



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 30, 2014 – 10:48 PM EDT

PDB ID : 4CCQ
Title : Crystal structure of the thioredoxin reductase from *Entamoeba histolytica* with NADP
Authors : Parsonage, D.; Kells, P.M.; Hirata, K.; Debnath, A.; Poole, L.B.; McKerrow, J.H.; Reed, S.L.; Podust, L.M.
Deposited on : 2013-10-24
Resolution : 1.50 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

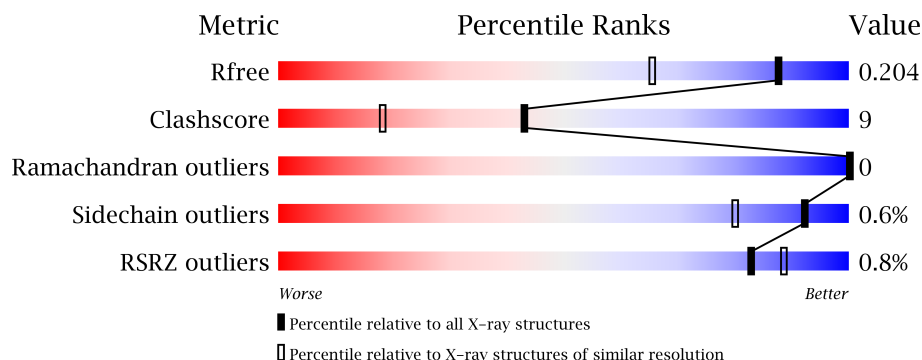
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable24103
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable24103

1 Overall quality at a glance

The reported resolution of this entry is 1.50 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}	66092	1513 (1.50-1.50)
Clashscore	79885	1768 (1.50-1.50)
Ramachandran outliers	78287	1720 (1.50-1.50)
Sidechain outliers	78261	1718 (1.50-1.50)
RSRZ outliers	66119	1514 (1.50-1.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	314	
1	B	314	

2 Entry composition i

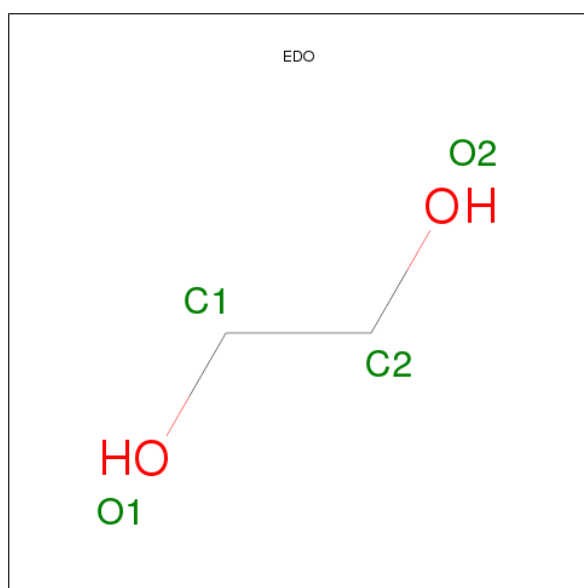
There are 6 unique types of molecules in this entry. The entry contains 6002 atoms, of which 0 are hydrogen and 0 are deuterium.

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 THIOREDOXIN REDUCTASE.

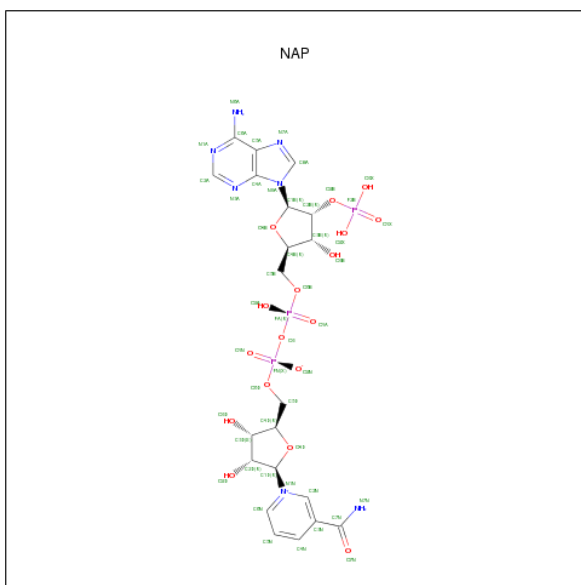
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	10	0
			2422	1529	410	465	18			
1	B	312	Total	C	N	O	S	0	19	0
			2506	1574	435	477	20			

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



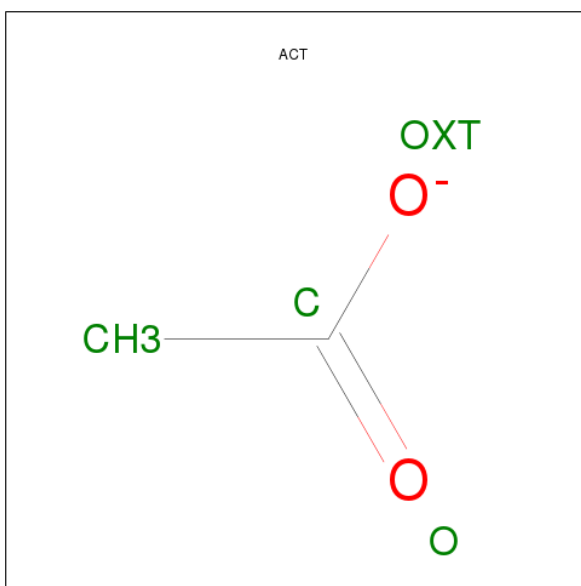
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	B	1	Total 40	C 15	N 6	O 16	P 3	0	0

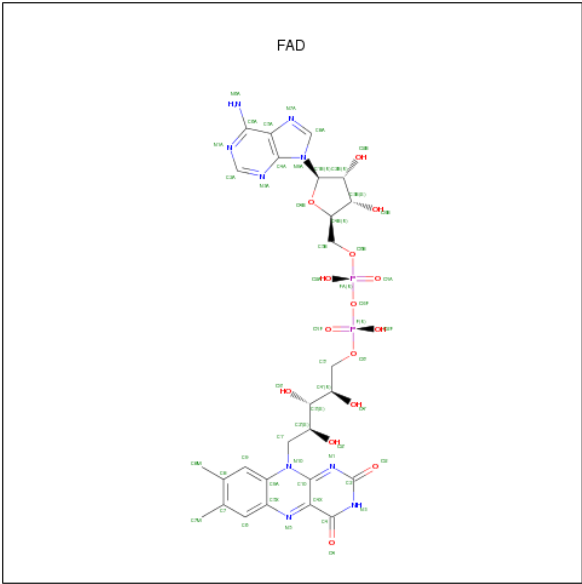
- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:

C₂₇H₃₃N₉O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 6 is water.

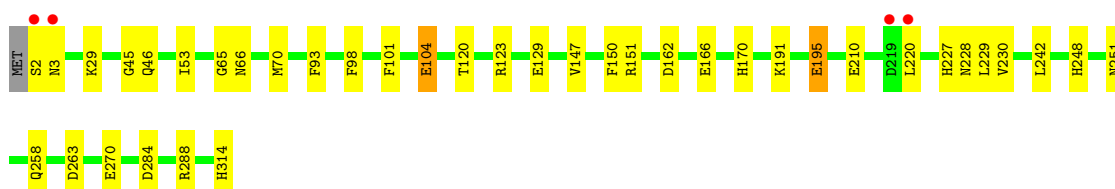
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	454	Total	O	0	0
			454	454		
6	B	410	Total	O	0	0
			410	410		

3 Residue-property plots i

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: THIOREDOXIN REDUCTASE

Chain A: 



• Molecule 1: THIOREDOXIN REDUCTASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.76Å 92.22Å 103.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	103.52 – 1.50 40.67 – 1.50	Depositor EDS
% Data completeness (in resolution range)	95.4 (103.52-1.50) 95.5 (40.67-1.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.172 , 0.207 0.169 , 0.204	Depositor DCC
R_{free} test set	4857 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 96720 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6002	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAP, EDO, FAD, ACT

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	1.25	5/2468 (0.2%)	1.11	6/3341 (0.2%)
1	B	1.17	1/2552 (0.0%)	1.16	8/3450 (0.2%)
All	All	1.21	6/5020 (0.1%)	1.14	14/6791 (0.2%)

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 (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	GLU	CD-OE2	-6.03	1.19	1.25
1	A	98	PHE	CD2-CE2	5.26	1.49	1.39
1	B	202	LYS	CE-NZ	5.26	1.62	1.49
1	A	104	GLU	CB-CG	-5.24	1.42	1.52
1	A	101	PHE	CE1-CZ	5.18	1.47	1.37
1	A	288	ARG	CG-CD	5.13	1.64	1.51

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	263[A]	ASP	CB-CG-OD2	13.93	130.83	118.30
1	B	263[B]	ASP	CB-CG-OD2	13.93	130.83	118.30
1	B	89[A]	ASP	CB-CG-OD1	11.76	128.89	118.30
1	B	89[B]	ASP	CB-CG-OD1	11.76	128.89	118.30
1	B	130	ASP	CB-CG-OD1	6.88	124.49	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	ASP	CB-CG-OD1	6.85	124.47	118.30
1	A	3	ASN	N-CA-C	-6.67	92.99	111.00
1	A	284	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	B	89[A]	ASP	CB-CG-OD2	-5.37	113.47	118.30
1	B	89[B]	ASP	CB-CG-OD2	-5.37	113.47	118.30
1	A	151	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	229[A]	LEU	CA-CB-CG	-5.07	103.63	115.30
1	A	229[B]	LEU	CA-CB-CG	-5.07	103.63	115.30
1	A	162	ASP	CB-CG-OD1	5.02	122.82	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	SER	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2422	0	2376	23	0
1	B	2506	0	2472	66	1
2	A	4	0	6	0	0
2	B	8	0	12	2	0
3	A	48	0	25	1	0
3	B	40	0	19	4	0
4	A	4	0	3	1	0
5	A	53	0	31	2	0
5	B	53	0	31	1	0
6	A	454	0	0	10	5
6	B	410	0	0	17	6
All	All	6002	0	4975	89	7

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (89) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:96[B]:GLN:NE2	1:B:97:PRO:HA	1.25	1.39
1:A:314:HIS:C	6:A:2221:HOH:O	1.92	1.05
1:B:193[A]:MET:HE3	1:B:196[A]:ARG:NH1	1.72	1.04
1:B:96[B]:GLN:NE2	1:B:97:PRO:CA	2.20	1.04
1:B:96[B]:GLN:HE21	1:B:97:PRO:CA	1.70	1.02
1:A:220:LEU:HB2	6:A:2334:HOH:O	1.59	0.99
1:B:291[B]:ARG:NH2	3:B:1315:NAP:O2N	1.98	0.95
1:B:124[A]:MET:HE3	1:B:244:TYR:CE1	2.03	0.93
1:B:193[A]:MET:CE	1:B:196[A]:ARG:NH1	2.31	0.92
1:B:96[B]:GLN:HE22	1:B:97:PRO:HA	1.29	0.87
1:B:248:HIS:ND1	1:B:291[B]:ARG:NE	2.22	0.87
1:B:134:GLN:HG3	6:B:2197:HOH:O	1.73	0.86
1:B:124[A]:MET:HE3	1:B:244:TYR:HE1	1.40	0.83
1:B:248:HIS:CE1	1:B:291[B]:ARG:HE	1.99	0.79
1:B:93:PHE:H	1:B:258:GLN:HE21	1.34	0.76
1:B:120[A]:THR:HG22	6:B:2050:HOH:O	1.85	0.74
1:B:96[B]:GLN:HE21	1:B:97:PRO:HA	0.92	0.74
1:B:193[A]:MET:HE3	1:B:196[A]:ARG:HH11	1.52	0.73
1:B:248:HIS:CE1	1:B:291[B]:ARG:NE	2.56	0.73
1:B:193[A]:MET:CE	1:B:196[A]:ARG:HH11	1.99	0.73
1:B:210:GLU:OE2	1:B:227:HIS:HD2	1.72	0.71
1:B:120[A]:THR:CG2	6:B:2050:HOH:O	2.38	0.71
1:A:120:THR:HB	1:A:251:ASN:HD21	1.55	0.71
1:B:93:PHE:H	1:B:258:GLN:NE2	1.89	0.69
1:B:273:LYS:HE3	6:B:2391:HOH:O	1.92	0.68
1:B:124[A]:MET:CE	1:B:244:TYR:CE1	2.78	0.67
1:B:248:HIS:ND1	1:B:291[B]:ARG:CZ	2.58	0.66
1:B:288[B]:ARG:NH2	6:B:2377:HOH:O	2.29	0.65
1:A:93:PHE:H	1:A:258:GLN:NE2	1.93	0.65
1:B:96[B]:GLN:HE21	1:B:96[B]:GLN:HA	1.62	0.63
1:A:93:PHE:H	1:A:258:GLN:HE21	1.45	0.62
1:A:210:GLU:OE2	1:A:227:HIS:HD2	1.82	0.61
1:B:263[A]:ASP:OD1	6:B:2347:HOH:O	2.16	0.61
1:B:124[A]:MET:CE	1:B:244:TYR:HE1	2.12	0.60
1:B:291[B]:ARG:HD2	6:B:2327:HOH:O	2.00	0.60
1:A:248:HIS:HD2	6:A:2361:HOH:O	1.85	0.58
1:B:70[B]:MET:CE	6:B:2104:HOH:O	2.51	0.58
1:B:50:THR:HG21	1:B:53:ILE:HD11	1.86	0.58
1:A:147:VAL:HG22	1:A:150:PHE:CD2	2.40	0.57
1:B:167[B]:GLU:OE2	6:B:2200:HOH:O	2.17	0.56
1:A:66:ASN:O	1:A:70[A]:MET:HG3	2.05	0.56
1:B:96[B]:GLN:HE21	1:B:96[B]:GLN:CA	2.19	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:104:GLU:HG2	6:A:2152:HOH:O	2.06	0.55
1:B:310:TRP:O	1:B:314:HIS:HD2	1.89	0.55
1:A:53:ILE:HD11	1:A:65:GLY:HA2	1.90	0.54
1:B:248:HIS:CE1	1:B:291[B]:ARG:HH21	2.25	0.53
1:B:227:HIS:HE1	6:B:2306:HOH:O	1.91	0.53
1:B:193[A]:MET:HE3	1:B:196[A]:ARG:HH12	1.65	0.53
1:B:185:ASP:HB3	1:B:208:ASN:HD21	1.74	0.53
1:B:192:THR:O	1:B:196[B]:ARG:HG3	2.09	0.52
1:A:170:HIS:HE1	6:A:2251:HOH:O	1.93	0.52
1:B:248:HIS:CE1	1:B:291[B]:ARG:NH2	2.78	0.52
1:A:46:GLN:HE21	5:A:1318:FAD:HM81	1.75	0.50
1:B:177:LYS:HE2	1:B:204:GLU:OE1	2.10	0.50
1:B:45:GLY:HA2	5:B:1318:FAD:O3B	2.12	0.50
1:B:66:ASN:O	1:B:70[A]:MET:HG3	2.12	0.50
1:A:166:GLU:CG	3:A:1316:NAP:O7N	2.60	0.49
1:B:248:HIS:HE1	1:B:291[B]:ARG:HH21	1.60	0.49
1:B:70[B]:MET:HE1	6:B:2104:HOH:O	2.11	0.49
1:A:191:LYS:O	1:A:195[B]:GLU:HG3	2.12	0.49
1:A:270[A]:GLU:HG2	4:A:1317:ACT:H1	1.95	0.49
1:B:291[B]:ARG:HH22	3:B:1315:NAP:PN	2.36	0.48
1:A:45:GLY:HA2	5:A:1318:FAD:O3B	2.13	0.48
1:B:70[B]:MET:HE3	6:B:2104:HOH:O	2.14	0.47
1:B:153:LYS:HE2	6:B:2212:HOH:O	2.15	0.47
1:B:248:HIS:CE1	1:B:291[B]:ARG:CZ	2.98	0.47
1:B:191[B]:LYS:HG2	1:B:195:GLU:OE2	2.15	0.47
1:B:291[B]:ARG:NH2	3:B:1315:NAP:PN	2.87	0.47
1:B:288[B]:ARG:HD3	6:B:2378:HOH:O	2.15	0.46
1:B:288[B]:ARG:CZ	6:B:2377:HOH:O	2.64	0.46
1:A:228:ASN:OD1	1:A:230:VAL:HG22	2.15	0.45
1:B:96[A]:GLN:NE2	6:B:2155:HOH:O	2.48	0.45
6:A:2039:HOH:O	1:B:192:THR:HB	2.14	0.45
1:A:263:ASP:HB2	6:A:2384:HOH:O	2.17	0.45
1:B:190:SER:OG	1:B:193[B]:MET:HG2	2.17	0.45
1:B:288[B]:ARG:NH1	6:B:2377:HOH:O	2.49	0.45
1:B:132:TYR:CZ	1:B:216:GLY:HA3	2.52	0.45
1:A:120:THR:CG2	6:A:2227:HOH:O	2.64	0.44
1:B:123:ARG:NH2	1:B:130:ASP:OD1	2.51	0.43
1:B:288[A]:ARG:O	1:B:291[A]:ARG:CZ	2.68	0.42
1:A:29:LYS:HE3	6:A:2040:HOH:O	2.18	0.42
1:B:193[A]:MET:HE1	1:B:196[A]:ARG:NH1	2.25	0.42
1:B:70[B]:MET:HE3	1:B:73:ARG:HB3	2.00	0.42
1:B:73:ARG:HD2	2:B:1316:EDO:H11	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:53:ILE:CD1	1:A:65:GLY:HA2	2.49	0.41
1:B:124[A]:MET:HG2	1:B:126:VAL:HG13	2.03	0.41
1:A:120:THR:HG23	6:A:2227:HOH:O	2.20	0.40
1:B:291[A]:ARG:NH2	3:B:1315:NAP:H52N	2.36	0.40
1:B:70[A]:MET:HA	2:B:1316:EDO:H22	2.02	0.40

All (7) 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	Distance(Å)	Clash(Å)
6:B:2034:HOH:O	6:B:2219:HOH:O[4_555]	1.31	0.89
6:A:2155:HOH:O	6:B:2101:HOH:O[3_544]	1.94	0.26
1:B:125:HIS:NE2	6:A:2384:HOH:O[3_554]	2.09	0.11
6:A:2384:HOH:O	6:B:2181:HOH:O[3_544]	2.11	0.09
6:B:2293:HOH:O	6:B:2403:HOH:O[3_554]	2.13	0.07
6:A:2235:HOH:O	6:B:2377:HOH:O[3_544]	2.18	0.02
6:A:2387:HOH:O	6:B:2330:HOH:O[3_544]	2.19	0.01

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	321/314 (102%)	313 (98%)	8 (2%)	0	100	100
1	B	329/314 (105%)	323 (98%)	6 (2%)	0	100	100
All	All	650/628 (104%)	636 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	256/251 (102%)	252 (98%)	4 (2%)	75	44
1	B	267/251 (106%)	267 (100%)	0	100	100
All	All	523/502 (104%)	519 (99%)	4 (1%)	92	73

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	ARG
1	A	195[A]	GLU
1	A	195[B]	GLU
1	A	242	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	61	ASN
1	A	194	GLN
1	A	208	ASN
1	A	227	HIS
1	A	248	HIS
1	A	251	ASN
1	A	258	GLN
1	A	292	GLN
1	B	61	ASN
1	B	66	ASN
1	B	194	GLN
1	B	208	ASN
1	B	227	HIS
1	B	258	GLN
1	B	314	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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	EDO	A	1315	-	3,3,3	0.74	0	2,2,2	1.07	0
3	NAP	A	1316	-	52,52,52	1.48	7 (13%)	80,80,80	1.47	14 (17%)
4	ACT	A	1317	-	1,3,3	1.39	0	0,3,3	0.00	-
5	FAD	A	1318	-	58,58,58	1.51	9 (15%)	85,89,89	1.98	17 (20%)
3	NAP	B	1315	-	42,43,52	1.63	8 (19%)	65,67,80	1.92	13 (20%)
2	EDO	B	1316	-	3,3,3	0.53	0	2,2,2	0.80	0
2	EDO	B	1317	-	3,3,3	0.90	0	2,2,2	0.12	0
5	FAD	B	1318	-	58,58,58	1.39	10 (17%)	85,89,89	1.76	17 (20%)

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	EDO	A	1315	-	-	0/1/1/1	0/0/0/0
3	NAP	A	1316	-	-	0/35/67/67	0/5/5/5
4	ACT	A	1317	-	-	0/0/0/0	0/0/0/0
5	FAD	A	1318	-	-	0/34/50/50	0/6/6/6
3	NAP	B	1315	-	-	0/27/59/67	0/4/4/5
2	EDO	B	1316	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	B	1317	-	-	0/1/1/1	0/0/0/0
5	FAD	B	1318	-	-	0/34/50/50	0/6/6/6

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1316	NAP	C4A-N9A	-4.62	1.31	1.37
3	B	1315	NAP	PN-O2N	4.56	1.59	1.48
5	A	1318	FAD	C4X-N5	4.48	1.40	1.33
3	A	1316	NAP	PN-O2N	4.22	1.58	1.48
5	B	1318	FAD	C4X-N5	4.08	1.40	1.33
5	A	1318	FAD	C1'-N10	3.90	1.52	1.48
3	A	1316	NAP	O4D-C1D	3.77	1.46	1.41
3	B	1315	NAP	C4A-N9A	-3.42	1.32	1.37
3	B	1315	NAP	PN-O3	3.41	1.67	1.60
5	B	1318	FAD	PA-O3P	3.11	1.65	1.59
5	B	1318	FAD	C5X-N5	3.08	1.40	1.35
5	A	1318	FAD	P-O2P	-3.05	1.41	1.55
5	A	1318	FAD	C10-N1	2.99	1.40	1.35
5	A	1318	FAD	C5X-N5	2.87	1.40	1.35
3	B	1315	NAP	C5A-C4A	2.79	1.46	1.40
5	B	1318	FAD	P-O2P	-2.75	1.43	1.55
3	B	1315	NAP	O4D-C1D	2.61	1.46	1.42
5	B	1318	FAD	C2A-N1A	2.59	1.39	1.33
5	A	1318	FAD	C8A-N9A	2.55	1.40	1.36
5	A	1318	FAD	C2A-N3A	2.48	1.36	1.32
5	A	1318	FAD	PA-O2A	-2.47	1.44	1.55
5	B	1318	FAD	C1'-N10	2.42	1.51	1.48
5	B	1318	FAD	C2A-N3A	2.41	1.36	1.32
3	B	1315	NAP	PN-O5D	2.33	1.67	1.59
3	A	1316	NAP	C2N-C3N	2.33	1.42	1.39
5	B	1318	FAD	P-O3P	2.30	1.64	1.59
3	B	1315	NAP	PN-O1N	2.27	1.54	1.48
3	A	1316	NAP	C3N-C7N	2.27	1.54	1.50
5	B	1318	FAD	C10-N1	2.25	1.39	1.35
3	B	1315	NAP	PA-O3	-2.22	1.56	1.59
3	A	1316	NAP	PN-O5D	2.21	1.67	1.59
5	A	1318	FAD	O4B-C1B	2.16	1.43	1.41
3	A	1316	NAP	C4N-C3N	2.15	1.43	1.39
5	B	1318	FAD	C4A-N3A	2.15	1.38	1.35

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1315	NAP	N3A-C2A-N1A	-6.90	122.82	128.89
5	A	1318	FAD	C6A-C5A-C4A	6.86	125.26	117.55
5	B	1318	FAD	C1'-N10-C9A	6.13	125.11	118.67
5	A	1318	FAD	C4A-C5A-N7A	-5.98	103.64	109.41
3	B	1315	NAP	C5A-C4A-N3A	-5.54	120.57	125.98
5	A	1318	FAD	C5A-C4A-N3A	-5.41	120.71	125.98
5	A	1318	FAD	C2-N1-C10	5.23	119.99	114.95
5	B	1318	FAD	N3A-C2A-N1A	-5.12	124.39	128.89
3	A	1316	NAP	C3D-C2D-C1D	4.74	108.35	100.92
3	B	1315	NAP	C4A-C5A-N7A	-4.38	105.17	109.41
5	B	1318	FAD	C5A-C4A-N3A	-4.33	121.76	125.98
5	A	1318	FAD	C5A-C4A-N9A	4.31	113.06	107.09
3	A	1316	NAP	C5A-C4A-N3A	-4.07	122.01	125.98
3	B	1315	NAP	O2N-PN-O3	-4.06	97.66	108.85
3	B	1315	NAP	N3A-C4A-N9A	4.00	132.26	125.39
5	B	1318	FAD	C1'-N10-C10	-3.87	114.61	118.82
5	A	1318	FAD	C1'-N10-C9A	3.81	122.67	118.67
5	A	1318	FAD	C8A-N9A-C4A	-3.77	103.90	106.96
5	A	1318	FAD	P-O3P-PA	3.71	142.19	131.93
3	A	1316	NAP	C4D-O4D-C1D	3.53	113.59	109.72
5	B	1318	FAD	C4'-C3'-C2'	-3.41	105.46	113.52
5	A	1318	FAD	C8A-N9A-C1B	3.38	132.52	126.15
5	B	1318	FAD	P-O3P-PA	3.36	141.25	131.93
5	B	1318	FAD	O2'-C2'-C1'	-3.17	102.33	109.88
5	A	1318	FAD	C10-C4X-N5	-3.13	119.20	122.57
3	B	1315	NAP	O2N-PN-O5D	3.12	120.96	108.19
5	B	1318	FAD	C4A-C5A-N7A	-3.06	106.45	109.41
3	A	1316	NAP	O4D-C1D-N1N	2.90	111.32	108.13
5	A	1318	FAD	O4B-C4B-C3B	-2.90	99.27	105.16
5	B	1318	FAD	C6A-C5A-C4A	2.90	120.80	117.55
3	A	1316	NAP	C6N-N1N-C2N	-2.71	118.25	121.79
5	B	1318	FAD	C5X-C9A-N10	2.65	119.59	117.63
3	B	1315	NAP	N6A-C6A-N1A	2.64	124.72	119.11
5	A	1318	FAD	C4X-N5-C5X	2.59	119.70	116.68
3	A	1316	NAP	N3A-C2A-N1A	-2.55	126.65	128.89
5	A	1318	FAD	O2'-C2'-C1'	-2.55	103.81	109.88
3	A	1316	NAP	O2N-PN-O3	-2.50	101.96	108.85
3	A	1316	NAP	O4D-C1D-C2D	-2.49	103.06	106.69
3	B	1315	NAP	C2A-N1A-C6A	2.46	123.14	118.76
5	B	1318	FAD	O3'-C3'-C4'	-2.41	102.66	108.73
5	B	1318	FAD	C5A-C6A-N6A	2.37	126.08	120.72
5	A	1318	FAD	C4'-C3'-C2'	-2.36	107.95	113.52
3	A	1316	NAP	O5B-C5B-C4B	-2.32	100.45	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1315	NAP	O2A-PA-O1A	2.30	124.86	112.14
5	B	1318	FAD	C5'-C4'-C3'	-2.28	107.75	112.06
3	B	1315	NAP	O2B-C2B-C1B	-2.22	101.44	110.13
5	A	1318	FAD	C8A-N7A-C5A	2.22	110.46	103.58
5	B	1318	FAD	N3A-C4A-N9A	2.21	129.18	125.39
5	B	1318	FAD	O5'-C5'-C4'	-2.19	103.13	109.23
3	B	1315	NAP	C2A-N3A-C4A	2.17	119.50	113.27
3	A	1316	NAP	O2N-PN-O1N	2.15	125.12	118.70
3	A	1316	NAP	N6A-C6A-N1A	2.14	123.65	119.11
3	B	1315	NAP	O3X-P2B-O2B	-2.12	100.97	107.09
3	A	1316	NAP	O2A-PA-O3	2.11	115.15	105.14
3	A	1316	NAP	O7N-C7N-N7N	-2.10	119.59	122.59
5	B	1318	FAD	O2B-C2B-C1B	-2.08	104.90	111.49
3	B	1315	NAP	O2A-PA-O3	2.07	114.97	105.14
5	A	1318	FAD	C5'-C4'-C3'	-2.04	108.21	112.06
5	B	1318	FAD	O3B-C3B-C4B	2.04	117.07	111.07
3	A	1316	NAP	N3A-C4A-N9A	2.04	128.89	125.39
5	A	1318	FAD	C1B-N9A-C4A	-2.01	123.16	126.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	313/314 (99%)	-0.36	4 (1%) 74 81	9, 16, 28, 41	0
1	B	312/314 (99%)	-0.36	1 (0%) 91 94	10, 16, 30, 38	0
All	All	625/628 (99%)	-0.36	5 (0%) 83 88	9, 16, 30, 41	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	SER	4.1
1	A	3	ASN	2.8
1	A	220	LEU	2.8
1	B	36	PHE	2.4
1	A	219	ASP	2.2

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	EDO	B	1316	4/4	0.12	1.73	32,36,37,43	0
2	EDO	A	1315	4/4	0.10	1.64	24,24,24,25	0
2	EDO	B	1317	4/4	0.11	0.97	37,38,38,40	0
5	FAD	B	1318	53/53	0.10	0.78	14,18,26,28	0
4	ACT	A	1317	4/4	0.09	0.31	47,48,48,48	0
3	NAP	A	1316	48/48	0.07	0.11	11,18,50,58	0
3	NAP	B	1315	40/48	0.08	-0.17	14,20,52,52	0
5	FAD	A	1318	53/53	0.06	-0.63	10,14,21,22	0

6.5 Other polymers ⓘ

There are no such residues in this entry.