Full wwPDB X-ray Structure Validation Report (i)

Sep 29, 2024 – 09:05 AM EDT

PDB ID	:	3UZ0
Title	:	Crystal Structure of SpoIIIAH and SpoIIQ Complex
Authors	:	Maehigashi, T.; Meisner, J.; Dunham, C.M.
Deposited on	:	2011-12-07
Resolution	:	2.82 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:

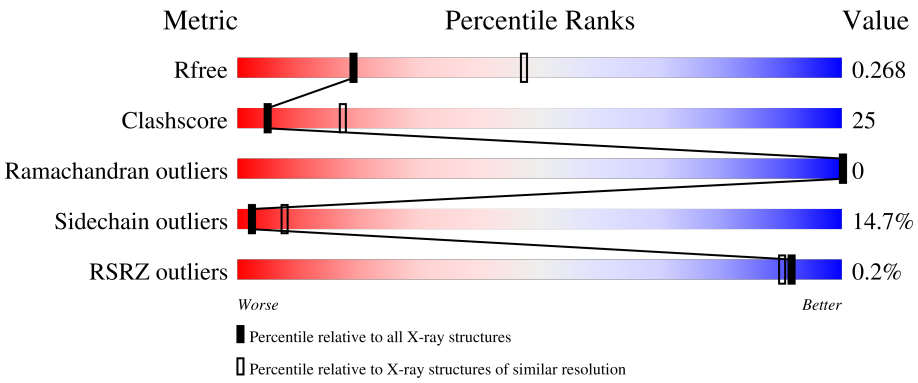
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.82 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}	164625	4293 (2.84-2.80)
Clashscore	180529	4801 (2.84-2.80)
Ramachandran outliers	177936	4739 (2.84-2.80)
Sidechain outliers	177891	4741 (2.84-2.80)
RSRZ outliers	164620	4295 (2.84-2.80)

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	133	53% 29% • 14%
1	C	133	52% 26% 7% 15%
2	B	152	% 47% 39% 7% 6%
2	D	152	40% 45% 9% 6%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3892 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 Stage III sporulation protein AH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	114	870	536	140	192	2	0	0	0
1	C	113	858	529	137	190	2	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	GLY	-	expression tag	UNP P49785
A	87	SER	-	expression tag	UNP P49785
A	88	HIS	-	expression tag	UNP P49785
A	89	MSE	-	expression tag	UNP P49785
C	86	GLY	-	expression tag	UNP P49785
C	87	SER	-	expression tag	UNP P49785
C	88	HIS	-	expression tag	UNP P49785
C	89	MSE	-	expression tag	UNP P49785

- Molecule 2 is a protein called Stage II sporulation protein Q.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
2	B	143	1067	665	174	225	3	0	0	0
2	D	143	1055	660	174	218	3	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	69	GLY	-	expression tag	UNP P71044
B	70	SER	-	expression tag	UNP P71044
B	71	HIS	-	expression tag	UNP P71044
B	72	MSE	-	expression tag	UNP P71044

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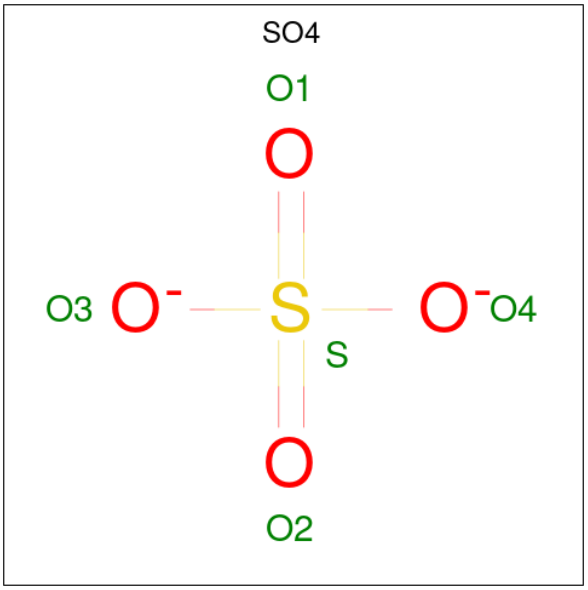
Chain	Residue	Modelled	Actual	Comment	Reference
D	69	GLY	-	expression tag	UNP P71044
D	70	SER	-	expression tag	UNP P71044
D	71	HIS	-	expression tag	UNP P71044
D	72	MSE	-	expression tag	UNP P71044

- 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
3	C	1	5	4	1	0	0
3	D	1	5	4	1	0	0

- Molecule 4 is water.

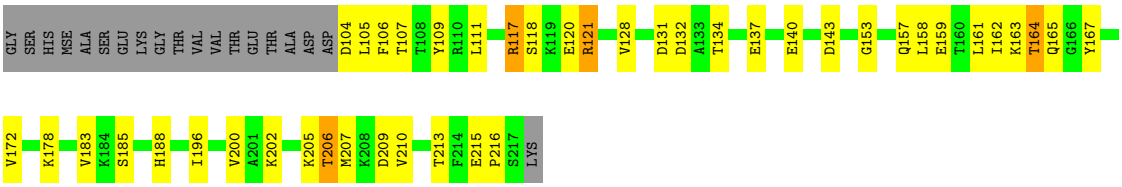
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	5	5	5	0	0
4	B	9	9	9	0	0
4	C	4	4	4	0	0
4	D	9	9	9	0	0

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.

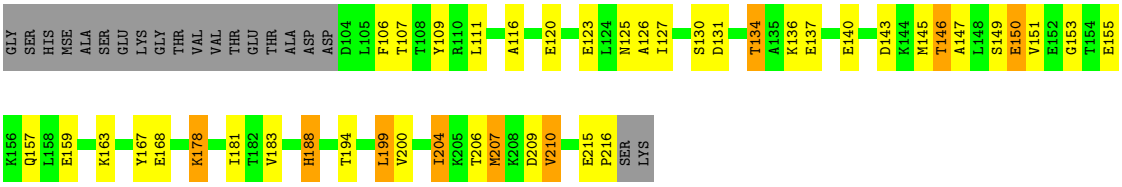
• Molecule 1: Stage III sporulation protein AH

Chain A:



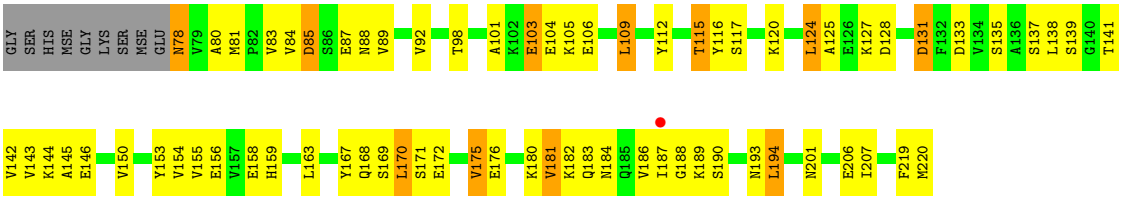
• Molecule 1: Stage III sporulation protein AH

Chain C:



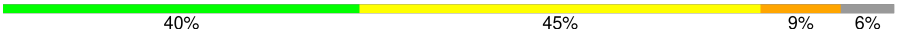
• Molecule 2: Stage II sporulation protein Q

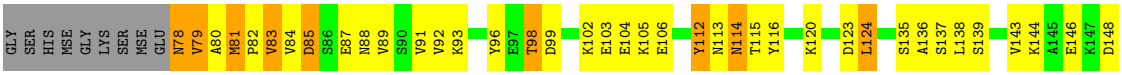
Chain B:



• Molecule 2: Stage II sporulation protein Q

Chain D:





P149
Y153
V154
V155
E156
V157
E158
H159
A160
D161
G162
L163
V166
Y167
Q168
S169
L170
V173
S174
V175
E176
Q177
K182
Q183
N184
Q185
V186
I187
G188
K192
N193
S196
E197
D198
S199
G200
N201
H202
V203
E206
I207
R208
K209
D210
G211
V212
A213
M214
N215
P216
F219
M220

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.37Å 64.49Å 176.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.40 – 2.82 36.40 – 2.82	Depositor EDS
% Data completeness (in resolution range)	98.8 (36.40-2.82) 98.7 (36.40-2.82)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.86 (at 2.72Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.208 , 0.270 0.209 , 0.268	Depositor DCC
R_{free} test set	1266 reflections (7.10%)	wwPDB-VP
Wilson B-factor (Å^2)	75.3	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 65.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.045 for k,h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3892	wwPDB-VP
Average B, all atoms (Å^2)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% 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	0.50	0/874	0.64	0/1177
1	C	0.46	0/862	0.66	0/1163
2	B	0.47	0/1079	0.68	0/1459
2	D	0.46	0/1067	0.70	1/1443 (0.1%)
All	All	0.47	0/3882	0.67	1/5242 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	112	TYR	N-CA-C	-5.09	97.27	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	206	THR	Peptide

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	870	0	845	25	0
1	C	858	0	822	30	0
2	B	1067	0	1008	58	0
2	D	1055	0	998	79	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
4	A	5	0	0	0	0
4	B	9	0	0	0	0
4	C	4	0	0	0	0
4	D	9	0	0	0	0
All	All	3892	0	3673	185	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 25.

All (185) 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:C:199:LEU:HD12	1:C:200:VAL:N	1.86	0.90
1:C:178:LYS:HG2	1:C:209:ASP:HB3	1.55	0.88
2:D:81:MSE:HG3	2:D:82:PRO:HD2	1.55	0.87
2:D:139:SER:HB3	2:D:183:GLN:N	1.90	0.86
2:D:84:VAL:HG23	2:D:85:ASP:H	1.40	0.84
2:B:172:GLU:O	2:B:188:GLY:HA3	1.77	0.83
2:B:142:VAL:HG21	2:B:175:VAL:HG21	1.62	0.81
2:D:139:SER:HB3	2:D:183:GLN:H	1.44	0.80
1:C:125:ASN:OD1	1:C:145:MSE:HE1	1.81	0.80
2:B:175:VAL:HG23	2:B:176:GLU:N	1.96	0.79
2:B:187:ILE:HG22	2:B:188:GLY:H	1.48	0.77
1:C:134:THR:HG22	1:C:137:GLU:H	1.49	0.77
1:C:178:LYS:CG	1:C:209:ASP:HB3	2.15	0.76
2:D:83:VAL:HG21	2:D:89:VAL:HG11	1.68	0.76
1:A:117:ARG:HH11	1:A:117:ARG:HG2	1.51	0.76
2:B:83:VAL:HG11	2:B:89:VAL:HG21	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:139:SER:HB3	2:B:183:GLN:N	2.03	0.74
2:B:127:LYS:O	2:B:128:ASP:OD1	2.06	0.72
2:D:175:VAL:HG22	2:D:176:GLU:N	2.05	0.71
2:B:80:ALA:HB3	2:B:137:SER:HA	1.71	0.71
2:D:79:VAL:HG11	2:D:219:PHE:HB3	1.74	0.70
1:A:106:PHE:O	1:A:109:TYR:HB2	1.92	0.69
2:B:175:VAL:CG2	2:B:176:GLU:N	2.55	0.69
2:D:124:LEU:HB2	2:D:203:VAL:HG12	1.74	0.69
2:B:183:GLN:HG2	2:B:184:ASN:ND2	2.08	0.68
1:C:123:GLU:O	1:C:127:ILE:HD13	1.94	0.68
2:B:112:TYR:CE2	2:D:197:GLU:HG2	2.29	0.67
1:A:134:THR:OG1	1:A:137:GLU:HG3	1.95	0.67
2:D:124:LEU:HB2	2:D:203:VAL:CG1	2.24	0.67
2:D:143:VAL:CG2	2:D:156:GLU:HG2	2.24	0.67
2:D:79:VAL:CG1	2:D:219:PHE:HB3	2.26	0.66
2:B:101:ALA:O	2:B:104:GLU:HG2	1.96	0.66
2:D:154:VAL:HG23	2:D:168:GLN:HG2	1.78	0.66
2:D:98:THR:O	2:D:105:LYS:HE3	1.96	0.65
1:A:153:GLY:O	1:A:157:GLN:HG3	1.98	0.63
2:D:139:SER:H	2:D:183:GLN:HB2	1.63	0.63
2:B:120:LYS:O	2:B:206:GLU:HG2	1.98	0.63
2:D:120:LYS:O	2:D:206:GLU:HG2	1.99	0.63
2:D:175:VAL:CG2	2:D:176:GLU:H	2.11	0.62
2:D:175:VAL:HG22	2:D:176:GLU:H	1.65	0.62
1:A:137:GLU:HA	1:A:140:GLU:OE1	2.00	0.62
2:B:87:GLU:O	2:B:88:ASN:HB2	2.00	0.61
2:D:175:VAL:CG2	2:D:176:GLU:N	2.63	0.61
2:B:84:VAL:HG23	2:B:85:ASP:H	1.66	0.61
2:D:187:ILE:HG13	2:D:187:ILE:O	2.01	0.60
2:B:171:SER:HB3	2:B:190:SER:O	2.02	0.59
1:A:213:THR:OG1	2:B:115:THR:HB	2.02	0.59
2:B:139:SER:HB3	2:B:183:GLN:H	1.64	0.59
1:A:161:LEU:O	1:A:164:THR:HB	2.03	0.58
2:D:174:SER:H	2:D:188:GLY:H	1.50	0.58
2:D:79:VAL:HA	2:D:138:LEU:HD13	1.84	0.58
2:D:84:VAL:HG23	2:D:85:ASP:N	2.16	0.58
2:D:81:MSE:HG3	2:D:82:PRO:CD	2.30	0.58
2:D:85:ASP:OD1	2:D:85:ASP:C	2.42	0.57
2:D:102:LYS:C	2:D:104:GLU:H	2.08	0.57
1:C:126:ALA:O	1:C:130:SER:HB2	2.04	0.57
2:D:84:VAL:CG1	2:D:135:SER:HB3	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:84:VAL:HG23	2:B:85:ASP:N	2.20	0.56
2:D:183:GLN:HG2	2:D:184:ASN:HD22	1.70	0.56
2:B:150:VAL:HA	2:B:194:LEU:HD22	1.87	0.56
2:B:98:THR:HA	2:B:105:LYS:HE2	1.88	0.56
2:D:87:GLU:O	2:D:88:ASN:HB2	2.06	0.56
2:B:138:LEU:O	2:B:159:HIS:CE1	2.58	0.56
1:C:147:ALA:HA	1:C:150:GLU:HG3	1.88	0.55
1:C:140:GLU:HA	1:C:140:GLU:OE1	2.05	0.55
2:D:183:GLN:HG2	2:D:184:ASN:ND2	2.21	0.55
1:A:104:ASP:HB3	1:A:107:THR:HB	1.89	0.55
2:D:144:LYS:HA	2:D:177:GLN:HG2	1.89	0.55
1:C:134:THR:HG22	1:C:137:GLU:N	2.20	0.55
2:D:163:LEU:HD23	2:D:209:LYS:HD3	1.88	0.54
1:A:188:HIS:HD2	2:B:116:TYR:OH	1.90	0.54
2:D:143:VAL:O	2:D:144:LYS:HB3	2.06	0.54
2:B:78:ASN:HD22	2:B:78:ASN:N	2.06	0.54
1:C:151:VAL:O	1:C:155:GLU:HG3	2.08	0.54
2:D:173:VAL:HA	2:D:188:GLY:HA2	1.89	0.54
1:C:116:ALA:O	1:C:120:GLU:HG3	2.08	0.53
2:B:175:VAL:HG23	2:B:176:GLU:H	1.70	0.53
2:B:183:GLN:HG2	2:B:184:ASN:HD22	1.73	0.53
1:A:158:LEU:O	1:A:162:ILE:HG13	2.08	0.53
1:A:167:TYR:CD1	1:A:183:VAL:HG11	2.44	0.53
2:B:143:VAL:CG2	2:B:158:GLU:HB2	2.38	0.53
2:D:93:LYS:HB3	2:D:123:ASP:HB2	1.90	0.53
1:A:205:LYS:O	1:A:206:THR:HB	2.09	0.52
2:D:169:SER:HB2	2:D:193:ASN:ND2	2.24	0.52
1:C:167:TYR:CD1	1:C:183:VAL:HG11	2.44	0.52
2:B:103:GLU:H	2:B:106:GLU:HG3	1.75	0.52
2:D:167:TYR:CD2	2:D:167:TYR:N	2.78	0.52
1:A:205:LYS:O	1:A:206:THR:CB	2.58	0.52
2:B:139:SER:HB3	2:B:182:LYS:HA	1.92	0.52
2:B:81:MSE:HA	2:B:220:MSE:HE2	1.92	0.52
1:C:163:LYS:HE3	1:C:168:GLU:O	2.11	0.51
2:B:131:ASP:OD1	2:B:131:ASP:C	2.49	0.51
2:B:153:TYR:O	2:B:170:LEU:HB2	2.11	0.51
1:C:146:THR:O	1:C:149:SER:HB3	2.11	0.51
2:D:96:TYR:CE2	2:D:105:LYS:HB3	2.46	0.50
2:B:139:SER:HB2	2:B:181:VAL:O	2.11	0.50
1:C:188:HIS:NE2	2:D:114:ASN:HB2	2.27	0.50
1:A:178:LYS:HG3	1:A:209:ASP:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:143:VAL:HG13	2:D:158:GLU:HB2	1.94	0.50
2:D:197:GLU:O	2:D:199:SER:O	2.29	0.50
2:B:169:SER:HB2	2:B:193:ASN:ND2	2.27	0.50
2:B:219:PHE:CD1	2:B:219:PHE:N	2.80	0.49
2:D:103:GLU:H	2:D:106:GLU:HG3	1.77	0.49
1:C:204:ILE:HG13	1:C:204:ILE:O	2.03	0.49
2:D:208:ARG:HA	2:D:212:VAL:O	2.12	0.49
2:D:163:LEU:HD22	2:D:163:LEU:HA	1.67	0.49
2:D:206:GLU:OE2	2:D:208:ARG:NH2	2.46	0.49
2:D:143:VAL:HG22	2:D:156:GLU:HG2	1.95	0.48
2:D:136:ALA:O	2:D:184:ASN:N	2.38	0.48
2:D:81:MSE:CG	2:D:82:PRO:HD2	2.37	0.48
2:D:84:VAL:HG13	2:D:135:SER:HB3	1.95	0.48
2:D:182:LYS:O	2:D:185:GLN:HB3	2.14	0.48
2:B:187:ILE:HG22	2:B:188:GLY:N	2.23	0.48
2:D:84:VAL:HG12	2:D:135:SER:HB3	1.96	0.47
2:D:173:VAL:HA	2:D:188:GLY:CA	2.43	0.47
1:A:117:ARG:HG2	1:A:117:ARG:NH1	2.19	0.47
2:D:214:MSE:O	2:D:216:PRO:HD3	2.14	0.47
2:B:127:LYS:O	2:B:128:ASP:CG	2.52	0.47
2:B:112:TYR:OH	2:D:192:LYS:HB3	2.15	0.47
2:D:175:VAL:HG12	2:D:187:ILE:HB	1.97	0.47
2:B:106:GLU:HA	2:B:109:LEU:CD2	2.44	0.47
1:C:134:THR:O	1:C:137:GLU:HB2	2.14	0.47
2:D:82:PRO:O	2:D:135:SER:N	2.47	0.47
1:C:107:THR:O	1:C:111:LEU:HD12	2.15	0.46
2:D:80:ALA:HB3	2:D:137:SER:CB	2.45	0.46
2:B:150:VAL:O	2:B:193:ASN:HB2	2.15	0.46
2:D:200:GLY:O	2:D:202:HIS:HD2	1.98	0.46
1:C:188:HIS:O	2:D:116:TYR:OH	2.20	0.46
1:C:199:LEU:HD12	1:C:199:LEU:C	2.34	0.46
2:B:131:ASP:CG	2:B:189:LYS:HD3	2.36	0.46
1:C:155:GLU:O	1:C:159:GLU:HG3	2.15	0.46
2:D:153:TYR:HB2	2:D:170:LEU:O	2.16	0.46
2:B:141:THR:HG22	2:B:180:LYS:HD2	1.97	0.46
2:B:138:LEU:HA	2:B:183:GLN:HB2	1.98	0.46
1:A:206:THR:O	1:A:206:THR:HG23	2.14	0.46
2:B:137:SER:O	2:B:138:LEU:C	2.54	0.46
1:C:134:THR:HG22	1:C:137:GLU:HG3	1.97	0.45
2:D:92:VAL:HG23	2:D:93:LYS:H	1.81	0.45
2:B:139:SER:CB	2:B:182:LYS:HA	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:MSE:HE2	1:A:210:VAL:HG21	1.99	0.45
2:B:145:ALA:O	2:B:146:GLU:HB3	2.16	0.45
2:B:194:LEU:HD12	2:B:194:LEU:HA	1.71	0.45
2:B:167:TYR:N	2:B:167:TYR:CD2	2.86	0.44
2:B:143:VAL:HG21	2:B:158:GLU:HB2	1.98	0.44
2:B:154:VAL:HG12	2:B:168:GLN:HG2	1.98	0.44
1:C:106:PHE:O	1:C:109:TYR:N	2.49	0.44
1:A:196:ILE:O	1:A:200:VAL:HG23	2.18	0.44
2:D:160:ALA:O	2:D:161:ASP:HB2	2.18	0.44
2:D:78:ASN:N	2:D:78:ASN:HD22	2.16	0.43
2:B:154:VAL:O	2:B:154:VAL:HG23	2.18	0.43
2:D:219:PHE:N	2:D:219:PHE:CD1	2.86	0.43
2:B:175:VAL:HG23	2:B:176:GLU:O	2.18	0.43
2:D:153:TYR:HB2	2:D:170:LEU:HB2	1.99	0.43
2:D:155:VAL:HG13	2:D:167:TYR:HB2	2.00	0.43
2:B:131:ASP:OD1	2:B:189:LYS:HD3	2.17	0.43
2:D:102:LYS:O	2:D:104:GLU:N	2.48	0.43
2:D:139:SER:HB3	2:D:182:LYS:HA	2.01	0.43
2:D:174:SER:H	2:D:188:GLY:N	2.14	0.42
2:D:80:ALA:O	2:D:220:MSE:HE2	2.18	0.42
2:D:139:SER:CB	2:D:183:GLN:H	2.25	0.42
1:C:153:GLY:O	1:C:157:GLN:HG3	2.20	0.42
1:C:199:LEU:HD12	1:C:200:VAL:H	1.78	0.42
1:A:215:GLU:HA	1:A:216:PRO:HD3	1.84	0.42
1:C:215:GLU:OE1	1:C:216:PRO:HD2	2.20	0.42
2:D:148:ASP:HA	2:D:149:PRO:HD2	1.80	0.42
2:D:154:VAL:CG2	2:D:168:GLN:HG2	2.49	0.42
2:D:162:GLY:HA3	2:D:210:ASP:OD1	2.20	0.42
2:B:144:LYS:HB3	2:B:156:GLU:HB3	2.02	0.41
2:B:133:ASP:HB3	2:B:186:VAL:HG13	2.02	0.41
1:A:188:HIS:CD2	2:B:116:TYR:OH	2.71	0.41
2:D:168:GLN:O	2:D:169:SER:HB3	2.20	0.41
2:D:182:LYS:O	2:D:185:GLN:CB	2.69	0.41
1:C:207:MSE:HE1	1:C:210:VAL:HG11	2.02	0.41
1:A:165:GLN:H	1:A:165:GLN:HG2	1.74	0.41
2:B:135:SER:HB2	2:B:184:ASN:HA	2.03	0.41
1:A:120:GLU:O	1:A:121:ARG:C	2.59	0.41
1:A:164:THR:O	1:A:164:THR:HG23	2.21	0.41
1:A:159:GLU:HG3	1:A:172:VAL:HG23	2.03	0.41
1:A:159:GLU:O	1:A:163:LYS:HG3	2.21	0.40
1:C:134:THR:HG23	1:C:136:LYS:H	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:124:LEU:CB	2:D:203:VAL:HG12	2.48	0.40
2:D:166:VAL:HG21	2:D:208:ARG:NH1	2.37	0.40
2:B:124:LEU:O	2:B:125:ALA:HB2	2.21	0.40
1:C:181:ILE:HD11	1:C:200:VAL:HG21	2.03	0.40
2:D:200:GLY:O	2:D:202:HIS:CD2	2.75	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	112/133 (84%)	108 (96%)	4 (4%)	0	100	100
1	C	111/133 (84%)	100 (90%)	11 (10%)	0	100	100
2	B	141/152 (93%)	128 (91%)	13 (9%)	0	100	100
2	D	141/152 (93%)	133 (94%)	8 (6%)	0	100	100
All	All	505/570 (89%)	469 (93%)	36 (7%)	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	95/111 (86%)	82 (86%)	13 (14%)	3	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	C	92/111 (83%)	80 (87%)	12 (13%)	3 10
2	B	116/125 (93%)	99 (85%)	17 (15%)	2 8
2	D	113/125 (90%)	94 (83%)	19 (17%)	1 5
All	All	416/472 (88%)	355 (85%)	61 (15%)	2 8

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	LEU
1	A	111	LEU
1	A	117	ARG
1	A	118	SER
1	A	121	ARG
1	A	128	VAL
1	A	131	ASP
1	A	132	ASP
1	A	143	ASP
1	A	164	THR
1	A	185	SER
1	A	202	LYS
1	A	206	THR
2	B	78	ASN
2	B	85	ASP
2	B	92	VAL
2	B	103	GLU
2	B	109	LEU
2	B	115	THR
2	B	117	SER
2	B	124	LEU
2	B	131	ASP
2	B	155	VAL
2	B	163	LEU
2	B	170	LEU
2	B	175	VAL
2	B	181	VAL
2	B	194	LEU
2	B	201	ASN
2	B	207	ILE
1	C	131	ASP
1	C	134	THR
1	C	143	ASP

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Mol	Chain	Res	Type
1	C	146	THR
1	C	150	GLU
1	C	178	LYS
1	C	188	HIS
1	C	194	THR
1	C	199	LEU
1	C	204	ILE
1	C	207	MSE
1	C	210	VAL
2	D	78	ASN
2	D	79	VAL
2	D	81	MSE
2	D	83	VAL
2	D	85	ASP
2	D	91	VAL
2	D	98	THR
2	D	99	ASP
2	D	112	TYR
2	D	113	ASN
2	D	114	ASN
2	D	115	THR
2	D	124	LEU
2	D	146	GLU
2	D	155	VAL
2	D	163	LEU
2	D	174	SER
2	D	177	GLN
2	D	196	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	188	HIS
2	B	114	ASN
2	B	184	ASN
2	D	184	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	1	-	4,4,4	0.33	0	6,6,6	0.32	0
3	SO4	C	2	-	4,4,4	0.26	0	6,6,6	0.40	0
3	SO4	D	3	-	4,4,4	0.28	0	6,6,6	0.20	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	112/133 (84%)	-0.57	0	100	100	51, 73, 105, 123	0
1	C	111/133 (83%)	-0.51	0	100	100	52, 77, 117, 129	0
2	B	140/152 (92%)	-0.31	1 (0%)	84	79	49, 82, 113, 120	0
2	D	140/152 (92%)	-0.37	0	100	100	49, 78, 111, 121	0
All	All	503/570 (88%)	-0.43	1 (0%)	92	90	49, 78, 113, 129	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	187	ILE	2.2

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	A	1	5/5	0.80	0.12	90,94,118,130	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SO4	C	2	5/5	0.81	0.10	70,90,120,130	0
3	SO4	D	3	5/5	0.85	0.12	81,97,140,146	0

6.5 Other polymers (i)

There are no such residues in this entry.