Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 05:18 AM EDT

PDB ID	:	5ODW
Title	:	Structure of the FpvAI-pyocin S2 complex
Authors	:	White, P.; Joshi, A.; Kleanthous, C.
Deposited on	:	2017-07-06
Resolution	:	2.80 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

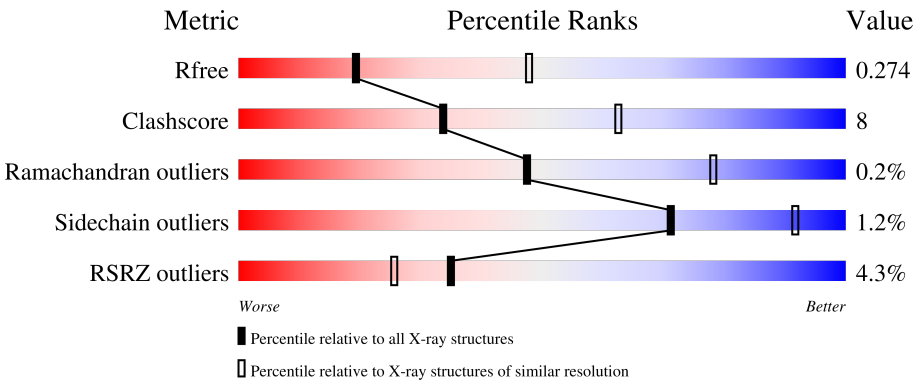
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of 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 <=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	815	% 80% 14% 5%
1	B	815	4% 75% 16% 9%
2	C	217	10% 74% 15% 10%
2	D	217	11% 76% 16% 9%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 15094 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 Ferripyoverdine receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	771	6105	3840	1044	1209	12	0	0	0
1	B	740	5874	3696	1008	1158	12	0	0	0

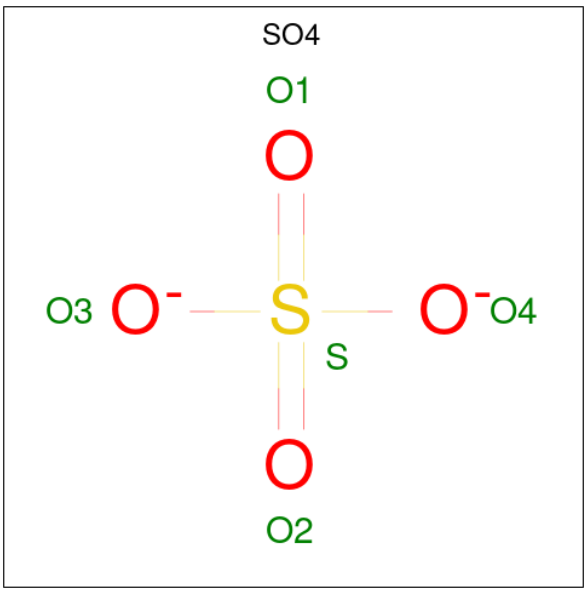
- Molecule 2 is a protein called Pyocin-S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	195	1520	950	264	301	5	0	0	0
2	D	198	1545	966	267	307	5	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	210	LEU	-	expression tag	UNP Q06584
C	211	GLU	-	expression tag	UNP Q06584
C	212	HIS	-	expression tag	UNP Q06584
C	213	HIS	-	expression tag	UNP Q06584
C	214	HIS	-	expression tag	UNP Q06584
C	215	HIS	-	expression tag	UNP Q06584
C	216	HIS	-	expression tag	UNP Q06584
C	217	HIS	-	expression tag	UNP Q06584
D	210	LEU	-	expression tag	UNP Q06584
D	211	GLU	-	expression tag	UNP Q06584
D	212	HIS	-	expression tag	UNP Q06584
D	213	HIS	-	expression tag	UNP Q06584
D	214	HIS	-	expression tag	UNP Q06584
D	215	HIS	-	expression tag	UNP Q06584
D	216	HIS	-	expression tag	UNP Q06584
D	217	HIS	-	expression tag	UNP Q06584

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



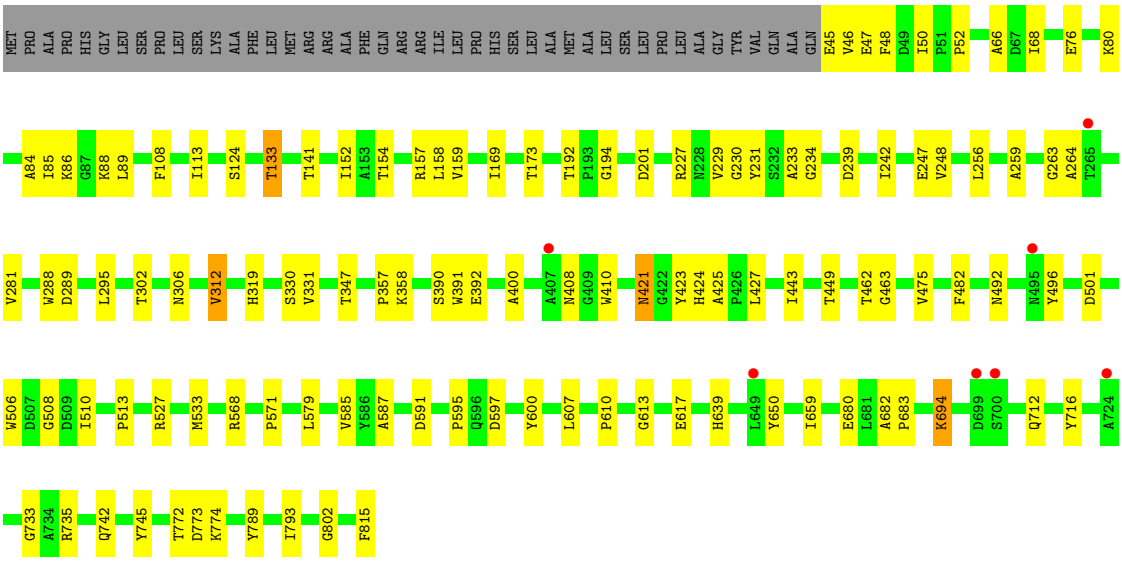
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	O	S			
3	A	1	5	4	1		0	0
			Total	O	S			
3	A	1	5	4	1		0	0
			Total	O	S			
3	A	1	5	4	1		0	0
			Total	O	S			
3	A	1	5	4	1		0	0
			Total	O	S			
3	A	1	5	4	1		0	0
			Total	O	S			
3	B	1	5	4	1		0	0
			Total	O	S			
3	B	1	5	4	1		0	0
			Total	O	S			
3	B	1	5	4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

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

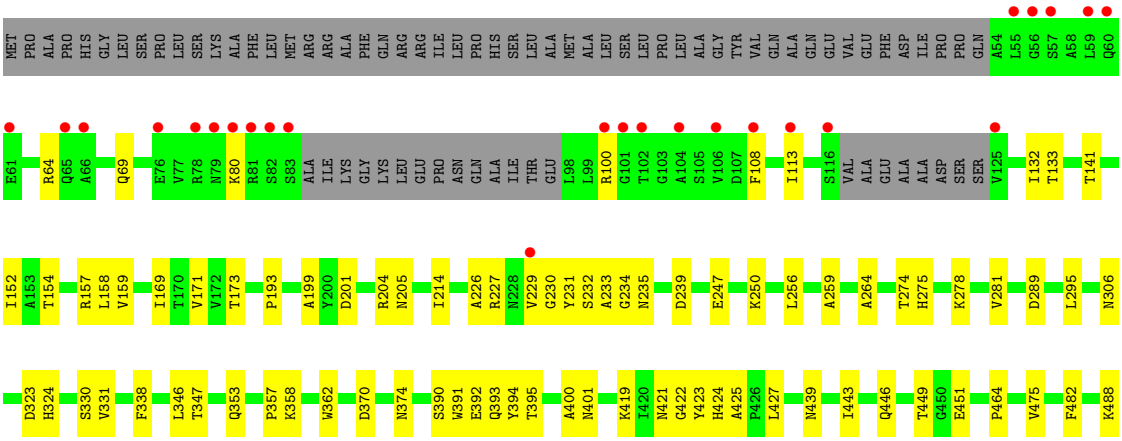
• Molecule 1: Ferripyoverdine receptor

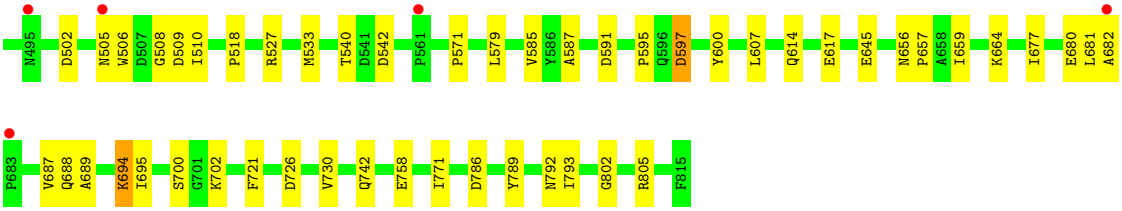
Chain A:



• Molecule 1: Ferripyoverdine receptor

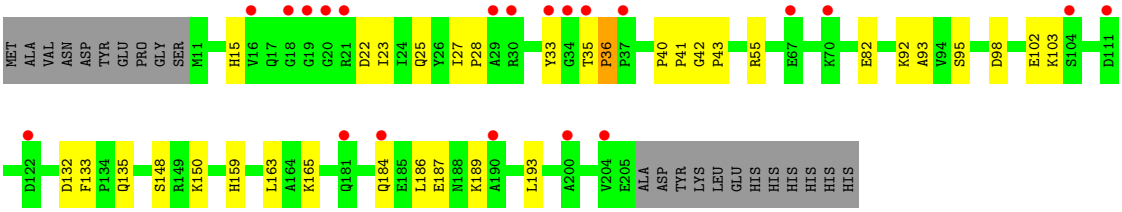
Chain B:





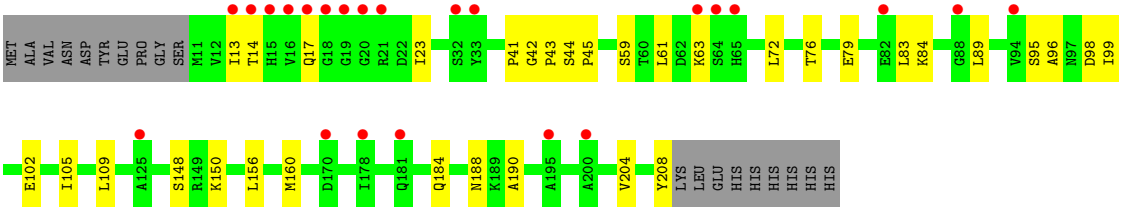
- Molecule 2: Pyocin-S2

Chain C:



- Molecule 2: Pyocin-S2

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	115.52Å 209.39Å 215.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.97 – 2.80 47.97 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.97-2.80) 98.6 (47.97-2.80)	Depositor EDS
R_{merge}	0.35	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.61 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.212 , 0.262 0.227 , 0.274	Depositor DCC
R_{free} test set	3094 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å^2)	36.4	Xtriage
Anisotropy	0.603	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 51.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.018 for -h,-l,-k	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	15094	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	0.48	0/6257	0.72	0/8502
1	B	0.47	0/6020	0.71	0/8175
2	C	0.45	0/1545	0.65	0/2079
2	D	0.44	0/1571	0.65	0/2115
All	All	0.47	0/15393	0.70	0/20871

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	6105	0	5780	92	0
1	B	5874	0	5555	110	0
2	C	1520	0	1525	35	0
2	D	1545	0	1542	32	0
3	A	25	0	0	0	0
3	B	25	0	0	0	0
All	All	15094	0	14402	250	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 8.

All (250) 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:306:ASN:HD21	1:B:306:ASN:ND2	1.00	1.48
1:A:306:ASN:ND2	1:B:306:ASN:HD21	0.98	1.45
1:B:227:ARG:NH2	1:B:234:GLY:H	1.22	1.33
2:D:95:SER:OG	2:D:98:ASP:OD2	1.53	1.26
1:A:742:GLN:NE2	1:A:793:ILE:O	1.65	1.26
2:D:89:LEU:HD12	2:D:105:ILE:HD11	1.21	1.17
1:A:227:ARG:NH1	1:A:234:GLY:H	1.43	1.16
1:A:227:ARG:HH12	1:A:234:GLY:N	1.46	1.11
1:B:227:ARG:NH2	1:B:234:GLY:N	1.96	1.11
1:B:227:ARG:HH12	1:B:235:ASN:HB2	1.18	1.04
1:A:462:THR:HG22	1:A:475:VAL:HG22	1.40	1.00
2:D:72:LEU:O	2:D:76:THR:HG22	1.61	0.99
1:A:194:GLY:HA2	1:A:712:GLN:HE21	1.26	0.99
1:B:230:GLY:HA3	2:D:43:PRO:CD	1.92	0.99
1:A:76:GLU:O	1:A:80:LYS:NZ	1.98	0.96
1:B:227:ARG:HH21	1:B:234:GLY:H	1.02	0.96
1:B:230:GLY:HA3	2:D:43:PRO:HD3	1.49	0.95
1:A:774:LYS:H	1:A:774:LYS:HD2	1.35	0.92
1:A:227:ARG:HH12	1:A:234:GLY:H	0.92	0.89
1:A:194:GLY:HA2	1:A:712:GLN:NE2	1.87	0.89
1:B:227:ARG:NH1	1:B:235:ASN:HB2	1.86	0.89
1:B:227:ARG:HH22	1:B:235:ASN:H	1.24	0.85
1:A:229:VAL:HB	2:C:41:PRO:HB2	1.58	0.84
1:B:227:ARG:HH22	1:B:234:GLY:N	1.67	0.83
2:D:89:LEU:HD12	2:D:105:ILE:CD1	2.06	0.83
1:A:694:LYS:HD3	1:A:694:LYS:O	1.80	0.81
1:B:193:PRO:O	1:B:250:LYS:NZ	2.14	0.81
2:C:133:PHE:CE2	2:C:163:LEU:CD2	2.65	0.80
1:B:227:ARG:HH22	1:B:234:GLY:H	1.16	0.79
1:B:742:GLN:OE1	1:B:793:ILE:O	2.00	0.79
1:A:597:ASP:HB2	1:A:600:TYR:HD1	1.48	0.79
1:B:230:GLY:HA3	2:D:43:PRO:HD2	1.66	0.78
2:D:99:ILE:HD11	2:D:190:ALA:HA	1.66	0.77
2:D:89:LEU:CD1	2:D:105:ILE:HD11	2.08	0.77
1:A:227:ARG:NH1	1:A:233:ALA:H	1.84	0.75
1:A:50:ILE:HD12	1:A:85:ILE:HD11	1.69	0.75
1:A:774:LYS:HD2	1:A:774:LYS:N	2.02	0.74
2:D:204:VAL:O	2:D:208:TYR:HD2	1.70	0.74
1:B:227:ARG:HH12	1:B:235:ASN:CB	1.99	0.73
1:B:227:ARG:HH22	1:B:235:ASN:N	1.86	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:680:GLU:HG3	1:A:682:ALA:O	1.89	0.73
2:C:133:PHE:CE2	2:C:163:LEU:HD23	2.23	0.73
1:B:540:THR:HG22	1:B:542:ASP:H	1.53	0.73
1:A:194:GLY:CA	1:A:712:GLN:NE2	2.52	0.72
1:A:600:TYR:O	1:A:607:LEU:HD12	1.90	0.71
2:C:27:ILE:HD12	2:C:27:ILE:N	2.05	0.71
1:B:645:GLU:OE1	1:B:664:LYS:HD3	1.90	0.71
1:B:227:ARG:HE	1:B:232:SER:HA	1.54	0.71
1:B:233:ALA:HB1	1:B:421:ASN:HB3	1.74	0.69
1:B:680:GLU:HG3	1:B:682:ALA:O	1.92	0.68
1:A:230:GLY:HA3	2:C:42:GLY:HA3	1.75	0.68
1:A:597:ASP:HB2	1:A:600:TYR:CD1	2.29	0.68
2:C:132:ASP:OD1	2:C:135:GLN:HB2	1.94	0.67
1:B:199:ALA:HB2	1:B:205:ASN:OD1	1.93	0.67
1:A:66:ALA:HB3	1:A:68:ILE:HG22	1.76	0.66
1:B:227:ARG:NH2	1:B:235:ASN:H	1.94	0.65
1:A:306:ASN:ND2	1:B:306:ASN:ND2	1.81	0.64
2:D:79:GLU:O	2:D:83:LEU:HG	1.97	0.64
1:A:152:ILE:HD13	1:A:169:ILE:HG12	1.80	0.64
1:B:152:ILE:HD13	1:B:169:ILE:HG12	1.80	0.64
1:B:227:ARG:HH21	1:B:233:ALA:N	1.95	0.63
1:B:597:ASP:HB2	1:B:600:TYR:HD1	1.64	0.63
1:B:681:LEU:HD11	1:B:687:VAL:HG21	1.80	0.62
1:A:52:PRO:HG3	1:A:84:ALA:HB2	1.80	0.62
1:A:680:GLU:CG	1:A:682:ALA:O	2.48	0.61
1:B:700:SER:OG	1:B:702:LYS:HG2	2.00	0.61
2:C:184:GLN:O	2:C:187:GLU:HG2	2.00	0.61
1:B:347:THR:HB	1:B:401:ASN:HB2	1.82	0.60
1:B:201:ASP:OD1	1:B:204:ARG:HB2	2.00	0.60
1:B:591:ASP:OD1	1:B:614:GLN:HG3	2.02	0.60
2:C:15:HIS:O	2:C:23:ILE:HG13	2.02	0.60
1:B:152:ILE:HG22	1:B:247:GLU:OE1	2.01	0.60
1:B:199:ALA:O	1:B:792:ASN:HB2	2.01	0.60
2:C:15:HIS:HE1	2:C:25:GLN:OE1	1.83	0.59
1:B:502:ASP:OD2	1:B:505:ASN:HB3	2.02	0.59
1:A:227:ARG:NH1	1:A:233:ALA:N	2.51	0.59
1:B:694:LYS:HG2	1:B:695:ILE:N	2.18	0.59
1:B:227:ARG:HH21	1:B:234:GLY:N	1.81	0.58
1:B:132:ILE:HD11	1:B:464:PRO:HG3	1.86	0.58
1:A:694:LYS:HD3	1:A:694:LYS:C	2.23	0.57
1:A:410:TRP:CD2	1:A:463:GLY:HA3	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:TYR:CE1	2:D:43:PRO:HG3	2.40	0.57
1:B:656:ASN:OD1	1:B:657:PRO:HD2	2.04	0.56
1:B:239:ASP:HB3	1:B:331:VAL:HG21	1.86	0.56
1:B:680:GLU:CG	1:B:682:ALA:O	2.53	0.56
2:C:95:SER:HB3	2:C:98:ASP:HB2	1.88	0.56
2:D:84:LYS:HG2	2:D:105:ILE:HD13	1.88	0.56
1:A:410:TRP:CE2	1:A:463:GLY:HA3	2.40	0.56
1:A:68:ILE:HD11	1:A:113:ILE:HG23	1.89	0.55
1:B:571:PRO:HD2	1:B:591:ASP:HB3	1.88	0.55
1:B:232:SER:OG	1:B:393:GLN:OE1	2.25	0.54
2:C:103:LYS:HA	2:C:186:LEU:HD13	1.89	0.54
2:C:159:HIS:O	2:C:163:LEU:HG	2.07	0.54
2:D:59:SER:O	2:D:63:LYS:HG2	2.07	0.54
1:B:201:ASP:HB3	1:B:362:TRP:O	2.08	0.54
1:A:194:GLY:C	1:A:712:GLN:HE22	2.10	0.54
1:A:133:THR:HG21	1:A:159:VAL:HG21	1.88	0.54
1:A:239:ASP:HB3	1:A:331:VAL:HG11	1.90	0.53
1:B:338:PHE:CZ	1:B:346:LEU:HD23	2.44	0.53
1:B:439:ASN:ND2	1:B:502:ASP:OD1	2.40	0.53
2:D:84:LYS:HD3	2:D:102:GLU:HG3	1.89	0.53
1:B:687:VAL:HG12	1:B:688:GLN:N	2.24	0.53
1:A:302:THR:OG1	1:A:306:ASN:OD1	2.10	0.53
2:C:133:PHE:CE2	2:C:163:LEU:HD21	2.44	0.53
1:A:745:TYR:CD1	2:C:33:TYR:HB2	2.44	0.53
1:A:774:LYS:H	1:A:774:LYS:CD	2.13	0.53
1:B:786:ASP:HB2	1:B:805:ARG:HB2	1.91	0.53
1:B:227:ARG:NH2	1:B:233:ALA:H	2.06	0.52
2:D:83:LEU:HD11	2:D:109:LEU:HD13	1.90	0.52
1:B:227:ARG:NH2	1:B:233:ALA:N	2.57	0.52
1:B:323:ASP:OD1	1:B:324:HIS:N	2.42	0.52
1:A:158:LEU:HD21	1:A:475:VAL:HG23	1.92	0.52
1:B:597:ASP:HB2	1:B:600:TYR:CD1	2.45	0.52
1:A:391:TRP:CZ3	1:A:427:LEU:HD11	2.45	0.51
1:A:613:GLY:HA2	1:A:639:HIS:O	2.10	0.51
1:A:242:ILE:HG22	1:A:312:VAL:HG22	1.92	0.51
2:C:133:PHE:CD2	2:C:163:LEU:HD23	2.45	0.51
2:C:189:LYS:O	2:C:193:LEU:HB2	2.11	0.51
1:A:230:GLY:HA3	2:C:43:PRO:HD2	1.93	0.51
1:A:475:VAL:O	1:A:533:MET:HA	2.10	0.51
1:B:281:VAL:HG22	1:B:295:LEU:HD13	1.94	0.50
1:B:424:HIS:HA	1:B:449:THR:HG22	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:424:HIS:HA	1:A:449:THR:HG22	1.93	0.50
1:A:745:TYR:CG	2:C:33:TYR:HB2	2.47	0.50
1:B:587:ALA:HA	1:B:617:GLU:O	2.12	0.50
1:A:587:ALA:HA	1:A:617:GLU:O	2.12	0.50
1:B:475:VAL:O	1:B:533:MET:HA	2.13	0.49
1:B:157:ARG:HE	1:B:256:LEU:HD22	1.77	0.49
2:D:188:ASN:N	2:D:188:ASN:HD22	2.10	0.49
1:A:194:GLY:CA	1:A:712:GLN:HE21	2.08	0.48
1:A:48:PHE:O	1:A:86:LYS:HA	2.13	0.48
1:A:158:LEU:CD2	1:A:475:VAL:HG23	2.44	0.48
1:B:687:VAL:CG1	1:B:688:GLN:N	2.76	0.48
2:C:132:ASP:OD1	2:C:135:GLN:OE1	2.31	0.48
1:A:259:ALA:HB1	1:A:595:PRO:HD3	1.96	0.48
1:B:391:TRP:CZ3	1:B:427:LEU:HD11	2.49	0.48
1:B:677:ILE:CG2	1:B:689:ALA:HB3	2.44	0.48
2:C:22:ASP:OD2	2:C:165:LYS:HG2	2.13	0.48
2:D:156:LEU:O	2:D:160:MET:HG2	2.14	0.48
1:B:230:GLY:CA	2:D:43:PRO:HD3	2.33	0.47
1:A:392:GLU:O	1:A:423:TYR:HA	2.14	0.47
1:B:677:ILE:HG23	1:B:689:ALA:HB3	1.96	0.47
2:C:15:HIS:CE1	2:C:25:GLN:OE1	2.66	0.47
1:B:721:PHE:O	1:B:726:ASP:HA	2.14	0.47
1:A:659:ILE:HG23	1:A:659:ILE:O	2.13	0.47
1:B:227:ARG:NE	1:B:233:ALA:H	2.13	0.47
1:A:492:ASN:HD22	2:C:55:ARG:HH22	1.61	0.47
1:A:390:SER:O	1:A:425:ALA:HA	2.15	0.47
1:A:571:PRO:HD2	1:A:591:ASP:HB3	1.95	0.47
1:B:579:LEU:HD11	1:B:585:VAL:HG13	1.96	0.47
1:A:152:ILE:HB	1:A:154:THR:HG22	1.96	0.47
1:A:568:ARG:NH2	1:A:610:PRO:HD2	2.30	0.46
1:A:568:ARG:CZ	1:A:610:PRO:HG2	2.45	0.46
1:A:192:THR:HG21	1:A:248:VAL:HG11	1.96	0.46
1:B:227:ARG:CZ	1:B:233:ALA:H	2.29	0.46
2:C:33:TYR:CG	2:C:33:TYR:O	2.69	0.46
2:D:96:ALA:O	2:D:99:ILE:HG12	2.14	0.46
1:A:501:ASP:OD1	1:A:501:ASP:N	2.49	0.46
2:D:17:GLN:CD	2:D:17:GLN:H	2.18	0.46
1:B:171:VAL:HG22	1:B:247:GLU:HG2	1.98	0.46
2:C:27:ILE:HD12	2:C:27:ILE:H	1.78	0.46
1:A:227:ARG:NH1	1:A:234:GLY:N	2.19	0.46
1:B:392:GLU:O	1:B:423:TYR:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:506:TRP:CZ2	1:B:508:GLY:HA2	2.50	0.46
1:A:154:THR:HB	1:A:247:GLU:OE1	2.15	0.46
1:A:229:VAL:CB	2:C:41:PRO:HB2	2.38	0.46
1:A:347:THR:O	1:A:400:ALA:HA	2.16	0.46
1:A:281:VAL:HG22	1:A:295:LEU:CD1	2.46	0.45
1:A:772:THR:HG22	1:A:774:LYS:HD2	1.98	0.45
1:A:263:GLY:O	1:A:264:ALA:HB2	2.16	0.45
2:C:35:THR:HA	2:C:36:PRO:HD3	1.82	0.45
1:B:152:ILE:HB	1:B:154:THR:HG22	1.98	0.45
1:B:230:GLY:HA3	2:D:42:GLY:HA3	1.99	0.45
1:B:789:TYR:CZ	1:B:802:GLY:HA3	2.51	0.45
2:D:95:SER:OG	2:D:98:ASP:CG	2.44	0.45
1:A:157:ARG:HH11	1:A:256:LEU:HD22	1.82	0.45
1:A:789:TYR:CZ	1:A:802:GLY:HA3	2.52	0.44
1:B:259:ALA:HB1	1:B:595:PRO:HD3	1.99	0.44
1:B:362:TRP:CH2	1:B:446:GLN:HB2	2.52	0.44
1:B:390:SER:O	1:B:425:ALA:HA	2.17	0.44
1:B:659:ILE:HG23	1:B:659:ILE:O	2.17	0.44
1:B:338:PHE:CE2	1:B:346:LEU:HD23	2.52	0.44
2:D:184:GLN:O	2:D:188:ASN:ND2	2.50	0.44
1:A:46:VAL:HG12	1:A:47:GLU:N	2.33	0.44
1:A:408:ASN:OD1	1:A:410:TRP:HB2	2.18	0.44
2:D:14:THR:HA	2:D:23:ILE:O	2.18	0.44
1:A:579:LEU:HD11	1:A:585:VAL:HG13	1.99	0.44
1:B:230:GLY:CA	2:D:43:PRO:HD2	2.42	0.44
2:C:33:TYR:O	2:C:33:TYR:CD2	2.70	0.44
1:B:357:PRO:O	1:B:358:LYS:HD2	2.17	0.44
1:B:229:VAL:HB	2:D:41:PRO:HB2	1.99	0.44
1:B:227:ARG:HH12	1:B:235:ASN:H	1.64	0.44
2:C:40:PRO:HA	2:C:41:PRO:HD3	1.90	0.44
2:D:95:SER:CB	2:D:98:ASP:OD2	2.60	0.43
1:A:482:PHE:CE2	1:A:527:ARG:HG2	2.54	0.43
1:B:158:LEU:HD21	1:B:475:VAL:HG23	2.01	0.43
1:B:656:ASN:HA	1:B:657:PRO:HD3	1.74	0.43
1:A:357:PRO:O	1:A:358:LYS:HD2	2.19	0.43
1:A:506:TRP:CZ2	1:A:508:GLY:HA2	2.53	0.43
1:B:64:ARG:O	1:B:275:HIS:CE1	2.71	0.43
1:B:158:LEU:CD2	1:B:475:VAL:HG23	2.49	0.43
1:B:108:PHE:HA	1:B:113:ILE:HG22	2.00	0.43
1:B:347:THR:O	1:B:400:ALA:HA	2.18	0.43
1:B:330:SER:O	1:B:353:GLN:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:ASP:OD1	1:B:374:ASN:N	2.52	0.43
1:B:394:TYR:CZ	1:B:422:GLY:HA3	2.53	0.43
1:B:443:ILE:HG13	1:B:510:ILE:HD13	1.99	0.43
1:B:227:ARG:HH22	1:B:234:GLY:CA	2.27	0.43
1:B:600:TYR:O	1:B:607:LEU:HD12	2.19	0.43
1:A:141:THR:HB	1:A:173:THR:HB	2.01	0.42
1:A:194:GLY:C	1:A:712:GLN:NE2	2.72	0.42
1:B:681:LEU:CD1	1:B:687:VAL:HG21	2.48	0.42
1:A:443:ILE:HG13	1:A:510:ILE:HD13	2.01	0.42
1:B:214:ILE:HG12	1:B:264:ALA:HB3	1.99	0.42
1:B:656:ASN:O	1:B:659:ILE:O	2.38	0.42
1:B:141:THR:HB	1:B:173:THR:HB	2.00	0.42
1:A:45:GLU:HG2	1:A:88:LYS:HG2	2.01	0.42
2:C:35:THR:O	2:C:35:THR:HG23	2.20	0.42
1:B:69:GLN:CG	1:B:133:THR:HG23	2.50	0.42
1:A:108:PHE:HA	1:A:113:ILE:HG22	2.02	0.41
1:A:650:TYR:OH	1:A:659:ILE:HG23	2.20	0.41
2:D:89:LEU:CD1	2:D:105:ILE:CD1	2.83	0.41
1:A:682:ALA:HB1	1:A:683:PRO:HD2	2.01	0.41
2:C:82:GLU:HA	2:C:82:GLU:OE1	2.20	0.41
1:A:52:PRO:CG	1:A:84:ALA:HB2	2.50	0.41
2:D:44:SER:HA	2:D:45:PRO:HD3	1.93	0.41
1:A:231:TYR:CE2	2:C:43:PRO:HD3	2.56	0.41
1:A:774:LYS:N	1:A:774:LYS:CD	2.73	0.41
1:B:488:LYS:HD3	1:B:518:PRO:HG2	2.02	0.41
1:B:742:GLN:CD	1:B:793:ILE:O	2.59	0.41
2:D:148:SER:HB2	2:D:150:LYS:HG2	2.02	0.41
1:A:233:ALA:HB1	1:A:421:ASN:HB3	2.03	0.41
1:B:69:GLN:NE2	1:B:133:THR:HG21	2.35	0.41
1:B:482:PHE:CE2	1:B:527:ARG:HG2	2.56	0.41
1:A:716:TYR:CD1	1:A:733:GLY:HA3	2.56	0.41
1:A:772:THR:HG22	1:A:773:ASP:N	2.36	0.40
1:A:281:VAL:HG23	1:A:815:PHE:CZ	2.56	0.40
1:B:419:LYS:HD3	1:B:421:ASN:OD1	2.21	0.40
2:C:92:LYS:H	2:C:92:LYS:HG2	1.65	0.40
2:C:93:ALA:HB3	2:C:102:GLU:OE1	2.21	0.40
1:A:288:TRP:O	1:A:319:HIS:HB2	2.21	0.40
2:D:13:ILE:HG22	2:D:14:THR:N	2.36	0.40
1:A:496:TYR:CG	1:A:513:PRO:HB3	2.56	0.40
1:B:274:THR:HG21	1:B:278:LYS:HE2	2.03	0.40
1:B:395:THR:HG22	1:B:421:ASN:ND2	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:THR:HG22	1:B:542:ASP:N	2.30	0.40
1:B:133:THR:HG21	1:B:159:VAL:HG11	2.04	0.40
1:B:394:TYR:OH	1:B:451:GLU:HG3	2.21	0.40
2:C:27:ILE:N	2:C:27:ILE:CD1	2.76	0.40
2:C:148:SER:HB2	2:C:150:LYS:HG2	2.02	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	769/815 (94%)	741 (96%)	28 (4%)	0	100	100
1	B	734/815 (90%)	708 (96%)	25 (3%)	1 (0%)	51	81
2	C	193/217 (89%)	187 (97%)	4 (2%)	2 (1%)	15	44
2	D	196/217 (90%)	190 (97%)	6 (3%)	0	100	100
All	All	1892/2064 (92%)	1826 (96%)	63 (3%)	3 (0%)	47	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	28	PRO
1	B	226	ALA
2	C	36	PRO

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	649/684 (95%)	639 (98%)	10 (2%)	65	89
1	B	624/684 (91%)	615 (99%)	9 (1%)	67	90
2	C	163/182 (90%)	163 (100%)	0	100	100
2	D	165/182 (91%)	164 (99%)	1 (1%)	86	96
All	All	1601/1732 (92%)	1581 (99%)	20 (1%)	71	92

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	LEU
1	A	124	SER
1	A	133	THR
1	A	201	ASP
1	A	289	ASP
1	A	312	VAL
1	A	330	SER
1	A	421	ASN
1	A	694	LYS
1	A	735	ARG
1	B	80	LYS
1	B	100	ARG
1	B	289	ASP
1	B	509	ASP
1	B	597	ASP
1	B	694	LYS
1	B	730	VAL
1	B	758	GLU
1	B	771	ILE
2	D	61	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	712	GLN
1	B	69	GLN
1	B	306	ASN
1	B	712	GLN

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Mol	Chain	Res	Type
2	C	15	HIS
2	D	175	GLN
2	D	184	GLN
2	D	188	ASN

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)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	SO4	A	904	-	4,4,4	0.34	0	6,6,6	0.22	0
3	SO4	B	903	-	4,4,4	0.14	0	6,6,6	0.26	0
3	SO4	A	902	-	4,4,4	0.21	0	6,6,6	0.32	0
3	SO4	A	901	-	4,4,4	0.17	0	6,6,6	0.30	0
3	SO4	B	905	-	4,4,4	0.16	0	6,6,6	0.19	0
3	SO4	A	905	-	4,4,4	0.19	0	6,6,6	0.15	0
3	SO4	B	902	-	4,4,4	0.25	0	6,6,6	0.47	0
3	SO4	B	904	-	4,4,4	0.15	0	6,6,6	0.10	0
3	SO4	B	901	-	4,4,4	0.22	0	6,6,6	0.26	0
3	SO4	A	903	-	4,4,4	0.33	0	6,6,6	0.22	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	771/815 (94%)	-0.05	7 (0%)	84	80	8, 29, 51, 76	0
1	B	740/815 (90%)	0.12	30 (4%)	37	27	14, 31, 65, 93	0
2	C	195/217 (89%)	0.75	21 (10%)	5	3	27, 52, 85, 105	0
2	D	198/217 (91%)	0.76	23 (11%)	4	2	28, 55, 89, 113	0
All	All	1904/2064 (92%)	0.18	81 (4%)	35	25	8, 33, 71, 113	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	19	GLY	6.4
2	D	18	GLY	6.4
2	C	34	GLY	6.3
2	D	17	GLN	5.7
2	C	18	GLY	5.6
1	B	79	ASN	5.3
2	C	29	ALA	5.2
2	D	21	ARG	4.9
2	D	20	GLY	4.6
2	C	20	GLY	4.3
1	B	113	ILE	4.0
1	B	60	GLN	3.8
1	B	108	PHE	3.6
1	B	116	SER	3.6
1	A	407	ALA	3.6
2	D	14	THR	3.5
2	D	15	HIS	3.5
2	C	30	ARG	3.4
2	C	204	VAL	3.3
1	B	78	ARG	3.3
1	B	83	SER	3.3

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Mol	Chain	Res	Type	RSRZ
2	D	63	LYS	3.2
1	B	229	VAL	3.2
2	C	35	THR	3.1
2	C	19	GLY	2.8
1	B	55	LEU	2.8
1	B	101	GLY	2.8
2	C	184	GLN	2.7
1	B	100	ARG	2.7
1	B	81	ARG	2.7
2	C	111	ASP	2.7
1	A	699	ASP	2.7
2	D	170	ASP	2.6
1	A	649	LEU	2.6
1	B	102	THR	2.5
1	B	125	VAL	2.5
1	B	82	SER	2.5
1	B	104	ALA	2.5
2	D	94	VAL	2.5
1	A	495	ASN	2.5
1	B	106	VAL	2.5
2	C	122	ASP	2.5
2	D	64	SER	2.4
2	D	178	ILE	2.4
2	C	70	LYS	2.4
2	D	82	GLU	2.4
1	A	724	ALA	2.4
1	B	65	GLN	2.3
2	D	33	TYR	2.3
1	B	59	LEU	2.3
1	B	505	ASN	2.3
2	C	33	TYR	2.3
1	B	56	GLY	2.3
2	C	181	GLN	2.3
2	D	16	VAL	2.2
2	D	125	ALA	2.2
1	B	61	GLU	2.2
2	C	104	SER	2.2
2	D	13	ILE	2.2
2	D	32	SER	2.2
2	D	195	ALA	2.2
2	D	88	GLY	2.2
1	B	682	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	57	SER	2.2
1	B	561	PRO	2.2
2	C	16	VAL	2.2
2	D	200	ALA	2.1
1	B	80	LYS	2.1
2	C	67	GLU	2.1
2	D	181	GLN	2.1
1	B	66	ALA	2.1
1	A	265	THR	2.1
2	C	21	ARG	2.1
1	B	76	GLU	2.1
2	D	65	HIS	2.1
1	A	700	SER	2.0
1	B	683	PRO	2.0
2	C	190	ALA	2.0
2	C	200	ALA	2.0
2	C	37	PRO	2.0
1	B	495	ASN	2.0

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	SO4	B	905	5/5	0.88	0.24	106,106,107,107	0
3	SO4	A	904	5/5	0.92	0.21	67,68,69,70	0
3	SO4	A	905	5/5	0.93	0.24	82,84,85,85	0
3	SO4	B	904	5/5	0.94	0.20	72,73,74,75	0
3	SO4	B	903	5/5	0.94	0.24	52,53,54,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SO4	A	902	5/5	0.95	0.20	53,58,61,61	0
3	SO4	B	902	5/5	0.96	0.17	54,55,55,55	0
3	SO4	B	901	5/5	0.96	0.17	54,55,56,58	0
3	SO4	A	903	5/5	0.97	0.17	52,55,56,57	0
3	SO4	A	901	5/5	0.98	0.14	53,56,58,59	0

6.5 Other polymers (i)

There are no such residues in this entry.