



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:41 PM UTC

PDB ID : 5GWV / pdb_00005gww
Title : Structure of MoeN5-Sso7d fusion protein in complex with a substrate analogue
Authors : Ko, T.-P.; Guo, R.-T.; Chen, C.-C.
Deposited on : 2016-09-14
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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

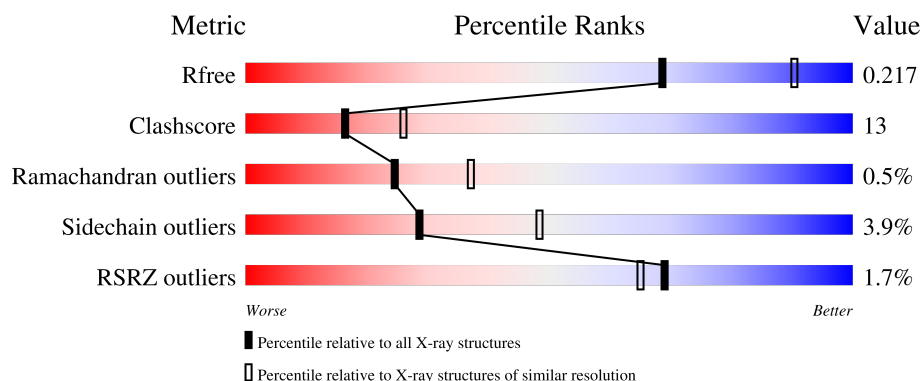
1 Overall quality at a glance

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	343	2% 59% 15% • 23%
1	B	343	76% 18% • •
1	C	343	57% 16% • 24%
1	D	343	4% 72% 23% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10016 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 MoeN5,DNA-binding protein 7d.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2013	1243	374	385	11			
1	B	332	Total	C	N	O	S	0	0	0
			2533	1571	463	485	14			
1	C	262	Total	C	N	O	S	0	0	0
			1998	1234	372	381	11			
1	D	333	Total	C	N	O	S	0	0	0
			2544	1579	466	485	14			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP A0A010
A	-12	ALA	-	expression tag	UNP A0A010
A	-11	HIS	-	expression tag	UNP A0A010
A	-10	HIS	-	expression tag	UNP A0A010
A	-9	HIS	-	expression tag	UNP A0A010
A	-8	HIS	-	expression tag	UNP A0A010
A	-7	HIS	-	expression tag	UNP A0A010
A	-6	HIS	-	expression tag	UNP A0A010
A	-5	VAL	-	expression tag	UNP A0A010
A	-4	ASP	-	expression tag	UNP A0A010
A	-3	ASP	-	expression tag	UNP A0A010
A	-2	ASP	-	expression tag	UNP A0A010
A	-1	ASP	-	expression tag	UNP A0A010
A	0	LYS	-	expression tag	UNP A0A010
A	261	ALA	-	linker	UNP A0A010
A	262	GLY	-	linker	UNP A0A010
A	263	ALA	-	linker	UNP A0A010
A	264	GLY	-	linker	UNP A0A010
A	265	ALA	-	linker	UNP A0A010
B	-13	MET	-	expression tag	UNP A0A010
B	-12	ALA	-	expression tag	UNP A0A010

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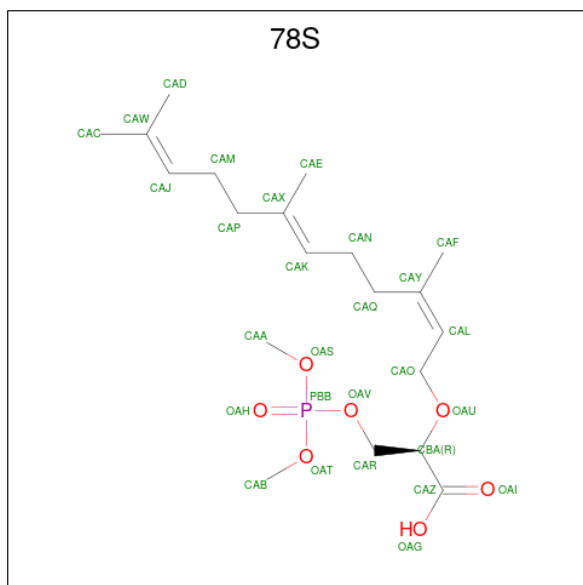
Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	HIS	-	expression tag	UNP A0A010
B	-10	HIS	-	expression tag	UNP A0A010
B	-9	HIS	-	expression tag	UNP A0A010
B	-8	HIS	-	expression tag	UNP A0A010
B	-7	HIS	-	expression tag	UNP A0A010
B	-6	HIS	-	expression tag	UNP A0A010
B	-5	VAL	-	expression tag	UNP A0A010
B	-4	ASP	-	expression tag	UNP A0A010
B	-3	ASP	-	expression tag	UNP A0A010
B	-2	ASP	-	expression tag	UNP A0A010
B	-1	ASP	-	expression tag	UNP A0A010
B	0	LYS	-	expression tag	UNP A0A010
B	261	ALA	-	linker	UNP A0A010
B	262	GLY	-	linker	UNP A0A010
B	263	ALA	-	linker	UNP A0A010
B	264	GLY	-	linker	UNP A0A010
B	265	ALA	-	linker	UNP A0A010
C	-13	MET	-	expression tag	UNP A0A010
C	-12	ALA	-	expression tag	UNP A0A010
C	-11	HIS	-	expression tag	UNP A0A010
C	-10	HIS	-	expression tag	UNP A0A010
C	-9	HIS	-	expression tag	UNP A0A010
C	-8	HIS	-	expression tag	UNP A0A010
C	-7	HIS	-	expression tag	UNP A0A010
C	-6	HIS	-	expression tag	UNP A0A010
C	-5	VAL	-	expression tag	UNP A0A010
C	-4	ASP	-	expression tag	UNP A0A010
C	-3	ASP	-	expression tag	UNP A0A010
C	-2	ASP	-	expression tag	UNP A0A010
C	-1	ASP	-	expression tag	UNP A0A010
C	0	LYS	-	expression tag	UNP A0A010
C	261	ALA	-	linker	UNP A0A010
C	262	GLY	-	linker	UNP A0A010
C	263	ALA	-	linker	UNP A0A010
C	264	GLY	-	linker	UNP A0A010
C	265	ALA	-	linker	UNP A0A010
D	-13	MET	-	expression tag	UNP A0A010
D	-12	ALA	-	expression tag	UNP A0A010
D	-11	HIS	-	expression tag	UNP A0A010
D	-10	HIS	-	expression tag	UNP A0A010
D	-9	HIS	-	expression tag	UNP A0A010
D	-8	HIS	-	expression tag	UNP A0A010

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	HIS	-	expression tag	UNP A0A010
D	-6	HIS	-	expression tag	UNP A0A010
D	-5	VAL	-	expression tag	UNP A0A010
D	-4	ASP	-	expression tag	UNP A0A010
D	-3	ASP	-	expression tag	UNP A0A010
D	-2	ASP	-	expression tag	UNP A0A010
D	-1	ASP	-	expression tag	UNP A0A010
D	0	LYS	-	expression tag	UNP A0A010
D	261	ALA	-	linker	UNP A0A010
D	262	GLY	-	linker	UNP A0A010
D	263	ALA	-	linker	UNP A0A010
D	264	GLY	-	linker	UNP A0A010
D	265	ALA	-	linker	UNP A0A010

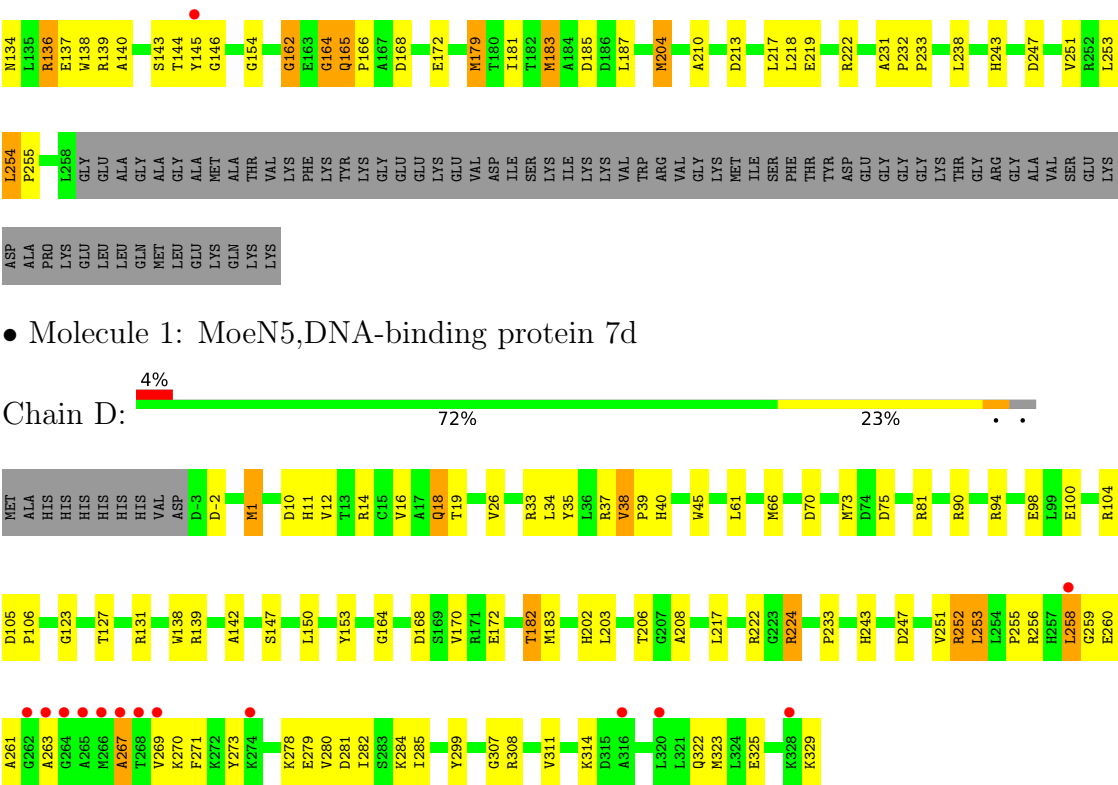
- Molecule 2 is (2R)-3-dimethoxyphosphoryloxy-2-[(2Z,6E)-3,7,11-trimethyldodeca-2,6,10-trienoxy]propanoic acid (CCD ID: 78S) (formula: C₂₀H₃₅O₇P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	O	P	0	0
			28	20	7	1		
2	D	1	Total	C	O	P	0	0
			18	10	7	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	223	Total 223	O 223	0	0
3	B	243	Total 243	O 243	0	0
3	C	175	Total 175	O 175	0	0
3	D	241	Total 241	O 241	0	0



• Molecule 1: MoeN5,DNA-binding protein 7d

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	137.94Å 217.70Å 104.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.40 25.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.6 (25.00-2.40) 95.2 (25.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.39Å)	Xtriage
Refinement program	CNS 1.21	Depositor
R, R_{free}	0.174 , 0.216 0.174 , 0.217	Depositor DCC
R_{free} test set	2976 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.8	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	10016	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 78S

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	0.79	1/2045 (0.0%)	1.06	9/2780 (0.3%)
1	B	0.75	1/2572 (0.0%)	1.01	7/3478 (0.2%)
1	C	0.79	2/2030 (0.1%)	1.04	4/2759 (0.1%)
1	D	0.74	2/2583 (0.1%)	1.06	10/3489 (0.3%)
All	All	0.76	6/9230 (0.1%)	1.04	30/12506 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	D	0	1
All	All	0	2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	MET	SD-CE	7.65	1.98	1.79
1	C	183	MET	SD-CE	6.50	1.95	1.79
1	D	66	MET	SD-CE	6.41	1.95	1.79
1	A	16	VAL	CA-CB	5.29	1.61	1.54
1	D	26	VAL	CA-CB	5.24	1.60	1.54
1	C	204	MET	SD-CE	-5.04	1.67	1.79

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	-2	ASP	N-CA-C	-10.52	99.84	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	253	LEU	N-CA-C	9.94	121.87	111.14
1	A	253	LEU	N-CA-C	7.86	121.35	111.69
1	A	33	ARG	N-CA-C	7.68	119.34	110.97
1	C	83	GLU	CB-CA-C	-7.40	98.97	110.81
1	B	253	LEU	N-CA-C	6.68	119.55	111.40
1	D	147	SER	N-CA-C	6.65	118.33	111.14
1	D	258	LEU	N-CA-C	-6.47	99.09	109.96
1	B	285	ILE	N-CA-C	6.23	117.34	108.93
1	B	265	ALA	N-CA-C	6.23	120.04	110.20
1	A	163	GLU	N-CA-C	6.15	118.46	110.53
1	A	104	ARG	N-CA-C	-6.09	104.55	111.07
1	D	33	ARG	N-CA-C	6.04	117.53	111.07
1	B	251	VAL	N-CA-C	5.97	116.61	110.82
1	C	104	ARG	N-CA-C	-5.83	104.83	111.07
1	A	147	SER	N-CA-C	5.82	117.29	111.07
1	B	190	TYR	N-CA-C	-5.79	104.88	111.07
1	A	164	GLY	N-CA-C	-5.78	105.59	115.08
1	D	38	VAL	CA-C-N	-5.70	113.74	119.56
1	D	38	VAL	C-N-CA	-5.70	113.74	119.56
1	D	278	LYS	N-CA-C	5.54	117.67	109.69
1	C	253	LEU	N-CA-C	5.53	120.01	112.04
1	A	251	VAL	N-CA-C	5.53	116.18	110.82
1	A	188	THR	N-CA-C	-5.42	106.72	113.55
1	B	163	GLU	N-CA-C	5.32	118.35	110.48
1	D	26	VAL	N-CA-C	-5.25	105.38	110.42
1	D	252	ARG	N-CA-C	5.15	116.75	111.03
1	C	154	GLY	N-CA-C	-5.12	106.56	112.50
1	B	104	ARG	N-CA-C	-5.07	105.66	111.14
1	A	122	GLY	N-CA-C	-5.03	106.63	113.37

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	TYR	Sidechain
1	D	153	TYR	Sidechain

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	2013	0	1996	49	0
1	B	2533	0	2532	54	0
1	C	1998	0	1983	55	0
1	D	2544	0	2554	83	0
2	D	46	0	0	1	0
3	A	223	0	0	9	0
3	B	243	0	0	10	0
3	C	175	0	0	4	0
3	D	241	0	0	8	0
All	All	10016	0	9065	237	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 13.

All (237) 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:C:136:ARG:HB3	1:C:136:ARG:HH11	1.07	1.10
1:D:325:GLU:HB2	1:D:329:LYS:HD3	1.48	0.92
1:C:136:ARG:HB3	1:C:136:ARG:NH1	1.86	0.88
1:A:139:ARG:HD3	1:A:179:MET:HE1	1.59	0.85
1:C:131:ARG:HH11	1:C:131:ARG:HG2	1.42	0.82
1:C:136:ARG:HH11	1:C:136:ARG:CB	1.89	0.82
1:C:1:MET:HE2	3:C:489:HOH:O	1.79	0.81
1:D:131:ARG:HG2	1:D:131:ARG:HH11	1.48	0.79
1:C:219:GLU:OE1	1:C:222:ARG:NE	2.16	0.79
1:B:222:ARG:HH11	1:B:243:HIS:HD2	1.29	0.77
1:D:202:HIS:O	1:D:206:THR:HG22	1.84	0.77
1:B:284:LYS:HE3	1:B:284:LYS:HA	1.67	0.76
1:A:38:VAL:HG12	1:A:39:PRO:HD3	1.69	0.75
1:A:186:ASP:OD1	1:A:200:LEU:N	2.19	0.75
1:D:75:ASP:HB2	1:D:81:ARG:NH1	2.01	0.75
1:D:222:ARG:HH11	1:D:243:HIS:HD2	1.35	0.74
1:B:106:PRO:HA	3:B:527:HOH:O	1.88	0.73
1:A:192:ARG:HB2	3:A:403:HOH:O	1.88	0.72
1:B:295:ILE:HD12	1:B:324:LEU:HD11	1.70	0.72
1:B:187:LEU:HD21	1:B:204:MET:HE1	1.70	0.72
1:C:222:ARG:HH11	1:C:243:HIS:HD2	1.38	0.72
1:D:123:GLY:O	1:D:127:THR:HG23	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ALA:HB2	3:B:539:HOH:O	1.89	0.70
1:C:247:ASP:O	1:C:251:VAL:HG12	1.92	0.69
1:D:267:ALA:HB1	1:D:282:ILE:HB	1.73	0.69
1:D:70:ASP:OD1	3:D:601:HOH:O	2.10	0.68
1:C:222:ARG:NH1	1:C:243:HIS:HD2	1.90	0.68
1:B:273:TYR:HB3	1:B:278:LYS:HD2	1.76	0.68
1:D:222:ARG:NH1	1:D:243:HIS:HD2	1.91	0.67
1:A:94:ARG:HD2	1:A:98:GLU:OE2	1.95	0.67
1:D:259:GLY:HA2	1:D:322:GLN:HG3	1.77	0.67
1:D:1:MET:CE	1:D:37:ARG:HG3	2.24	0.67
1:C:187:LEU:HD21	1:C:204:MET:HE1	1.77	0.65
1:D:247:ASP:O	1:D:251:VAL:HG22	1.97	0.65
1:D:61:LEU:HD13	1:D:94:ARG:HG3	1.78	0.65
1:D:94:ARG:HD2	1:D:98:GLU:OE2	1.96	0.64
1:A:61:LEU:HD13	1:A:94:ARG:HG3	1.80	0.64
1:B:43:THR:HA	3:B:445:HOH:O	1.97	0.64
1:D:105:ASP:OD2	1:D:106:PRO:HD2	1.98	0.64
1:D:267:ALA:HB1	1:D:282:ILE:CG1	2.28	0.64
1:B:164:GLY:O	3:B:401:HOH:O	2.16	0.63
1:C:73:MET:HE1	1:C:124:GLN:NE2	2.14	0.63
1:C:222:ARG:HH11	1:C:243:HIS:CD2	2.16	0.63
1:D:282:ILE:HA	1:D:285:ILE:HD13	1.80	0.63
1:A:187:LEU:HD21	1:A:204:MET:HE1	1.79	0.63
1:B:115:GLN:NE2	3:B:404:HOH:O	2.32	0.63
1:B:303:GLY:C	1:B:305:LYS:H	2.06	0.62
1:A:195:GLU:O	1:A:196:ARG:HD2	2.00	0.62
1:A:192:ARG:N	3:A:403:HOH:O	2.31	0.62
1:D:10:ASP:O	1:D:14:ARG:HG3	2.00	0.62
1:A:200:LEU:HG	1:A:204:MET:HE2	1.82	0.61
1:A:189:ASP:HB3	1:A:195:GLU:OE1	1.99	0.61
1:A:107:LYS:O	1:A:107:LYS:HD3	2.01	0.61
1:C:44:GLU:HG2	1:C:238:LEU:HG	1.83	0.61
1:D:1:MET:HG2	1:D:40:HIS:CD2	2.35	0.61
1:D:138:TRP:CH2	1:D:182:THR:HG22	2.36	0.60
1:B:266:MET:HE3	1:B:266:MET:HA	1.83	0.60
1:D:104:ARG:HD3	3:D:743:HOH:O	2.01	0.59
1:C:145:TYR:CD1	1:C:146:GLY:N	2.70	0.59
1:D:271:PHE:HE2	1:D:273:TYR:HB2	1.67	0.59
1:D:269:VAL:O	1:D:279:GLU:HA	2.02	0.59
1:B:301:GLU:CD	1:B:305:LYS:HE3	2.28	0.59
1:B:321:LEU:O	1:B:324:LEU:HB2	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:224:ARG:HG2	1:D:224:ARG:HH21	1.67	0.59
1:D:94:ARG:HD3	1:D:98:GLU:HG3	1.84	0.59
1:D:271:PHE:CE2	1:D:273:TYR:HB2	2.38	0.59
1:D:142:ALA:HB1	1:D:182:THR:HG21	1.85	0.58
1:B:273:TYR:CE2	1:B:274:LYS:HD2	2.38	0.58
1:D:38:VAL:HG12	1:D:39:PRO:HD3	1.84	0.58
1:C:12:VAL:HG12	1:C:29:THR:HG21	1.84	0.58
1:B:179:MET:HE2	1:B:217:LEU:HD11	1.86	0.58
1:D:131:ARG:HG2	1:D:131:ARG:NH1	2.13	0.57
1:C:38:VAL:HG12	1:C:39:PRO:HD3	1.85	0.57
1:C:14:ARG:HD3	3:C:467:HOH:O	2.05	0.56
1:D:325:GLU:HB2	1:D:329:LYS:CD	2.29	0.56
1:D:267:ALA:HB1	1:D:282:ILE:CB	2.35	0.56
1:A:131:ARG:HG3	1:A:197:ASP:HB3	1.87	0.56
1:A:90:ARG:NH1	1:A:94:ARG:HB2	2.20	0.56
1:C:131:ARG:HH11	1:C:131:ARG:CG	2.18	0.56
1:D:325:GLU:HA	1:D:329:LYS:HB3	1.88	0.56
1:A:115:GLN:HG3	3:A:457:HOH:O	2.06	0.56
1:C:107:LYS:O	1:C:107:LYS:HD3	2.05	0.55
1:D:1:MET:HE3	1:D:37:ARG:HG3	1.88	0.55
1:A:187:LEU:CD2	1:A:204:MET:HE1	2.37	0.55
1:D:224:ARG:HG2	3:D:619:HOH:O	2.06	0.55
1:A:33:ARG:HG2	1:A:33:ARG:HH11	1.70	0.55
1:A:33:ARG:HG2	1:A:33:ARG:NH1	2.22	0.55
1:C:187:LEU:HD21	1:C:204:MET:CE	2.37	0.55
1:D:18:GLN:HG2	1:D:18:GLN:O	2.07	0.55
1:A:158:ALA:HA	1:A:165:GLN:HG2	1.88	0.55
1:A:187:LEU:HD21	1:A:204:MET:CE	2.37	0.55
1:D:222:ARG:HH11	1:D:243:HIS:CD2	2.20	0.54
1:B:104:ARG:HG3	1:B:105:ASP:N	2.22	0.54
1:D:1:MET:HG2	1:D:40:HIS:HD2	1.71	0.54
1:C:46:THR:HG21	1:C:51:ARG:HG2	1.89	0.54
1:B:187:LEU:HD21	1:B:204:MET:CE	2.37	0.54
1:C:81:ARG:HD3	1:D:81:ARG:NH2	2.23	0.54
1:D:252:ARG:NH1	1:D:253:LEU:HD21	2.23	0.54
1:A:138:TRP:CH2	1:A:183:MET:HG2	2.43	0.53
1:A:55:VAL:HG22	1:A:102:LEU:HD13	1.91	0.53
1:C:1:MET:SD	1:C:1:MET:C	2.92	0.53
1:A:179:MET:HE2	1:A:221:LEU:HD21	1.91	0.53
1:B:319:GLU:CD	1:B:319:GLU:H	2.17	0.53
1:B:139:ARG:HG3	1:B:179:MET:HE1	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:ARG:HD3	1:B:243:HIS:CD2	2.44	0.53
1:D:168:ASP:O	1:D:172:GLU:HG3	2.09	0.52
1:D:34:LEU:HD13	3:D:652:HOH:O	2.09	0.52
1:C:81:ARG:NH2	3:C:407:HOH:O	2.40	0.52
1:B:200:LEU:HG	1:B:204:MET:HE2	1.92	0.52
1:C:219:GLU:CD	1:C:222:ARG:HE	2.18	0.52
1:C:183:MET:HE1	1:C:218:LEU:HG	1.92	0.51
1:D:251:VAL:HG23	1:D:252:ARG:N	2.25	0.51
1:B:222:ARG:HH11	1:B:243:HIS:CD2	2.19	0.51
1:C:22:SER:N	1:C:83:GLU:OE1	2.38	0.51
1:C:19:THR:O	1:C:19:THR:HG22	2.10	0.51
1:D:206:THR:HG23	1:D:208:ALA:H	1.76	0.51
1:D:256:ARG:NH1	3:D:603:HOH:O	2.42	0.51
1:B:263:ALA:CB	3:B:539:HOH:O	2.52	0.50
1:C:73:MET:HE1	1:C:124:GLN:HE22	1.77	0.50
1:D:1:MET:HE1	1:D:37:ARG:HG3	1.93	0.50
1:D:105:ASP:OD2	1:D:106:PRO:CD	2.60	0.50
1:D:285:ILE:HG22	1:D:323:MET:HE1	1.94	0.50
1:A:14:ARG:HD2	3:A:531:HOH:O	2.12	0.50
1:D:90:ARG:HD3	1:D:90:ARG:C	2.37	0.49
1:A:192:ARG:CB	3:A:403:HOH:O	2.54	0.49
1:D:12:VAL:HG22	1:D:61:LEU:HD23	1.95	0.49
1:B:179:MET:CE	1:B:217:LEU:HD11	2.41	0.49
1:C:101:SER:HB2	3:D:699:HOH:O	2.12	0.49
1:D:75:ASP:HB2	1:D:81:ARG:HH12	1.72	0.49
1:A:195:GLU:C	1:A:196:ARG:HD2	2.38	0.49
1:B:229:VAL:HG13	1:B:235:ALA:O	2.13	0.49
1:B:131:ARG:HG3	1:B:197:ASP:HB3	1.95	0.48
1:C:104:ARG:HH11	1:C:162:GLY:HA2	1.78	0.48
1:C:166:PRO:HD3	1:C:233:PRO:HD2	1.96	0.48
1:C:210:ALA:HB3	1:C:213:ASP:OD2	2.14	0.48
1:B:272:LYS:HA	1:B:276:GLU:O	2.14	0.47
1:D:307:GLY:C	1:D:308:ARG:HG3	2.39	0.47
1:D:203:LEU:HA	1:D:206:THR:CG2	2.43	0.47
1:D:35:TYR:OH	2:D:501:78S:CAF	2.63	0.47
1:B:179:MET:HE3	1:B:217:LEU:HD21	1.96	0.47
1:B:301:GLU:OE1	1:B:305:LYS:HE3	2.14	0.47
1:C:131:ARG:CG	1:C:131:ARG:NH1	2.77	0.47
1:C:139:ARG:O	1:C:143:SER:HB2	2.14	0.47
1:A:163:GLU:O	3:A:401:HOH:O	2.21	0.46
1:D:267:ALA:HB1	1:D:282:ILE:HG13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:10:ASP:HB2	3:D:701:HOH:O	2.15	0.46
1:D:75:ASP:CG	1:D:75:ASP:O	2.59	0.46
1:C:254:LEU:N	1:C:255:PRO:CD	2.79	0.46
1:C:12:VAL:HG12	1:C:29:THR:CG2	2.45	0.45
1:D:280:VAL:HG21	1:D:299:TYR:CE2	2.51	0.45
1:B:295:ILE:CD1	1:B:324:LEU:HD11	2.44	0.45
1:A:38:VAL:CG1	1:A:39:PRO:HD3	2.44	0.45
1:B:271:PHE:HB3	1:B:311:VAL:CG1	2.46	0.45
1:C:138:TRP:CH2	1:C:183:MET:HG2	2.52	0.45
1:A:115:GLN:CG	3:A:457:HOH:O	2.63	0.45
1:C:134:ASN:OD1	1:C:137:GLU:HG3	2.15	0.45
1:D:164:GLY:O	1:D:233:PRO:HG3	2.16	0.45
1:D:75:ASP:OD2	1:D:81:ARG:HD2	2.17	0.45
1:B:272:LYS:HG3	1:B:276:GLU:O	2.16	0.45
1:D:183:MET:CE	1:D:217:LEU:HD23	2.46	0.45
1:D:255:PRO:O	1:D:258:LEU:O	2.35	0.45
1:B:201:ALA:HB1	1:B:257:HIS:CE1	2.52	0.45
1:D:251:VAL:HG23	1:D:252:ARG:H	1.80	0.45
1:D:259:GLY:O	1:D:260:GLU:HG2	2.16	0.45
1:A:44:GLU:O	1:A:44:GLU:HG3	2.17	0.44
1:A:76:ASP:OD1	1:A:76:ASP:N	2.50	0.44
1:C:90:ARG:HD3	1:C:90:ARG:C	2.42	0.44
1:D:269:VAL:HG12	1:D:270:LYS:N	2.32	0.44
1:C:38:VAL:CG1	1:C:39:PRO:HD3	2.47	0.44
1:D:183:MET:HE3	1:D:217:LEU:HD23	2.00	0.44
1:B:256:ARG:NH1	3:B:405:HOH:O	2.33	0.44
1:B:187:LEU:HD23	1:B:201:ALA:HB2	1.99	0.44
1:B:9:ARG:HD2	3:B:434:HOH:O	2.17	0.44
1:D:70:ASP:HA	1:D:73:MET:HE3	2.00	0.44
1:B:239:VAL:HB	1:B:240:PRO:HD3	1.99	0.44
1:D:138:TRP:HH2	1:D:182:THR:HG22	1.80	0.43
1:C:131:ARG:HG2	1:C:131:ARG:NH1	2.18	0.43
1:D:11:HIS:NE2	1:D:61:LEU:HD11	2.33	0.43
1:D:38:VAL:CG1	1:D:39:PRO:HD3	2.47	0.43
1:B:104:ARG:CG	1:B:105:ASP:N	2.81	0.43
1:D:281:ASP:HB3	1:D:284:LYS:HG2	2.01	0.43
1:A:33:ARG:HD3	3:A:422:HOH:O	2.19	0.43
1:A:165:GLN:HE21	1:A:165:GLN:HB2	1.71	0.43
1:C:168:ASP:O	1:C:172:GLU:HG3	2.18	0.43
1:C:231:ALA:HA	1:C:232:PRO:HD3	1.84	0.43
1:D:19:THR:HG22	1:D:19:THR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:ALA:CB	1:D:182:THR:HG21	2.48	0.43
1:A:90:ARG:C	1:A:90:ARG:HD3	2.43	0.43
1:B:261:ALA:C	1:B:263:ALA:H	2.25	0.43
1:B:286:LYS:HG3	1:B:299:TYR:HA	2.00	0.43
1:A:158:ALA:CB	1:A:165:GLN:HG2	2.49	0.43
1:A:201:ALA:HB1	1:A:257:HIS:CE1	2.54	0.43
1:C:34:LEU:HD11	1:C:181:ILE:HD12	2.01	0.43
1:B:19:THR:HG22	1:B:19:THR:O	2.19	0.43
1:B:267:ALA:O	1:B:282:ILE:HG12	2.19	0.42
1:B:222:ARG:HD3	1:B:243:HIS:HD2	1.83	0.42
1:C:164:GLY:N	3:C:404:HOH:O	2.38	0.42
1:D:271:PHE:HB3	1:D:311:VAL:CG1	2.49	0.42
1:A:62:ASP:O	1:A:65:SER:HB3	2.20	0.42
1:D:271:PHE:HB3	1:D:311:VAL:HG12	2.01	0.42
1:A:25:LEU:HD22	1:A:83:GLU:HB3	2.02	0.42
1:B:72:LEU:HD21	1:B:85:ALA:HB2	2.01	0.42
1:C:25:LEU:HD22	1:C:83:GLU:HB3	2.00	0.42
1:D:34:LEU:CD1	3:D:652:HOH:O	2.67	0.42
1:A:134:ASN:OD1	1:A:137:GLU:HG3	2.20	0.42
1:B:90:ARG:HD3	1:B:90:ARG:C	2.44	0.42
1:B:269:VAL:HG21	1:B:285:ILE:HD11	2.01	0.42
1:B:303:GLY:O	1:B:305:LYS:N	2.51	0.42
1:D:12:VAL:O	1:D:16:VAL:HG23	2.20	0.42
1:A:193:ASN:HB3	1:A:195:GLU:HG3	2.01	0.42
1:D:75:ASP:CB	1:D:81:ARG:NH1	2.79	0.42
1:A:131:ARG:HG2	1:A:131:ARG:HH11	1.85	0.41
1:C:140:ALA:O	1:C:144:THR:HG23	2.20	0.41
1:A:192:ARG:HA	3:A:428:HOH:O	2.19	0.41
1:C:179:MET:CE	1:C:217:LEU:HD11	2.49	0.41
1:C:139:ARG:HE	1:C:179:MET:HE1	1.86	0.41
1:C:165:GLN:HE21	1:C:165:GLN:HB3	1.58	0.41
1:D:150:LEU:HD12	1:D:150:LEU:HA	1.90	0.41
1:B:243:HIS:HE1	3:B:553:HOH:O	2.02	0.41
1:B:163:GLU:HA	3:B:472:HOH:O	2.19	0.41
1:C:139:ARG:HG3	1:C:179:MET:HE1	2.02	0.41
1:A:81:ARG:HH11	1:A:81:ARG:HG2	1.85	0.41
1:C:100:GLU:OE1	1:D:100:GLU:OE1	2.39	0.41
1:A:81:ARG:NH2	1:B:72:LEU:O	2.54	0.41
1:A:84:LEU:HD23	1:A:84:LEU:HA	1.95	0.41
1:B:32:LEU:HD22	1:B:35:TYR:CD2	2.55	0.41
1:C:46:THR:CG2	1:C:51:ARG:HG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:TRP:CE3	1:D:170:VAL:HG22	2.56	0.41
1:D:203:LEU:C	1:D:206:THR:HG22	2.46	0.41
1:B:105:ASP:CG	1:B:106:PRO:HD2	2.46	0.40
1:A:205:ARG:NH1	1:D:131:ARG:HD3	2.36	0.40
1:A:239:VAL:N	1:A:240:PRO:HD2	2.37	0.40
1:C:32:LEU:HB3	1:C:35:TYR:CD2	2.56	0.40
1:D:104:ARG:HH21	1:D:104:ARG:HG3	1.86	0.40
1:A:138:TRP:CZ3	1:A:183:MET:HG2	2.56	0.40
1:A:150:LEU:HD12	1:A:150:LEU:HA	1.86	0.40
1:B:321:LEU:O	1:B:325:GLU:OE1	2.39	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	262/343 (76%)	256 (98%)	6 (2%)	0	100	100
1	B	330/343 (96%)	313 (95%)	16 (5%)	1 (0%)	36	50
1	C	260/343 (76%)	251 (96%)	7 (3%)	2 (1%)	16	25
1	D	331/343 (96%)	311 (94%)	17 (5%)	3 (1%)	14	22
All	All	1183/1372 (86%)	1131 (96%)	46 (4%)	6 (0%)	24	37

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	260	GLU
1	C	162	GLY
1	D	263	ALA
1	D	261	ALA
1	D	267	ALA

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Mol	Chain	Res	Type
1	C	164	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	208/270 (77%)	196 (94%)	12 (6%)	18	32
1	B	260/270 (96%)	248 (95%)	12 (5%)	24	41
1	C	206/270 (76%)	200 (97%)	6 (3%)	37	60
1	D	261/270 (97%)	255 (98%)	6 (2%)	44	66
All	All	935/1080 (87%)	899 (96%)	36 (4%)	28	48

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	ASP
1	A	18	GLN
1	A	29	THR
1	A	38	VAL
1	A	106	PRO
1	A	107	LYS
1	A	139	ARG
1	A	163	GLU
1	A	165	GLN
1	A	186	ASP
1	A	197	ASP
1	A	254	LEU
1	B	18	GLN
1	B	38	VAL
1	B	131	ARG
1	B	186	ASP
1	B	206	THR
1	B	254	LEU
1	B	266	MET
1	B	269	VAL

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Mol	Chain	Res	Type
1	B	284	LYS
1	B	288	VAL
1	B	293	LYS
1	B	314	LYS
1	C	52	ARG
1	C	136	ARG
1	C	165	GLN
1	C	179	MET
1	C	185	ASP
1	C	254	LEU
1	D	1	MET
1	D	18	GLN
1	D	139	ARG
1	D	182	THR
1	D	224	ARG
1	D	314	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	115	GLN
1	A	202	HIS
1	B	11	HIS
1	B	193	ASN
1	B	202	HIS
1	B	243	HIS
1	C	115	GLN
1	C	119	HIS
1	C	124	GLN
1	C	165	GLN
1	C	212	GLN
1	C	243	HIS
1	D	165	GLN
1	D	193	ASN
1	D	243	HIS
1	D	322	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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](#)

2 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$
2	78S	D	501	-	27,27,27	2.93	3 (11%)	30,34,34	2.13	8 (26%)
2	78S	D	502	-	17,17,27	2.01	1 (5%)	18,22,34	2.02	3 (16%)

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	78S	D	501	-	-	11/34/34/34	-
2	78S	D	502	-	-	8/22/22/34	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	78S	CAK-CAX	8.88	1.53	1.33
2	D	501	78S	CAL-CAY	8.82	1.53	1.33
2	D	502	78S	CAL-CAY	7.03	1.53	1.32
2	D	501	78S	CAJ-CAW	7.03	1.53	1.32

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	502	78S	CAO-CAL-CAY	-7.17	115.17	126.76
2	D	501	78S	CAO-CAL-CAY	-6.51	115.53	126.20
2	D	501	78S	CAN-CAK-CAX	-5.41	115.24	127.62
2	D	501	78S	CAM-CAJ-CAW	-3.77	115.09	127.64
2	D	501	78S	CAF-CAY-CAQ	3.24	120.84	115.23
2	D	501	78S	OAI-CAZ-CBA	-2.84	116.72	122.85
2	D	502	78S	OAI-CAZ-CBA	-2.59	117.27	122.85
2	D	502	78S	CAF-CAY-CAQ	2.38	120.06	114.59
2	D	501	78S	CAO-OAU-CBA	-2.32	108.97	113.15
2	D	501	78S	CAE-CAX-CAP	2.28	119.19	115.23
2	D	501	78S	CAD-CAW-CAC	2.24	119.75	114.59

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	501	78S	CAJ-CAM-CAP-CAX
2	D	501	78S	CAR-CBA-OAU-CAO
2	D	501	78S	OAG-CAZ-CBA-OAU
2	D	501	78S	OAV-CAR-CBA-OAU
2	D	501	78S	OAV-CAR-CBA-CAZ
2	D	502	78S	CAR-CBA-OAU-CAO
2	D	502	78S	OAI-CAZ-CBA-OAU
2	D	502	78S	OAG-CAZ-CBA-OAU
2	D	501	78S	CAA-OAS-PBB-OAV
2	D	501	78S	CAA-OAS-PBB-OAH
2	D	502	78S	CAA-OAS-PBB-OAH
2	D	502	78S	CAZ-CBA-OAU-CAO
2	D	501	78S	CAB-OAT-PBB-OAH
2	D	502	78S	OAV-CAR-CBA-OAU
2	D	501	78S	OAI-CAZ-CBA-OAU
2	D	502	78S	CBA-CAR-OAV-PBB
2	D	501	78S	CAB-OAT-PBB-OAV
2	D	502	78S	OAV-CAR-CBA-CAZ
2	D	501	78S	CAN-CAQ-CAY-CAF

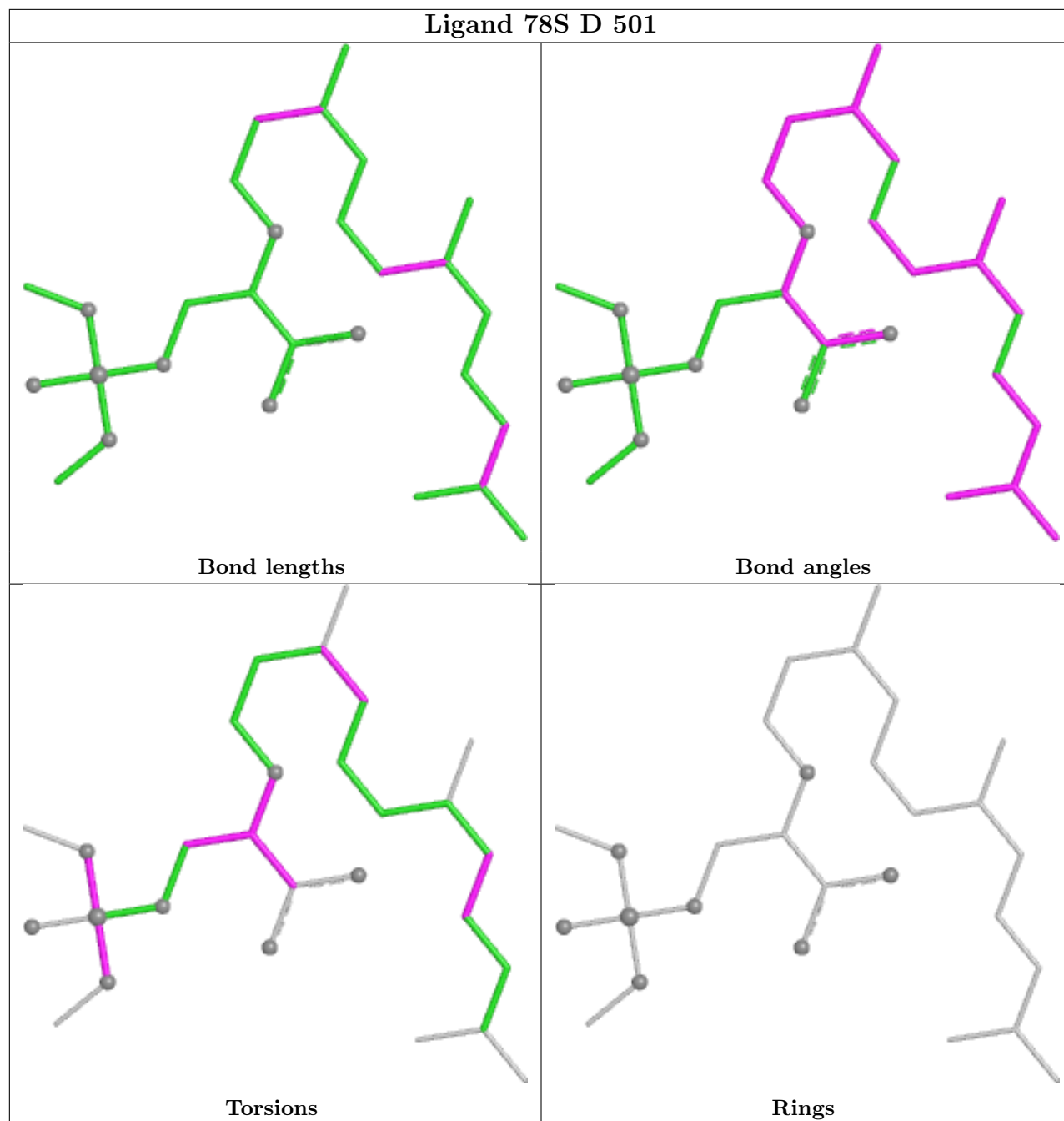
There are no ring outliers.

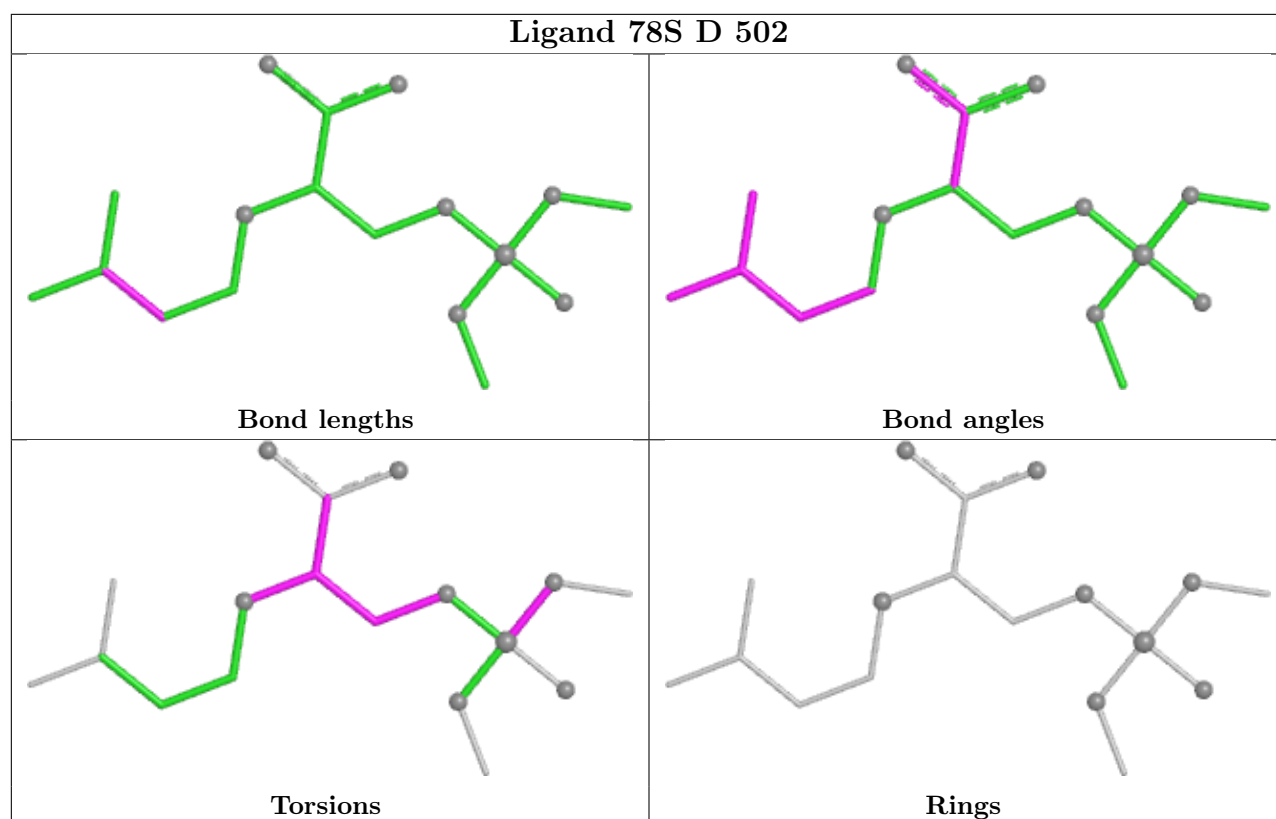
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	78S	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand 78S D 501





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	264/343 (76%)	-0.61	6 (2%) 61 57	27, 39, 72, 118	0
1	B	332/343 (96%)	-0.50	0 100 100	27, 43, 87, 96	0
1	C	262/343 (76%)	-0.64	1 (0%) 88 86	24, 41, 58, 98	0
1	D	333/343 (97%)	-0.32	13 (3%) 43 39	25, 42, 121, 143	0
All	All	1191/1372 (86%)	-0.51	20 (1%) 69 65	24, 41, 96, 143	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	263	ALA	3.6
1	A	164	GLY	3.5
1	D	265	ALA	3.1
1	D	262	GLY	2.5
1	D	264	GLY	2.5
1	A	190	TYR	2.5
1	D	269	VAL	2.5
1	A	194	GLY	2.4
1	D	258	LEU	2.4
1	A	192	ARG	2.4
1	D	266	MET	2.3
1	D	328	LYS	2.3
1	D	268	THR	2.2
1	A	198	GLY	2.2
1	C	145	TYR	2.2
1	D	274	LYS	2.2
1	A	-5	VAL	2.1
1	D	320	LEU	2.1
1	D	267	ALA	2.0
1	D	316	ALA	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](#)

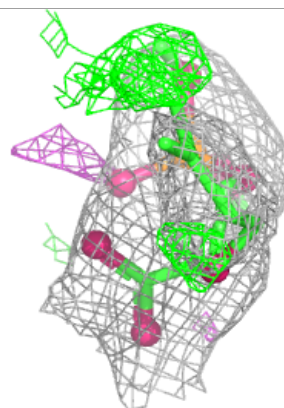
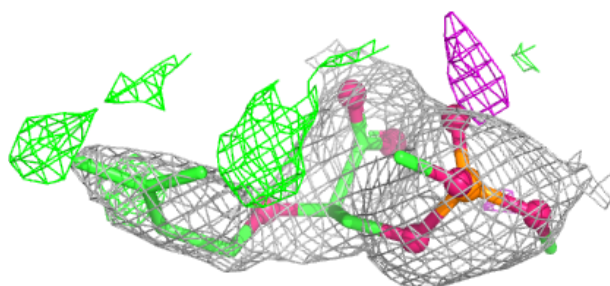
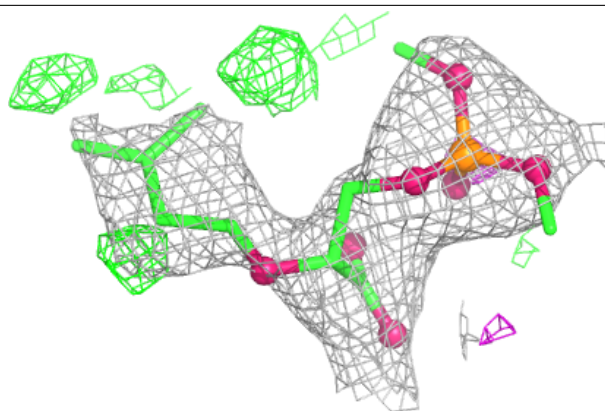
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	78S	D	502	18/28	0.72	0.19	100,108,111,112	0
2	78S	D	501	28/28	0.82	0.19	49,83,116,118	0

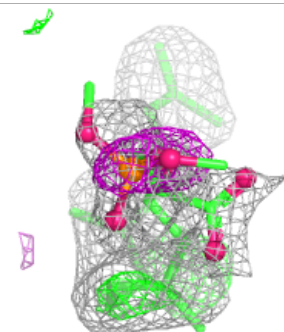
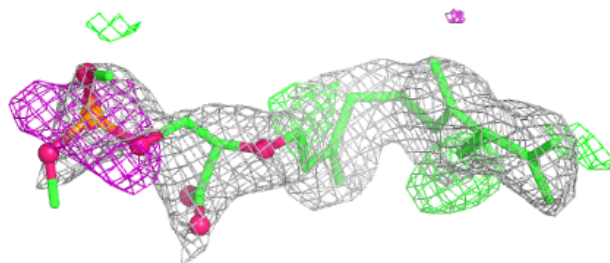
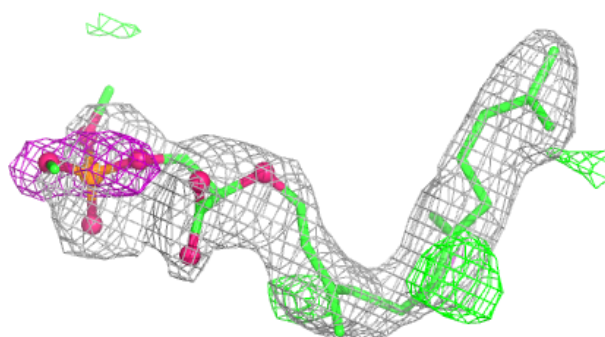
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 78S D 502:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around 78S D 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.