



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:31 AM GMT

PDB ID : 2DBV  
Title : GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE MUTANT  
WITH ASP 32 REPLACED BY GLY, LEU 187 REPLACED BY ALA, AND  
PRO 188 REPLACED BY SER COMPLEXED WITH NADP+  
Authors : Didierjean, C.; Rahuel-Clermont, S.; Vitoux, B.; Dideberg, O.; Branlant, G.;  
Aubry, A.  
Deposited on : 1996-12-19  
Resolution : 2.20 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](mailto: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.

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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 : 1.7 (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

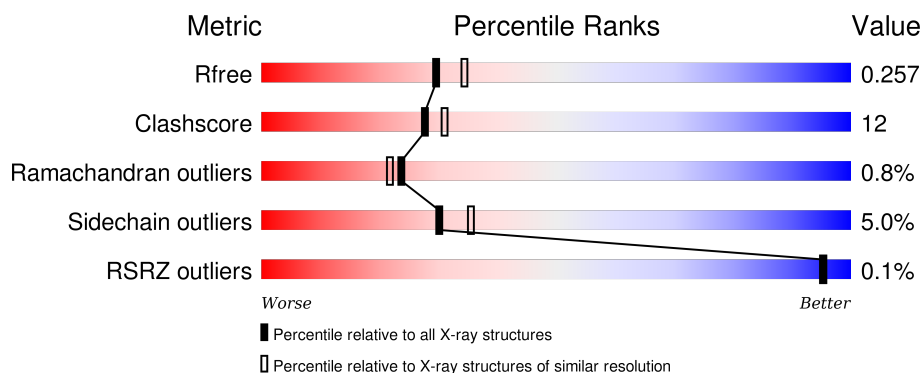
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*





The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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. 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	O	334	 72% 26% •
1	P	334	 69% 27% •
1	Q	334	 74% 25% •
1	R	334	 72% 27% •

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	O	339	-	-	-	X
2	SO4	R	339	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10693 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.

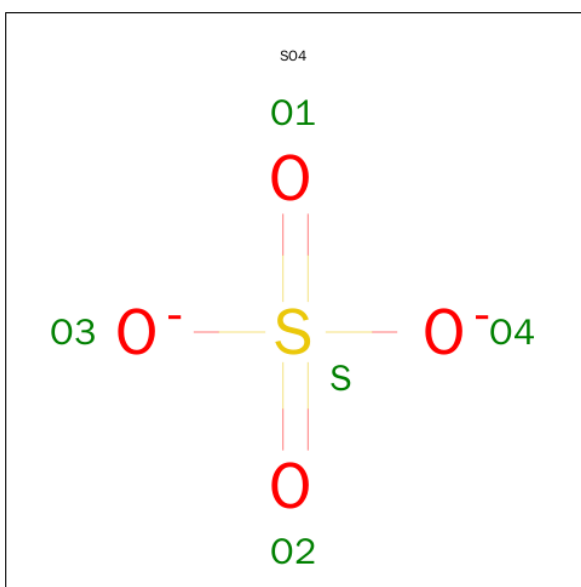
- Molecule 1 is a protein called GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	O	334	Total	C	N	O	S	0	0	0
			2517	1575	445	488	9			
1	P	334	Total	C	N	O	S	0	0	0
			2517	1575	445	488	9			
1	Q	334	Total	C	N	O	S	0	0	0
			2517	1575	445	488	9			
1	R	334	Total	C	N	O	S	0	0	0
			2517	1575	445	488	9			

There are 12 discrepancies between the modelled and reference sequences:

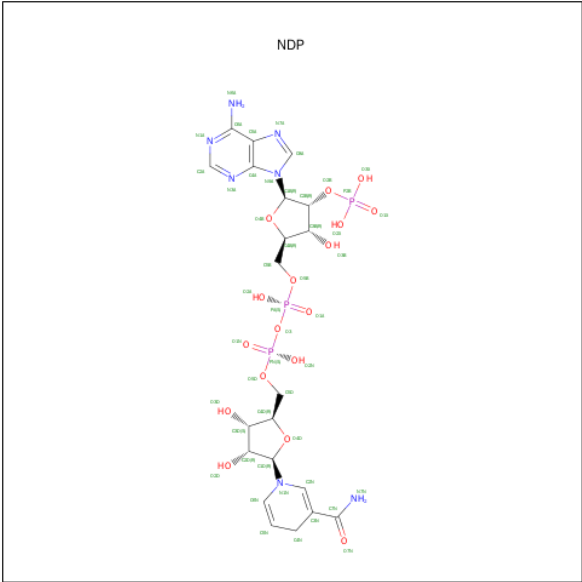
Chain	Residue	Modelled	Actual	Comment	Reference
O	32	GLY	ASP	ENGINEERED	UNP P00362
O	187	ALA	LEU	ENGINEERED	UNP P00362
O	188	SER	PRO	ENGINEERED	UNP P00362
P	32	GLY	ASP	ENGINEERED	UNP P00362
P	187	ALA	LEU	ENGINEERED	UNP P00362
P	188	SER	PRO	ENGINEERED	UNP P00362
Q	32	GLY	ASP	ENGINEERED	UNP P00362
Q	187	ALA	LEU	ENGINEERED	UNP P00362
Q	188	SER	PRO	ENGINEERED	UNP P00362
R	32	GLY	ASP	ENGINEERED	UNP P00362
R	187	ALA	LEU	ENGINEERED	UNP P00362
R	188	SER	PRO	ENGINEERED	UNP P00362

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	O	1	Total	O	S	0	0
			5	4	1		
2	O	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	Q	1	Total	O	S	0	0
			5	4	1		
2	Q	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C<sub>21</sub>H<sub>30</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



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

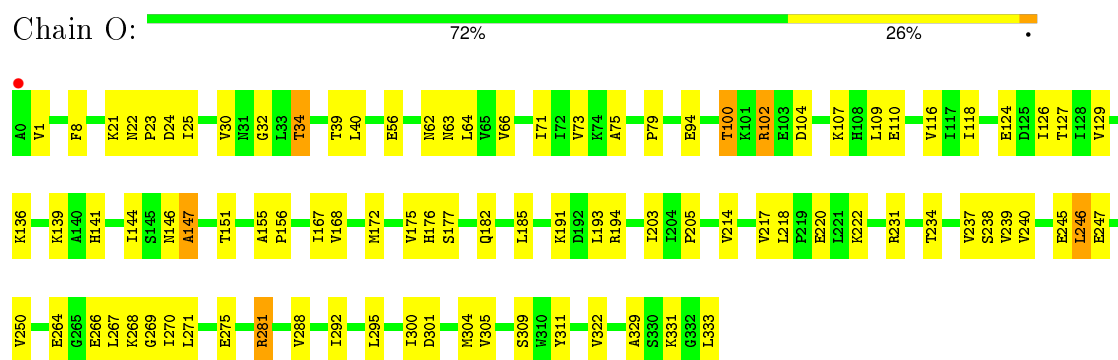
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	O	96	Total	O	0	0
			96	96		
4	P	106	Total	O	0	0
			106	106		
4	Q	87	Total	O	0	0
			87	87		
4	R	104	Total	O	0	0
			104	104		

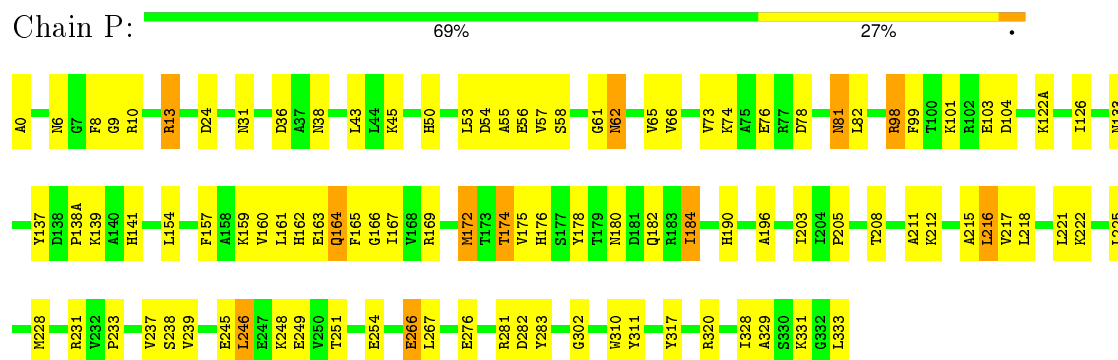
### 3 Residue-property plots

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.

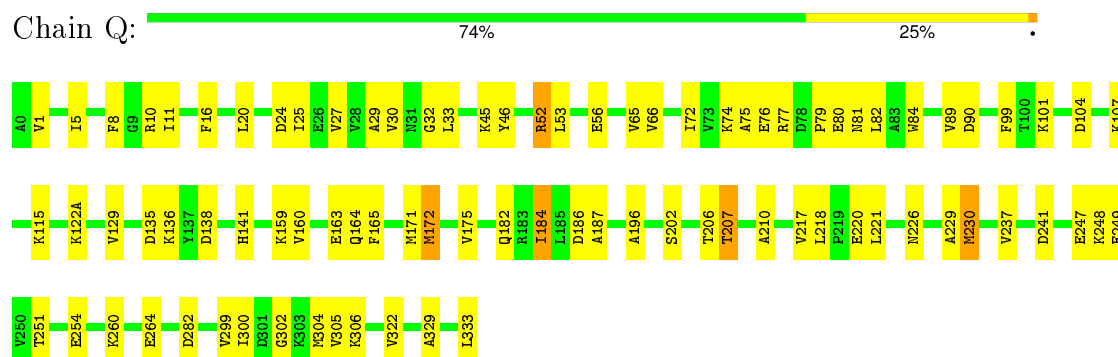
#### • Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE



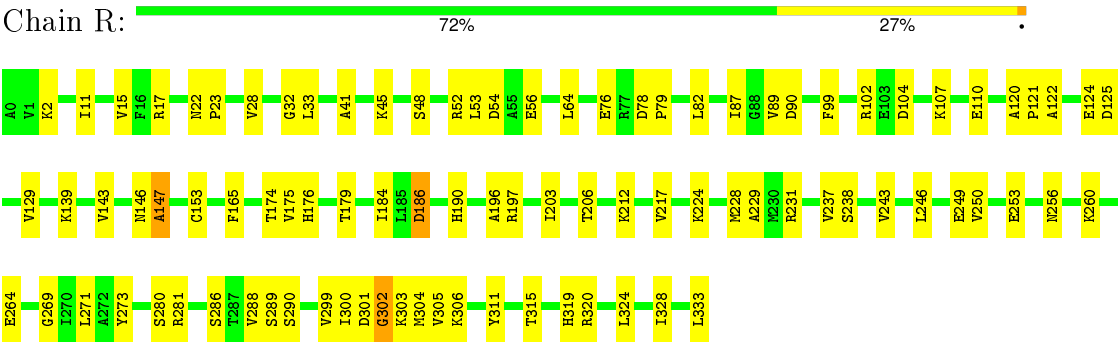
#### • Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE



#### • Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE



● Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.37Å 124.27Å 82.45Å 90.00° 108.93° 90.00°	Depositor
Resolution (Å)	8.00 – 2.20 8.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	78.8 (8.00-2.20) 83.9 (8.00-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.77 (at 2.21Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.195 , 0.265 0.197 , 0.257	Depositor DCC
$R_{free}$ test set	6665 reflections (10.17%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.3	Xtriage
Anisotropy	0.147	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 36.9	EDS
Estimated twinning fraction	0.440 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 65530 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10693	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 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	O	0.56	0/2552	0.80	2/3460 (0.1%)
1	P	0.55	0/2552	0.82	4/3460 (0.1%)
1	Q	0.55	0/2552	0.81	2/3460 (0.1%)
1	R	0.57	0/2552	0.81	1/3460 (0.0%)
All	All	0.56	0/10208	0.81	9/13840 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	203	ILE	N-CA-C	-6.56	93.29	111.00
1	Q	175	VAL	N-CA-C	-6.01	94.77	111.00
1	P	203	ILE	N-CA-C	-5.93	94.99	111.00
1	P	61	GLY	N-CA-C	-5.61	99.08	113.10
1	O	175	VAL	N-CA-C	-5.57	95.96	111.00
1	O	203	ILE	N-CA-C	-5.49	96.18	111.00
1	P	233	PRO	N-CA-C	5.33	125.96	112.10
1	Q	10	ARG	NE-CZ-NH2	5.25	122.92	120.30
1	P	175	VAL	N-CA-C	-5.08	97.28	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	O	2517	0	2563	57	0
1	P	2517	0	2563	71	0
1	Q	2517	0	2563	68	0
1	R	2517	0	2563	64	0
2	O	10	0	0	0	0
2	P	10	0	0	0	0
2	Q	10	0	0	1	0
2	R	10	0	0	0	0
3	O	48	0	26	2	0
3	P	48	0	26	3	0
3	Q	48	0	26	2	0
3	R	48	0	26	2	0
4	O	96	0	0	5	0
4	P	106	0	0	8	0
4	Q	87	0	0	2	0
4	R	104	0	0	2	0
All	All	10693	0	10356	243	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 12.

All (243) 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:O:300:ILE:HG22	1:O:304:MET:HB3	1.46	0.94
1:R:2:LYS:HE3	1:R:28:VAL:HG11	1.50	0.93
1:P:50:HIS:HD2	4:P:373:HOH:O	1.59	0.85
1:O:1:VAL:HG21	1:O:329:ALA:HB1	1.59	0.85
1:P:162:HIS:CD2	1:P:221:LEU:HD21	2.21	0.76
1:Q:76:GLU:HG2	1:Q:81:ASN:HB2	1.69	0.74
1:O:300:ILE:CG2	1:O:304:MET:HB3	2.18	0.74
1:Q:226:ASN:HB2	1:R:300:ILE:HD11	1.72	0.72
1:O:136:LYS:HE2	1:O:136:LYS:HA	1.70	0.72
1:R:17:ARG:HG2	1:R:53:LEU:HD13	1.72	0.71
1:O:8:PHE:HD2	1:O:40:LEU:HD22	1.55	0.71
1:Q:251:THR:OG1	1:Q:254:GLU:HG3	1.91	0.71
1:Q:25:ILE:HD13	1:Q:322:VAL:HG13	1.71	0.70
1:R:184:ILE:HD12	1:R:184:ILE:H	1.55	0.70
1:Q:165:PHE:HA	1:Q:248:LYS:HD3	1.73	0.69
1:Q:187:ALA:O	1:Q:196:ALA:HB1	1.92	0.69
1:P:176:HIS:HB3	1:P:231:ARG:HD3	1.75	0.68
1:P:98:ARG:HD3	4:P:409:HOH:O	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:184:ILE:HD11	1:Q:182:GLN:O	1.94	0.68
1:R:260:LYS:HE3	1:R:264:GLU:OE1	1.96	0.66
1:P:282:ASP:HB3	1:R:52:ARG:NH2	2.10	0.66
1:Q:136:LYS:HA	1:Q:136:LYS:HE2	1.77	0.65
1:O:1:VAL:HB	1:O:25:ILE:HG22	1.79	0.65
1:P:266:GLU:HG2	1:P:267:LEU:HG	1.78	0.64
1:R:32:GLY:O	1:R:33:LEU:HD23	1.98	0.64
1:O:56:GLU:O	1:O:66:VAL:HA	1.98	0.63
1:Q:80:GLU:HG2	1:Q:107:LYS:HD3	1.79	0.63
1:Q:82:LEU:HD13	1:Q:84:TRP:CZ2	2.34	0.62
1:P:218:LEU:HD13	1:P:221:LEU:HD12	1.80	0.62
1:Q:76:GLU:CG	1:Q:81:ASN:HB2	2.30	0.61
1:Q:226:ASN:CB	1:R:300:ILE:HD11	2.29	0.61
1:Q:160:VAL:O	1:Q:164:GLN:HB2	1.99	0.61
1:P:78:ASP:HB3	1:P:81:ASN:OD1	1.99	0.61
1:R:315:THR:O	1:R:319:HIS:HD2	1.83	0.61
1:Q:202:SER:OG	1:R:281:ARG:HG2	2.01	0.61
1:O:129:VAL:HG23	1:O:217:VAL:HG11	1.82	0.61
1:Q:129:VAL:HG23	1:Q:217:VAL:HG11	1.83	0.61
1:Q:76:GLU:HG2	1:Q:81:ASN:O	2.01	0.60
1:O:32:GLY:O	1:O:75:ALA:HA	2.03	0.59
1:P:9:GLY:HA3	3:P:336:NDP:O5B	2.03	0.59
1:Q:101:LYS:HG2	1:Q:122(A):LYS:HD2	1.85	0.58
1:Q:136:LYS:HE2	1:Q:136:LYS:CA	2.34	0.58
1:P:320:ARG:NE	1:P:320:ARG:HA	2.17	0.58
1:Q:306:LYS:HE2	1:R:228:MET:HG2	1.87	0.57
1:Q:11:ILE:HD11	3:Q:336:NDP:H42N	1.87	0.57
1:R:249:GLU:HA	1:R:302:GLY:O	2.05	0.56
1:O:266:GLU:HG3	1:O:267:LEU:HG	1.86	0.56
1:P:163:GLU:HB2	1:P:164:GLN:NE2	2.20	0.56
1:R:79:PRO:HD2	1:R:107:LYS:NZ	2.21	0.56
1:R:176:HIS:O	1:R:231:ARG:HA	2.06	0.56
1:Q:16:PHE:CE1	1:Q:20:LEU:HD21	2.41	0.56
1:R:260:LYS:HD3	1:R:273:TYR:CD2	2.41	0.56
1:O:79:PRO:HG2	1:O:107:LYS:HD3	1.87	0.56
1:O:126:ILE:HD12	1:O:141:HIS:CE1	2.41	0.56
1:R:122:ALA:HB3	1:R:124:GLU:HB3	1.89	0.55
1:O:8:PHE:CD2	1:O:40:LEU:HD22	2.38	0.55
1:P:62:ASN:O	1:P:73:VAL:HB	2.07	0.55
1:P:180:ASN:HA	1:Q:184:ILE:HD12	1.89	0.55
1:Q:84:TRP:HB3	1:Q:89:VAL:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:300:ILE:HG22	1:Q:304:MET:HB3	1.88	0.54
1:R:165:PHE:CD1	1:R:250:VAL:HG11	2.43	0.54
1:R:76:GLU:HB2	1:R:82:LEU:CD2	2.38	0.54
1:P:10:ARG:NH1	1:Q:186:ASP:HB2	2.23	0.54
1:R:11:ILE:O	1:R:15:VAL:HG23	2.08	0.53
1:P:53:LEU:HD23	1:P:57:VAL:CG2	2.39	0.53
1:P:101:LYS:HB3	1:P:103:GLU:OE1	2.09	0.53
1:O:240:VAL:HG23	1:O:311:TYR:CE1	2.43	0.53
1:P:162:HIS:CD2	1:P:167:ILE:H	2.26	0.53
1:R:129:VAL:HG23	1:R:217:VAL:HG11	1.91	0.53
1:Q:115:LYS:NZ	1:Q:141:HIS:O	2.41	0.53
1:P:58:SER:OG	1:P:65:VAL:HB	2.09	0.53
1:P:101:LYS:HE3	4:P:345:HOH:O	2.08	0.53
1:R:90:ASP:HB3	1:R:333:LEU:HD23	1.92	0.52
1:Q:329:ALA:HA	1:Q:333:LEU:HD13	1.91	0.52
1:P:320:ARG:HA	1:P:320:ARG:HE	1.75	0.52
1:P:0:ALA:N	1:P:24:ASP:O	2.43	0.52
1:P:329:ALA:HA	1:P:333:LEU:HD22	1.90	0.52
1:O:177:SER:HB3	1:O:234:THR:O	2.10	0.51
1:P:281:ARG:NH1	4:P:429:HOH:O	2.41	0.51
1:P:282:ASP:HB3	1:R:52:ARG:HH21	1.72	0.51
1:P:76:GLU:HB2	1:P:82:LEU:HD23	1.93	0.51
1:R:99:PHE:HD1	1:R:104:ASP:HB3	1.74	0.51
1:Q:305:VAL:HG22	1:Q:306:LYS:N	2.25	0.51
3:O:336:NDP:H1B	4:O:423:HOH:O	2.10	0.51
1:P:74:LYS:HE3	1:P:82:LEU:O	2.11	0.51
1:Q:300:ILE:CG2	1:Q:304:MET:HB3	2.41	0.50
1:Q:5:ILE:HB	1:Q:30:VAL:HG12	1.91	0.50
1:P:211:ALA:HB3	4:P:350:HOH:O	2.11	0.50
1:O:269:GLY:O	1:O:288:VAL:HG12	2.12	0.50
1:O:194:ARG:HD2	1:O:205:PRO:O	2.12	0.50
1:Q:206:THR:HG23	1:Q:229:ALA:HB3	1.92	0.50
1:R:146:ASN:O	1:R:147:ALA:HB3	2.12	0.50
1:R:315:THR:O	1:R:319:HIS:CD2	2.65	0.50
1:P:161:LEU:O	1:P:165:PHE:HB2	2.12	0.49
1:R:328:ILE:HG22	1:R:333:LEU:HD11	1.95	0.49
1:P:36:ASP:OD2	1:P:38:ASN:HB2	2.12	0.49
1:O:301:ASP:HB2	1:P:169:ARG:HD3	1.94	0.49
1:P:190:HIS:HB3	1:P:196:ALA:HB2	1.94	0.49
1:Q:207:THR:HG22	4:Q:381:HOH:O	2.11	0.49
4:Q:406:HOH:O	1:R:228:MET:HE3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:29:ALA:HB2	1:Q:72:ILE:HB	1.94	0.49
1:Q:99:PHE:HB3	1:Q:104:ASP:HB3	1.95	0.49
1:O:264:GLU:HA	1:O:268:LYS:HD3	1.95	0.49
1:R:212:LYS:HD3	4:R:355:HOH:O	2.12	0.49
1:P:6:ASN:OD1	1:P:31:ASN:HB3	2.12	0.49
1:P:138(A):PRO:HG2	1:P:139:LYS:HD2	1.95	0.49
1:Q:29:ALA:CB	1:Q:72:ILE:HB	2.42	0.49
1:R:238:SER:HB2	1:R:311:TYR:CZ	2.48	0.48
1:R:190:HIS:HB3	1:R:196:ALA:HB2	1.95	0.48
1:Q:56:GLU:O	1:Q:66:VAL:HA	2.13	0.48
1:P:238:SER:HB2	1:P:311:TYR:CZ	2.48	0.48
1:R:76:GLU:HB2	1:R:82:LEU:HD21	1.95	0.48
1:O:94:GLU:HG3	4:O:357:HOH:O	2.12	0.48
1:P:249:GLU:HA	1:P:302:GLY:O	2.14	0.48
1:R:41:ALA:HB2	1:R:64:LEU:HD22	1.96	0.48
1:Q:33:LEU:HD11	3:Q:336:NDP:C5A	2.43	0.48
1:P:56:GLU:O	1:P:66:VAL:HA	2.15	0.47
1:Q:79:PRO:HA	1:Q:82:LEU:HD12	1.95	0.47
1:O:144:ILE:HD12	1:O:144:ILE:H	1.79	0.47
1:Q:74:LYS:HE2	1:Q:82:LEU:O	2.15	0.47
1:P:76:GLU:HB2	1:P:82:LEU:CD2	2.44	0.47
1:O:100:THR:HG22	1:O:118:ILE:HG21	1.95	0.47
1:P:43:LEU:HA	4:P:344:HOH:O	2.15	0.47
1:Q:249:GLU:HA	1:Q:302:GLY:O	2.15	0.47
1:O:167:ILE:HG12	1:O:246:LEU:CD1	2.45	0.47
1:R:120:ALA:HB2	3:R:336:NDP:O3D	2.14	0.47
1:P:103:GLU:H	1:P:103:GLU:CD	2.17	0.47
1:R:186:ASP:HA	1:R:196:ALA:O	2.15	0.47
1:O:62:ASN:O	1:O:73:VAL:HB	2.15	0.47
1:Q:202:SER:HB3	1:R:280:SER:OG	2.15	0.47
1:R:107:LYS:O	1:R:110:GLU:HB2	2.14	0.47
1:P:216:LEU:N	1:P:216:LEU:HD23	2.30	0.47
1:O:155:ALA:HB3	1:O:156:PRO:HD3	1.97	0.47
1:R:269:GLY:O	1:R:288:VAL:HG12	2.15	0.47
1:O:107:LYS:O	1:O:110:GLU:HB2	2.15	0.46
1:Q:45:LYS:O	1:Q:52:ARG:HA	2.16	0.46
1:O:151:THR:HG23	1:O:214:VAL:HG22	1.96	0.46
1:O:146:ASN:O	1:O:147:ALA:HB3	2.15	0.46
1:Q:82:LEU:HD13	1:Q:84:TRP:HZ2	1.77	0.46
1:R:299:VAL:HG13	1:R:304:MET:O	2.15	0.46
1:P:281:ARG:HG2	1:R:48:SER:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:172:MET:HA	1:Q:241:ASP:O	2.16	0.46
1:P:99:PHE:HB3	1:P:104:ASP:HB3	1.97	0.46
1:O:104:ASP:O	1:O:107:LYS:HD2	2.15	0.46
1:O:271:LEU:HD11	1:O:292:ILE:HG12	1.97	0.46
1:O:168:VAL:HG22	1:O:247:GLU:HG3	1.98	0.46
1:O:176:HIS:HA	1:O:238:SER:HB3	1.98	0.45
1:O:182:GLN:HB3	4:O:360:HOH:O	2.16	0.45
1:R:289:SER:OG	1:R:320:ARG:HD2	2.16	0.45
1:R:271:LEU:HD12	1:R:290:SER:HB3	1.98	0.45
1:O:239:VAL:HG23	1:O:309:SER:O	2.14	0.45
1:P:283:TYR:CE2	1:P:310:TRP:CD1	3.05	0.45
1:O:264:GLU:HA	1:O:268:LYS:CD	2.46	0.45
1:R:243:VAL:HA	1:R:305:VAL:O	2.15	0.45
1:P:178:TYR:HA	1:P:182:GLN:OE1	2.17	0.45
1:Q:218:LEU:HB3	1:Q:221:LEU:HD12	1.98	0.45
1:Q:210:ALA:HB2	2:Q:339:SO4:O1	2.17	0.45
1:R:249:GLU:HG2	1:R:303:LYS:HG3	1.98	0.45
1:P:159:LYS:O	1:P:163:GLU:HG3	2.17	0.45
1:O:270:ILE:O	1:O:288:VAL:HB	2.17	0.45
1:O:64:LEU:HB2	1:O:71:ILE:HB	1.99	0.45
1:Q:226:ASN:HB2	1:R:300:ILE:CD1	2.44	0.45
1:P:166:GLY:O	1:P:246:LEU:HA	2.17	0.45
1:O:245:GLU:HG3	1:P:245:GLU:OE2	2.17	0.45
1:R:87:ILE:HG13	1:R:89:VAL:HG23	1.99	0.44
1:R:125:ASP:HB2	1:R:143:VAL:O	2.17	0.44
1:O:22:ASN:HD22	1:O:322:VAL:HG12	1.82	0.44
1:P:101:LYS:HG3	1:P:122(A):LYS:HE2	1.99	0.44
1:O:1:VAL:HG21	1:O:329:ALA:CB	2.38	0.44
1:P:281:ARG:NH2	4:P:385:HOH:O	2.51	0.44
1:O:144:ILE:N	1:O:144:ILE:HD12	2.32	0.44
1:Q:52:ARG:HG2	1:Q:52:ARG:HH11	1.82	0.44
1:O:331:LYS:HA	1:O:331:LYS:HD3	1.64	0.44
1:Q:90:ASP:HB3	1:Q:333:LEU:HD23	1.99	0.43
1:R:305:VAL:HG22	1:R:306:LYS:N	2.33	0.43
1:P:154:LEU:HA	1:P:157:PHE:CE2	2.52	0.43
1:P:174:THR:HA	1:P:239:VAL:O	2.18	0.43
1:P:246:LEU:N	1:P:246:LEU:HD13	2.33	0.43
1:Q:159:LYS:O	1:Q:163:GLU:HG3	2.19	0.43
1:Q:165:PHE:O	1:Q:247:GLU:HB2	2.18	0.43
1:Q:16:PHE:CE2	1:Q:27:VAL:HG11	2.53	0.43
1:R:120:ALA:HB1	1:R:121:PRO:HD2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:109:LEU:HD21	1:O:116:VAL:HG23	2.01	0.43
1:O:176:HIS:HB3	1:O:231:ARG:HD3	2.00	0.43
1:R:129:VAL:CG2	1:R:217:VAL:HG11	2.48	0.43
1:O:222:LYS:C	1:O:222:LYS:HD3	2.39	0.43
1:O:238:SER:HB2	1:O:311:TYR:CZ	2.53	0.43
1:R:153:CYS:HG	1:R:311:TYR:HD1	1.67	0.43
1:P:137:TYR:HB3	4:P:393:HOH:O	2.19	0.43
1:Q:45:LYS:HD2	1:Q:46:TYR:CZ	2.53	0.43
1:O:34:THR:CG2	1:O:39:THR:HB	2.49	0.43
1:O:295:LEU:HG	4:O:350:HOH:O	2.18	0.42
1:Q:230:MET:CE	1:R:175:VAL:HG21	2.48	0.42
1:P:215:ALA:HB2	1:P:222:LYS:HA	2.00	0.42
1:R:41:ALA:O	1:R:45:LYS:HB2	2.19	0.42
1:Q:230:MET:HE2	1:R:175:VAL:HG21	2.01	0.42
1:Q:1:VAL:HG12	1:Q:24:ASP:O	2.18	0.42
1:Q:32:GLY:O	1:Q:75:ALA:HA	2.19	0.42
1:R:206:THR:HG23	1:R:229:ALA:HB3	2.02	0.42
1:P:54:ASP:O	1:P:55:ALA:HB2	2.20	0.42
1:P:205:PRO:HB3	1:P:228:MET:HE1	2.01	0.42
1:P:10:ARG:HG2	3:P:336:NDP:O2A	2.19	0.42
1:R:146:ASN:HD22	1:R:324:LEU:HD22	1.85	0.42
1:O:193:LEU:HD23	1:O:193:LEU:HA	1.87	0.42
1:O:32:GLY:HA2	3:O:336:NDP:N3A	2.34	0.42
1:Q:25:ILE:CD1	1:Q:322:VAL:HG13	2.45	0.41
1:P:160:VAL:O	1:P:164:GLN:NE2	2.53	0.41
1:Q:20:LEU:HD12	1:Q:53:LEU:HD11	2.02	0.41
1:Q:138:ASP:H	1:Q:141:HIS:CE1	2.36	0.41
1:O:167:ILE:HG12	1:O:246:LEU:HD12	2.02	0.41
1:P:251:THR:H	1:P:254:GLU:HB2	1.85	0.41
1:R:179:THR:OG1	1:R:231:ARG:NH2	2.53	0.41
1:P:133:ASN:ND2	1:P:217:VAL:HA	2.35	0.41
1:Q:33:LEU:CD2	1:Q:77:ARG:HB3	2.49	0.41
1:Q:52:ARG:NH1	1:Q:52:ARG:HG2	2.35	0.41
1:Q:260:LYS:O	1:Q:264:GLU:HG3	2.20	0.41
1:R:174:THR:HG23	1:R:174:THR:O	2.20	0.41
1:R:22:ASN:HA	1:R:23:PRO:HD2	1.92	0.41
1:P:126:ILE:HD12	1:P:141:HIS:CE1	2.55	0.41
1:P:162:HIS:HD2	1:P:167:ILE:H	1.66	0.41
1:O:270:ILE:HG22	1:O:271:LEU:N	2.35	0.41
1:Q:138:ASP:HB3	1:Q:141:HIS:CE1	2.56	0.41
4:O:427:HOH:O	1:Q:52:ARG:HD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:185:LEU:HD23	1:O:185:LEU:HA	1.93	0.41
1:O:281:ARG:HD3	1:O:281:ARG:HA	1.90	0.41
1:Q:282:ASP:CG	1:R:197:ARG:HH22	2.23	0.41
1:Q:16:PHE:CD2	1:Q:27:VAL:HG11	2.55	0.41
1:O:102:ARG:HG3	1:O:124:GLU:HA	2.01	0.41
1:P:9:GLY:O	1:P:13:ARG:HB2	2.20	0.41
1:R:11:ILE:HD11	3:R:336:NDP:H42N	2.03	0.41
1:O:71:ILE:N	1:O:71:ILE:HD12	2.36	0.41
1:P:137:TYR:OH	1:P:328:ILE:HG23	2.21	0.41
1:P:45:LYS:HB2	1:P:45:LYS:HE2	1.90	0.41
1:R:253:GLU:H	1:R:253:GLU:CD	2.23	0.41
1:P:10:ARG:HH11	1:Q:186:ASP:HB2	1.86	0.40
1:P:208:THR:HG22	1:P:228:MET:HA	2.03	0.40
1:R:56:GLU:HA	4:R:419:HOH:O	2.20	0.40
1:P:317:TYR:O	1:P:320:ARG:HB2	2.22	0.40
1:P:31:ASN:ND2	3:P:336:NDP:H2A	2.36	0.40
1:R:99:PHE:CD1	1:R:104:ASP:HB3	2.55	0.40
1:O:139:LYS:HD3	1:O:139:LYS:HA	1.90	0.40
1:R:90:ASP:CB	1:R:333:LEU:HD23	2.51	0.40
1:P:154:LEU:HD22	1:P:172:MET:SD	2.62	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	O	332/334 (99%)	311 (94%)	18 (5%)	3 (1%)	21	19
1	P	332/334 (99%)	310 (93%)	20 (6%)	2 (1%)	30	29
1	Q	332/334 (99%)	306 (92%)	25 (8%)	1 (0%)	46	50
1	R	332/334 (99%)	314 (95%)	14 (4%)	4 (1%)	16	12
All	All	1328/1336 (99%)	1241 (93%)	77 (6%)	10 (1%)	24	22

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	147	ALA
1	O	237	VAL
1	P	237	VAL
1	Q	237	VAL
1	R	147	ALA
1	R	237	VAL
1	R	186	ASP
1	P	266	GLU
1	R	302	GLY
1	O	23	PRO

### 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	O	270/270 (100%)	252 (93%)	18 (7%)	20	21
1	P	270/270 (100%)	254 (94%)	16 (6%)	24	27
1	Q	270/270 (100%)	259 (96%)	11 (4%)	37	45
1	R	270/270 (100%)	261 (97%)	9 (3%)	45	56
All	All	1080/1080 (100%)	1026 (95%)	54 (5%)	30	35

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	21	LYS
1	O	24	ASP
1	O	30	VAL
1	O	34	THR
1	O	63	ASN
1	O	100	THR
1	O	102	ARG
1	O	127	THR
1	O	172	MET
1	O	191	LYS

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Mol	Chain	Res	Type
1	O	218	LEU
1	O	220	GLU
1	O	246	LEU
1	O	250	VAL
1	O	275	GLU
1	O	281	ARG
1	O	305	VAL
1	O	333	LEU
1	P	8	PHE
1	P	13	ARG
1	P	62	ASN
1	P	81	ASN
1	P	98	ARG
1	P	164	GLN
1	P	172	MET
1	P	174	THR
1	P	184	ILE
1	P	212	LYS
1	P	216	LEU
1	P	225	LEU
1	P	246	LEU
1	P	248	LYS
1	P	276	GLU
1	P	331	LYS
1	Q	8	PHE
1	Q	52	ARG
1	Q	65	VAL
1	Q	135	ASP
1	Q	171	MET
1	Q	172	MET
1	Q	184	ILE
1	Q	207	THR
1	Q	220	GLU
1	Q	230	MET
1	Q	299	VAL
1	R	54	ASP
1	R	78	ASP
1	R	102	ARG
1	R	139	LYS
1	R	224	LYS
1	R	246	LEU
1	R	256	ASN

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Mol	Chain	Res	Type
1	R	286	SER
1	R	301	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	O	146	ASN
1	O	152	ASN
1	O	164	GLN
1	O	256	ASN
1	P	38	ASN
1	P	50	HIS
1	P	63	ASN
1	P	146	ASN
1	P	152	ASN
1	P	162	HIS
1	P	164	GLN
1	P	256	ASN
1	Q	256	ASN
1	R	63	ASN
1	R	81	ASN
1	R	146	ASN
1	R	152	ASN
1	R	256	ASN
1	R	319	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules 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

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	NDP	O	336	-	42,52,52	1.36	4 (9%)	55,80,80	2.27	12 (21%)
2	SO4	O	338	-	4,4,4	1.39	0	6,6,6	0.57	0
2	SO4	O	339	-	4,4,4	1.24	1 (25%)	6,6,6	0.38	0
3	NDP	P	336	-	42,52,52	1.38	5 (11%)	55,80,80	2.14	10 (18%)
2	SO4	P	338	-	4,4,4	1.92	2 (50%)	6,6,6	0.74	0
2	SO4	P	339	-	4,4,4	1.03	0	6,6,6	1.02	1 (16%)
3	NDP	Q	336	-	42,52,52	1.26	4 (9%)	55,80,80	2.34	17 (30%)
2	SO4	Q	338	-	4,4,4	1.98	1 (25%)	6,6,6	0.30	0
2	SO4	Q	339	-	4,4,4	1.30	0	6,6,6	0.37	0
3	NDP	R	336	-	42,52,52	1.36	5 (11%)	55,80,80	1.82	12 (21%)
2	SO4	R	338	-	4,4,4	1.90	2 (50%)	6,6,6	0.25	0
2	SO4	R	339	-	4,4,4	1.32	0	6,6,6	0.54	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	NDP	O	336	-	-	0/30/77/77	0/5/5/5
2	SO4	O	338	-	-	0/0/0/0	0/0/0/0
2	SO4	O	339	-	-	0/0/0/0	0/0/0/0
3	NDP	P	336	-	-	0/30/77/77	0/5/5/5
2	SO4	P	338	-	-	0/0/0/0	0/0/0/0
2	SO4	P	339	-	-	0/0/0/0	0/0/0/0
3	NDP	Q	336	-	-	0/30/77/77	0/5/5/5
2	SO4	Q	338	-	-	0/0/0/0	0/0/0/0
2	SO4	Q	339	-	-	0/0/0/0	0/0/0/0
3	NDP	R	336	-	-	0/30/77/77	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	R	338	-	-	0/0/0/0	0/0/0/0
2	SO4	R	339	-	-	0/0/0/0	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	R	336	NDP	C4N-C5N	-4.64	1.39	1.49
3	O	336	NDP	C4N-C5N	-4.37	1.39	1.49
3	Q	336	NDP	C4N-C5N	-4.16	1.40	1.49
3	P	336	NDP	C4N-C5N	-2.89	1.42	1.49
2	Q	338	SO4	O1-S	-2.51	1.38	1.47
2	R	338	SO4	O4-S	-2.35	1.38	1.47
2	P	338	SO4	O1-S	-2.29	1.39	1.47
2	O	339	SO4	O1-S	-2.08	1.40	1.47
2	P	338	SO4	O2-S	-2.04	1.40	1.47
2	R	338	SO4	O2-S	-2.01	1.40	1.47
3	R	336	NDP	P2B-O2X	2.20	1.62	1.54
3	Q	336	NDP	C2N-C3N	2.23	1.40	1.34
3	R	336	NDP	C2N-C3N	2.24	1.40	1.34
3	P	336	NDP	P2B-O3X	2.31	1.63	1.54
3	Q	336	NDP	P2B-O1X	2.64	1.59	1.51
3	Q	336	NDP	C6N-C5N	2.70	1.38	1.33
3	O	336	NDP	C2N-C3N	2.73	1.41	1.34
3	R	336	NDP	C6N-C5N	2.86	1.38	1.33
3	P	336	NDP	C2N-C3N	3.19	1.42	1.34
3	P	336	NDP	P2B-O1X	3.34	1.62	1.51
3	O	336	NDP	C6N-C5N	3.53	1.40	1.33
3	R	336	NDP	P2B-O1X	3.67	1.63	1.51
3	O	336	NDP	P2B-O1X	3.68	1.63	1.51
3	P	336	NDP	C6N-C5N	3.90	1.40	1.33

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	336	NDP	N3A-C2A-N1A	-9.54	121.59	128.89
3	P	336	NDP	N3A-C2A-N1A	-8.88	122.10	128.89
3	O	336	NDP	N3A-C2A-N1A	-8.78	122.17	128.89
3	O	336	NDP	C4B-O4B-C1B	-8.62	100.25	109.72
3	R	336	NDP	N3A-C2A-N1A	-7.85	122.89	128.89
3	Q	336	NDP	C4B-O4B-C1B	-5.99	103.14	109.72
3	P	336	NDP	C4B-O4B-C1B	-5.59	103.58	109.72
3	Q	336	NDP	O4B-C4B-C5B	-4.64	92.73	109.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	336	NDP	O4B-C1B-C2B	-4.13	99.13	106.60
3	Q	336	NDP	C4N-C5N-C6N	-3.18	117.34	122.58
3	P	336	NDP	C1B-N9A-C4A	-3.16	122.18	126.94
3	R	336	NDP	O2X-P2B-O1X	-3.13	100.50	110.58
3	Q	336	NDP	C3N-C2N-N1N	-2.93	118.94	123.14
3	O	336	NDP	O3D-C3D-C4D	-2.76	102.76	111.05
3	R	336	NDP	O4B-C1B-C2B	-2.65	101.81	106.60
3	O	336	NDP	O2X-P2B-O1X	-2.60	102.20	110.58
3	Q	336	NDP	C1B-N9A-C4A	-2.59	123.03	126.94
3	O	336	NDP	O5D-PN-O1N	-2.46	100.07	109.62
3	R	336	NDP	C4N-C5N-C6N	-2.40	118.63	122.58
3	R	336	NDP	C3N-C2N-N1N	-2.37	119.75	123.14
3	P	336	NDP	P2B-O2B-C2B	-2.35	115.93	121.56
3	Q	336	NDP	C3B-C2B-C1B	-2.30	98.28	102.73
3	R	336	NDP	C3D-C2D-C1D	-2.30	96.79	101.40
3	R	336	NDP	O5D-PN-O1N	-2.27	100.79	109.62
3	O	336	NDP	C1D-N1N-C6N	-2.21	115.86	120.81
3	O	336	NDP	C3N-C2N-N1N	-2.11	120.11	123.14
3	O	336	NDP	C1B-N9A-C4A	-2.08	123.80	126.94
3	Q	336	NDP	O5B-C5B-C4B	-2.07	101.49	109.12
3	O	336	NDP	P2B-O2B-C2B	-2.06	116.63	121.56
3	Q	336	NDP	C3D-C2D-C1D	-2.04	97.30	101.40
3	R	336	NDP	O5B-C5B-C4B	-2.03	101.65	109.12
2	P	339	SO4	O2-S-O1	2.06	116.04	109.50
3	Q	336	NDP	O2A-PA-O3	2.10	114.62	105.09
3	R	336	NDP	O4D-C1D-N1N	2.23	112.78	108.07
3	Q	336	NDP	O3B-C3B-C4B	2.23	117.75	111.05
3	P	336	NDP	C2A-N1A-C6A	2.26	122.81	118.77
3	R	336	NDP	C4A-C5A-N7A	2.27	111.57	109.48
3	P	336	NDP	O2A-PA-O3	2.49	116.38	105.09
3	Q	336	NDP	C2D-C3D-C4D	2.59	107.94	102.61
3	P	336	NDP	O2B-P2B-O1X	2.61	113.63	107.11
3	Q	336	NDP	PN-O3-PA	2.69	140.29	132.73
3	R	336	NDP	O2B-P2B-O1X	2.74	113.96	107.11
3	P	336	NDP	C5N-C4N-C3N	2.78	120.19	112.52
3	O	336	NDP	C5N-C4N-C3N	2.79	120.21	112.52
3	P	336	NDP	O3-PA-O5B	2.81	110.38	102.94
3	Q	336	NDP	C2A-N1A-C6A	2.99	124.12	118.77
3	R	336	NDP	C5N-C4N-C3N	3.70	122.72	112.52
3	Q	336	NDP	C4A-C5A-N7A	3.77	112.95	109.48
3	O	336	NDP	O3X-P2B-O2X	3.80	121.84	107.38
3	Q	336	NDP	C5N-C4N-C3N	3.87	123.18	112.52

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Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	O	336	NDP	C4A-C5A-N7A	4.84	113.94	109.48
3	P	336	NDP	C4A-C5A-N7A	5.45	114.49	109.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	O	336	NDP	2	0
3	P	336	NDP	3	0
3	Q	336	NDP	2	0
2	Q	339	SO4	1	0
3	R	336	NDP	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 [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, 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	O	334/334 (100%)	-0.83	1 (0%) 94 94	7, 22, 44, 58	0
1	P	334/334 (100%)	-0.81	0 100 100	7, 22, 46, 55	0
1	Q	334/334 (100%)	-0.80	0 100 100	7, 23, 44, 60	0
1	R	334/334 (100%)	-0.83	0 100 100	6, 21, 42, 58	0
All	All	1336/1336 (100%)	-0.82	1 (0%) 95 95	6, 22, 44, 60	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	O	0	ALA	2.7

### 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, 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	O	339	5/5	0.90	0.17	5.42	54,56,58,59	0
2	SO4	R	339	5/5	0.94	0.14	3.62	64,65,67,68	0
2	SO4	P	339	5/5	0.96	0.10	1.94	43,47,49,50	0
2	SO4	Q	339	5/5	0.97	0.10	1.08	62,63,64,67	0
3	NDP	Q	336	48/48	0.96	0.09	0.07	6,17,24,29	0
3	NDP	O	336	48/48	0.97	0.08	-0.29	13,22,36,43	0
3	NDP	R	336	48/48	0.97	0.08	-0.43	9,21,30,32	0
3	NDP	P	336	48/48	0.97	0.08	-0.49	10,19,31,34	0
2	SO4	O	338	5/5	0.97	0.12	-	40,40,47,48	0
2	SO4	Q	338	5/5	0.99	0.12	-	34,34,38,40	0
2	SO4	R	338	5/5	0.97	0.13	-	49,54,56,56	0
2	SO4	P	338	5/5	0.98	0.08	-	40,44,46,47	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.