Full wwPDB X-ray Structure Validation Report (i)

Sep 13, 2023 – 11:05 AM EDT

PDB ID	:	4PSM
Title	:	Crystal structure of pfuThermo-DBP-RP1 (crystal form II)
Authors	:	Gahlei, H.; von Moeller, H.; Eppers, D.; Loll, B.; Wahl, M.C.
Deposited on	:	2014-03-07
Resolution	:	2.43 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:

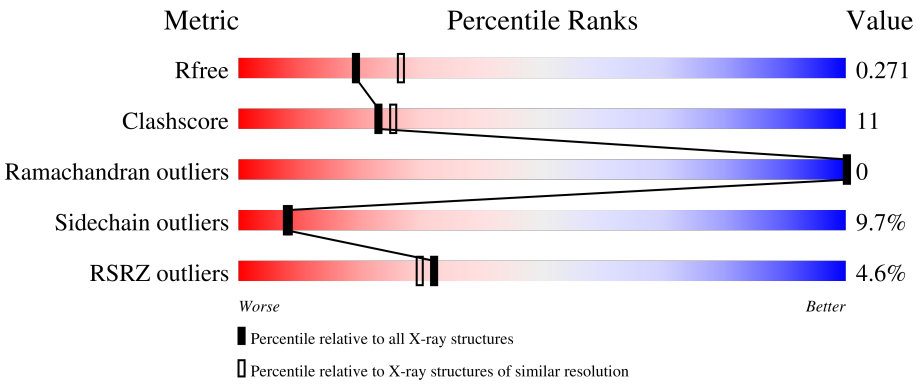
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.43 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	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

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	149	% 66% 32% ••
1	B	149	5% 62% 32% • •
1	C	149	9% 62% 32% • •
1	D	149	3% 65% 32% ••

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4883 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.

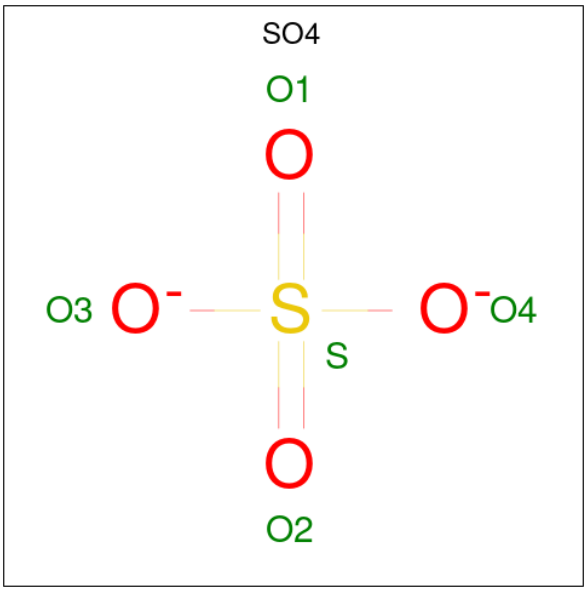
- Molecule 1 is a protein called ssDNA binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	147	1204	766	198	234	6	6	1	0
1	B	146	1196	761	197	231	7	0	1	0
1	C	146	1191	757	197	231	6	0	0	0
1	D	147	1204	766	198	234	6	6	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q8U208
A	1	ALA	-	expression tag	UNP Q8U208
B	0	GLY	-	expression tag	UNP Q8U208
B	1	ALA	-	expression tag	UNP Q8U208
C	0	GLY	-	expression tag	UNP Q8U208
C	1	ALA	-	expression tag	UNP Q8U208
D	0	GLY	-	expression tag	UNP Q8U208
D	1	ALA	-	expression tag	UNP Q8U208

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	22	22	22	0	0
3	B	14	14	14	0	0
3	C	10	10	10	0	0
3	D	22	22	22	0	0

3 Residue-property plots (i)

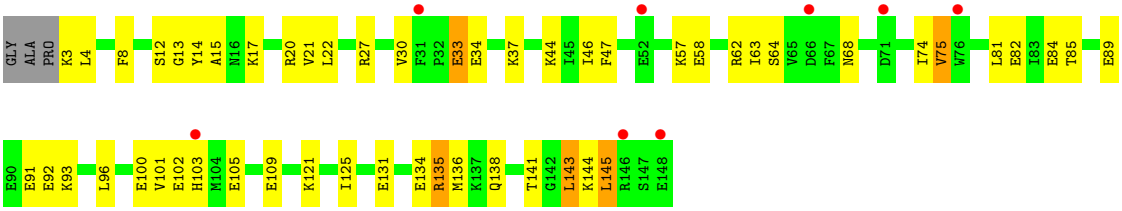
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: ssDNA binding protein

Chain A:

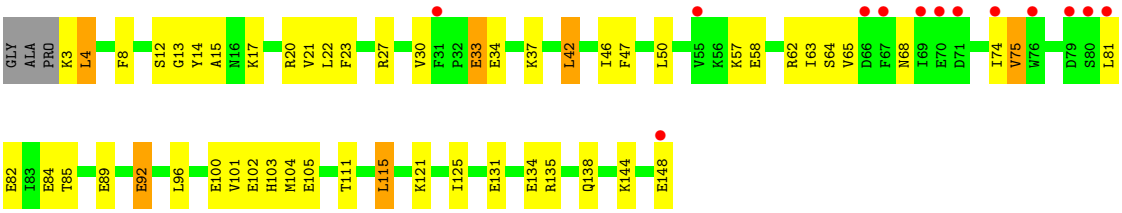
• Molecule 1: ssDNA binding protein

Chain B:



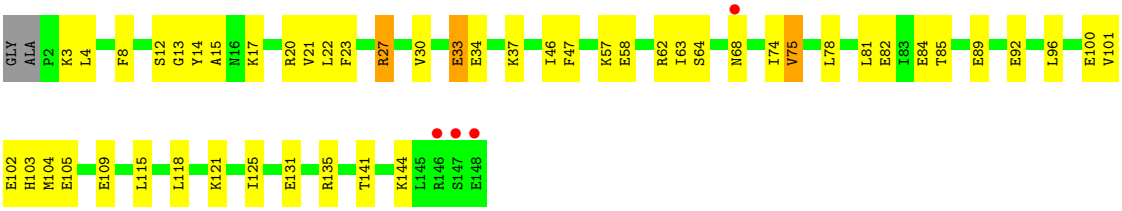
• Molecule 1: ssDNA binding protein

Chain C:



• Molecule 1: ssDNA binding protein

Chain D:


GLY
ALA
P2
K3
L4
F8
S12
G13
Y14
A15
N16
K17
R20
V21
L22
F23
R27
V30
E33
E34
K37
I46
F47
K57
E58
R62
I63
S64
N68
I74
V75
L78
L81
E82
I83
E84
T85
E89
E92
L96
E100
V101
E102
H103
M104
E105
E109
L115
L118
K121
I125
E131
R135
T141
K144
L145
R146
S147
E148

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	61.25Å 199.30Å 114.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.41 – 2.43 33.41 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.0 (33.41-2.43) 99.4 (33.41-2.43)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.06 (at 2.42Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.223 , 0.276 0.221 , 0.271	Depositor DCC
R_{free} test set	1341 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	37.5	Xtriage
Anisotropy	0.851	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 64.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4883	wwPDB-VP
Average B, all atoms (Å^2)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.5043e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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	0.33	0/1214	0.78	6/1609 (0.4%)
1	B	0.32	0/1205	0.76	5/1596 (0.3%)
1	C	0.32	0/1197	0.79	7/1586 (0.4%)
1	D	0.33	0/1214	0.75	6/1609 (0.4%)
All	All	0.32	0/4830	0.77	24/6400 (0.4%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	135	ARG	NE-CZ-NH2	-12.42	114.09	120.30
1	B	135	ARG	NE-CZ-NH2	-11.68	114.46	120.30
1	A	135	ARG	NE-CZ-NH1	11.62	126.11	120.30
1	D	135	ARG	NE-CZ-NH1	-11.55	114.53	120.30
1	C	135	ARG	NE-CZ-NH1	-11.43	114.59	120.30
1	B	135	ARG	NE-CZ-NH1	11.25	125.93	120.30
1	C	135	ARG	NE-CZ-NH2	11.09	125.85	120.30
1	D	135	ARG	NE-CZ-NH2	11.07	125.83	120.30
1	C	27	ARG	NE-CZ-NH2	-10.54	115.03	120.30
1	D	27	ARG	NE-CZ-NH2	-10.44	115.08	120.30
1	D	27	ARG	NE-CZ-NH1	10.34	125.47	120.30
1	C	27	ARG	NE-CZ-NH1	10.27	125.44	120.30
1	A	27	ARG	NE-CZ-NH1	-10.18	115.21	120.30
1	B	27	ARG	NE-CZ-NH1	-10.15	115.23	120.30
1	A	27	ARG	NE-CZ-NH2	9.73	125.17	120.30
1	B	27	ARG	NE-CZ-NH2	9.62	125.11	120.30
1	A	135	ARG	CD-NE-CZ	5.76	131.66	123.60
1	D	135	ARG	CD-NE-CZ	5.58	131.41	123.60
1	C	135	ARG	CD-NE-CZ	5.55	131.37	123.60
1	B	135	ARG	CD-NE-CZ	5.50	131.30	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	4	LEU	CA-CB-CG	5.46	127.87	115.30
1	C	4	LEU	CA-CB-CG	5.38	127.67	115.30
1	D	27	ARG	CD-NE-CZ	5.16	130.82	123.60
1	C	27	ARG	CD-NE-CZ	5.01	130.62	123.60

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	1204	0	1242	33	0
1	B	1196	0	1237	41	0
1	C	1191	0	1228	35	0
1	D	1204	0	1242	32	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	22	0	0	0	0
3	B	14	0	0	0	0
3	C	10	0	0	0	1
3	D	22	0	0	0	0
All	All	4883	0	4949	111	1

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 11.

All (111) 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:125:ILE:HD12	1:D:125:ILE:HD12	1.55	0.87
1:C:46:ILE:HD13	1:C:63:ILE:HD13	1.65	0.77
1:B:125:ILE:HD12	1:C:125:ILE:HD12	1.65	0.76
1:D:46:ILE:HD13	1:D:63:ILE:HD13	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:ILE:HD13	1:A:63:ILE:HD13	1.66	0.75
1:B:46:ILE:HD13	1:B:63:ILE:HD13	1.68	0.75
1:A:125:ILE:HD11	1:D:121:LYS:HB3	1.69	0.74
1:A:121:LYS:HB3	1:D:125:ILE:HD11	1.70	0.73
1:B:125:ILE:HD11	1:C:121:LYS:HB3	1.71	0.70
1:B:121:LYS:HB3	1:C:125:ILE:HD11	1.76	0.66
1:C:14:TYR:OH	1:C:47:PHE:HB2	1.98	0.64
1:B:12:SER:HB3	1:B:57:LYS:HB2	1.80	0.63
1:A:14:TYR:OH	1:A:47:PHE:HB2	1.99	0.63
1:B:109:GLU:HG2	1:D:141:THR:HB	1.82	0.62
1:D:14:TYR:OH	1:D:47:PHE:HB2	2.00	0.62
1:A:12:SER:HB3	1:A:57:LYS:HB2	1.81	0.61
1:B:14:TYR:OH	1:B:47:PHE:HB2	2.00	0.61
1:D:12:SER:HB3	1:D:57:LYS:HB2	1.83	0.61
1:C:12:SER:HB3	1:C:57:LYS:HB2	1.81	0.60
1:B:13:GLY:HA3	1:B:102:GLU:OE2	2.00	0.60
1:C:13:GLY:HA3	1:C:102:GLU:OE2	2.03	0.59
1:D:13:GLY:HA3	1:D:102:GLU:OE2	2.02	0.59
1:C:115:LEU:HD11	1:D:115:LEU:HD21	1.83	0.59
1:A:13:GLY:HA3	1:A:102:GLU:OE2	2.05	0.56
1:A:104:MSE:CE	1:B:20:ARG:HH21	2.19	0.55
1:B:144:LYS:NZ	1:B:144:LYS:HB2	2.22	0.55
1:A:57:LYS:HD2	2:A:201:SO4:O4	2.06	0.55
1:D:144:LYS:HB2	1:D:144:LYS:NZ	2.23	0.54
1:A:23:PHE:CG	1:C:138:GLN:HB3	2.43	0.54
1:C:144:LYS:HB2	1:C:144:LYS:NZ	2.24	0.53
1:A:93:LYS:HE3	1:B:91:GLU:OE2	2.08	0.53
1:A:144:LYS:HB2	1:A:144:LYS:NZ	2.24	0.52
1:C:20:ARG:HH21	1:D:104:MSE:CE	2.23	0.51
1:B:92:GLU:H	1:B:92:GLU:CD	2.13	0.51
1:C:42:LEU:HD22	1:C:46:ILE:HG13	1.93	0.51
1:B:4:LEU:C	1:B:4:LEU:HD23	2.32	0.50
1:B:143:LEU:O	1:B:145:LEU:HD13	2.12	0.49
1:D:4:LEU:C	1:D:4:LEU:HD23	2.33	0.49
1:B:125:ILE:CD1	1:C:121:LYS:HB3	2.40	0.48
1:C:57:LYS:HG3	1:C:58:GLU:OE2	2.13	0.48
1:B:134:GLU:O	1:B:138:GLN:HG3	2.13	0.48
1:A:17:LYS:O	1:A:21:VAL:HG23	2.13	0.48
1:B:138:GLN:HB3	1:D:23:PHE:CG	2.48	0.48
1:B:141:THR:HB	1:D:109:GLU:HG2	1.94	0.48
1:C:42:LEU:HD11	1:C:65:VAL:HG11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:LYS:HG3	1:D:58:GLU:OE2	2.14	0.48
1:A:138:GLN:HB3	1:C:23:PHE:CG	2.48	0.48
1:B:57:LYS:HG3	1:B:58:GLU:OE2	2.13	0.48
1:A:121:LYS:HB3	1:D:125:ILE:CD1	2.41	0.47
1:A:125:ILE:CD1	1:D:121:LYS:HB3	2.43	0.47
1:B:96:LEU:O	1:B:100:GLU:HG3	2.14	0.47
1:A:57:LYS:HG3	1:A:58:GLU:OE2	2.15	0.47
1:C:104:MSE:CE	1:D:20:ARG:HH21	2.28	0.47
1:B:136:MSE:HG2	1:C:111:THR:HG21	1.96	0.46
1:B:17:LYS:O	1:B:21:VAL:HG23	2.14	0.46
1:C:3:LYS:HB2	1:C:3:LYS:HE3	1.56	0.46
1:D:96:LEU:O	1:D:100:GLU:HG3	2.14	0.46
1:C:96:LEU:O	1:C:100:GLU:HG3	2.14	0.46
1:C:101:VAL:O	1:C:105:GLU:HG3	2.15	0.46
1:A:96:LEU:HG	1:B:8:PHE:HE2	1.80	0.46
1:A:104:MSE:HE2	1:B:20:ARG:HH21	1.79	0.46
1:D:17:LYS:O	1:D:21:VAL:HG23	2.16	0.45
1:D:3:LYS:HE3	1:D:3:LYS:HB2	1.57	0.45
1:C:17:LYS:O	1:C:21:VAL:HG23	2.17	0.45
1:C:62:ARG:HD3	1:C:84:GLU:OE1	2.17	0.45
1:D:62:ARG:HD3	1:D:84:GLU:OE1	2.17	0.45
1:B:22:LEU:HD11	1:B:74:ILE:HD13	1.99	0.45
1:B:64:SER:O	1:B:81:LEU:HD12	2.16	0.45
1:A:36:VAL:HG11	1:C:148:GLU:HG2	1.98	0.45
1:B:8:PHE:CE1	1:B:62:ARG:HD2	2.52	0.45
1:B:3:LYS:HE3	1:B:3:LYS:HB2	1.57	0.45
1:B:75:VAL:O	1:B:75:VAL:CG2	2.65	0.44
1:B:101:VAL:O	1:B:105:GLU:HG3	2.18	0.44
1:C:64:SER:O	1:C:81:LEU:HD12	2.17	0.43
1:C:14:TYR:CG	1:C:15:ALA:N	2.86	0.43
1:B:135:ARG:HG3	1:D:27:ARG:NH1	2.34	0.43
1:A:64:SER:O	1:A:81:LEU:HD12	2.18	0.43
1:A:91:GLU:OE2	1:B:93:LYS:HE3	2.18	0.43
1:B:121:LYS:HB3	1:C:125:ILE:CD1	2.47	0.43
1:A:101:VAL:O	1:A:105:GLU:HG3	2.18	0.43
1:C:75:VAL:O	1:C:75:VAL:CG2	2.65	0.43
1:B:14:TYR:CG	1:B:15:ALA:N	2.86	0.43
1:C:92:GLU:CD	1:C:92:GLU:H	2.20	0.43
1:A:96:LEU:O	1:A:100:GLU:HG3	2.19	0.43
1:B:62:ARG:HD3	1:B:84:GLU:OE1	2.19	0.43
1:C:20:ARG:HH21	1:D:104:MSE:HE2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:PHE:CE1	1:C:62:ARG:HD2	2.54	0.43
1:D:101:VAL:O	1:D:105:GLU:HG3	2.19	0.43
1:A:8:PHE:CE1	1:A:62:ARG:HD2	2.54	0.42
1:D:64:SER:O	1:D:81:LEU:HD12	2.18	0.42
1:B:33:GLU:CD	1:B:33:GLU:H	2.21	0.42
1:A:3:LYS:HE3	1:A:3:LYS:HB2	1.57	0.42
1:D:14:TYR:CG	1:D:15:ALA:N	2.87	0.42
1:A:62:ARG:HD3	1:A:84:GLU:OE1	2.19	0.41
1:B:30:VAL:HG13	1:B:34:GLU:HB3	2.02	0.41
1:D:8:PHE:CE1	1:D:62:ARG:HD2	2.55	0.41
1:D:22:LEU:HD11	1:D:74:ILE:HD13	2.02	0.41
1:C:33:GLU:CD	1:C:33:GLU:H	2.23	0.41
1:A:30:VAL:HG13	1:A:34:GLU:HB3	2.03	0.41
1:A:96:LEU:HD23	1:B:8:PHE:CE2	2.56	0.41
1:C:30:VAL:HG13	1:C:34:GLU:HB3	2.01	0.41
1:C:134:GLU:O	1:C:138:GLN:HG3	2.21	0.41
1:B:44:LYS:HE3	1:B:44:LYS:HB3	1.96	0.41
1:A:14:TYR:CG	1:A:15:ALA:N	2.89	0.41
1:A:96:LEU:CG	1:B:8:PHE:HE2	2.34	0.41
1:D:33:GLU:CD	1:D:33:GLU:H	2.23	0.40
1:D:75:VAL:O	1:D:75:VAL:CG2	2.69	0.40
1:A:8:PHE:CE2	1:B:96:LEU:HD23	2.55	0.40
1:A:75:VAL:O	1:A:75:VAL:CG2	2.68	0.40
1:D:30:VAL:HG13	1:D:34:GLU:HB3	2.03	0.40
1:C:22:LEU:HD11	1:C:74:ILE:HD13	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:306:HOH:O	3:C:306:HOH:O[4_565]	1.81	0.39

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	146/149 (98%)	143 (98%)	3 (2%)	0	100	100
1	B	145/149 (97%)	143 (99%)	2 (1%)	0	100	100
1	C	144/149 (97%)	142 (99%)	2 (1%)	0	100	100
1	D	146/149 (98%)	144 (99%)	2 (1%)	0	100	100
All	All	581/596 (98%)	572 (98%)	9 (2%)	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	133/126 (106%)	118 (89%)	15 (11%)	6	5
1	B	132/126 (105%)	121 (92%)	11 (8%)	11	12
1	C	131/126 (104%)	117 (89%)	14 (11%)	6	6
1	D	133/126 (106%)	120 (90%)	13 (10%)	8	8
All	All	529/504 (105%)	476 (90%)	53 (10%)	8	7

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	33	GLU
1	A	37	LYS
1	A	68	ASN
1	A	75	VAL
1	A	78	LEU
1	A	82	GLU
1	A	85	THR
1	A	89	GLU
1	A	92[A]	GLU

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Mol	Chain	Res	Type
1	A	92[B]	GLU
1	A	103	HIS
1	A	118	LEU
1	A	131	GLU
1	A	143	LEU
1	B	33	GLU
1	B	37	LYS
1	B	68	ASN
1	B	75	VAL
1	B	82	GLU
1	B	85	THR
1	B	89	GLU
1	B	103	HIS
1	B	131	GLU
1	B	143	LEU
1	B	145	LEU
1	C	4	LEU
1	C	33	GLU
1	C	37	LYS
1	C	42	LEU
1	C	50	LEU
1	C	68	ASN
1	C	75	VAL
1	C	82	GLU
1	C	85	THR
1	C	89	GLU
1	C	92	GLU
1	C	103	HIS
1	C	115	LEU
1	C	131	GLU
1	D	33	GLU
1	D	37	LYS
1	D	68	ASN
1	D	75	VAL
1	D	78	LEU
1	D	82	GLU
1	D	85	THR
1	D	89	GLU
1	D	92[A]	GLU
1	D	92[B]	GLU
1	D	103	HIS
1	D	118	LEU

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Mol	Chain	Res	Type
1	D	131	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	A	201	-	4,4,4	0.16	0	6,6,6	0.07	0
2	SO4	B	201	-	4,4,4	0.12	0	6,6,6	0.13	0
2	SO4	D	201	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	C	201	-	4,4,4	0.12	0	6,6,6	0.13	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	SO4	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	141/149 (94%)	0.05	1 (0%)	87	87	28, 49, 86, 108	0
1	B	140/149 (93%)	0.36	8 (5%)	23	20	34, 60, 100, 109	0
1	C	140/149 (93%)	0.52	13 (9%)	8	6	33, 63, 102, 109	0
1	D	141/149 (94%)	0.05	4 (2%)	53	49	30, 50, 88, 110	0
All	All	562/596 (94%)	0.24	26 (4%)	32	30	28, 55, 96, 110	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	76	TRP	5.9
1	C	69	ILE	3.7
1	C	67	PHE	3.4
1	C	66	ASP	3.4
1	B	66	ASP	3.3
1	C	31	PHE	3.1
1	C	71	ASP	3.0
1	C	55	VAL	2.9
1	C	148	GLU	2.9
1	D	147	SER	2.8
1	B	76	TRP	2.8
1	C	70	GLU	2.8
1	D	146	ARG	2.8
1	C	74	ILE	2.7
1	B	148	GLU	2.7
1	B	31	PHE	2.7
1	C	81	LEU	2.7
1	B	146	ARG	2.6
1	C	79	ASP	2.6
1	B	71	ASP	2.5
1	A	148	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	80	SER	2.2
1	B	103	HIS	2.1
1	B	52	GLU	2.1
1	D	148	GLU	2.0
1	D	68	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
2	SO4	C	201	5/5	0.88	0.13	66,99,156,178	0
2	SO4	B	201	5/5	0.89	0.15	79,85,159,160	0
2	SO4	A	201	5/5	0.91	0.18	53,132,142,163	0
2	SO4	D	201	5/5	0.95	0.21	43,63,81,81	5

6.5 Other polymers (i)

There are no such residues in this entry.