



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:57 pm GMT

PDB ID : 1NQ5
Title : Glyceraldehyde-3-Phosphate Dehydrogenase Mutant With Cys 149 Replaced By Ser Complexed With Nad+
Authors : Didierjean, C.; Corbier, C.; Fatih, M.; Favier, F.; Boschi-Muller, S.; Branlant, G.; Aubry, A.
Deposited on : 2003-01-21
Resolution : 2.11 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	:	trunk28683
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

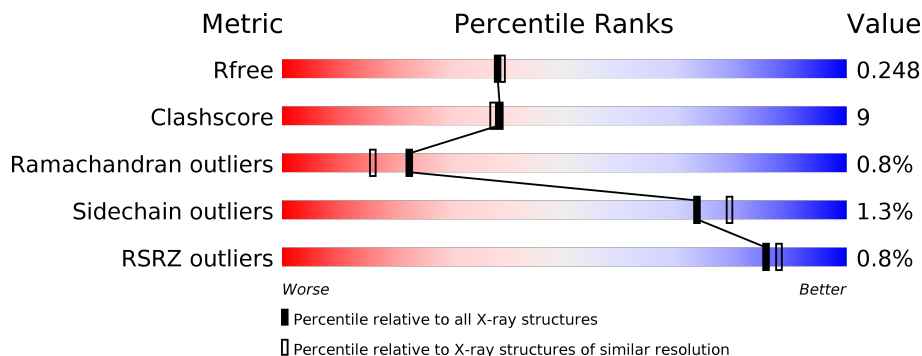
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	4988 (2.14-2.10)
Clashscore	112137	5557 (2.14-2.10)
Ramachandran outliers	110173	5504 (2.14-2.10)
Sidechain outliers	110143	5505 (2.14-2.10)
RSRZ outliers	101464	5021 (2.14-2.10)

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	334	<div> <div style="width: 2%; background-color: red;"></div> <div style="width: 82%; background-color: green;"></div> <div style="width: 17%; background-color: yellow;"></div> <div style="width: 1%; background-color: orange;"></div> </div> <div>82% 17% .</div>
1	C	334	<div> <div style="width: 78%; background-color: green;"></div> <div style="width: 21%; background-color: yellow;"></div> <div style="width: 1%; background-color: orange;"></div> </div> <div>78% 21% .</div>
1	O	334	<div> <div style="width: 2%; background-color: red;"></div> <div style="width: 77%; background-color: green;"></div> <div style="width: 22%; background-color: yellow;"></div> <div style="width: 1%; background-color: orange;"></div> </div> <div>2% 77% 22% .</div>
1	Q	334	<div> <div style="width: 84%; background-color: green;"></div> <div style="width: 16%; background-color: yellow;"></div> </div> <div>84% 16% .</div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10900 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.

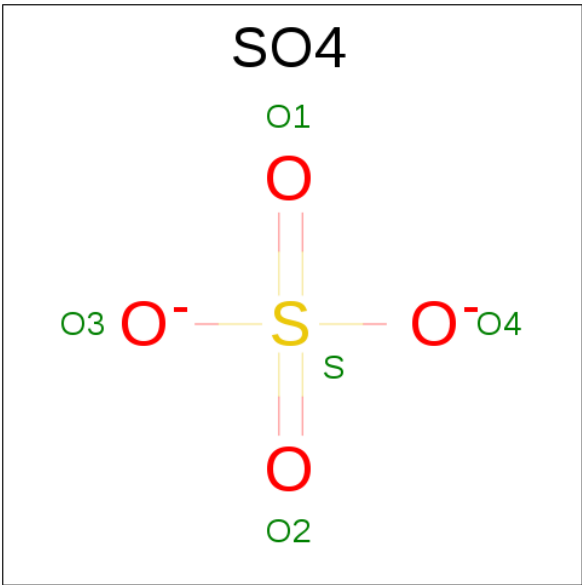
- Molecule 1 is a protein called Glyceraldehyde 3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	O	334	Total	C	N	O	S	0	0	0
			2525	1582	445	490	8			
1	Q	334	Total	C	N	O	S	0	0	0
			2525	1582	445	490	8			
1	A	334	Total	C	N	O	S	0	0	0
			2525	1582	445	490	8			
1	C	334	Total	C	N	O	S	0	0	0
			2525	1582	445	490	8			

There are 4 discrepancies between the modelled and reference sequences:

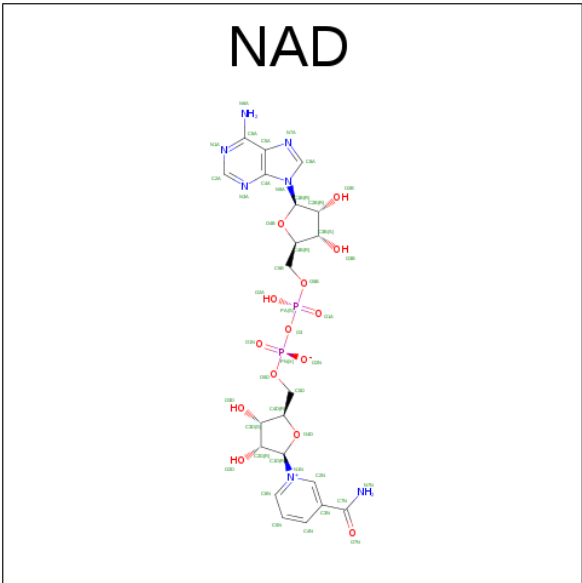
Chain	Residue	Modelled	Actual	Comment	Reference
O	149	SER	CYS	ENGINEERED	UNP P00362
Q	149	SER	CYS	ENGINEERED	UNP P00362
A	149	SER	CYS	ENGINEERED	UNP P00362
C	149	SER	CYS	ENGINEERED	UNP P00362

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



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

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

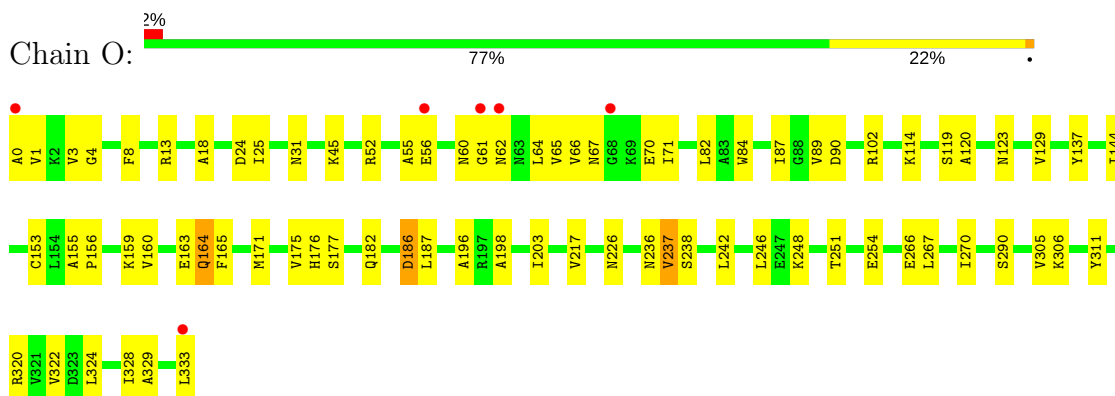
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	154	Total	O	0	0
			154	154		
4	C	159	Total	O	0	0
			159	159		
4	O	130	Total	O	0	0
			130	130		
4	Q	161	Total	O	0	0
			161	161		

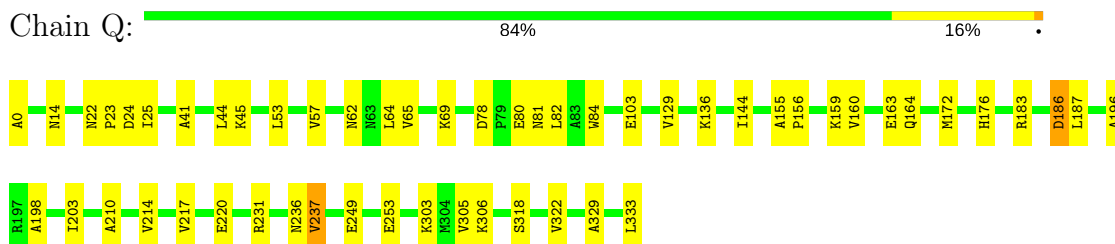
3 Residue-property plots [i](#)

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

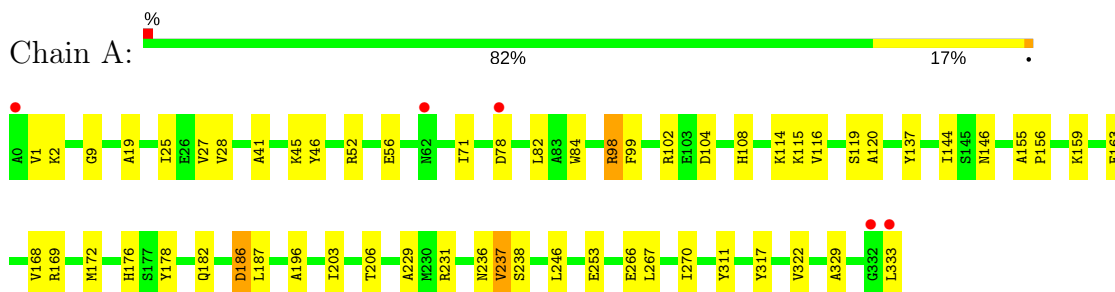
- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase



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A0	V1	G9	N14	R17	K21	D24	I25	L44	K45	H50	L53	E56	V57	M62	M63	L64	E70	I71	I72	V73	D78	L82	A83	M84	I128	V129	M130	G131	Y137	I144	C153	L154	A155	P156	K159	V160	E163	Q164	M171
M172	H176	L185	D186	L187	A196	I203	A210	V214	V217	N226	G227	R231	N236	V237	L246	E249	V252	E253	K260	E264	G265	E266	L267	I270	S290	T297	K303	E314	S318	H319	R320	A326	A329	G332	L333				

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.48Å 87.91Å 119.92Å 90.00° 119.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.11 8.00 – 2.11	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.11) 99.3 (8.00-2.11)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.11Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.198 , 0.249 0.197 , 0.248	Depositor DCC
R_{free} test set	7204 reflections (10.12%)	DCC
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 54.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10900	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.01% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.31	0/2561	0.63	1/3474 (0.0%)
1	C	0.31	0/2561	0.63	1/3474 (0.0%)
1	O	0.30	0/2561	0.62	1/3474 (0.0%)
1	Q	0.30	0/2561	0.62	1/3474 (0.0%)
All	All	0.30	0/10244	0.62	4/13896 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	203	ILE	N-CA-C	-5.68	95.67	111.00
1	C	203	ILE	N-CA-C	-5.60	95.88	111.00
1	A	203	ILE	N-CA-C	-5.25	96.82	111.00
1	O	203	ILE	N-CA-C	-5.20	96.97	111.00

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	2525	0	2572	43	0
1	C	2525	0	2572	52	0
1	O	2525	0	2572	50	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Q	2525	0	2572	34	0
2	A	5	0	0	0	0
2	C	5	0	0	0	0
2	O	5	0	0	0	0
2	Q	5	0	0	0	0
3	A	44	0	26	0	0
3	C	44	0	26	1	0
3	O	44	0	26	0	0
3	Q	44	0	26	0	0
4	A	154	0	0	1	0
4	C	159	0	0	2	0
4	O	130	0	0	3	0
4	Q	161	0	0	6	0
All	All	10900	0	10392	178	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 9.

All (178) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:VAL:HG21	1:A:329:ALA:HB1	1.48	0.94
1:C:329:ALA:HA	1:C:333:LEU:HD11	1.55	0.88
1:A:25:ILE:HD13	1:A:322:VAL:HG13	1.55	0.87
1:O:165:PHE:HA	1:O:248:LYS:HD2	1.58	0.84
1:A:1:VAL:HG21	1:A:329:ALA:CB	2.10	0.81
1:Q:329:ALA:HA	1:Q:333:LEU:HD21	1.63	0.80
1:C:9:GLY:HA3	3:C:7336:NAD:H52A	1.65	0.78
1:O:248:LYS:HE3	4:O:1469:HOH:O	1.86	0.76
1:O:129:VAL:HG23	1:O:217:VAL:HG11	1.68	0.76
1:O:45:LYS:O	1:O:45:LYS:HD3	1.88	0.74
1:O:0:ALA:HB3	1:O:24:ASP:O	1.90	0.72
1:A:98:ARG:HA	1:A:98:ARG:HH11	1.55	0.70
1:C:25:ILE:HD11	1:C:326:ALA:HB2	1.75	0.68
1:Q:25:ILE:HD13	1:Q:322:VAL:HG13	1.76	0.67
1:A:82:LEU:HD13	1:A:84:TRP:CZ2	2.29	0.67
1:C:210:ALA:O	1:C:214:VAL:HG23	1.94	0.67
1:C:329:ALA:HA	1:C:333:LEU:CD1	2.26	0.66
1:C:45:LYS:HE2	1:C:57:VAL:HB	1.78	0.65
1:A:27:VAL:O	1:A:71:ILE:HD12	1.95	0.65
1:A:41:ALA:O	1:A:45:LYS:HG3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:87:ILE:HG13	1:O:89:VAL:HG23	1.78	0.64
1:O:144:ILE:N	1:O:144:ILE:HD12	2.13	0.64
1:C:129:VAL:HG23	1:C:217:VAL:HG11	1.80	0.62
1:C:1:VAL:HG21	1:C:329:ALA:HB1	1.82	0.62
1:Q:82:LEU:HD13	1:Q:84:TRP:CZ2	2.35	0.62
1:O:1:VAL:HG21	1:O:329:ALA:HB1	1.82	0.61
1:A:102:ARG:HG2	1:A:102:ARG:HH11	1.65	0.61
1:O:82:LEU:HD13	1:O:84:TRP:CZ2	2.35	0.60
1:A:187:LEU:O	1:A:196:ALA:HB1	2.00	0.60
1:Q:103:GLU:HG2	4:Q:3390:HOH:O	2.00	0.60
1:C:128:ILE:HD11	1:C:137:TYR:HB2	1.81	0.60
1:A:159:LYS:O	1:A:163:GLU:HG3	2.01	0.60
1:C:187:LEU:O	1:C:196:ALA:HB1	2.02	0.59
1:C:332:GLY:C	1:C:333:LEU:HD12	2.21	0.59
1:C:82:LEU:HD13	1:C:84:TRP:CZ2	2.38	0.59
1:C:62:ASN:O	1:C:73:VAL:HB	2.03	0.59
1:A:19:ALA:HA	1:A:25:ILE:HD11	1.85	0.59
1:Q:187:LEU:O	1:Q:196:ALA:HB1	2.03	0.58
1:O:155:ALA:HB3	1:O:156:PRO:HD3	1.84	0.58
1:Q:155:ALA:HB3	1:Q:156:PRO:HD3	1.86	0.58
1:C:0:ALA:HB3	1:C:24:ASP:O	2.03	0.58
1:O:324:LEU:O	1:O:328:ILE:HG13	2.04	0.57
1:C:44:LEU:HG	1:C:53:LEU:HD22	1.87	0.57
1:O:159:LYS:O	1:O:163:GLU:HG3	2.04	0.57
1:Q:41:ALA:HB2	1:Q:64:LEU:HD22	1.87	0.56
1:O:129:VAL:CG2	1:O:217:VAL:HG11	2.35	0.56
1:O:25:ILE:HD13	1:O:322:VAL:HG13	1.88	0.55
1:O:25:ILE:CD1	1:O:322:VAL:HG13	2.36	0.55
1:A:270:ILE:N	1:A:270:ILE:HD12	2.21	0.55
1:A:102:ARG:CG	1:A:102:ARG:HH11	2.20	0.55
1:A:114:LYS:CB	1:A:333:LEU:HD11	2.37	0.55
1:Q:136:LYS:HE2	4:Q:3406:HOH:O	2.06	0.55
1:C:253:GLU:H	1:C:253:GLU:CD	2.10	0.55
1:Q:253:GLU:H	1:Q:253:GLU:CD	2.10	0.55
1:C:144:ILE:HD12	1:C:144:ILE:N	2.22	0.54
1:O:90:ASP:OD1	1:O:114:LYS:HE2	2.08	0.54
1:A:25:ILE:CD1	1:A:322:VAL:HG13	2.32	0.54
1:A:46:TYR:HE1	1:A:52:ARG:NH1	2.05	0.54
1:A:266:GLU:HG3	1:A:267:LEU:HG	1.91	0.53
1:O:236:ASN:O	1:O:237:VAL:HB	2.08	0.53
1:C:159:LYS:O	1:C:163:GLU:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:160:VAL:O	1:O:164:GLN:HB2	2.07	0.53
1:A:114:LYS:HB2	1:A:333:LEU:HD11	1.90	0.53
1:Q:159:LYS:O	1:Q:163:GLU:HG3	2.08	0.53
1:Q:210:ALA:O	1:Q:214:VAL:HG23	2.09	0.53
1:C:1:VAL:HG21	1:C:329:ALA:CB	2.40	0.52
1:C:236:ASN:O	1:C:237:VAL:HB	2.09	0.52
1:A:2:LYS:HD3	1:A:28:VAL:HG11	1.92	0.52
1:Q:0:ALA:HB3	1:Q:24:ASP:O	2.08	0.52
1:C:266:GLU:HG3	1:C:267:LEU:HG	1.92	0.52
1:C:176:HIS:HB3	1:C:231:ARG:HD3	1.92	0.52
1:Q:41:ALA:HB2	1:Q:64:LEU:CD2	2.40	0.51
1:Q:176:HIS:HB3	1:Q:231:ARG:HD3	1.93	0.51
1:A:119:SER:O	1:A:120:ALA:HB2	2.11	0.50
1:O:90:ASP:HB3	1:O:333:LEU:HD23	1.93	0.50
1:A:155:ALA:HB3	1:A:156:PRO:HD3	1.93	0.50
1:A:98:ARG:HA	1:A:98:ARG:NH1	2.25	0.50
1:A:176:HIS:HB3	1:A:231:ARG:HD3	1.93	0.50
1:O:187:LEU:O	1:O:196:ALA:HB1	2.10	0.50
1:Q:144:ILE:HD12	1:Q:144:ILE:N	2.26	0.50
1:O:266:GLU:HG3	1:O:267:LEU:HG	1.94	0.50
1:C:24:ASP:C	1:C:25:ILE:HD12	2.32	0.49
1:C:128:ILE:CD1	1:C:137:TYR:HB2	2.43	0.49
1:O:186:ASP:HA	1:O:196:ALA:O	2.12	0.48
1:Q:249:GLU:HG3	4:Q:3447:HOH:O	2.13	0.48
1:O:64:LEU:HB2	1:O:71:ILE:HB	1.95	0.48
1:O:56:GLU:O	1:O:66:VAL:HA	2.14	0.48
1:Q:236:ASN:O	1:Q:237:VAL:HB	2.14	0.48
1:Q:78:ASP:OD1	1:Q:80:GLU:HG2	2.13	0.48
1:O:45:LYS:C	1:O:45:LYS:HD3	2.34	0.47
1:Q:25:ILE:CD1	1:Q:322:VAL:HG13	2.44	0.47
1:A:137:TYR:HE1	1:A:144:ILE:HD11	1.79	0.47
1:O:60:ASN:HB2	1:O:65:VAL:CG2	2.45	0.47
1:C:154:LEU:HD23	1:C:214:VAL:HG21	1.97	0.47
1:O:175:VAL:HG23	4:O:1381:HOH:O	2.14	0.47
1:C:186:ASP:HA	1:C:196:ALA:O	2.14	0.47
1:C:155:ALA:HB3	1:C:156:PRO:HD3	1.97	0.47
1:Q:129:VAL:HG23	1:Q:217:VAL:HG11	1.97	0.47
1:Q:80:GLU:HG3	1:Q:81:ASN:ND2	2.31	0.47
1:O:84:TRP:HB3	1:O:89:VAL:HB	1.98	0.46
1:A:102:ARG:CG	1:A:102:ARG:NH1	2.78	0.46
1:A:168:VAL:HG12	1:A:169:ARG:HG2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:PHE:HB3	1:A:104:ASP:HB3	1.98	0.46
1:C:176:HIS:O	1:C:231:ARG:HA	2.16	0.46
1:Q:186:ASP:HA	1:Q:196:ALA:O	2.16	0.46
1:Q:44:LEU:HG	1:Q:53:LEU:HD22	1.98	0.46
1:C:303:LYS:NZ	4:C:7465:HOH:O	2.42	0.45
1:Q:69:LYS:HE3	1:C:56:GLU:HG2	1.97	0.45
1:O:182:GLN:HB3	4:O:1396:HOH:O	2.16	0.45
1:C:154:LEU:HD23	1:C:214:VAL:CG2	2.47	0.45
1:C:17:ARG:HH12	1:C:50:HIS:CG	2.34	0.45
1:O:65:VAL:HG22	1:O:70:GLU:HG3	1.98	0.45
1:C:260:LYS:O	1:C:264:GLU:HG3	2.16	0.45
1:O:251:THR:OG1	1:O:254:GLU:HG3	2.17	0.45
1:A:178:TYR:HA	1:A:182:GLN:OE1	2.17	0.45
1:A:108:HIS:HB2	1:A:116:VAL:HG21	2.00	0.44
1:A:186:ASP:HA	1:A:196:ALA:O	2.17	0.44
1:C:14:ASN:HB3	1:C:318:SER:OG	2.17	0.44
1:O:242:LEU:O	1:O:306:LYS:HA	2.17	0.44
1:Q:62:ASN:O	1:Q:62:ASN:OD1	2.36	0.44
1:C:333:LEU:N	1:C:333:LEU:HD12	2.32	0.44
1:O:3:VAL:HG12	1:O:4:GLY:N	2.32	0.44
1:A:19:ALA:CB	1:A:25:ILE:HD11	2.48	0.44
1:C:185:LEU:O	1:C:186:ASP:C	2.55	0.44
1:C:64:LEU:HB2	1:C:71:ILE:HB	1.98	0.44
1:C:129:VAL:CG2	1:C:217:VAL:HG11	2.47	0.44
1:A:1:VAL:CG2	1:A:329:ALA:HB1	2.34	0.44
1:O:31:ASN:ND2	1:O:82:LEU:HD21	2.32	0.44
1:Q:14:ASN:HB3	1:Q:318:SER:OG	2.18	0.44
1:A:9:GLY:HA2	4:A:5342:HOH:O	2.17	0.43
1:A:253:GLU:CD	1:A:253:GLU:H	2.21	0.43
1:Q:183:ARG:HD2	1:Q:187:LEU:HB2	2.00	0.43
1:C:236:ASN:OD1	1:C:314:GLU:HG3	2.19	0.43
1:C:131:GLY:HA3	1:C:270:ILE:CD1	2.48	0.43
1:A:115:LYS:HG2	1:A:333:LEU:HD11	2.01	0.43
1:O:102:ARG:HD2	1:O:123:ASN:O	2.17	0.43
1:O:18:ALA:HB3	1:O:322:VAL:HG21	2.00	0.43
1:C:249:GLU:OE1	1:C:303:LYS:HE2	2.18	0.43
1:Q:22:ASN:HA	1:Q:23:PRO:HD2	1.94	0.43
1:A:236:ASN:O	1:A:237:VAL:HB	2.18	0.43
1:O:270:ILE:N	1:O:270:ILE:HD12	2.34	0.43
1:O:238:SER:HB2	1:O:311:TYR:CZ	2.54	0.43
1:O:119:SER:O	1:O:120:ALA:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:65:VAL:HG12	4:Q:3368:HOH:O	2.19	0.43
1:O:176:HIS:ND1	1:O:177:SER:O	2.41	0.42
1:Q:160:VAL:O	1:Q:164:GLN:HB2	2.19	0.42
1:C:160:VAL:O	1:C:164:GLN:HB2	2.19	0.42
1:Q:62:ASN:HB2	4:Q:3364:HOH:O	2.20	0.42
1:Q:303:LYS:HE3	4:Q:3447:HOH:O	2.20	0.42
1:A:146:ASN:HB3	1:A:317:TYR:OH	2.20	0.42
1:A:114:LYS:HB3	1:A:333:LEU:HD11	2.01	0.42
1:C:130:MET:O	1:C:270:ILE:HD13	2.20	0.42
1:O:153:CYS:HA	1:O:290:SER:HB2	2.02	0.42
1:A:206:THR:HG23	1:A:229:ALA:HB3	2.01	0.42
1:C:320:ARG:HA	1:C:320:ARG:NE	2.35	0.42
1:O:305:VAL:HG22	1:O:306:LYS:N	2.34	0.42
1:C:236:ASN:O	1:C:237:VAL:CB	2.69	0.41
1:A:238:SER:HB2	1:A:311:TYR:CZ	2.54	0.41
1:Q:45:LYS:HE2	1:Q:57:VAL:HB	2.01	0.41
1:C:153:CYS:HA	1:C:290:SER:HB2	2.02	0.41
1:O:55:ALA:HB1	1:O:67:ASN:ND2	2.35	0.41
1:A:82:LEU:HD13	1:A:84:TRP:CE2	2.55	0.41
1:C:17:ARG:HD2	4:C:7346:HOH:O	2.20	0.41
1:C:154:LEU:CD2	1:C:214:VAL:HG21	2.51	0.41
1:O:329:ALA:HA	1:O:333:LEU:CD1	2.51	0.41
1:C:172:MET:HG3	1:C:227:GLY:HA3	2.03	0.41
1:O:144:ILE:CD1	1:O:144:ILE:N	2.82	0.41
1:A:25:ILE:O	1:A:25:ILE:HG13	2.20	0.41
1:C:252:VAL:HG13	1:C:297:THR:CG2	2.51	0.41
1:C:21:LYS:NZ	1:C:319:HIS:HE1	2.19	0.41
1:O:45:LYS:HE2	1:O:52:ARG:HD3	2.02	0.41
1:Q:305:VAL:HG22	1:Q:306:LYS:N	2.35	0.41
1:O:137:TYR:CZ	1:O:328:ILE:HG23	2.55	0.40
1:O:8:PHE:O	1:O:13:ARG:NH1	2.50	0.40
1:A:236:ASN:O	1:A:237:VAL:CB	2.68	0.40
1:O:320:ARG:NE	1:O:320:ARG:HA	2.35	0.40
1:C:171:MET:HA	1:C:226:ASN:O	2.22	0.40
1:O:171:MET:HA	1:O:226:ASN:O	2.21	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	332/334 (99%)	315 (95%)	15 (4%)	2 (1%)	28	23
1	C	332/334 (99%)	313 (94%)	17 (5%)	2 (1%)	28	23
1	O	332/334 (99%)	316 (95%)	12 (4%)	4 (1%)	15	9
1	Q	332/334 (99%)	313 (94%)	16 (5%)	3 (1%)	20	13
All	All	1328/1336 (99%)	1257 (95%)	60 (4%)	11 (1%)	22	16

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	186	ASP
1	O	237	VAL
1	Q	186	ASP
1	Q	237	VAL
1	A	186	ASP
1	A	237	VAL
1	C	186	ASP
1	C	237	VAL
1	O	198	ALA
1	Q	198	ALA
1	O	61	GLY

5.3.2 Protein sidechains [i](#)

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	272/272 (100%)	267 (98%)	5 (2%)	64	69
1	C	272/272 (100%)	268 (98%)	4 (2%)	70	75
1	O	272/272 (100%)	269 (99%)	3 (1%)	78	82
1	Q	272/272 (100%)	270 (99%)	2 (1%)	87	90
All	All	1088/1088 (100%)	1074 (99%)	14 (1%)	73	79

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	62	ASN
1	O	164	GLN
1	O	246	LEU
1	Q	172	MET
1	Q	220	GLU
1	A	56	GLU
1	A	78	ASP
1	A	98	ARG
1	A	172	MET
1	A	246	LEU
1	C	70	GLU
1	C	78	ASP
1	C	164	GLN
1	C	246	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	O	62	ASN
1	O	63	ASN
1	O	81	ASN
1	Q	38	ASN
1	Q	42	HIS
1	Q	62	ASN
1	Q	63	ASN
1	Q	81	ASN
1	Q	146	ASN
1	Q	319	HIS
1	A	63	ASN
1	A	81	ASN
1	A	146	ASN

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Mol	Chain	Res	Type
1	C	63	ASN
1	C	164	GLN
1	C	319	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](#)

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
3	NAD	A	5336	-	41,48,48	1.09	4 (9%)	43,73,73	1.94	4 (9%)
2	SO4	A	5338	-	4,4,4	0.33	0	6,6,6	0.08	0
3	NAD	C	7336	-	41,48,48	1.14	5 (12%)	43,73,73	2.29	9 (20%)
2	SO4	C	7338	-	4,4,4	0.34	0	6,6,6	0.08	0
3	NAD	O	1336	-	41,48,48	1.04	2 (4%)	43,73,73	2.18	5 (11%)
2	SO4	O	1338	-	4,4,4	0.33	0	6,6,6	0.08	0
3	NAD	Q	3336	-	41,48,48	1.08	4 (9%)	43,73,73	2.16	5 (11%)
2	SO4	Q	3338	-	4,4,4	0.33	0	6,6,6	0.06	0

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
3	NAD	A	5336	-	-	0/22/62/62	0/5/5/5
2	SO4	A	5338	-	-	0/0/0/0	0/0/0/0
3	NAD	C	7336	-	-	0/22/62/62	0/5/5/5
2	SO4	C	7338	-	-	0/0/0/0	0/0/0/0
3	NAD	O	1336	-	-	0/22/62/62	0/5/5/5
2	SO4	O	1338	-	-	0/0/0/0	0/0/0/0
3	NAD	Q	3336	-	-	0/22/62/62	0/5/5/5
2	SO4	Q	3338	-	-	0/0/0/0	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	7336	NAD	C5A-N7A	-2.25	1.31	1.39
3	A	5336	NAD	C5A-N7A	-2.11	1.32	1.39
3	Q	3336	NAD	C5A-N7A	-2.05	1.32	1.39
3	C	7336	NAD	C2A-N3A	2.27	1.36	1.32
3	C	7336	NAD	O4D-C1D	2.30	1.44	1.41
3	C	7336	NAD	O4B-C1B	2.45	1.44	1.41
3	Q	3336	NAD	O4D-C1D	2.52	1.44	1.41
3	Q	3336	NAD	C3N-C7N	2.53	1.54	1.50
3	A	5336	NAD	O4B-C1B	2.61	1.44	1.41
3	A	5336	NAD	C6N-N1N	2.72	1.42	1.35
3	Q	3336	NAD	C6N-N1N	2.76	1.42	1.35
3	O	1336	NAD	O4D-C1D	2.84	1.45	1.41
3	C	7336	NAD	C6N-N1N	3.07	1.43	1.35
3	A	5336	NAD	O4D-C1D	3.11	1.45	1.41
3	O	1336	NAD	C6N-N1N	3.14	1.43	1.35

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	3336	NAD	N3A-C2A-N1A	-11.77	118.60	128.86
3	O	1336	NAD	N3A-C2A-N1A	-11.75	118.62	128.86
3	C	7336	NAD	N3A-C2A-N1A	-11.55	118.80	128.86
3	A	5336	NAD	N3A-C2A-N1A	-10.39	119.81	128.86
3	C	7336	NAD	C4B-O4B-C1B	-4.38	105.11	109.77
3	C	7336	NAD	O5B-PA-O1A	-2.88	97.64	109.25
3	O	1336	NAD	C1B-N9A-C4A	-2.43	122.44	126.64
3	C	7336	NAD	C3N-C7N-N7N	-2.40	115.03	117.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	7336	NAD	C4D-O4D-C1D	2.04	111.94	109.77
3	Q	3336	NAD	O2D-C2D-C3D	2.05	118.38	111.83
3	O	1336	NAD	O2D-C2D-C3D	2.08	118.48	111.83
3	Q	3336	NAD	C5N-C4N-C3N	2.27	123.02	120.35
3	A	5336	NAD	C5N-C4N-C3N	2.34	123.11	120.35
3	C	7336	NAD	O4B-C4B-C5B	2.36	117.38	109.40
3	A	5336	NAD	C4A-C5A-N7A	2.52	111.84	109.41
3	C	7336	NAD	C5N-C4N-C3N	2.59	123.39	120.35
3	Q	3336	NAD	C2A-N1A-C6A	2.67	123.44	118.77
3	C	7336	NAD	C2A-N1A-C6A	2.69	123.48	118.77
3	A	5336	NAD	C2A-N1A-C6A	2.70	123.49	118.77
3	O	1336	NAD	C2A-N1A-C6A	2.70	123.49	118.77
3	C	7336	NAD	C4A-C5A-N7A	2.79	112.11	109.41
3	Q	3336	NAD	C4A-C5A-N7A	4.00	113.27	109.41
3	O	1336	NAD	C4A-C5A-N7A	4.19	113.46	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	7336	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 [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	334/334 (100%)	-0.33	5 (1%) 74 77	15, 22, 36, 50	0
1	C	334/334 (100%)	-0.41	0 100 100	14, 23, 34, 42	0
1	O	334/334 (100%)	-0.26	6 (1%) 69 73	11, 24, 41, 51	0
1	Q	334/334 (100%)	-0.36	0 100 100	14, 23, 35, 43	0
All	All	1336/1336 (100%)	-0.34	11 (0%) 86 88	11, 23, 36, 51	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	O	68	GLY	3.8
1	O	0	ALA	3.7
1	A	333	LEU	3.3
1	A	62	ASN	3.2
1	O	61	GLY	3.2
1	A	0	ALA	3.0
1	O	56	GLU	2.8
1	A	78	ASP	2.7
1	O	333	LEU	2.6
1	A	332	GLY	2.6
1	O	62	ASN	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 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NAD	C	7336	44/44	0.97	0.08	-0.55	14,18,22,25	0
3	NAD	Q	3336	44/44	0.97	0.07	-0.61	14,18,20,21	0
3	NAD	O	1336	44/44	0.97	0.07	-0.70	14,19,22,25	0
3	NAD	A	5336	44/44	0.98	0.07	-1.10	14,21,24,25	0
2	SO4	A	5338	5/5	0.87	0.23	-	74,74,75,75	0
2	SO4	Q	3338	5/5	0.66	0.34	-	97,97,98,98	0
2	SO4	C	7338	5/5	0.89	0.23	-	69,69,70,70	0
2	SO4	O	1338	5/5	0.90	0.24	-	60,60,62,62	0

6.5 Other polymers

There are no such residues in this entry.