



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 11:36 pm BST

PDB ID : 1YEP
Title : Structural and biochemical analysis of the link between enzymatic activity and oligomerization in AhpC, a bacterial peroxiredoxin.
Authors : Parsonage, D.; Youngblood, D.S.; Sarma, G.N.; Wood, Z.A.; Karplus, P.A.; Poole, L.B.
Deposited on : 2004-12-28
Resolution : 2.50 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 : 2.11
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) : 2.11

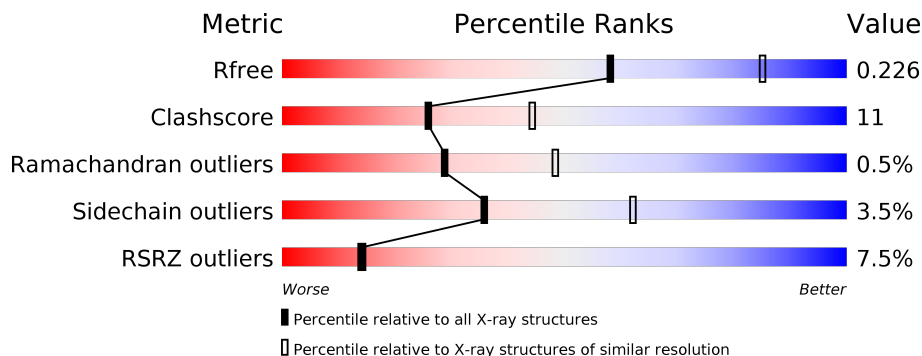
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	186	<div style="display: flex; align-items: center;"> <div style="width: 9%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 73%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 17%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">9% 73% 17% • 10%</p>
1	B	186	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 76%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 76% 12% • 11%</p>
1	C	186	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 67%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">6% 67% 21% 12%</p>
1	D	186	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 70%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 19%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">4% 70% 19% • 11%</p>
1	E	186	<div style="display: flex; align-items: center;"> <div style="width: 11%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 58%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">11% 58% 27% • • 11%</p>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	C	803	-	-	-	X
2	SO4	E	805	-	-	-	X

2 Entry composition [i](#)

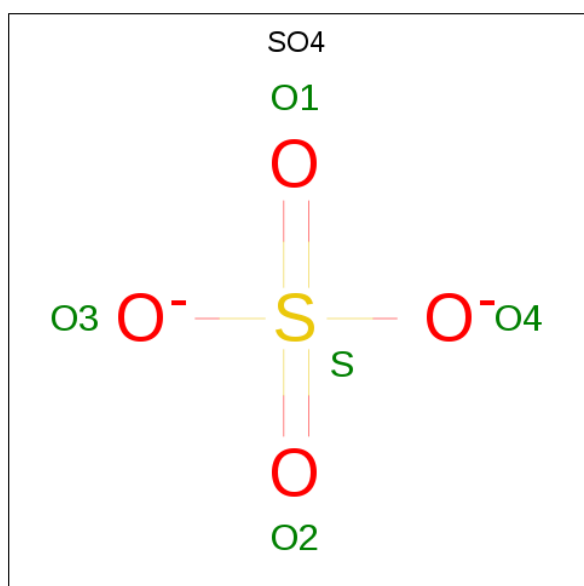
There are 3 unique types of molecules in this entry. The entry contains 6866 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alkyl hydroperoxide reductase subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	168	Total 1320	C 840	N 220	O 256	S 4	0	0	0
1	B	165	Total 1299	C 826	N 216	O 253	S 4	0	0	0
1	C	163	Total 1286	C 818	N 214	O 251	S 3	0	0	0
1	D	166	Total 1306	C 831	N 217	O 254	S 4	0	0	0
1	E	165	Total 1299	C 826	N 216	O 253	S 4	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

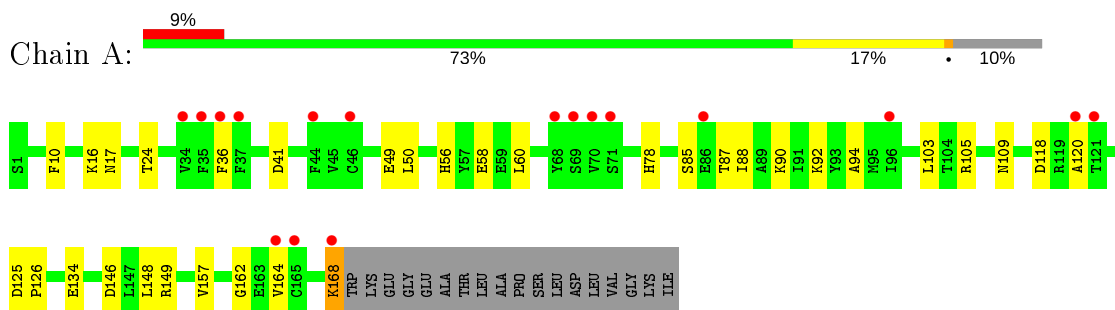
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	77	Total	O	0	0
			77	77		
3	B	88	Total	O	0	0
			88	88		
3	C	67	Total	O	0	0
			67	67		
3	D	58	Total	O	0	0
			58	58		
3	E	21	Total	O	0	0
			21	21		

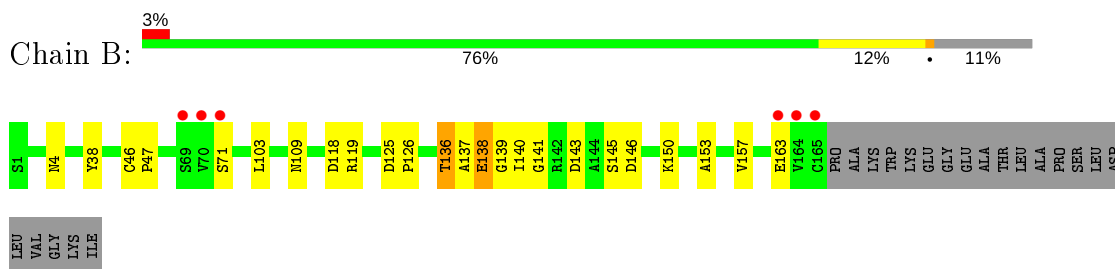
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

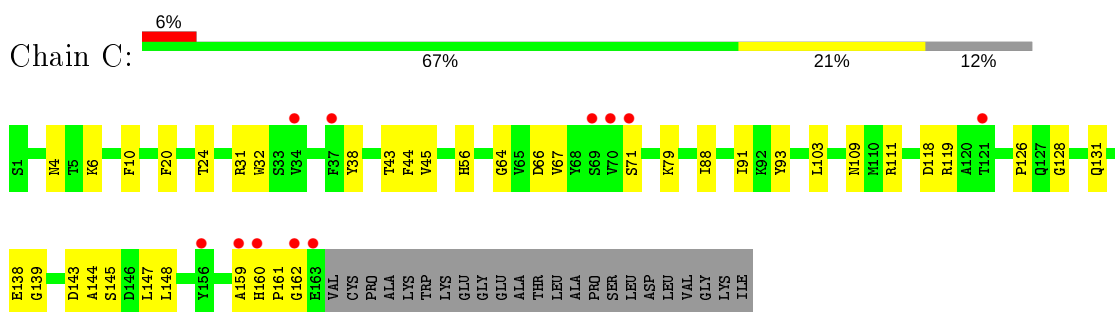
- Molecule 1: Alkyl hydroperoxide reductase subunit C



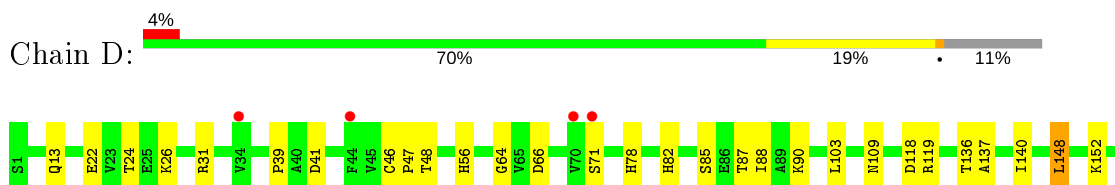
- Molecule 1: Alkyl hydroperoxide reductase subunit C

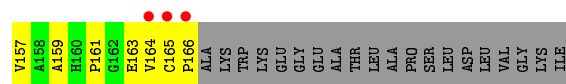


- Molecule 1: Alkyl hydroperoxide reductase subunit C



- Molecule 1: Alkyl hydroperoxide reductase subunit C





- Molecule 1: Alkyl hydroperoxide reductase subunit C



GLY
LYS
ILE

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	127.47Å 171.63Å 135.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 24.99 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.50) 95.4 (24.99-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.41Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.178 , 0.226 0.178 , 0.226	Depositor DCC
R_{free} test set	5128 reflections (8.90%)	wwPDB-VP
Wilson B-factor (Å ²)	49.4	Xtrriage
Anisotropy	0.145	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6866	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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	0/1351	0.67	0/1831
1	B	0.49	0/1329	0.67	0/1801
1	C	0.46	0/1316	0.65	1/1783 (0.1%)
1	D	0.44	0/1337	0.64	0/1813
1	E	0.37	0/1329	0.60	0/1801
All	All	0.45	0/6662	0.65	1/9029 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	111	ARG	N-CA-C	-5.00	97.49	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1320	0	1276	22	0
1	B	1299	0	1251	17	0
1	C	1286	0	1238	22	0
1	D	1306	0	1259	29	0
1	E	1299	0	1253	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
2	E	5	0	0	0	0
3	A	77	0	0	1	0
3	B	88	0	0	1	0
3	C	67	0	0	0	0
3	D	58	0	0	1	0
3	E	21	0	0	0	0
All	All	6866	0	6277	135	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 11.

All (135) 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:119:ARG:HE	1:C:138:GLU:HG2	1.04	1.11
1:C:119:ARG:NE	1:C:138:GLU:HG2	1.85	0.91
1:E:111:ARG:HH11	1:E:118:ASP:HA	1.40	0.87
1:E:119:ARG:HE	1:E:138:GLU:HG2	1.47	0.79
1:D:148:LEU:HD22	1:D:152:LYS:NZ	1.98	0.79
1:D:31:ARG:HD3	1:D:66:ASP:OD2	1.88	0.74
1:C:56:HIS:HD2	1:C:148:LEU:HD12	1.53	0.74
1:D:148:LEU:HD22	1:D:152:LYS:HZ2	1.58	0.67
1:E:149:ARG:HH11	1:E:149:ARG:HB3	1.61	0.66
1:B:119:ARG:HD3	3:B:2004:HOH:O	1.97	0.64
1:E:39:PRO:HG3	1:E:111:ARG:HH12	1.63	0.63
1:E:36:PHE:HZ	1:E:49:GLU:HB3	1.63	0.63
1:C:109:ASN:ND2	1:C:118:ASP:HB2	2.13	0.62
1:C:139:GLY:HA3	1:D:157:VAL:HG11	1.81	0.62
1:B:109:ASN:ND2	1:B:118:ASP:HB2	2.15	0.62
1:D:164:VAL:HG12	1:D:165:CYS:N	2.14	0.62
1:C:109:ASN:HD21	1:C:118:ASP:HB2	1.64	0.62
1:E:149:ARG:HB3	1:E:149:ARG:NH1	2.15	0.61
1:E:147:LEU:O	1:E:151:ILE:HG13	2.01	0.61
1:B:136:THR:HG21	1:B:140:ILE:O	2.01	0.61
1:A:87:THR:O	1:A:90:LYS:HG2	2.00	0.61
1:D:109:ASN:ND2	1:D:118:ASP:HB2	2.15	0.60
1:C:6:LYS:HA	1:C:128:GLY:O	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:39:PRO:HG2	3:D:4064:HOH:O	2.01	0.60
1:A:17:ASN:HD21	1:A:92:LYS:HA	1.67	0.59
1:D:148:LEU:HD22	1:D:152:LYS:HZ1	1.69	0.56
1:E:56:HIS:HD2	1:E:148:LEU:HD12	1.71	0.56
1:A:85:SER:HB3	1:A:88:ILE:HB	1.88	0.56
1:B:109:ASN:HD21	1:B:118:ASP:HB2	1.71	0.56
1:E:159:ALA:O	1:E:161:PRO:HD3	2.06	0.56
1:D:31:ARG:NH2	1:D:64:GLY:HA2	2.21	0.55
1:C:56:HIS:HD2	1:C:148:LEU:CD1	2.20	0.55
1:B:163:GLU:HA	1:B:163:GLU:OE1	2.08	0.54
1:A:109:ASN:ND2	1:A:118:ASP:HB2	2.23	0.54
1:D:87:THR:O	1:D:90:LYS:HG3	2.08	0.53
1:E:30:GLY:HA2	1:E:126:PRO:HB3	1.91	0.53
1:A:168:LYS:O	1:A:168:LYS:HE3	2.09	0.53
1:D:137:ALA:HB3	1:D:140:ILE:HD12	1.91	0.53
1:D:109:ASN:HD21	1:D:118:ASP:HB2	1.73	0.53
1:E:82:HIS:HA	1:E:88:ILE:HG22	1.91	0.53
1:C:160:HIS:N	1:C:161:PRO:HD3	2.24	0.53
1:E:119:ARG:HE	1:E:138:GLU:HA	1.74	0.52
1:E:101:GLY:O	1:E:105:ARG:HG3	2.09	0.52
1:E:164:VAL:HG12	1:E:165:CYS:N	2.24	0.52
1:A:16:LYS:HG2	1:A:94:ALA:HA	1.92	0.52
1:C:31:ARG:HD3	1:C:66:ASP:OD2	2.09	0.52
1:E:86:GLU:HA	1:E:86:GLU:OE1	2.10	0.51
1:E:161:PRO:O	1:E:162:GLY:C	2.48	0.51
1:B:153:ALA:O	1:B:157:VAL:HG23	2.11	0.51
1:D:82:HIS:HA	1:D:88:ILE:HG22	1.92	0.51
1:E:119:ARG:NE	1:E:138:GLU:HG2	2.22	0.51
1:A:146:ASP:OD1	1:A:149:ARG:NH1	2.36	0.51
1:C:67:VAL:O	1:C:93:TYR:HB2	2.11	0.51
1:B:119:ARG:HB2	1:B:136:THR:O	2.10	0.51
1:C:88:ILE:HG23	1:C:91:ILE:HD12	1.93	0.51
1:E:82:HIS:O	1:E:82:HIS:ND1	2.44	0.51
1:E:24:THR:HG23	1:E:27:ASP:OD1	2.11	0.51
1:C:10:PHE:O	1:C:24:THR:HA	2.11	0.50
1:A:56:HIS:CD2	1:A:148:LEU:HD12	2.46	0.50
1:C:143:ASP:OD2	1:C:145:SER:HB2	2.11	0.50
1:E:119:ARG:HB2	1:E:136:THR:O	2.12	0.50
1:D:164:VAL:CG1	1:D:165:CYS:N	2.76	0.49
1:E:151:ILE:O	1:E:155:GLN:HB2	2.11	0.49
1:E:149:ARG:HH22	1:E:150:LYS:HE3	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:ASP:HB2	1:E:126:PRO:CD	2.42	0.49
1:C:44:PHE:HB3	1:D:164:VAL:HG13	1.95	0.49
1:A:36:PHE:HZ	1:A:49:GLU:HB3	1.78	0.49
1:B:146:ASP:OD2	1:B:150:LYS:HE2	2.13	0.48
1:E:140:ILE:N	1:E:140:ILE:HD12	2.28	0.48
1:A:56:HIS:HD2	1:A:148:LEU:HD12	1.79	0.48
1:D:47:PRO:HB3	1:D:88:ILE:HD11	1.94	0.48
1:A:162:GLY:C	1:A:164:VAL:H	2.17	0.48
1:B:125:ASP:HB2	1:B:126:PRO:CD	2.44	0.48
1:E:6:LYS:HA	1:E:128:GLY:O	2.13	0.48
1:C:38:TYR:CZ	1:C:71:SER:HB3	2.49	0.48
1:C:32:TRP:CE2	1:C:126:PRO:HD3	2.48	0.47
1:B:119:ARG:HE	1:B:138:GLU:CD	2.18	0.46
1:D:39:PRO:HD3	1:D:119:ARG:HG2	1.96	0.46
1:B:137:ALA:HB3	1:B:140:ILE:HD12	1.97	0.46
1:E:49:GLU:O	1:E:53:VAL:HG23	2.15	0.46
1:A:10:PHE:O	1:A:24:THR:HA	2.16	0.46
1:E:59:GLU:O	1:E:62:LYS:HB3	2.16	0.46
1:C:43:THR:OG1	1:C:45:VAL:HG22	2.16	0.46
1:E:36:PHE:CZ	1:E:49:GLU:HB3	2.47	0.45
1:E:140:ILE:HD12	1:E:140:ILE:H	1.81	0.45
1:D:46:CYS:C	1:D:48:THR:H	2.19	0.45
1:C:31:ARG:NH2	1:C:64:GLY:HA2	2.32	0.45
1:E:111:ARG:HD2	1:E:114:GLU:OE1	2.17	0.45
1:D:136:THR:HB	1:D:140:ILE:HB	1.98	0.44
1:D:46:CYS:O	1:D:48:THR:HG22	2.17	0.44
1:E:82:HIS:CD2	1:E:88:ILE:O	2.70	0.44
1:A:168:LYS:C	1:A:168:LYS:HE3	2.38	0.44
1:E:63:LEU:HD21	1:E:152:LYS:HD3	1.99	0.44
1:A:105:ARG:NH1	3:A:1032:HOH:O	2.39	0.44
1:A:109:ASN:HD21	1:A:118:ASP:HB2	1.82	0.44
1:A:125:ASP:HB2	1:A:126:PRO:CD	2.48	0.44
1:D:24:THR:OG1	1:D:26:LYS:HB2	2.17	0.44
1:E:125:ASP:HB2	1:E:126:PRO:HD2	1.98	0.44
1:A:16:LYS:HG2	1:A:94:ALA:CA	2.48	0.44
1:E:119:ARG:NE	1:E:138:GLU:HA	2.32	0.44
1:E:136:THR:HG21	1:E:140:ILE:O	2.18	0.43
1:A:58:GLU:HA	1:A:58:GLU:OE1	2.19	0.43
1:E:38:TYR:CZ	1:E:71:SER:HB2	2.53	0.43
1:B:136:THR:CG2	1:B:140:ILE:O	2.66	0.43
1:D:85:SER:HB3	1:D:88:ILE:HB	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:109:ASN:ND2	1:E:118:ASP:HB2	2.34	0.43
1:E:47:PRO:HB3	1:E:88:ILE:HD11	2.01	0.43
1:B:46:CYS:HA	1:B:47:PRO:HD3	1.76	0.42
1:D:165:CYS:HA	1:D:166:PRO:HD3	1.93	0.42
1:E:153:ALA:O	1:E:157:VAL:HG23	2.19	0.42
1:E:15:PHE:HE1	1:E:79:LYS:HE3	1.84	0.42
1:E:26:LYS:HD2	1:E:26:LYS:H	1.85	0.42
1:C:20:PHE:CZ	1:C:79:LYS:HD2	2.54	0.42
1:A:41:ASP:OD2	1:A:78:HIS:ND1	2.40	0.42
1:C:144:ALA:O	1:C:147:LEU:HB3	2.20	0.42
1:D:56:HIS:CD2	1:D:148:LEU:HD12	2.54	0.42
1:E:144:ALA:O	1:E:147:LEU:HB3	2.19	0.42
1:B:143:ASP:OD1	1:B:145:SER:HB2	2.20	0.42
1:D:56:HIS:HD2	1:D:148:LEU:HD12	1.84	0.42
1:E:15:PHE:CD1	1:E:79:LYS:HB2	2.55	0.42
1:E:56:HIS:CD2	1:E:148:LEU:HD12	2.53	0.41
1:D:159:ALA:C	1:D:161:PRO:HD3	2.41	0.41
1:A:120:ALA:HA	1:A:134:GLU:O	2.21	0.41
1:E:39:PRO:HG3	1:E:111:ARG:NH1	2.32	0.41
1:D:41:ASP:OD2	1:D:78:HIS:ND1	2.43	0.41
1:E:25:GLU:HG2	1:E:26:LYS:N	2.34	0.41
1:E:46:CYS:C	1:E:48:THR:H	2.24	0.41
1:D:119:ARG:NH1	1:D:136:THR:OG1	2.53	0.41
1:A:149:ARG:NH2	1:B:141:GLY:O	2.53	0.40
1:D:13:GLN:HE22	1:D:22:GLU:HB2	1.87	0.40
1:C:159:ALA:C	1:C:160:HIS:ND1	2.75	0.40
1:A:157:VAL:HG21	1:B:139:GLY:HA3	2.03	0.40
1:E:149:ARG:HH22	1:E:150:LYS:CE	2.34	0.40
1:E:57:TYR:CE1	1:E:93:TYR:HB3	2.56	0.40
1:B:38:TYR:CZ	1:B:71:SER:HB3	2.57	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	166/186 (89%)	158 (95%)	8 (5%)	0	100	100
1	B	163/186 (88%)	160 (98%)	3 (2%)	0	100	100
1	C	161/186 (87%)	154 (96%)	6 (4%)	1 (1%)	25	43
1	D	164/186 (88%)	159 (97%)	5 (3%)	0	100	100
1	E	163/186 (88%)	146 (90%)	14 (9%)	3 (2%)	8	14
All	All	817/930 (88%)	777 (95%)	36 (4%)	4 (0%)	29	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	162	GLY
1	C	162	GLY
1	E	161	PRO
1	E	164	VAL

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	140/154 (91%)	136 (97%)	4 (3%)	42	69
1	B	138/154 (90%)	134 (97%)	4 (3%)	42	69
1	C	136/154 (88%)	133 (98%)	3 (2%)	52	77
1	D	139/154 (90%)	135 (97%)	4 (3%)	42	69
1	E	138/154 (90%)	129 (94%)	9 (6%)	17	33
All	All	691/770 (90%)	667 (96%)	24 (4%)	36	62

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU

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Mol	Chain	Res	Type
1	A	60	LEU
1	A	103	LEU
1	A	168	LYS
1	B	4	ASN
1	B	103	LEU
1	B	136	THR
1	B	138	GLU
1	C	4	ASN
1	C	103	LEU
1	C	131	GLN
1	D	71	SER
1	D	103	LEU
1	D	148	LEU
1	D	163	GLU
1	E	2	LEU
1	E	24	THR
1	E	27	ASP
1	E	43	THR
1	E	87	THR
1	E	103	LEU
1	E	136	THR
1	E	155	GLN
1	E	161	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	56	HIS
1	A	109	ASN
1	A	127	GLN
1	A	131	GLN
1	A	155	GLN
1	B	109	ASN
1	B	155	GLN
1	C	56	HIS
1	C	109	ASN
1	C	127	GLN
1	C	155	GLN
1	D	13	GLN
1	D	109	ASN
1	D	155	GLN

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Mol	Chain	Res	Type
1	D	160	HIS
1	E	109	ASN
1	E	131	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

9 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	C	803	-	4,4,4	0.26	0	6,6,6	0.07	0
2	SO4	B	812	-	4,4,4	0.27	0	6,6,6	0.11	0
2	SO4	B	802	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	A	811	-	4,4,4	0.25	0	6,6,6	0.08	0
2	SO4	A	801	-	4,4,4	0.26	0	6,6,6	0.06	0
2	SO4	D	814	-	4,4,4	0.27	0	6,6,6	0.06	0
2	SO4	E	805	-	4,4,4	0.26	0	6,6,6	0.08	0
2	SO4	D	804	-	4,4,4	0.27	0	6,6,6	0.05	0
2	SO4	C	813	-	4,4,4	0.25	0	6,6,6	0.05	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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	168/186 (90%)	-0.02	17 (10%) 7 6	27, 44, 76, 100	0
1	B	165/186 (88%)	-0.13	6 (3%) 42 46	27, 41, 69, 135	0
1	C	163/186 (87%)	-0.07	11 (6%) 17 18	29, 48, 78, 111	0
1	D	166/186 (89%)	-0.01	7 (4%) 36 39	34, 50, 80, 129	0
1	E	165/186 (88%)	0.39	21 (12%) 3 3	42, 70, 107, 144	0
All	All	827/930 (88%)	0.03	62 (7%) 14 14	27, 50, 92, 144	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	VAL	7.5
1	E	165	CYS	7.0
1	D	165	CYS	7.0
1	B	165	CYS	6.0
1	D	166	PRO	5.1
1	E	159	ALA	4.6
1	E	164	VAL	4.6
1	E	163	GLU	4.5
1	C	162	GLY	4.4
1	D	44	PHE	4.2
1	D	164	VAL	3.7
1	E	70	VAL	3.7
1	E	44	PHE	3.5
1	E	34	VAL	3.5
1	A	168	LYS	3.5
1	A	70	VAL	3.4
1	C	156	TYR	3.3
1	E	162	GLY	3.3
1	A	36	PHE	3.2
1	A	96	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	44	PHE	2.9
1	A	35	PHE	2.8
1	B	70	VAL	2.7
1	A	69	SER	2.7
1	E	71	SER	2.7
1	E	36	PHE	2.7
1	A	165	CYS	2.6
1	A	68	TYR	2.6
1	E	156	TYR	2.6
1	E	96	ILE	2.6
1	D	71	SER	2.6
1	E	46	CYS	2.5
1	E	86	GLU	2.5
1	A	46	CYS	2.4
1	B	71	SER	2.4
1	E	69	SER	2.4
1	C	34	VAL	2.4
1	E	35	PHE	2.4
1	C	159	ALA	2.4
1	E	160	HIS	2.4
1	A	71	SER	2.4
1	A	34	VAL	2.3
1	C	163	GLU	2.3
1	E	62	LYS	2.3
1	C	121	THR	2.3
1	A	86	GLU	2.3
1	D	34	VAL	2.3
1	A	164	VAL	2.3
1	E	58	GLU	2.3
1	A	37	PHE	2.3
1	C	70	VAL	2.2
1	B	163	GLU	2.2
1	E	37	PHE	2.2
1	C	69	SER	2.2
1	D	70	VAL	2.1
1	B	69	SER	2.1
1	C	71	SER	2.1
1	A	120	ALA	2.1
1	A	121	THR	2.0
1	C	37	PHE	2.0
1	C	160	HIS	2.0
1	E	161	PRO	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	E	805	5/5	0.62	0.49	179,179,180,181	0
2	SO4	C	803	5/5	0.67	0.52	176,176,177,177	0
2	SO4	A	801	5/5	0.71	0.38	163,164,165,166	0
2	SO4	D	814	5/5	0.73	0.33	153,154,155,156	0
2	SO4	D	804	5/5	0.77	0.34	163,163,164,165	0
2	SO4	B	802	5/5	0.81	0.40	143,144,147,147	0
2	SO4	C	813	5/5	0.82	0.34	140,140,141,142	0
2	SO4	B	812	5/5	0.91	0.21	105,106,108,113	0
2	SO4	A	811	5/5	0.92	0.25	151,151,151,151	0

6.5 Other polymers [i](#)

There are no such residues in this entry.