



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:43 PM GMT

PDB ID : 1YE6
Title : Crystal structure of the Lys-274 to Arg mutant of *Candida tenuis* xylose reductase (AKR2B5) bound to NADP+
Authors : Leitgeb, S.; Petschacher, B.; Wilson, D.K.; Nidetzky, B.
Deposited on : 2004-12-28
Resolution : 2.30 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

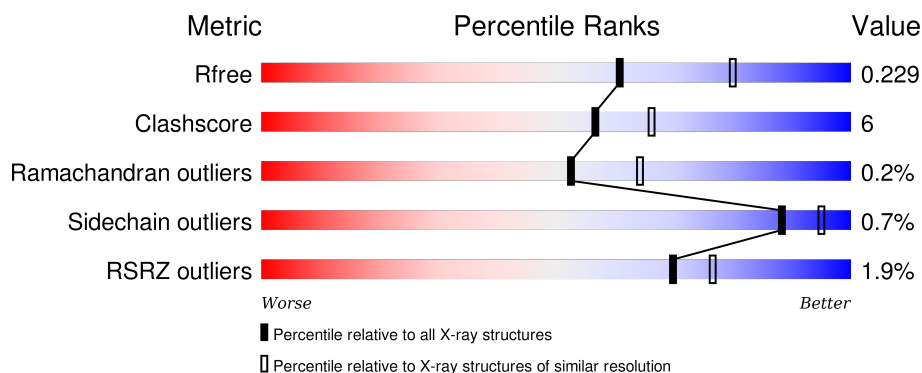
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	322	<div> <div>2%</div> <div>84%</div> <div>15%</div> <div>..</div> </div>
1	B	322	<div> <div>2%</div> <div>85%</div> <div>14%</div> <div>.</div> </div>
1	C	322	<div> <div>2%</div> <div>84%</div> <div>15%</div> <div>..</div> </div>
1	D	322	<div> <div>%</div> <div>88%</div> <div>11%</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	B	901	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11116 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 NAD(P)H-dependent D-xylose reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2533	1645	421	463	4			
1	B	319	Total	C	N	O	S	0	0	0
			2533	1645	421	463	4			
1	C	319	Total	C	N	O	S	0	0	0
			2533	1645	421	463	4			
1	D	319	Total	C	N	O	S	0	0	0
			2533	1645	421	463	4			

There are 4 discrepancies between the modelled and reference sequences:

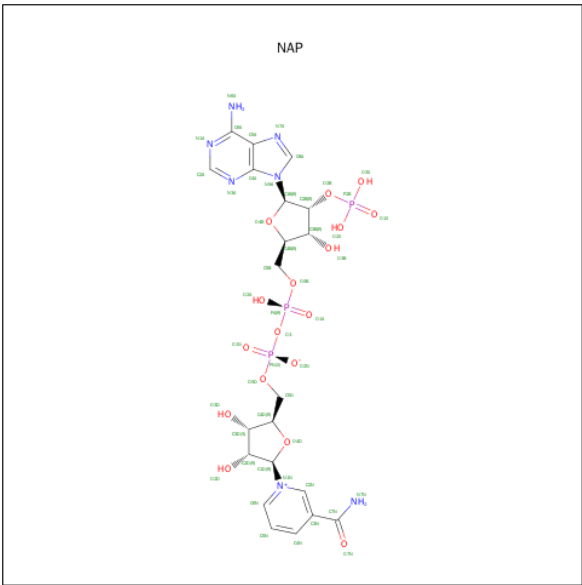
Chain	Residue	Modelled	Actual	Comment	Reference
A	274	ARG	LYS	ENGINEERED	UNP O74237
B	274	ARG	LYS	ENGINEERED	UNP O74237
C	274	ARG	LYS	ENGINEERED	UNP O74237
D	274	ARG	LYS	ENGINEERED	UNP O74237

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



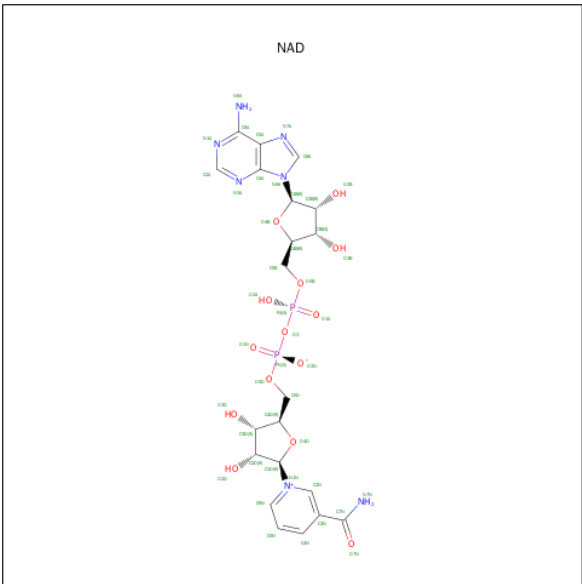
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

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

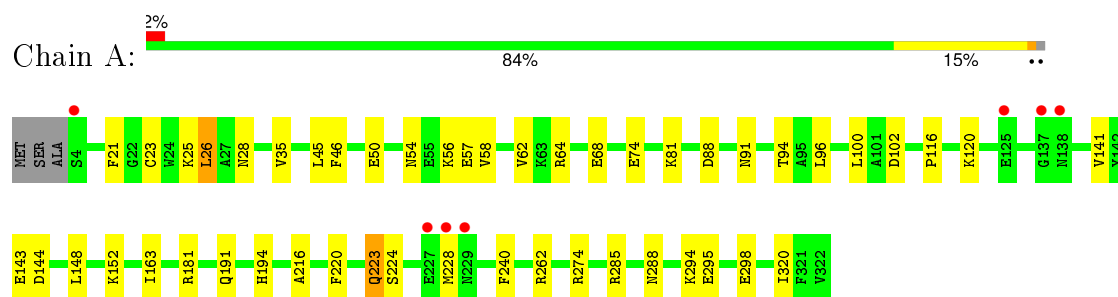
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	182	Total	O	0	0
			182	182		
5	B	167	Total	O	0	0
			167	167		
5	C	187	Total	O	0	0
			187	187		
5	D	224	Total	O	0	0
			224	224		

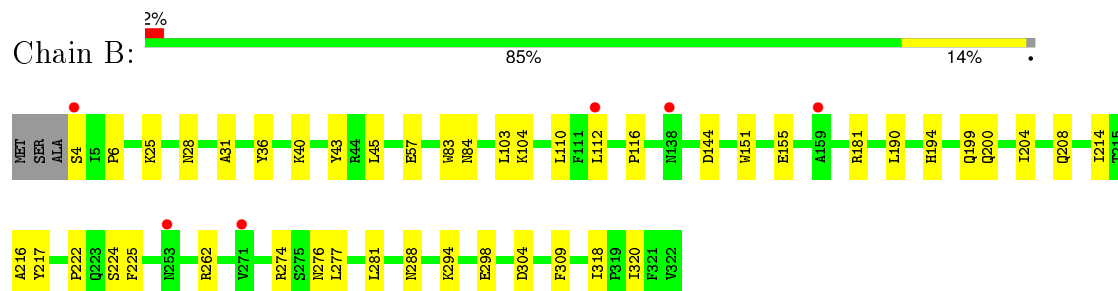
3 Residue-property plots [i](#)

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.

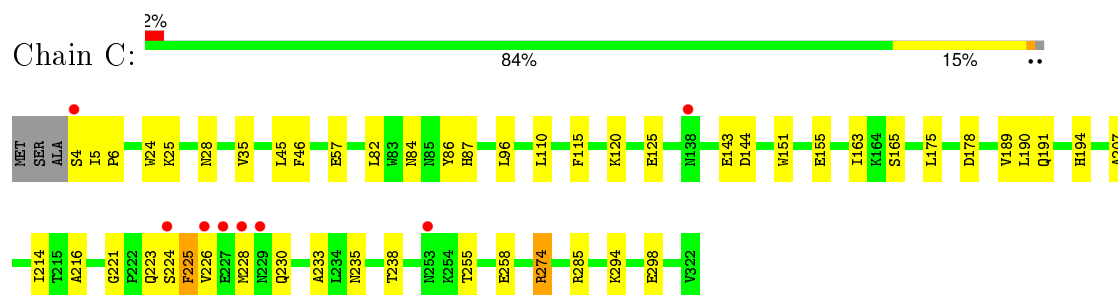
- Molecule 1: NAD(P)H-dependent D-xylose reductase



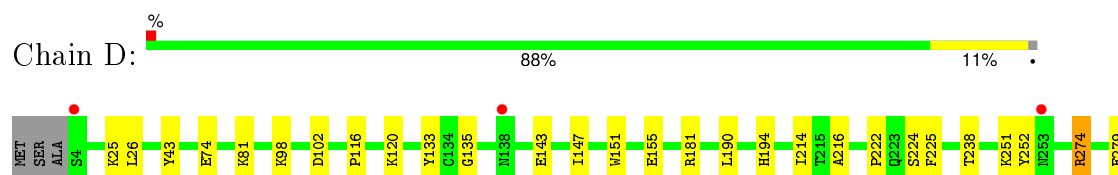
- Molecule 1: NAD(P)H-dependent D-xylose reductase



- Molecule 1: NAD(P)H-dependent D-xylose reductase



- Molecule 1: NAD(P)H-dependent D-xylose reductase



L280	L281	L282	Q283	Q284	R285	R294	E298	F309	T318	P319	I320	F321	Y322
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	180.00Å 127.99Å 80.07Å 90.00° 90.22° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 29.86 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.30) 99.7 (29.86-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.51 (at 2.31Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.192 , 0.229 0.193 , 0.229	Depositor DCC
R_{free} test set	3921 reflections (4.89%)	DCC
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.9	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 80192 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11116	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NAP, 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	0.34	0/2599	0.55	0/3531
1	B	0.34	0/2599	0.57	0/3531
1	C	0.34	0/2599	0.56	0/3531
1	D	0.35	0/2599	0.58	0/3531
All	All	0.34	0/10396	0.57	0/14124

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	217	TYR	Sidechain

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	2533	0	2525	35	0
1	B	2533	0	2525	29	0
1	C	2533	0	2525	40	0
1	D	2533	0	2525	26	0
2	A	5	0	0	0	0
2	B	15	0	0	0	0
2	C	5	0	0	0	0
2	D	15	0	0	0	0
3	A	48	0	25	5	0
3	C	48	0	25	6	0
4	B	44	0	26	1	0
4	D	44	0	26	2	0
5	A	182	0	0	2	0
5	B	167	0	0	1	0
5	C	187	0	0	2	0
5	D	224	0	0	3	1
All	All	11116	0	10202	126	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (126) 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:224:SER:HA	3:A:1000:NAP:H51A	1.52	0.91
1:C:224:SER:HA	3:C:1002:NAP:H51A	1.55	0.85
1:A:223:GLN:HG2	5:A:1092:HOH:O	1.76	0.84
1:C:224:SER:O	3:C:1002:NAP:H52A	1.79	0.83
1:C:224:SER:O	1:C:225:PHE:HB2	1.83	0.78
1:C:294:LYS:O	1:C:298:GLU:HG3	1.88	0.73
1:C:224:SER:O	1:C:225:PHE:CB	2.39	0.70
1:B:194:HIS:HB3	1:B:216:ALA:HB1	1.78	0.66
1:C:223:GLN:HE22	1:C:238:THR:HA	1.59	0.66
1:D:74:GLU:HG3	5:D:1046:HOH:O	1.97	0.64
1:D:26:LEU:HA	1:D:274:ARG:NH2	2.13	0.63
1:B:318:ILE:HG22	1:B:320:ILE:HG12	1.80	0.62
1:B:25:LYS:HE3	5:B:1074:HOH:O	1.99	0.62
1:A:220:PHE:HB3	1:A:223:GLN:HB2	1.82	0.62
1:A:224:SER:HA	3:A:1000:NAP:C5B	2.29	0.61
1:C:226:VAL:HG21	1:C:233:ALA:HB2	1.82	0.61
1:D:279:GLU:O	1:D:283:GLN:HG3	2.02	0.60
1:C:224:SER:HA	3:C:1002:NAP:C5B	2.29	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:LYS:O	1:B:298:GLU:HG3	2.03	0.59
1:A:294:LYS:O	1:A:298:GLU:HG3	2.02	0.59
1:A:28:ASN:O	1:A:57:GLU:HG2	2.02	0.58
1:A:262:ARG:NH1	1:A:288:ASN:OD1	2.36	0.58
1:D:26:LEU:HA	1:D:274:ARG:HH21	1.69	0.58
1:D:274:ARG:O	4:D:1003:NAD:H8A	2.04	0.56
1:B:110:LEU:HD21	1:B:112:LEU:HD21	1.87	0.56
1:C:45:LEU:HD23	1:C:45:LEU:C	2.26	0.55
1:A:74:GLU:CD	1:A:74:GLU:H	2.10	0.55
1:A:223:GLN:HG3	1:A:240:PHE:CD1	2.43	0.54
1:B:194:HIS:HB3	1:B:216:ALA:CB	2.38	0.54
1:D:25:LYS:NZ	1:D:224:SER:O	2.42	0.53
1:C:194:HIS:HB3	1:C:216:ALA:HB1	1.90	0.53
1:B:28:ASN:O	1:B:57:GLU:HG2	2.09	0.52
1:A:94:THR:HG21	5:A:1022:HOH:O	2.08	0.52
1:A:58:VAL:O	1:A:62:VAL:HG23	2.10	0.52
1:D:120:LYS:HA	1:D:143:GLU:CG	2.40	0.51
1:D:294:LYS:O	1:D:298:GLU:HG3	2.10	0.51
1:B:190:LEU:HB3	1:B:214:ILE:HD13	1.93	0.51
1:A:116:PRO:HG2	1:A:320:ILE:HD12	1.93	0.51
1:D:43:TYR:CE1	1:D:281:LEU:HD22	2.46	0.50
1:A:148:LEU:HG	1:A:152:LYS:HE2	1.93	0.50
1:C:28:ASN:O	1:C:57:GLU:HG2	2.12	0.50
1:D:238:THR:HG22	5:D:1128:HOH:O	2.12	0.50
1:B:199:GLN:NE2	1:B:304:ASP:HB2	2.27	0.50
1:C:255:THR:OG1	1:C:258:GLU:HG3	2.11	0.49
1:B:204:ILE:O	1:B:208:GLN:HG3	2.12	0.49
1:B:116:PRO:HG2	1:B:320:ILE:HD12	1.94	0.49
1:B:274:ARG:O	4:B:1001:NAD:H8A	2.12	0.49
1:D:120:LYS:HA	1:D:143:GLU:HG3	1.93	0.49
1:A:144:ASP:HB3	1:B:181:ARG:O	2.12	0.49
1:B:45:LEU:HD23	1:B:45:LEU:C	2.33	0.49
1:D:274:ARG:HD3	5:D:1083:HOH:O	2.13	0.48
1:C:223:GLN:NE2	1:C:238:THR:HA	2.28	0.48
1:A:194:HIS:HB3	1:A:216:ALA:HB1	1.95	0.48
1:C:45:LEU:HD23	1:C:46:PHE:N	2.29	0.48
1:A:23:CYS:HA	1:A:26:LEU:HD22	1.96	0.48
1:B:25:LYS:NZ	1:B:224:SER:O	2.47	0.48
1:A:96:LEU:HD22	1:A:163:ILE:HD11	1.96	0.48
1:A:64:ARG:O	1:A:68:GLU:HG3	2.14	0.47
1:D:190:LEU:HB3	1:D:214:ILE:HD13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:194:HIS:HB3	1:D:216:ALA:HB1	1.97	0.46
1:D:98:LYS:HE3	1:D:102:ASP:OD2	2.14	0.46
1:A:181:ARG:O	1:B:144:ASP:HB3	2.15	0.46
1:B:4:SER:O	1:B:6:PRO:HD3	2.15	0.46
1:C:189:VAL:HG12	1:C:190:LEU:N	2.30	0.46
1:C:120:LYS:HE3	5:C:1159:HOH:O	2.15	0.46
1:D:251:LYS:HE3	1:D:252:TYR:CZ	2.51	0.46
1:D:116:PRO:HG2	1:D:320:ILE:HD12	1.98	0.46
1:C:144:ASP:HB3	1:D:181:ARG:O	2.16	0.46
1:C:125:GLU:H	1:C:125:GLU:CD	2.19	0.46
1:C:82:LEU:HD11	1:C:86:TYR:HB2	1.98	0.45
1:C:223:GLN:OE1	1:C:238:THR:HG22	2.17	0.45
1:A:120:LYS:HD2	1:A:141:VAL:HG11	1.98	0.45
1:A:120:LYS:HA	1:A:143:GLU:CG	2.46	0.45
1:A:274:ARG:N	3:A:1000:NAP:O2A	2.48	0.45
1:B:31:ALA:HB3	1:B:57:GLU:HB3	1.98	0.45
1:C:274:ARG:N	3:C:1002:NAP:O2A	2.47	0.44
1:C:221:GLY:O	1:C:224:SER:HB2	2.17	0.44
1:A:96:LEU:O	1:A:100:LEU:HG	2.18	0.44
1:C:84:ASN:OD1	1:C:115:PHE:HB2	2.17	0.44
1:C:110:LEU:HA	1:C:165:SER:O	2.17	0.43
1:D:225:PHE:CD1	1:D:225:PHE:N	2.86	0.43
1:C:189:VAL:CG1	1:C:190:LEU:N	2.81	0.43
1:B:222:PRO:HB3	1:B:309:PHE:CZ	2.53	0.43
1:D:133:TYR:CZ	1:D:135:GLY:HA2	2.53	0.43
1:B:36:TYR:CE2	1:B:40:LYS:HD2	2.54	0.43
1:A:54:ASN:O	1:A:58:VAL:HG23	2.19	0.43
1:A:50:GLU:HG3	1:A:81:LYS:O	2.19	0.43
1:C:5:ILE:HD11	1:C:285:ARG:HD3	2.01	0.43
1:D:318:ILE:HG22	1:D:320:ILE:HG12	2.00	0.43
1:A:274:ARG:O	3:A:1000:NAP:H8A	2.19	0.43
1:B:151:TRP:O	1:B:155:GLU:HG3	2.19	0.43
1:B:83:TRP:CG	1:B:84:ASN:N	2.87	0.43
1:A:295:GLU:CD	1:A:295:GLU:H	2.21	0.42
1:C:24:TRP:O	1:C:25:LYS:HB2	2.18	0.42
1:A:191:GLN:OE1	3:A:1000:NAP:H2N	2.19	0.42
1:C:4:SER:O	1:C:6:PRO:HD3	2.18	0.42
1:D:43:TYR:CE1	1:D:285:ARG:HD2	2.55	0.42
1:B:43:TYR:CE1	1:B:281:LEU:HD22	2.54	0.42
1:A:45:LEU:HD23	1:A:45:LEU:C	2.40	0.42
1:C:175:LEU:HD12	1:C:175:LEU:HA	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:151:TRP:O	1:D:155:GLU:HG3	2.20	0.42
1:C:120:LYS:HA	1:C:143:GLU:HG3	2.00	0.42
1:A:56:LYS:HA	1:A:102:ASP:O	2.20	0.42
1:D:222:PRO:HB3	1:D:309:PHE:CZ	2.54	0.42
1:A:88:ASP:HB3	1:A:91:ASN:ND2	2.35	0.42
1:C:96:LEU:HD22	1:C:163:ILE:HD11	2.02	0.42
1:B:199:GLN:O	1:B:200:GLN:C	2.58	0.41
1:C:25:LYS:HA	1:C:25:LYS:HD3	1.90	0.41
1:B:276:ASN:O	1:B:277:LEU:HD23	2.21	0.41
1:C:194:HIS:HB3	1:C:216:ALA:CB	2.50	0.41
1:B:262:ARG:NH1	1:B:288:ASN:OD1	2.53	0.41
1:C:274:ARG:O	3:C:1002:NAP:H8A	2.20	0.41
1:C:87:HIS:HB2	5:C:1004:HOH:O	2.20	0.41
1:A:88:ASP:HB3	1:A:91:ASN:HD22	1.85	0.41
1:D:81:LYS:HE2	4:D:1003:NAD:H71N	1.85	0.41
1:C:35:VAL:HG13	1:C:46:PHE:CE2	2.56	0.41
1:B:36:TYR:CZ	1:B:40:LYS:HD2	2.55	0.41
1:A:25:LYS:HD3	1:A:25:LYS:HA	1.77	0.41
1:B:103:LEU:O	1:B:104:LYS:HB2	2.21	0.40
1:C:207:ALA:HB3	1:C:214:ILE:HD11	2.03	0.40
1:C:178:ASP:OD1	1:D:147:ILE:HG12	2.22	0.40
1:B:225:PHE:N	1:B:225:PHE:CD1	2.89	0.40
1:C:191:GLN:OE1	3:C:1002:NAP:H2N	2.21	0.40
1:A:21:PHE:HE2	1:A:26:LEU:HD21	1.86	0.40
1:A:35:VAL:HG13	1:A:46:PHE:CE2	2.56	0.40
1:C:151:TRP:O	1:C:155:GLU:HG3	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:1226:HOH:O	5:D:1226:HOH:O[2_657]	1.90	0.30

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	317/322 (98%)	307 (97%)	10 (3%)	0	100	100
1	B	317/322 (98%)	307 (97%)	10 (3%)	0	100	100
1	C	317/322 (98%)	307 (97%)	8 (2%)	2 (1%)	30	36
1	D	317/322 (98%)	308 (97%)	9 (3%)	0	100	100
All	All	1268/1288 (98%)	1229 (97%)	37 (3%)	2 (0%)	52	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	225	PHE
1	C	228	MET

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	268/270 (99%)	264 (98%)	4 (2%)	72	85
1	B	268/270 (99%)	268 (100%)	0	100	100
1	C	268/270 (99%)	265 (99%)	3 (1%)	80	90
1	D	268/270 (99%)	267 (100%)	1 (0%)	93	97
All	All	1072/1080 (99%)	1064 (99%)	8 (1%)	88	95

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	223	GLN
1	A	228	MET
1	A	285	ARG
1	C	230	GLN
1	C	235	ASN

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Mol	Chain	Res	Type
1	C	274	ARG
1	D	274	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	317	ASN
1	B	34	GLN
1	B	91	ASN
1	B	223	GLN
1	C	34	GLN
1	C	223	GLN
1	C	276	ASN
1	C	317	ASN
1	D	34	GLN
1	D	253	ASN
1	D	317	ASN

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](#)

12 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	NAP	A	1000	-	42,52,52	1.77	7 (16%)	54,80,80	1.72	10 (18%)
2	SO4	A	904	-	4,4,4	0.17	0	6,6,6	0.13	0
4	NAD	B	1001	-	38,48,48	1.98	10 (26%)	47,73,73	1.88	8 (17%)
2	SO4	B	901	-	4,4,4	0.21	0	6,6,6	0.09	0
2	SO4	B	905	-	4,4,4	0.17	0	6,6,6	0.08	0
2	SO4	B	907	-	4,4,4	0.23	0	6,6,6	0.08	0
3	NAP	C	1002	-	42,52,52	1.85	9 (21%)	54,80,80	1.82	12 (22%)
2	SO4	C	903	-	4,4,4	0.21	0	6,6,6	0.09	0
4	NAD	D	1003	-	38,48,48	1.97	8 (21%)	47,73,73	1.87	8 (17%)
2	SO4	D	902	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	D	906	-	4,4,4	0.23	0	6,6,6	0.06	0
2	SO4	D	908	-	4,4,4	0.16	0	6,6,6	0.07	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	NAP	A	1000	-	-	0/27/67/67	0/5/5/5
2	SO4	A	904	-	-	0/0/0/0	0/0/0/0
4	NAD	B	1001	-	-	0/22/62/62	0/5/5/5
2	SO4	B	901	-	-	0/0/0/0	0/0/0/0
2	SO4	B	905	-	-	0/0/0/0	0/0/0/0
2	SO4	B	907	-	-	0/0/0/0	0/0/0/0
3	NAP	C	1002	-	-	0/27/67/67	0/5/5/5
2	SO4	C	903	-	-	0/0/0/0	0/0/0/0
4	NAD	D	1003	-	-	0/22/62/62	0/5/5/5
2	SO4	D	902	-	-	0/0/0/0	0/0/0/0
2	SO4	D	906	-	-	0/0/0/0	0/0/0/0
2	SO4	D	908	-	-	0/0/0/0	0/0/0/0

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	NAP	PA-O1A	-2.96	1.40	1.51
3	C	1002	NAP	PA-O1A	-2.83	1.40	1.51
4	D	1003	NAD	PA-O1A	-2.74	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1001	NAD	PA-O1A	-2.69	1.41	1.51
4	D	1003	NAD	O4B-C4B	2.03	1.49	1.45
4	D	1003	NAD	C4A-N3A	2.05	1.38	1.35
3	C	1002	NAP	C6N-C5N	2.05	1.43	1.38
4	B	1001	NAD	C5N-C4N	2.07	1.43	1.38
4	B	1001	NAD	C6N-C5N	2.08	1.43	1.38
3	A	1000	NAP	C4A-N3A	2.09	1.38	1.35
3	C	1002	NAP	C3D-C4D	2.26	1.59	1.53
4	B	1001	NAD	O4B-C4B	2.30	1.50	1.45
3	C	1002	NAP	C4A-N3A	2.42	1.39	1.35
4	B	1001	NAD	C4A-N3A	2.60	1.39	1.35
3	C	1002	NAP	C6N-N1N	3.26	1.44	1.35
4	B	1001	NAD	C6N-N1N	3.32	1.44	1.35
3	A	1000	NAP	C6N-N1N	3.45	1.44	1.35
4	D	1003	NAD	C2A-N1A	3.51	1.40	1.33
4	D	1003	NAD	C6N-N1N	3.60	1.45	1.35
3	A	1000	NAP	C2A-N1A	3.64	1.40	1.33
3	C	1002	NAP	C2A-N1A	3.67	1.40	1.33
3	A	1000	NAP	C2A-N3A	3.86	1.39	1.32
3	C	1002	NAP	C2A-N3A	3.90	1.39	1.32
4	D	1003	NAD	C2A-N3A	3.97	1.39	1.32
4	B	1001	NAD	C2A-N1A	4.08	1.41	1.33
4	B	1001	NAD	C2A-N3A	4.30	1.39	1.32
3	A	1000	NAP	C4N-C3N	4.42	1.46	1.39
3	C	1002	NAP	C4N-C3N	4.50	1.47	1.39
4	B	1001	NAD	C4N-C3N	4.58	1.47	1.39
4	D	1003	NAD	C4N-C3N	4.76	1.47	1.39
3	A	1000	NAP	O4D-C1D	4.85	1.47	1.41
4	B	1001	NAD	O4D-C1D	5.72	1.48	1.41
3	C	1002	NAP	O4D-C1D	5.82	1.48	1.41
4	D	1003	NAD	O4D-C1D	5.95	1.48	1.41

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1002	NAP	N3A-C2A-N1A	-7.81	122.91	128.89
4	B	1001	NAD	N3A-C2A-N1A	-7.62	123.06	128.89
4	D	1003	NAD	N3A-C2A-N1A	-7.59	123.08	128.89
3	A	1000	NAP	N3A-C2A-N1A	-7.55	123.12	128.89
4	D	1003	NAD	C1B-N9A-C4A	-3.85	121.14	126.94
3	C	1002	NAP	C1B-N9A-C4A	-3.82	121.18	126.94
4	B	1001	NAD	C1B-N9A-C4A	-3.81	121.20	126.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1000	NAP	C1B-N9A-C4A	-3.35	121.88	126.94
4	D	1003	NAD	O7N-C7N-C3N	-3.02	116.29	119.59
3	C	1002	NAP	O4B-C1B-N9A	-2.87	102.09	108.10
3	A	1000	NAP	O7N-C7N-C3N	-2.80	116.53	119.59
4	B	1001	NAD	O7N-C7N-C3N	-2.71	116.63	119.59
3	A	1000	NAP	O3B-C3B-C4B	-2.65	103.09	111.05
3	C	1002	NAP	O7N-C7N-C3N	-2.62	116.72	119.59
3	C	1002	NAP	O3B-C3B-C4B	-2.35	104.00	111.05
3	C	1002	NAP	O5B-C5B-C4B	2.05	116.69	109.12
3	C	1002	NAP	O3X-P2B-O2X	2.13	115.48	107.38
3	A	1000	NAP	O3X-P2B-O2X	2.14	115.51	107.38
4	D	1003	NAD	O4B-C4B-C5B	2.20	117.17	109.32
3	A	1000	NAP	C4B-O4B-C1B	2.32	112.26	109.72
4	B	1001	NAD	O2A-PA-O1A	2.38	125.45	112.53
3	A	1000	NAP	O2A-PA-O1A	2.46	125.86	112.53
4	D	1003	NAD	O2A-PA-O1A	2.46	125.88	112.53
3	C	1002	NAP	O2A-PA-O1A	2.47	125.91	112.53
4	B	1001	NAD	O4B-C4B-C5B	2.49	118.22	109.32
3	A	1000	NAP	O4D-C1D-N1N	2.72	111.12	108.13
4	B	1001	NAD	C3N-C7N-N7N	2.83	120.91	117.82
3	C	1002	NAP	C3N-C7N-N7N	2.83	120.91	117.82
4	D	1003	NAD	C2N-C3N-C4N	2.93	121.55	118.29
4	B	1001	NAD	C2N-C3N-C4N	2.95	121.57	118.29
3	C	1002	NAP	C2N-C3N-C4N	2.99	121.61	118.29
3	A	1000	NAP	C3N-C7N-N7N	3.00	121.10	117.82
3	A	1000	NAP	C2N-C3N-C4N	3.02	121.65	118.29
3	C	1002	NAP	C4B-O4B-C1B	3.06	113.08	109.72
3	C	1002	NAP	O4D-C1D-N1N	3.18	111.62	108.13
4	D	1003	NAD	C3N-C7N-N7N	3.43	121.57	117.82
4	D	1003	NAD	O4D-C1D-N1N	3.60	112.08	108.13
4	B	1001	NAD	O4D-C1D-N1N	3.95	112.47	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1000	NAP	5	0
4	B	1001	NAD	1	0
3	C	1002	NAP	6	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1003	NAD	2	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	319/322 (99%)	-0.02	7 (2%) 65 73	17, 26, 40, 56	0
1	B	319/322 (99%)	0.01	6 (1%) 70 76	16, 24, 36, 47	0
1	C	319/322 (99%)	-0.03	8 (2%) 61 70	13, 22, 40, 61	0
1	D	319/322 (99%)	-0.28	3 (0%) 85 89	12, 20, 32, 44	0
All	All	1276/1288 (99%)	-0.08	24 (1%) 70 76	12, 23, 37, 61	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	4	SER	5.1
1	C	228	MET	5.0
1	C	226	VAL	4.7
1	C	229	ASN	4.3
1	A	228	MET	3.5
1	A	229	ASN	3.5
1	A	4	SER	3.4
1	B	138	ASN	3.3
1	C	4	SER	3.3
1	D	4	SER	3.3
1	A	125	GLU	3.2
1	C	138	ASN	3.1
1	C	253	ASN	3.0
1	C	227	GLU	2.8
1	A	138	ASN	2.7
1	D	138	ASN	2.5
1	C	224	SER	2.3
1	A	137	GLY	2.3
1	B	253	ASN	2.3
1	B	112	LEU	2.3
1	A	227	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	159	ALA	2.1
1	D	253	ASN	2.1
1	B	271	VAL	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 [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(Å ²)	Q<0.9
2	SO4	B	901	5/5	0.94	0.30	5.51	74,75,75,75	0
2	SO4	D	906	5/5	0.94	0.18	0.84	66,66,67,67	0
3	NAP	A	1000	48/48	0.94	0.14	-0.22	19,26,37,41	0
4	NAD	B	1001	44/44	0.96	0.14	-0.28	16,26,29,33	0
3	NAP	C	1002	48/48	0.93	0.13	-0.33	15,30,37,39	0
4	NAD	D	1003	44/44	0.97	0.12	-0.37	17,22,26,27	0
2	SO4	D	902	5/5	0.95	0.23	-	63,63,64,64	0
2	SO4	C	903	5/5	0.96	0.23	-	63,63,64,64	0
2	SO4	B	907	5/5	0.88	0.25	-	78,78,79,79	0
2	SO4	B	905	5/5	0.99	0.12	-	42,42,42,43	0
2	SO4	A	904	5/5	0.98	0.17	-	46,46,47,47	0
2	SO4	D	908	5/5	0.94	0.28	-	66,67,67,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.