



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:53 AM UTC

PDB ID : 3CKC / pdb_00003ckc
Title : B. thetaiotaomicron SusD
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Deposited on : 2008-03-14
Resolution : 1.50 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

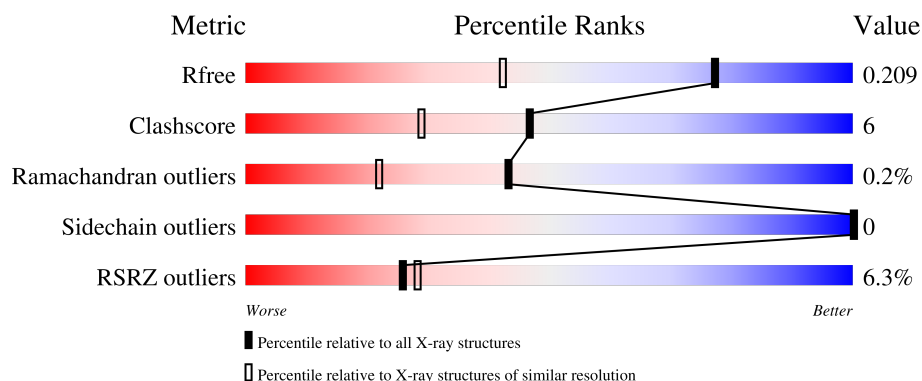
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	527	9% 83% 12% 5%
1	B	527	3% 83% 12% 5%

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	A	903	-	-	X	-
3	EDO	B	910	-	-	X	-
5	PEG	A	800	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9347 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 SusD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			4027	2534	692	782	19			
1	B	500	Total	C	N	O	S	0	0	0
			4021	2531	691	780	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	GLY	-	expression tag	UNP Q8A1G2
A	100	LYS	GLU	SEE REMARK 999	UNP Q8A1G2
B	25	GLY	-	expression tag	UNP Q8A1G2
B	100	LYS	GLU	SEE REMARK 999	UNP Q8A1G2

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



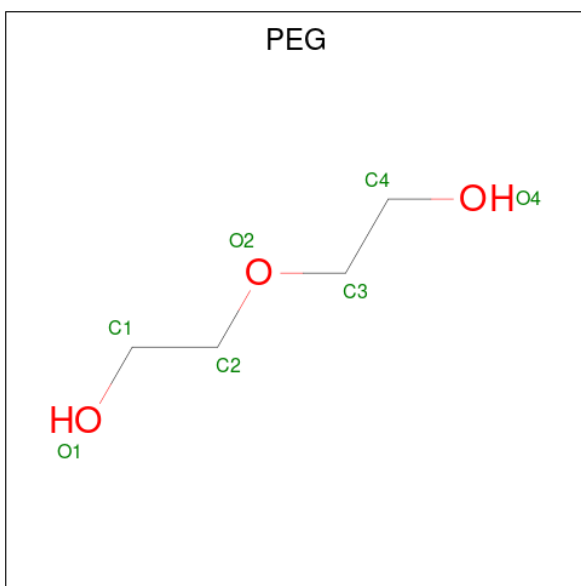
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total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

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

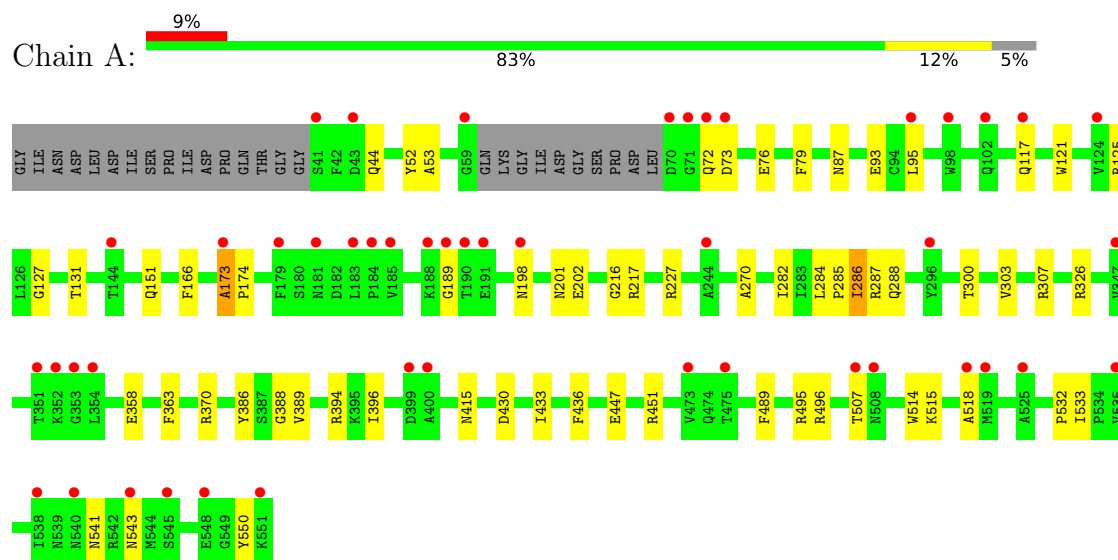
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	558	Total 558	O 558	0	0
6	B	665	Total 665	O 665	0	0

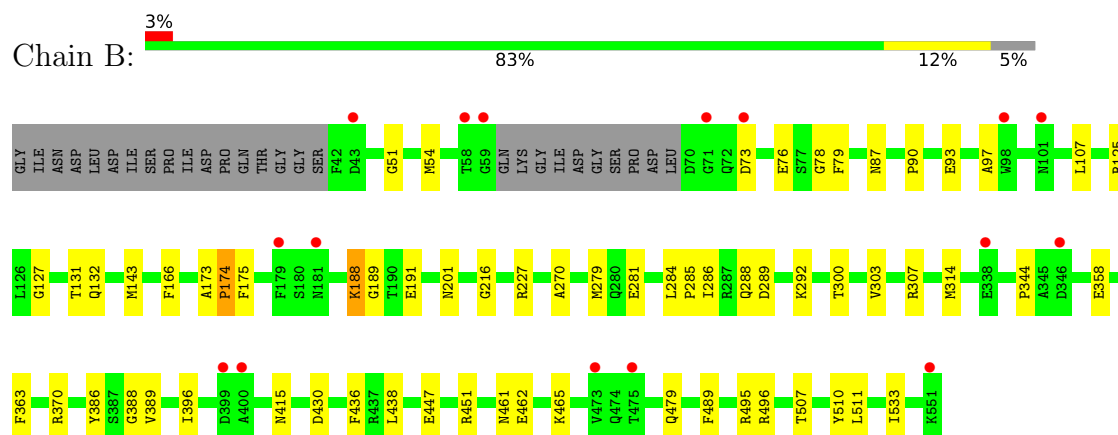
3 Residue-property plots [i](#)

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: SusD



• Molecule 1: SusD



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.20Å 68.06Å 83.05Å 111.09° 93.17° 109.17°	Depositor
Resolution (Å)	50.00 – 1.50 50.00 – 1.50	Depositor EDS
% Data completeness (in resolution range)	96.2 (50.00-1.50) 96.2 (50.00-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	17.69 (at 1.47Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.192 , 0.210 0.192 , 0.209	Depositor DCC
R_{free} test set	18043 reflections (9.47%)	wwPDB-VP
Wilson B-factor (Å ²)	14.2	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.6	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	0.96	EDS
Total number of atoms	9347	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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: PEG, CA, EDO, MES, CSO

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	0.32	0/4111	0.83	10/5561 (0.2%)
1	B	0.34	0/4105	0.86	14/5553 (0.3%)
All	All	0.33	0/8216	0.84	24/11114 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	533	ILE	N-CA-C	-7.70	101.41	108.95
1	A	396	ILE	N-CA-C	-7.65	105.23	111.81
1	A	533	ILE	N-CA-C	-7.29	102.36	108.63
1	B	396	ILE	N-CA-C	-7.29	105.54	111.81
1	B	174	PRO	N-CA-C	-7.17	98.83	110.74
1	B	97	ALA	N-CA-C	6.22	118.06	111.28
1	B	286	ILE	N-CA-C	-6.16	98.33	107.51
1	B	175	PHE	N-CA-C	6.10	119.58	109.46
1	A	507	THR	N-CA-C	6.09	118.76	110.35
1	B	370	ARG	N-CA-C	-5.87	102.62	110.55
1	B	78	GLY	N-CA-C	5.85	118.66	111.93
1	A	370	ARG	N-CA-C	-5.72	102.83	110.55
1	B	489	PHE	N-CA-C	5.66	120.18	112.88
1	A	286	ILE	N-CA-C	-5.58	99.19	107.51
1	B	270	ALA	N-CA-C	5.50	119.09	112.38
1	B	507	THR	N-CA-C	5.49	117.92	110.35
1	A	489	PHE	N-CA-C	5.47	119.94	112.88
1	B	188	LYS	N-CA-C	5.31	117.94	110.14
1	B	462	GLU	N-CA-C	-5.28	105.60	111.36
1	B	300	THR	N-CA-C	-5.23	105.58	111.28
1	A	174	PRO	N-CA-C	-5.08	102.00	112.47
1	A	300	THR	N-CA-C	-5.08	105.75	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	270	ALA	N-CA-C	5.04	118.53	112.38
1	A	173	ALA	N-CA-C	5.04	114.61	108.11

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	4027	0	3835	41	0
1	B	4021	0	3830	44	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	16	0	23	8	0
3	B	32	0	47	11	0
4	A	12	0	13	0	0
5	A	7	0	10	4	0
5	B	7	0	10	2	0
6	A	558	0	0	4	0
6	B	665	0	0	11	0
All	All	9347	0	7768	90	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 (90) 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)
3:A:903:EDO:H21	1:B:358:GLU:HG2	1.38	1.05
1:A:287:ARG:HB3	5:A:800:PEG:H32	1.53	0.88
1:A:44:GLN:HE21	1:A:151:GLN:HE22	1.24	0.85
1:A:72:GLN:HG2	1:A:76:GLU:HB2	1.62	0.81
1:B:279:MET:HE1	3:B:909:EDO:H21	1.64	0.80
1:A:288:GLN:O	5:A:800:PEG:H41	1.84	0.77
1:A:166:PHE:HB3	1:A:173:ALA:HB2	1.67	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:ASN:HD22	1:B:227:ARG:HH12	1.36	0.74
1:B:73:ASP:O	1:B:76:GLU:HG2	1.88	0.72
1:B:510:TYR:HA	3:B:910:EDO:H12	1.70	0.72
1:A:95:LEU:HG	1:A:326:ARG:HG2	1.72	0.72
1:A:217:ARG:HG3	3:A:901:EDO:H22	1.71	0.70
3:A:903:EDO:H22	6:A:1464:HOH:O	1.92	0.70
3:A:903:EDO:H21	1:B:358:GLU:CG	2.20	0.69
1:A:201:ASN:HD22	1:A:227:ARG:HH12	1.41	0.69
1:A:44:GLN:HE21	1:A:151:GLN:NE2	1.90	0.68
1:B:288:GLN:O	5:B:801:PEG:H41	1.95	0.66
3:A:903:EDO:H11	1:B:358:GLU:HB3	1.76	0.66
1:B:279:MET:CE	3:B:909:EDO:H21	2.29	0.62
1:B:279:MET:HE2	6:B:1430:HOH:O	1.98	0.62
1:A:198:ASN:O	1:A:202:GLU:HG3	2.02	0.59
1:B:166:PHE:HB3	1:B:173:ALA:HB2	1.85	0.58
1:A:447:GLU:O	1:A:451:ARG:HG2	2.03	0.57
1:A:93:GLU:HG3	1:A:495:ARG:HG2	1.86	0.56
1:A:415:ASN:ND2	1:A:430:ASP:H	2.04	0.55
1:B:188:LYS:HB2	6:B:1850:HOH:O	2.04	0.55
1:B:511:LEU:H	3:B:910:EDO:C1	2.20	0.55
3:B:908:EDO:O1	5:B:801:PEG:H42	2.08	0.53
1:B:447:GLU:O	1:B:451:ARG:HG2	2.08	0.53
1:A:284:LEU:HB3	1:A:436:PHE:HB2	1.91	0.53
1:B:93:GLU:HG3	1:B:495:ARG:HG2	1.89	0.53
5:A:800:PEG:H31	6:A:1451:HOH:O	2.09	0.52
1:B:284:LEU:HB3	1:B:436:PHE:HB2	1.91	0.52
1:A:541:ASN:OD1	1:A:543:ASN:HB2	2.10	0.52
1:A:415:ASN:HD21	1:A:430:ASP:H	1.58	0.51
1:B:479:GLN:NE2	6:B:1447:HOH:O	2.44	0.51
1:A:117:GLN:HB2	6:A:1149:HOH:O	2.11	0.51
1:B:216:GLY:HA2	1:B:285:PRO:HG2	1.93	0.50
1:A:216:GLY:HA2	1:A:285:PRO:HG2	1.94	0.50
1:B:415:ASN:ND2	1:B:430:ASP:H	2.10	0.49
1:A:72:GLN:HG2	1:A:76:GLU:CB	2.39	0.49
1:B:90:PRO:HB3	1:B:107:LEU:O	2.13	0.49
1:A:52:TYR:O	3:A:902:EDO:H11	2.14	0.48
1:B:127:GLY:HA2	3:B:906:EDO:H12	1.95	0.48
1:B:143:MET:HE3	6:B:1809:HOH:O	2.12	0.48
1:A:95:LEU:HG	1:A:326:ARG:CG	2.40	0.48
1:B:461:ASN:O	1:B:465:LYS:HG3	2.14	0.48
1:B:51:GLY:HA2	1:B:54:MET:HE2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:VAL:O	1:A:307:ARG:HG2	2.15	0.47
1:A:127:GLY:O	1:A:131:THR:HG23	2.15	0.47
1:B:314:MET:HE2	6:B:1538:HOH:O	2.16	0.45
1:B:415:ASN:HD21	1:B:430:ASP:H	1.64	0.45
1:B:479:GLN:HB2	6:B:1680:HOH:O	2.15	0.45
1:A:286:ILE:HA	3:A:902:EDO:H22	1.99	0.45
1:A:358:GLU:H	1:A:358:GLU:CD	2.25	0.45
1:A:287:ARG:NE	5:A:800:PEG:O4	2.49	0.45
1:B:289:ASP:OD2	3:B:908:EDO:H11	2.17	0.45
1:A:117:GLN:HG2	1:A:121:TRP:CE2	2.52	0.44
1:B:281:GLU:HG3	1:B:438:LEU:HB3	1.99	0.44
1:A:282:ILE:HG21	1:A:285:PRO:HB3	2.00	0.44
1:B:303:VAL:O	1:B:307:ARG:HG2	2.18	0.44
1:B:344:PRO:HD3	6:B:1627:HOH:O	2.18	0.44
1:B:201:ASN:HD22	1:B:227:ARG:NH1	2.11	0.44
1:A:53:ALA:HA	3:A:902:EDO:H11	2.00	0.43
1:A:515:LYS:HA	6:A:1069:HOH:O	2.18	0.43
1:A:386:TYR:CZ	1:A:388:GLY:HA3	2.54	0.43
1:B:363:PHE:CE1	1:B:389:VAL:HG11	2.54	0.43
1:B:386:TYR:CZ	1:B:388:GLY:HA3	2.54	0.43
1:A:166:PHE:CB	1:A:173:ALA:HB2	2.43	0.43
1:B:79:PHE:HA	1:B:125:ARG:HG2	2.01	0.43
3:B:911:EDO:H22	6:B:1784:HOH:O	2.19	0.43
1:B:87:ASN:HD22	1:B:496:ARG:HB3	1.84	0.42
1:B:191:GLU:HG3	6:B:1720:HOH:O	2.19	0.42
1:B:289:ASP:HB3	1:B:292:LYS:HB2	2.00	0.42
1:A:363:PHE:CE1	1:A:389:VAL:HG11	2.54	0.42
1:B:358:GLU:CD	1:B:358:GLU:H	2.28	0.42
1:A:514:TRP:HA	1:A:518:ALA:O	2.20	0.42
1:B:314:MET:HE1	6:B:1872:HOH:O	2.19	0.41
1:B:54:MET:HE1	1:B:132:GLN:HB2	2.03	0.41
1:B:174:PRO:HG2	3:B:906:EDO:O1	2.21	0.41
1:A:87:ASN:HD22	1:A:496:ARG:HD3	1.86	0.41
3:B:910:EDO:H11	6:B:1930:HOH:O	2.21	0.41
1:A:79:PHE:HA	1:A:125:ARG:HG2	2.03	0.41
1:A:394:ARG:HH11	1:A:394:ARG:HG2	1.86	0.41
1:B:511:LEU:H	3:B:910:EDO:H12	1.85	0.40
1:A:286:ILE:HB	1:A:433:ILE:HB	2.03	0.40
1:A:87:ASN:HD22	1:A:87:ASN:HA	1.72	0.40
1:B:127:GLY:O	1:B:131:THR:HG23	2.20	0.40
1:A:73:ASP:N	1:A:76:GLU:HG3	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:532:PRO:HB3	1:A:550:TYR:CG	2.55	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	496/527 (94%)	481 (97%)	14 (3%)	1 (0%)	43	22
1	B	495/527 (94%)	483 (98%)	11 (2%)	1 (0%)	43	22
All	All	991/1054 (94%)	964 (97%)	25 (2%)	2 (0%)	43	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	189	GLY
1	A	189	GLY

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	419/440 (95%)	419 (100%)	0	100	100
1	B	418/440 (95%)	418 (100%)	0	100	100
All	All	837/880 (95%)	837 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	ASN
1	A	102	GLN
1	A	117	GLN
1	A	151	GLN
1	A	198	ASN
1	A	201	ASN
1	A	295	ASN
1	A	415	ASN
1	A	468	ASN
1	A	479	GLN
1	A	540	ASN
1	A	543	ASN
1	B	44	GLN
1	B	85	ASN
1	B	87	ASN
1	B	151	GLN
1	B	201	ASN
1	B	275	ASN
1	B	415	ASN
1	B	468	ASN
1	B	479	GLN
1	B	540	ASN
1	B	543	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSO	B	322	1	3,6,7	0.62	0	1,6,8	1.61	0
1	CSO	A	322	1	3,6,7	0.61	0	1,6,8	1.45	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
1	CSO	B	322	1	-	0/1/5/7	-
1	CSO	A	322	1	-	0/1/5/7	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 2 are monoatomic - leaving 15 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	EDO	A	907	2	3,3,3	0.75	0	2,2,2	0.27	0
5	PEG	A	800	2	6,6,6	0.95	0	5,5,5	2.39	3 (60%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	B	909	-	3,3,3	0.70	0	2,2,2	0.39	0
5	PEG	B	801	2	6,6,6	0.95	0	5,5,5	2.37	3 (60%)
3	EDO	A	901	-	3,3,3	0.66	0	2,2,2	0.44	0
3	EDO	A	902	-	3,3,3	0.79	0	2,2,2	0.28	0
3	EDO	B	906	-	3,3,3	0.66	0	2,2,2	0.41	0
3	EDO	B	904	-	3,3,3	0.67	0	2,2,2	0.40	0
3	EDO	B	908	2	3,3,3	0.75	0	2,2,2	0.29	0
3	EDO	B	910	-	3,3,3	0.68	0	2,2,2	0.41	0
3	EDO	B	911	-	3,3,3	0.69	0	2,2,2	0.35	0
3	EDO	A	903	-	3,3,3	0.69	0	2,2,2	0.26	0
3	EDO	B	905	-	3,3,3	0.68	0	2,2,2	0.40	0
3	EDO	B	912	-	3,3,3	0.73	0	2,2,2	0.35	0
4	MES	A	1000	-	12,12,12	1.16	1 (8%)	15,16,16	1.00	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
3	EDO	A	907	2	-	1/1/1/1	-
5	PEG	A	800	2	-	2/4/4/4	-
3	EDO	B	909	-	-	1/1/1/1	-
5	PEG	B	801	2	-	2/4/4/4	-
3	EDO	A	901	-	-	0/1/1/1	-
3	EDO	A	902	-	-	1/1/1/1	-
3	EDO	B	906	-	-	0/1/1/1	-
3	EDO	B	904	-	-	0/1/1/1	-
3	EDO	B	908	2	-	1/1/1/1	-
3	EDO	B	910	-	-	0/1/1/1	-
3	EDO	B	911	-	-	0/1/1/1	-
3	EDO	A	903	-	-	1/1/1/1	-
3	EDO	B	905	-	-	0/1/1/1	-
3	EDO	B	912	-	-	0/1/1/1	-
4	MES	A	1000	-	-	0/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1000	MES	C8-S	2.92	1.81	1.77

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	800	PEG	C3-O2-C2	3.15	127.05	113.26
5	B	801	PEG	C3-O2-C2	3.08	126.74	113.26
5	A	800	PEG	O2-C3-C4	2.89	122.86	110.11
5	B	801	PEG	O2-C2-C1	2.86	122.71	110.11
5	A	800	PEG	O2-C2-C1	2.82	122.53	110.11
5	B	801	PEG	O2-C3-C4	2.81	122.51	110.11

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	801	PEG	O2-C3-C4-O4
3	B	909	EDO	O1-C1-C2-O2
5	A	800	PEG	O2-C3-C4-O4
5	B	801	PEG	C1-C2-O2-C3
5	A	800	PEG	C1-C2-O2-C3
3	A	902	EDO	O1-C1-C2-O2
3	B	908	EDO	O1-C1-C2-O2
3	A	903	EDO	O1-C1-C2-O2
3	A	907	EDO	O1-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	800	PEG	4	0
3	B	909	EDO	2	0
5	B	801	PEG	2	0
3	A	901	EDO	1	0
3	A	902	EDO	3	0
3	B	906	EDO	2	0
3	B	908	EDO	2	0
3	B	910	EDO	4	0
3	B	911	EDO	1	0
3	A	903	EDO	4	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

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	500/527 (94%)	0.66	47 (9%)	14 14	9, 17, 26, 34	0
1	B	499/527 (94%)	0.23	16 (3%)	50 54	8, 14, 23, 29	0
All	All	999/1054 (94%)	0.44	63 (6%)	26 28	8, 15, 25, 34	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	71	GLY	5.8
1	A	70	ASP	5.7
1	B	98	TRP	5.0
1	A	59	GLY	4.6
1	B	73	ASP	4.3
1	B	551	LYS	4.2
1	A	73	ASP	4.2
1	B	59	GLY	4.0
1	A	144	THR	3.8
1	A	551	LYS	3.8
1	A	354	LEU	3.5
1	A	181	ASN	3.3
1	A	41	SER	3.3
1	A	548	GLU	3.1
1	B	58	THR	3.1
1	A	188	LYS	3.0
1	A	352	LYS	3.0
1	A	72	GLN	2.9
1	A	538	ILE	2.8
1	B	475	THR	2.7
1	A	518	ALA	2.7
1	A	189	GLY	2.7
1	A	351	THR	2.7
1	A	399	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	525	ALA	2.6
1	A	519	MET	2.6
1	B	181	ASN	2.6
1	A	98	TRP	2.5
1	A	400	ALA	2.5
1	B	399	ASP	2.4
1	A	185	VAL	2.4
1	A	545	SER	2.4
1	B	400	ALA	2.3
1	A	124	VAL	2.3
1	A	508	ASN	2.3
1	A	117	GLN	2.3
1	A	173	ALA	2.3
1	A	347	VAL	2.2
1	A	198	ASN	2.2
1	A	475	THR	2.2
1	A	102	GLN	2.2
1	B	71	GLY	2.2
1	A	473	VAL	2.2
1	A	535	VAL	2.2
1	A	296	TYR	2.1
1	A	540	ASN	2.1
1	A	543	ASN	2.1
1	A	43	ASP	2.1
1	A	190	THR	2.1
1	B	101	ASN	2.1
1	A	191	GLU	2.1
1	A	353	GLY	2.1
1	B	338	GLU	2.1
1	B	43	ASP	2.1
1	A	95	LEU	2.1
1	A	507	THR	2.0
1	B	179	PHE	2.0
1	A	244	ALA	2.0
1	A	183	LEU	2.0
1	B	346	ASP	2.0
1	A	184	PRO	2.0
1	A	179	PHE	2.0
1	B	473	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSO	A	322	7/8	0.82	0.15	20,21,30,31	0
1	CSO	B	322	7/8	0.94	0.09	16,16,22,28	0

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
4	MES	A	1000	12/12	0.69	0.20	31,33,39,39	0
3	EDO	B	910	4/4	0.76	0.19	23,23,26,29	0
5	PEG	B	801	7/7	0.77	0.15	20,24,29,30	0
3	EDO	B	909	4/4	0.79	0.16	28,28,29,30	0
3	EDO	A	903	4/4	0.80	0.21	24,25,26,27	0
3	EDO	B	911	4/4	0.80	0.19	27,28,28,29	0
3	EDO	A	901	4/4	0.81	0.19	27,28,28,28	0
3	EDO	B	906	4/4	0.84	0.16	28,28,29,29	0
5	PEG	A	800	7/7	0.85	0.14	18,22,28,30	0
3	EDO	B	908	4/4	0.86	0.23	22,25,27,30	0
3	EDO	A	907	4/4	0.89	0.20	19,23,25,26	0
3	EDO	B	904	4/4	0.90	0.13	24,24,25,27	0
3	EDO	B	912	4/4	0.91	0.11	24,25,25,26	0
3	EDO	A	902	4/4	0.91	0.19	22,23,23,23	0
3	EDO	B	905	4/4	0.96	0.07	18,20,21,21	0
2	CA	B	700	1/1	0.98	0.12	18,18,18,18	0
2	CA	A	600	1/1	0.98	0.11	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.