Full wwPDB X-ray Structure Validation Report (i)

Oct 17, 2023 – 09:29 PM EDT

PDB ID	:	2DP4
Title	:	Crystal structure of the complex formed between proteinase K and a human lactoferrin fragment at 2.9 A resolution
Authors	:	Singh, A.K.; Singh, N.; Sharma, S.; Bhushan, A.; Singh, T.P.
Deposited on	:	2006-05-05
Resolution	:	2.90 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 (i) 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 (i)) were used in the production of this report:

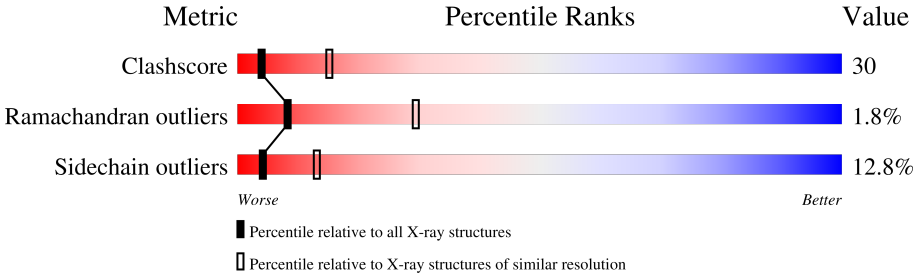
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	279	59% 35% 5%
2	I	8	12% 38% 50%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2201 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 Proteinase K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	279	2029	1247	357	415	10	0	0	0

- Molecule 2 is a protein called 8-mer peptide from Lactotransferrin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	I	8	60	33	11	16	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	E	107	107	107	0	0
3	I	5	5	5	0	0

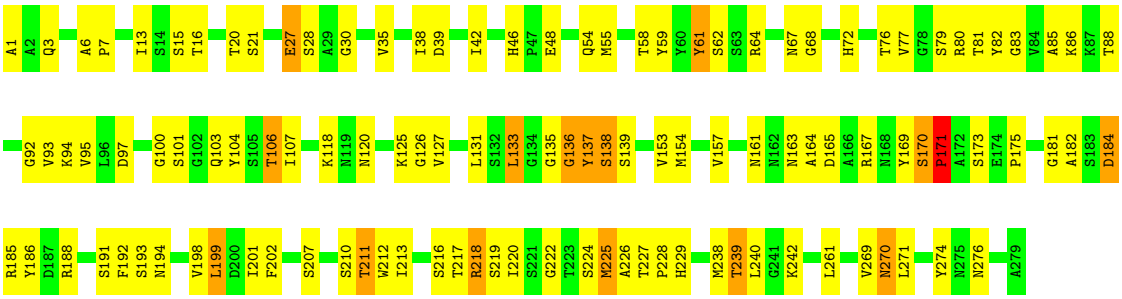
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: Proteinase K

Chain E: 59% 35% 5%



• Molecule 2: 8-mer peptide from Lactotransferrin

Chain I: 12% 38% 50%

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	68.40Å 68.40Å 108.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	11.80 – 2.90	Depositor
% Data completeness (in resolution range)	92.8 (11.80-2.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.193 , 0.235	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2201	wwPDB-VP
Average B, all atoms (Å^2)	14.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

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	E	0.55	0/2068	0.99	9/2810 (0.3%)
2	I	1.15	0/59	2.72	5/76 (6.6%)
All	All	0.58	0/2127	1.07	14/2886 (0.5%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	170	SER	C-N-CD	-23.74	68.37	120.60
1	E	170	SER	C-N-CA	11.37	169.76	122.00
1	E	137	TYR	CB-CG-CD2	8.98	126.39	121.00
1	E	136	GLY	C-N-CA	7.56	140.61	121.70
1	E	137	TYR	CB-CG-CD1	-7.34	116.59	121.00
1	E	211	THR	CA-CB-CG2	-7.17	102.36	112.40
1	E	171	PRO	CA-N-CD	-6.66	102.17	111.50
2	I	2	ASP	C-N-CA	-6.21	106.17	121.70
2	I	4	GLN	N-CA-C	6.01	127.23	111.00
1	E	137	TYR	N-CA-CB	5.81	121.05	110.60
2	I	3	GLU	N-CA-C	5.68	126.34	111.00
1	E	81	THR	N-CA-CB	-5.46	99.92	110.30
2	I	3	GLU	CA-C-N	-5.16	105.86	117.20
2	I	5	GLY	N-CA-C	-5.05	100.47	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	E	2029	0	1935	111	0
2	I	60	0	51	28	0
3	E	107	0	0	8	0
3	I	5	0	0	3	0
All	All	2201	0	1986	123	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 30.

All (123) 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)
2:I:3:GLU:HG3	2:I:4:GLN:H	1.23	1.00
1:E:30:GLY:HA2	1:E:239:THR:HG21	1.52	0.89
1:E:224:SER:OG	2:I:2:ASP:HB3	1.74	0.88
1:E:58:THR:HG21	1:E:62:SER:O	1.77	0.84
1:E:212:TRP:CD2	1:E:218:ARG:HD3	2.12	0.84
1:E:261:LEU:H	1:E:270:ASN:HD21	1.27	0.83
1:E:185:ARG:HD2	3:E:354:HOH:O	1.86	0.75
1:E:218:ARG:NH2	2:I:6:GLU:HB2	2.03	0.74
1:E:46:HIS:HD2	1:E:48:GLU:H	1.34	0.73
1:E:35:VAL:HG11	1:E:77:VAL:HG11	1.71	0.72
2:I:3:GLU:CG	2:I:4:GLN:H	1.99	0.69
1:E:64:ARG:HD3	3:E:358:HOH:O	1.92	0.69
1:E:220:ILE:HB	2:I:4:GLN:O	1.92	0.68
1:E:207:SER:HB3	3:E:306:HOH:O	1.95	0.67
1:E:97:ASP:HB3	1:E:101:SER:O	1.95	0.66
2:I:3:GLU:HG3	2:I:4:GLN:N	2.05	0.66
2:I:4:GLN:O	2:I:4:GLN:CG	2.44	0.65
1:E:64:ARG:HH11	1:E:64:ARG:HG3	1.61	0.65
1:E:30:GLY:HA2	1:E:239:THR:CG2	2.27	0.65
1:E:100:GLY:HA3	2:I:8:LYS:C	2.17	0.63
1:E:42:ILE:HD12	1:E:92:GLY:HA3	1.81	0.63
1:E:133:LEU:C	1:E:133:LEU:HD22	2.21	0.61
1:E:68:GLY:HA2	1:E:213:ILE:HG23	1.83	0.61
1:E:185:ARG:CD	3:E:354:HOH:O	2.45	0.60
1:E:173:SER:HA	1:E:198:VAL:HG21	1.83	0.60
1:E:125:LYS:CG	1:E:239:THR:HA	2.32	0.60
1:E:135:GLY:O	1:E:169:TYR:HD1	1.84	0.60
1:E:38:ILE:HB	1:E:131:LEU:HD23	1.82	0.60
1:E:163:ASN:ND2	1:E:191:SER:O	2.36	0.59
1:E:38:ILE:HG22	1:E:95:VAL:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:269:VAL:HB	1:E:271:LEU:HD12	1.86	0.58
1:E:220:ILE:CG2	2:I:5:GLY:HA2	2.33	0.58
1:E:30:GLY:CA	1:E:239:THR:HG21	2.31	0.57
1:E:72:HIS:NE2	1:E:210:SER:HB3	2.20	0.57
2:I:6:GLU:O	2:I:7:ASN:HB2	2.03	0.57
1:E:58:THR:HG23	1:E:94:LYS:HB3	1.87	0.57
2:I:3:GLU:CG	2:I:4:GLN:N	2.66	0.56
1:E:165:ASP:OD1	1:E:167:ARG:HD2	2.05	0.56
1:E:186:TYR:HD2	1:E:188:ARG:NH2	2.05	0.55
1:E:46:HIS:CD2	1:E:48:GLU:H	2.20	0.55
1:E:225:MET:SD	1:E:225:MET:N	2.80	0.54
1:E:192:PHE:CE1	1:E:222:GLY:HA2	2.43	0.54
1:E:72:HIS:CD2	1:E:210:SER:HB3	2.43	0.54
1:E:61:TYR:CD1	1:E:61:TYR:N	2.67	0.54
1:E:79:SER:H	1:E:83:GLY:HA3	1.72	0.54
1:E:201:ILE:HG13	1:E:202:PHE:N	2.22	0.53
1:E:82:TYR:CD2	1:E:211:THR:CG2	2.92	0.53
1:E:131:LEU:HD22	1:E:133:LEU:HD11	1.91	0.53
1:E:227:THR:N	1:E:228:PRO:HD2	2.25	0.52
1:E:42:ILE:HD12	1:E:92:GLY:CA	2.40	0.52
2:I:3:GLU:CD	3:I:40:HOH:O	2.47	0.52
1:E:184:ASP:C	1:E:184:ASP:OD1	2.48	0.52
2:I:3:GLU:HG2	3:I:40:HOH:O	2.09	0.52
1:E:93:VAL:HG12	1:E:95:VAL:HG13	1.92	0.52
1:E:137:TYR:O	1:E:138:SER:HB2	2.11	0.51
1:E:72:HIS:CE1	1:E:210:SER:HB3	2.46	0.51
1:E:3:GLN:HE22	1:E:86:LYS:NZ	2.09	0.51
1:E:76:THR:O	1:E:83:GLY:HA2	2.11	0.50
1:E:165:ASP:CG	1:E:167:ARG:HD2	2.32	0.50
1:E:125:LYS:HG2	1:E:239:THR:HA	1.94	0.50
1:E:125:LYS:HG2	1:E:239:THR:HB	1.92	0.50
1:E:126:GLY:HA3	1:E:238:MET:HE3	1.93	0.50
1:E:58:THR:HG22	1:E:59:TYR:H	1.77	0.49
1:E:136:GLY:HA2	1:E:169:TYR:HA	1.95	0.49
1:E:220:ILE:HG22	2:I:5:GLY:HA2	1.95	0.49
1:E:64:ARG:HG3	1:E:64:ARG:NH1	2.26	0.49
1:E:133:LEU:HA	2:I:1:GLY:N	2.28	0.49
1:E:137:TYR:CG	1:E:138:SER:N	2.81	0.48
1:E:212:TRP:CG	1:E:218:ARG:HD3	2.47	0.48
2:I:1:GLY:O	2:I:8:LYS:HB2	2.13	0.48
1:E:103:GLN:HB2	1:E:106:THR:HG23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:226:ALA:O	1:E:229:HIS:HB2	2.14	0.48
1:E:126:GLY:HA3	1:E:238:MET:CE	2.44	0.48
1:E:131:LEU:HB2	1:E:157:VAL:HG12	1.95	0.48
1:E:169:TYR:HB3	3:E:298:HOH:O	2.14	0.47
1:E:182:ALA:HB3	1:E:193:SER:HB2	1.95	0.47
1:E:85:ALA:O	1:E:88:THR:HG22	2.15	0.47
1:E:224:SER:OG	2:I:2:ASP:CB	2.57	0.46
1:E:58:THR:HG23	1:E:94:LYS:HD3	1.96	0.46
1:E:133:LEU:C	1:E:133:LEU:CD2	2.85	0.46
1:E:213:ILE:C	1:E:213:ILE:HD12	2.35	0.45
1:E:220:ILE:HG21	2:I:5:GLY:HA2	1.98	0.45
1:E:100:GLY:CA	2:I:8:LYS:C	2.83	0.45
1:E:164:ALA:H	1:E:194:ASN:ND2	2.14	0.45
1:E:220:ILE:HG21	2:I:5:GLY:CA	2.47	0.45
1:E:1:ALA:HB2	1:E:27:GLU:OE1	2.16	0.44
1:E:104:TYR:OH	1:E:135:GLY:HA3	2.16	0.44
1:E:100:GLY:HA3	2:I:8:LYS:HG2	2.00	0.44
1:E:212:TRP:N	1:E:216:SER:O	2.36	0.44
1:E:133:LEU:HD22	1:E:133:LEU:O	2.17	0.44
2:I:4:GLN:O	2:I:4:GLN:HG3	2.15	0.44
1:E:15:SER:HB2	3:E:366:HOH:O	2.17	0.44
1:E:222:GLY:C	1:E:224:SER:H	2.20	0.43
1:E:38:ILE:HG22	1:E:95:VAL:CG2	2.46	0.43
1:E:82:TYR:CD2	1:E:211:THR:HG23	2.53	0.43
1:E:82:TYR:HD2	1:E:211:THR:CG2	2.31	0.43
1:E:202:PHE:HB2	1:E:271:LEU:O	2.19	0.43
1:E:240:LEU:HD12	1:E:240:LEU:HA	1.80	0.43
1:E:270:ASN:ND2	3:E:317:HOH:O	2.52	0.43
1:E:217:THR:O	1:E:218:ARG:HB3	2.18	0.43
1:E:54:GLN:HG2	1:E:55:MET:O	2.19	0.43
1:E:6:ALA:HB1	1:E:7:PRO:CD	2.49	0.42
1:E:173:SER:O	1:E:175:PRO:HD3	2.19	0.42
1:E:13:ILE:O	1:E:274:TYR:HD1	2.03	0.42
1:E:3:GLN:HE22	1:E:86:LYS:HZ1	1.66	0.42
2:I:3:GLU:CG	3:I:40:HOH:O	2.67	0.42
1:E:201:ILE:HG13	1:E:202:PHE:H	1.83	0.42
1:E:127:VAL:HG12	1:E:153:VAL:HG13	2.01	0.42
1:E:100:GLY:HA2	2:I:8:LYS:O	2.19	0.42
1:E:222:GLY:C	1:E:224:SER:N	2.74	0.42
2:I:7:ASN:CG	2:I:8:LYS:H	2.24	0.41
1:E:3:GLN:NE2	1:E:86:LYS:NZ	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:LYS:HG2	1:E:239:THR:CB	2.51	0.41
1:E:210:SER:O	1:E:217:THR:HB	2.21	0.41
1:E:186:TYR:CD2	1:E:188:ARG:NH2	2.86	0.41
1:E:103:GLN:O	1:E:107:ILE:HG12	2.21	0.41
1:E:67:ASN:ND2	2:I:8:LYS:O	2.54	0.41
1:E:80:ARG:NH2	3:E:367:HOH:O	2.41	0.41
1:E:20:THR:HG22	1:E:21:SER:N	2.36	0.40
1:E:181:GLY:HA3	1:E:199:LEU:HD21	2.03	0.40
1:E:218:ARG:HH22	2:I:6:GLU:HB2	1.82	0.40
1:E:161:ASN:HA	1:E:192:PHE:O	2.21	0.40
1:E:161:ASN:ND2	2:I:2:ASP:OD1	2.54	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	E	277/279 (99%)	256 (92%)	19 (7%)	2 (1%)	22 54
2	I	6/8 (75%)	2 (33%)	1 (17%)	3 (50%)	0 0
All	All	283/287 (99%)	258 (91%)	20 (7%)	5 (2%)	8 29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	138	SER
1	E	171	PRO
2	I	2	ASP
2	I	4	GLN
2	I	7	ASN

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	E	213/213 (100%)	191 (90%)	22 (10%)	7 22
2	I	6/6 (100%)	0	6 (100%)	0 0
All	All	219/219 (100%)	191 (87%)	28 (13%)	4 13

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	16	THR
1	E	27	GLU
1	E	28	SER
1	E	39	ASP
1	E	61	TYR
1	E	106	THR
1	E	118	LYS
1	E	120	ASN
1	E	133	LEU
1	E	139	SER
1	E	154	MET
1	E	170	SER
1	E	171	PRO
1	E	184	ASP
1	E	199	LEU
1	E	218	ARG
1	E	219	SER
1	E	225	MET
1	E	239	THR
1	E	242	LYS
1	E	270	ASN
1	E	276	ASN
2	I	2	ASP
2	I	3	GLU
2	I	4	GLN
2	I	6	GLU
2	I	7	ASN

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Mol	Chain	Res	Type
2	I	8	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	3	GLN
1	E	5	ASN
1	E	31	GLN
1	E	46	HIS
1	E	54	GLN
1	E	67	ASN
1	E	119	ASN
1	E	120	ASN
1	E	162	ASN
1	E	168	ASN
1	E	194	ASN
1	E	270	ASN
2	I	4	GLN

5.3.3 RNA (i)

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.