



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:43 AM GMT

PDB ID : 2JD0
Title : X-RAY STRUCTURE OF MUTANT 1-DEOXY-D-XYLULOSE 5-PHOSPHATE REDUCTOISOMERASE, DXR, RV2870C, FROM MYCOBACTERIUM TUBERCULOSIS, IN COMPLEX WITH NADPH
Authors : Henriksson, L.M.; Unge, T.; Jones, T.A.; Mowbray, S.L.
Deposited on : 2007-01-04
Resolution : 2.30 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>

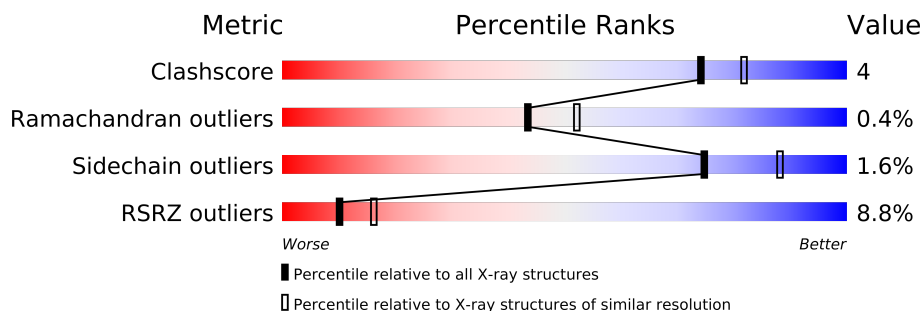
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.30 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))
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (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. 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	398	
1	B	398	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5907 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.

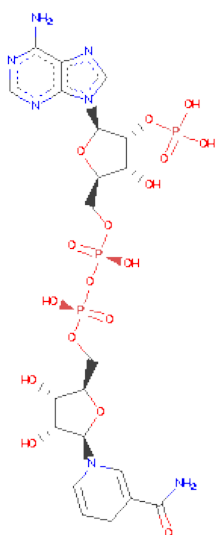
- Molecule 1 is a protein called 1-DEOXY-D-XYLULOSE 5-PHOSPHATE REDUCTOISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2781	1742	504	526	9			
1	B	379	Total	C	N	O	S	0	2	0
			2800	1753	509	528	10			

There are 4 discrepancies between the modelled and reference sequences:

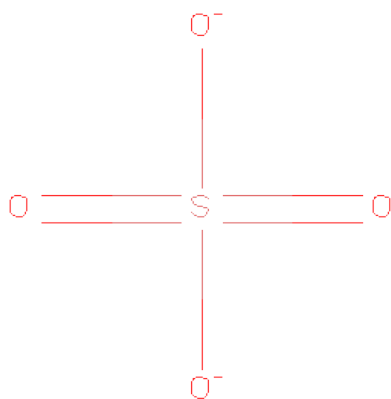
Chain	Residue	Modelled	Actual	Comment	Reference
A	151	ASN	ASP	ENGINEERED MUTATION	UNP P64012
A	222	GLN	GLU	ENGINEERED MUTATION	UNP P64012
B	151	ASN	ASP	ENGINEERED MUTATION	UNP P64012
B	222	GLN	GLU	ENGINEERED MUTATION	UNP P64012

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



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

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



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

- Molecule 4 is water.

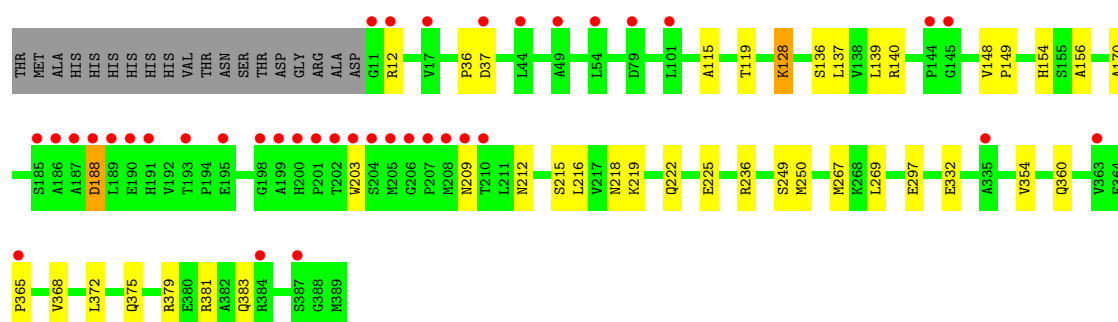
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	115	Total	O	0	0
			115	115		
4	B	148	Total	O	0	0
			148	148		

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.

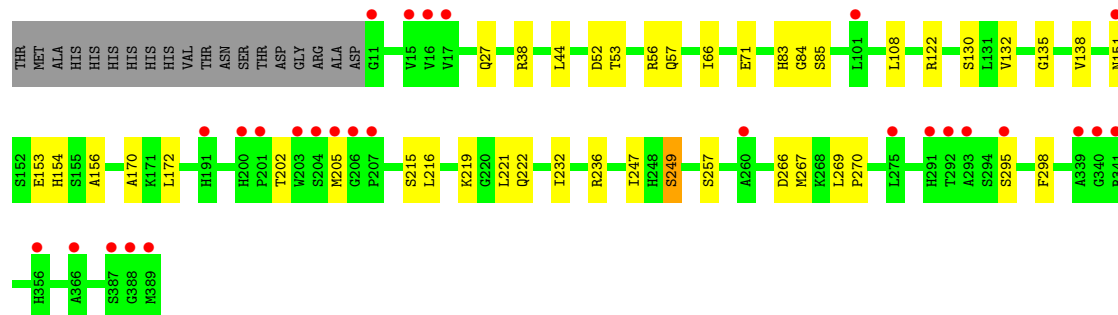
• Molecule 1: 1-DEOXY-D-XYLULOSE 5-PHOSPHATE REDUCTOISOMERASE

Chain A: 



• Molecule 1: 1-DEOXY-D-XYLULOSE 5-PHOSPHATE REDUCTOISOMERASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.63Å 65.61Å 86.02Å 90.00° 102.02° 90.00°	Depositor
Resolution (Å)	35.00 – 2.30 24.08 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (35.00-2.30) 100.0 (24.08-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.37 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.255 0.198 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.940	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 21.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 32928 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5907	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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: NDP, 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	0.43	0/2835	0.55	0/3873
1	B	0.43	0/2854	0.55	0/3897
All	All	0.43	0/5689	0.55	0/7770

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2781	0	2772	23	0
1	B	2800	0	2792	25	0
2	A	48	0	26	0	0
3	A	10	0	0	0	0
3	B	5	0	0	0	0
4	A	115	0	0	2	0
4	B	148	0	0	2	0
All	All	5907	0	5590	46	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 4.

All (46) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:71:GLU:OE2	1:B:83:HIS:HD2	1.53	0.90
1:B:71:GLU:OE2	1:B:83:HIS:CD2	2.37	0.76
1:A:219:LYS:NZ	1:A:222:GLN:HE22	1.90	0.69
1:B:151:ASN:HD21	1:B:153:GLU:HB2	1.60	0.67
1:B:151:ASN:HD22	1:B:154:HIS:H	1.50	0.59
1:A:219:LYS:HZ3	1:A:222:GLN:HE22	1.50	0.58
1:A:269:LEU:HD21	1:B:257:SER:HB2	1.85	0.57
1:A:297:GLU:HG2	1:B:295:SER:HB3	1.87	0.56
1:A:375:GLN:O	1:A:379:ARG:HG3	2.07	0.53
1:A:188:ASP:N	1:A:188:ASP:OD2	2.41	0.53
1:A:140:ARG:NH2	1:A:365:PRO:O	2.42	0.53
1:A:203:TRP:CZ3	1:A:267:MET:HG2	2.43	0.52
1:B:219:LYS:NZ	1:B:222:GLN:HE22	2.07	0.51
1:A:332:GLU:HG3	1:A:379:ARG:HA	1.94	0.49
1:B:172:LEU:HD11	1:B:232:ILE:HG13	1.94	0.49
1:B:247:ILE:HD12	1:B:298:PHE:HZ	1.78	0.48
1:A:170:ALA:O	1:A:236:ARG:HD2	2.14	0.48
1:A:148:VAL:HG13	4:A:2005:HOH:O	2.13	0.48
1:B:170:ALA:O	1:B:236:ARG:HD2	2.14	0.47
1:B:151:ASN:ND2	1:B:154:HIS:H	2.12	0.47
1:A:354:VAL:HG22	1:A:381:ARG:HG3	1.97	0.47
1:B:151:ASN:HB2	4:B:2065:HOH:O	2.15	0.47
1:A:360:GLN:HG2	4:A:2104:HOH:O	2.14	0.46
1:A:149:PRO:HB2	1:A:154:HIS:HB2	1.96	0.46
1:B:156:ALA:HB3	1:B:249:SER:HB2	1.98	0.46
1:B:53:THR:O	1:B:57:GLN:HG3	2.16	0.46
1:A:128:LYS:HE3	1:A:225:GLU:OE1	2.16	0.45
1:B:38[B]:ARG:HA	1:B:38[B]:ARG:HD2	1.87	0.45
1:B:27:GLN:HE22	1:B:267[B]:MET:CE	2.29	0.45
1:B:71:GLU:HG2	1:B:84:GLY:HA2	1.98	0.44
1:A:212:ASN:HB3	1:A:218:ASN:HB2	1.99	0.44
1:A:136:SER:HA	1:A:139:LEU:HB2	1.98	0.44
1:A:12:ARG:HD2	1:A:36:PRO:O	2.18	0.44
1:A:156:ALA:HB3	1:A:249:SER:HB2	2.00	0.43
1:B:85:SER:HA	4:B:2033:HOH:O	2.18	0.43
1:B:44:LEU:HB2	1:B:66:ILE:HG22	2.01	0.42
1:A:215:SER:O	1:A:216:LEU:HB2	2.19	0.42
1:B:215:SER:O	1:B:216:LEU:HB2	2.20	0.42
1:B:130:SER:O	1:B:138:VAL:HG21	2.19	0.42
1:A:219:LYS:HZ2	1:A:222:GLN:HE22	1.63	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:132:VAL:HG11	1:B:221:LEU:HD22	2.02	0.42
1:B:269:LEU:HB3	1:B:270:PRO:CD	2.50	0.41
1:B:52:ASP:O	1:B:56:ARG:HG3	2.21	0.41
1:A:368:VAL:O	1:A:372:LEU:HG	2.21	0.41
1:A:115:ALA:O	1:A:119:THR:HG23	2.21	0.41
1:B:202:THR:HB	1:B:205:MET:HB3	2.02	0.40

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	377/398 (95%)	368 (98%)	9 (2%)	0	100	100
1	B	379/398 (95%)	369 (97%)	7 (2%)	3 (1%)	27	30
All	All	756/796 (95%)	737 (98%)	16 (2%)	3 (0%)	43	52

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	249	SER
1	B	135	GLY
1	B	266	ASP

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	282/298 (95%)	275 (98%)	7 (2%)	60	77

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	284/298 (95%)	282 (99%)	2 (1%)	91	97
All	All	566/596 (95%)	557 (98%)	9 (2%)	75	88

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASP
1	A	128	LYS
1	A	137	LEU
1	A	188	ASP
1	A	209	ASN
1	A	250	MET
1	A	383	GLN
1	B	108	LEU
1	B	122	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	ASN
1	A	191	HIS
1	A	209	ASN
1	A	222	GLN
1	A	375	GLN
1	B	27	GLN
1	B	83	HIS
1	B	209	ASN
1	B	222	GLN
1	B	383	GLN

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

4 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	NDP	A	1390	-	52,52,52	1.45	4 (7%)	80,80,80	2.17	14 (17%)
3	SO4	A	1391	-	4,4,4	0.25	0	6,6,6	0.20	0
3	SO4	A	1392	-	4,4,4	0.06	0	6,6,6	0.10	0
3	SO4	B	1390	-	4,4,4	0.27	0	6,6,6	0.37	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
2	NDP	A	1390	-	-	0/35/77/77	0/3/5/5
3	SO4	A	1391	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1392	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1390	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1390	NDP	PN-O3	-7.74	1.45	1.59
2	A	1390	NDP	C2A-N3A	2.95	1.38	1.32
2	A	1390	NDP	C2A-N1A	2.34	1.38	1.33
2	A	1390	NDP	O4D-C1D	2.21	1.47	1.42

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1390	NDP	N3A-C2A-N1A	-11.14	119.40	128.71
2	A	1390	NDP	PN-O3-PA	8.81	157.52	131.68
2	A	1390	NDP	O4B-C1B-N9A	5.90	113.93	108.44
2	A	1390	NDP	O3-PN-O1N	4.63	144.54	111.28
2	A	1390	NDP	O3-PN-O5D	-4.33	84.05	103.41
2	A	1390	NDP	N3A-C4A-N9A	3.61	131.94	125.43
2	A	1390	NDP	C4N-C3N-C2N	-2.84	118.22	121.68
2	A	1390	NDP	O2N-PN-O3	-2.52	93.20	105.14
2	A	1390	NDP	C2A-N3A-C4A	2.17	120.18	114.01
2	A	1390	NDP	C5A-C4A-N3A	-2.15	121.02	125.70
2	A	1390	NDP	C6N-N1N-C2N	2.12	122.21	119.44
2	A	1390	NDP	C2N-C3N-C7N	2.09	123.83	118.49
2	A	1390	NDP	N7A-C8A-N9A	-2.07	108.50	114.36
2	A	1390	NDP	O4D-C1D-N1N	2.02	112.34	108.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	379/398 (95%)	0.60	38 (10%) 8 13	13, 26, 59, 79	0
1	B	379/398 (95%)	0.28	28 (7%) 14 21	12, 22, 43, 56	0
All	All	758/796 (95%)	0.44	66 (8%) 10 16	12, 24, 49, 79	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	203	TRP	11.6
1	A	201	PRO	8.8
1	A	202	THR	7.9
1	A	200	HIS	7.8
1	A	199	ALA	7.2
1	A	11	GLY	6.6
1	A	204	SER	6.5
1	A	206	GLY	6.3
1	B	389	MET	5.6
1	A	185	SER	4.8
1	A	205	MET	4.6
1	A	191	HIS	4.3
1	B	205	MET	4.2
1	B	206	GLY	3.9
1	A	186	ALA	3.9
1	A	335	ALA	3.9
1	B	356	HIS	3.8
1	B	388	GLY	3.8
1	A	207	PRO	3.8
1	B	339	ALA	3.8
1	A	189	LEU	3.6
1	B	203	TRP	3.4
1	B	387	SER	3.3
1	B	291	HIS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	195	GLU	3.2
1	B	201	PRO	3.2
1	A	208	MET	3.2
1	A	190	GLU	3.0
1	A	198	GLY	3.0
1	A	209	ASN	3.0
1	B	15	VAL	2.9
1	B	204	SER	2.9
1	A	384	ARG	2.9
1	B	340	GLY	2.9
1	B	191	HIS	2.8
1	B	17	VAL	2.8
1	B	11	GLY	2.8
1	A	188	ASP	2.8
1	A	187	ALA	2.7
1	B	293	ALA	2.7
1	B	260	ALA	2.6
1	B	200	HIS	2.6
1	A	17	VAL	2.5
1	A	365	PRO	2.5
1	A	12	ARG	2.5
1	B	151	ASN	2.5
1	B	292	THR	2.5
1	B	366	ALA	2.4
1	A	387	SER	2.3
1	A	49	ALA	2.3
1	B	16	VAL	2.2
1	B	295	SER	2.2
1	B	101	LEU	2.2
1	A	144	PRO	2.2
1	A	44	LEU	2.1
1	A	54	LEU	2.1
1	A	363	VAL	2.1
1	A	79	ASP	2.1
1	B	207	PRO	2.1
1	A	101	LEU	2.1
1	B	275	LEU	2.1
1	A	193	THR	2.1
1	A	145	GLY	2.1
1	A	210	THR	2.1
1	A	37	ASP	2.1
1	B	341	ARG	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
2	NDP	A	1390	48/48	0.35	1.53	45,60,85,86	21
3	SO4	B	1390	5/5	0.13	0.31	29,29,30,30	0
3	SO4	A	1392	5/5	0.11	-0.06	36,37,37,38	0
3	SO4	A	1391	5/5	0.11	-0.56	31,33,33,33	0

6.5 Other polymers

There are no such residues in this entry.