



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 04:04 AM GMT

PDB ID : 4JR9  
Title : Crystal structure of nitrate/nitrite exchanger NarK  
Authors : Zheng, H.; Wisedchaisri, G.; Gonen, T.  
Deposited on : 2013-03-21  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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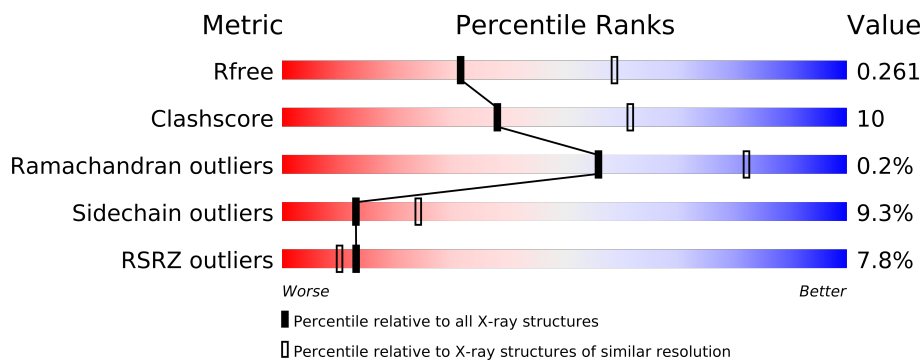
The following versions of software and data (see [references](#)) were used in the production of this report:

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  $\geq 3$ , 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	466	
2	H	217	
3	L	211	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6337 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 Nitrite extrusion protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	1	0
			3093	2058	493	522	20			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P10903
A	-1	SER	-	EXPRESSION TAG	UNP P10903
A	0	HIS	-	EXPRESSION TAG	UNP P10903

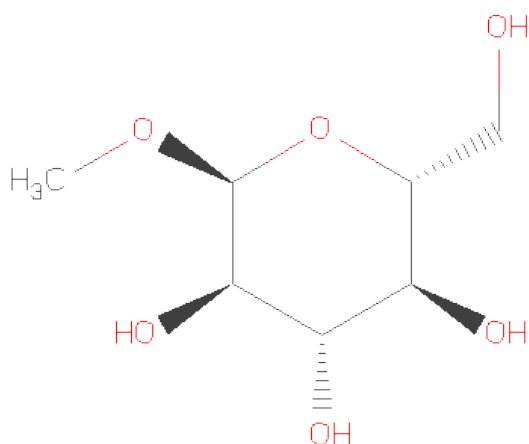
- Molecule 2 is a protein called Immunoglobulin Gamma-2a, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	213	Total	C	N	O	S	0	0	0
			1592	1016	261	308	7			

- Molecule 3 is a protein called Immunoglobulin Kappa, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1594	1001	261	326	6			

- Molecule 4 is SUGAR (METHYL-ALPHA-D-GLUCOPYRANOSIDE) (three-letter code: GYP) (formula: C<sub>7</sub>H<sub>14</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	7	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total	O	0	0
			5	5		
5	H	22	Total	O	0	0
			22	22		
5	L	18	Total	O	0	0
			18	18		



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.27Å 86.55Å 144.79Å 90.00° 115.41° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 43.27 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.60) 99.5 (43.27-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.96 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.229 , 0.261 0.229 , 0.261	Depositor DCC
$R_{free}$ test set	2406 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	56.1	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 43.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 47445 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6337	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: GYP

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	0.52	6/3181 (0.2%)	0.56	2/4322 (0.0%)
2	H	0.65	2/1636 (0.1%)	0.75	2/2234 (0.1%)
3	L	0.56	3/1630 (0.2%)	0.65	1/2226 (0.0%)
All	All	0.57	11/6447 (0.2%)	0.64	5/8782 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	162	TRP	CD2-CE2	6.26	1.48	1.41
1	A	219	TRP	CD2-CE2	5.58	1.48	1.41
2	H	50	TRP	CD2-CE2	5.28	1.47	1.41
3	L	147	TRP	CD2-CE2	5.27	1.47	1.41
1	A	257	TRP	CD2-CE2	5.23	1.47	1.41
3	L	35	TRP	CD2-CE2	5.23	1.47	1.41
1	A	39	TRP	CD2-CE2	5.22	1.47	1.41
2	H	108	TRP	CD2-CE2	5.21	1.47	1.41
1	A	52	TRP	CD2-CE2	5.08	1.47	1.41
1	A	217	TRP	CD2-CE2	5.06	1.47	1.41
1	A	454	TRP	CD2-CE2	5.05	1.47	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	157	LEU	CA-CB-CG	7.25	131.98	115.30
3	L	95	LEU	CA-CB-CG	6.63	130.56	115.30
2	H	83	LEU	CB-CG-CD2	-6.13	100.58	111.00
1	A	46	LEU	CA-CB-CG	6.05	129.22	115.30
1	A	315	LEU	CA-CB-CG	5.78	128.60	115.30

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	3093	0	3128	40	0
2	H	1592	0	1541	52	0
3	L	1594	0	1494	35	0
4	A	13	0	14	1	0
5	A	5	0	0	0	0
5	H	22	0	0	1	0
5	L	18	0	0	1	0
All	All	6337	0	6177	123	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 10.

All (123) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:100:TRP:O	2:H:100:TRP:HE3	1.21	1.24
2:H:50:TRP:CE2	2:H:100:TRP:HD1	1.72	1.06
2:H:37:VAL:HG21	2:H:105:MET:HE1	1.37	1.04
2:H:100:TRP:O	2:H:100:TRP:CE3	2.12	1.01
2:H:18:VAL:HG23	2:H:86:LEU:HD11	1.39	0.99
2:H:50:TRP:CE2	2:H:100:TRP:CD1	2.51	0.98
2:H:34:MET:CE	2:H:96:CYS:HB2	1.95	0.96
2:H:37:VAL:CG2	2:H:105:MET:HE1	1.96	0.95
2:H:18:VAL:CG2	2:H:86:LEU:HD11	2.01	0.91
2:H:34:MET:HE1	2:H:96:CYS:HB2	1.53	0.90
2:H:35:ASN:HD22	2:H:47:TRP:HE1	1.21	0.87
2:H:50:TRP:CD2	2:H:100:TRP:HD1	1.95	0.83
1:A:36:ARG:HH21	1:A:235:ASN:HD21	1.28	0.82
1:A:312:SER:HB3	1:A:371:GLN:HG2	1.61	0.81
1:A:119:LEU:HB3	1:A:136:ILE:HG21	1.65	0.77
3:L:132:VAL:HG22	3:L:177:THR:HG22	1.66	0.75
3:L:5:THR:HG23	5:L:302:HOH:O	1.92	0.69
2:H:97:ALA:HB3	2:H:105:MET:HE2	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:260:SER:HB3	1:A:413:ILE:HG21	1.75	0.68
2:H:37:VAL:CG2	2:H:105:MET:CE	2.70	0.68
2:H:35:ASN:ND2	2:H:47:TRP:HE1	1.92	0.68
2:H:55:ASN:HB3	2:H:57:TYR:HB2	1.79	0.64
2:H:100:TRP:C	2:H:100:TRP:HE3	2.00	0.64
1:A:326:ILE:HD12	1:A:450:VAL:HG21	1.79	0.64
2:H:50:TRP:CZ2	2:H:100:TRP:CD1	2.87	0.63
1:A:312:SER:CB	1:A:371:GLN:HG2	2.28	0.63
2:H:50:TRP:CD2	2:H:100:TRP:CD1	2.84	0.63
2:H:34:MET:HE1	2:H:96:CYS:CB	2.27	0.62
3:L:89:LEU:HD21	3:L:95:LEU:HD13	1.83	0.61
3:L:112:PRO:HG3	3:L:143:ILE:HD11	1.82	0.60
1:A:335:LEU:HD23	1:A:439:MET:SD	2.41	0.60
3:L:11:LEU:HD11	3:L:19:VAL:HG13	1.83	0.60
1:A:181:MET:HG3	1:A:219:TRP:CH2	2.37	0.59
2:H:52:ASN:HB3	2:H:55:ASN:HB2	1.84	0.59
2:H:100:TRP:CZ3	3:L:91:TYR:O	2.56	0.59
2:H:100:TRP:C	2:H:100:TRP:CE3	2.76	0.58
2:H:141:VAL:O	2:H:187:THR:HA	2.04	0.58
1:A:182:GLN:HB2	1:A:295:PHE:CB	2.34	0.58
3:L:34:VAL:HG13	3:L:91:TYR:CE1	2.38	0.57
3:L:196:THR:HG22	3:L:203:PRO:HB3	1.85	0.57
2:H:100:TRP:HZ3	3:L:91:TYR:O	1.87	0.57
3:L:148:LYS:HB2	3:L:192:THR:HB	1.86	0.57
1:A:160:LYS:HG3	1:A:397:ARG:HA	1.87	0.56
3:L:34:VAL:HG23	3:L:89:LEU:HB3	1.87	0.56
1:A:381:THR:HG21	1:A:398:GLU:HB2	1.87	0.56
1:A:399:ALA:O	1:A:403:THR:HG23	2.06	0.55
3:L:162:TRP:N	3:L:162:TRP:CD1	2.75	0.55
1:A:178:VAL:O	1:A:182:GLN:HG2	2.06	0.55
1:A:358:LEU:HA	1:A:361:THR:HG22	1.89	0.55
2:H:168:VAL:HG22	2:H:186:VAL:CG1	2.38	0.54
2:H:58:THR:CG2	2:H:70:PHE:HB2	2.38	0.54
1:A:37:ASN:HD21	1:A:235:ASN:HD22	1.55	0.53
3:L:143:ILE:HG22	3:L:162:TRP:CZ3	2.43	0.53
1:A:271:ILE:HG22	1:A:418:GLY:HA3	1.90	0.53
3:L:38:GLN:HG2	3:L:44:PRO:HG3	1.90	0.53
2:H:17:THR:HA	2:H:83:LEU:O	2.08	0.53
2:H:55:ASN:CB	2:H:57:TYR:HB2	2.38	0.53
2:H:159:TRP:HZ3	2:H:215:ILE:HD13	1.73	0.53
3:L:78:LEU:HD11	3:L:103:LEU:HD21	1.89	0.53
1:A:381:THR:CG2	1:A:398:GLU:HB2	2.40	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:L:143:ILE:HG22	3:L:162:TRP:CH2	2.46	0.51
1:A:75:PHE:O	1:A:78:THR:HG22	2.11	0.51
2:H:91:THR:HG23	2:H:115:THR:HA	1.93	0.51
3:L:140:PRO:O	3:L:197:HIS:HE1	1.94	0.50
3:L:79:GLN:NE2	3:L:81:GLU:OE2	2.38	0.50
3:L:80:PRO:HA	3:L:105:ILE:HD12	1.93	0.49
2:H:55:ASN:CB	2:H:57:TYR:H	2.25	0.49
1:A:312:SER:OG	1:A:317:GLY:HA2	2.12	0.49
2:H:47:TRP:CH2	2:H:49:GLY:HA2	2.49	0.48
2:H:68:PHE:CE2	2:H:83:LEU:HD21	2.48	0.48
1:A:308:GLY:O	1:A:312:SER:HB2	2.13	0.48
2:H:143:LEU:HD22	2:H:215:ILE:HG21	1.96	0.48
2:H:189:THR:O	2:H:192:THR:HG22	2.14	0.48
2:H:97:ALA:CB	2:H:105:MET:HE2	2.41	0.47
3:L:81:GLU:HG3	3:L:81:GLU:H	1.38	0.47
3:L:114:VAL:HG22	3:L:135:LEU:HD22	1.95	0.47
3:L:107:ARG:HD3	3:L:108:ALA:O	2.14	0.47
3:L:35:TRP:CZ3	3:L:88:CYS:HB3	2.50	0.47
2:H:12:LYS:HG3	2:H:18:VAL:HG22	1.96	0.47
1:A:147:PHE:O	1:A:151:MET:HB2	2.15	0.46
1:A:182:GLN:HE21	1:A:298:PRO:HG2	1.81	0.46
1:A:17:ASP:H	2:H:102:ASN:HD21	1.63	0.46
2:H:55:ASN:HB3	2:H:57:TYR:H	1.81	0.46
1:A:188:VAL:HG21	1:A:218:ILE:HB	1.97	0.46
1:A:36:ARG:NH2	1:A:235:ASN:HD21	2.06	0.46
1:A:55:PHE:HA	1:A:58:VAL:HB	1.97	0.46
2:H:153:GLU:HG3	5:H:316:HOH:O	2.14	0.46
1:A:290:ILE:HG13	1:A:294:ALA:HB2	1.98	0.45
1:A:80:LEU:N	1:A:81:PRO:HD2	2.32	0.45
3:L:29:ILE:HG21	3:L:90:GLN:HG3	1.99	0.45
2:H:34:MET:HE2	2:H:96:CYS:HB2	1.93	0.45
3:L:141:LYS:HD2	3:L:172:TYR:CE1	2.52	0.45
2:H:36:TRP:HB3	2:H:48:MET:CE	2.47	0.45
1:A:21:GLU:OE2	1:A:160:LYS:HG2	2.16	0.45
1:A:122:ALA:HB1	1:A:128:THR:HG21	1.98	0.45
1:A:289:GLN:HG2	1:A:289:GLN:O	2.18	0.44
3:L:34:VAL:CG2	3:L:89:LEU:HB3	2.47	0.44
1:A:181:MET:HG3	1:A:219:TRP:HH2	1.81	0.44
1:A:393:GLU:HB3	1:A:397:ARG:NH2	2.32	0.44
2:H:76:ALA:O	2:H:78:THR:HG23	2.17	0.44
2:H:36:TRP:HB3	2:H:48:MET:HE1	1.99	0.43
3:L:35:TRP:CH2	3:L:88:CYS:HB3	2.52	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:119:ALA:HB3	2:H:151:PHE:CE2	2.53	0.43
2:H:58:THR:HG22	2:H:70:PHE:HB2	2.01	0.43
1:A:271:ILE:HA	1:A:271:ILE:HD13	1.90	0.43
1:A:258:ILE:HD13	1:A:456:VAL:HG21	2.01	0.43
3:L:35:TRP:CE2	3:L:73:PHE:HB2	2.54	0.43
3:L:3:LEU:HB2	3:L:26:SER:HB3	2.00	0.43
1:A:420:PHE:HB3	4:A:501:GYP:H2	2.01	0.42
3:L:186:GLU:HA	3:L:210:ARG:NH1	2.34	0.42
2:H:142:THR:HA	2:H:186:VAL:O	2.20	0.41
1:A:300:ILE:HD13	1:A:300:ILE:HA	1.84	0.41
2:H:69:SER:HB2	2:H:82:GLN:HB3	2.02	0.41
2:H:47:TRP:CE2	3:L:95:LEU:HG	2.56	0.41
3:L:149:ILE:HD12	3:L:154:ARG:HD3	2.02	0.41
1:A:120:GLY:HA3	1:A:217:TRP:NE1	2.36	0.41
3:L:47:LEU:HD23	3:L:58:ILE:HD12	2.03	0.41
1:A:272:GLY:HA2	1:A:418:GLY:HA2	2.03	0.40
1:A:179:SER:OG	1:A:299:PHE:HB2	2.21	0.40
2:H:50:TRP:NE1	2:H:100:TRP:CD1	2.89	0.40
2:H:142:THR:HB	2:H:187:THR:HG22	2.02	0.40
3:L:194:GLU:CG	3:L:205:VAL:HG22	2.52	0.40
3:L:190:SER:HB3	3:L:209:ASN:OD1	2.21	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	404/466 (87%)	393 (97%)	11 (3%)	0	100	100
2	H	209/217 (96%)	206 (99%)	2 (1%)	1 (0%)	38	67
3	L	209/211 (99%)	197 (94%)	11 (5%)	1 (0%)	38	67
All	All	822/894 (92%)	796 (97%)	24 (3%)	2 (0%)	56	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	99	SER
3	L	92	ASN

### 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	315/364 (86%)	292 (93%)	23 (7%)	20	38
2	H	174/182 (96%)	154 (88%)	20 (12%)	8	15
3	L	177/190 (93%)	158 (89%)	19 (11%)	10	17
All	All	666/736 (90%)	604 (91%)	62 (9%)	13	24

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	46	LEU
1	A	53	MET
1	A	70	THR
1	A	78	THR
1	A	119	LEU
1	A	126	THR
1	A	181	MET
1	A	189	VAL
1	A	216	SER
1	A	256	LEU
1	A	261	LEU
1	A	284	GLN
1	A	312	SER
1	A	315	LEU
1	A	326	ILE
1	A	328	MET
1	A	344	GLN
1	A	355	PHE
1	A	371	GLN
1	A	409	PHE
1	A	416	ILE

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Mol	Chain	Res	Type
1	A	431	LEU
2	H	5	VAL
2	H	57	TYR
2	H	58	THR
2	H	83	LEU
2	H	100	TRP
2	H	102	ASN
2	H	121	THR
2	H	141	VAL
2	H	142	THR
2	H	148	LYS
2	H	156	THR
2	H	157	LEU
2	H	163	SER
2	H	181	THR
2	H	186	VAL
2	H	187	THR
2	H	195	SER
2	H	201	ASN
2	H	208	SER
2	H	209	THR
3	L	7	SER
3	L	12	SER
3	L	15	VAL
3	L	30	THR
3	L	33	ILE
3	L	38	GLN
3	L	79	GLN
3	L	81	GLU
3	L	89	LEU
3	L	90	GLN
3	L	95	LEU
3	L	105	ILE
3	L	107	ARG
3	L	113	THR
3	L	124	LEU
3	L	162	TRP
3	L	174	MET
3	L	186	GLU
3	L	201	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	182	GLN
1	A	235	ASN
1	A	284	GLN
1	A	344	GLN
1	A	371	GLN
2	H	35	ASN
2	H	102	ASN
3	L	27	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 ⓘ

1 ligand is 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	$\# Z  > 2$
4	GYP	A	501	-	13,13,13	0.72	1 (7%)	18,18,18	1.56	3 (16%)

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
4	GYP	A	501	-	-	0/4/24/24	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	501	GYP	O1-C1	2.07	1.43	1.40

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501	GYP	O1-C1-C2	4.18	113.30	108.18
4	A	501	GYP	C3-C4-C5	-3.07	104.72	110.20
4	A	501	GYP	C1-C2-C3	2.53	114.92	110.00

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	409/466 (87%)	0.50	49 (11%) 5 4	39, 89, 122, 155	0
2	H	213/217 (98%)	-0.28	2 (0%) 81 82	30, 43, 72, 129	0
3	L	211/211 (100%)	0.09	13 (6%) 20 17	29, 48, 79, 107	0
All	All	833/894 (93%)	0.20	64 (7%) 13 10	29, 61, 116, 155	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	132	VAL	7.6
1	A	335	LEU	6.0
1	A	130	TYR	5.4
1	A	121	PHE	5.3
1	A	336	PHE	5.2
3	L	196	THR	5.2
1	A	315	LEU	4.9
3	L	203	PRO	4.7
2	H	137	THR	4.6
1	A	12	GLY	4.3
1	A	343	GLY	4.3
1	A	370	PHE	4.1
1	A	257	TRP	4.0
1	A	457	TYR	4.0
2	H	132	VAL	4.0
1	A	65	VAL	3.9
1	A	259	MET	3.8
1	A	67	PHE	3.7
1	A	69	PHE	3.7
3	L	199	THR	3.7
1	A	314	ARG	3.6
1	A	238	ALA	3.6
3	L	1	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	258	ILE	3.5
1	A	337	LEU	3.5
1	A	129	PRO	3.5
1	A	68	ASN	3.5
1	A	376	ILE	3.4
1	A	77	LEU	3.3
3	L	209	ASN	3.3
3	L	108	ALA	3.1
1	A	63	PRO	3.1
1	A	379	LYS	3.1
1	A	261	LEU	3.0
1	A	456	VAL	3.0
3	L	195	ALA	3.0
1	A	404	ALA	3.0
1	A	440	LYS	2.9
1	A	443	LEU	2.9
1	A	133	PHE	2.9
1	A	62	LEU	2.7
1	A	405	ALA	2.7
1	A	334	LEU	2.6
1	A	447	ILE	2.6
3	L	205	VAL	2.5
3	L	200	SER	2.4
1	A	349	MET	2.4
3	L	145	VAL	2.4
1	A	131	SER	2.3
1	A	273	PHE	2.3
3	L	167	SER	2.3
1	A	393	GLU	2.3
1	A	215	ALA	2.2
1	A	219	TRP	2.2
1	A	437	GLY	2.1
3	L	197	HIS	2.1
1	A	212	LEU	2.1
1	A	128	THR	2.1
3	L	194	GLU	2.1
1	A	127	SER	2.1
1	A	444	ILE	2.1
1	A	319	ARG	2.0
1	A	330	ILE	2.0
1	A	213	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	GYP	A	501	13/13	0.18	0.06	83,91,98,107	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.