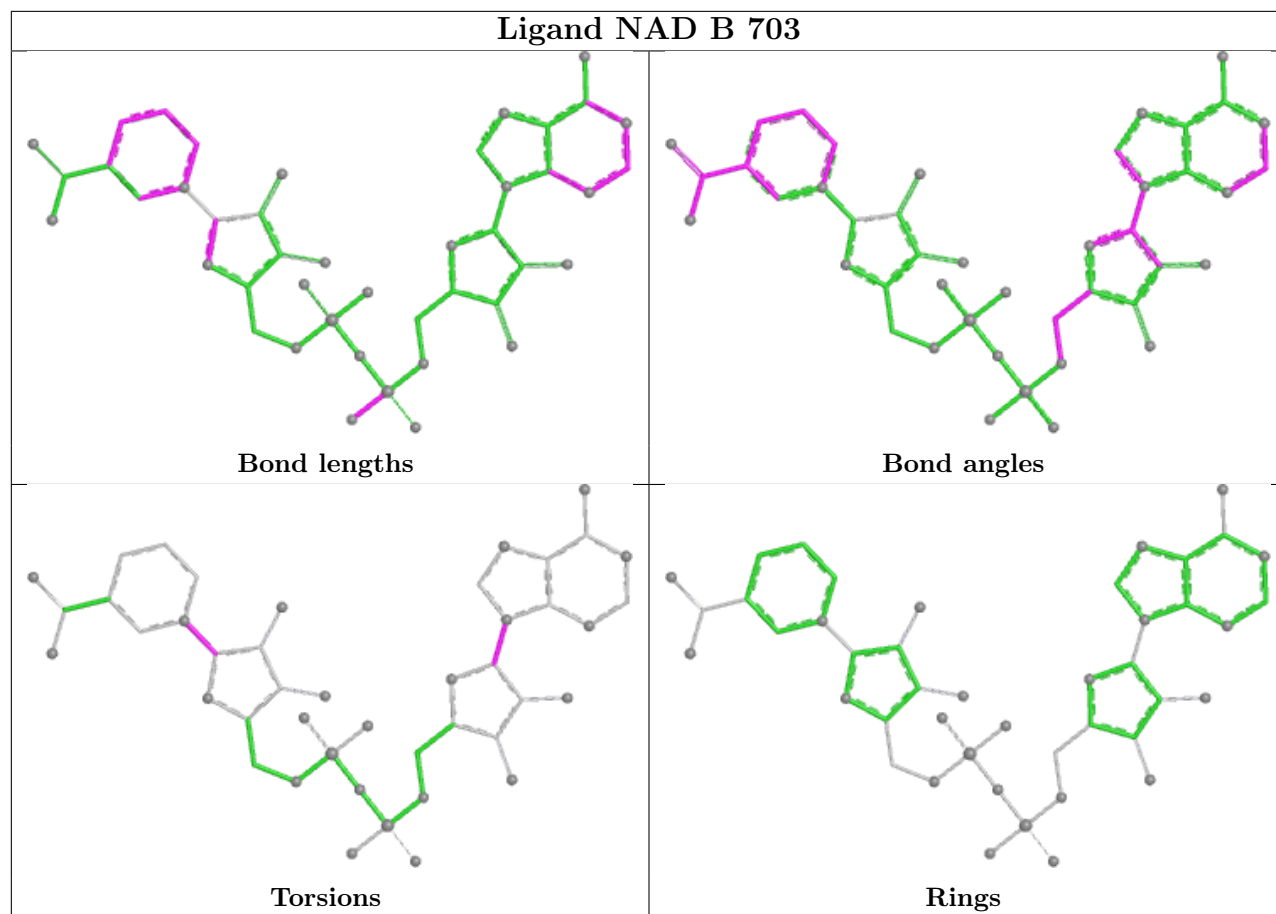
There are no ring outliers.

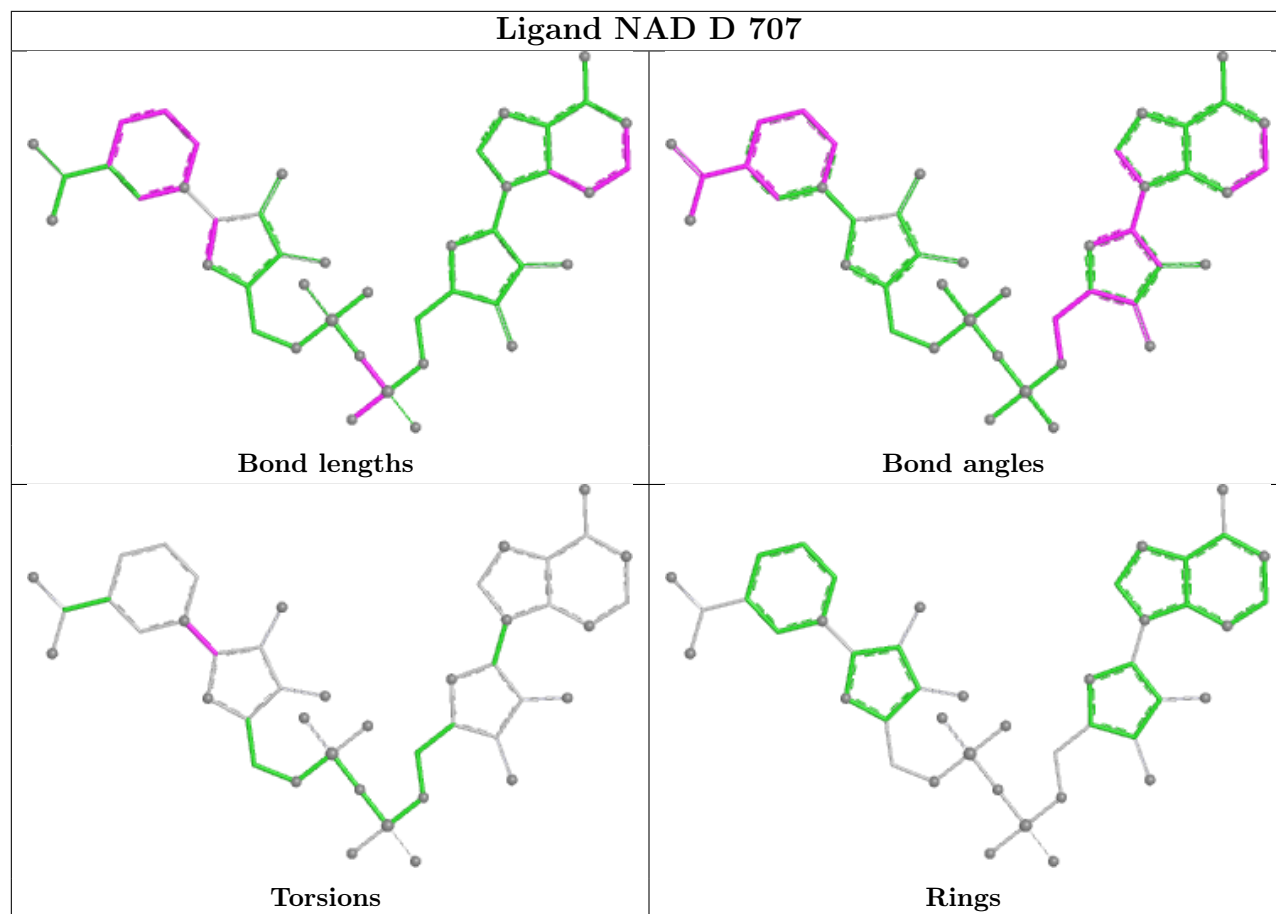
7 monomers are involved in 11 short contacts:

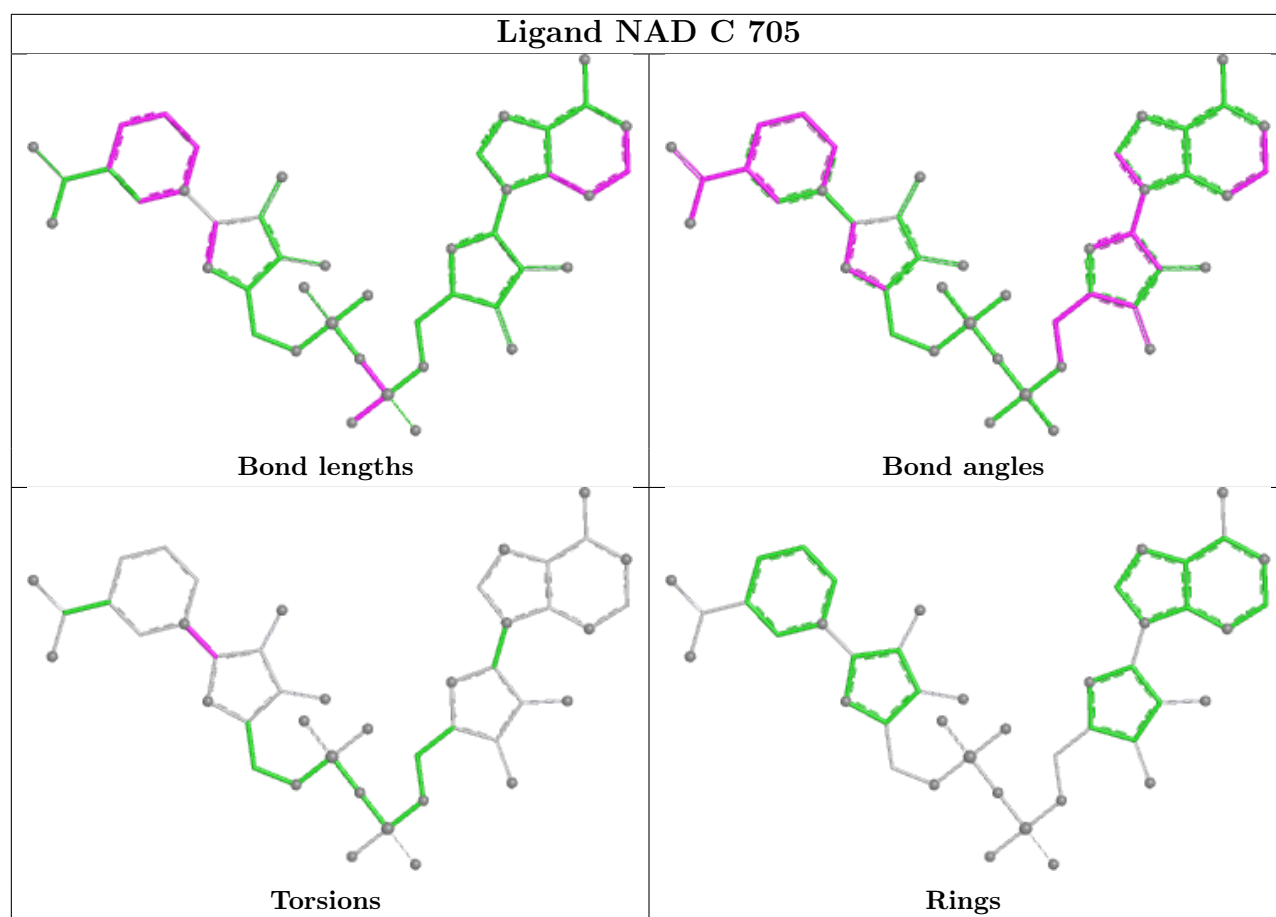
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	NAD	1	0
3	C	706	OXQ	2	0
2	B	703	NAD	1	0
2	D	707	NAD	1	0
3	D	708	OXQ	1	0
3	A	702	OXQ	1	0
2	C	705	NAD	4	0

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

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	329/329 (100%)	0.05	8 (2%) 59 69	9, 14, 29, 48	0
1	B	329/329 (100%)	0.02	8 (2%) 59 69	8, 14, 29, 44	0
1	C	326/329 (99%)	-0.01	8 (2%) 58 68	9, 15, 26, 43	0
1	D	326/329 (99%)	0.04	10 (3%) 51 61	9, 15, 29, 45	0
All	All	1310/1316 (99%)	0.02	34 (2%) 57 67	8, 15, 29, 48	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	14	PRO	8.7
1	A	335	GLY	6.6
1	B	83	TYR	6.4
1	B	335	GLY	6.3
1	A	144	CYS	5.6
1	A	83	TYR	4.2
1	B	103(C)	GLY	3.9
1	D	14	PRO	3.8
1	B	14	PRO	3.8
1	A	14	PRO	3.6
1	D	331	LEU	3.6
1	D	328	VAL	3.4
1	D	83	TYR	3.2
1	A	333	ALA	3.2
1	D	279	TYR	3.2
1	D	56	LYS	3.2
1	D	329	ALA	3.1
1	C	83	TYR	3.0
1	A	56	LYS	2.9
1	B	56	LYS	2.9
1	C	332	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	56	LYS	2.6
1	D	304	GLU	2.6
1	C	331	LEU	2.4
1	C	330	ALA	2.4
1	D	330	ALA	2.4
1	A	318	SER	2.4
1	B	280	GLY	2.2
1	C	239	ILE	2.2
1	D	269	ILE	2.2
1	A	60	GLU	2.2
1	B	333	ALA	2.2
1	C	303	ILE	2.2
1	B	328	VAL	2.1

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.

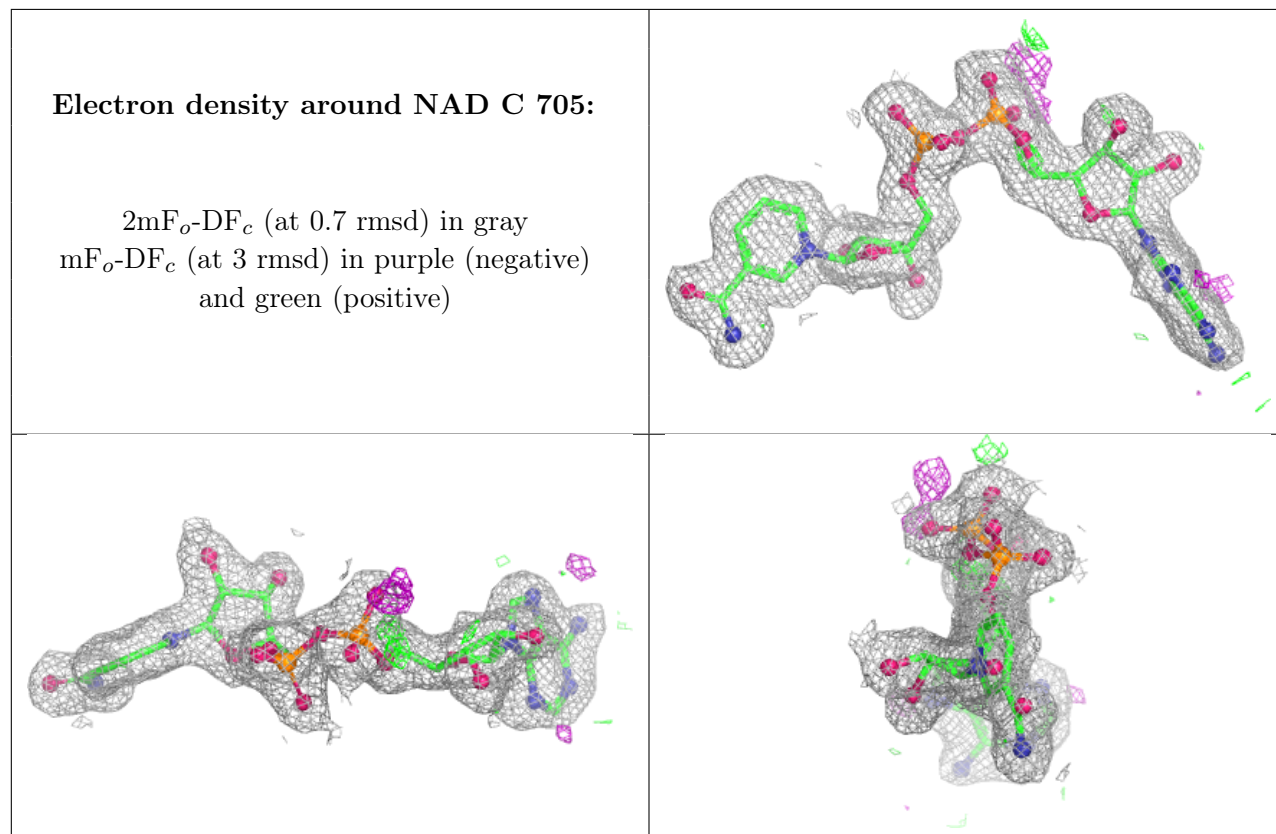
6.4 Ligands [i](#)

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	OXQ	D	708	9/9	0.92	0.10	14,16,17,19	0
3	OXQ	B	704	9/9	0.94	0.07	13,14,15,15	0
3	OXQ	C	706	9/9	0.95	0.07	12,13,15,16	0
3	OXQ	A	702	9/9	0.95	0.06	12,14,15,16	0
2	NAD	C	705	44/44	0.96	0.07	11,15,22,26	0
2	NAD	D	707	44/44	0.96	0.06	9,15,23,26	0
2	NAD	A	701	44/44	0.96	0.07	9,13,25,31	0
2	NAD	B	703	44/44	0.97	0.06	11,13,24,26	0

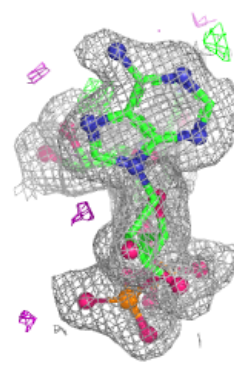
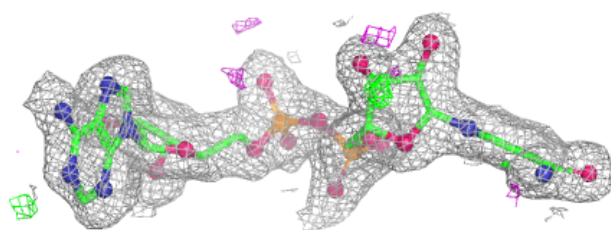
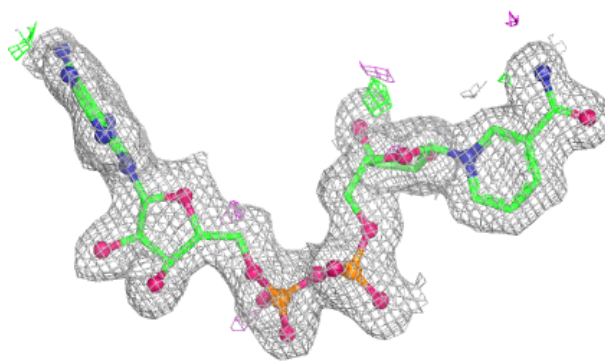
The following is a graphical depiction of the model fit to experimental electron density of all

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.

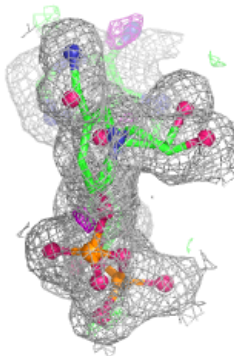
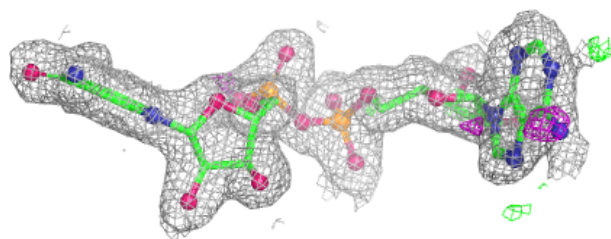
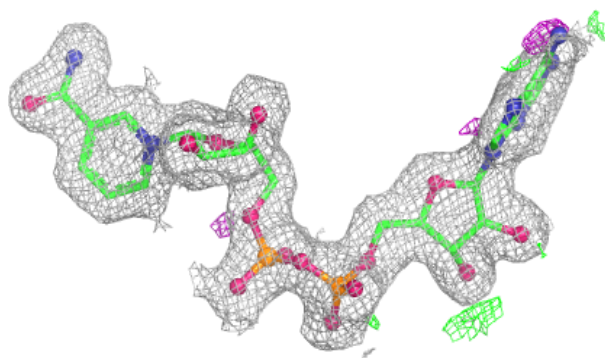


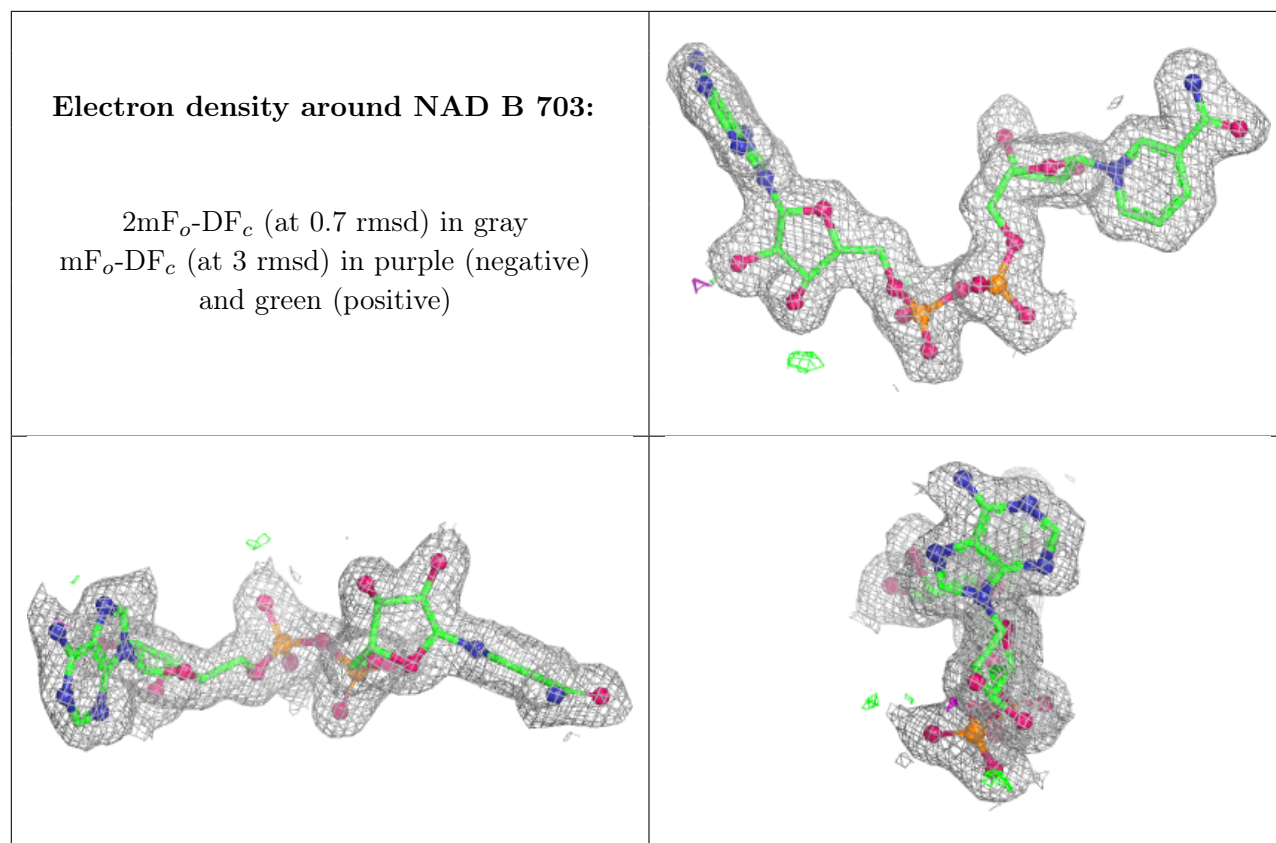
Electron density around NAD D 707:

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

**Electron density around NAD A 701:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.