



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:00 PM UTC

PDB ID : 6IGB / pdb_00006igb
Title : the structure of Pseudomonas aeruginosa Periplasmic gluconolactonase, PpgL
Authors : Song, Y.J.; Shen, Y.L.; Wang, K.L.; Li, T.; Zhu, Y.B.; Li, C.C.; He, L.H.;
Zhao, N.L.; Zhao, C.; Yang, J.; Huang, Q.; Mu, X.Y.
Deposited on : 2018-09-25
Resolution : 1.65 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

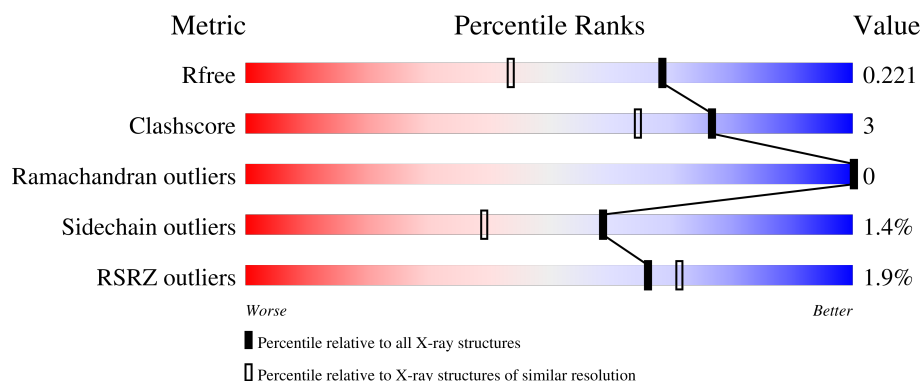
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	364	0% 73% 24% .
1	B	364	2% 83% 15% .

2 Entry composition [i](#)

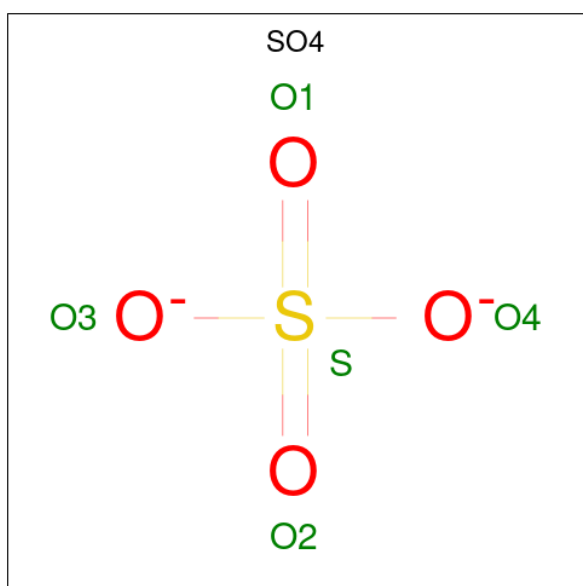
There are 4 unique types of molecules in this entry. The entry contains 5859 atoms, of which 6 are hydrogens and 0 are deuteriums.

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 Periplasmic gluconolactonase, PpgL.

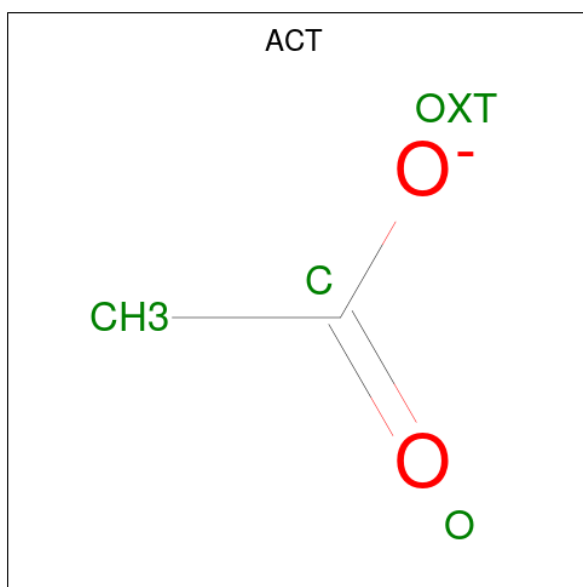
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2765	1734	503	527	1			
1	B	363	Total	C	N	O	S	0	0	0
			2765	1734	503	527	1			

- 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		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	B	1	Total	C	H	O	0	0
			7	2	3	2		

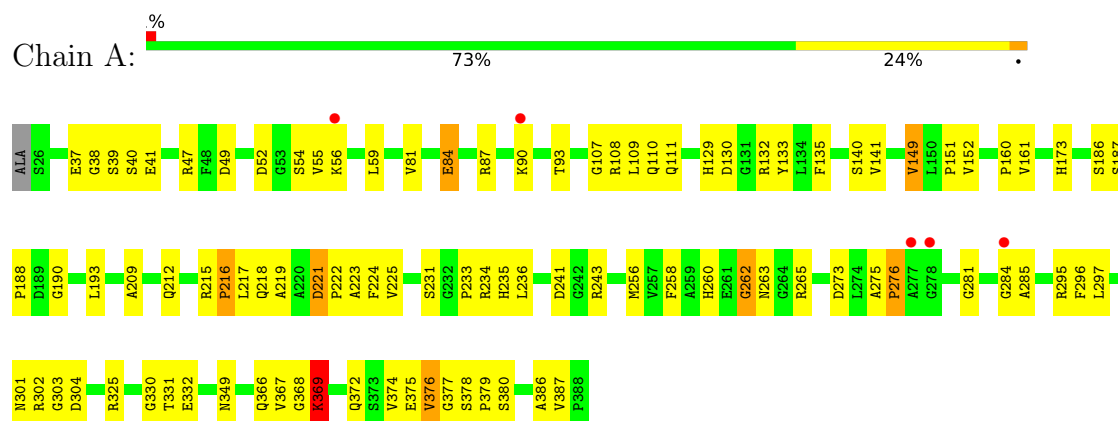
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	170	Total	O	0	0
			170	170		
4	B	135	Total	O	0	0
			135	135		

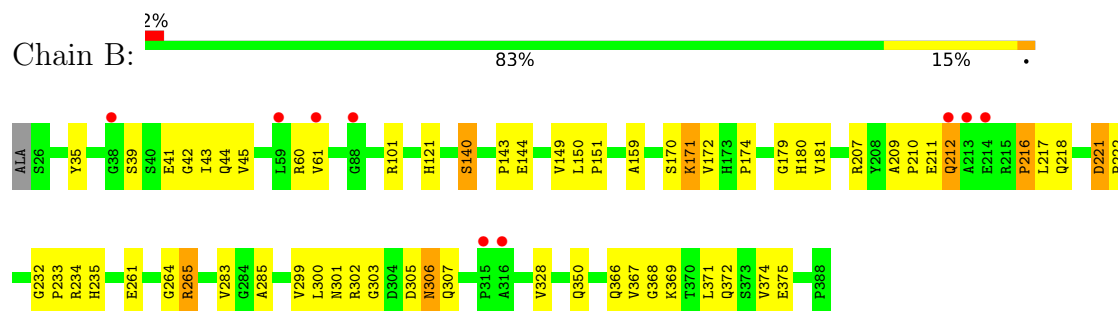
3 Residue-property plots

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: Periplasmic gluconolactonase, PpgL



- Molecule 1: Periplasmic gluconolactonase, PpgL



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.76Å 91.76Å 170.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.70 – 1.65 38.70 – 1.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (38.70-1.65) 100.0 (38.70-1.65)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 1.65Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.196 , 0.222 0.197 , 0.221	Depositor DCC
R_{free} test set	2000 reflections (2.27%)	wwPDB-VP
Wilson B-factor (Å ²)	15.9	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5859	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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: ACT, 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	1.49	74/2832 (2.6%)	0.99	11/3852 (0.3%)
1	B	1.13	39/2832 (1.4%)	0.85	10/3852 (0.3%)
All	All	1.32	113/5664 (2.0%)	0.93	21/7704 (0.3%)

All (113) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	378	SER	C-O	-15.70	1.16	1.23
1	B	216	PRO	N-CD	11.38	1.63	1.47
1	A	332	GLU	C-O	-10.26	1.19	1.23
1	B	179	GLY	C-O	-8.44	1.14	1.23
1	A	81	VAL	C-O	-8.16	1.16	1.24
1	A	331	THR	C-O	-7.94	1.13	1.23
1	B	235	HIS	C-O	-7.94	1.14	1.24
1	A	215	ARG	C-O	-7.85	1.14	1.24
1	B	374	VAL	C-O	-7.77	1.16	1.24
1	B	299	VAL	C-O	-7.72	1.16	1.24
1	B	172	VAL	C-O	-7.64	1.15	1.24
1	A	366	GLN	C-O	-7.39	1.14	1.23
1	A	235	HIS	C-O	-7.36	1.14	1.24
1	B	305	ASP	C-O	-7.20	1.16	1.24
1	A	108	ARG	C-O	-7.13	1.15	1.23
1	B	150	LEU	C-O	-7.12	1.16	1.24
1	A	377	GLY	C-O	-7.10	1.18	1.24
1	A	187	SER	C-O	-7.08	1.14	1.23
1	A	302	ARG	C-O	-7.08	1.15	1.24
1	A	330	GLY	C-O	-6.98	1.15	1.23
1	A	110	GLN	C-O	-6.92	1.16	1.24
1	B	367	VAL	C-O	-6.90	1.16	1.24
1	A	374	VAL	C-O	-6.87	1.16	1.24
1	A	39	SER	C-O	-6.84	1.15	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	216	PRO	C-O	-6.70	1.15	1.24
1	A	209	ALA	C-O	-6.67	1.16	1.24
1	A	107	GLY	C-O	-6.67	1.14	1.24
1	B	302	ARG	C-O	-6.62	1.15	1.24
1	B	180	HIS	C-O	-6.61	1.16	1.23
1	A	152	VAL	C-O	-6.57	1.17	1.24
1	A	217	LEU	C-O	-6.53	1.16	1.24
1	B	375	GLU	C-O	-6.50	1.16	1.24
1	A	219	ALA	C-O	-6.48	1.16	1.23
1	B	233	PRO	C-O	-6.48	1.16	1.23
1	B	181	VAL	C-O	-6.44	1.16	1.24
1	A	186	SER	C-O	-6.39	1.15	1.23
1	A	303	GLY	C-O	-6.37	1.16	1.24
1	B	307	GLN	C-O	-6.31	1.15	1.23
1	A	231	SER	C-O	-6.31	1.15	1.24
1	A	233	PRO	C-O	-6.31	1.16	1.23
1	A	218	GLN	C-O	-6.31	1.16	1.24
1	A	173	HIS	C-O	-6.30	1.15	1.24
1	A	133	TYR	C-O	-6.29	1.16	1.23
1	A	225	VAL	C-O	-6.29	1.17	1.23
1	B	372	GLN	C-O	-6.29	1.16	1.23
1	A	149	VAL	C-O	-6.28	1.17	1.24
1	B	207	ARG	C-O	-6.25	1.16	1.24
1	B	234	ARG	C-O	-6.21	1.16	1.24
1	A	379	PRO	C-O	-6.18	1.17	1.23
1	B	302	ARG	N-CA	-6.09	1.42	1.46
1	B	218	GLN	C-O	-6.08	1.16	1.24
1	A	151	PRO	C-O	-6.01	1.17	1.23
1	A	376	VAL	C-O	-5.99	1.17	1.23
1	B	232	GLY	C-O	-5.99	1.15	1.24
1	A	47	ARG	C-O	-5.98	1.16	1.24
1	A	212	GLN	C-O	-5.98	1.15	1.23
1	A	188	PRO	C-O	-5.97	1.16	1.24
1	A	380	SER	C-O	-5.91	1.16	1.24
1	A	38	GLY	C-O	-5.90	1.16	1.23
1	A	302	ARG	N-CA	-5.90	1.42	1.46
1	A	372	GLN	C-O	-5.83	1.16	1.23
1	B	368	GLY	C-O	-5.81	1.16	1.23
1	A	93	THR	C-O	-5.77	1.16	1.24
1	A	111	GLN	C-O	-5.75	1.17	1.23
1	B	42	GLY	C-O	-5.75	1.16	1.23
1	A	160	PRO	C-O	-5.72	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	234	ARG	C-O	-5.70	1.16	1.23
1	B	301	ASN	C-O	-5.70	1.17	1.24
1	A	297	LEU	C-O	-5.69	1.17	1.24
1	B	209	ALA	CA-C	-5.67	1.45	1.52
1	A	276	PRO	C-O	-5.63	1.16	1.23
1	A	281	GLY	C-O	-5.60	1.16	1.23
1	B	366	GLN	C-O	-5.59	1.17	1.23
1	B	217	LEU	C-O	-5.58	1.17	1.24
1	A	295	ARG	C-O	-5.58	1.16	1.24
1	A	296	PHE	C-O	-5.57	1.16	1.23
1	B	303	GLY	C-O	-5.55	1.17	1.24
1	A	224	PHE	C-O	-5.54	1.17	1.23
1	A	369	LYS	C-O	-5.54	1.17	1.23
1	A	273	ASP	C-O	-5.51	1.16	1.23
1	A	236	LEU	C-O	-5.49	1.17	1.24
1	A	90	LYS	CA-C	-5.49	1.46	1.52
1	B	369	LYS	C-O	-5.48	1.17	1.23
1	A	129	HIS	C-O	-5.45	1.17	1.24
1	A	375	GLU	C-O	-5.44	1.17	1.24
1	A	367	VAL	C-O	-5.43	1.18	1.24
1	A	223	ALA	C-O	-5.42	1.17	1.24
1	B	265	ARG	C-O	-5.41	1.17	1.24
1	A	52	ASP	C-O	-5.40	1.17	1.24
1	A	84	GLU	C-O	-5.39	1.17	1.23
1	A	59	LEU	C-O	-5.37	1.17	1.24
1	A	387	VAL	C-O	-5.33	1.17	1.24
1	B	300	LEU	C-O	-5.32	1.17	1.23
1	B	371	LEU	C-O	-5.29	1.17	1.24
1	A	304	ASP	C-O	-5.27	1.16	1.23
1	A	349	ASN	C-O	-5.27	1.17	1.23
1	B	45	VAL	C-O	-5.26	1.18	1.24
1	A	190	GLY	C-O	-5.26	1.17	1.24
1	B	43	ILE	C-O	-5.25	1.18	1.24
1	A	40	SER	C-O	-5.22	1.17	1.23
1	A	55	VAL	C-O	-5.21	1.18	1.24
1	B	264	GLY	C-O	-5.21	1.16	1.23
1	A	386	ALA	C-O	-5.19	1.17	1.23
1	B	261	GLU	C-O	-5.19	1.17	1.24
1	B	174	PRO	N-CD	5.17	1.54	1.47
1	A	368	GLY	C-O	-5.14	1.17	1.23
1	A	109	LEU	C-O	-5.14	1.17	1.24
1	A	161	VAL	C-O	-5.13	1.19	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	209	ALA	C-O	-5.13	1.18	1.24
1	B	221	ASP	C-O	-5.07	1.21	1.25
1	A	54	SER	C-O	-5.05	1.17	1.23
1	A	49	ASP	C-O	-5.04	1.17	1.24
1	A	260	HIS	C-O	-5.01	1.18	1.24

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	LYS	N-CA-C	-7.51	94.87	108.02
1	B	302	ARG	N-CA-C	6.90	114.44	108.78
1	A	221	ASP	C-N-CD	6.52	134.94	120.60
1	A	87	ARG	N-CA-C	6.32	118.98	111.71
1	B	306	ASN	N-CA-C	6.29	120.33	111.52
1	B	216	PRO	CA-N-CD	-6.21	103.31	112.00
1	B	221	ASP	C-N-CD	6.17	134.19	120.60
1	B	211	GLU	N-CA-C	-6.13	105.90	113.19
1	A	285	ALA	N-CA-C	-5.96	99.99	109.76
1	A	275	ALA	CA-C-N	-5.89	114.55	120.21
1	A	275	ALA	C-N-CA	-5.89	114.55	120.21
1	B	307	GLN	N-CA-C	5.68	118.33	109.52
1	A	187	SER	N-CA-C	-5.35	102.50	110.20
1	A	262	GLY	N-CA-C	-5.30	104.97	112.81
1	A	129	HIS	N-CA-C	5.28	118.97	112.54
1	B	140	SER	N-CA-C	5.24	121.96	110.80
1	B	39	SER	N-CA-C	5.22	119.49	113.18
1	A	378	SER	O-C-N	5.16	123.90	121.53
1	B	285	ALA	N-CA-C	-5.13	102.89	110.48
1	A	295	ARG	N-CA-C	5.11	119.57	113.23
1	B	283	VAL	CB-CA-C	-5.10	102.54	110.69

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	2765	0	2679	17	0
1	B	2765	0	2679	15	1
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	4	3	3	1	0
3	B	4	3	3	0	0
4	A	170	0	0	4	1
4	B	135	0	0	0	0
All	All	5853	6	5364	32	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 3.

All (32) 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:A:37:GLU:HG2	4:A:538:HOH:O	1.83	0.78
1:A:37:GLU:CG	4:A:538:HOH:O	2.40	0.66
1:B:41:GLU:HG2	1:B:61:VAL:HG11	1.83	0.59
1:B:210:PRO:C	1:B:212:GLN:H	2.10	0.59
1:B:151:PRO:HB2	1:B:159:ALA:HB3	1.89	0.54
1:B:41:GLU:CG	1:B:61:VAL:HG11	2.38	0.54
1:A:284:GLY:O	1:A:301:ASN:HA	2.08	0.54
1:B:149:VAL:HG11	1:B:216:PRO:HG2	1.92	0.52
1:B:210:PRO:C	1:B:212:GLN:N	2.67	0.52
1:B:306:ASN:O	1:B:328:VAL:HG22	2.11	0.49
1:A:276:PRO:HD3	1:A:325:ARG:CZ	2.45	0.47
1:B:41:GLU:OE2	1:B:41:GLU:N	2.48	0.46
1:B:143:PRO:HA	1:B:144:GLU:HA	1.78	0.46
1:B:121:HIS:HB2	1:B:140:SER:H	1.81	0.46
1:A:149:VAL:HG11	1:A:216:PRO:HG2	1.99	0.43
1:A:284:GLY:O	1:A:301:ASN:OD1	2.36	0.43
1:A:221:ASP:HA	1:A:222:PRO:C	2.43	0.43
1:B:170:SER:C	1:B:171:LYS:HG2	2.44	0.43
1:A:256:MET:HE2	1:A:258:PHE:CZ	2.54	0.43
1:B:35:TYR:CE2	1:B:350:GLN:NE2	2.88	0.42
1:A:135:PHE:CD2	1:A:193:LEU:HD13	2.54	0.42
1:B:151:PRO:HD2	1:B:159:ALA:HB3	2.02	0.42
1:A:140:SER:O	1:A:141:VAL:C	2.63	0.42
1:A:130:ASP:OD2	1:A:132:ARG:HD2	2.19	0.41
1:A:262:GLY:O	1:A:263:ASN:HB2	2.20	0.41
1:B:221:ASP:HA	1:B:222:PRO:HA	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ASP:OD2	1:A:243:ARG:HD3	2.21	0.41
1:A:84:GLU:OE2	3:A:402:ACT:H1	2.21	0.41
1:B:35:TYR:CD2	1:B:350:GLN:NE2	2.89	0.41
1:A:37:GLU:HG3	4:A:538:HOH:O	2.17	0.40
1:A:222:PRO:HD2	1:A:265:ARG:HG2	2.03	0.40
1:A:369:LYS:HG3	4:A:628:HOH:O	2.22	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:B:60:ARG:NH1	1:B:265:ARG:NH1[5_455]	2.05	0.15
4:A:583:HOH:O	4:A:618:HOH:O[8_554]	2.14	0.06

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	361/364 (99%)	354 (98%)	7 (2%)	0	100	100
1	B	361/364 (99%)	351 (97%)	10 (3%)	0	100	100
All	All	722/728 (99%)	705 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	294/294 (100%)	290 (99%)	4 (1%)	59	39
1	B	294/294 (100%)	290 (99%)	4 (1%)	59	39
All	All	588/588 (100%)	580 (99%)	8 (1%)	59	39

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLU
1	A	56	LYS
1	A	369	LYS
1	A	376	VAL
1	B	44	GLN
1	B	101	ARG
1	B	171	LYS
1	B	212	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	HIS
1	A	83	ASN
1	A	114	GLN
1	A	116	GLN
1	A	129	HIS
1	B	83	ASN
1	B	114	GLN
1	B	350	GLN
1	B	366	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACT	A	402	-	3,3,3	1.48	0	3,3,3	1.60	1 (33%)
2	SO4	A	401	-	4,4,4	0.21	0	6,6,6	0.25	0
3	ACT	B	402	-	3,3,3	1.14	0	3,3,3	3.51	2 (66%)
2	SO4	B	401	-	4,4,4	0.29	0	6,6,6	0.33	0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	B	402	ACT	O-C-CH3	-4.60	103.68	122.53
3	B	402	ACT	OXT-C-CH3	3.90	131.41	115.05
3	A	402	ACT	OXT-C-O	-2.40	113.13	122.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	ACT	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	363/364 (99%)	-0.33	5 (1%) 73 78	9, 16, 28, 46	0
1	B	363/364 (99%)	-0.00	9 (2%) 58 64	11, 20, 38, 46	0
All	All	726/728 (99%)	-0.17	14 (1%) 66 71	9, 18, 36, 46	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	284	GLY	3.9
1	A	56	LYS	3.3
1	B	61	VAL	3.1
1	A	277	ALA	2.9
1	B	316	ALA	2.9
1	B	213	ALA	2.8
1	B	88	GLY	2.3
1	A	90	LYS	2.3
1	B	214	GLU	2.2
1	B	315	PRO	2.2
1	B	59	LEU	2.1
1	A	278	GLY	2.1
1	B	212	GLN	2.1
1	B	38	GLY	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
3	ACT	A	402	4/4	0.80	0.14	21,25,31,42	0
3	ACT	B	402	4/4	0.80	0.19	17,31,33,44	0
2	SO4	A	401	5/5	0.98	0.10	19,19,27,38	0
2	SO4	B	401	5/5	0.98	0.07	20,28,35,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.