



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 08:04 PM EDT

PDB ID : 7B5G / pdb_00007b5g
Title : Crystal structure of E.coli LexA in complex with nanobody NbSOS3(Nb14527)
Authors : Maso, L.; Vascon, F.; Chinellato, M.; Pardon, E.; Steyaert, J.; Angelini, A.;
Tondi, D.; Cendron, L.
Deposited on : 2020-12-03
Resolution : 2.40 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

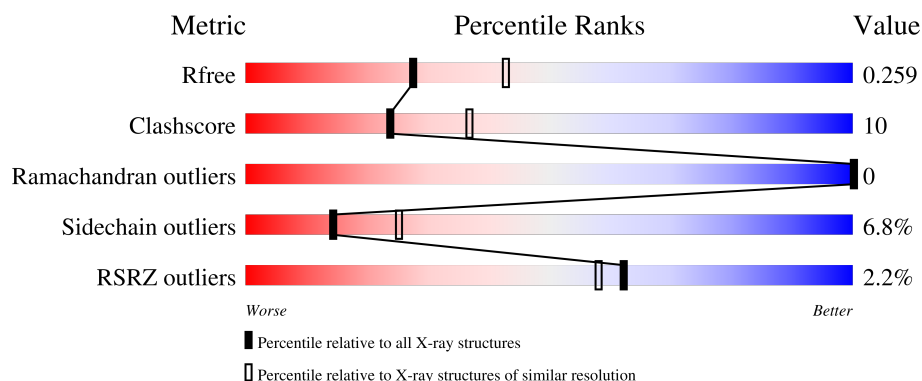
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	202	0% 50% 13% • 35%
1	C	202	49% 15% 36%
1	E	202	2% 50% 13% • 36%
1	G	202	47% 15% • 36%
2	B	132	2% 82% 11% 7%

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Mol	Chain	Length	Quality of chain
2	D	132	% 72% 21% • 6%
2	F	132	4% 76% 19% 5%
2	H	132	4% 63% 27% • 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7961 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 LexA repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	1	0
			1035	653	182	197	3			
1	C	130	Total	C	N	O	S	0	2	0
			1026	649	182	192	3			
1	E	130	Total	C	N	O	S	0	0	0
			1010	638	179	190	3			
1	G	129	Total	C	N	O	S	0	1	0
			1008	638	179	188	3			

- Molecule 2 is a protein called Nanobody Nb14527, NbSOS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	123	Total	C	N	O	S	0	0	0
			925	564	171	186	4			
2	D	124	Total	C	N	O	S	0	0	0
			935	570	174	187	4			
2	F	125	Total	C	N	O	S	0	0	0
			945	576	177	188	4			
2	H	124	Total	C	N	O	S	0	0	0
			935	570	174	187	4			

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



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

- Molecule 4 is water.

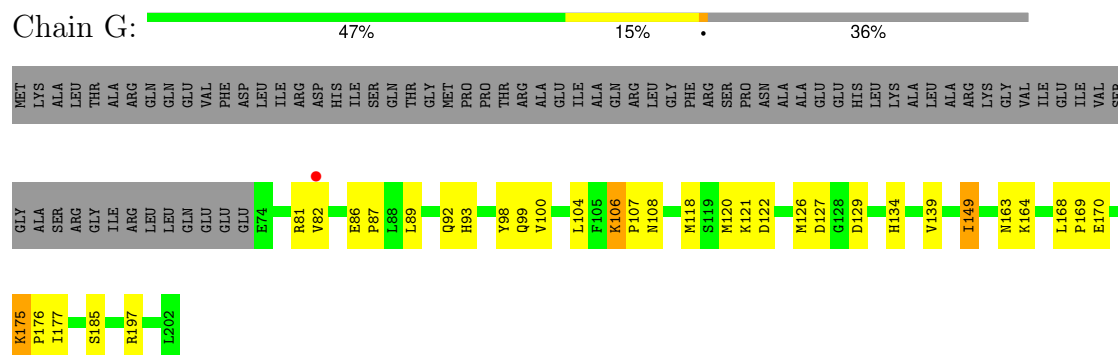
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	23	Total O 23 23	0	0
4	B	19	Total O 19 19	0	0
4	C	12	Total O 12 12	0	0
4	D	12	Total O 12 12	0	0
4	E	14	Total O 14 14	0	0
4	F	20	Total O 20 20	0	0

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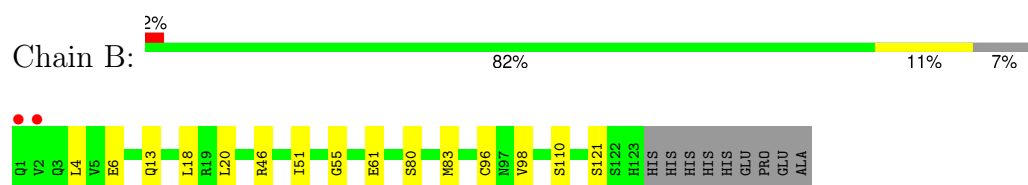
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	9	Total	O	0	0
			9	9		
4	H	9	Total	O	0	0
			9	9		

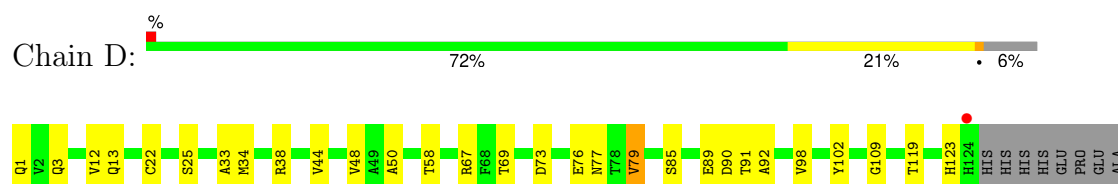
- Molecule 1: LexA repressor



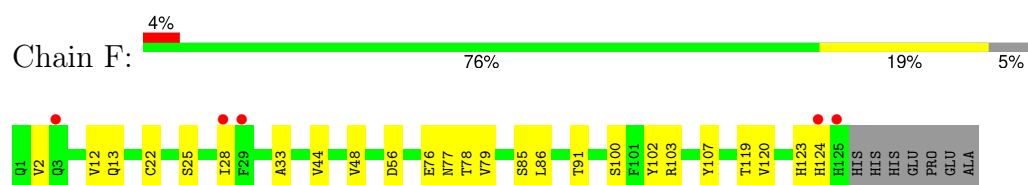
- Molecule 2: Nanobody Nb14527, NbSOS3



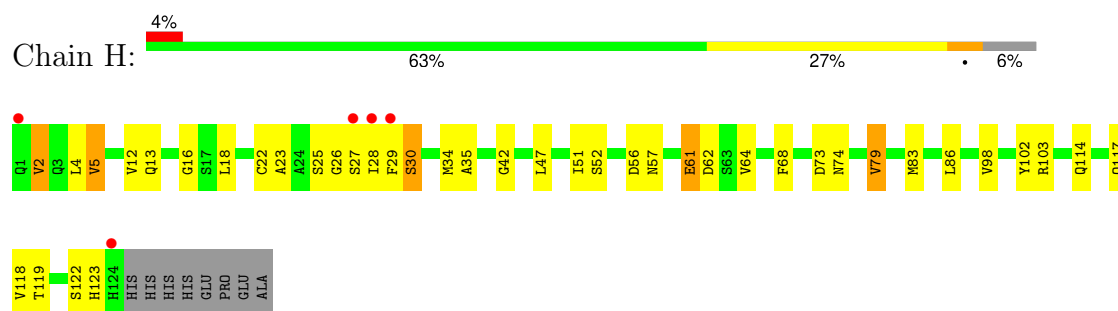
- Molecule 2: Nanobody Nb14527, NbSOS3



- Molecule 2: Nanobody Nb14527, NbSOS3



- Molecule 2: Nanobody Nb14527, NbSOS3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.07Å 57.08Å 122.51Å 84.53° 83.21° 68.43°	Depositor
Resolution (Å)	49.61 – 2.40 49.61 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.0 (49.61-2.40) 98.1 (49.61-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.206 , 0.252 0.210 , 0.259	Depositor DCC
R_{free} test set	2360 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7961	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% 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: EDO

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	0.56	0/1050	1.03	1/1419 (0.1%)
1	C	0.53	0/1041	0.98	1/1406 (0.1%)
1	E	0.53	0/1025	1.04	2/1385 (0.1%)
1	G	0.50	0/1023	1.02	0/1383
2	B	0.54	0/941	1.08	0/1273
2	D	0.53	0/952	0.98	1/1288 (0.1%)
2	F	0.57	0/963	1.05	0/1303
2	H	0.57	0/952	1.02	0/1288
All	All	0.54	0/7947	1.03	5/10745 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	123	HIS	N-CA-C	-6.85	105.70	112.97
1	E	74	GLU	CB-CG-CD	5.74	122.36	112.60
1	A	171	ASN	N-CA-C	-5.21	99.70	110.80
1	C	99	GLN	CB-CA-C	5.14	118.23	109.84
1	E	110	ASP	N-CA-C	-5.14	107.56	113.88

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	1035	0	1050	29	0
1	C	1026	0	1050	21	0
1	E	1010	0	1030	21	0
1	G	1008	0	1032	22	0
2	B	925	0	875	8	0
2	D	935	0	882	27	0
2	F	945	0	889	14	0
2	H	935	0	882	28	0
3	A	8	0	12	0	0
3	C	16	0	24	1	0
4	A	23	0	0	1	0
4	B	19	0	0	0	0
4	C	12	0	0	0	0
4	D	12	0	0	2	0
4	E	14	0	0	0	0
4	F	20	0	0	0	0
4	G	9	0	0	0	0
4	H	9	0	0	0	0
All	All	7961	0	7726	158	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 10.

All (158) 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:201:TRP:HZ3	2:H:13:GLN:HB3	1.04	1.15
2:D:34:MET:HE1	2:D:98:VAL:HG22	1.35	1.08
1:A:201:TRP:CZ3	2:H:13:GLN:HB3	1.93	1.03
1:C:124:GLY:HA2	1:C:126:MET:HE2	1.50	0.94
1:E:129:ASP:OD1	1:E:197:ARG:HD3	1.72	0.90
1:A:71:GLU:OE1	1:A:97:HIS:HE1	1.55	0.89
1:C:160:LYS:HG2	1:C:165:VAL:HG22	1.53	0.88
2:H:102:TYR:CE1	2:H:103:ARG:HG3	2.13	0.83
1:G:86:GLU:HG3	1:G:87:PRO:HD2	1.62	0.81
2:D:34:MET:HE1	2:D:98:VAL:CG2	2.12	0.79
2:H:52:SER:HB3	2:H:57:ASN:HB2	1.65	0.79
1:A:71:GLU:OE1	1:A:97:HIS:CE1	2.37	0.78
2:H:5:VAL:HG12	2:H:23:ALA:HB3	1.68	0.76
2:H:30:SER:HA	2:H:74:ASN:ND2	2.01	0.74
2:D:34:MET:CE	2:D:98:VAL:HA	2.17	0.73
2:H:61:GLU:HG3	2:H:62:ASP:N	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:3:GLN:HB2	2:D:25:SER:OG	1.88	0.73
2:D:34:MET:HE2	2:D:98:VAL:HA	1.69	0.73
1:A:129:ASP:OD1	1:A:197:ARG:HD3	1.89	0.73
1:C:140:ARG:HB3	3:C:302:EDO:H21	1.72	0.71
1:A:201:TRP:HZ3	2:H:13:GLN:CB	1.96	0.70
2:D:12:VAL:HG22	2:D:13:GLN:H	1.58	0.69
1:E:74:GLU:HG3	1:E:97:HIS:CE1	2.27	0.68
1:E:78:LEU:HD13	1:E:98:TYR:HE2	1.58	0.68
2:F:12:VAL:HG21	2:F:86:LEU:HD13	1.76	0.68
2:H:4:LEU:HD11	2:H:98:VAL:HG23	1.76	0.67
1:G:118:MET:HE3	1:G:121:LYS:HG3	1.77	0.66
2:H:12:VAL:HG11	2:H:86:LEU:HD13	1.77	0.65
1:G:82:VAL:HA	1:G:93:HIS:CD2	2.31	0.65
1:C:129:ASP:OD1	1:C:197:ARG:HD3	1.98	0.64
1:E:74:GLU:O	1:E:97:HIS:HB3	1.98	0.63
1:E:171:ASN:OD1	1:E:173:GLU:HG2	1.98	0.63
1:G:149:ILE:HD12	1:G:149:ILE:N	2.13	0.63
1:G:129:ASP:OD1	1:G:197:ARG:HD3	2.00	0.61
1:A:108:ASN:HD22	1:A:108:ASN:H	1.48	0.61
2:D:3:GLN:HB2	2:D:25:SER:HG	1.66	0.61
1:G:98:TYR:O	1:G:100:VAL:HG23	2.01	0.60
1:G:118:MET:HE1	1:G:126:MET:HE3	1.82	0.60
1:C:121[A]:LYS:HD3	1:C:171:ASN:ND2	2.16	0.59
1:C:129:ASP:OD2	1:C:197:ARG:NH1	2.35	0.59
1:C:160:LYS:HG2	1:C:165:VAL:CG2	2.28	0.59
1:C:124:GLY:HA2	1:C:126:MET:CE	2.28	0.58
2:B:20:LEU:HG	2:B:83:MET:HE2	1.84	0.58
1:C:129:ASP:CG	1:C:197:ARG:HH11	2.11	0.58
1:G:82:VAL:HA	1:G:93:HIS:HD2	1.67	0.57
2:D:12:VAL:HG22	2:D:13:GLN:N	2.19	0.57
2:D:109:GLY:HA3	4:D:303:HOH:O	2.04	0.57
2:H:16:GLY:O	2:H:86:LEU:HD12	2.05	0.57
2:H:30:SER:HA	2:H:74:ASN:HD22	1.69	0.56
1:E:118:MET:O	1:E:171:ASN:ND2	2.30	0.56
1:A:108:ASN:H	1:A:108:ASN:ND2	2.05	0.55
2:H:83:MET:HE1	2:H:118:VAL:HG21	1.89	0.55
1:A:101:ASP:OD1	1:A:102:PRO:HD2	2.07	0.55
2:F:22:CYS:HB3	2:F:79:VAL:HG12	1.89	0.55
2:F:12:VAL:HG21	2:F:86:LEU:CD1	2.38	0.54
2:H:29:PHE:HE1	2:H:34:MET:HE2	1.71	0.54
2:H:114:GLN:H	2:H:114:GLN:CD	2.14	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:149:ILE:HD12	1:E:149:ILE:N	2.23	0.54
2:D:38:ARG:HG2	2:D:48:VAL:HG22	1.88	0.53
2:H:61:GLU:HG3	2:H:62:ASP:H	1.69	0.52
1:G:106:LYS:HA	1:G:107:PRO:C	2.33	0.52
1:E:76:LEU:O	1:E:97:HIS:HA	2.10	0.52
2:F:33:ALA:HB3	2:F:100:SER:HA	1.90	0.52
2:D:91:THR:HG23	2:D:119:THR:HA	1.91	0.51
2:D:34:MET:HE1	2:D:98:VAL:HA	1.91	0.51
2:D:34:MET:HE2	2:D:34:MET:HA	1.91	0.51
1:G:89:LEU:HD23	2:H:102:TYR:CD1	2.45	0.51
1:G:120:MET:HA	1:G:170:GLU:HB2	1.93	0.51
1:C:92:GLN:H	1:C:92:GLN:CD	2.19	0.50
2:D:12:VAL:CG2	2:D:13:GLN:H	2.23	0.50
1:A:82:VAL:HG23	1:A:93:HIS:CD2	2.46	0.50
1:A:78:LEU:HD12	1:A:98:TYR:HE2	1.77	0.50
1:E:123:ILE:HD12	1:E:125:ILE:HD12	1.93	0.49
1:E:123:ILE:CD1	1:E:125:ILE:HD12	2.41	0.49
1:A:165:VAL:O	1:A:179:VAL:HG12	2.13	0.49
1:E:74:GLU:HG3	1:E:97:HIS:ND1	2.28	0.49
2:D:22:CYS:HB3	2:D:79:VAL:HG13	1.94	0.49
1:E:88:LEU:HD11	1:E:153:VAL:HG11	1.93	0.49
1:G:86:GLU:HG3	1:G:87:PRO:CD	2.39	0.49
2:H:22:CYS:HB3	2:H:79:VAL:HG13	1.94	0.49
2:D:67:ARG:NH2	2:D:90:ASP:OD2	2.46	0.48
2:D:34:MET:HE1	2:D:98:VAL:CB	2.43	0.48
1:A:146:VAL:HG12	1:A:189:GLU:HB2	1.96	0.48
1:E:89:LEU:HD23	2:F:102:TYR:CD2	2.49	0.48
1:A:119:SER:HB3	4:A:410:HOH:O	2.14	0.48
1:A:82:VAL:HG23	1:A:93:HIS:HD2	1.80	0.47
2:F:22:CYS:O	2:F:78:THR:HA	2.14	0.47
1:C:124:GLY:CA	1:C:126:MET:HE2	2.32	0.47
1:A:125:ILE:HG12	1:A:195[B]:VAL:HG21	1.96	0.47
1:E:131:LEU:HD11	1:E:192:ALA:HB1	1.97	0.47
1:A:78:LEU:C	1:A:78:LEU:HD23	2.40	0.47
1:A:120:MET:HE3	1:A:123:ILE:HD11	1.96	0.46
1:A:79:VAL:HG21	1:A:89:LEU:HD22	1.97	0.46
1:A:82:VAL:HA	1:A:93:HIS:CD2	2.51	0.46
2:D:38:ARG:HG2	2:D:48:VAL:CG2	2.45	0.46
1:A:129:ASP:CG	1:A:197:ARG:HH11	2.24	0.46
1:E:89:LEU:HD22	2:F:107:TYR:CZ	2.50	0.46
1:E:108:ASN:O	1:E:134:HIS:ND1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:168:LEU:HD23	1:G:176:PRO:HB3	1.97	0.45
1:C:80:GLY:HA3	1:C:95:GLU:OE2	2.15	0.45
1:C:141:ASN:HD21	1:C:160:LYS:H	1.64	0.45
1:E:78:LEU:HD13	1:E:98:TYR:CE2	2.45	0.45
2:H:68:PHE:CE1	2:H:83:MET:HB3	2.51	0.45
1:A:160:LYS:HG2	1:A:165:VAL:HG22	1.98	0.45
1:A:129:ASP:OD2	1:A:197:ARG:NH1	2.49	0.45
1:C:75:GLY:O	1:C:97:HIS:HD2	2.00	0.45
1:A:85:GLY:HA2	2:B:110:SER:OG	2.17	0.45
2:D:91:THR:O	2:D:92:ALA:HB2	2.17	0.44
2:H:117:GLN:NE2	2:H:119:THR:HG23	2.32	0.44
1:A:200:ASP:OD1	1:A:200:ASP:N	2.51	0.44
1:G:81:ARG:O	1:G:93:HIS:CD2	2.71	0.44
1:A:202:LEU:C	2:H:123:HIS:CD2	2.95	0.44
1:C:131:LEU:HD11	1:C:192:ALA:HB1	1.99	0.44
1:C:150:ASP:OD2	1:C:184:GLN:HG2	2.18	0.44
1:G:92:GLN:HG2	1:G:93:HIS:HD1	1.82	0.44
1:E:75:GLY:O	2:F:103:ARG:NH1	2.48	0.43
1:G:168:LEU:HA	1:G:169:PRO:HD2	1.89	0.43
2:H:30:SER:HA	2:H:74:ASN:HD21	1.81	0.43
1:E:107:PRO:HG2	1:E:193:VAL:HG11	2.00	0.43
1:G:108:ASN:O	1:G:134:HIS:HB2	2.18	0.43
1:G:163:ASN:CG	2:H:42:GLY:HA2	2.44	0.43
1:A:108:ASN:ND2	1:A:108:ASN:N	2.67	0.43
2:B:51:ILE:HG12	2:B:55:GLY:HA2	2.00	0.43
1:C:77:PRO:HG3	2:D:102:TYR:CE2	2.54	0.43
2:F:12:VAL:CG2	2:F:86:LEU:HD13	2.46	0.43
2:H:73:ASP:OD1	2:H:73:ASP:C	2.62	0.43
2:D:33:ALA:O	2:D:34:MET:HE2	2.19	0.43
2:B:46:ARG:NH1	2:B:61:GLU:OE1	2.39	0.42
1:C:83:ALA:O	1:C:84:ALA:C	2.62	0.42
2:D:33:ALA:O	2:D:34:MET:CE	2.67	0.42
1:E:74:GLU:O	1:E:97:HIS:CB	2.65	0.42
2:B:4:LEU:HD11	2:B:98:VAL:HG23	2.01	0.42
2:H:2:VAL:HG23	2:H:27:SER:HB2	2.02	0.42
2:B:18:LEU:HD23	2:B:83:MET:CE	2.50	0.42
2:B:18:LEU:HD23	2:B:83:MET:HE1	2.01	0.42
1:E:107:PRO:HG2	1:E:193:VAL:CG1	2.49	0.42
2:D:1:GLN:HA	2:D:1:GLN:OE1	2.20	0.42
2:F:91:THR:HG23	2:F:119:THR:HA	2.01	0.42
2:H:2:VAL:HA	2:H:26:GLY:HA3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:ASN:ND2	1:C:160:LYS:H	2.18	0.41
2:D:73:ASP:OD2	2:D:76:GLU:HB2	2.20	0.41
1:G:175:LYS:HB2	1:G:175:LYS:HE2	1.69	0.41
2:D:12:VAL:CG2	2:D:13:GLN:N	2.83	0.41
1:C:146:VAL:HG22	1:C:155:VAL:HG22	2.02	0.41
2:F:22:CYS:O	2:F:78:THR:HG22	2.21	0.41
1:G:104:LEU:HD23	1:G:104:LEU:HA	1.97	0.41
1:C:201:TRP:O	2:F:123:HIS:HD2	2.03	0.41
2:F:12:VAL:O	2:F:120:VAL:HA	2.21	0.41
2:F:76:GLU:O	2:F:77:ASN:HB2	2.21	0.41
2:H:18:LEU:HB3	2:H:83:MET:HE3	2.03	0.41
1:G:126:MET:O	1:G:127:ASP:C	2.63	0.40
1:A:106:LYS:HA	1:A:107:PRO:HA	1.84	0.40
1:A:146:VAL:HG22	1:A:155:VAL:HG22	2.02	0.40
2:D:109:GLY:C	4:D:303:HOH:O	2.64	0.40
2:B:6:GLU:HG3	2:B:96:CYS:HB2	2.02	0.40
2:D:50:ALA:O	2:D:58:THR:HA	2.21	0.40
1:G:122:ASP:HB2	1:G:170:GLU:O	2.21	0.40
2:H:35:ALA:HB1	2:H:47:LEU:CD1	2.51	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	131/202 (65%)	123 (94%)	8 (6%)	0	100	100
1	C	130/202 (64%)	124 (95%)	6 (5%)	0	100	100
1	E	128/202 (63%)	121 (94%)	7 (6%)	0	100	100
1	G	128/202 (63%)	117 (91%)	11 (9%)	0	100	100
2	B	121/132 (92%)	118 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	122/132 (92%)	115 (94%)	7 (6%)	0	100	100
2	F	123/132 (93%)	120 (98%)	3 (2%)	0	100	100
2	H	122/132 (92%)	118 (97%)	4 (3%)	0	100	100
All	All	1005/1336 (75%)	956 (95%)	49 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	114/169 (68%)	106 (93%)	8 (7%)	14	24
1	C	113/169 (67%)	106 (94%)	7 (6%)	16	29
1	E	111/169 (66%)	104 (94%)	7 (6%)	16	29
1	G	111/169 (66%)	103 (93%)	8 (7%)	13	23
2	B	98/106 (92%)	95 (97%)	3 (3%)	35	57
2	D	99/106 (93%)	93 (94%)	6 (6%)	17	30
2	F	100/106 (94%)	91 (91%)	9 (9%)	9	15
2	H	99/106 (93%)	88 (89%)	11 (11%)	6	9
All	All	845/1100 (77%)	786 (93%)	59 (7%)	14	24

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	LYS
1	A	108	ASN
1	A	152	GLU
1	A	165	VAL
1	A	195[A]	VAL
1	A	195[B]	VAL
1	A	200	ASP
1	A	202	LEU

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Mol	Chain	Res	Type
2	B	13	GLN
2	B	80	SER
2	B	121	SER
1	C	95	GLU
1	C	104	LEU
1	C	178	VAL
1	C	185	SER
1	C	195[A]	VAL
1	C	195[B]	VAL
1	C	202	LEU
2	D	44	VAL
2	D	69	THR
2	D	77	ASN
2	D	79	VAL
2	D	85	SER
2	D	89	GLU
1	E	73	GLU
1	E	88	LEU
1	E	106	LYS
1	E	108	ASN
1	E	152	GLU
1	E	195	VAL
1	E	202	LEU
2	F	2	VAL
2	F	13	GLN
2	F	25	SER
2	F	28	ILE
2	F	44	VAL
2	F	48	VAL
2	F	56	ASP
2	F	85	SER
2	F	124	HIS
1	G	99	GLN
1	G	106	LYS
1	G	139	VAL
1	G	149	ILE
1	G	164	LYS
1	G	175	LYS
1	G	177	ILE
1	G	185	SER
2	H	2	VAL
2	H	5	VAL

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Mol	Chain	Res	Type
2	H	25	SER
2	H	28	ILE
2	H	30	SER
2	H	51	ILE
2	H	56	ASP
2	H	61	GLU
2	H	64	VAL
2	H	79	VAL
2	H	122	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	HIS
1	A	108	ASN
2	B	77	ASN
2	B	82	GLN
2	B	117	GLN
1	C	92	GLN
1	C	97	HIS
2	D	74	ASN
2	D	77	ASN
1	E	161	GLN
2	F	3	GLN
2	F	13	GLN
2	F	82	GLN
2	F	123	HIS
1	G	161	GLN
2	H	13	GLN
2	H	117	GLN
2	H	123	HIS
2	H	124	HIS

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

6 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	EDO	A	302	-	3,3,3	0.07	0	2,2,2	0.17	0
3	EDO	C	301	-	3,3,3	0.08	0	2,2,2	0.08	0
3	EDO	C	303	-	3,3,3	0.14	0	2,2,2	0.26	0
3	EDO	A	301	-	3,3,3	0.11	0	2,2,2	0.11	0
3	EDO	C	304	-	3,3,3	0.12	0	2,2,2	0.22	0
3	EDO	C	302	-	3,3,3	0.13	0	2,2,2	0.53	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	302	-	-	0/1/1/1	-
3	EDO	C	301	-	-	1/1/1/1	-
3	EDO	C	303	-	-	0/1/1/1	-
3	EDO	A	301	-	-	1/1/1/1	-
3	EDO	C	304	-	-	1/1/1/1	-
3	EDO	C	302	-	-	0/1/1/1	-

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
3	A	301	EDO	O1-C1-C2-O2
3	C	304	EDO	O1-C1-C2-O2
3	C	301	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	302	EDO	1	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	132/202 (65%)	-0.27	2 (1%) 72 68	13, 36, 71, 109	1 (0%)
1	C	130/202 (64%)	-0.20	1 (0%) 82 80	12, 39, 65, 92	2 (1%)
1	E	130/202 (64%)	0.08	5 (3%) 44 40	26, 48, 78, 104	0
1	G	129/202 (63%)	0.11	1 (0%) 82 80	23, 48, 76, 98	1 (0%)
2	B	123/132 (93%)	-0.44	2 (1%) 70 66	22, 33, 56, 91	0
2	D	124/132 (93%)	0.15	1 (0%) 82 80	28, 45, 76, 106	0
2	F	125/132 (94%)	0.05	5 (4%) 42 38	24, 40, 83, 106	0
2	H	124/132 (93%)	0.19	5 (4%) 42 38	31, 49, 82, 104	0
All	All	1017/1336 (76%)	-0.04	22 (2%) 62 58	12, 43, 78, 109	4 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	124	HIS	4.2
1	E	92	GLN	3.6
2	B	1	GLN	3.4
2	F	29	PHE	3.0
2	F	125	HIS	2.8
2	H	27	SER	2.8
2	H	28	ILE	2.7
2	F	124	HIS	2.6
2	F	28	ILE	2.6
1	C	73	GLU	2.6
1	E	90	ALA	2.6
2	H	29	PHE	2.3
1	E	73	GLU	2.3
1	A	81	ARG	2.2
1	G	82	VAL	2.2
2	H	124	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	94	ILE	2.1
1	E	182	ARG	2.1
1	A	164	LYS	2.1
2	H	1	GLN	2.1
2	F	3	GLN	2.1
2	B	2	VAL	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	EDO	A	301	4/4	0.74	0.19	62,66,69,74	0
3	EDO	C	301	4/4	0.78	0.27	56,61,63,66	0
3	EDO	C	302	4/4	0.79	0.16	48,51,52,55	0
3	EDO	A	302	4/4	0.82	0.17	52,57,63,65	0
3	EDO	C	303	4/4	0.91	0.15	51,53,58,58	0
3	EDO	C	304	4/4	0.92	0.11	57,58,64,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.