



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 09:25 PM GMT

PDB ID : 2005
Title : Structure of transhydrogenase (dI.H₂NADH)₂(dIII.NADP⁺)₁asymmetric complex
Authors : Bhakta, T.; Jackson, J.B.
Deposited on : 2007-01-25
Resolution : 2.60 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>

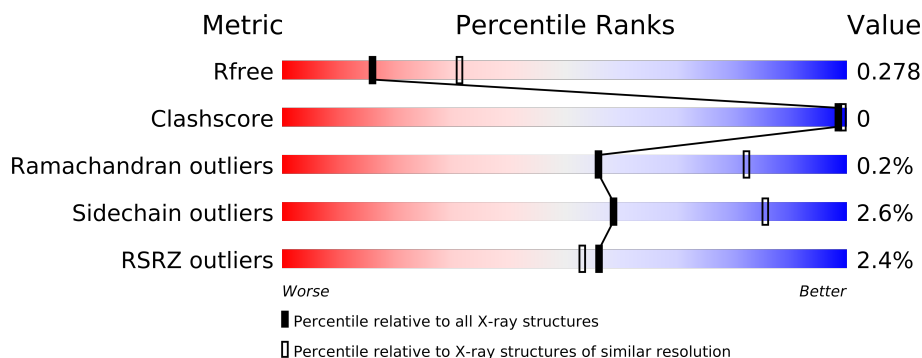
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.60 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	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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	384	
1	B	384	
2	C	174	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SO4	A	385	-	X
3	SO4	A	386	-	X
3	SO4	B	385	-	X
3	SO4	B	386	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7082 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.

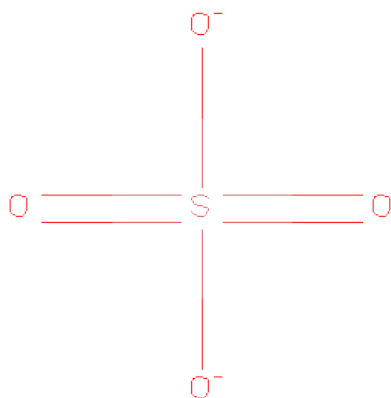
- Molecule 1 is a protein called NAD(P) transhydrogenase subunit alpha part 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	0	0
			2794	1763	481	532	18			
1	B	380	Total	C	N	O	S	0	0	0
			2794	1763	481	532	18			

- Molecule 2 is a protein called NAD(P) transhydrogenase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	174	Total	C	N	O	S	0	0	0
			1311	830	217	253	11			

- Molecule 3 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



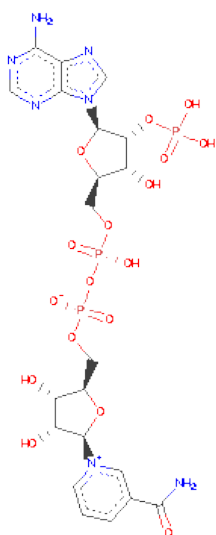
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

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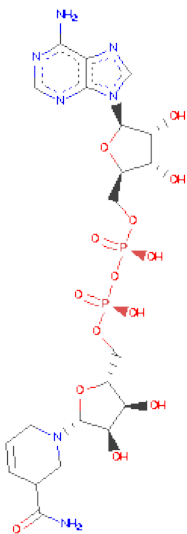
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



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

- Molecule 5 is 1,4,5,6-TETRAHYDRONICOTINAMIDEADENINE DINUCLEOTIDE (three-letter code: TXD) (formula: $C_{21}H_{31}N_7O_{14}P_2$).



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

- Molecule 6 is water.

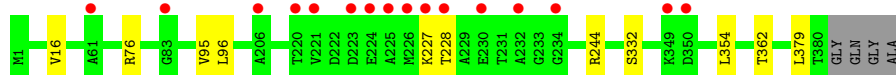
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	10	Total	O	0	0
			10	10		
6	B	8	Total	O	0	0
			8	8		
6	C	9	Total	O	0	0
			9	9		

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 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) transhydrogenase subunit alpha part 1

Chain A: 



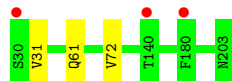
- Molecule 1: NAD(P) transhydrogenase subunit alpha part 1

Chain B: 



- Molecule 2: NAD(P) transhydrogenase subunit beta

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.20Å 101.03Å 131.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	80.06 – 2.60 43.61 – 2.60	Depositor EDS
% Data completeness (in resolution range)	91.9 (80.06-2.60) 91.9 (43.61-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.219 , 0.275 0.226 , 0.278	Depositor DCC
R_{free} test set	1396 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 12.9	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 27507 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7082	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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, TXD, SO4

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	0.40	0/2831	0.58	0/3837
1	B	0.39	0/2831	0.65	2/3837 (0.1%)
2	C	0.40	0/1334	0.55	0/1803
All	All	0.39	0/6996	0.60	2/9477 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	68	ALA	N-CA-C	-6.66	93.02	111.00
1	B	127	ARG	NE-CZ-NH1	5.09	122.85	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	66	SER	Peptide
1	B	67	GLN	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	2794	0	860	1	0
1	B	2794	0	769	3	0
2	C	1311	0	5	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
4	C	48	0	25	0	0
5	A	44	0	28	0	0
5	B	44	0	28	1	0
6	A	10	0	0	0	0
6	B	8	0	0	0	0
6	C	9	0	0	0	0
All	All	7082	0	1715	4	0

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 0.

All (4) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:16:VAL:N	1:A:332:SER:HG	2.03	0.57
1:B:14:ASP:N	1:B:325:SER:HG	2.02	0.57
1:B:127:ARG:NH1	5:B:600:TXD:H6N1	2.32	0.44
1:B:106:ARG:N	1:B:107:PRO:CD	2.84	0.41

There are no symmetry-related clashes.

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	378/384 (98%)	362 (96%)	15 (4%)	1 (0%)	50	77
1	B	378/384 (98%)	365 (97%)	12 (3%)	1 (0%)	50	77
2	C	172/174 (99%)	166 (96%)	6 (4%)	0	100	100
All	All	928/942 (98%)	893 (96%)	33 (4%)	2 (0%)	56	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	379	LEU
1	B	225	ALA

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	295/296 (100%)	287 (97%)	8 (3%)	57	85
1	B	295/296 (100%)	287 (97%)	8 (3%)	57	85
2	C	138/138 (100%)	135 (98%)	3 (2%)	64	89
All	All	728/730 (100%)	709 (97%)	19 (3%)	59	85

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	95	VAL
1	A	96	LEU
1	A	227	LYS
1	A	228	THR
1	A	244	ARG
1	A	354	LEU
1	A	362	THR
1	B	67	GLN
1	B	130	ARG
1	B	137	LEU
1	B	147	ARG
1	B	199	MET

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Mol	Chain	Res	Type
1	B	244	ARG
1	B	289	VAL
1	B	350	ASP
2	C	31	VAL
2	C	61	GLN
2	C	72	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	300	ASN

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 ⓘ

7 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	SO4	A	385	-	4,4,4	0.19	0	6,6,6	0.08	0
3	SO4	A	386	-	4,4,4	0.23	0	6,6,6	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	TXD	A	500	-	48,48,48	2.58	7 (14%)	73,73,73	2.85	14 (19%)
3	SO4	B	385	-	4,4,4	0.19	0	6,6,6	0.10	0
3	SO4	B	386	-	4,4,4	0.20	0	6,6,6	0.12	0
5	TXD	B	600	-	48,48,48	2.59	6 (12%)	73,73,73	2.89	16 (21%)
4	NAP	C	400	-	52,52,52	1.31	3 (5%)	80,80,80	1.64	9 (11%)

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	SO4	A	385	-	-	0/0/0/0	0/0/0/0
3	SO4	A	386	-	-	0/0/0/0	0/0/0/0
5	TXD	A	500	-	1/1/13/15	0/30/72/72	0/3/5/5
3	SO4	B	385	-	-	0/0/0/0	0/0/0/0
3	SO4	B	386	-	-	0/0/0/0	0/0/0/0
5	TXD	B	600	-	1/1/13/15	0/30/72/72	0/3/5/5
4	NAP	C	400	-	-	0/35/67/67	0/3/5/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	600	TXD	C2N-C3N	-13.13	1.39	1.52
5	A	500	TXD	C2N-C3N	-13.06	1.39	1.52
5	A	500	TXD	C2N-N1N	-7.14	1.35	1.47
4	C	400	NAP	O7N-C7N	7.00	1.40	1.24
5	B	600	TXD	C2N-N1N	-6.95	1.35	1.47
5	B	600	TXD	C6N-C5N	-5.33	1.38	1.49
5	A	500	TXD	C4N-C5N	5.00	1.46	1.32
5	A	500	TXD	C6N-C5N	-4.82	1.39	1.49
5	B	600	TXD	C6N-N1N	-4.70	1.34	1.47
5	B	600	TXD	C4N-C5N	4.59	1.45	1.32
5	A	500	TXD	C6N-N1N	-4.35	1.35	1.47
4	C	400	NAP	C2A-N3A	3.10	1.38	1.32
5	B	600	TXD	C4A-N9A	-2.71	1.33	1.37
5	A	500	TXD	C4A-N9A	-2.52	1.34	1.37
4	C	400	NAP	C2A-N1A	2.37	1.38	1.33
5	A	500	TXD	C3N-C7N	-2.32	1.50	1.53

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	500	TXD	C3N-C2N-N1N	15.67	121.58	109.74
5	B	600	TXD	C3N-C2N-N1N	15.15	121.19	109.74
4	C	400	NAP	N3A-C2A-N1A	-10.81	119.67	128.71
5	B	600	TXD	C6N-N1N-C2N	8.82	122.36	109.61
5	A	500	TXD	C2N-C3N-C4N	8.57	121.14	108.75
5	A	500	TXD	N3A-C2A-N1A	-8.38	121.70	128.71
5	B	600	TXD	N3A-C2A-N1A	-8.24	121.82	128.71
5	A	500	TXD	C6N-N1N-C2N	7.93	121.07	109.61
5	B	600	TXD	C2N-C3N-C4N	7.30	119.31	108.75
5	B	600	TXD	C3N-C7N-N7N	5.02	120.13	116.38
5	A	500	TXD	C5N-C6N-N1N	4.38	122.28	110.92
5	B	600	TXD	C5N-C6N-N1N	4.21	121.82	110.92
5	A	500	TXD	N3A-C4A-N9A	4.20	133.01	125.43
5	B	600	TXD	N3A-C4A-N9A	4.18	132.98	125.43
5	B	600	TXD	C2N-C3N-C7N	3.88	122.00	112.19
4	C	400	NAP	O4D-C1D-N1N	3.76	111.80	107.95
5	B	600	TXD	PA-O3-PN	-3.65	120.97	131.68
4	C	400	NAP	N3A-C4A-N9A	3.52	131.79	125.43
5	A	500	TXD	O4'-C1'-N9A	3.31	111.52	108.44
5	A	500	TXD	C2N-C3N-C7N	3.18	120.24	112.19
5	B	600	TXD	C4N-C3N-C7N	3.08	118.69	110.04
5	A	500	TXD	C4N-C3N-C7N	3.06	118.61	110.04
5	A	500	TXD	C5A-C4A-N3A	-2.83	119.55	125.70
5	B	600	TXD	O4'-C1'-N9A	2.72	110.97	108.44
5	A	500	TXD	PA-O3-PN	-2.69	123.79	131.68
5	A	500	TXD	C2A-N3A-C4A	2.68	121.65	114.01
4	C	400	NAP	C6N-N1N-C2N	-2.61	119.10	122.04
5	B	600	TXD	C5A-C4A-N3A	-2.60	120.03	125.70
5	B	600	TXD	O7N-C7N-C3N	-2.56	119.20	121.26
5	B	600	TXD	C2A-N3A-C4A	2.52	121.20	114.01
5	A	500	TXD	C3N-C4N-C5N	-2.43	112.06	122.25
5	B	600	TXD	C8A-N9A-C4A	2.42	108.75	106.90
4	C	400	NAP	C5A-C4A-N3A	-2.26	120.78	125.70
4	C	400	NAP	C2A-N3A-C4A	2.24	120.38	114.01
5	A	500	TXD	O4'A-C1'A-N1N	2.19	112.01	108.78
5	B	600	TXD	C4'-O4'-C1'	-2.17	107.40	109.75
4	C	400	NAP	C4A-C5A-N7A	-2.05	107.77	109.52
4	C	400	NAP	C2D-C1D-N1N	2.04	117.31	113.86
4	C	400	NAP	N7A-C8A-N9A	-2.04	108.59	114.36

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	500	TXD	C3N
5	B	600	TXD	C3N

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	380/384 (98%)	0.09	16 (4%) 35 31	22, 42, 66, 77	0
1	B	380/384 (98%)	-0.08	4 (1%) 77 79	19, 38, 58, 69	0
2	C	174/174 (100%)	0.06	3 (1%) 67 66	23, 40, 56, 63	0
All	All	934/942 (99%)	0.02	23 (2%) 56 52	19, 40, 61, 77	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	140	THR	4.0
1	A	224	GLU	3.5
1	A	226	MET	3.5
1	B	80	ALA	3.3
1	A	228	THR	3.2
2	C	30	SER	3.0
1	A	221	VAL	2.9
1	A	350	ASP	2.7
1	A	83	GLY	2.6
1	A	227	LYS	2.6
1	A	232	ALA	2.4
1	A	220	THR	2.4
2	C	180	PHE	2.3
1	A	225	ALA	2.3
1	A	349	LYS	2.2
1	A	230	GLU	2.2
1	A	223	ASP	2.2
1	A	61	ALA	2.1
1	B	349	LYS	2.1
1	B	59	SER	2.1
1	B	224	GLU	2.1
1	A	234	GLY	2.0
1	A	206	ALA	2.0

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(Å ²)	Q<0.9
3	SO4	B	386	5/5	0.23	8.10	85,85,85,85	0
3	SO4	B	385	5/5	0.26	4.71	76,76,77,77	0
3	SO4	A	386	5/5	0.24	4.57	79,80,80,80	0
3	SO4	A	385	5/5	0.29	4.57	82,83,83,83	0
4	NAP	C	400	48/48	0.15	-0.53	21,29,31,32	0
5	TXD	B	600	44/44	0.13	-0.87	21,26,31,32	0
5	TXD	A	500	44/44	0.13	-1.00	34,35,37,38	0

6.5 Other polymers ⓘ

There are no such residues in this entry.