



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:34 pm GMT

PDB ID : 1WPQ
Title : Ternary Complex Of Glycerol 3-phosphate Dehydrogenase 1 with NAD and dihydroxyactone
Authors : Ou, X.; Han, X.; Rao, Z.
Deposited on : 2004-09-10
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

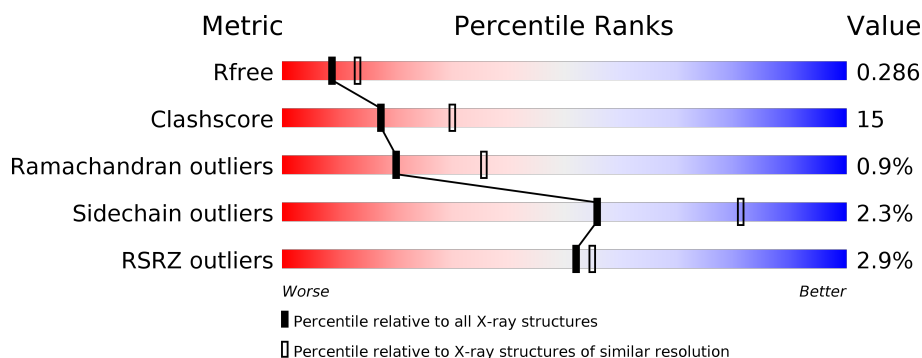
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	349	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>23%</div> <div>.</div> </div> </div>
1	B	349	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>26%</div> <div>.</div> </div> </div>

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
2	SO4	A	1001	-	-	-	X
2	SO4	A	1002	-	-	-	X
3	13P	A	2002	-	-	X	X
3	13P	B	2001	-	-	X	X

2 Entry composition [i](#)

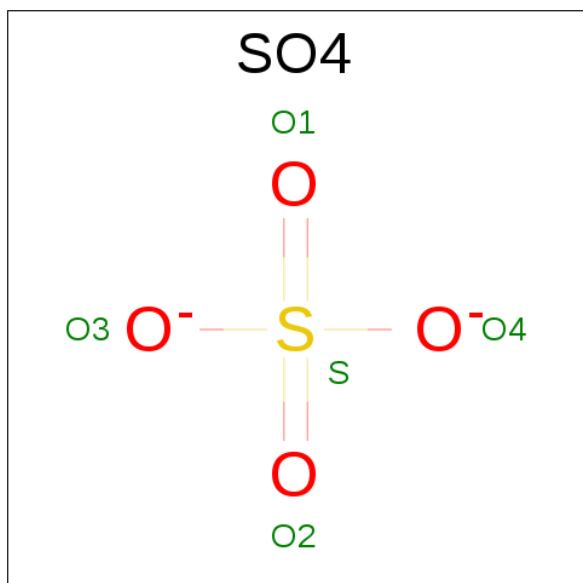
There are 5 unique types of molecules in this entry. The entry contains 5993 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycerol-3-phosphate dehydrogenase [NAD⁺], cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	0	0
			2624	1670	445	490	19			
1	B	348	Total	C	N	O	S	0	0	0
			2624	1670	445	490	19			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



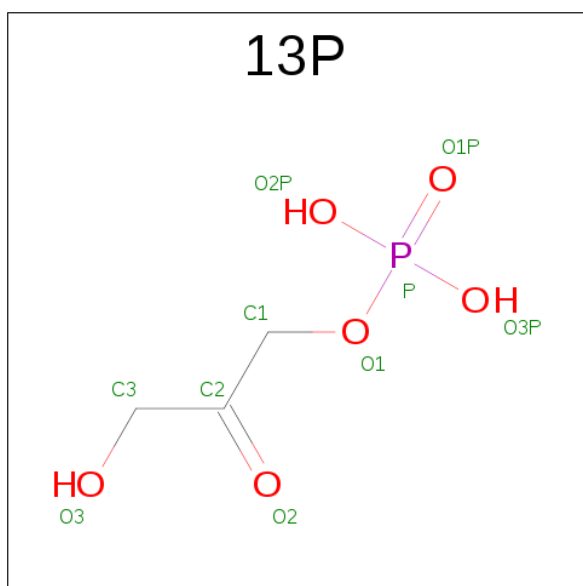
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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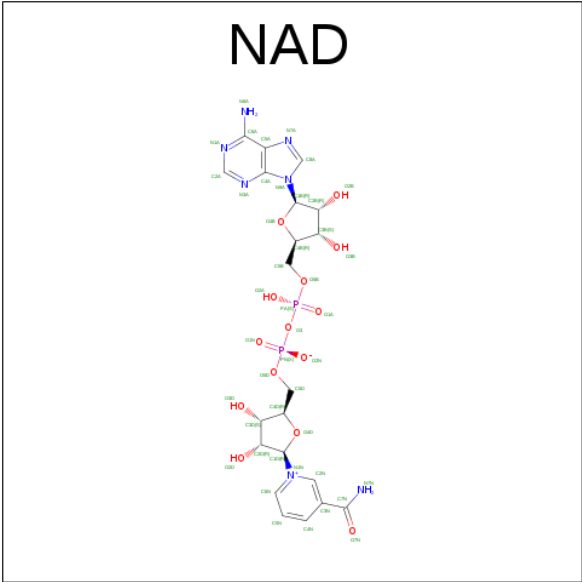
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1,3-DIHYDROXYACETONEPHOSPHATE (three-letter code: 13P) (formula: $C_3H_7O_6P$).



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

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

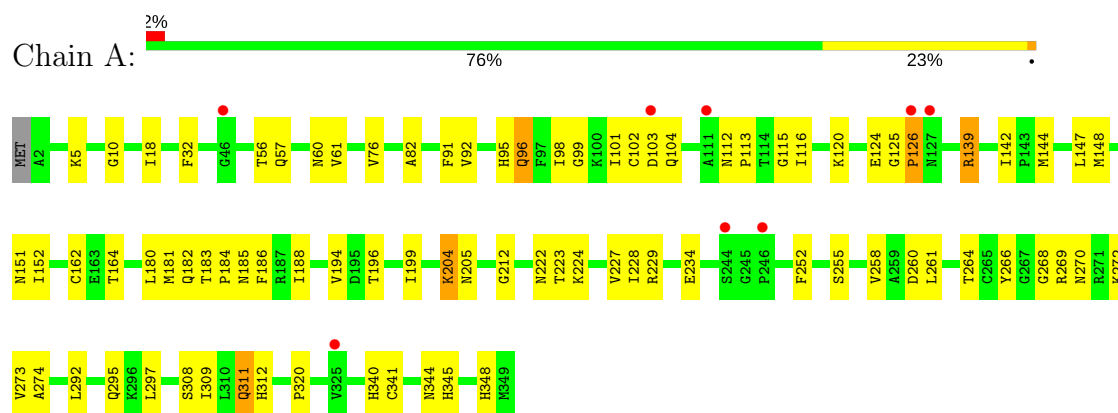
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	322	Total	O	0	0
			322	322		
5	B	280	Total	O	0	0
			280	280		

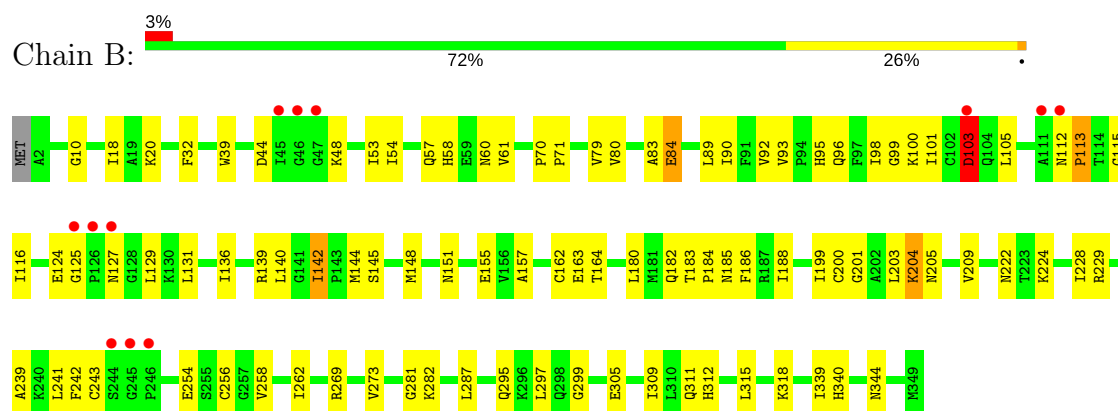
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Glycerol-3-phosphate dehydrogenase [NAD+], cytoplasmic



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4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	116.00Å 116.00Å 153.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 49.15 – 2.46	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 93.7 (49.15-2.46)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.45Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.258 , 0.294 0.256 , 0.286	Depositor DCC
R_{free} test set	1745 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 23.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5993	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 81.80 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.2580e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 13P, SO4, 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	0/2670	0.62	1/3608 (0.0%)
1	B	0.37	0/2670	0.63	1/3608 (0.0%)
All	All	0.36	0/5340	0.62	2/7216 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	113	PRO	N-CA-CB	5.71	110.15	103.30
1	A	113	PRO	N-CA-CB	5.23	109.58	103.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2624	0	2662	84	0
1	B	2624	0	2662	89	0
2	A	30	0	0	1	0
2	B	5	0	0	0	0
3	A	10	0	5	9	0
3	B	10	0	5	5	0
4	A	44	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	44	0	26	1	0
5	A	322	0	0	11	0
5	B	280	0	0	16	0
All	All	5993	0	5385	166	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 15.

All (166) 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:100:LYS:O	1:B:103:ASP:HB2	1.49	1.12
1:A:204:LYS:HZ1	3:A:2002:13P:H32	1.16	1.09
1:A:269:ARG:HH11	1:A:295:GLN:NE2	1.58	0.99
1:A:183:THR:HG22	1:A:185:ASN:H	1.26	0.99
1:B:269:ARG:HH11	1:B:295:GLN:HE21	1.14	0.95
1:A:269:ARG:HH11	1:A:295:GLN:HE21	1.05	0.93
1:A:96:GLN:HE21	1:A:96:GLN:H	1.17	0.90
1:A:204:LYS:NZ	3:A:2002:13P:H32	1.88	0.87
1:A:139:ARG:HG2	1:A:139:ARG:HH11	1.44	0.82
1:B:269:ARG:HH11	1:B:295:GLN:NE2	1.77	0.81
1:B:183:THR:HG22	1:B:185:ASN:H	1.46	0.81
1:A:204:LYS:HZ1	3:A:2002:13P:C3	1.97	0.77
1:B:79:VAL:HG22	1:B:105:LEU:HD21	1.66	0.77
1:A:223:THR:O	1:A:227:VAL:HG23	1.89	0.71
1:B:142:ILE:HD11	1:B:144:MET:HE3	1.73	0.70
1:A:103:ASP:HB2	1:A:139:ARG:NH2	2.07	0.69
1:B:124:GLU:HG3	5:B:3124:HOH:O	1.93	0.68
1:A:152:ILE:HD13	5:A:3197:HOH:O	1.94	0.67
1:A:124:GLU:HG3	1:A:309:ILE:HG12	1.76	0.67
1:A:212:GLY:HA3	1:A:274:ALA:HB3	1.75	0.67
1:A:269:ARG:NH1	1:A:295:GLN:HE21	1.84	0.67
1:A:124:GLU:OE2	1:A:312:HIS:HD2	1.78	0.66
1:B:124:GLU:OE2	1:B:312:HIS:HD2	1.78	0.66
1:A:297:LEU:HB2	5:A:3070:HOH:O	1.96	0.66
1:B:182:GLN:HE21	1:B:188:ILE:H	1.43	0.65
1:B:269:ARG:HB2	3:B:2001:13P:O2P	1.96	0.65
1:B:115:GLY:HA3	1:B:144:MET:HE1	1.77	0.65
1:B:129:LEU:HD21	1:B:131:LEU:HD23	1.79	0.64
1:A:264:THR:HG23	3:A:2002:13P:O1P	1.97	0.64
1:B:115:GLY:HA3	1:B:144:MET:CE	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:THR:HB	1:A:186:PHE:HB3	1.81	0.62
1:A:164:THR:HG22	1:A:188:ILE:HG12	1.82	0.62
1:A:182:GLN:NE2	5:A:3002:HOH:O	2.32	0.61
1:A:102:CYS:HB3	1:A:139:ARG:HD3	1.83	0.61
1:B:32:PHE:CZ	1:B:180:LEU:HB2	2.36	0.60
1:A:96:GLN:NE2	1:A:96:GLN:H	1.96	0.60
1:B:339:ILE:HB	5:B:3149:HOH:O	2.01	0.60
1:B:140:LEU:HB2	1:B:142:ILE:HD12	1.84	0.59
1:B:163:GLU:HG3	5:B:3006:HOH:O	2.02	0.59
1:A:340:HIS:CE1	1:A:344:ASN:ND2	2.71	0.59
1:A:183:THR:HG23	1:A:184:PRO:HD2	1.84	0.58
1:B:241:LEU:HD22	1:B:242:PHE:CE1	2.39	0.58
1:B:199:ILE:HG12	1:B:203:LEU:CD1	2.34	0.58
1:B:183:THR:HB	1:B:186:PHE:HB3	1.86	0.57
1:B:99:GLY:O	5:B:3050:HOH:O	2.17	0.57
1:B:269:ARG:NH1	1:B:295:GLN:HB3	2.19	0.57
1:A:228:ILE:HG23	1:A:261:LEU:HG	1.87	0.57
1:A:139:ARG:HG2	1:A:139:ARG:NH1	2.18	0.56
1:B:142:ILE:HD13	1:B:142:ILE:H	1.71	0.56
1:A:32:PHE:CZ	1:A:180:LEU:HB2	2.40	0.56
1:A:348:HIS:HE1	1:B:254:GLU:OE2	1.89	0.55
1:B:183:THR:HG23	1:B:184:PRO:HD2	1.87	0.55
1:A:196:THR:O	1:A:199:ILE:HG22	2.07	0.55
1:B:340:HIS:CE1	1:B:344:ASN:HD21	2.26	0.54
1:B:124:GLU:OE2	1:B:312:HIS:CD2	2.60	0.54
1:B:10:GLY:HA2	1:B:93:VAL:HG12	1.90	0.54
1:B:98:ILE:HG12	1:B:136:ILE:HD11	1.90	0.53
1:B:281:GLY:HA2	5:B:3134:HOH:O	2.08	0.53
1:B:241:LEU:HD22	1:B:242:PHE:CD1	2.43	0.53
1:A:340:HIS:CE1	1:A:344:ASN:HD21	2.27	0.53
1:B:183:THR:HG21	5:B:3125:HOH:O	2.08	0.53
1:A:120:LYS:HD2	3:A:2002:13P:H31	1.91	0.53
1:A:269:ARG:NH1	1:A:295:GLN:HB3	2.24	0.52
1:B:53:ILE:HG23	1:B:57:GLN:OE1	2.10	0.52
1:B:142:ILE:CD1	1:B:144:MET:HE3	2.40	0.52
1:A:234:GLU:OE1	1:A:320:PRO:HD2	2.10	0.52
1:B:204:LYS:HZ1	3:B:2001:13P:H31	1.76	0.51
1:B:95:HIS:O	1:B:98:ILE:HG22	2.10	0.51
1:A:273:VAL:HG21	1:A:297:LEU:HD11	1.93	0.51
1:A:116:ILE:HD11	1:A:147:LEU:HB2	1.92	0.51
1:B:54:ILE:O	1:B:58:HIS:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:GLU:HG3	1:A:309:ILE:CG1	2.40	0.51
1:A:204:LYS:HD2	1:A:205:ASN:N	2.26	0.51
1:A:152:ILE:N	1:B:222:ASN:HD21	2.09	0.51
1:A:152:ILE:HG12	1:B:222:ASN:OD1	2.11	0.50
1:A:292:LEU:HB3	1:A:295:GLN:HB2	1.92	0.50
1:A:124:GLU:OE2	1:A:312:HIS:CD2	2.62	0.49
1:A:252:PHE:O	1:A:258:VAL:HG23	2.12	0.49
1:A:115:GLY:HA3	1:A:144:MET:HE1	1.94	0.49
1:A:269:ARG:NH1	1:A:295:GLN:NE2	2.43	0.49
1:A:204:LYS:HD2	1:A:204:LYS:C	2.34	0.49
1:A:120:LYS:HE2	1:A:148:MET:SD	2.52	0.48
1:A:229:ARG:HD3	1:B:163:GLU:OE2	2.13	0.48
1:B:224:LYS:O	1:B:228:ILE:HG13	2.13	0.48
1:A:152:ILE:HA	5:A:3197:HOH:O	2.14	0.48
1:A:204:LYS:CE	3:A:2002:13P:H32	2.43	0.48
1:A:10:GLY:HA3	1:A:92:VAL:O	2.13	0.48
1:B:273:VAL:HG11	1:B:297:LEU:HD11	1.94	0.48
1:A:151:ASN:O	4:A:3001:NAD:H4N	2.14	0.48
1:B:201:GLY:O	1:B:204:LYS:HG3	2.14	0.48
1:A:120:LYS:HA	5:A:3120:HOH:O	2.13	0.47
1:B:164:THR:HG22	1:B:188:ILE:HG12	1.96	0.47
1:B:48:LYS:HB2	1:B:53:ILE:HD11	1.95	0.47
1:A:268:GLY:HA3	3:A:2002:13P:O1P	2.14	0.47
1:A:269:ARG:HG3	3:A:2002:13P:O2P	2.14	0.47
1:B:61:VAL:HG22	5:B:3096:HOH:O	2.13	0.47
1:B:84:GLU:HG2	5:B:3048:HOH:O	2.14	0.47
1:A:229:ARG:NH1	1:B:163:GLU:OE2	2.47	0.47
1:B:148:MET:HB3	5:B:3007:HOH:O	2.14	0.47
1:B:90:ILE:N	1:B:90:ILE:HD12	2.30	0.47
1:B:125:GLY:C	1:B:127:ASN:N	2.68	0.47
1:A:224:LYS:HD2	1:A:266:TYR:CZ	2.50	0.46
1:A:5:LYS:HB3	5:A:3199:HOH:O	2.14	0.46
1:B:269:ARG:HD3	1:B:295:GLN:NE2	2.29	0.46
1:A:103:ASP:HB2	1:A:139:ARG:CZ	2.45	0.46
1:B:151:ASN:O	4:B:3002:NAD:H4N	2.14	0.46
1:A:272:LYS:HE2	2:A:1003:SO4:O2	2.16	0.46
1:A:142:ILE:HD11	1:A:144:MET:CE	2.46	0.46
1:A:255:SER:N	1:B:229:ARG:HE	2.14	0.46
1:B:204:LYS:NZ	3:B:2001:13P:H31	2.31	0.46
1:B:269:ARG:H	3:B:2001:13P:P	2.39	0.46
1:B:305:GLU:O	1:B:309:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:THR:HG23	1:A:184:PRO:CD	2.47	0.45
1:B:44:ASP:HB2	5:B:3240:HOH:O	2.14	0.45
1:B:140:LEU:HB2	1:B:142:ILE:CD1	2.46	0.45
1:B:79:VAL:HG13	1:B:80:VAL:N	2.31	0.45
1:A:61:VAL:HG22	5:A:3210:HOH:O	2.16	0.45
1:B:125:GLY:O	1:B:127:ASN:N	2.50	0.45
1:A:125:GLY:O	1:A:126:PRO:C	2.55	0.45
1:B:315:LEU:HD22	1:B:318:LYS:HG3	1.99	0.45
1:A:18:ILE:HG23	5:A:3043:HOH:O	2.16	0.45
1:A:92:VAL:O	1:A:92:VAL:HG13	2.17	0.44
1:B:157:ALA:HA	5:B:3065:HOH:O	2.18	0.44
1:B:199:ILE:HG12	1:B:203:LEU:HD12	2.00	0.44
1:B:39:TRP:CZ3	1:B:101:ILE:HD11	2.52	0.44
1:A:162:CYS:HB2	5:A:3036:HOH:O	2.17	0.44
1:A:120:LYS:HE3	4:A:3001:NAD:H6N	1.99	0.44
1:A:183:THR:HB	1:A:186:PHE:CB	2.47	0.43
1:A:56:THR:HG22	1:A:57:GLN:NE2	2.33	0.43
1:B:239:ALA:O	1:B:243:CYS:HB2	2.17	0.43
1:A:270:ASN:ND2	3:A:2002:13P:O3P	2.52	0.43
1:A:269:ARG:NH2	5:A:3037:HOH:O	2.36	0.43
1:B:162:CYS:HB2	5:B:3008:HOH:O	2.19	0.43
1:A:18:ILE:HD12	1:A:181:MET:HE1	2.01	0.43
1:B:125:GLY:C	1:B:127:ASN:H	2.20	0.43
1:A:91:PHE:HE1	1:A:144:MET:HE1	1.83	0.42
1:A:340:HIS:HB2	5:A:3217:HOH:O	2.18	0.42
1:B:205:ASN:O	1:B:209:VAL:HG23	2.19	0.42
1:B:182:GLN:NE2	1:B:188:ILE:H	2.13	0.42
1:B:205:ASN:HB3	1:B:299:GLY:HA2	2.02	0.42
1:B:100:LYS:HA	5:B:3050:HOH:O	2.18	0.42
1:B:224:LYS:HA	5:B:3118:HOH:O	2.17	0.42
1:A:98:ILE:HG23	1:A:99:GLY:N	2.33	0.42
1:A:101:ILE:O	1:A:104:GLN:HB2	2.20	0.42
1:B:129:LEU:HD21	1:B:131:LEU:CD2	2.46	0.42
1:B:116:ILE:HA	1:B:145:SER:O	2.20	0.42
1:B:83:ALA:HB1	1:B:89:LEU:HD21	2.01	0.42
1:A:222:ASN:ND2	1:B:155:GLU:HB2	2.35	0.41
1:A:204:LYS:HE2	1:A:260:ASP:OD2	2.20	0.41
1:B:200:CYS:HB3	1:B:256:CYS:O	2.19	0.41
1:B:258:VAL:O	1:B:262:ILE:HG13	2.20	0.41
1:A:181:MET:HB2	1:A:188:ILE:CD1	2.51	0.41
1:B:204:LYS:HZ1	3:B:2001:13P:C3	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:PRO:HA	1:B:71:PRO:HD3	1.95	0.41
1:A:308:SER:O	1:A:311:GLN:HG3	2.21	0.41
1:B:151:ASN:N	5:B:3012:HOH:O	2.49	0.41
1:B:311:GLN:HG3	1:B:312:HIS:N	2.35	0.41
1:B:92:VAL:O	1:B:92:VAL:HG22	2.21	0.41
1:A:341:CYS:O	1:A:345:HIS:HB2	2.21	0.41
1:A:76:VAL:HB	1:A:82:ALA:HB2	2.03	0.41
1:B:113:PRO:O	1:B:142:ILE:HB	2.20	0.41
1:B:224:LYS:HB2	1:B:224:LYS:HE3	1.92	0.41
1:A:273:VAL:CB	1:A:297:LEU:HD11	2.51	0.40
1:B:20:LYS:HE3	5:B:3222:HOH:O	2.21	0.40
1:B:241:LEU:C	1:B:241:LEU:HD23	2.42	0.40
1:B:282:LYS:HB2	1:B:287:LEU:HD21	2.03	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	346/349 (99%)	330 (95%)	13 (4%)	3 (1%)	20	36
1	B	346/349 (99%)	320 (92%)	23 (7%)	3 (1%)	20	36
All	All	692/698 (99%)	650 (94%)	36 (5%)	6 (1%)	20	36

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	103	ASP
1	B	112	ASN
1	A	126	PRO
1	B	84	GLU
1	A	112	ASN

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Mol	Chain	Res	Type
1	A	194	VAL

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	282/284 (99%)	276 (98%)	6 (2%)	59	83
1	B	282/284 (99%)	275 (98%)	7 (2%)	53	79
All	All	564/568 (99%)	551 (98%)	13 (2%)	56	81

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	95	HIS
1	A	96	GLN
1	A	139	ARG
1	A	204	LYS
1	A	311	GLN
1	B	18	ILE
1	B	60	ASN
1	B	96	GLN
1	B	103	ASP
1	B	139	ARG
1	B	142	ILE
1	B	204	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	57	GLN
1	A	60	ASN
1	A	81	GLN

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Mol	Chain	Res	Type
1	A	96	GLN
1	A	108	HIS
1	A	151	ASN
1	A	185	ASN
1	A	192	GLN
1	A	295	GLN
1	A	311	GLN
1	A	312	HIS
1	A	344	ASN
1	A	348	HIS
1	B	25	ASN
1	B	60	ASN
1	B	81	GLN
1	B	182	GLN
1	B	192	GLN
1	B	295	GLN
1	B	312	HIS
1	B	340	HIS
1	B	344	ASN
1	B	348	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 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	SO4	A	1001	-	4,4,4	0.77	0	6,6,6	0.15	0
2	SO4	A	1002	-	4,4,4	0.94	0	6,6,6	0.14	0
2	SO4	A	1003	-	4,4,4	0.99	0	6,6,6	0.23	0
2	SO4	A	1004	-	4,4,4	1.09	0	6,6,6	0.20	0
2	SO4	A	1006	-	4,4,4	1.10	0	6,6,6	0.25	0
2	SO4	A	1007	-	4,4,4	0.94	0	6,6,6	0.19	0
3	13P	A	2002	-	9,9,9	3.68	2 (22%)	10,12,12	1.55	2 (20%)
4	NAD	A	3001	-	41,48,48	2.59	11 (26%)	43,73,73	2.53	11 (25%)
2	SO4	B	1005	-	4,4,4	1.01	0	6,6,6	0.23	0
3	13P	B	2001	-	9,9,9	2.32	1 (11%)	10,12,12	1.61	2 (20%)
4	NAD	B	3002	-	41,48,48	2.66	11 (26%)	43,73,73	2.58	11 (25%)

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	SO4	A	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1004	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1006	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1007	-	-	0/0/0/0	0/0/0/0
3	13P	A	2002	-	-	0/7/8/8	0/0/0/0
4	NAD	A	3001	-	-	0/22/62/62	0/5/5/5
2	SO4	B	1005	-	-	0/0/0/0	0/0/0/0
3	13P	B	2001	-	-	0/7/8/8	0/0/0/0
4	NAD	B	3002	-	-	0/22/62/62	0/5/5/5

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2002	13P	C1-C2	-3.10	1.44	1.50
4	A	3001	NAD	O3D-C3D	-2.70	1.36	1.43
4	B	3002	NAD	C2B-C1B	-2.35	1.49	1.53
4	B	3002	NAD	O3D-C3D	-2.19	1.37	1.43
4	A	3001	NAD	PA-O2A	-2.06	1.44	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	3002	NAD	O4D-C1D	2.26	1.44	1.41
4	A	3001	NAD	C2A-N1A	2.48	1.38	1.33
4	B	3002	NAD	C3N-C7N	2.51	1.54	1.50
4	A	3001	NAD	O4D-C1D	2.52	1.44	1.41
4	B	3002	NAD	C2A-N1A	2.59	1.38	1.33
4	A	3001	NAD	C3N-C7N	2.62	1.54	1.50
4	A	3001	NAD	C4A-N3A	3.23	1.40	1.35
4	A	3001	NAD	C2A-N3A	3.29	1.37	1.32
4	B	3002	NAD	C4A-N3A	3.30	1.40	1.35
4	B	3002	NAD	C2A-N3A	3.65	1.38	1.32
4	A	3001	NAD	C6N-N1N	4.65	1.47	1.35
4	B	3002	NAD	C6N-C5N	4.67	1.48	1.38
4	A	3001	NAD	C6N-C5N	4.88	1.49	1.38
4	B	3002	NAD	C6N-N1N	5.33	1.49	1.35
3	B	2001	13P	O1-C1	6.31	1.47	1.43
4	A	3001	NAD	C5N-C4N	6.78	1.51	1.38
4	B	3002	NAD	C5N-C4N	7.12	1.52	1.38
3	A	2002	13P	O1-C1	10.11	1.49	1.43
4	A	3001	NAD	C4N-C3N	10.28	1.56	1.39
4	B	3002	NAD	C4N-C3N	10.59	1.56	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	3002	NAD	N3A-C2A-N1A	-11.73	118.64	128.86
4	A	3001	NAD	N3A-C2A-N1A	-11.57	118.78	128.86
4	B	3002	NAD	O7N-C7N-N7N	-3.80	117.18	122.58
4	A	3001	NAD	C5N-C4N-C3N	-3.59	116.13	120.35
4	B	3002	NAD	C5N-C4N-C3N	-3.34	116.42	120.35
4	A	3001	NAD	O7N-C7N-N7N	-3.17	118.08	122.58
4	B	3002	NAD	C1B-N9A-C4A	-2.58	122.18	126.64
4	B	3002	NAD	C3N-C2N-N1N	-2.29	118.12	120.43
4	B	3002	NAD	O2B-C2B-C1B	-2.28	104.48	111.61
4	A	3001	NAD	C1B-N9A-C4A	-2.21	122.82	126.64
4	A	3001	NAD	O4D-C4D-C5D	-2.19	102.01	109.40
4	A	3001	NAD	C3N-C2N-N1N	-2.18	118.24	120.43
4	B	3002	NAD	C2D-C3D-C4D	2.25	107.00	102.62
4	B	3002	NAD	O2A-PA-O1A	2.28	124.07	112.28
4	A	3001	NAD	O2A-PA-O1A	2.32	124.28	112.28
4	A	3001	NAD	C2D-C3D-C4D	2.37	107.24	102.62
3	B	2001	13P	O2-C2-C1	2.52	124.51	120.52
3	A	2002	13P	O2P-P-O1P	2.55	120.50	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	3001	NAD	O5B-PA-O1A	2.61	119.77	109.25
3	B	2001	13P	O2P-P-O1P	2.62	120.73	110.50
4	B	3002	NAD	O5B-PA-O1A	2.77	120.41	109.25
3	A	2002	13P	O2-C2-C1	2.88	125.07	120.52
4	A	3001	NAD	C2N-C3N-C4N	4.51	123.40	118.26
4	B	3002	NAD	C2N-C3N-C4N	4.71	123.63	118.26
4	A	3001	NAD	C3N-C7N-N7N	5.54	124.09	117.77
4	B	3002	NAD	C3N-C7N-N7N	5.84	124.44	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1003	SO4	1	0
3	A	2002	13P	9	0
4	A	3001	NAD	2	0
3	B	2001	13P	5	0
4	B	3002	NAD	1	0

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	348/349 (99%)	0.10	8 (2%) 61 63	28, 28, 28, 28	0
1	B	348/349 (99%)	0.05	12 (3%) 46 48	28, 28, 28, 28	0
All	All	696/698 (99%)	0.07	20 (2%) 52 55	28, 28, 28, 28	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	127	ASN	4.6
1	B	127	ASN	4.3
1	B	46	GLY	3.7
1	B	111	ALA	3.5
1	B	103	ASP	3.4
1	A	103	ASP	3.4
1	A	126	PRO	3.3
1	A	244	SER	3.2
1	B	246	PRO	3.1
1	B	126	PRO	3.0
1	B	112	ASN	3.0
1	A	111	ALA	2.8
1	B	45	ILE	2.7
1	B	47	GLY	2.6
1	A	246	PRO	2.5
1	B	244	SER	2.4
1	A	325	VAL	2.2
1	A	46	GLY	2.1
1	B	125	GLY	2.1
1	B	245	GLY	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	13P	B	2001	10/10	0.88	0.27	5.68	28,28,28,28	0
2	SO4	A	1001	5/5	0.92	0.19	4.22	28,28,28,28	0
2	SO4	A	1002	5/5	0.91	0.20	3.80	28,28,28,28	0
3	13P	A	2002	10/10	0.92	0.25	2.69	28,28,28,28	0
4	NAD	B	3002	44/44	0.92	0.18	1.67	28,28,28,28	0
4	NAD	A	3001	44/44	0.93	0.18	1.22	28,28,28,28	0
2	SO4	B	1005	5/5	0.95	0.29	-	28,28,28,28	0
2	SO4	A	1003	5/5	0.95	0.25	-	28,28,28,28	0
2	SO4	A	1004	5/5	0.88	0.46	-	28,28,28,28	0
2	SO4	A	1006	5/5	0.84	0.40	-	28,28,28,28	0
2	SO4	A	1007	5/5	0.90	0.39	-	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.