



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2023 – 11:44 AM EDT

PDB ID : 5ES5
Title : Crystal structure of the initiation module of LgrA in the "open" and "closed" adenylation states
Authors : Reimer, J.M.; Aloise, M.N.; Schmeing, T.M.
Deposited on : 2015-11-16
Resolution : 2.80 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

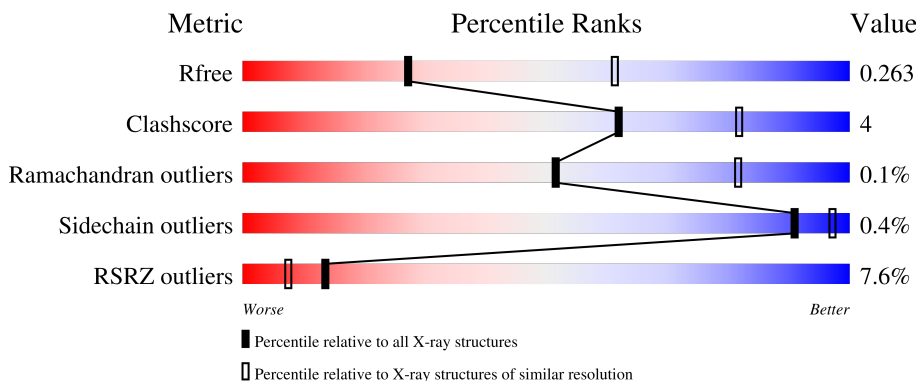
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	776	 5% 81% 7% 12%
1	B	776	 9% 78% 10% 12%

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 21585 atoms, of which 10679 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 Linear gramicidin synthetase subunit A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	683	Total	C	H	N	O	S	0	0	0
			10764	3471	5324	930	1014	25			
1	B	685	Total	C	H	N	O	S	0	0	0
			10802	3476	5355	930	1016	25			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	initiating methionine	UNP Q70LM7
B	2	GLY	-	expression tag	UNP Q70LM7
B	768	ALA	-	expression tag	UNP Q70LM7
B	769	ALA	-	expression tag	UNP Q70LM7
B	770	ALA	-	expression tag	UNP Q70LM7
B	771	GLU	-	expression tag	UNP Q70LM7
B	772	ASN	-	expression tag	UNP Q70LM7
B	773	LEU	-	expression tag	UNP Q70LM7
B	774	TYR	-	expression tag	UNP Q70LM7
B	775	PHE	-	expression tag	UNP Q70LM7
B	776	GLN	-	expression tag	UNP Q70LM7
A	1	MET	-	initiating methionine	UNP Q70LM7
A	2	GLY	-	expression tag	UNP Q70LM7
A	768	ALA	-	expression tag	UNP Q70LM7
A	769	ALA	-	expression tag	UNP Q70LM7
A	770	ALA	-	expression tag	UNP Q70LM7
A	771	GLU	-	expression tag	UNP Q70LM7
A	772	ASN	-	expression tag	UNP Q70LM7
A	773	LEU	-	expression tag	UNP Q70LM7
A	774	TYR	-	expression tag	UNP Q70LM7
A	775	PHE	-	expression tag	UNP Q70LM7
A	776	GLN	-	expression tag	UNP Q70LM7

- Molecule 2 is SULFATE ION (three-letter code: 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		

- Molecule 3 is water.

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

VAL
ALA
PRO
THR
THR
ASP
GLU
LEU
GLU
GLN
GLN
LEU
LEU
ALA
GLN
VAL
TRP
SER
HIS
SER
VAL
VAL
LEU
GLY
ILE
PRO
GLN
MET
GLY
ILE
ASP
ASP
HIS
PHE
LEU
LEU
ARG
GLY
GLY
GLY
ASP
SER
SER
ILE
LYS
VAL
MET
GLN
GLN
LEU
ILE
HIS
GLN
LEU
LYS
ASN
ILE
GLY
LEU
LEU
SER
TYR
ARG
ASP
GLN
LEU
PHE

THR
HIS
PRO
THR
THR
ILE
ARG
GLN
LEU
LYS
ARG
LEU
LEU
THR
GLN
GLU
GLN
ALA
SER
ALA
ALA
ALA
GLU
ASN
LEU
TYR
PHE
GLN

4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	278.84Å 278.84Å 82.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.05 – 2.80 46.05 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.05-2.80) 92.3 (46.05-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.81Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.227 , 0.263 0.227 , 0.263	Depositor DCC
R_{free} test set	2961 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	77.0	Xtrriage
Anisotropy	0.313	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 86.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.014 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21585	wwPDB-VP
Average B, all atoms (Å ²)	148.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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	0.25	0/5569	0.43	0/7564
1	B	0.26	0/5576	0.44	0/7574
All	All	0.26	0/11145	0.44	0/15138

There are no bond length outliers.

There are no bond angle outliers.

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	5440	5324	5357	31	2
1	B	5447	5355	5362	64	1
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	5	0	0	0	0
3	B	4	0	0	0	0
All	All	10906	10679	10719	95	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 4.

All (95) 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:B:611:GLU:C	1:B:612:LEU:HD23	1.64	1.17
1:A:199:LYS:O	1:A:510:ARG:NH1	2.03	0.91
1:A:512:ARG:NH2	1:A:561:GLU:OE2	2.14	0.80
1:B:611:GLU:O	1:B:612:LEU:HD23	1.85	0.75
1:B:231:TYR:O	1:B:235:ASN:ND2	2.23	0.72
1:B:612:LEU:HD23	1:B:612:LEU:N	2.03	0.71
1:B:367:ASN:HB3	1:B:487:THR:HG21	1.73	0.71
1:B:605:MET:HA	1:B:610:ILE:HD12	1.76	0.67
1:B:488:ILE:HG22	1:B:489:MET:H	1.61	0.66
1:B:80:ASN:OD1	1:B:89:SER:N	2.34	0.60
1:B:454:GLN:N	1:B:454:GLN:OE1	2.34	0.60
1:B:486:CYS:O	1:B:487:THR:OG1	2.20	0.59
1:A:75:SER:OG	1:A:79:TRP:O	2.18	0.59
1:B:585:ASP:OD1	1:B:586:ASN:N	2.36	0.59
1:B:605:MET:HE1	1:B:611:GLU:O	2.04	0.58
1:B:427:ASN:OD1	1:B:431:ASN:ND2	2.37	0.58
1:B:209:PHE:O	1:B:213:VAL:HG23	2.04	0.57
1:A:78:PRO:O	1:A:81:LYS:HE3	2.05	0.56
1:B:391:SER:OG	1:B:398:SER:HA	2.05	0.56
1:A:71:ASN:HB2	1:A:103:ILE:HD11	1.86	0.56
1:B:374:TRP:CG	1:B:508:ILE:HG23	2.40	0.56
1:B:89:SER:O	1:B:96:LYS:NZ	2.37	0.56
1:A:442:THR:OG1	1:A:464:ASP:HB2	2.06	0.55
1:B:363:ILE:HD12	1:B:487:THR:HG22	1.89	0.55
1:A:464:ASP:OD1	1:A:672:LYS:NZ	2.28	0.54
1:B:277:GLY:O	1:B:340:ASN:ND2	2.37	0.54
1:B:75:SER:OG	1:B:79:TRP:O	2.27	0.52
1:B:606:ASN:N	1:B:606:ASN:OD1	2.41	0.52
1:B:332:ASN:OD1	1:B:334:ASN:N	2.39	0.52
1:B:345:LEU:HD11	1:B:540:ARG:NH1	2.25	0.52
1:A:80:ASN:OD1	1:A:89:SER:N	2.43	0.51
1:B:602:PRO:O	1:B:606:ASN:OD1	2.29	0.51
1:B:209:PHE:CE2	1:B:235:ASN:OD1	2.64	0.51
1:B:374:TRP:CD1	1:B:508:ILE:HG23	2.47	0.50
1:B:605:MET:CE	1:B:611:GLU:O	2.60	0.49
1:A:619:GLU:OE2	1:A:625:LYS:NZ	2.39	0.49
1:A:104:ASP:OD2	1:A:110:GLY:N	2.45	0.48
1:B:479:ASN:OD1	1:B:480:GLY:N	2.47	0.48
1:B:453:ASN:OD1	1:B:455:SER:N	2.46	0.48
1:B:508:ILE:O	1:B:511:THR:OG1	2.25	0.48
1:B:362:MET:HE2	1:B:540:ARG:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:PHE:O	1:A:457:ARG:NH1	2.45	0.48
1:B:607:HIS:HB3	1:B:610:ILE:HG12	1.95	0.48
1:B:213:VAL:CG1	1:B:235:ASN:HD22	2.26	0.47
1:B:213:VAL:CG2	1:B:235:ASN:ND2	2.77	0.47
1:B:520:LEU:HD23	1:B:520:LEU:N	2.29	0.47
1:B:604:LEU:O	1:B:610:ILE:HD13	2.14	0.47
1:B:345:LEU:HD12	1:B:345:LEU:H	1.79	0.47
1:B:246:ARG:NH2	1:B:275:LYS:O	2.48	0.46
1:A:170:ARG:NH2	1:A:177:ASN:OD1	2.48	0.46
1:A:80:ASN:OD1	1:A:89:SER:CA	2.63	0.46
1:B:482:GLY:HA3	1:B:489:MET:HA	1.98	0.46
1:A:385:ASP:OD2	1:A:457:ARG:NH2	2.49	0.46
1:A:512:ARG:HH22	1:A:561:GLU:CD	2.14	0.46
1:B:605:MET:HA	1:B:610:ILE:CD1	2.45	0.46
1:A:147:ASN:OD1	1:A:152:ARG:NH2	2.41	0.46
1:A:614:THR:HG21	1:A:673:VAL:HG21	1.97	0.45
1:B:442:THR:HB	1:B:464:ASP:HB3	1.98	0.45
1:B:236:GLU:O	1:B:240:GLN:HG3	2.17	0.45
1:B:532:PHE:CE2	1:B:565:ARG:HG2	2.51	0.45
1:A:469:ILE:HD12	1:A:495:VAL:HG11	1.99	0.45
1:A:630:TYR:HD1	1:A:662:MET:HE1	1.81	0.45
1:B:3:ARG:NH1	1:B:43:GLN:O	2.50	0.45
1:A:37:ALA:HA	1:A:42:LEU:HD11	1.98	0.44
1:A:509:ASP:O	1:A:511:THR:HG23	2.18	0.44
1:B:262:LYS:HE2	1:B:391:SER:O	2.17	0.44
1:B:480:GLY:HA3	1:B:489:MET:SD	2.57	0.44
1:B:508:ILE:HG22	1:B:509:ASP:N	2.33	0.44
1:B:595:ILE:HG22	1:B:596:GLU:N	2.33	0.44
1:A:76:LEU:HD23	1:A:95:PRO:HB2	2.00	0.44
1:B:396:ASP:OD1	1:B:397:GLY:N	2.51	0.44
1:A:103:ILE:HG22	1:A:104:ASP:N	2.34	0.43
1:B:612:LEU:C	1:B:613:THR:HG23	2.39	0.43
1:B:73:HIS:O	1:B:75:SER:N	2.44	0.43
1:B:471:ARG:HD2	1:B:475:TYR:CE1	2.53	0.43
1:B:223:VAL:N	1:B:412:LEU:O	2.39	0.42
1:B:607:HIS:CG	1:B:608:PRO:HD2	2.55	0.42
1:A:604:LEU:HB3	1:A:613:THR:HG21	2.01	0.42
1:B:261:ASP:OD2	1:B:310:HIS:NE2	2.50	0.42
1:B:80:ASN:OD1	1:B:89:SER:CA	2.68	0.42
1:B:508:ILE:CG2	1:B:509:ASP:N	2.83	0.41
1:A:202:ASP:OD1	1:A:202:ASP:N	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:569:ARG:NH1	1:A:599:GLU:OE1	2.54	0.41
1:B:213:VAL:HG11	1:B:235:ASN:HD22	1.84	0.41
1:A:116:GLU:OE2	1:A:143:ARG:NH2	2.50	0.41
1:A:485:GLU:OE1	1:A:485:GLU:N	2.53	0.41
1:B:294:TYR:OH	1:B:545:ARG:NH2	2.54	0.41
1:B:604:LEU:O	1:B:610:ILE:CD1	2.69	0.41
1:A:70:ILE:HD12	1:A:70:ILE:N	2.36	0.41
1:A:69:ILE:HG22	1:A:103:ILE:HG13	2.03	0.41
1:B:114:VAL:HG11	1:B:144:GLU:HG3	2.03	0.41
1:A:239:ASN:OD1	1:A:275:LYS:HE3	2.21	0.40
1:B:38:GLN:N	1:B:38:GLN:OE1	2.54	0.40
1:B:481:TYR:C	1:B:489:MET:HG2	2.42	0.40
1:B:71:ASN:HB2	1:B:103:ILE:HD11	2.03	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)
1:A:637:ILE:O	1:B:682:GLN:NE2[3_765]	2.06	0.14
1:A:220:VAL:O	1:A:341:ARG:HH22[5_535]	1.51	0.09

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	681/776 (88%)	660 (97%)	20 (3%)	1 (0%)	51 81
1	B	683/776 (88%)	654 (96%)	28 (4%)	1 (0%)	51 81
All	All	1364/1552 (88%)	1314 (96%)	48 (4%)	2 (0%)	51 81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	488	ILE
1	A	488	ILE

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	583/666 (88%)	583 (100%)	0	100	100
1	B	583/666 (88%)	578 (99%)	5 (1%)	78	94
All	All	1166/1332 (88%)	1161 (100%)	5 (0%)	91	97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	80	ASN
1	B	442	THR
1	B	472	ASN
1	B	606	ASN
1	B	612	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	235	ASN
1	B	450	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	801	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	B	801	-	4,4,4	0.13	0	6,6,6	0.07	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	683/776 (88%)	0.48	35 (5%) 28 19	73, 117, 177, 194	0
1	B	685/776 (88%)	0.71	69 (10%) 7 4	86, 148, 192, 233	0
All	All	1368/1552 (88%)	0.59	104 (7%) 13 7	73, 139, 183, 233	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	ARG	6.3
1	B	655	ILE	6.1
1	B	646	LEU	5.9
1	B	589	LYS	5.8
1	A	37	ALA	5.3
1	B	659	PHE	5.0
1	B	197	GLY	4.8
1	B	654	MET	4.6
1	B	593	TYR	4.6
1	B	652	ASP	4.4
1	B	656	PRO	4.4
1	B	590	ILE	4.3
1	B	657	THR	4.3
1	A	42	LEU	4.2
1	B	650	LEU	4.1
1	B	592	GLY	4.0
1	B	591	ARG	3.8
1	A	57	LEU	3.8
1	B	624	ARG	3.7
1	B	603	PHE	3.6
1	B	584	LEU	3.6
1	A	196	ARG	3.6
1	B	595	ILE	3.5
1	B	636	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	599	GLU	3.4
1	B	649	ASP	3.3
1	B	609	LEU	3.3
1	B	634	PRO	3.3
1	B	196	ARG	3.3
1	B	13	LYS	3.3
1	B	642	LEU	3.2
1	A	683	ALA	3.2
1	B	631	TYR	3.2
1	A	34	LYS	3.2
1	A	48	ILE	3.1
1	A	654	MET	3.1
1	A	197	GLY	3.1
1	A	226	GLY	3.1
1	B	629	GLY	3.1
1	A	620	GLN	3.1
1	A	627	LEU	3.0
1	B	176	LEU	3.0
1	B	522	LEU	3.0
1	A	10	PHE	3.0
1	B	235	ASN	3.0
1	B	604	LEU	3.0
1	B	620	GLN	2.9
1	A	609	LEU	2.9
1	A	36	HIS	2.9
1	B	651	PRO	2.9
1	B	600	ILE	2.9
1	A	228	SER	2.8
1	A	32	GLN	2.8
1	B	198	GLU	2.8
1	B	199	LYS	2.7
1	A	17	VAL	2.7
1	A	35	VAL	2.7
1	B	541	GLY	2.7
1	A	372	CYS	2.7
1	B	484	THR	2.7
1	B	625	LYS	2.7
1	B	338	HIS	2.7
1	A	4	ILE	2.6
1	B	653	TYR	2.6
1	A	126	LEU	2.6
1	B	637	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	596	GLU	2.5
1	B	419	ARG	2.5
1	A	38	GLN	2.5
1	A	61	ILE	2.4
1	A	13	LYS	2.4
1	A	60	GLU	2.4
1	A	486	CYS	2.4
1	A	68	ARG	2.4
1	B	635	GLU	2.3
1	B	645	TRP	2.3
1	B	456	PHE	2.3
1	B	663	LYS	2.2
1	B	208	LEU	2.2
1	B	514	LEU	2.2
1	B	201	ILE	2.2
1	B	289	GLY	2.2
1	B	565	ARG	2.2
1	B	588	VAL	2.2
1	B	304	LEU	2.1
1	A	18	VAL	2.1
1	B	55	TYR	2.1
1	A	626	TYR	2.1
1	B	626	TYR	2.1
1	A	632	VAL	2.1
1	B	388	LEU	2.1
1	B	648	ASN	2.1
1	B	18	VAL	2.1
1	A	550	ALA	2.1
1	A	290	GLU	2.1
1	B	259	MET	2.1
1	B	628	VAL	2.1
1	B	317	ALA	2.0
1	B	393	PHE	2.0
1	B	676	ARG	2.0
1	B	438	SER	2.0
1	B	318	CYS	2.0
1	A	33	GLU	2.0
1	B	623	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	B	801	5/5	0.90	0.14	175,176,186,186	0
2	SO4	A	801	5/5	0.91	0.16	143,151,164,166	0

6.5 Other polymers [i](#)

There are no such residues in this entry.