



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:19 AM UTC

PDB ID : 4S37 / pdb_00004s37
Title : Crystal structure of R2 pyocin membrane-piercing spike
Authors : Browning, C.B.; Leiman, P.G.; Shneider, M.M.
Deposited on : 2015-01-26
Resolution : 2.20 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

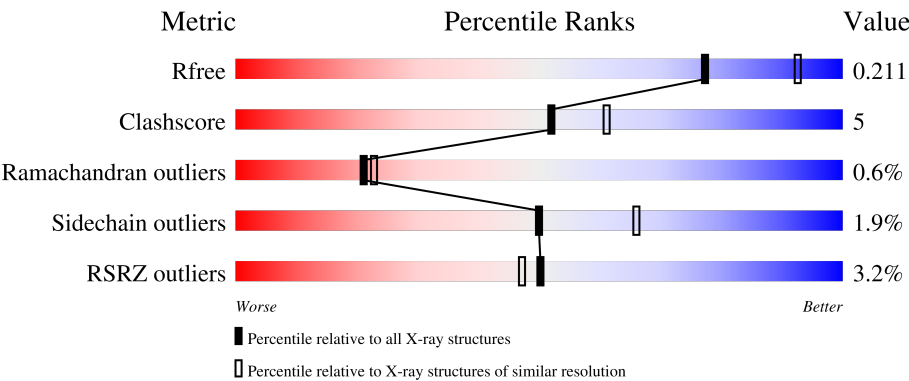
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	185	% 82%9%9%
1	B	185	2% 84%9%6%
1	C	185	2% 79%13%7%
1	D	185	% 80%13%6%
1	E	185	% 81%12%7%

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Mol	Chain	Length	Quality of chain
1	F	185	
1	G	185	
1	H	185	
1	I	185	
1	J	185	
1	K	185	
1	L	185	
1	M	185	
1	N	185	
1	O	185	
1	P	185	
1	Q	185	
1	R	185	

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	B	202	-	-	X	-
2	SO4	D	202	-	-	X	-
2	SO4	F	202	-	-	X	-
2	SO4	M	201	-	-	X	-
2	SO4	M	202	-	-	X	-
3	CL	G	205	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 26421 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 Phage baseplate protein.

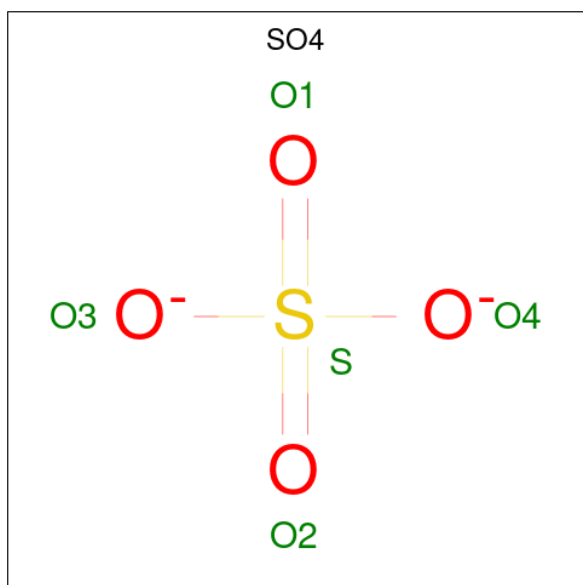
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	6	0
			1287	806	233	243	5			
1	B	174	Total	C	N	O	S	0	9	0
			1347	844	244	254	5			
1	C	172	Total	C	N	O	S	0	9	0
			1346	838	248	255	5			
1	D	173	Total	C	N	O	S	0	11	0
			1360	849	245	260	6			
1	E	172	Total	C	N	O	S	0	10	0
			1343	837	243	258	5			
1	F	175	Total	C	N	O	S	0	7	0
			1348	842	243	257	6			
1	G	175	Total	C	N	O	S	0	6	0
			1342	837	242	258	5			
1	H	176	Total	C	N	O	S	7	10	0
			1374	854	247	268	5			
1	I	175	Total	C	N	O	S	0	6	0
			1340	839	241	255	5			
1	J	175	Total	C	N	O	S	0	5	0
			1331	833	238	255	5			
1	K	176	Total	C	N	O	S	0	9	0
			1368	853	246	263	6			
1	L	175	Total	C	N	O	S	0	6	0
			1338	835	239	259	5			
1	M	172	Total	C	N	O	S	8	3	0
			1293	809	233	246	5			
1	N	171	Total	C	N	O	S	0	7	0
			1322	822	245	250	5			
1	O	174	Total	C	N	O	S	0	2	0
			1294	809	231	249	5			
1	P	166	Total	C	N	O	S	0	0	0
			1217	765	216	232	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	166	Total	C	N	O	S	8	2	0
			1231	774	218	236	3			
1	R	169	Total	C	N	O	S	0	3	0
			1269	793	231	241	4			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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	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	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		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		
2	M	1	Total	O	S	0	0
			5	4	1		
2	M	1	Total	O	S	0	0
			5	4	1		
2	M	1	Total	O	S	0	0
			5	4	1		
2	N	1	Total	O	S	0	0
			5	4	1		
2	O	1	Total	O	S	0	0
			5	4	1		

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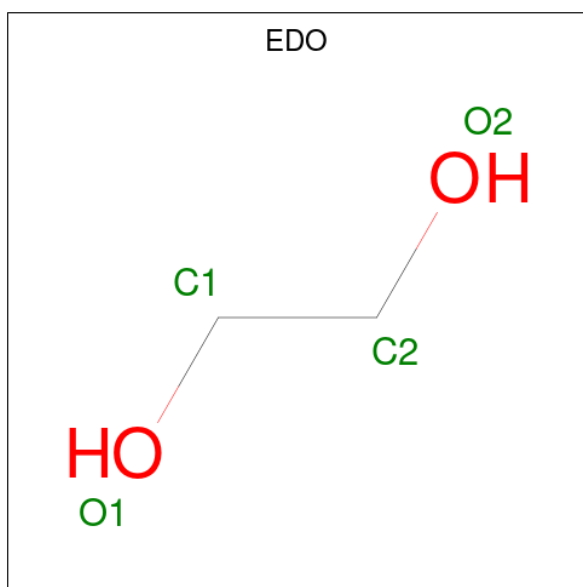
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	O	1	Total	O	S	0	0
			5	4	1		
2	O	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		
3	D	2	Total	Cl	0	0
			2	2		
3	E	2	Total	Cl	0	0
			2	2		
3	F	1	Total	Cl	0	0
			1	1		
3	G	2	Total	Cl	0	0
			2	2		
3	H	2	Total	Cl	0	0
			2	2		
3	I	2	Total	Cl	0	0
			2	2		
3	J	2	Total	Cl	0	0
			2	2		
3	K	2	Total	Cl	0	0
			2	2		
3	L	1	Total	Cl	0	0
			1	1		
3	O	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	K	1	Total C O 4 2 2	0	0
4	L	1	Total C O 4 2 2	0	0
4	N	1	Total C O 4 2 2	0	0

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

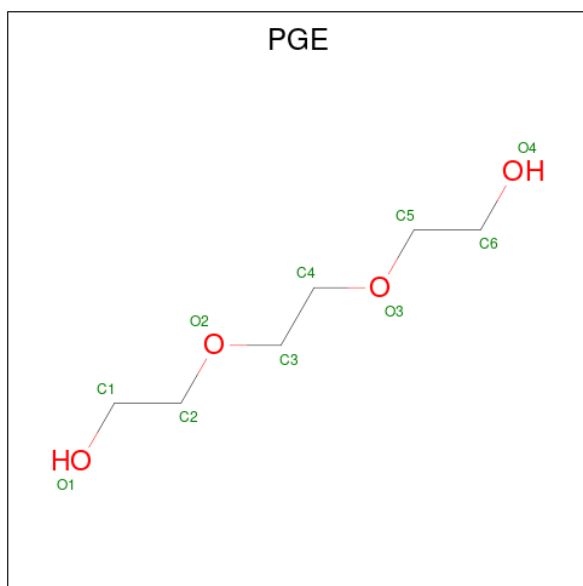
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Ca 1 1	0	0
5	D	1	Total Ca 1 1	0	0
5	F	1	Total Ca 1 1	0	0
5	H	1	Total Ca 1 1	0	0
5	J	1	Total Ca 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	O	1	Total	Ca	0	0
			1	1		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	151	Total	O	0	0
			151	151		
7	B	166	Total	O	0	0
			166	166		
7	C	133	Total	O	0	0
			133	133		
7	D	173	Total	O	0	0
			173	173		
7	E	147	Total	O	0	0
			147	147		
7	F	149	Total	O	0	0
			149	149		
7	G	174	Total	O	0	0
			174	174		

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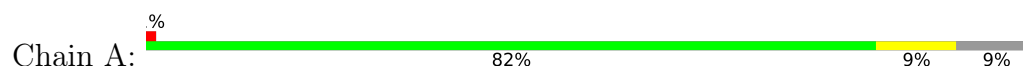
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	H	147	Total 147	O 147	0	0
7	I	138	Total 138	O 138	0	0
7	J	159	Total 159	O 159	0	0
7	K	140	Total 140	O 140	0	0
7	L	149	Total 149	O 149	0	0
7	M	105	Total 105	O 105	0	0
7	N	111	Total 111	O 111	0	0
7	O	106	Total 106	O 106	0	0
7	P	100	Total 100	O 100	0	0
7	Q	90	Total 90	O 90	0	0
7	R	95	Total 95	O 95	0	0

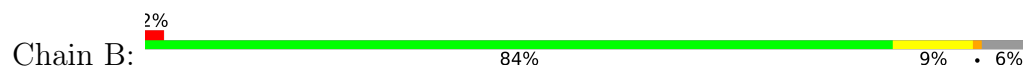
3 Residue-property plots [i](#)

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.

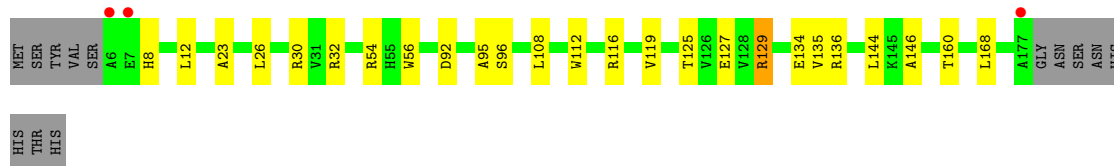
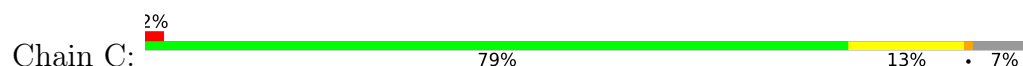
- Molecule 1: Phage baseplate protein



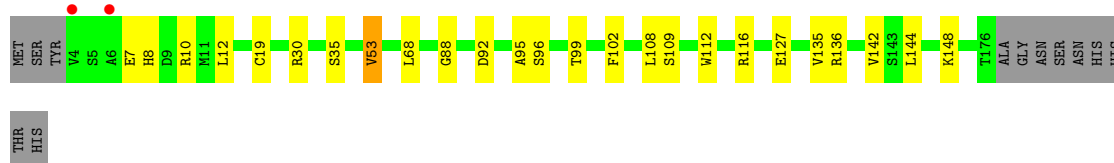
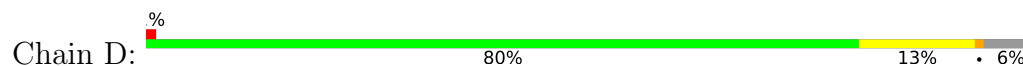
- Molecule 1: Phage baseplate protein



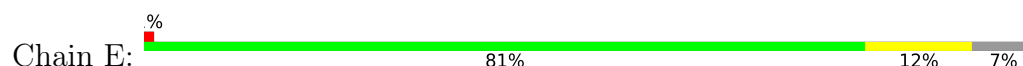
- Molecule 1: Phage baseplate protein



- Molecule 1: Phage baseplate protein

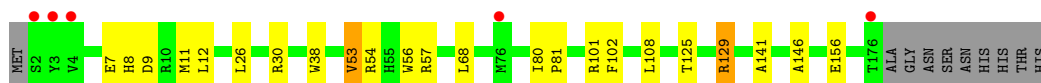
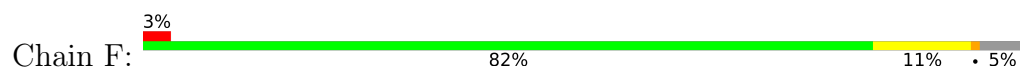


- Molecule 1: Phage baseplate protein

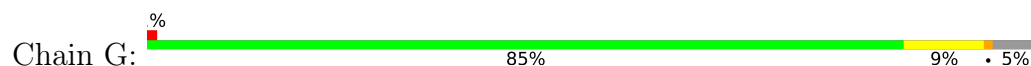




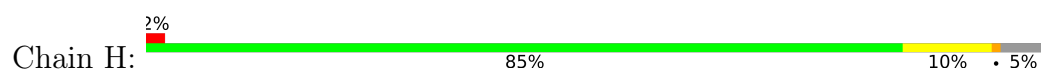
- Molecule 1: Phage baseplate protein



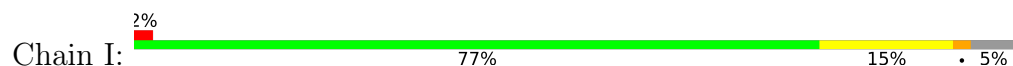
- Molecule 1: Phage baseplate protein



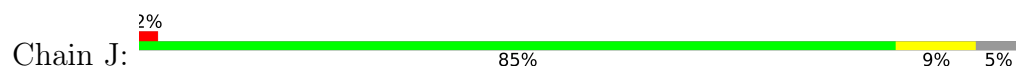
- Molecule 1: Phage baseplate protein



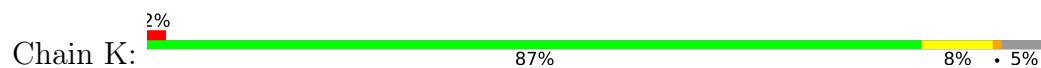
- Molecule 1: Phage baseplate protein



- Molecule 1: Phage baseplate protein

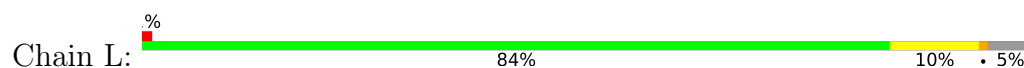


- Molecule 1: Phage baseplate protein

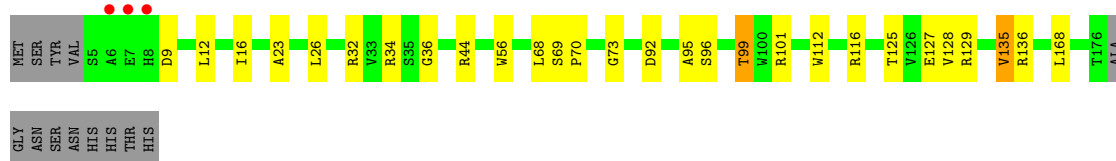
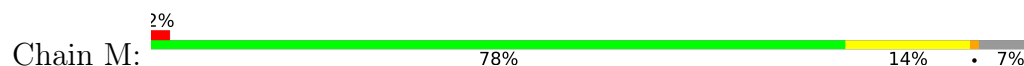




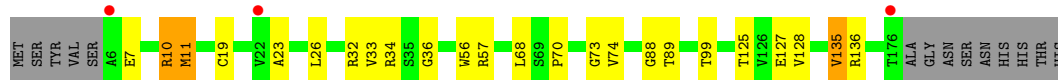
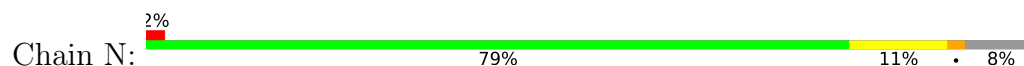
- Molecule 1: Phage baseplate protein



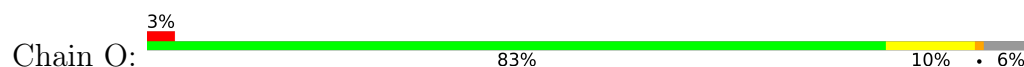
- Molecule 1: Phage baseplate protein



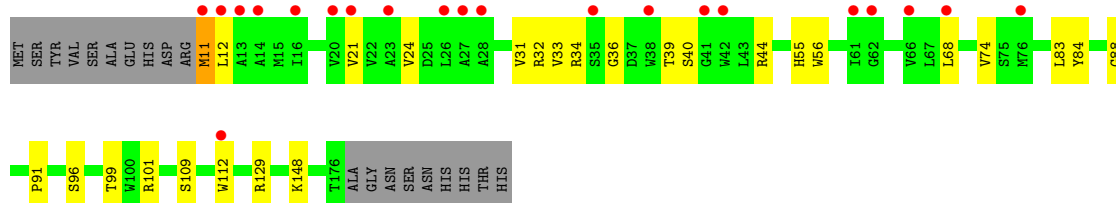
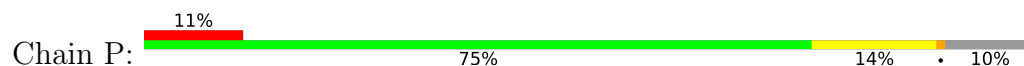
- Molecule 1: Phage baseplate protein



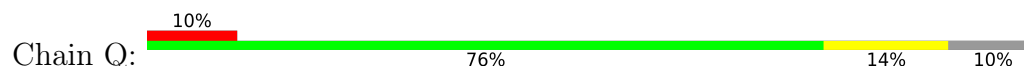
- Molecule 1: Phage baseplate protein

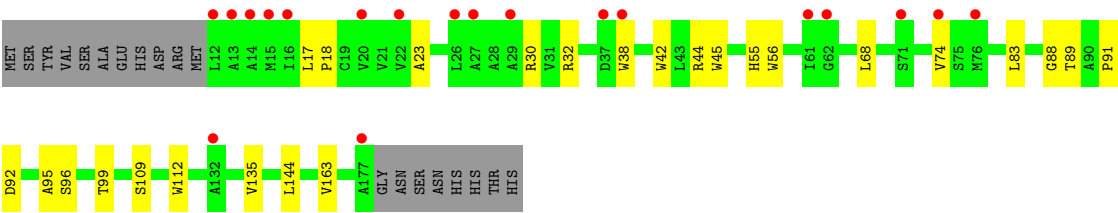


- Molecule 1: Phage baseplate protein

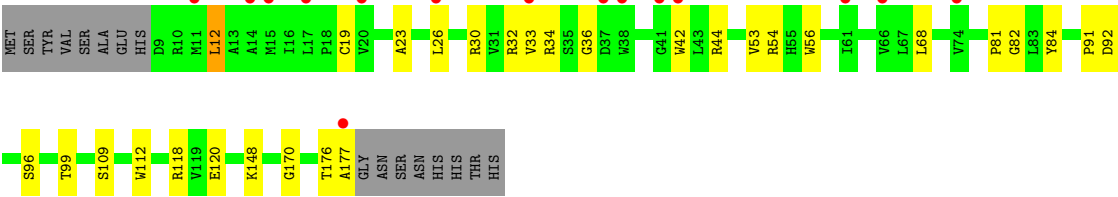


- Molecule 1: Phage baseplate protein





● Molecule 1: Phage baseplate protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.94Å 137.47Å 164.76Å 90.00° 106.56° 90.00°	Depositor
Resolution (Å)	49.19 – 2.20 49.19 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.19-2.20) 99.2 (49.19-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.163 , 0.208 0.166 , 0.211	Depositor DCC
R_{free} test set	10160 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 62.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.077 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	26421	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% 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, CL, EDO, PGE, CA

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	0/1314	0.73	0/1788
1	B	0.33	0/1384	0.72	0/1883
1	C	0.34	0/1374	0.74	0/1868
1	D	0.33	0/1388	0.71	0/1889
1	E	0.34	0/1371	0.73	0/1866
1	F	0.33	0/1377	0.71	1/1874 (0.1%)
1	G	0.35	0/1371	0.71	0/1866
1	H	0.40	0/1403	0.73	0/1910
1	I	0.35	0/1372	0.75	1/1867 (0.1%)
1	J	0.34	0/1360	0.74	0/1852
1	K	0.33	0/1397	0.73	1/1902 (0.1%)
1	L	0.35	0/1367	0.76	0/1861
1	M	0.30	0/1321	0.72	0/1798
1	N	0.29	0/1350	0.72	1/1835 (0.1%)
1	O	0.30	0/1322	0.76	0/1801
1	P	0.29	0/1244	0.72	0/1696
1	Q	0.29	0/1258	0.73	0/1716
1	R	0.30	0/1296	0.72	0/1764
All	All	0.33	0/24269	0.73	4/33036 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	53	VAL	CB-CA-C	-5.33	103.47	110.98
1	F	141	ALA	N-CA-C	5.05	116.20	108.52
1	I	53	VAL	CB-CA-C	-5.03	104.33	111.21
1	N	99	THR	N-CA-C	5.02	116.57	108.34

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1287	0	1276	12	0
1	B	1347	0	1341	17	0
1	C	1346	0	1324	19	0
1	D	1360	0	1338	23	0
1	E	1343	0	1316	20	0
1	F	1348	0	1324	20	0
1	G	1342	0	1312	22	0
1	H	1374	0	1334	17	0
1	I	1340	0	1320	25	0
1	J	1331	0	1307	12	0
1	K	1368	0	1337	10	0
1	L	1338	0	1305	15	0
1	M	1293	0	1274	22	0
1	N	1322	0	1299	16	0
1	O	1294	0	1268	15	0
1	P	1217	0	1204	20	0
1	Q	1231	0	1215	19	0
1	R	1269	0	1254	25	0
2	A	5	0	0	0	0
2	B	15	0	0	5	0
2	C	15	0	0	0	0
2	D	10	0	0	3	0
2	E	5	0	0	0	0
2	F	10	0	0	3	0
2	G	10	0	0	1	0
2	H	15	0	0	0	0
2	I	5	0	0	0	0
2	J	15	0	0	0	0
2	K	10	0	0	0	0
2	L	10	0	0	0	0
2	M	15	0	0	4	0
2	N	5	0	0	0	0
2	O	15	0	0	1	0
2	P	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	R	10	0	0	0	0
3	A	1	0	0	0	0
3	C	1	0	0	1	0
3	D	2	0	0	1	0
3	E	2	0	0	0	0
3	F	1	0	0	1	0
3	G	2	0	0	3	0
3	H	2	0	0	1	0
3	I	2	0	0	0	0
3	J	2	0	0	1	0
3	K	2	0	0	1	0
3	L	1	0	0	0	0
3	O	1	0	0	1	0
4	B	4	0	6	0	0
4	D	4	0	6	0	0
4	F	4	0	6	1	0
4	G	4	0	6	0	0
4	K	4	0	6	0	0
4	L	4	0	6	0	0
4	N	4	0	6	0	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
5	F	1	0	0	0	0
5	H	1	0	0	0	0
5	J	1	0	0	0	0
5	O	1	0	0	0	0
6	H	10	0	14	0	0
7	A	151	0	0	2	1
7	B	166	0	0	3	0
7	C	133	0	0	1	1
7	D	173	0	0	3	0
7	E	147	0	0	1	0
7	F	149	0	0	2	0
7	G	174	0	0	2	0
7	H	147	0	0	1	1
7	I	138	0	0	3	0
7	J	159	0	0	3	0
7	K	140	0	0	1	0
7	L	149	0	0	3	0
7	M	105	0	0	0	0
7	N	111	0	0	0	0
7	O	106	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	P	100	0	0	3	0
7	Q	90	0	0	1	1
7	R	95	0	0	0	0
All	All	26421	0	23404	242	2

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 5.

All (242) 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:30[B]:ARG:NH2	3:C:204:CL:CL	2.47	0.84
1:F:30:ARG:NE	3:F:205:CL:CL	2.46	0.81
1:G:30:ARG:NE	3:G:204:CL:CL	2.49	0.81
1:H:30:ARG:NE	3:H:207:CL:CL	2.52	0.80
1:D:136[B]:ARG:NH2	2:D:202:SO4:S	2.56	0.79
3:G:205:CL:CL	7:J:446:HOH:O	2.38	0.78
3:G:205:CL:CL	7:J:447:HOH:O	2.42	0.75
1:I:3:TYR:N	7:I:414:HOH:O	2.23	0.71
1:B:134:GLU:OE2	1:B:136[B]:ARG:NH1	2.26	0.68
1:G:136[B]:ARG:NH2	7:G:411:HOH:O	2.27	0.68
1:Q:83:LEU:HD13	1:R:44:ARG:HH21	1.59	0.68
1:B:32:ARG:NH1	7:B:413:HOH:O	2.28	0.67
1:B:136[B]:ARG:NH2	2:B:202:SO4:O3	2.28	0.67
1:M:128:VAL:HB	1:M:135:VAL:HG22	1.76	0.67
1:D:148:LYS:NZ	1:E:156:GLU:OE2	2.28	0.66
1:L:127:GLU:OE2	7:L:431:HOH:O	2.13	0.66
1:A:152:GLU:OE2	7:A:434:HOH:O	2.13	0.66
1:B:136[B]:ARG:NH2	2:B:202:SO4:S	2.65	0.65
1:N:128:VAL:HB	1:N:135:VAL:HG13	1.78	0.65
1:M:12:LEU:HD21	1:N:11:MET:HG2	1.79	0.65
1:J:8:HIS:NE2	1:L:9:ASP:OD1	2.28	0.64
1:K:89:THR:O	1:L:54:ARG:HD2	1.97	0.64
1:L:101:ARG:NH1	7:L:444:HOH:O	2.15	0.64
1:O:30:ARG:NE	3:O:205:CL:CL	2.59	0.63
1:D:30:ARG:NE	3:D:206:CL:CL	2.69	0.63
1:G:54:ARG:HD2	1:I:89:THR:O	1.99	0.62
1:D:127:GLU:OE1	1:D:136[B]:ARG:NH1	2.32	0.62
1:I:91:PRO:O	1:I:101[A]:ARG:NH2	2.33	0.62
1:A:108[B]:LEU:HD23	1:A:119:VAL:HG22	1.80	0.62
1:M:125:THR:HG21	1:M:136[A]:ARG:HH11	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:HIS:NE2	1:F:9:ASP:OD1	2.27	0.61
1:N:34:ARG:NH1	1:N:36:GLY:O	2.33	0.61
1:I:7:GLU:HG3	1:I:10:ARG:HH12	1.66	0.60
1:B:136[B]:ARG:NH2	2:B:202:SO4:O1	2.31	0.60
1:N:89:THR:O	1:O:54:ARG:HD2	2.02	0.60
1:I:30[A]:ARG:NH1	1:I:42:TRP:O	2.34	0.60
1:L:7:GLU:HG3	1:L:10:ARG:HH21	1.66	0.59
1:P:91:PRO:O	1:P:101:ARG:NH2	2.35	0.59
1:A:56:TRP:CD1	1:C:92:ASP:HB3	2.39	0.58
1:B:23:ALA:HB3	1:B:32:ARG:HB2	1.84	0.57
1:P:129:ARG:HB3	1:R:120:GLU:HG2	1.86	0.57
1:K:116:ARG:HD3	1:K:118:ARG:NE	2.20	0.57
1:D:136[B]:ARG:NH2	2:D:202:SO4:O4	2.37	0.57
1:R:30:ARG:NH2	1:R:42:TRP:O	2.38	0.56
1:N:7:GLU:HA	1:N:10:ARG:HD3	1.87	0.56
1:P:12:LEU:HD11	1:R:12:LEU:HD21	1.87	0.56
1:F:156:GLU:OE2	7:F:438:HOH:O	2.18	0.56
1:G:53:VAL:HG13	1:I:57:ARG:HD3	1.87	0.56
1:Q:30:ARG:NH1	1:Q:42:TRP:O	2.37	0.56
1:L:136[B]:ARG:NH1	7:L:414:HOH:O	2.35	0.56
1:E:102:PHE:HE1	1:E:108[B]:LEU:HD23	1.70	0.56
1:G:92:ASP:HB3	1:H:56:TRP:CD1	2.41	0.56
1:H:91:PRO:O	1:H:101[A]:ARG:NH1	2.40	0.55
1:J:56:TRP:CD1	1:L:92:ASP:HB3	2.41	0.55
1:D:95:ALA:HB2	1:E:26:LEU:HD11	1.89	0.55
1:Q:45:TRP:O	7:Q:215:HOH:O	2.18	0.55
1:R:34:ARG:NH1	1:R:36:GLY:O	2.40	0.55
1:P:44:ARG:NH1	1:R:84:TYR:O	2.40	0.54
1:M:92:ASP:HB3	1:N:56:TRP:CD1	2.42	0.54
1:Q:68:LEU:HD21	1:R:68:LEU:HD13	1.89	0.54
1:C:129:ARG:HH11	1:C:134:GLU:HB2	1.72	0.54
1:P:56:TRP:HB2	1:R:91:PRO:HG2	1.89	0.54
1:R:23:ALA:HB3	1:R:32:ARG:HB2	1.88	0.54
1:I:99:THR:HG23	1:I:109[A]:SER:HB3	1.90	0.54
1:F:102:PHE:HE2	1:F:108[B]:LEU:HD23	1.73	0.54
1:P:99:THR:HG23	1:P:109:SER:HB3	1.89	0.54
1:Q:23:ALA:HB3	1:Q:32:ARG:HB3	1.89	0.53
1:G:129:ARG:NH1	2:G:202:SO4:O3	2.41	0.53
1:C:96:SER:HA	1:C:112:TRP:CZ2	2.43	0.53
1:B:92:ASP:HB3	1:C:56:TRP:CD1	2.44	0.53
1:E:6:ALA:N	7:E:443:HOH:O	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:96:SER:HA	1:O:112:TRP:CZ2	2.43	0.53
1:D:53:VAL:HG13	1:F:57:ARG:HD3	1.91	0.52
1:H:8:HIS:O	1:H:12:LEU:HB2	2.09	0.52
1:N:127:GLU:OE2	1:N:136[B]:ARG:NH1	2.42	0.52
1:G:129:ARG:HD2	1:G:133:SER:O	2.09	0.52
1:I:8:HIS:O	1:I:12:LEU:HB2	2.10	0.52
1:M:26:LEU:HD11	1:O:95:ALA:HB2	1.91	0.52
1:G:68:LEU:HD21	1:H:68:LEU:HD13	1.91	0.52
1:G:125[B]:THR:HG23	1:I:116:ARG:HG3	1.92	0.51
1:A:57:ARG:HD3	1:B:53[A]:VAL:HG13	1.92	0.51
1:M:101:ARG:NE	2:M:202:SO4:O2	2.37	0.51
1:Q:91:PRO:HG2	1:R:56:TRP:HB2	1.91	0.51
1:J:68:LEU:HD13	1:L:68:LEU:HD21	1.93	0.51
1:P:68:LEU:HD13	1:R:68:LEU:HD21	1.93	0.51
1:A:136[A]:ARG:NH1	7:A:422:HOH:O	2.40	0.51
1:P:32:ARG:NH2	1:P:40:SER:O	2.44	0.51
1:G:116:ARG:HG3	1:H:125[B]:THR:HG23	1.92	0.51
1:D:136[A]:ARG:NH1	7:D:409:HOH:O	2.44	0.50
1:D:136[B]:ARG:NH2	2:D:202:SO4:O2	2.37	0.50
1:P:96:SER:HA	1:P:112:TRP:CZ2	2.46	0.50
1:H:96:SER:HA	1:H:112:TRP:CZ2	2.47	0.50
1:M:95:ALA:HB2	1:N:26:LEU:HD11	1.93	0.50
1:M:96:SER:HA	1:M:112:TRP:CZ2	2.46	0.50
1:Q:92:ASP:HB3	1:R:56:TRP:CG	2.46	0.50
1:P:11:MET:N	7:P:393:HOH:O	2.43	0.50
1:P:148:LYS:NZ	7:P:368:HOH:O	2.44	0.50
1:L:96:SER:HA	1:L:112:TRP:CZ2	2.46	0.50
1:C:125[A]:THR:HG21	1:C:136[A]:ARG:HH21	1.75	0.50
1:I:96:SER:HA	1:I:112:TRP:CZ2	2.47	0.50
1:D:102:PHE:HE2	1:D:108[B]:LEU:HD23	1.77	0.49
1:A:102:PHE:HE1	1:A:108[B]:LEU:HD12	1.78	0.49
1:O:7:GLU:O	1:O:11:MET:HB2	2.13	0.49
1:Q:96:SER:HA	1:Q:112:TRP:CZ2	2.46	0.49
1:G:96:SER:HA	1:G:112:TRP:CZ2	2.47	0.49
1:R:118:ARG:HG2	1:R:120:GLU:HG3	1.95	0.49
1:M:101:ARG:NH2	2:M:202:SO4:O3	2.46	0.49
1:D:135:VAL:HG22	1:D:144:LEU:HD22	1.95	0.49
1:P:34:ARG:HB3	1:P:39:THR:HG22	1.95	0.49
1:H:68:LEU:HD21	1:I:68:LEU:HD13	1.94	0.48
1:P:56:TRP:CG	1:R:92:ASP:HB3	2.48	0.48
1:M:68:LEU:HD21	1:N:68:LEU:HD13	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:23:ALA:HB3	1:N:32[B]:ARG:HB2	1.96	0.48
1:B:136[A]:ARG:NH2	2:B:202:SO4:O1	2.34	0.48
1:P:91:PRO:HG2	1:Q:56:TRP:HB2	1.95	0.48
1:H:92:ASP:HB3	1:I:56:TRP:CD1	2.49	0.48
1:R:99:THR:HG23	1:R:109[A]:SER:HB2	1.95	0.48
1:D:68:LEU:HD13	1:F:68:LEU:HD21	1.96	0.48
1:J:56:TRP:CG	1:L:92:ASP:HB3	2.49	0.48
1:M:136[B]:ARG:NH2	2:M:201:SO4:O4	2.47	0.48
1:M:116:ARG:HG3	1:N:125:THR:HG23	1.95	0.48
1:J:26:LEU:HD11	1:L:95:ALA:HB2	1.96	0.47
1:C:135:VAL:HG22	1:C:144:LEU:HD22	1.97	0.47
1:Q:95:ALA:HB2	1:R:26:LEU:HD11	1.96	0.47
1:B:54:ARG:NH2	2:B:204:SO4:O1	2.33	0.47
1:Q:135:VAL:HG22	1:Q:144:LEU:HD22	1.96	0.47
1:E:57:ARG:HD3	1:F:53:VAL:HG13	1.96	0.47
1:H:95:ALA:HB2	1:I:26:LEU:HD11	1.96	0.47
1:J:91:PRO:HG2	1:K:56:TRP:HB2	1.97	0.47
3:J:205:CL:CL	7:J:434:HOH:O	2.58	0.47
1:K:10:ARG:NH2	7:K:411:HOH:O	2.47	0.47
1:D:116:ARG:HG3	1:E:125[A]:THR:HG23	1.96	0.47
1:E:19[B]:CYS:SG	1:E:35:SER:HB2	2.55	0.47
1:B:136[A]:ARG:NH1	7:B:436:HOH:O	2.34	0.46
1:Q:99:THR:HG23	1:Q:109:SER:HB3	1.96	0.46
1:M:68:LEU:HD13	1:O:68:LEU:HD21	1.96	0.46
1:D:7:GLU:OE2	1:D:10:ARG:NH1	2.49	0.46
1:E:116:ARG:HG3	1:F:125:THR:HG23	1.98	0.46
1:I:136:ARG:NH2	7:I:392:HOH:O	2.48	0.46
1:F:54:ARG:NH2	2:F:203:SO4:O1	2.43	0.46
1:F:129:ARG:NH1	2:F:202:SO4:O3	2.48	0.46
1:C:23:ALA:HB3	1:C:32:ARG:HB2	1.97	0.46
1:D:19[B]:CYS:SG	1:D:35:SER:HB2	2.56	0.46
1:G:30:ARG:NH2	1:G:42:TRP:O	2.46	0.46
1:B:142:VAL:HG23	1:C:146:ALA:HB2	1.97	0.46
1:C:127:GLU:OE2	1:C:136[B]:ARG:NH2	2.49	0.46
1:G:56:TRP:CD1	1:I:92:ASP:HB3	2.51	0.46
1:J:96:SER:HA	1:J:112:TRP:CZ2	2.50	0.46
1:N:57:ARG:HD3	1:O:53:VAL:HG13	1.98	0.46
1:G:142:VAL:HG23	1:H:146:ALA:HB2	1.98	0.45
1:J:125[B]:THR:HG22	1:J:138:SER:HB3	1.97	0.45
1:J:135:VAL:HG22	1:J:144:LEU:HD22	1.97	0.45
1:M:34:ARG:NH1	1:M:36:GLY:O	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:201:SO4:O2	1:O:118:ARG:NH2	2.48	0.45
1:O:6:ALA:HA	1:O:9:ASP:HB2	1.97	0.45
1:I:99:THR:HG23	1:I:109[B]:SER:HB2	1.96	0.45
1:M:9:ASP:HA	1:M:12:LEU:HB3	1.98	0.45
1:D:92:ASP:HB3	1:E:56:TRP:CG	2.52	0.45
1:M:56:TRP:CD1	1:O:92:ASP:HB3	2.50	0.45
1:R:96:SER:HA	1:R:112:TRP:CZ2	2.51	0.45
1:D:96:SER:HA	1:D:112:TRP:CZ2	2.52	0.45
1:E:118:ARG:NH1	2:F:202:SO4:O4	2.49	0.45
1:R:176:THR:HA	1:R:177:ALA:HA	1.53	0.45
1:Q:92:ASP:HB3	1:R:56:TRP:CD1	2.52	0.45
1:A:76[A]:MET:SD	1:H:10:ARG:HG2	2.57	0.45
1:N:19[B]:CYS:HB3	1:N:33:VAL:HB	1.98	0.45
1:E:95:ALA:HB2	1:F:26:LEU:HD11	1.98	0.45
1:P:24:VAL:HG22	1:P:31:VAL:HG12	1.98	0.45
1:P:84:TYR:OH	1:Q:55:HIS:O	2.28	0.45
1:F:8:HIS:O	1:F:12:LEU:HB2	2.17	0.44
7:D:359:HOH:O	1:E:129:ARG:HG2	2.16	0.44
1:M:69:SER:HA	1:M:70:PRO:HD2	1.89	0.44
1:B:99:THR:HG23	1:B:109[A]:SER:HB2	2.00	0.44
1:A:15:MET:SD	1:C:12:LEU:HD22	2.57	0.44
1:J:92:ASP:HB3	1:K:56:TRP:CD1	2.52	0.44
1:G:57:ARG:HD3	1:H:53:VAL:HG13	1.99	0.44
1:J:46:HIS:CE1	1:L:66:VAL:HG12	2.52	0.44
1:R:19:CYS:HB2	1:R:33:VAL:HB	2.00	0.44
1:H:101[B]:ARG:NH2	7:H:414:HOH:O	2.27	0.44
1:F:38:TRP:CG	4:F:201:EDO:H22	2.53	0.44
1:M:127:GLU:OE1	1:M:136[B]:ARG:NH1	2.50	0.44
1:A:125:THR:HG23	1:C:116:ARG:HG3	2.00	0.43
1:I:23:ALA:HB3	1:I:32:ARG:HB2	1.99	0.43
1:B:92:ASP:HB3	1:C:56:TRP:CG	2.53	0.43
7:B:333:HOH:O	1:C:125[B]:THR:HG22	2.18	0.43
1:F:101[B]:ARG:NH2	7:F:448:HOH:O	2.39	0.43
1:M:23:ALA:HB3	1:M:32:ARG:HB2	1.99	0.43
1:Q:163:VAL:O	1:R:170:GLY:N	2.47	0.43
1:Q:89:THR:OG1	1:R:54:ARG:NH1	2.52	0.43
1:A:166:ASP:H	1:C:160:THR:HG22	1.84	0.43
1:K:12:LEU:HD11	1:L:12:LEU:HD12	2.01	0.43
1:K:34:ARG:HD3	3:K:205:CL:CL	2.56	0.43
1:H:80:ILE:HA	1:H:81:PRO:HD3	1.86	0.43
1:I:34:ARG:HD3	7:I:399:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:99[A]:THR:HG23	1:K:109[A]:SER:HB3	2.01	0.42
1:M:99:THR:HG23	1:N:56:TRP:HB3	2.00	0.42
1:O:136:ARG:NH1	2:O:202:SO4:O4	2.46	0.42
1:D:68:LEU:HD21	1:E:68:LEU:HD13	2.01	0.42
1:G:68:LEU:HD13	1:I:68:LEU:HD21	2.01	0.42
1:E:142:VAL:HG23	1:F:146:ALA:HB2	2.02	0.42
1:N:23:ALA:HB3	1:N:32[A]:ARG:HB2	2.02	0.42
1:O:23:ALA:HB3	1:O:32:ARG:HB2	2.01	0.42
1:B:89:THR:O	1:C:54[A]:ARG:HD2	2.18	0.42
1:O:30:ARG:NH2	7:O:388:HOH:O	2.51	0.42
1:E:92:ASP:HB3	1:F:56:TRP:CD1	2.54	0.42
1:F:80:ILE:HA	1:F:81:PRO:HD3	1.89	0.42
1:L:135:VAL:HG22	1:L:144:LEU:HD22	2.02	0.42
1:D:12:LEU:HG	1:E:15:MET:SD	2.60	0.42
1:F:7:GLU:O	1:F:11:MET:HB2	2.20	0.42
1:G:146:ALA:HB2	1:I:142:VAL:HG23	2.01	0.42
1:P:21:VAL:HA	1:P:33:VAL:HG12	2.01	0.42
1:P:56:TRP:CD1	1:R:92:ASP:HB3	2.55	0.42
1:Q:89:THR:O	1:R:54:ARG:HD2	2.20	0.42
1:E:92:ASP:HB3	1:F:56:TRP:CG	2.55	0.41
1:H:92:ASP:HB3	1:I:56:TRP:CG	2.54	0.41
1:B:80:ILE:HA	1:B:81:PRO:HD3	1.90	0.41
1:E:135:VAL:HG22	1:E:144:LEU:HD22	2.02	0.41
1:G:92:ASP:HB3	1:H:56:TRP:CG	2.55	0.41
1:O:7:GLU:HA	1:O:10:ARG:NH2	2.36	0.41
1:G:26:LEU:HD11	1:I:95:ALA:HB2	2.01	0.41
1:G:54:ARG:HD3	1:I:88:GLY:HA2	2.01	0.41
1:J:129:ARG:HD2	1:J:134:GLU:HB2	2.03	0.41
1:Q:17:LEU:HA	1:Q:18:PRO:HD2	1.95	0.41
1:C:108[A]:LEU:HD13	1:C:119:VAL:HG22	2.03	0.41
1:G:53:VAL:CG1	1:I:57:ARG:HD3	2.49	0.41
1:M:44:ARG:NH2	1:O:83:LEU:HB3	2.36	0.41
7:D:346:HOH:O	1:E:125[A]:THR:HG22	2.20	0.41
1:K:7:GLU:OE2	1:K:10:ARG:NH1	2.54	0.41
1:M:16:ILE:HG22	1:N:70:PRO:HG3	2.03	0.41
1:Q:38:TRP:CH2	1:Q:74:VAL:HB	2.56	0.41
1:B:95:ALA:HB2	1:C:26:LEU:HD11	2.02	0.40
1:D:53:VAL:CG1	1:F:57:ARG:HD3	2.50	0.40
1:D:99:THR:HG23	1:D:109[B]:SER:HB3	2.03	0.40
1:G:125[B]:THR:HG22	7:G:378:HOH:O	2.21	0.40
1:I:11:MET:O	1:I:15:MET:HG3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:81:PRO:HA	1:R:82:GLY:HA2	1.81	0.40
1:A:26:LEU:HD11	1:C:95:ALA:HB2	2.02	0.40
1:A:125:THR:HG22	7:C:348:HOH:O	2.21	0.40
1:K:9:ASP:HB3	1:L:8:HIS:NE2	2.37	0.40
1:P:34:ARG:NH2	1:P:36:GLY:O	2.55	0.40
1:D:142:VAL:HG23	1:E:146:ALA:HB2	2.03	0.40
1:P:55:HIS:HD2	7:P:337:HOH:O	2.04	0.40

All (2) 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)
7:H:347:HOH:O	7:Q:283:HOH:O[1_645]	2.17	0.03
7:A:354:HOH:O	7:C:387:HOH:O[2_646]	2.18	0.02

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	172/185 (93%)	167 (97%)	4 (2%)	1 (1%)	21	23
1	B	181/185 (98%)	175 (97%)	5 (3%)	1 (1%)	21	23
1	C	179/185 (97%)	172 (96%)	7 (4%)	0	100	100
1	D	182/185 (98%)	177 (97%)	4 (2%)	1 (0%)	24	27
1	E	180/185 (97%)	174 (97%)	5 (3%)	1 (1%)	21	23
1	F	180/185 (97%)	172 (96%)	8 (4%)	0	100	100
1	G	179/185 (97%)	173 (97%)	6 (3%)	0	100	100
1	H	184/185 (100%)	176 (96%)	7 (4%)	1 (0%)	24	27
1	I	179/185 (97%)	171 (96%)	6 (3%)	2 (1%)	11	10
1	J	178/185 (96%)	172 (97%)	4 (2%)	2 (1%)	11	10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	183/185 (99%)	175 (96%)	8 (4%)	0	100	100
1	L	179/185 (97%)	173 (97%)	5 (3%)	1 (1%)	21	23
1	M	173/185 (94%)	165 (95%)	7 (4%)	1 (1%)	21	23
1	N	176/185 (95%)	168 (96%)	6 (3%)	2 (1%)	11	10
1	O	174/185 (94%)	165 (95%)	7 (4%)	2 (1%)	11	10
1	P	164/185 (89%)	151 (92%)	12 (7%)	1 (1%)	21	23
1	Q	166/185 (90%)	157 (95%)	8 (5%)	1 (1%)	21	23
1	R	170/185 (92%)	163 (96%)	7 (4%)	0	100	100
All	All	3179/3330 (96%)	3046 (96%)	116 (4%)	17 (0%)	21	27

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	73	GLY
1	N	73	GLY
1	P	88	GLY
1	I	81	PRO
1	J	81	PRO
1	O	88	GLY
1	Q	88	GLY
1	D	88	GLY
1	E	88	GLY
1	J	88	GLY
1	L	88	GLY
1	N	88	GLY
1	B	88	GLY
1	I	88	GLY
1	O	81	PRO
1	A	88	GLY
1	H	88	GLY

5.3.2 Protein sidechains ⓘ

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	134/142 (94%)	130 (97%)	4 (3%)	36	49
1	B	141/142 (99%)	136 (96%)	5 (4%)	32	43
1	C	139/142 (98%)	136 (98%)	3 (2%)	45	61
1	D	143/142 (101%)	142 (99%)	1 (1%)	76	87
1	E	140/142 (99%)	140 (100%)	0	100	100
1	F	141/142 (99%)	139 (99%)	2 (1%)	59	75
1	G	140/142 (99%)	139 (99%)	1 (1%)	76	87
1	H	144/142 (101%)	142 (99%)	2 (1%)	59	75
1	I	139/142 (98%)	135 (97%)	4 (3%)	37	51
1	J	139/142 (98%)	139 (100%)	0	100	100
1	K	143/142 (101%)	139 (97%)	4 (3%)	38	52
1	L	140/142 (99%)	137 (98%)	3 (2%)	47	63
1	M	134/142 (94%)	130 (97%)	4 (3%)	36	49
1	N	137/142 (96%)	133 (97%)	4 (3%)	37	51
1	O	134/142 (94%)	132 (98%)	2 (2%)	57	73
1	P	126/142 (89%)	123 (98%)	3 (2%)	43	58
1	Q	127/142 (89%)	126 (99%)	1 (1%)	73	85
1	R	131/142 (92%)	128 (98%)	3 (2%)	44	59
All	All	2472/2556 (97%)	2426 (98%)	46 (2%)	50	66

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	53[A]	VAL
1	A	53[B]	VAL
1	A	156	GLU
1	B	4	VAL
1	B	53[A]	VAL
1	B	53[B]	VAL
1	B	74	VAL
1	B	129	ARG
1	C	8	HIS
1	C	129	ARG
1	C	168	LEU
1	D	53	VAL
1	F	53	VAL

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Mol	Chain	Res	Type
1	F	129	ARG
1	G	53	VAL
1	H	53	VAL
1	H	129	ARG
1	I	7	GLU
1	I	10	ARG
1	I	12	LEU
1	I	74	VAL
1	K	7	GLU
1	K	96	SER
1	K	120	GLU
1	K	129	ARG
1	L	12	LEU
1	L	161	LEU
1	L	168	LEU
1	M	99	THR
1	M	129	ARG
1	M	135	VAL
1	M	168	LEU
1	N	10	ARG
1	N	11	MET
1	N	74	VAL
1	N	135	VAL
1	O	10	ARG
1	O	53	VAL
1	P	11	MET
1	P	74	VAL
1	P	83	LEU
1	Q	44	ARG
1	R	12	LEU
1	R	53	VAL
1	R	148	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	64	GLN

5.3.3 RNA ⓘ

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](#)

Of 68 ligands modelled in this entry, 25 are monoatomic - leaving 43 for Mogul analysis.

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	O	201	-	4,4,4	0.25	0	6,6,6	0.06	0
2	SO4	P	201	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	D	203	-	4,4,4	0.25	0	6,6,6	0.06	0
4	EDO	L	201	-	3,3,3	0.43	0	2,2,2	0.35	0
2	SO4	M	203	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	F	202	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	B	204	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	J	201	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	G	202	-	4,4,4	0.23	0	6,6,6	0.06	0
4	EDO	K	201	-	3,3,3	0.41	0	2,2,2	0.38	0
2	SO4	C	202	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	F	203	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	G	203	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	K	202	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	M	201	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	A	201	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	O	202	-	4,4,4	0.25	0	6,6,6	0.15	0
6	PGE	H	201	-	9,9,9	0.32	0	8,8,8	0.27	0
4	EDO	F	201	-	3,3,3	0.44	0	2,2,2	0.34	0
4	EDO	G	201	-	3,3,3	0.43	0	2,2,2	0.37	0
2	SO4	M	202	-	4,4,4	0.23	0	6,6,6	0.08	0
4	EDO	N	201	-	3,3,3	0.43	0	2,2,2	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	O	203	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	J	202	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	H	203	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	J	203	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	H	204	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	R	202	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	N	202	-	4,4,4	0.23	0	6,6,6	0.06	0
4	EDO	B	201	-	3,3,3	0.43	0	2,2,2	0.38	0
2	SO4	E	201	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	I	201	-	4,4,4	0.24	0	6,6,6	0.09	0
4	EDO	D	201	-	3,3,3	0.44	0	2,2,2	0.34	0
2	SO4	D	202	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	C	201	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	B	202	-	4,4,4	0.23	0	6,6,6	0.11	0
2	SO4	K	203	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	C	203	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	L	202	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	L	203	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	R	201	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	H	202	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	B	203	-	4,4,4	0.24	0	6,6,6	0.04	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	L	201	-	-	0/1/1/1	-
4	EDO	D	201	-	-	0/1/1/1	-
6	PGE	H	201	-	-	2/7/7/7	-
4	EDO	F	201	-	-	0/1/1/1	-
4	EDO	G	201	-	-	0/1/1/1	-
4	EDO	N	201	-	-	0/1/1/1	-
4	EDO	K	201	-	-	0/1/1/1	-
4	EDO	B	201	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	H	201	PGE	C1-C2-O2-C3
6	H	201	PGE	C3-C4-O3-C5

There are no ring outliers.

10 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	202	SO4	2	0
2	B	204	SO4	1	0
2	G	202	SO4	1	0
2	F	203	SO4	1	0
2	M	201	SO4	2	0
2	O	202	SO4	1	0
4	F	201	EDO	1	0
2	M	202	SO4	2	0
2	D	202	SO4	3	0
2	B	202	SO4	4	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 ⓘ

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	168/185 (90%)	-0.48	1 (0%) 85 83	16, 37, 59, 99	6 (3%)
1	B	174/185 (94%)	-0.48	3 (1%) 69 66	14, 36, 69, 135	9 (5%)
1	C	172/185 (92%)	-0.46	3 (1%) 69 66	14, 36, 59, 130	9 (5%)
1	D	173/185 (93%)	-0.52	2 (1%) 76 74	14, 35, 70, 139	11 (6%)
1	E	172/185 (92%)	-0.37	2 (1%) 76 74	13, 37, 72, 135	10 (5%)
1	F	175/185 (94%)	-0.39	5 (2%) 53 50	14, 37, 80, 182	7 (4%)
1	G	175/185 (94%)	-0.53	1 (0%) 85 83	15, 35, 64, 102	6 (3%)
1	H	176/185 (95%)	-0.31	3 (1%) 69 66	15, 43, 66, 92	10 (5%)
1	I	175/185 (94%)	-0.40	4 (2%) 61 58	14, 38, 86, 144	6 (3%)
1	J	175/185 (94%)	-0.46	3 (1%) 69 66	15, 37, 76, 106	5 (2%)
1	K	176/185 (95%)	-0.43	3 (1%) 69 66	15, 37, 81, 119	9 (5%)
1	L	175/185 (94%)	-0.50	2 (1%) 78 76	15, 36, 68, 134	6 (3%)
1	M	171/185 (92%)	0.17	3 (1%) 67 64	21, 65, 101, 173	2 (1%)
1	N	171/185 (92%)	0.13	3 (1%) 67 64	18, 57, 103, 138	7 (4%)
1	O	174/185 (94%)	-0.00	5 (2%) 53 50	31, 52, 109, 175	2 (1%)
1	P	166/185 (89%)	0.74	21 (12%) 8 6	24, 97, 162, 196	0
1	Q	166/185 (89%)	0.64	19 (11%) 10 7	20, 87, 160, 178	2 (1%)
1	R	169/185 (91%)	0.61	15 (8%) 15 13	26, 89, 155, 190	3 (1%)
All	All	3103/3330 (93%)	-0.18	98 (3%) 50 47	13, 42, 131, 196	110 (3%)

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	6	ALA	5.7
1	Q	12	LEU	5.0
1	E	177	ALA	5.0

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Mol	Chain	Res	Type	RSRZ
1	C	177	ALA	4.8
1	E	6	ALA	4.6
1	H	178	ALA	4.3
1	O	4	VAL	4.2
1	R	14	ALA	4.2
1	J	76[A]	MET	3.7
1	D	4	VAL	3.6
1	P	12	LEU	3.6
1	Q	22	VAL	3.5
1	J	176	THR	3.4
1	Q	177	ALA	3.4
1	H	76[A]	MET	3.4
1	F	3	TYR	3.3
1	I	3	TYR	3.3
1	P	26	LEU	3.2
1	K	177	ALA	3.2
1	O	76[A]	MET	3.2
1	P	27	ALA	3.1
1	M	8	HIS	3.1
1	O	14	ALA	3.1
1	L	4	VAL	3.1
1	I	4	VAL	2.9
1	O	12	LEU	2.9
1	B	4	VAL	2.9
1	Q	132	ALA	2.9
1	O	177	ALA	2.8
1	B	177	ALA	2.8
1	R	177	ALA	2.8
1	R	15	MET	2.8
1	K	4	VAL	2.8
1	Q	62	GLY	2.8
1	P	68	LEU	2.8
1	N	6	ALA	2.8
1	D	6	ALA	2.7
1	F	2	SER	2.7
1	K	2	SER	2.7
1	Q	61	ILE	2.7
1	P	38	TRP	2.7
1	P	112	TRP	2.7
1	P	42	TRP	2.6
1	M	6	ALA	2.6
1	Q	27	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	2	SER	2.6
1	I	177	ALA	2.6
1	Q	29	ALA	2.6
1	P	16	ILE	2.6
1	P	35	SER	2.6
1	R	66	VAL	2.5
1	Q	15	MET	2.5
1	F	176	THR	2.5
1	F	4	VAL	2.5
1	L	2	SER	2.5
1	B	8	HIS	2.5
1	N	176	THR	2.4
1	P	28	ALA	2.4
1	P	66	VAL	2.4
1	P	41	GLY	2.4
1	P	62	GLY	2.4
1	P	14	ALA	2.4
1	Q	13	ALA	2.4
1	A	9	ASP	2.3
1	Q	16	ILE	2.3
1	R	33	VAL	2.3
1	R	11	MET	2.3
1	R	61	ILE	2.3
1	Q	20	VAL	2.3
1	Q	26	LEU	2.3
1	F	76[A]	MET	2.3
1	C	7	GLU	2.3
1	P	21	VAL	2.3
1	R	74	VAL	2.3
1	P	76	MET	2.3
1	Q	38	TRP	2.3
1	P	20	VAL	2.3
1	Q	76	MET	2.2
1	M	7	GLU	2.2
1	P	23	ALA	2.2
1	Q	71	SER	2.2
1	R	26	LEU	2.2
1	P	11	MET	2.2
1	N	22	VAL	2.2
1	Q	37	ASP	2.2
1	J	2	SER	2.2
1	Q	74	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	R	38	TRP	2.1
1	Q	14	ALA	2.1
1	I	76[A]	MET	2.1
1	R	17	LEU	2.1
1	R	41	GLY	2.1
1	H	12	LEU	2.1
1	P	61	ILE	2.1
1	R	37	ASP	2.1
1	R	20	VAL	2.1
1	P	13	ALA	2.0
1	R	42	TRP	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 oligosaccharides 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	R	201	5/5	0.32	0.14	198,200,201,203	0
2	SO4	H	202	5/5	0.46	0.12	148,150,152,152	0
2	SO4	J	202	5/5	0.52	0.13	174,174,175,176	0
2	SO4	P	201	5/5	0.62	0.11	146,148,151,152	0
2	SO4	F	202	5/5	0.68	0.15	171,172,172,173	0
2	SO4	H	203	5/5	0.69	0.14	137,137,138,140	0
2	SO4	O	201	5/5	0.70	0.16	133,140,141,145	0
2	SO4	M	203	5/5	0.71	0.11	144,144,145,146	0
2	SO4	I	201	5/5	0.72	0.12	132,134,136,138	0
2	SO4	O	203	5/5	0.72	0.11	129,129,131,132	0
2	SO4	K	203	5/5	0.73	0.12	126,127,129,129	0
2	SO4	N	202	5/5	0.74	0.15	176,178,179,181	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	201	5/5	0.75	0.12	110,112,116,121	0
3	CL	I	202	1/1	0.75	0.14	97,97,97,97	0
2	SO4	C	201	5/5	0.76	0.14	148,149,149,150	0
2	SO4	C	202	5/5	0.76	0.12	139,140,142,143	0
2	SO4	B	203	5/5	0.76	0.13	130,132,133,135	0
2	SO4	F	203	5/5	0.76	0.13	107,108,111,111	0
2	SO4	H	204	5/5	0.77	0.11	97,99,103,106	0
2	SO4	J	203	5/5	0.77	0.10	95,96,100,107	0
2	SO4	C	203	5/5	0.78	0.11	109,110,113,114	0
2	SO4	G	202	5/5	0.78	0.14	138,140,144,145	0
2	SO4	G	203	5/5	0.79	0.13	117,124,127,128	0
2	SO4	M	201	5/5	0.80	0.14	167,168,170,171	0
2	SO4	M	202	5/5	0.80	0.15	180,180,181,182	0
2	SO4	E	201	5/5	0.80	0.12	148,151,153,154	0
3	CL	O	205	1/1	0.80	0.15	108,108,108,108	0
2	SO4	R	202	5/5	0.81	0.12	142,144,146,147	0
2	SO4	L	203	5/5	0.82	0.12	112,117,119,119	0
2	SO4	J	201	5/5	0.82	0.14	149,150,151,154	0
2	SO4	D	203	5/5	0.82	0.10	99,100,101,103	0
4	EDO	N	201	4/4	0.83	0.15	92,96,96,98	0
5	CA	J	204	1/1	0.83	0.20	94,94,94,94	0
2	SO4	K	202	5/5	0.84	0.14	167,169,171,171	0
3	CL	A	202	1/1	0.84	0.22	106,106,106,106	0
5	CA	B	205	1/1	0.85	0.11	121,121,121,121	0
3	CL	D	206	1/1	0.85	0.14	91,91,91,91	0
5	CA	O	204	1/1	0.85	0.10	113,113,113,113	0
3	CL	H	206	1/1	0.86	0.10	99,99,99,99	0
4	EDO	B	201	4/4	0.86	0.18	60,61,62,68	0
2	SO4	B	202	5/5	0.87	0.12	67,89,97,106	0
2	SO4	B	204	5/5	0.87	0.08	81,81,85,89	0
3	CL	I	203	1/1	0.87	0.14	85,85,85,85	0
6	PGE	H	201	10/10	0.87	0.15	70,73,78,78	0
3	CL	H	207	1/1	0.88	0.13	92,92,92,92	0
3	CL	F	205	1/1	0.89	0.18	88,88,88,88	0
4	EDO	G	201	4/4	0.89	0.15	55,62,63,63	0
4	EDO	L	201	4/4	0.89	0.12	55,55,59,62	0
3	CL	E	203	1/1	0.89	0.11	95,95,95,95	0
5	CA	D	204	1/1	0.90	0.10	99,99,99,99	0
3	CL	K	205	1/1	0.90	0.15	83,83,83,83	0
4	EDO	D	201	4/4	0.91	0.09	46,51,55,58	0
3	CL	E	202	1/1	0.91	0.12	93,93,93,93	0
4	EDO	K	201	4/4	0.91	0.15	59,64,73,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	G	204	1/1	0.91	0.15	89,89,89,89	0
3	CL	J	206	1/1	0.91	0.14	96,96,96,96	0
4	EDO	F	201	4/4	0.92	0.10	65,68,70,72	0
3	CL	L	204	1/1	0.93	0.10	74,74,74,74	0
3	CL	K	204	1/1	0.93	0.09	85,85,85,85	0
3	CL	C	204	1/1	0.94	0.10	92,92,92,92	0
2	SO4	O	202	5/5	0.95	0.07	57,63,77,79	0
3	CL	D	205	1/1	0.95	0.12	79,79,79,79	0
2	SO4	D	202	5/5	0.96	0.13	54,55,58,73	0
2	SO4	L	202	5/5	0.96	0.08	58,61,62,62	0
5	CA	H	205	1/1	0.97	0.15	69,69,69,69	0
3	CL	J	205	1/1	0.97	0.08	76,76,76,76	0
5	CA	F	204	1/1	0.98	0.15	72,72,72,72	0
3	CL	G	205	1/1	0.99	0.11	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.