



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 07:31 AM GMT

PDB ID : 4L9O
Title : Crystal Structure of the Sec13-Sec16 blade-inserted complex from *Pichia pastoris*
Authors : McMahon, C.; Jeffrey, P.D.; Hughson, F.M.
Deposited on : 2013-06-18
Resolution : 1.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

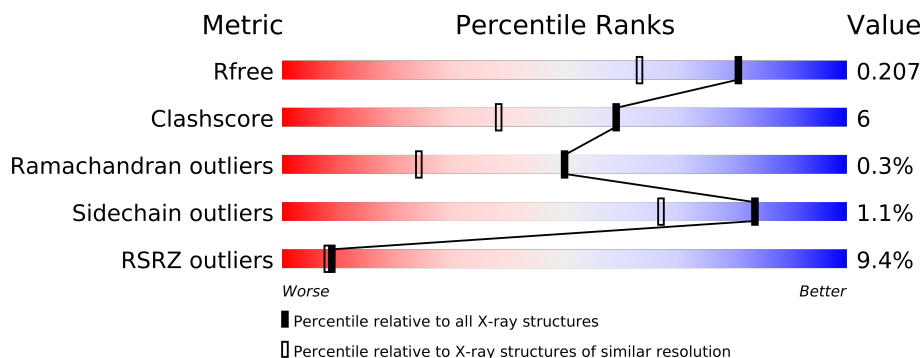
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.60 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}	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	349	
1	B	349	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	EDO	A	2408	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10256 atoms, of which 4900 are hydrogens and 0 are deuterium.

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 Protein transport protein SEC13, COPII coat assembly protein SEC16 fusion protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	330	Total	C	H	N	O	S	0	0	0
			5097	1649	2508	453	479	8			
1	B	311	Total	C	H	N	O	S	0	0	0
			4809	1561	2366	424	451	7			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1026	GLY	-	EXPRESSION TAG	UNP P53024
A	1027	SER	-	EXPRESSION TAG	UNP P53024
A	1028	HIS	-	EXPRESSION TAG	UNP P53024
A	1029	MET	-	EXPRESSION TAG	UNP P53024
A	1077	GLY	-	LINKER	UNP P53024
A	1078	SER	-	LINKER	UNP P53024
A	1079	GLY	-	LINKER	UNP P53024
A	1080	GLY	-	LINKER	UNP P53024
A	1081	GLY	-	LINKER	UNP P53024
A	1082	SER	-	LINKER	UNP P53024
A	1083	GLY	-	LINKER	UNP P53024
A	1084	GLY	-	LINKER	UNP P53024
A	1085	GLY	-	LINKER	UNP P53024
A	1086	SER	-	LINKER	UNP P53024
B	1026	GLY	-	EXPRESSION TAG	UNP P53024
B	1027	SER	-	EXPRESSION TAG	UNP P53024
B	1028	HIS	-	EXPRESSION TAG	UNP P53024
B	1029	MET	-	EXPRESSION TAG	UNP P53024
B	1077	GLY	-	LINKER	UNP P53024
B	1078	SER	-	LINKER	UNP P53024
B	1079	GLY	-	LINKER	UNP P53024
B	1080	GLY	-	LINKER	UNP P53024
B	1081	GLY	-	LINKER	UNP P53024
B	1082	SER	-	LINKER	UNP P53024

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1083	GLY	-	LINKER	UNP P53024
B	1084	GLY	-	LINKER	UNP P53024
B	1085	GLY	-	LINKER	UNP P53024
B	1086	SER	-	LINKER	UNP P53024

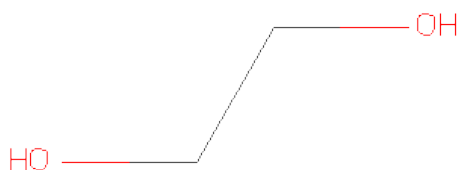
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	2	Total Ca 2 2	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H O 8 2 4 2	0	0
4	A	1	Total C H O 8 2 4 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is water.

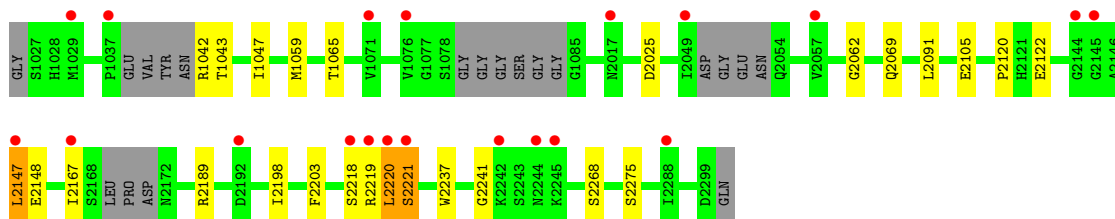
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	171	Total	O	0	0
			171	171		
5	B	128	Total	O	0	0
			128	128		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

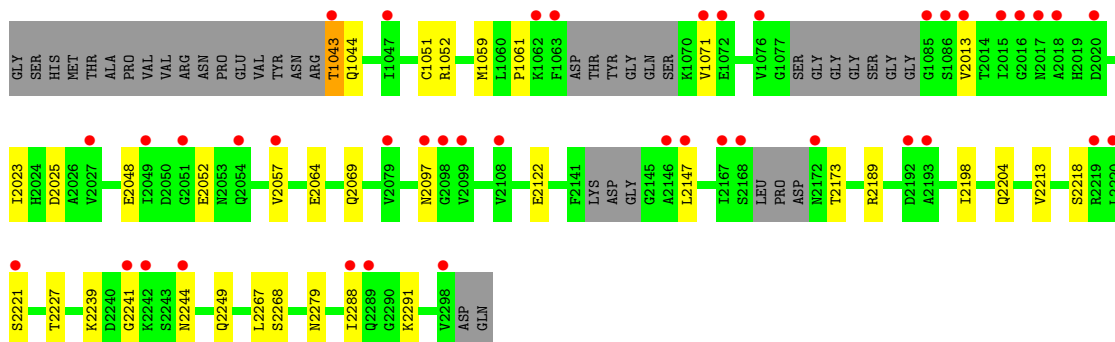
- Molecule 1: Protein transport protein SEC13, COPII coat assembly protein SEC16 fusion protein

Chain A: 



- Molecule 1: Protein transport protein SEC13, COPII coat assembly protein SEC16 fusion protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.51Å 49.32Å 90.90Å 90.00° 111.70° 90.00°	Depositor
Resolution (Å)	44.42 – 1.60 44.42 – 1.56	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.42-1.60) 98.9 (44.42-1.56)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 1.55Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.173 , 0.205 0.176 , 0.207	Depositor DCC
R_{free} test set	4155 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 42.8	EDS
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 82785 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10256	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, EDO, CL

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	0.54	0/2657	0.70	0/3607
1	B	0.48	0/2507	0.66	0/3403
All	All	0.51	0/5164	0.68	0/7010

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2589	2508	11	26	0
1	B	2443	2366	10	37	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
4	A	20	26	0	3	0
5	A	171	0	0	1	0
5	B	128	0	0	6	0
All	All	5356	4900	21	61	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (61) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:2147:LEU:HD22	1:A:2148:GLU:HG3	1.75	0.69
1:B:1052:ARG:HA	1:B:2267:LEU:HD22	1.76	0.68
1:B:2218:SER:O	5:B:2628:HOH:O	2.12	0.66
1:B:1059:MET:HE1	1:B:2023:ILE:HD12	1.77	0.66
1:B:1059:MET:HE2	1:B:1071:VAL:HG11	1.78	0.66
1:B:2239:LYS:NZ	1:B:2244:ASN:HA	2.11	0.64
1:A:1042:ARG:HG3	1:A:1043:THR:HG23	1.80	0.63
1:B:1059:MET:CE	1:B:2023:ILE:HD12	2.30	0.62
1:A:2120:PRO:HB3	1:A:2167:ILE:CD1	2.31	0.61
1:B:1059:MET:HG2	1:B:1071:VAL:HG11	1.82	0.60
1:B:1044:GLN:HG3	1:B:1061:PRO:HG3	1.82	0.59
1:B:1059:MET:HE1	1:B:2023:ILE:CD1	2.33	0.59
1:B:2048:GLU:HG3	1:B:2057:VAL:CG2	2.33	0.59
1:B:1059:MET:HE1	1:B:2023:ILE:HB	1.85	0.59
1:A:2203:PHE:HB3	1:A:2237:TRP:CH2	2.39	0.58
1:B:2025:ASP:OD2	1:B:2069:GLN:HG3	2.04	0.58
1:B:1059:MET:CG	1:B:1071:VAL:HG11	2.34	0.57
1:B:1052:ARG:HA	1:B:2267:LEU:CD2	2.35	0.57
1:A:2167:ILE:O	1:A:2219:ARG:HG2	2.04	0.57
1:B:2064:GLU:OE1	5:B:2525:HOH:O	2.17	0.57
1:A:2275:SER:OG	4:A:2408:EDO:C1	2.52	0.57
1:B:1059:MET:HE1	1:B:2023:ILE:CB	2.35	0.56
1:A:2120:PRO:HB3	1:A:2167:ILE:HD11	1.88	0.56
1:A:2221:SER:HB2	1:A:2241:GLY:HA3	1.92	0.51
1:B:1061:PRO:HD2	1:B:2279:ASN:OD1	2.11	0.51
1:A:2189:ARG:CZ	1:A:2198:ILE:HD11	2.41	0.51
1:B:2173:THR:CG2	1:B:2189:ARG:HG3	2.40	0.50
1:B:2267:LEU:HD12	1:B:2267:LEU:N	2.27	0.50
1:B:2013:VAL:CG2	1:B:2052:GLU:HA	2.41	0.50
1:B:2221:SER:OG	1:B:2241:GLY:HA3	2.13	0.48
1:A:2120:PRO:CB	1:A:2167:ILE:HD11	2.42	0.48
1:A:2268:SER:HB2	1:B:2122:GLU:HG3	1.95	0.47
1:B:2204:GLN:CD	5:B:2551:HOH:O	2.51	0.47
1:A:2147:LEU:H	1:A:2147:LEU:CD1	2.28	0.47
1:A:2189:ARG:NH1	5:A:2617:HOH:O	2.32	0.47
1:B:2213:VAL:HG12	1:B:2227:THR:HG22	1.97	0.47
1:A:2120:PRO:HB3	1:A:2167:ILE:HD12	1.97	0.47
1:B:2198:ILE:C	1:B:2198:ILE:HD12	2.36	0.46
1:B:1059:MET:HG2	1:B:1071:VAL:CG1	2.45	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:2147:LEU:N	1:B:2147:LEU:HD12	2.31	0.46
1:B:1059:MET:CE	1:B:1071:VAL:HG11	2.45	0.46
1:A:2218:SER:OG	1:A:2220:LEU:CD2	2.64	0.46
1:B:1044:GLN:CG	1:B:1061:PRO:HG3	2.46	0.45
1:B:1059:MET:CG	1:B:1071:VAL:CG1	2.94	0.45
1:B:1043:THR:CG2	5:B:2627:HOH:O	2.64	0.45
1:A:1047:ILE:HD12	1:A:1059:MET:CE	2.47	0.45
1:A:2147:LEU:H	1:A:2147:LEU:HD13	1.81	0.44
1:A:2275:SER:OG	4:A:2408:EDO:H11	2.18	0.44
1:A:2218:SER:OG	1:A:2220:LEU:CD1	2.66	0.44
1:B:1043:THR:HG22	5:B:2627:HOH:O	2.16	0.44
1:B:2288:ILE:HG23	5:B:2611:HOH:O	2.17	0.44
1:A:2025:ASP:OD2	1:A:2069:GLN:HG3	2.18	0.44
1:A:2091:LEU:HD22	1:A:2105:GLU:HG2	2.01	0.43
1:B:2249:GLN:OE1	1:B:2291:LYS:HD3	2.19	0.43
1:A:2122:GLU:OE2	1:B:1052:ARG:NH1	2.53	0.42
1:A:2275:SER:CB	4:A:2408:EDO:H11	2.49	0.41
1:B:2147:LEU:CD1	1:B:2147:LEU:N	2.84	0.41
1:A:2220:LEU:HD22	1:A:2220:LEU:O	2.21	0.40
1:B:2048:GLU:HG3	1:B:2057:VAL:HG21	2.04	0.40
1:A:2091:LEU:CD2	1:A:2105:GLU:HG2	2.52	0.40
1:A:1065:THR:O	1:A:1065:THR:HG22	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	320/349 (92%)	309 (97%)	9 (3%)	2 (1%)	33	10
1	B	301/349 (86%)	286 (95%)	15 (5%)	0	100	100
All	All	621/698 (89%)	595 (96%)	24 (4%)	2 (0%)	50	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2221	SER
1	A	2062	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	277/289 (96%)	275 (99%)	2 (1%)	91	81
1	B	260/289 (90%)	256 (98%)	4 (2%)	76	53
All	All	537/578 (93%)	531 (99%)	6 (1%)	84	67

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2147	LEU
1	A	2220	LEU
1	B	1043	THR
1	B	1051	CYS
1	B	2097	ASN
1	B	2268	SER

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

Mol	Chain	Res	Type
1	A	2172	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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	$\# Z > 2$
4	EDO	A	2405	2	3,3,3	0.79	0	2,2,2	0.67	0
4	EDO	A	2406	2	3,3,3	0.60	0	2,2,2	0.48	0
4	EDO	A	2407	-	3,3,3	0.44	0	2,2,2	0.84	0
4	EDO	A	2408	-	3,3,3	0.51	0	2,2,2	0.25	0
4	EDO	A	2409	-	3,3,3	0.54	0	2,2,2	0.40	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	A	2405	2	-	0/1/1/1	0/0/0/0
4	EDO	A	2406	2	-	0/1/1/1	0/0/0/0
4	EDO	A	2407	-	-	0/1/1/1	0/0/0/0
4	EDO	A	2408	-	-	0/1/1/1	0/0/0/0
4	EDO	A	2409	-	-	0/1/1/1	0/0/0/0

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.

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 ⓘ

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	330/349 (94%)	0.18	20 (6%) 21 19	11, 18, 43, 59	0
1	B	311/349 (89%)	0.44	41 (13%) 4 3	13, 23, 52, 62	0
All	All	641/698 (91%)	0.31	61 (9%) 9 8	11, 20, 48, 62	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2219	ARG	7.2
1	A	2288	ILE	6.5
1	B	1085	GLY	6.4
1	B	2099	VAL	5.8
1	B	2242	LYS	5.5
1	B	2017	ASN	4.9
1	B	2057	VAL	4.9
1	B	2288	ILE	4.8
1	B	2098	GLY	4.6
1	B	2015	ILE	4.4
1	A	2220	LEU	4.4
1	B	2298	VAL	4.2
1	B	2167	ILE	4.1
1	B	2219	ARG	3.8
1	B	1043	THR	3.8
1	B	2097	ASN	3.7
1	A	2244	ASN	3.6
1	A	1071	VAL	3.6
1	B	2168	SER	3.6
1	B	1062	LYS	3.5
1	B	1076	VAL	3.5
1	B	2020	ASP	3.4
1	B	2192	ASP	3.3
1	B	2172	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	2220	LEU	3.2
1	B	2013	VAL	3.2
1	B	1071	VAL	3.0
1	A	2221	SER	3.0
1	B	2049	ILE	2.9
1	B	1063	PHE	2.9
1	B	2147	LEU	2.8
1	A	2242	LYS	2.8
1	B	1047	ILE	2.8
1	B	2146	ALA	2.8
1	A	2057	VAL	2.8
1	A	2144	GLY	2.8
1	B	2079	VAL	2.7
1	B	2016	GLY	2.7
1	A	2218	SER	2.6
1	B	1072	GLU	2.6
1	A	2049	ILE	2.5
1	A	2147	LEU	2.4
1	B	2221	SER	2.4
1	B	2244	ASN	2.4
1	B	2108	VAL	2.4
1	B	2193	ALA	2.4
1	A	2017	ASN	2.4
1	B	2241	GLY	2.3
1	B	2054	GLN	2.2
1	A	1037	PRO	2.2
1	A	2192	ASP	2.2
1	A	1076	VAL	2.2
1	A	2167	ILE	2.2
1	A	2245	LYS	2.2
1	B	1086	SER	2.2
1	A	1029	MET	2.2
1	A	2145	GLY	2.2
1	B	2051	GLY	2.1
1	B	2027	VAL	2.1
1	B	2018	ALA	2.0
1	B	2289	GLN	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	EDO	A	2408	4/4	0.14	2.21	19,26,32,37	0
4	EDO	A	2409	4/4	0.12	0.66	24,31,38,38	0
4	EDO	A	2405	4/4	0.09	0.19	19,21,24,24	0
2	CA	A	2402	1/1	0.08	0.00	20,20,20,20	0
2	CA	A	2401	1/1	0.08	-0.21	11,11,11,11	0
2	CA	B	2401	1/1	0.06	-0.30	15,15,15,15	0
4	EDO	A	2407	4/4	0.09	-0.44	18,27,28,33	0
4	EDO	A	2406	4/4	0.07	-0.65	14,16,19,19	0
3	CL	A	2403	1/1	0.07	-0.70	18,18,18,18	0
3	CL	A	2404	1/1	0.06	-0.97	21,21,21,21	0

6.5 Other polymers ⓘ

There are no such residues in this entry.