



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 05:16 AM BST

PDB ID : 3VIB
Title : Structural basis for multidrug recognition and antimicrobial resistance by MTRR, an efflux pump regulator from *Neisseria Gonorrhoeae*
Authors : Kumaraswami, M.; Shafer, W.M.; Brennan, R.G.
Deposited on : 2011-09-27
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

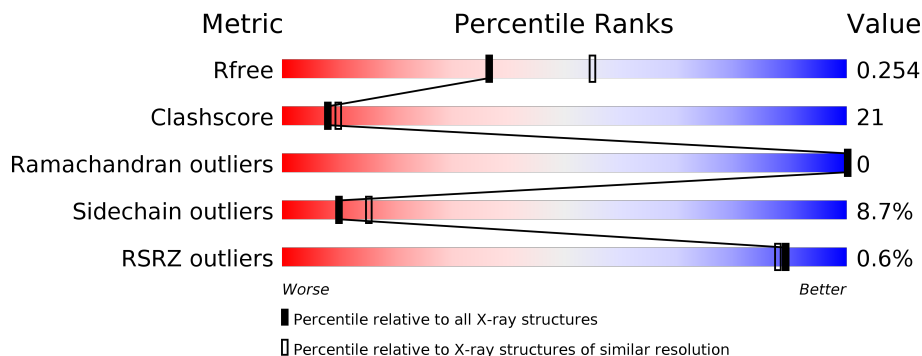
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	210	 67% 21% 8%
1	B	210	 60% 29% 8%
1	C	210	 57% 35% 8%
1	D	210	 2% 53% 32% 7% 8%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	C	212	-	-	X	-
3	CXS	A	212	-	-	X	X
3	CXS	C	211	-	-	-	X
3	CXS	D	211	-	-	-	X

2 Entry composition [i](#)

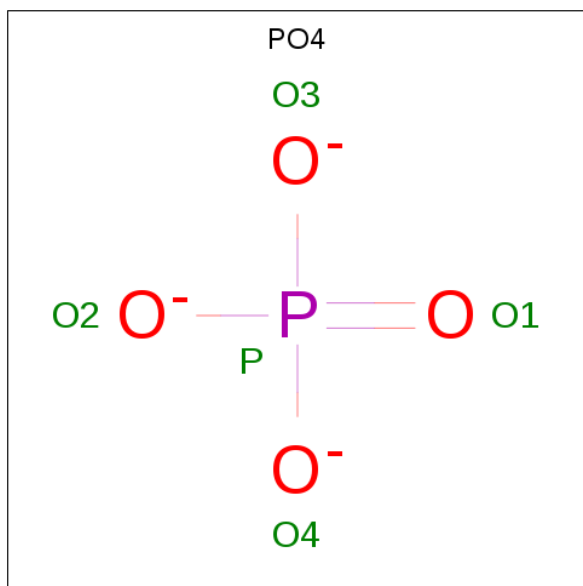
There are 4 unique types of molecules in this entry. The entry contains 6864 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MtrR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	193	Total 1570	C 1004	N 278	O 281	S 7	0	0	0
1	B	194	Total 1566	C 1001	N 279	O 279	S 7	0	0	0
1	C	201	Total 1601	C 1022	N 283	O 289	S 7	0	0	0
1	D	194	Total 1576	C 1007	N 279	O 283	S 7	0	0	0

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



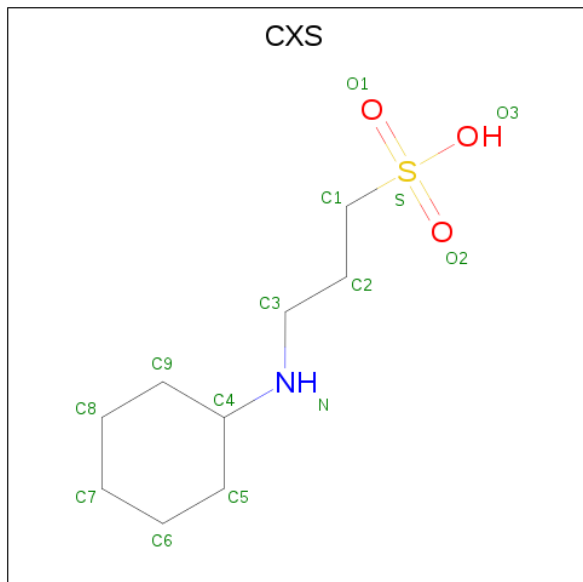
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	Total 5	O 4	P 1	0	0
2	B	1	Total 5	O 4	P 1	0	0

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

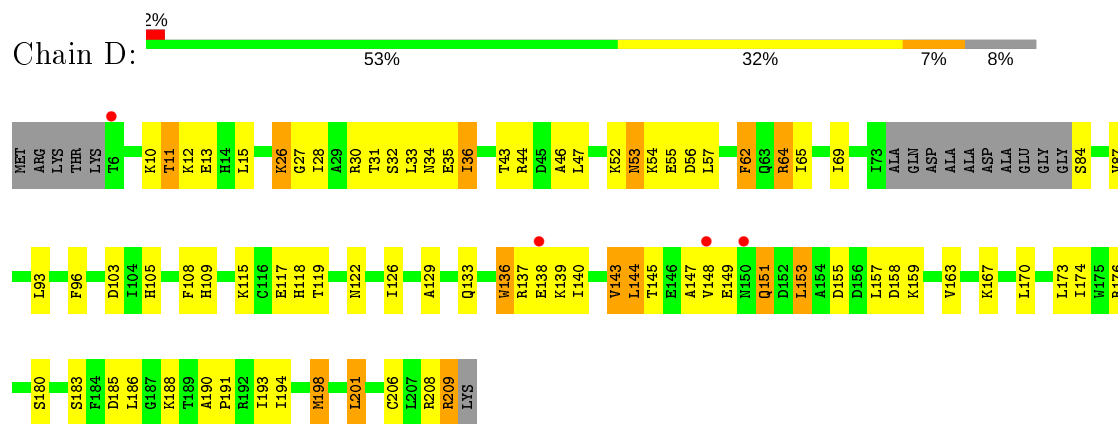
- Molecule 3 is 3-CYCLOHEXYL-1-PROPYLSULFONIC ACID (three-letter code: CXS) (formula: C₉H₁₉NO₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
3	B	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
3	C	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
3	D	1	Total	C	N	O	S	0	0
			14	9	1	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	124	Total 124	O 124	0	0
4	B	125	Total 125	O 125	0	0
4	C	79	Total 79	O 79	0	0
4	D	132	Total 132	O 132	0	0



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	218.30Å 84.60Å 58.10Å 90.00° 103.90° 90.00°	Depositor
Resolution (Å)	45.46 – 2.40 45.46 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (45.46-2.40) 95.0 (45.46-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.00Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.212 , 0.256 0.209 , 0.254	Depositor DCC
R_{free} test set	6680 reflections (10.13%)	wwPDB-VP
Wilson B-factor (Å ²)	42.6	Xtrriage
Anisotropy	0.372	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.022 for -h-2*1,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6864	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: PO4, CXS

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	0.37	0/1604	0.58	0/2165
1	B	0.36	0/1600	0.57	0/2161
1	C	0.36	0/1635	0.60	0/2211
1	D	0.39	0/1610	0.59	0/2175
All	All	0.37	0/6449	0.59	0/8712

There are no bond length outliers.

There are no bond angle outliers.

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	A	1570	0	1546	47	0
1	B	1566	0	1536	67	0
1	C	1601	0	1554	80	0
1	D	1576	0	1542	90	0
2	A	5	0	0	1	0
2	B	10	0	0	0	0
2	C	10	0	0	2	0
2	D	10	0	0	1	0
3	A	14	0	19	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	14	0	19	4	0
3	C	14	0	19	1	0
3	D	14	0	19	2	0
4	A	124	0	0	3	0
4	B	125	0	0	12	0
4	C	79	0	0	1	0
4	D	132	0	0	7	0
All	All	6864	0	6254	270	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 21.

All (270) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:TRP:CZ3	3:A:212:CXS:H62	1.23	1.62
1:A:136:TRP:CZ3	3:A:212:CXS:C6	2.17	1.26
1:D:137:ARG:HG2	1:D:167:LYS:HE3	1.36	1.07
1:D:34:ASN:OD1	1:D:44:ARG:HD3	1.62	1.00
1:D:10:LYS:HA	4:D:449:HOH:O	1.60	0.99
1:A:136:TRP:CH2	3:A:212:CXS:H62	2.01	0.95
1:A:156:ASP:HB3	1:A:208:ARG:HD2	1.50	0.93
1:A:136:TRP:HZ3	3:A:212:CXS:H62	1.15	0.92
1:D:137:ARG:HG2	1:D:167:LYS:CE	1.99	0.91
1:D:118:HIS:HA	1:D:126:ILE:HD11	1.52	0.91
1:C:192:ARG:HB2	1:D:206:CYS:SG	2.12	0.89
1:C:136:TRP:HZ3	1:C:139:LYS:HE2	1.38	0.88
1:B:170:LEU:HD13	3:B:213:CXS:H91	1.56	0.87
1:C:53:ASN:ND2	1:C:56:ASP:H	1.77	0.83
1:B:27:GLY:O	1:B:31:THR:HG23	1.79	0.82
1:A:136:TRP:CE3	3:A:212:CXS:H62	2.10	0.82
1:C:145:THR:HA	1:C:159:LYS:NZ	1.99	0.78
1:D:137:ARG:CG	1:D:167:LYS:HE3	2.15	0.76
1:A:114:LEU:HD11	1:A:178:PHE:HB3	1.68	0.75
1:B:130:ARG:HD2	4:B:443:HOH:O	1.87	0.75
1:C:102:ASN:HD22	1:C:102:ASN:C	1.90	0.75
1:C:190:ALA:HB3	1:C:191:PRO:HD3	1.69	0.75
1:B:26:LYS:NZ	1:B:30:ARG:HG3	2.02	0.75
1:A:135:ILE:HG13	1:A:136:TRP:N	2.01	0.74
1:D:96:PHE:CE2	1:D:173:LEU:HD12	2.22	0.74
1:D:36:ILE:HD13	1:D:47:LEU:HD11	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:ARG:O	1:C:30:ARG:HG3	1.88	0.74
1:B:70:GLU:HA	1:B:73:ILE:HG22	1.70	0.73
1:D:133:GLN:NE2	4:D:605:HOH:O	2.21	0.73
1:B:43:THR:HG23	1:B:46:ALA:H	1.53	0.73
1:C:32:SER:OG	1:C:35:GLU:HG3	1.88	0.73
1:C:145:THR:HA	1:C:159:LYS:HZ1	1.54	0.73
1:B:209:ARG:HG3	1:B:209:ARG:O	1.88	0.72
1:A:13:GLU:O	1:A:17:LEU:HG	1.90	0.71
1:A:132:HIS:O	1:A:136:TRP:CD1	2.45	0.70
1:A:136:TRP:CH2	3:A:212:CXS:C6	2.68	0.70
1:C:64:ARG:NH1	1:C:104:ILE:HD13	2.06	0.70
1:B:84:SER:O	1:B:87:VAL:HG22	1.92	0.70
1:D:32:SER:OG	1:D:35:GLU:HG3	1.91	0.70
1:D:159:LYS:O	1:D:163:VAL:HG23	1.91	0.69
1:B:109:HIS:HE1	4:B:398:HOH:O	1.74	0.69
1:A:132:HIS:O	1:A:136:TRP:HD1	1.75	0.69
1:D:139:LYS:O	1:D:143:VAL:HG12	1.94	0.68
1:C:118:HIS:HA	1:C:126:ILE:HD11	1.76	0.68
1:D:173:LEU:HD11	1:D:186:LEU:HD11	1.73	0.67
1:D:118:HIS:HA	1:D:126:ILE:CD1	2.24	0.67
1:D:145:THR:O	1:D:149:GLU:HG3	1.94	0.67
1:B:30:ARG:O	1:B:30:ARG:HG3	1.94	0.66
1:B:124:ALA:HB3	4:B:459:HOH:O	1.95	0.66
1:B:70:GLU:HA	1:B:73:ILE:CG2	2.25	0.66
1:D:43:THR:HG23	1:D:46:ALA:H	1.60	0.66
1:D:208:ARG:O	1:D:209:ARG:HB2	1.95	0.66
1:B:181:GLY:O	4:B:487:HOH:O	2.14	0.65
1:C:118:HIS:HA	1:C:126:ILE:CD1	2.26	0.65
1:D:173:LEU:HD11	1:D:186:LEU:CD1	2.27	0.65
1:D:133:GLN:O	1:D:137:ARG:HG3	1.97	0.64
1:C:171:ASP:OD1	3:C:211:CXS:O2	2.15	0.64
1:A:69:ILE:HD12	1:A:136:TRP:HH2	1.63	0.64
1:B:26:LYS:HB2	1:B:31:THR:HG22	1.80	0.64
1:D:12:LYS:HD2	4:D:334:HOH:O	1.99	0.63
1:B:170:LEU:CD1	3:B:213:CXS:H91	2.29	0.62
1:C:136:TRP:HA	1:C:136:TRP:CE3	2.34	0.62
1:C:53:ASN:HD22	1:C:56:ASP:H	1.47	0.62
1:C:15:LEU:HD21	1:C:42:VAL:HG21	1.81	0.62
1:C:36:ILE:HD13	1:C:57:LEU:HD23	1.82	0.61
1:C:37:ALA:HB2	1:C:47:LEU:HD23	1.82	0.61
1:B:138:GLU:HG3	4:B:469:HOH:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:53:ASN:O	1:D:56:ASP:OD1	2.18	0.61
1:C:115:LYS:HE3	1:D:118:HIS:O	2.00	0.61
1:D:137:ARG:NH2	4:D:605:HOH:O	2.33	0.60
1:D:105:HIS:O	1:D:109:HIS:HD2	1.84	0.60
1:D:117:GLU:HG3	1:D:119:THR:HG23	1.84	0.60
1:C:119:THR:HB	1:D:30:ARG:NH2	2.17	0.60
1:C:158:ASP:CG	1:C:161:THR:HG23	2.22	0.59
1:B:62:PHE:HE1	1:B:136:TRP:HZ3	1.51	0.59
1:C:121:GLN:OE1	1:D:30:ARG:HD3	2.03	0.59
1:C:136:TRP:CZ3	1:C:139:LYS:HE2	2.28	0.59
1:B:12:LYS:NZ	1:B:50:HIS:HD2	2.01	0.59
1:B:43:THR:HG22	1:B:46:ALA:HB2	1.85	0.58
1:C:55:GLU:HG2	1:C:124:ALA:HB1	1.84	0.58
1:C:102:ASN:HD22	1:C:103:ASP:N	2.00	0.58
1:B:26:LYS:HE3	1:B:35:GLU:OE1	2.03	0.58
1:B:10:LYS:N	4:B:426:HOH:O	2.36	0.58
1:C:136:TRP:HZ3	1:C:139:LYS:CE	2.14	0.58
1:B:100:GLN:HG3	1:B:100:GLN:O	2.02	0.58
1:D:129:ALA:O	1:D:133:GLN:HG3	2.04	0.57
1:B:137:ARG:HG3	1:B:163:VAL:HG11	1.85	0.57
1:C:136:TRP:HA	1:C:136:TRP:HE3	1.69	0.57
1:B:173:LEU:HD13	1:B:193:ILE:HD12	1.87	0.57
1:C:70:GLU:CD	1:C:139:LYS:HE3	2.25	0.57
1:C:53:ASN:ND2	1:C:56:ASP:HB2	2.20	0.57
1:D:53:ASN:ND2	1:D:55:GLU:HB2	2.20	0.56
1:C:46:ALA:O	1:C:49:TRP:HB3	2.06	0.56
1:D:30:ARG:HD2	4:D:646:HOH:O	2.04	0.56
1:D:185:ASP:OD1	1:D:188:LYS:HD3	2.05	0.56
1:B:26:LYS:HZ2	1:B:30:ARG:HG3	1.70	0.56
1:A:27:GLY:HA2	1:A:115:LYS:HB3	1.87	0.55
1:C:176:ARG:HH12	1:D:137:ARG:NH2	2.05	0.55
1:C:64:ARG:HH12	1:C:104:ILE:CD1	2.19	0.55
1:A:137:ARG:HD3	2:A:211:PO4:O3	2.06	0.55
1:D:147:ALA:HB3	1:D:153:LEU:HD22	1.87	0.55
1:C:70:GLU:OE2	1:C:139:LYS:HE3	2.07	0.55
1:B:156:ASP:O	1:B:209:ARG:N	2.30	0.55
1:D:147:ALA:CB	1:D:153:LEU:HD22	2.36	0.55
1:C:64:ARG:NH1	1:C:104:ILE:CD1	2.70	0.55
1:C:15:LEU:HD21	1:C:42:VAL:CG2	2.37	0.54
1:D:93:LEU:HD23	1:D:194:ILE:HD12	1.90	0.54
1:D:198:MET:HE3	1:D:201:LEU:HD23	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:ALA:HB3	1:B:191:PRO:HD3	1.90	0.53
1:D:140:ILE:O	1:D:143:VAL:HG13	2.08	0.53
1:D:84:SER:O	1:D:87:VAL:HG22	2.08	0.53
1:A:137:ARG:HG3	1:A:163:VAL:HG11	1.90	0.53
1:C:137:ARG:HB2	1:C:167:LYS:HE3	1.90	0.53
1:C:59:ASP:OD1	1:C:132:HIS:HE1	1.92	0.53
1:A:29:ALA:HB2	1:A:122:ASN:HD21	1.73	0.53
1:A:131:LYS:NZ	1:A:131:LYS:HB3	2.23	0.53
1:D:137:ARG:HG2	1:D:167:LYS:HE2	1.89	0.53
1:D:153:LEU:HG	1:D:157:LEU:HD22	1.90	0.53
1:D:47:LEU:C	1:D:47:LEU:HD23	2.28	0.53
1:A:136:TRP:CE3	3:A:212:CXS:C5	2.91	0.53
1:B:129:ALA:O	1:B:133:GLN:HG3	2.10	0.52
1:B:210:LYS:C	4:B:456:HOH:O	2.47	0.52
1:A:136:TRP:CE3	3:A:212:CXS:C6	2.84	0.52
1:C:65:ILE:HG13	1:C:108:PHE:CD2	2.44	0.52
1:D:208:ARG:HH11	1:D:208:ARG:HG3	1.74	0.52
1:D:198:MET:CE	1:D:201:LEU:HD23	2.39	0.52
1:A:136:TRP:CZ3	3:A:212:CXS:C5	2.88	0.52
1:B:105:HIS:O	1:B:109:HIS:HD2	1.92	0.51
1:B:114:LEU:HD11	1:B:178:PHE:HB3	1.92	0.51
1:C:102:ASN:ND2	1:C:104:ILE:H	2.07	0.51
1:B:43:THR:CG2	1:B:46:ALA:H	2.23	0.51
1:D:140:ILE:HA	1:D:143:VAL:CG1	2.39	0.51
1:C:166:ILE:O	1:C:170:LEU:HG	2.11	0.51
1:C:176:ARG:HH12	1:D:137:ARG:HH21	1.58	0.51
1:C:16:MET:HE1	1:C:60:ALA:HB1	1.93	0.51
1:A:126:ILE:O	1:A:130:ARG:HG2	2.12	0.50
1:B:136:TRP:NE1	3:B:213:CXS:H61	2.27	0.50
1:D:54:LYS:NZ	2:D:213:PO4:O4	2.45	0.50
1:C:102:ASN:C	1:C:102:ASN:ND2	2.62	0.49
1:A:122:ASN:O	1:A:126:ILE:HG12	2.12	0.49
1:B:122:ASN:ND2	4:B:472:HOH:O	2.45	0.49
1:B:137:ARG:CG	1:B:163:VAL:HG11	2.41	0.49
1:C:154:ALA:C	1:C:156:ASP:H	2.16	0.49
1:D:144:LEU:O	1:D:148:VAL:HG13	2.11	0.49
1:D:28:ILE:O	1:D:54:LYS:NZ	2.44	0.49
1:A:32:SER:OG	1:A:35:GLU:HG3	2.12	0.49
1:C:137:ARG:NE	2:C:212:PO4:O3	2.43	0.49
1:D:138:GLU:HG2	1:D:139:LYS:N	2.28	0.49
1:C:33:LEU:HD21	1:C:54:LYS:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:ARG:HG3	1:D:209:ARG:NH1	2.29	0.48
1:C:148:VAL:HG23	1:C:149:GLU:N	2.29	0.48
1:C:133:GLN:HG3	2:C:212:PO4:O2	2.14	0.48
1:C:173:LEU:HD11	1:C:190:ALA:HA	1.96	0.48
1:D:136:TRP:HA	1:D:136:TRP:CE3	2.49	0.48
1:A:128:ILE:O	1:A:131:LYS:HG2	2.14	0.48
1:D:208:ARG:NH1	1:D:208:ARG:HG3	2.29	0.48
1:D:69:ILE:HD13	3:D:211:CXS:H61	1.96	0.48
1:A:167:LYS:HB3	1:B:176:ARG:HD3	1.97	0.47
4:B:463:HOH:O	1:C:44:ARG:HG2	2.14	0.47
1:C:53:ASN:HD21	1:C:56:ASP:HB2	1.80	0.47
1:D:96:PHE:HE2	1:D:173:LEU:HD12	1.77	0.47
1:A:172:GLY:O	1:A:176:ARG:HB2	2.14	0.47
1:B:66:CYS:O	1:B:70:GLU:HG2	2.15	0.47
1:D:170:LEU:O	1:D:174:ILE:HG13	2.15	0.47
1:A:44:ARG:HG2	4:A:455:HOH:O	2.14	0.47
1:B:182:GLU:HB2	4:B:450:HOH:O	2.15	0.47
1:C:43:THR:HG23	1:C:46:ALA:H	1.80	0.46
1:D:159:LYS:NZ	4:D:621:HOH:O	2.49	0.46
1:D:36:ILE:HD12	1:D:57:LEU:HD13	1.97	0.46
1:D:52:LYS:N	1:D:56:ASP:OD2	2.47	0.46
1:A:176:ARG:HG3	1:B:168:SER:HA	1.97	0.46
1:B:43:THR:HG22	1:B:46:ALA:CB	2.44	0.46
1:D:158:ASP:HB2	1:D:209:ARG:HG2	1.97	0.46
1:B:182:GLU:C	4:B:487:HOH:O	2.53	0.46
1:C:85:TRP:CZ2	1:C:153:LEU:HD21	2.51	0.46
1:A:176:ARG:HD2	1:B:167:LYS:HB3	1.98	0.46
1:D:27:GLY:HA2	1:D:115:LYS:HB3	1.98	0.46
1:C:149:GLU:C	1:C:151:GLN:H	2.20	0.45
1:C:53:ASN:HD21	1:C:56:ASP:H	1.62	0.45
1:C:51:PHE:CD1	1:C:57:LEU:HD13	2.52	0.45
1:D:190:ALA:HB3	1:D:191:PRO:HD3	1.97	0.45
1:D:173:LEU:HD23	1:D:193:ILE:HD12	1.99	0.45
1:A:72:CYS:C	1:A:73:ILE:HG13	2.36	0.45
1:C:12:LYS:HE3	1:C:51:PHE:CZ	2.52	0.45
1:A:8:ALA:O	1:A:11:THR:HG22	2.16	0.45
1:C:25:ARG:NH2	4:C:592:HOH:O	2.47	0.44
1:A:114:LEU:CD1	1:A:178:PHE:HB3	2.44	0.44
1:D:26:LYS:O	1:D:30:ARG:NE	2.50	0.44
1:D:26:LYS:HA	1:D:30:ARG:NH2	2.32	0.44
1:C:62:PHE:CD2	1:C:108:PHE:HZ	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:48:TYR:HA	1:C:51:PHE:O	2.18	0.44
1:C:188:LYS:O	1:C:192:ARG:HD2	2.18	0.44
1:C:121:GLN:HG3	1:D:30:ARG:NH1	2.33	0.44
1:D:64:ARG:O	1:D:64:ARG:HG3	2.17	0.44
1:B:100:GLN:NE2	4:B:519:HOH:O	2.50	0.44
1:A:131:LYS:O	1:A:135:ILE:HG23	2.18	0.43
3:B:213:CXS:H51	3:B:213:CXS:H32	1.86	0.43
1:C:36:ILE:HG21	1:C:57:LEU:HD21	1.99	0.43
1:B:70:GLU:CA	1:B:73:ILE:HG22	2.45	0.43
1:D:11:THR:O	1:D:15:LEU:HG	2.18	0.43
1:D:31:THR:O	1:D:54:LYS:HD3	2.17	0.43
1:C:130:ARG:HA	1:C:133:GLN:HB3	1.99	0.43
1:C:36:ILE:HD13	1:C:57:LEU:CD2	2.47	0.43
1:D:209:ARG:HH11	1:D:209:ARG:CG	2.31	0.43
1:B:144:LEU:O	1:B:148:VAL:HG23	2.19	0.43
1:A:206:CYS:O	1:B:192:ARG:HD3	2.19	0.43
1:C:107:LYS:O	1:C:111:ILE:HG13	2.19	0.43
1:A:51:PHE:HB3	1:A:57:LEU:HD22	1.99	0.43
1:C:16:MET:CE	1:C:60:ALA:HB1	2.48	0.43
1:A:131:LYS:NZ	1:A:131:LYS:CB	2.82	0.43
1:C:18:ALA:HB1	1:C:39:ALA:HB3	1.99	0.43
1:D:167:LYS:NZ	3:D:211:CXS:O2	2.47	0.43
1:A:29:ALA:HB2	1:A:122:ASN:ND2	2.34	0.43
1:B:12:LYS:HZ1	1:B:50:HIS:HD2	1.67	0.43
1:B:59:ASP:HB2	1:B:128:ILE:HD13	2.00	0.43
1:B:173:LEU:HD13	1:B:193:ILE:CD1	2.48	0.43
1:B:98:ARG:C	1:B:100:GLN:N	2.72	0.43
1:B:158:ASP:HA	1:B:209:ARG:HB3	2.00	0.43
1:A:199:ASP:HB3	4:A:278:HOH:O	2.18	0.42
1:B:204:HIS:HA	1:B:205:PRO:HD3	1.77	0.42
1:C:119:THR:HB	1:D:30:ARG:HH22	1.82	0.42
1:B:204:HIS:ND1	1:B:206:CYS:HB2	2.34	0.42
1:D:36:ILE:HD12	1:D:57:LEU:CD1	2.49	0.42
1:D:140:ILE:HA	1:D:143:VAL:HG12	2.00	0.42
1:B:177:TRP:CZ2	1:B:182:GLU:HG2	2.55	0.42
1:A:192:ARG:HH21	1:B:161:THR:HG21	1.85	0.42
1:D:53:ASN:C	1:D:53:ASN:HD22	2.23	0.42
1:C:145:THR:HA	1:C:159:LYS:HZ3	1.83	0.42
1:D:53:ASN:ND2	1:D:53:ASN:N	2.68	0.42
1:A:18:ALA:O	1:A:22:THR:HG22	2.20	0.42
1:C:154:ALA:C	1:C:156:ASP:N	2.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:138:GLU:CG	1:D:139:LYS:N	2.83	0.42
1:A:70:GLU:C	1:A:72:CYS:H	2.22	0.42
1:A:147:ALA:HB3	1:A:153:LEU:HD12	2.02	0.41
1:D:180:SER:OG	1:D:183:SER:HB3	2.20	0.41
1:D:62:PHE:CD2	1:D:108:PHE:HZ	2.38	0.41
1:B:70:GLU:OE2	1:B:139:LYS:NZ	2.47	0.41
1:C:158:ASP:OD1	1:C:161:THR:HG23	2.20	0.41
1:A:93:LEU:O	1:A:97:GLU:HG3	2.20	0.41
1:B:55:GLU:HB3	1:B:128:ILE:HD12	2.02	0.41
1:B:13:GLU:O	1:B:17:LEU:HG	2.19	0.41
1:B:208:ARG:HA	1:B:208:ARG:HD3	1.78	0.41
1:D:209:ARG:NH1	1:D:209:ARG:CG	2.83	0.41
1:B:128:ILE:O	1:B:131:LYS:HG2	2.20	0.41
1:C:15:LEU:HD22	1:C:47:LEU:HD22	2.02	0.41
1:B:12:LYS:HZ2	1:B:50:HIS:HD2	1.68	0.41
1:B:62:PHE:HE1	1:B:136:TRP:CZ3	2.35	0.41
1:C:58:PHE:CD2	1:C:125:VAL:HG13	2.56	0.41
1:D:155:ASP:HB3	4:D:661:HOH:O	2.20	0.41
1:B:27:GLY:HA2	1:B:115:LYS:HB3	2.01	0.41
1:D:32:SER:HG	1:D:35:GLU:HG3	1.84	0.41
1:C:24:TYR:O	1:C:115:LYS:HE2	2.21	0.41
1:D:65:ILE:HG13	1:D:108:PHE:CD2	2.56	0.41
1:A:204:HIS:CG	1:A:205:PRO:HD2	2.56	0.41
1:B:136:TRP:O	1:B:140:ILE:HG13	2.21	0.41
1:C:27:GLY:HA2	1:C:115:LYS:HB3	2.02	0.41
1:D:136:TRP:HE3	1:D:136:TRP:HA	1.86	0.41
1:D:140:ILE:O	1:D:143:VAL:CG1	2.68	0.41
1:B:55:GLU:HB3	1:B:128:ILE:CD1	2.51	0.40
1:C:204:HIS:HA	1:C:205:PRO:HD2	1.88	0.40
1:A:63:GLN:HG3	4:A:246:HOH:O	2.21	0.40
1:C:112:LEU:HA	1:C:112:LEU:HD23	1.94	0.40
1:D:151:GLN:HE21	1:D:151:GLN:HB3	1.62	0.40
1:A:117:GLU:HA	1:B:114:LEU:O	2.21	0.40
1:B:37:ALA:HB2	1:B:47:LEU:HD12	2.02	0.40
1:C:53:ASN:C	1:C:53:ASN:HD22	2.24	0.40
1:D:145:THR:HA	1:D:159:LYS:HE2	2.03	0.40
1:D:198:MET:HA	1:D:198:MET:CE	2.52	0.40
1:C:121:GLN:CD	1:D:30:ARG:HD3	2.42	0.40
1:D:26:LYS:C	1:D:30:ARG:HH21	2.24	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	189/210 (90%)	184 (97%)	5 (3%)	0	100	100
1	B	190/210 (90%)	185 (97%)	5 (3%)	0	100	100
1	C	197/210 (94%)	190 (96%)	7 (4%)	0	100	100
1	D	190/210 (90%)	187 (98%)	3 (2%)	0	100	100
All	All	766/840 (91%)	746 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	164/179 (92%)	151 (92%)	13 (8%)	12	19
1	B	162/179 (90%)	151 (93%)	11 (7%)	16	25
1	C	163/179 (91%)	149 (91%)	14 (9%)	10	16
1	D	164/179 (92%)	145 (88%)	19 (12%)	5	7
All	All	653/716 (91%)	596 (91%)	57 (9%)	10	15

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	THR
1	A	26	LYS

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Mol	Chain	Res	Type
1	A	43	THR
1	A	62	PHE
1	A	70	GLU
1	A	73	ILE
1	A	103	ASP
1	A	130	ARG
1	A	131	LYS
1	A	138	GLU
1	A	149	GLU
1	A	176	ARG
1	A	192	ARG
1	B	20	LEU
1	B	44	ARG
1	B	62	PHE
1	B	70	GLU
1	B	100	GLN
1	B	103	ASP
1	B	122	ASN
1	B	136	TRP
1	B	176	ARG
1	B	206	CYS
1	B	209	ARG
1	C	20	LEU
1	C	38	GLN
1	C	47	LEU
1	C	53	ASN
1	C	62	PHE
1	C	93	LEU
1	C	94	HIS
1	C	102	ASN
1	C	112	LEU
1	C	133	GLN
1	C	136	TRP
1	C	161	THR
1	C	173	LEU
1	C	207	LEU
1	D	11	THR
1	D	13	GLU
1	D	26	LYS
1	D	33	LEU
1	D	36	ILE
1	D	53	ASN

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Mol	Chain	Res	Type
1	D	62	PHE
1	D	64	ARG
1	D	103	ASP
1	D	122	ASN
1	D	136	TRP
1	D	143	VAL
1	D	144	LEU
1	D	151	GLN
1	D	153	LEU
1	D	176	ARG
1	D	198	MET
1	D	201	LEU
1	D	209	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	HIS
1	A	94	HIS
1	A	105	HIS
1	A	200	ASN
1	B	38	GLN
1	B	50	HIS
1	B	109	HIS
1	B	122	ASN
1	B	200	ASN
1	C	53	ASN
1	C	63	GLN
1	C	102	ASN
1	C	132	HIS
1	C	133	GLN
1	C	200	ASN
1	D	53	ASN
1	D	109	HIS
1	D	122	ASN
1	D	133	GLN
1	D	150	ASN
1	D	151	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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	PO4	B	211	-	4,4,4	1.61	0	6,6,6	0.43	0
2	PO4	C	212	-	4,4,4	0.91	0	6,6,6	0.42	0
2	PO4	B	212	-	4,4,4	1.65	0	6,6,6	0.44	0
3	CXS	D	211	-	14,14,14	2.14	1 (7%)	18,18,18	1.35	5 (27%)
2	PO4	C	213	-	4,4,4	0.92	0	6,6,6	0.43	0
3	CXS	A	212	-	14,14,14	2.13	1 (7%)	18,18,18	1.35	5 (27%)
3	CXS	C	211	-	14,14,14	2.13	1 (7%)	18,18,18	1.35	5 (27%)
3	CXS	B	213	-	14,14,14	2.14	1 (7%)	18,18,18	1.35	5 (27%)
2	PO4	D	213	-	4,4,4	1.68	0	6,6,6	0.42	0
2	PO4	A	211	-	4,4,4	1.68	0	6,6,6	0.43	0
2	PO4	D	212	-	4,4,4	1.69	0	6,6,6	0.43	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	CXS	A	212	-	-	2/8/16/16	0/1/1/1
3	CXS	C	211	-	-	6/8/16/16	0/1/1/1
3	CXS	B	213	-	-	6/8/16/16	0/1/1/1
3	CXS	D	211	-	-	6/8/16/16	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	213	CXS	C1-S	-7.67	1.66	1.77
3	D	211	CXS	C1-S	-7.67	1.66	1.77
3	C	211	CXS	C1-S	-7.62	1.66	1.77
3	A	212	CXS	C1-S	-7.60	1.66	1.77

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	211	CXS	C3-N-C4	2.99	120.01	114.14
3	D	211	CXS	C3-N-C4	2.99	120.01	114.14
3	A	212	CXS	C3-N-C4	2.99	120.00	114.14
3	B	213	CXS	C3-N-C4	2.98	120.00	114.14
3	A	212	CXS	C2-C1-S	-2.46	109.47	113.25
3	D	211	CXS	C2-C1-S	-2.44	109.51	113.25
3	B	213	CXS	C2-C1-S	-2.43	109.52	113.25
3	C	211	CXS	C2-C1-S	-2.42	109.53	113.25
3	B	213	CXS	O3-S-C1	2.31	109.50	105.77
3	A	212	CXS	O3-S-C1	2.30	109.48	105.77
3	D	211	CXS	O3-S-C1	2.29	109.46	105.77
3	C	211	CXS	O3-S-C1	2.28	109.46	105.77
3	B	213	CXS	O1-S-C1	2.13	109.48	106.92
3	B	213	CXS	O2-S-C1	2.12	109.47	106.92
3	A	212	CXS	O1-S-C1	2.12	109.47	106.92
3	D	211	CXS	O1-S-C1	2.11	109.46	106.92
3	D	211	CXS	O2-S-C1	2.11	109.46	106.92
3	C	211	CXS	O1-S-C1	2.11	109.46	106.92
3	C	211	CXS	O2-S-C1	2.09	109.44	106.92
3	A	212	CXS	O2-S-C1	2.09	109.43	106.92

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	211	CXS	C2-C1-S-O2
3	A	212	CXS	C1-C2-C3-N
3	C	211	CXS	C2-C1-S-O1
3	C	211	CXS	C2-C1-S-O2
3	C	211	CXS	S-C1-C2-C3
3	B	213	CXS	C2-C1-S-O2
3	B	213	CXS	C5-C4-N-C3
3	D	211	CXS	C2-C1-S-O3
3	C	211	CXS	C2-C1-S-O3
3	D	211	CXS	S-C1-C2-C3
3	D	211	CXS	C5-C4-N-C3
3	C	211	CXS	C5-C4-N-C3
3	B	213	CXS	C2-C3-N-C4
3	D	211	CXS	C2-C1-S-O1
3	B	213	CXS	C2-C1-S-O1
3	B	213	CXS	C2-C1-S-O3
3	A	212	CXS	S-C1-C2-C3
3	D	211	CXS	C9-C4-N-C3
3	C	211	CXS	C9-C4-N-C3
3	B	213	CXS	C9-C4-N-C3

There are no ring outliers.

7 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	212	PO4	2	0
3	D	211	CXS	2	0
3	A	212	CXS	9	0
3	C	211	CXS	1	0
3	B	213	CXS	4	0
2	D	213	PO4	1	0
2	A	211	PO4	1	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, 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	193/210 (91%)	-0.34	1 (0%) 91 89	26, 45, 67, 81	0
1	B	194/210 (92%)	-0.33	0 100 100	27, 45, 64, 70	0
1	C	201/210 (95%)	-0.50	0 100 100	28, 47, 67, 96	0
1	D	194/210 (92%)	-0.26	4 (2%) 63 61	28, 46, 70, 94	0
All	All	782/840 (93%)	-0.36	5 (0%) 89 88	26, 46, 67, 96	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	6	THR	3.9
1	D	148	VAL	2.6
1	D	150	ASN	2.5
1	A	136	TRP	2.2
1	D	138	GLU	2.1

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 monosaccharides 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. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CXS	A	212	14/14	0.37	0.48	134,135,140,140	0
3	CXS	C	211	14/14	0.41	0.48	146,147,152,153	0
3	CXS	D	211	14/14	0.50	0.40	138,139,140,141	0
3	CXS	B	213	14/14	0.71	0.30	127,128,132,132	0
2	PO4	C	213	5/5	0.74	0.25	119,120,120,120	0
2	PO4	D	213	5/5	0.77	0.27	136,136,136,136	0
2	PO4	A	211	5/5	0.80	0.23	111,111,112,112	0
2	PO4	C	212	5/5	0.88	0.14	107,107,108,108	0
2	PO4	B	212	5/5	0.89	0.18	103,104,104,104	0
2	PO4	D	212	5/5	0.97	0.13	83,83,83,83	0
2	PO4	B	211	5/5	0.98	0.11	66,67,68,69	0

6.5 Other polymers

There are no such residues in this entry.