



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2024 – 07:42 PM EST

PDB ID : 4G65
Title : Potassium transporter peripheral membrane component (trkA) from *Vibrio vulnificus*
Authors : Filippova, E.V.; Wawrzak, Z.; Skarina, T.; Onopriyenko, O.; Edwards, A.; Savchenko, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2012-07-18
Resolution : 2.09 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

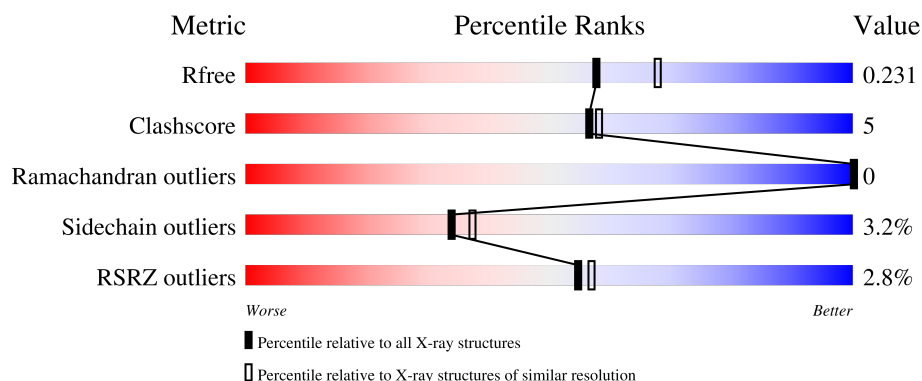
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

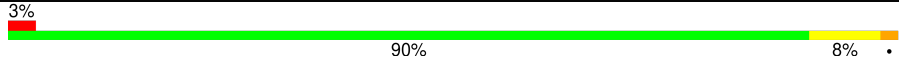
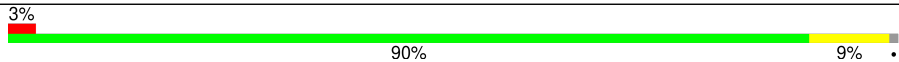
The reported resolution of this entry is 2.09 Å.

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}	164625	6234 (2.10-2.10)
Clashscore	180529	6893 (2.10-2.10)
Ramachandran outliers	177936	6839 (2.10-2.10)
Sidechain outliers	177891	6840 (2.10-2.10)
RSRZ outliers	164620	6234 (2.10-2.10)

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	461	
1	B	461	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7583 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.

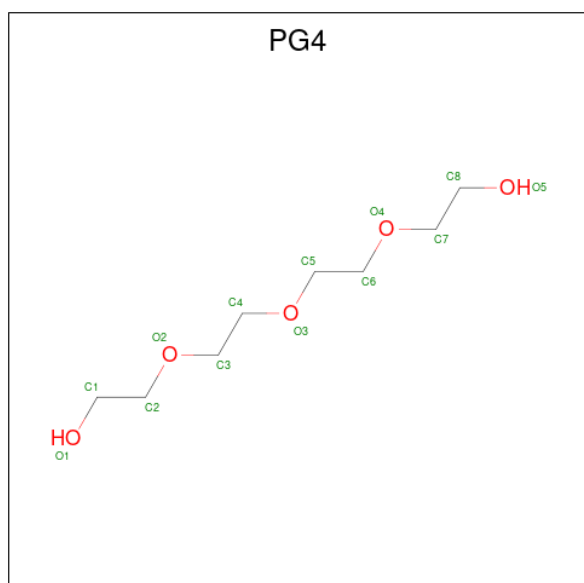
- Molecule 1 is a protein called Trk system potassium uptake protein trkA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	459	Total	C	N	O	S	Se	0	8	0
			3583	2246	633	690	2	12			
1	B	458	Total	C	N	O	S	Se	0	2	0
			3533	2218	622	680	2	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q8DDE6
A	-1	ASN	-	expression tag	UNP Q8DDE6
A	0	ALA	-	expression tag	UNP Q8DDE6
B	-2	SER	-	expression tag	UNP Q8DDE6
B	-1	ASN	-	expression tag	UNP Q8DDE6
B	0	ALA	-	expression tag	UNP Q8DDE6

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		

- 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	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	Cl	0	0
			7	7		
4	B	7	Total	Cl	0	0
			7	7		

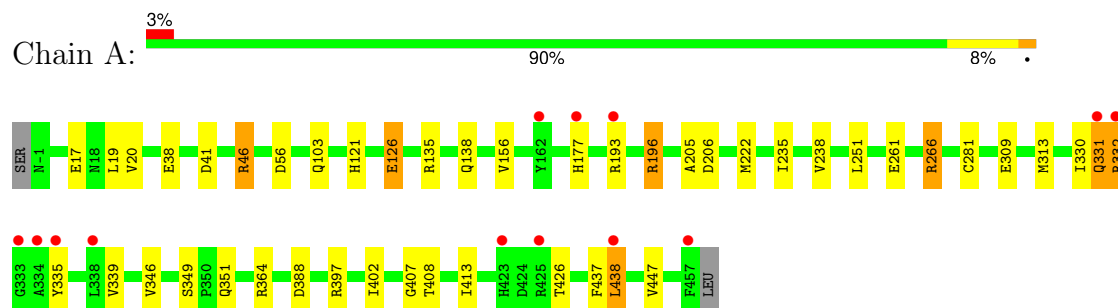
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	234	Total	O	0	6
			240	240		
5	B	192	Total	O	0	4
			196	196		

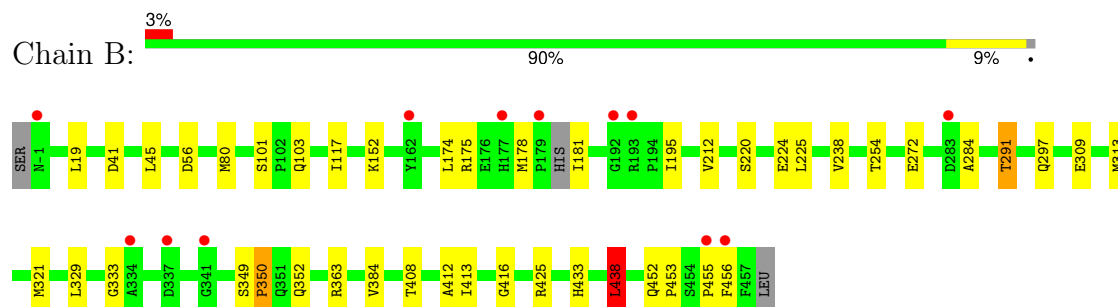
3 Residue-property plots [i](#)

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: Trk system potassium uptake protein trkA



- Molecule 1: Trk system potassium uptake protein trkA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	68.11Å 225.26Å 139.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.67 – 2.09 29.67 – 2.09	Depositor EDS
% Data completeness (in resolution range)	97.5 (29.67-2.09) 97.5 (29.67-2.09)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.179 , 0.228 0.186 , 0.231	Depositor DCC
R_{free} test set	3155 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7583	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

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

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, PG4, 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	0.76	1/3621 (0.0%)	0.88	6/4890 (0.1%)
1	B	0.66	0/3571	0.81	2/4823 (0.0%)
All	All	0.71	1/7192 (0.0%)	0.85	8/9713 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	GLU	CD-OE2	-5.17	1.20	1.25

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	388	ASP	CB-CG-OD1	5.61	123.35	118.30
1	A	388	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	46	ARG	NE-CZ-NH2	5.33	122.96	120.30
1	A	135	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	266	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	56	ASP	CB-CG-OD1	5.09	122.88	118.30
1	B	438	LEU	CA-CB-CG	5.03	126.88	115.30

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	3583	0	3619	36	0
1	B	3533	0	3573	28	0
2	A	7	0	9	4	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	7	0	0	0	0
4	B	7	0	0	1	0
5	A	240	0	0	3	0
5	B	196	0	0	0	0
All	All	7583	0	7201	65	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 5.

All (65) 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:437:PHE:C	1:A:438:LEU:HD23	1.82	0.98
1:B:80:MSE:CE	1:B:117:ILE:HD13	2.06	0.86
1:A:438:LEU:HD23	1:A:438:LEU:N	1.96	0.80
1:A:313[A]:MSE:HE3	1:A:313[A]:MSE:HA	1.67	0.76
1:A:156[B]:VAL:HG11	1:A:222:MSE:SE	2.36	0.76
1:B:80:MSE:HE2	1:B:117:ILE:HD13	1.68	0.75
1:A:332:ARG:NH1	5:A:817:HOH:O	2.22	0.72
1:A:438:LEU:HD11	1:A:447:VAL:HG21	1.73	0.68
1:A:46:ARG:HH21	1:A:46:ARG:HG3	1.60	0.66
1:B:408:THR:HG23	1:B:438:LEU:HD22	1.80	0.63
1:B:333:GLY:HA3	1:B:352:GLN:HG3	1.81	0.62
1:B:45:LEU:HD12	1:B:45:LEU:C	2.21	0.61
1:A:413:ILE:HD12	1:A:426:THR:HG21	1.83	0.61
1:B:384:VAL:HG22	1:B:433:HIS:CE1	2.36	0.60
1:B:333:GLY:CA	1:B:352:GLN:HG3	2.32	0.59
1:A:332:ARG:H	1:A:332:ARG:HD2	1.68	0.58
1:A:438:LEU:N	1:A:438:LEU:CD2	2.68	0.56
1:B:220:SER:O	1:B:224:GLU:HG2	2.05	0.56
1:A:103:GLN:NE2	5:A:780:HOH:O	2.40	0.55
1:A:438:LEU:HD11	1:A:447:VAL:CG2	2.37	0.54
1:B:212:VAL:HG11	1:B:412:ALA:HB2	1.89	0.54
1:B:152:LYS:HE3	1:B:416:GLY:O	2.08	0.54
1:A:332:ARG:HG2	1:A:335:TYR:CD1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:TYR:O	1:A:339:VAL:HG23	2.08	0.53
1:A:266:ARG:HG3	1:A:266:ARG:HH11	1.74	0.53
1:A:332:ARG:CG	1:A:335:TYR:CD1	2.95	0.50
1:A:332:ARG:HD2	1:A:332:ARG:N	2.27	0.49
1:A:126:GLU:OE2	5:A:786:HOH:O	2.20	0.49
1:B:291:THR:HG23	1:B:321:MSE:HE2	1.93	0.49
1:B:45:LEU:HD12	1:B:45:LEU:O	2.13	0.49
1:B:291:THR:HG23	1:B:321:MSE:CE	2.43	0.49
1:B:384:VAL:HG22	1:B:433:HIS:ND1	2.28	0.49
1:B:309:GLU:OE1	1:B:313:MSE:HE1	2.12	0.49
1:A:205:ALA:O	1:A:206:ASP:HB2	2.13	0.49
1:A:309:GLU:H	2:A:501:PG4:H41	1.77	0.49
1:A:332:ARG:HG2	1:A:335:TYR:CG	2.48	0.49
1:A:331:GLN:NE2	1:A:331:GLN:HA	2.29	0.48
1:B:333:GLY:HA3	1:B:352:GLN:CG	2.43	0.48
1:A:309:GLU:N	2:A:501:PG4:H41	2.28	0.48
2:A:501:PG4:H32	2:A:501:PG4:O1	2.13	0.48
1:B:352:GLN:HG2	4:B:505:CL:CL	2.50	0.47
1:B:453:PRO:HB2	1:B:455:PRO:HD2	1.97	0.47
1:A:397:ARG:HB2	1:A:402:ILE:HD11	1.95	0.47
1:B:363:ARG:NH2	1:B:384:VAL:HG21	2.30	0.46
1:B:452:GLN:HG3	1:B:456:PHE:CD2	2.50	0.46
1:B:238:VAL:HG11	1:B:284:ALA:HB1	1.97	0.46
1:A:103:GLN:HG2	1:B:56:ASP:OD1	2.17	0.44
1:B:174:LEU:HD22	1:B:178:MSE:HE2	1.99	0.44
1:B:178:MSE:HE1	1:B:225:LEU:HD23	2.00	0.44
1:A:17:GLU:O	1:A:20:VAL:HG22	2.18	0.43
2:A:501:PG4:O1	2:A:501:PG4:C3	2.65	0.43
1:B:178:MSE:HE1	1:B:225:LEU:CD2	2.49	0.43
1:A:235[A]:ILE:HD13	1:A:251:LEU:HD13	2.01	0.42
1:A:261:GLU:O	1:A:281:CYS:HA	2.19	0.42
1:A:408:THR:HG23	1:A:438:LEU:HD22	2.02	0.42
1:A:313[A]:MSE:HE3	1:A:313[A]:MSE:CA	2.42	0.42
1:B:329:LEU:HD11	1:B:350:PRO:HD3	2.01	0.42
1:B:101[B]:SER:OG	1:B:103:GLN:OE1	2.35	0.41
1:A:196:ARG:HG3	1:A:407:GLY:HA2	2.03	0.41
1:A:121:HIS:CE1	1:A:364:ARG:HH22	2.39	0.41
1:A:331:GLN:O	1:A:351:GLN:OE1	2.39	0.41
1:A:330:ILE:O	1:A:349:SER:HB2	2.21	0.41
1:A:332:ARG:H	1:A:332:ARG:CD	2.32	0.41
1:B:349:SER:HB3	1:B:352:GLN:HB2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:PHE:C	1:A:438:LEU:CD2	2.72	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	465/461 (101%)	454 (98%)	11 (2%)	0	100	100
1	B	456/461 (99%)	449 (98%)	7 (2%)	0	100	100
All	All	921/922 (100%)	903 (98%)	18 (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	386/369 (105%)	373 (97%)	13 (3%)	32	35
1	B	379/369 (103%)	367 (97%)	12 (3%)	34	37
All	All	765/738 (104%)	740 (97%)	25 (3%)	34	36

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	38	GLU
1	A	41	ASP
1	A	138[A]	GLN
1	A	138[B]	GLN
1	A	177	HIS
1	A	193	ARG
1	A	196	ARG
1	A	238	VAL
1	A	331	GLN
1	A	332	ARG
1	A	346	VAL
1	A	438	LEU
1	B	19	LEU
1	B	41	ASP
1	B	181	ILE
1	B	195	ILE
1	B	254	THR
1	B	272	GLU
1	B	291	THR
1	B	297	GLN
1	B	350	PRO
1	B	413	ILE
1	B	425	ARG
1	B	438	LEU

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

Mol	Chain	Res	Type
1	A	352	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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 17 ligands modelled in this entry, 14 are monoatomic - leaving 3 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	$\# Z > 2$
3	SO4	B	501	-	4,4,4	0.40	0	6,6,6	0.55	0
2	PG4	A	501	-	6,6,12	0.43	0	5,5,11	0.31	0
3	SO4	A	502	-	4,4,4	0.38	0	6,6,6	0.76	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	PG4	A	501	-	-	3/4/4/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	PG4	C1-C2-O2-C3
2	A	501	PG4	O1-C1-C2-O2
2	A	501	PG4	O2-C3-C4-O3

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	PG4	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	448/461 (97%)	-0.16	13 (2%) 54 55	6, 23, 47, 70	7 (1%)
1	B	447/461 (96%)	-0.10	12 (2%) 56 58	10, 29, 50, 79	2 (0%)
All	All	895/922 (97%)	-0.13	25 (2%) 55 57	6, 26, 49, 79	9 (1%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162[A]	TYR	5.4
1	A	457	PHE	4.6
1	B	455	PRO	3.9
1	B	179	PRO	3.8
1	A	335	TYR	3.4
1	A	162	TYR	3.3
1	B	456	PHE	3.3
1	A	338	LEU	3.0
1	A	332	ARG	2.8
1	A	177	HIS	2.8
1	A	331	GLN	2.5
1	A	425	ARG	2.5
1	A	423	HIS	2.4
1	B	177	HIS	2.4
1	B	192	GLY	2.3
1	A	333	GLY	2.3
1	B	193	ARG	2.2
1	A	334	ALA	2.2
1	B	337	ASP	2.2
1	B	334	ALA	2.2
1	A	438	LEU	2.1
1	B	283	ASP	2.1
1	A	193	ARG	2.1
1	B	-1	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	341	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 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	PG4	A	501	7/13	0.68	0.17	35,46,49,50	0
4	CL	A	505	1/1	0.84	0.15	48,48,48,48	0
4	CL	B	508	1/1	0.89	0.14	59,59,59,59	0
4	CL	B	505	1/1	0.90	0.10	49,49,49,49	0
4	CL	A	506	1/1	0.90	0.11	51,51,51,51	0
4	CL	B	506	1/1	0.93	0.08	47,47,47,47	0
4	CL	A	509	1/1	0.93	0.07	46,46,46,46	0
4	CL	B	504	1/1	0.94	0.08	29,29,29,29	1
4	CL	A	504	1/1	0.94	0.09	29,29,29,29	0
3	SO4	A	502	5/5	0.97	0.10	27,28,32,35	0
4	CL	A	507	1/1	0.97	0.08	30,30,30,30	0
3	SO4	B	501	5/5	0.97	0.07	27,30,32,36	0
4	CL	B	502	1/1	0.97	0.06	33,33,33,33	0
4	CL	A	503	1/1	0.98	0.04	36,36,36,36	0
4	CL	B	507	1/1	0.98	0.03	34,34,34,34	0
4	CL	A	508	1/1	0.98	0.07	28,28,28,28	0
4	CL	B	503	1/1	0.99	0.05	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.