



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:22 am GMT

PDB ID : 2WKE
Title : CRYSTAL STRUCTURE OF THE ACTINOMADURA R39 DD-PEPTIDASE INHIBITED BY 6-BETA-IODOPENICILLANATE
Authors : Sauvage, E.; Herman, R.; Kerff, F.; Charlier, P.
Deposited on : 2009-06-10
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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

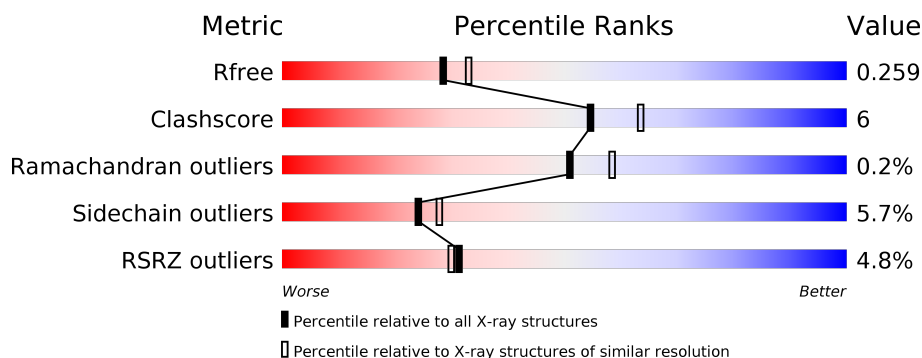
1 Overall quality at a glance

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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	466	<div> <div>6%</div> <div>86%</div> <div>12%</div> <div>.</div> </div>
1	B	466	<div> <div>5%</div> <div>87%</div> <div>12%</div> <div>.</div> </div>
1	C	466	<div> <div>4%</div> <div>88%</div> <div>10%</div> <div>.</div> </div>
1	D	466	<div> <div>5%</div> <div>85%</div> <div>13%</div> <div>.</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	603	-	-	-	X
3	SO4	A	605	-	-	X	-
3	SO4	B	603	-	-	-	X
3	SO4	B	605	-	-	X	-
3	SO4	C	603	-	-	-	X
3	SO4	C	605	-	-	X	-
3	SO4	D	605	-	-	X	-

2 Entry composition [i](#)

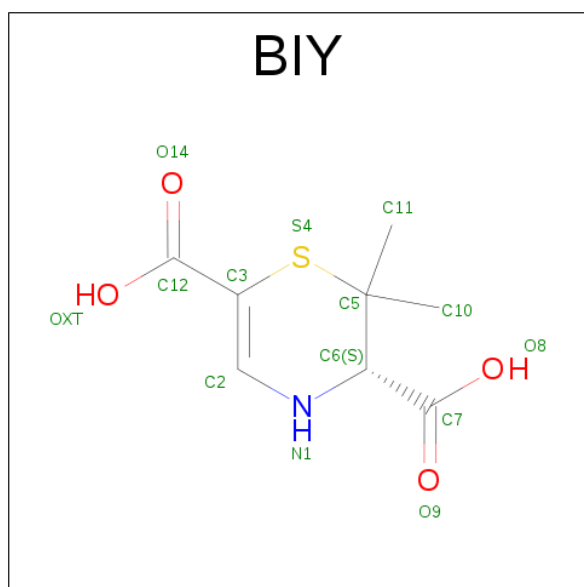
There are 5 unique types of molecules in this entry. The entry contains 14160 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-ALANYL-D-ALANINE CARBOXYPEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	0	0
			3353	2076	564	707	6			
1	B	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	C	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	D	466	Total	C	N	O	S	0	0	0
			3353	2076	564	707	6			

- Molecule 2 is (3S)-2,2-DIMETHYL-3,4-DIHYDRO-2H-1,4-THIAZINE-3,6-DICARBOXYLIC ACID (three-letter code: BIY) (formula: C₈H₁₁NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
2	B	1	Total	C	N	O	S	0	0
			13	8	1	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
2	D	1	Total	C	N	O	S	0	0
			13	8	1	3	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 4 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	3	Total Co 3 3	0	0
4	A	4	Total Co 4 4	0	0
4	D	4	Total Co 4 4	0	0
4	C	3	Total Co 3 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	170	Total O 170 170	0	0
5	B	143	Total O 143 143	0	0

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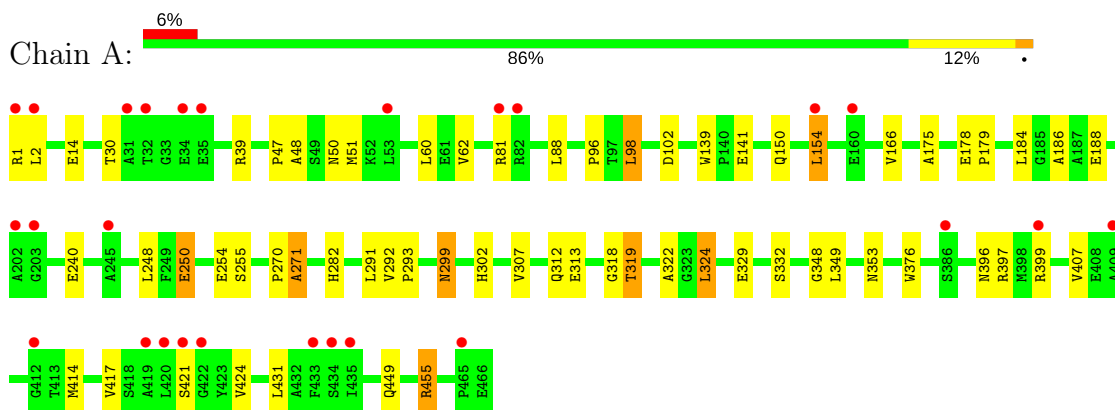
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	148	Total 148	O 148	0	0
5	D	139	Total 139	O 139	0	0

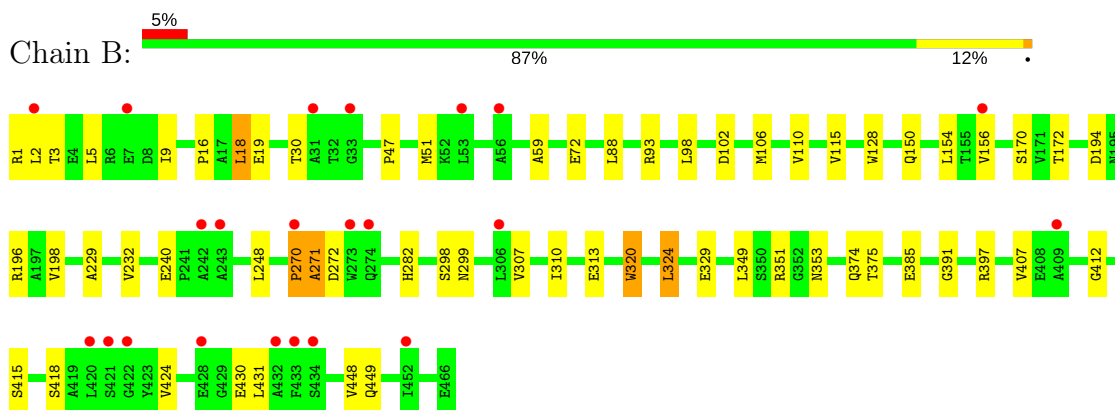
3 Residue-property plots [i](#)

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.

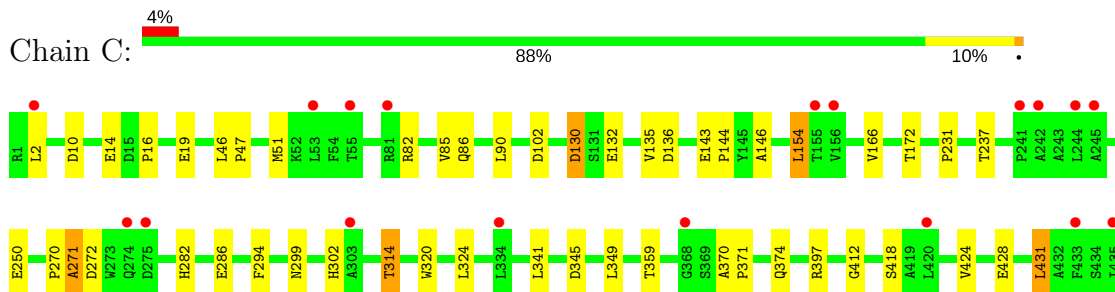
• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

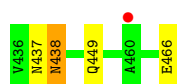


• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE



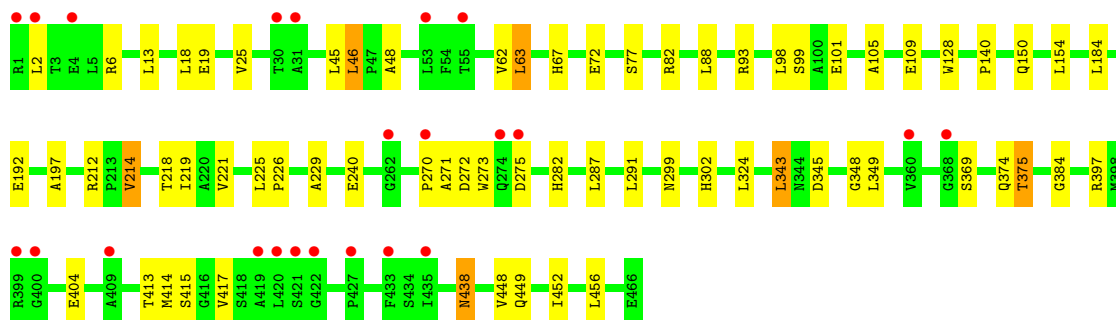
• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE





● Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

Chain D: 5% 85% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.98Å 92.69Å 107.13Å 90.00° 95.03° 90.00°	Depositor
Resolution (Å)	29.06 – 2.20 29.06 – 2.20	Depositor EDS
% Data completeness (in resolution range)	91.5 (29.06-2.20) 91.5 (29.06-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.206 , 0.261 0.206 , 0.259	Depositor DCC
R_{free} test set	4728 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14160	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.75% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CO, SO4, BIY

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	0.58	0/3412	0.69	2/4666 (0.0%)
1	B	0.56	0/3403	0.66	0/4656
1	C	0.54	0/3403	0.63	0/4656
1	D	0.53	0/3412	0.67	1/4666 (0.0%)
All	All	0.55	0/13630	0.66	3/18644 (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	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	343	LEU	CA-CB-CG	6.26	129.71	115.30
1	A	98	LEU	CA-CB-CG	5.38	127.68	115.30
1	A	154	LEU	CB-CG-CD2	5.03	119.56	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	318	GLY	Peptide

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	3353	0	3200	45	0
1	B	3344	0	3194	39	0
1	C	3344	0	3194	36	0
1	D	3353	0	3200	38	0
2	A	13	0	9	1	0
2	B	13	0	9	1	0
2	C	13	0	9	0	0
2	D	13	0	9	0	0
3	A	25	0	0	3	0
3	B	25	0	0	6	0
3	C	25	0	0	6	0
3	D	25	0	0	3	0
4	A	4	0	0	0	0
4	B	3	0	0	0	0
4	C	3	0	0	0	0
4	D	4	0	0	0	0
5	A	170	0	0	11	0
5	B	143	0	0	4	0
5	C	148	0	0	9	0
5	D	139	0	0	7	0
All	All	14160	0	12824	158	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 6.

All (158) 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:166:VAL:HG12	5:C:2101:HOH:O	1.23	1.38
1:D:282:HIS:NE2	3:D:605:SO4:O4	1.88	1.06
1:A:282:HIS:NE2	3:A:605:SO4:O1	1.98	0.95
1:B:51:MET:HE3	1:B:353:ASN:HB3	1.48	0.95
1:B:51:MET:CE	1:B:353:ASN:HB3	1.97	0.93
1:B:282:HIS:NE2	3:B:605:SO4:O3	2.09	0.86
1:D:287:LEU:HB3	5:D:2009:HOH:O	1.75	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:THR:HG21	5:A:2127:HOH:O	1.77	0.85
1:A:51:MET:HE2	1:A:353:ASN:HB3	1.58	0.85
1:A:270:PRO:O	1:A:271:ALA:HB3	1.77	0.83
1:C:102:ASP:OD2	3:C:605:SO4:O3	1.97	0.82
1:A:1:ARG:HH22	1:A:455:ARG:HH22	1.28	0.81
1:B:102:ASP:OD2	3:B:605:SO4:O3	1.99	0.81
1:C:282:HIS:NE2	3:C:605:SO4:O3	2.14	0.81
1:C:237:THR:HG22	5:C:2101:HOH:O	1.83	0.77
1:C:16:PRO:HA	1:C:19:GLU:OE2	1.86	0.76
1:A:319:THR:HG22	1:A:322:ALA:H	1.50	0.75
1:C:374:GLN:HG2	5:D:2091:HOH:O	1.88	0.74
1:A:270:PRO:O	1:A:271:ALA:CB	2.35	0.73
1:C:282:HIS:NE2	3:C:605:SO4:S	2.61	0.73
1:A:166:VAL:HG12	5:A:2097:HOH:O	1.90	0.72
1:A:150:GLN:HE22	1:A:240:GLU:H	1.38	0.71
1:D:150:GLN:NE2	1:D:240:GLU:H	1.88	0.71
1:A:51:MET:CE	1:A:353:ASN:HB3	2.21	0.70
1:D:282:HIS:CD2	3:D:605:SO4:O4	2.43	0.70
1:D:282:HIS:NE2	3:D:605:SO4:S	2.56	0.70
1:B:51:MET:HE3	1:B:353:ASN:CB	2.24	0.68
1:D:150:GLN:HE22	1:D:240:GLU:H	1.38	0.68
1:A:178:GLU:HG3	1:A:179:PRO:HD2	1.75	0.67
1:B:282:HIS:NE2	3:B:605:SO4:S	2.62	0.67
1:B:351:ARG:NH2	1:B:415:SER:O	2.28	0.66
1:C:166:VAL:CG1	5:C:2101:HOH:O	2.04	0.66
1:A:1:ARG:HG3	5:A:2002:HOH:O	1.94	0.66
1:B:229:ALA:O	5:B:2088:HOH:O	2.14	0.66
1:B:150:GLN:HE22	1:B:240:GLU:H	1.43	0.65
1:C:397:ARG:HH12	1:C:449:GLN:HE21	1.44	0.65
1:A:102:ASP:OD1	3:A:605:SO4:O1	2.14	0.65
1:B:150:GLN:NE2	1:B:240:GLU:H	1.94	0.65
1:B:270:PRO:O	1:B:271:ALA:CB	2.45	0.64
1:C:172:THR:HG22	1:C:231:PRO:HB3	1.79	0.64
1:C:237:THR:CG2	5:C:2101:HOH:O	2.43	0.63
1:D:77:SER:HB2	5:D:2014:HOH:O	1.98	0.62
1:A:1:ARG:HH22	1:A:455:ARG:NH2	1.95	0.62
1:C:102:ASP:OD2	3:C:605:SO4:S	2.57	0.62
1:B:375:THR:HG23	5:B:2119:HOH:O	1.98	0.62
1:A:139:TRP:HB3	1:A:141:GLU:OE1	2.00	0.61
1:A:150:GLN:NE2	1:A:240:GLU:H	1.99	0.60
1:D:384:GLY:HA3	1:D:404:GLU:HG3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:270:PRO:O	1:C:271:ALA:CB	2.50	0.60
1:D:99:SER:OG	1:D:101:GLU:HG2	2.00	0.60
1:C:102:ASP:CG	3:C:605:SO4:O3	2.39	0.60
1:B:59:ALA:HA	1:B:310:ILE:HD11	1.83	0.60
1:B:51:MET:CE	1:B:353:ASN:CB	2.77	0.60
1:B:397:ARG:HH12	1:B:449:GLN:HE21	1.48	0.60
1:B:270:PRO:O	1:B:271:ALA:HB3	2.02	0.59
1:C:282:HIS:CE1	3:C:605:SO4:O3	2.56	0.59
1:D:270:PRO:O	1:D:271:ALA:HB3	2.01	0.59
1:C:130:ASP:HB3	1:C:132:GLU:H	1.67	0.58
1:C:397:ARG:HH12	1:C:449:GLN:NE2	2.02	0.57
1:C:47:PRO:HG3	1:C:51:MET:CE	2.35	0.56
1:C:270:PRO:O	1:C:271:ALA:HB3	2.05	0.56
1:B:102:ASP:OD2	3:B:605:SO4:S	2.63	0.56
1:D:67:HIS:N	5:D:2009:HOH:O	2.39	0.55
1:A:141:GLU:HG2	5:A:2059:HOH:O	2.05	0.55
1:A:1:ARG:NH2	1:A:455:ARG:HH22	1.99	0.55
1:B:47:PRO:HG3	1:B:51:MET:HE1	1.88	0.55
1:D:197:ALA:HB2	1:D:221:VAL:HG12	1.88	0.55
1:A:102:ASP:OD1	3:A:605:SO4:S	2.64	0.55
1:D:6:ARG:HD3	5:D:2002:HOH:O	2.07	0.55
1:A:141:GLU:CG	5:A:2059:HOH:O	2.56	0.53
1:A:319:THR:CG2	1:A:322:ALA:H	2.20	0.53
1:D:72:GLU:OE1	1:D:93:ARG:NH2	2.41	0.52
1:A:166:VAL:CG1	5:A:2097:HOH:O	2.55	0.52
1:A:250:GLU:HG2	5:A:2103:HOH:O	2.10	0.52
1:C:370:ALA:HB1	1:C:371:PRO:HD2	1.92	0.50
1:A:299:ASN:ND2	1:A:302:HIS:H	2.11	0.49
2:B:500:BIY:H103	2:B:500:BIY:C2	2.43	0.48
1:B:407:VAL:HG22	1:B:424:VAL:HG22	1.95	0.48
1:D:48:ALA:O	1:D:348:GLY:HA3	2.13	0.48
1:C:286:GLU:HG3	5:C:2107:HOH:O	2.13	0.48
1:A:51:MET:HE2	1:A:353:ASN:CB	2.37	0.48
1:A:271:ALA:HA	5:A:2109:HOH:O	2.14	0.48
1:B:282:HIS:CD2	3:B:605:SO4:O3	2.67	0.48
1:B:93:ARG:HD3	1:B:128:TRP:CE3	2.49	0.47
1:C:90:LEU:HG	1:C:154:LEU:HD11	1.96	0.47
1:C:85:VAL:HG12	1:C:86:GLN:N	2.29	0.47
1:D:140:PRO:HB3	1:D:214:VAL:HG11	1.97	0.47
1:D:225:LEU:HD12	1:D:226:PRO:HD2	1.97	0.47
1:B:18:LEU:CD2	1:B:448:VAL:HG11	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:VAL:HG21	1:C:345:ASP:HB3	1.97	0.47
1:D:452:ILE:O	1:D:456:LEU:HG	2.14	0.47
1:A:81:ARG:HD3	1:B:313:GLU:OE1	2.14	0.47
1:D:18:LEU:HD11	1:D:448:VAL:HG11	1.96	0.47
1:A:397:ARG:HH12	1:A:449:GLN:HE21	1.63	0.47
1:C:466:GLU:N	5:C:2144:HOH:O	2.48	0.47
1:D:287:LEU:CB	5:D:2009:HOH:O	2.48	0.47
1:B:5:LEU:O	1:B:9:ILE:HG13	2.14	0.47
1:C:46:LEU:CD2	5:C:2120:HOH:O	2.63	0.46
1:D:414:MET:HG2	1:D:417:VAL:HB	1.97	0.46
2:A:500:BIY:O9	2:A:500:BIY:H103	2.15	0.46
1:A:60:LEU:HD11	1:A:291:LEU:HD11	1.97	0.46
1:C:424:VAL:HB	1:C:431:LEU:HB2	1.97	0.46
1:B:424:VAL:HB	1:B:431:LEU:HB2	1.98	0.46
1:A:184:LEU:HB2	1:A:188:GLU:HG2	1.98	0.46
1:B:156:VAL:HG21	1:B:248:LEU:HD12	1.97	0.46
1:D:397:ARG:HH12	1:D:449:GLN:HE21	1.63	0.45
1:D:45:LEU:HD22	1:D:438:ASN:HB2	1.97	0.45
1:B:1:ARG:HD2	1:D:2:LEU:HG	1.98	0.45
1:C:314:THR:CB	5:C:2111:HOH:O	2.64	0.45
1:B:385:GLU:O	1:B:391:GLY:HA3	2.15	0.45
1:B:110:VAL:HG12	1:B:115:VAL:HB	1.97	0.45
1:A:96:PRO:HB2	1:A:302:HIS:CD2	2.52	0.45
1:B:72:GLU:HG2	5:B:2026:HOH:O	2.17	0.45
1:B:397:ARG:HH12	1:B:449:GLN:NE2	2.14	0.45
1:C:294:PHE:HB2	1:C:302:HIS:HD2	1.81	0.45
1:B:307:VAL:HG11	1:B:324:LEU:HD13	1.99	0.45
1:A:47:PRO:HD2	5:A:2134:HOH:O	2.17	0.44
1:B:194:ASP:OD2	1:B:196:ARG:HD3	2.16	0.44
1:A:396:ASN:OD1	1:A:399:ARG:NH1	2.51	0.44
1:A:186:ALA:HB1	1:A:248:LEU:HD11	2.00	0.44
1:C:397:ARG:NH1	1:C:449:GLN:HE21	2.13	0.44
1:D:270:PRO:O	1:D:271:ALA:CB	2.66	0.44
1:C:146:ALA:HA	5:C:2101:HOH:O	2.18	0.44
1:D:192:GLU:HG3	1:D:218:THR:OG1	2.17	0.44
1:D:226:PRO:HG2	1:D:229:ALA:HB2	1.99	0.44
1:A:307:VAL:HG11	1:A:324:LEU:HD13	1.99	0.44
1:A:407:VAL:HG22	1:A:424:VAL:HG22	2.00	0.44
1:B:102:ASP:CG	3:B:605:SO4:O3	2.56	0.43
1:C:143:GLU:N	1:C:144:PRO:CD	2.82	0.43
1:B:320:TRP:O	1:B:324:LEU:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:ASP:O	1:C:14:GLU:HG2	2.19	0.43
1:C:412:GLY:O	1:C:418:SER:HA	2.19	0.43
1:B:170:SER:HA	1:B:232:VAL:O	2.19	0.43
1:D:271:ALA:C	1:D:273:TRP:H	2.22	0.43
1:D:13:LEU:HD11	1:D:25:VAL:HG21	2.00	0.42
1:D:397:ARG:NH1	1:D:449:GLN:HE21	2.16	0.42
1:A:1:ARG:HH12	1:A:455:ARG:HH12	1.65	0.42
1:D:93:ARG:HG2	1:D:128:TRP:CD2	2.54	0.42
1:A:414:MET:HB2	1:A:417:VAL:HB	2.01	0.42
1:C:437:ASN:C	1:C:438:ASN:HD22	2.23	0.42
1:B:106:MET:O	1:B:110:VAL:HG23	2.20	0.42
1:D:212:ARG:HB2	1:D:219:ILE:HD13	2.01	0.42
1:D:397:ARG:HH12	1:D:449:GLN:NE2	2.18	0.42
1:A:175:ALA:HB3	5:A:2085:HOH:O	2.20	0.42
1:D:105:ALA:O	1:D:109:GLU:HG3	2.20	0.42
1:A:292:VAL:HB	1:A:293:PRO:CD	2.50	0.41
1:A:299:ASN:HD22	1:A:299:ASN:C	2.24	0.41
1:D:375:THR:HG23	5:D:2117:HOH:O	2.19	0.41
1:D:345:ASP:OD1	1:D:345:ASP:C	2.59	0.41
1:D:62:VAL:HG12	1:D:63:LEU:HD13	2.00	0.41
1:C:341:LEU:HG	1:C:359:THR:HG21	2.02	0.41
1:D:46:LEU:N	1:D:46:LEU:HD23	2.36	0.41
1:B:412:GLY:O	1:B:418:SER:HA	2.20	0.41
1:A:271:ALA:HB2	5:A:2034:HOH:O	2.21	0.41
1:A:62:VAL:HG12	1:A:313:GLU:HG2	2.03	0.40
1:A:48:ALA:O	1:A:348:GLY:HA3	2.21	0.40
1:A:50:ASN:ND2	1:A:421:SER:OG	2.55	0.40
1:B:172:THR:HB	5:B:2019:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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	464/466 (100%)	450 (97%)	13 (3%)	1 (0%)	51	58
1	B	464/466 (100%)	444 (96%)	19 (4%)	1 (0%)	51	58
1	C	464/466 (100%)	444 (96%)	18 (4%)	2 (0%)	38	41
1	D	464/466 (100%)	447 (96%)	17 (4%)	0	100	100
All	All	1856/1864 (100%)	1785 (96%)	67 (4%)	4 (0%)	51	58

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	271	ALA
1	C	271	ALA
1	C	130	ASP
1	A	271	ALA

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	339/339 (100%)	319 (94%)	20 (6%)	23	26
1	B	338/339 (100%)	318 (94%)	20 (6%)	23	26
1	C	338/339 (100%)	324 (96%)	14 (4%)	35	44
1	D	339/339 (100%)	316 (93%)	23 (7%)	18	20
All	All	1354/1356 (100%)	1277 (94%)	77 (6%)	24	28

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	14	GLU
1	A	30	THR
1	A	39	ARG
1	A	88	LEU
1	A	98	LEU
1	A	154	LEU

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Mol	Chain	Res	Type
1	A	250	GLU
1	A	254	GLU
1	A	255	SER
1	A	299	ASN
1	A	312	GLN
1	A	319	THR
1	A	324	LEU
1	A	329	GLU
1	A	332	SER
1	A	349	LEU
1	A	376	TRP
1	A	431	LEU
1	A	455	ARG
1	B	2	LEU
1	B	3	THR
1	B	16	PRO
1	B	18	LEU
1	B	19	GLU
1	B	30	THR
1	B	88	LEU
1	B	98	LEU
1	B	154	LEU
1	B	198	VAL
1	B	270	PRO
1	B	272	ASP
1	B	298	SER
1	B	299	ASN
1	B	320	TRP
1	B	324	LEU
1	B	329	GLU
1	B	349	LEU
1	B	374	GLN
1	B	430	GLU
1	C	2	LEU
1	C	82	ARG
1	C	136	ASP
1	C	154	LEU
1	C	250	GLU
1	C	272	ASP
1	C	299	ASN
1	C	314	THR
1	C	320	TRP

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Mol	Chain	Res	Type
1	C	324	LEU
1	C	349	LEU
1	C	428	GLU
1	C	431	LEU
1	C	438	ASN
1	D	19	GLU
1	D	46	LEU
1	D	63	LEU
1	D	82	ARG
1	D	88	LEU
1	D	98	LEU
1	D	154	LEU
1	D	184	LEU
1	D	214	VAL
1	D	272	ASP
1	D	275	ASP
1	D	291	LEU
1	D	299	ASN
1	D	302	HIS
1	D	324	LEU
1	D	343	LEU
1	D	349	LEU
1	D	369	SER
1	D	374	GLN
1	D	375	THR
1	D	413	THR
1	D	415	SER
1	D	438	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	150	GLN
1	A	299	ASN
1	A	312	GLN
1	A	437	ASN
1	A	449	GLN
1	B	44	GLN
1	B	50	ASN
1	B	150	GLN
1	B	396	ASN

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Mol	Chain	Res	Type
1	B	437	ASN
1	B	449	GLN
1	C	50	ASN
1	C	299	ASN
1	C	302	HIS
1	C	437	ASN
1	C	449	GLN
1	D	44	GLN
1	D	50	ASN
1	D	150	GLN
1	D	299	ASN
1	D	302	HIS
1	D	366	GLN
1	D	437	ASN
1	D	449	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](#)

Of 38 ligands modelled in this entry, 14 are monoatomic - leaving 24 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	BIY	A	500	1	7,13,14	3.01	3 (42%)	7,19,21	2.19	2 (28%)
3	SO4	A	601	-	4,4,4	0.25	0	6,6,6	0.46	0
3	SO4	A	602	-	4,4,4	0.19	0	6,6,6	0.22	0
3	SO4	A	603	-	4,4,4	0.20	0	6,6,6	0.37	0
3	SO4	A	604	-	4,4,4	0.12	0	6,6,6	0.22	0
3	SO4	A	605	-	4,4,4	0.87	0	6,6,6	0.65	0
2	BIY	B	500	1	7,13,14	3.19	3 (42%)	7,19,21	1.09	0
3	SO4	B	601	-	4,4,4	0.15	0	6,6,6	0.39	0
3	SO4	B	602	-	4,4,4	0.20	0	6,6,6	0.19	0
3	SO4	B	603	-	4,4,4	0.20	0	6,6,6	0.20	0
3	SO4	B	604	-	4,4,4	0.16	0	6,6,6	0.23	0
3	SO4	B	605	-	4,4,4	0.37	0	6,6,6	0.42	0
2	BIY	C	500	1	7,13,14	2.21	4 (57%)	7,19,21	1.60	2 (28%)
3	SO4	C	601	-	4,4,4	0.11	0	6,6,6	0.37	0
3	SO4	C	602	-	4,4,4	0.13	0	6,6,6	0.18	0
3	SO4	C	603	-	4,4,4	0.17	0	6,6,6	0.18	0
3	SO4	C	604	-	4,4,4	0.14	0	6,6,6	0.35	0
3	SO4	C	605	-	4,4,4	0.54	0	6,6,6	0.33	0
2	BIY	D	500	1	7,13,14	4.70	4 (57%)	7,19,21	1.71	2 (28%)
3	SO4	D	601	-	4,4,4	0.15	0	6,6,6	0.27	0
3	SO4	D	602	-	4,4,4	0.15	0	6,6,6	0.20	0
3	SO4	D	603	-	4,4,4	0.32	0	6,6,6	0.91	0
3	SO4	D	604	-	4,4,4	0.19	0	6,6,6	0.40	0
3	SO4	D	605	-	4,4,4	0.35	0	6,6,6	0.26	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
2	BIY	A	500	1	-	0/0/22/24	0/0/1/1
3	SO4	A	601	-	-	0/0/0/0	0/0/0/0
3	SO4	A	602	-	-	0/0/0/0	0/0/0/0
3	SO4	A	603	-	-	0/0/0/0	0/0/0/0
3	SO4	A	604	-	-	0/0/0/0	0/0/0/0
3	SO4	A	605	-	-	0/0/0/0	0/0/0/0
2	BIY	B	500	1	-	0/0/22/24	0/0/1/1
3	SO4	B	601	-	-	0/0/0/0	0/0/0/0
3	SO4	B	602	-	-	0/0/0/0	0/0/0/0
3	SO4	B	603	-	-	0/0/0/0	0/0/0/0
3	SO4	B	604	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	605	-	-	0/0/0/0	0/0/0/0
2	BIY	C	500	1	-	0/0/22/24	0/0/1/1
3	SO4	C	601	-	-	0/0/0/0	0/0/0/0
3	SO4	C	602	-	-	0/0/0/0	0/0/0/0
3	SO4	C	603	-	-	0/0/0/0	0/0/0/0
3	SO4	C	604	-	-	0/0/0/0	0/0/0/0
3	SO4	C	605	-	-	0/0/0/0	0/0/0/0
2	BIY	D	500	1	-	0/0/22/24	0/0/1/1
3	SO4	D	601	-	-	0/0/0/0	0/0/0/0
3	SO4	D	602	-	-	0/0/0/0	0/0/0/0
3	SO4	D	603	-	-	0/0/0/0	0/0/0/0
3	SO4	D	604	-	-	0/0/0/0	0/0/0/0
3	SO4	D	605	-	-	0/0/0/0	0/0/0/0

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	BIY	C3-S4	-2.67	1.67	1.73
2	C	500	BIY	C3-S4	-2.41	1.67	1.73
2	A	500	BIY	C3-S4	-2.27	1.68	1.73
2	D	500	BIY	C3-S4	-2.03	1.68	1.73
2	C	500	BIY	C12-C3	2.01	1.48	1.45
2	D	500	BIY	C12-C3	2.99	1.49	1.45
2	A	500	BIY	C6-N1	3.17	1.50	1.45
2	C	500	BIY	C6-N1	3.36	1.50	1.45
2	C	500	BIY	C2-N1	3.57	1.39	1.34
2	B	500	BIY	C6-N1	3.78	1.50	1.45
2	D	500	BIY	C6-N1	3.86	1.51	1.45
2	A	500	BIY	C2-N1	6.80	1.44	1.34
2	B	500	BIY	C2-N1	6.81	1.44	1.34
2	D	500	BIY	C2-N1	11.22	1.51	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	BIY	O14-C12-C3	-4.71	115.20	124.94
2	C	500	BIY	O14-C12-C3	-2.87	119.00	124.94
2	A	500	BIY	C10-C5-C11	-2.87	106.27	109.63
2	D	500	BIY	C10-C5-C11	-2.86	106.28	109.63
2	C	500	BIY	C10-C5-C11	-2.44	106.77	109.63
2	D	500	BIY	C11-C5-S4	2.36	114.07	107.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	BIY	1	0
3	A	605	SO4	3	0
2	B	500	BIY	1	0
3	B	605	SO4	6	0
3	C	605	SO4	6	0
3	D	605	SO4	3	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	466/466 (100%)	0.13	26 (5%) 25 24	11, 19, 27, 32	0
1	B	466/466 (100%)	-0.00	22 (4%) 32 31	13, 21, 29, 35	0
1	C	466/466 (100%)	0.03	19 (4%) 38 36	11, 19, 27, 33	0
1	D	466/466 (100%)	0.20	23 (4%) 30 29	9, 23, 34, 39	0
All	All	1864/1864 (100%)	0.09	90 (4%) 31 30	9, 20, 31, 39	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	2	LEU	6.9
1	A	2	LEU	5.7
1	D	1	ARG	5.4
1	B	428	GLU	5.3
1	B	433	PHE	4.4
1	D	274	GLN	4.3
1	A	422	GLY	4.3
1	A	31	ALA	3.9
1	D	31	ALA	3.8
1	D	270	PRO	3.8
1	B	306	LEU	3.8
1	D	421	SER	3.8
1	D	419	ALA	3.7
1	B	421	SER	3.7
1	A	435	ILE	3.7
1	C	303	ALA	3.7
1	D	433	PHE	3.7
1	A	53	LEU	3.6
1	B	53	LEU	3.6
1	A	1	ARG	3.6
1	D	409	ALA	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	386	SER	3.5
1	A	245	ALA	3.5
1	B	422	GLY	3.4
1	C	55	THR	3.3
1	A	433	PHE	3.3
1	C	245	ALA	3.2
1	D	275	ASP	3.2
1	D	427	PRO	3.1
1	A	203	GLY	3.1
1	D	262	GLY	3.0
1	C	275	ASP	2.9
1	D	435	ILE	2.9
1	C	155	THR	2.9
1	C	81	ARG	2.9
1	A	35	GLU	2.9
1	A	421	SER	2.8
1	D	422	GLY	2.8
1	A	419	ALA	2.8
1	C	241	PRO	2.8
1	A	420	LEU	2.7
1	C	420	LEU	2.7
1	B	274	GLN	2.7
1	A	81	ARG	2.7
1	B	7	GLU	2.6
1	B	409	ALA	2.6
1	D	4	GLU	2.6
1	C	460	ALA	2.6
1	B	432	ALA	2.6
1	D	368	GLY	2.6
1	B	156	VAL	2.5
1	C	433	PHE	2.5
1	C	368	GLY	2.5
1	B	273	TRP	2.5
1	B	242	ALA	2.5
1	A	160	GLU	2.5
1	B	270	PRO	2.5
1	D	55	THR	2.5
1	A	412	GLY	2.5
1	A	409	ALA	2.4
1	A	434	SER	2.4
1	D	420	LEU	2.4
1	B	243	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	53	LEU	2.4
1	A	202	ALA	2.3
1	A	399	ARG	2.3
1	C	274	GLN	2.3
1	B	33	GLY	2.3
1	D	400	GLY	2.3
1	A	34	GLU	2.3
1	C	242	ALA	2.3
1	D	399	ARG	2.2
1	A	154	LEU	2.2
1	B	420	LEU	2.2
1	D	53	LEU	2.2
1	B	434	SER	2.2
1	C	334	LEU	2.2
1	B	56	ALA	2.1
1	D	360	VAL	2.1
1	A	32	THR	2.1
1	B	31	ALA	2.1
1	C	156	VAL	2.1
1	A	82	ARG	2.1
1	C	244	LEU	2.1
1	A	465	PRO	2.1
1	B	452	ILE	2.0
1	D	30	THR	2.0
1	C	435	ILE	2.0
1	B	2	LEU	2.0
1	D	2	LEU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	C	603	5/5	0.95	0.26	10.44	60,60,61,61	0
3	SO4	B	603	5/5	0.97	0.18	9.85	46,46,47,48	0
3	SO4	A	603	5/5	0.89	0.23	3.85	56,57,59,59	0
2	BIY	C	500	13/14	0.95	0.12	-0.44	32,35,39,39	0
3	SO4	A	605	5/5	0.93	0.12	-0.48	17,25,27,29	0
2	BIY	D	500	13/14	0.90	0.14	-0.56	26,37,39,40	0
2	BIY	A	500	13/14	0.92	0.13	-0.71	29,37,42,45	0
4	CO	A	612	1/1	0.94	0.10	-0.77	66,66,66,66	0
3	SO4	B	605	5/5	0.96	0.09	-0.83	30,32,34,35	0
4	CO	B	610	1/1	0.97	0.08	-1.16	48,48,48,48	0
2	BIY	B	500	13/14	0.95	0.09	-1.19	32,37,37,37	0
3	SO4	D	605	5/5	0.96	0.08	-1.31	35,37,37,40	0
3	SO4	C	605	5/5	0.99	0.07	-1.51	22,24,28,29	0
4	CO	A	611	1/1	0.98	0.05	-2.00	31,31,31,31	0
4	CO	D	611	1/1	0.99	0.06	-2.64	46,46,46,46	0
3	SO4	D	603	5/5	0.97	0.08	-2.76	28,29,31,32	0
4	CO	A	610	1/1	0.99	0.02	-3.25	33,33,33,33	0
4	CO	D	610	1/1	0.99	0.03	-3.58	35,35,35,35	0
4	CO	C	610	1/1	0.99	0.04	-4.59	47,47,47,47	0
4	CO	B	613	1/1	0.78	0.10	-	97,97,97,97	0
3	SO4	C	601	5/5	0.99	0.20	-	38,39,40,40	0
3	SO4	D	604	5/5	0.95	0.25	-	38,41,43,43	0
3	SO4	A	604	5/5	0.91	0.36	-	70,71,72,72	0
4	CO	C	612	1/1	0.97	0.11	-	67,67,67,67	0
4	CO	A	613	1/1	0.96	0.07	-	62,62,62,62	0
4	CO	C	613	1/1	0.89	0.14	-	95,95,95,95	0
4	CO	D	612	1/1	0.91	0.10	-	66,66,66,66	0
3	SO4	C	604	5/5	0.97	0.20	-	44,44,45,47	0
3	SO4	A	602	5/5	0.93	0.20	-	63,64,65,65	0
3	SO4	B	604	5/5	0.96	0.20	-	49,51,51,52	0
4	CO	B	612	1/1	0.97	0.09	-	67,67,67,67	0
3	SO4	D	601	5/5	0.98	0.18	-	52,52,53,54	0
3	SO4	A	601	5/5	0.98	0.21	-	32,34,36,37	0
3	SO4	D	602	5/5	0.91	0.20	-	66,66,66,67	0
3	SO4	B	602	5/5	0.95	0.16	-	55,55,56,57	0
4	CO	D	613	1/1	0.91	0.08	-	87,87,87,87	0
3	SO4	C	602	5/5	0.96	0.17	-	52,52,54,54	0
3	SO4	B	601	5/5	0.97	0.26	-	52,52,53,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.