



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:36 PM UTC

PDB ID : 1T7D / pdb_00001t7d
Title : Crystal structure of Escherichia coli type I signal peptidase in complex with a lipopeptide inhibitor
Authors : Paetzel, M.; Goodall, J.J.; Kania, M.; Dalbey, R.E.; Page, M.G.P.
Deposited on : 2004-05-09
Resolution : 2.47 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

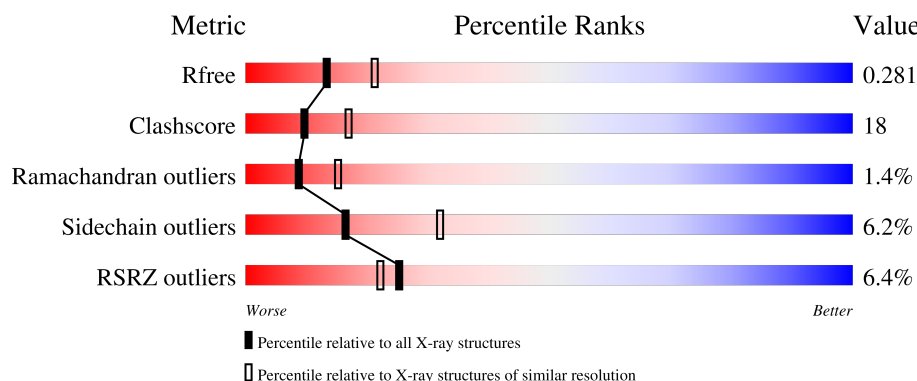
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	250	4% 58% 21% • 17%
1	B	250	7% 56% 28% • 11%
2	C	6	17% 67% 33%
2	D	6	17% 33% 67%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3766 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 SIGNAL PEPTIDASE I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1648	1058	274	309	7			
1	B	223	Total	C	N	O	S	0	0	0
			1755	1121	293	333	8			

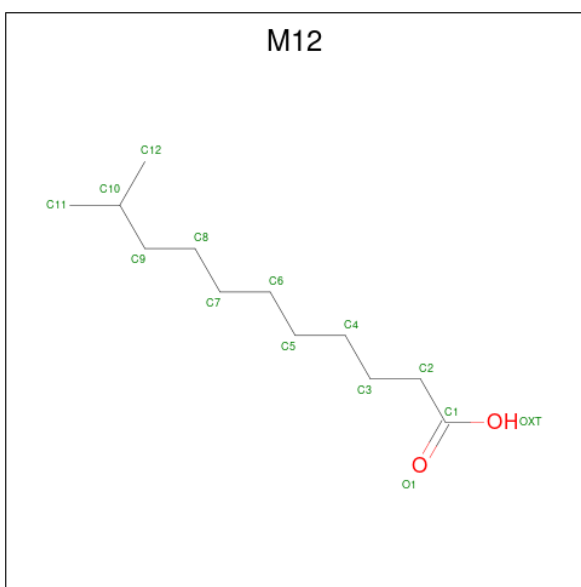
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	MET	-	initiating methionine	UNP P00803
B	74	MET	-	initiating methionine	UNP P00803

- Molecule 2 is a protein called ARYLOMYCIN A2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			46	30	6	10			
2	D	6	Total	C	N	O	0	0	0
			46	30	6	10			

- Molecule 3 is 10-METHYLUNDECANOIC ACID (CCD ID: M12) (formula: C₁₂H₂₄O₂).



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

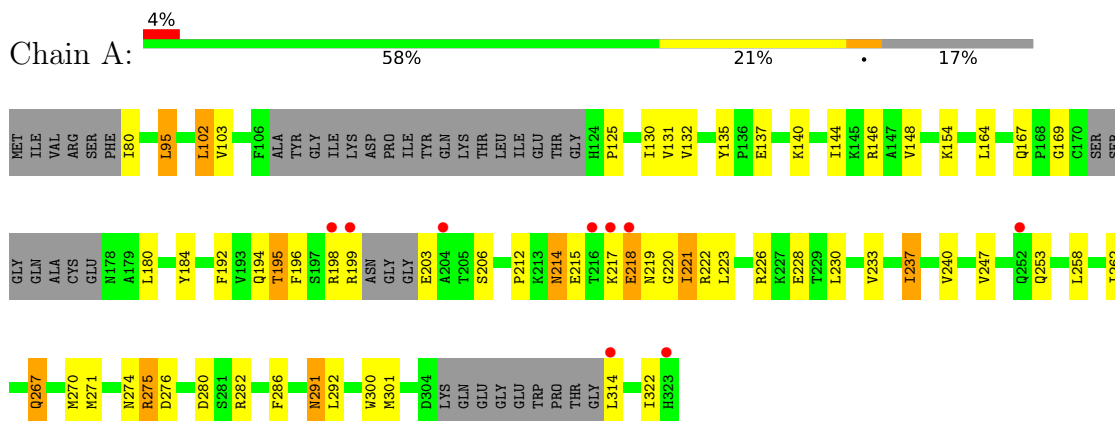
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	122	Total	O	0	0
			122	122		
4	B	135	Total	O	0	0
			135	135		
4	C	5	Total	O	0	0
			5	5		
4	D	3	Total	O	0	0
			3	3		

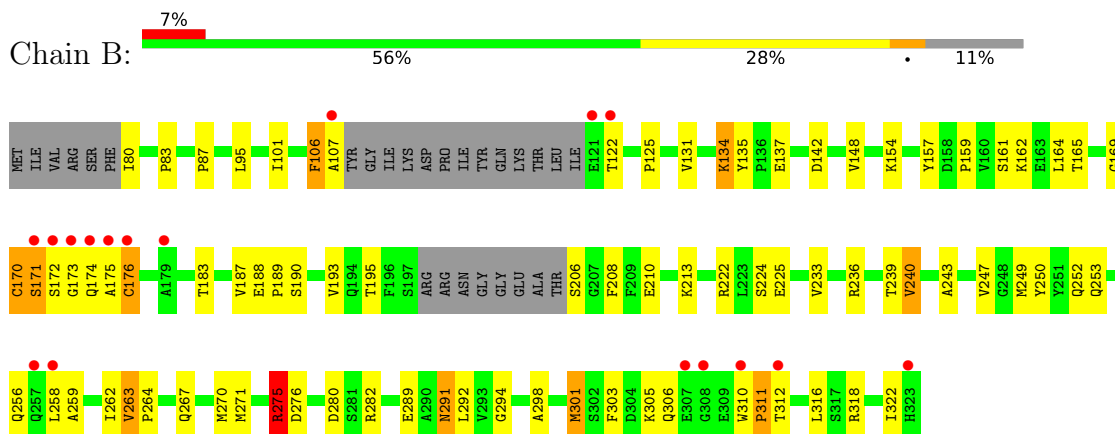
3 Residue-property plots

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.

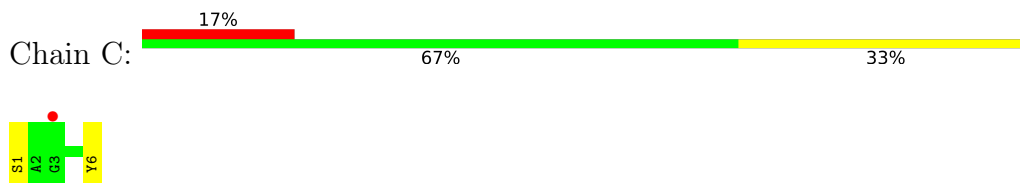
• Molecule 1: SIGNAL PEPTIDASE I



• Molecule 1: SIGNAL PEPTIDASE I



• Molecule 2: ARYLOMYCIN A2



• Molecule 2: ARYLOMYCIN A2





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	69.61Å 69.61Å 258.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.54 – 2.47 29.54 – 2.47	Depositor EDS
% Data completeness (in resolution range)	90.6 (29.54-2.47) 90.5 (29.54-2.47)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.48 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.283 0.231 , 0.281	Depositor DCC
R_{free} test set	2142 reflections (9.92%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.513	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.0	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	0.93	EDS
Total number of atoms	3766	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 6.54% 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: M12, DAL, DSE, 5PG

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	0.52	0/1689	0.98	8/2289 (0.3%)
1	B	0.48	0/1801	0.97	8/2443 (0.3%)
2	C	0.54	0/21	0.54	0/24
2	D	0.52	0/21	0.65	0/24
All	All	0.50	0/3532	0.98	16/4780 (0.3%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	ARG	N-CA-C	-8.36	87.58	111.00
1	A	275	ARG	N-CA-C	8.17	121.11	111.71
1	A	198	ARG	N-CA-C	6.80	119.83	108.13
1	A	140	LYS	N-CA-C	-6.64	105.14	113.18
1	B	176	CYS	N-CA-C	6.60	118.84	110.33
1	B	240	VAL	N-CA-C	-6.53	100.45	107.77
1	A	192	PHE	N-CA-C	6.46	120.06	109.72
1	A	198	ARG	CA-C-N	6.27	132.99	121.70
1	A	198	ARG	C-N-CA	6.27	132.99	121.70
1	A	274	ASN	N-CA-C	-6.04	98.05	108.56
1	B	311	PRO	N-CA-C	5.77	120.25	111.19
1	B	106	PHE	N-CA-C	-5.68	105.23	111.82
1	B	275	ARG	N-CA-C	5.55	122.62	110.80
1	B	210	GLU	N-CA-C	-5.11	101.44	109.25
1	B	298	ALA	N-CA-C	5.10	116.59	108.79
1	B	263	VAL	N-CA-C	5.01	113.86	108.95

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	1648	0	1606	52	0
1	B	1755	0	1701	67	0
2	C	46	0	36	1	0
2	D	46	0	36	3	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
4	A	122	0	0	5	0
4	B	135	0	0	6	0
4	C	5	0	0	0	0
4	D	3	0	0	0	0
All	All	3766	0	3379	122	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 18.

All (122) 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:253:GLN:HE22	1:A:262:ILE:H	1.05	1.01
1:B:253:GLN:HE22	1:B:262:ILE:H	1.07	0.96
1:A:322:ILE:HA	4:A:2122:HOH:O	1.78	0.82
1:A:218:GLU:H	1:A:218:GLU:CD	1.90	0.80
1:B:148:VAL:HG11	1:B:164:LEU:HD13	1.64	0.80
1:B:240:VAL:HG13	1:B:243:ALA:HB3	1.68	0.76
1:B:247:VAL:HG12	1:B:258:LEU:HD13	1.73	0.71
1:B:135:TYR:CE2	1:B:137:GLU:HB2	2.28	0.69
1:B:301:MET:HA	1:B:301:MET:HE2	1.74	0.68
1:B:240:VAL:HG12	1:B:276:ASP:HB2	1.74	0.68
1:B:253:GLN:NE2	1:B:262:ILE:H	1.89	0.68
1:B:253:GLN:HE22	1:B:262:ILE:N	1.86	0.68
1:B:225:GLU:OE1	1:B:236:ARG:HD2	1.96	0.66
1:A:154:LYS:HB3	1:A:167:GLN:HB2	1.77	0.66
1:B:134:LYS:HE2	1:B:142:ASP:OD1	1.96	0.66
1:B:291:ASN:HD22	1:B:291:ASN:N	1.93	0.66
1:B:240:VAL:HG12	1:B:276:ASP:CB	2.26	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HD13	1:A:300:TRP:HB2	1.79	0.65
1:A:212:PRO:C	1:A:214:ASN:H	2.02	0.64
1:A:291:ASN:HD22	1:A:291:ASN:N	1.95	0.64
1:A:271:MET:HE3	1:A:282:ARG:NE	2.14	0.63
1:B:170:CYS:HA	1:B:176:CYS:HB3	1.81	0.62
1:A:223:LEU:CD2	1:A:240:VAL:HG22	2.28	0.62
1:A:253:GLN:HE22	1:A:262:ILE:N	1.88	0.62
1:A:275:ARG:HG3	4:A:2105:HOH:O	1.99	0.62
1:A:135:TYR:CE2	1:A:137:GLU:HB3	2.37	0.60
1:A:267:GLN:HG2	1:A:286:PHE:CB	2.32	0.60
1:B:187:VAL:HG13	1:B:224:SER:HB3	1.84	0.60
1:A:130:ILE:HD11	1:A:146:ARG:NH2	2.17	0.59
1:B:131:VAL:HG11	1:B:292:LEU:HD22	1.85	0.59
1:B:107:ALA:HB2	1:B:122:THR:HA	1.84	0.58
1:A:80:ILE:N	1:A:80:ILE:HD12	2.19	0.58
1:B:301:MET:HE1	1:B:316:LEU:CD2	2.33	0.58
1:B:310:TRP:HB3	1:B:311:PRO:HD2	1.86	0.57
1:B:301:MET:HE1	1:B:316:LEU:HD23	1.85	0.57
1:A:212:PRO:C	1:A:214:ASN:N	2.63	0.57
1:A:130:ILE:HD11	1:A:146:ARG:CZ	2.35	0.56
1:A:130:ILE:CG2	1:A:144:ILE:HD11	2.36	0.56
1:A:226:ARG:HD3	4:A:2085:HOH:O	2.06	0.56
1:B:270:MET:O	1:B:280:ASP:HB2	2.06	0.56
1:B:125:PRO:HG2	1:B:292:LEU:HD13	1.88	0.55
1:B:174:GLN:HG2	1:B:175:ALA:H	1.71	0.55
1:A:131:VAL:HG11	1:A:292:LEU:HD22	1.87	0.55
1:B:174:GLN:O	1:B:176:CYS:N	2.40	0.54
1:A:154:LYS:HG2	1:A:262:ILE:CD1	2.38	0.54
1:A:130:ILE:HG22	1:A:144:ILE:HD11	1.89	0.54
1:A:233:VAL:HA	4:A:2089:HOH:O	2.08	0.54
1:A:267:GLN:HG2	1:A:286:PHE:HB3	1.88	0.54
1:A:195:THR:HG22	1:A:220:GLY:HA3	1.89	0.53
1:B:240:VAL:CG1	1:B:243:ALA:HB3	2.39	0.53
1:B:171:SER:C	1:B:173:GLY:H	2.16	0.52
1:A:223:LEU:HD21	1:A:240:VAL:HG22	1.91	0.52
1:A:154:LYS:HG2	1:A:262:ILE:HD13	1.91	0.52
1:A:148:VAL:HG11	1:A:164:LEU:HD13	1.92	0.51
1:B:174:GLN:HG2	1:B:175:ALA:N	2.26	0.51
1:B:154:LYS:HD2	4:B:2045:HOH:O	2.10	0.51
1:B:303:PHE:HB3	1:B:311:PRO:HB3	1.92	0.51
1:A:267:GLN:HG2	1:A:286:PHE:HB2	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ALA:HB3	1:B:122:THR:OG1	2.11	0.51
1:A:240:VAL:HG23	1:A:276:ASP:HB2	1.94	0.50
1:A:230:LEU:O	1:A:233:VAL:HG12	2.11	0.50
1:B:107:ALA:CB	1:B:122:THR:OG1	2.60	0.49
1:B:247:VAL:O	1:B:250:TYR:HB2	2.13	0.49
1:A:218:GLU:O	1:A:219:ASN:HB2	2.12	0.49
1:A:221:ILE:HD13	1:A:222:ARG:O	2.12	0.49
1:B:247:VAL:CG1	1:B:258:LEU:HD13	2.41	0.49
1:A:247:VAL:HG12	1:A:258:LEU:HD13	1.94	0.49
1:B:208:PHE:H	1:B:306:GLN:HE22	1.58	0.49
1:A:247:VAL:CG1	1:A:258:LEU:HB3	2.44	0.48
1:B:106:PHE:O	1:B:107:ALA:HB2	2.12	0.48
1:B:258:LEU:O	1:B:259:ALA:HB3	2.13	0.48
1:B:195:THR:O	1:B:206:SER:N	2.47	0.48
1:B:125:PRO:CG	1:B:292:LEU:HD13	2.44	0.48
1:A:102:LEU:HD13	1:A:300:TRP:CB	2.44	0.48
1:B:208:PHE:H	1:B:306:GLN:NE2	2.12	0.47
1:B:157:TYR:O	1:B:159:PRO:HD3	2.13	0.47
1:B:224:SER:HB2	1:B:239:THR:OG1	2.14	0.47
1:A:300:TRP:CD1	1:A:301:MET:HG3	2.50	0.47
1:B:206:SER:N	4:B:2078:HOH:O	2.47	0.46
1:B:169:GLY:C	1:B:171:SER:H	2.24	0.46
1:B:193:VAL:HG22	1:B:222:ARG:HA	1.96	0.46
1:A:218:GLU:CD	1:A:218:GLU:N	2.67	0.46
1:A:223:LEU:HD22	1:A:240:VAL:HG22	1.96	0.46
1:B:252:GLN:NE2	1:B:258:LEU:HD21	2.31	0.46
2:D:6:TYR:CD1	2:D:6:TYR:N	2.84	0.46
1:B:271:MET:HE3	1:B:282:ARG:NE	2.31	0.46
1:B:188:GLU:HG3	1:B:189:PRO:HD2	1.98	0.45
1:B:206:SER:N	4:B:2079:HOH:O	2.49	0.45
1:B:275:ARG:HG3	4:B:2114:HOH:O	2.16	0.45
1:A:103:VAL:HG21	1:A:132:VAL:HG21	1.99	0.45
1:A:130:ILE:N	1:A:130:ILE:HD12	2.31	0.45
1:A:203:GLU:N	4:A:2065:HOH:O	2.49	0.45
1:A:132:VAL:CG2	1:A:144:ILE:HD13	2.47	0.45
1:B:83:PRO:HA	1:B:101:ILE:O	2.17	0.45
2:D:4:5PG:HC1	2:D:5:ALA:N	2.32	0.45
1:B:263:VAL:HA	1:B:264:PRO:HD3	1.75	0.45
1:B:291:ASN:N	1:B:291:ASN:ND2	2.64	0.45
1:A:196:PHE:CD1	1:A:196:PHE:N	2.85	0.44
1:B:125:PRO:HG2	1:B:292:LEU:CD1	2.47	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:SER:CB	1:B:225:GLU:HB2	2.48	0.44
1:A:194:GLN:HE21	1:A:206:SER:HB3	1.82	0.44
1:B:310:TRP:CB	1:B:311:PRO:HD2	2.47	0.44
1:B:106:PHE:CD1	1:B:294:GLY:HA2	2.53	0.44
1:B:253:GLN:HB3	1:B:256:GLN:HB2	1.99	0.44
1:A:125:PRO:HG2	1:A:292:LEU:HD13	2.00	0.44
1:A:184:TYR:CE2	1:A:228:GLU:HG3	2.53	0.43
1:B:318:ARG:HH21	1:B:318:ARG:HG2	1.84	0.43
1:B:305:LYS:HG2	1:B:306:GLN:N	2.33	0.43
1:B:80:ILE:HD13	4:B:2022:HOH:O	2.18	0.43
1:A:95:LEU:HD11	1:A:322:ILE:HD12	2.00	0.42
1:A:267:GLN:HE21	1:A:267:GLN:HB2	1.47	0.42
1:A:253:GLN:NE2	1:A:262:ILE:H	1.90	0.42
1:B:161:SER:O	1:B:162:LYS:HB2	2.19	0.41
1:A:228:GLU:HB2	1:A:237:ILE:HD13	2.02	0.41
1:B:240:VAL:HG13	1:B:240:VAL:O	2.19	0.41
1:B:312:THR:O	1:B:312:THR:HG22	2.20	0.41
1:B:322:ILE:HA	4:B:2134:HOH:O	2.21	0.41
2:C:6:TYR:N	2:C:6:TYR:CD1	2.88	0.41
1:B:87:PRO:HG3	2:D:4:5PG:CD1	2.51	0.41
1:B:187:VAL:CG1	1:B:224:SER:HB3	2.50	0.41
1:A:270:MET:O	1:A:280:ASP:HB2	2.21	0.41
1:B:253:GLN:NE2	1:B:256:GLN:HG3	2.36	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	198/250 (79%)	184 (93%)	12 (6%)	2 (1%)	12	22
1	B	217/250 (87%)	193 (89%)	20 (9%)	4 (2%)	6	11

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	2/6 (33%)	2 (100%)	0	0	100	100
2	D	2/6 (33%)	2 (100%)	0	0	100	100
All	All	419/512 (82%)	381 (91%)	32 (8%)	6 (1%)	9	15

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	SER
1	A	169	GLY
1	B	275	ARG
1	A	215	GLU
1	B	172	SER
1	B	213	LYS

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	180/215 (84%)	168 (93%)	12 (7%)	15	29
1	B	191/215 (89%)	180 (94%)	11 (6%)	18	35
2	C	1/1 (100%)	1 (100%)	0	100	100
2	D	1/1 (100%)	1 (100%)	0	100	100
All	All	373/432 (86%)	350 (94%)	23 (6%)	16	32

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LEU
1	A	102	LEU
1	A	180	LEU
1	A	195	THR
1	A	214	ASN
1	A	217	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	218	GLU
1	A	221	ILE
1	A	237	ILE
1	A	267	GLN
1	A	291	ASN
1	A	314	LEU
1	B	95	LEU
1	B	134	LYS
1	B	165	THR
1	B	170	CYS
1	B	183	THR
1	B	233	VAL
1	B	249	MET
1	B	267	GLN
1	B	289	GLU
1	B	291	ASN
1	B	301	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	124	HIS
1	A	167	GLN
1	A	194	GLN
1	A	214	ASN
1	A	244	GLN
1	A	252	GLN
1	A	253	GLN
1	A	257	GLN
1	A	267	GLN
1	A	277	ASN
1	A	291	ASN
1	B	124	HIS
1	B	194	GLN
1	B	244	GLN
1	B	252	GLN
1	B	253	GLN
1	B	277	ASN
1	B	291	ASN
1	B	306	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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
2	5PG	D	4	2	11,12,13	0.64	0	10,15,17	0.57	0
2	5PG	C	4	2	11,12,13	0.67	0	10,15,17	0.77	0
2	DSE	C	1	3,2	5,6,7	0.48	0	5,6,8	1.76	1 (20%)
2	DSE	D	1	3,2	5,6,7	0.47	0	5,6,8	1.80	1 (20%)

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	5PG	D	4	2	-	3/5/8/10	0/1/1/1
2	5PG	C	4	2	-	4/5/8/10	0/1/1/1
2	DSE	C	1	3,2	-	1/3/6/8	-
2	DSE	D	1	3,2	-	0/3/6/8	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	DSE	CN-N-CA	3.49	124.13	113.70
2	D	1	DSE	CN-N-CA	3.46	124.05	113.70

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	DSE	O-C-CA-CB
2	C	4	5PG	N-CA-CB-CC1
2	C	4	5PG	N-CA-CB-CC2
2	C	4	5PG	C-CA-CB-CC1
2	C	4	5PG	C-CA-CB-CC2
2	D	4	5PG	C-CA-CB-CC1
2	D	4	5PG	C-CA-CB-CC2
2	D	4	5PG	N-CA-CB-CC1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	5PG	2	0

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	$\# Z > 2$
3	M12	C	0	2	1,2,13	0.85	0	0,1,14	-	-
3	M12	D	0	2	1,2,13	0.82	0	0,1,14	-	-

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.

5.7 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	208/250 (83%)	0.31	9 (4%) 40 36	28, 46, 79, 85	0
1	B	223/250 (89%)	0.40	17 (7%) 20 17	23, 47, 89, 112	0
2	C	3/6 (50%)	1.05	1 (33%) 1 1	40, 40, 43, 59	0
2	D	3/6 (50%)	1.51	1 (33%) 1 1	41, 41, 43, 61	0
All	All	437/512 (85%)	0.36	28 (6%) 25 22	23, 46, 82, 112	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	199	ARG	5.7
1	B	121	GLU	3.7
2	D	3	GLY	3.6
1	A	198	ARG	3.5
1	B	174	GLN	3.2
2	C	3	GLY	3.2
1	B	107	ALA	3.2
1	B	171	SER	3.1
1	B	175	ALA	3.0
1	A	217	LYS	3.0
1	A	323	HIS	2.9
1	B	323	HIS	2.8
1	B	122	THR	2.8
1	B	310	TRP	2.7
1	B	258	LEU	2.6
1	A	216	THR	2.6
1	B	179	ALA	2.6
1	B	312	THR	2.5
1	B	172	SER	2.5
1	B	308	GLY	2.4
1	B	307	GLU	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	173	GLY	2.2
1	A	252	GLN	2.2
1	A	218	GLU	2.2
1	B	257	GLN	2.1
1	A	314	LEU	2.1
1	B	176	CYS	2.0
1	A	204	ALA	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
2	DSE	C	1	7/8	0.63	0.26	71,73,74,76	5
2	DAL	C	2	5/6	0.69	0.24	64,66,67,69	5
2	DSE	D	1	7/8	0.70	0.23	73,76,78,78	5
2	DAL	D	2	5/6	0.85	0.20	65,66,67,70	5
2	5PG	D	4	12/13	0.93	0.10	42,44,49,52	0
2	5PG	C	4	12/13	0.96	0.08	41,42,51,51	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
3	M12	D	0	3/14	0.50	0.31	77,77,77,77	3
3	M12	C	0	3/14	0.54	0.22	73,73,73,73	3

6.5 Other polymers [i](#)

There are no such residues in this entry.