Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 12:08 AM EDT

PDB ID	:	1JRK
Title	:	Crystal Structure of a Nudix Protein from Pyrobaculum aerophilum Reveals a Dimer with Intertwined Beta Sheets
Authors	:	Wang, S.; Mura, C.; Sawaya, M.R.; Cascio, D.; Eisenberg, D.
Deposited on	:	2001-08-13
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 (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:

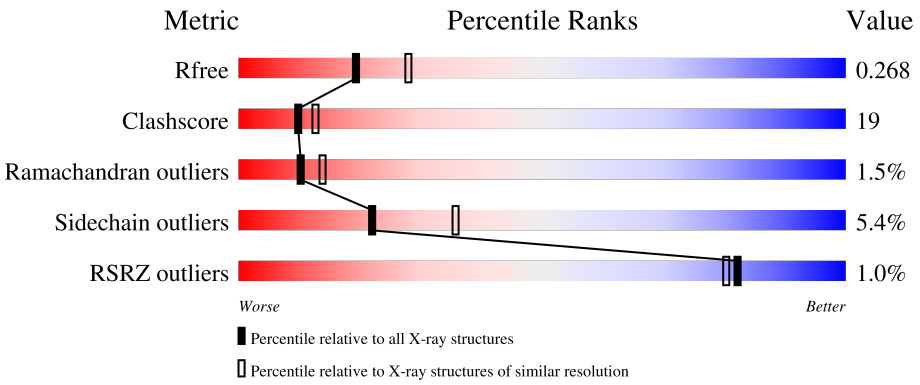
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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 <=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	156	67% 26% • • 5%
1	B	156	% 58% 33% • 6%
1	C	156	3% 62% 29% • 5%
1	D	156	% 69% 24% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MPD	C	205	-	-	X	-
2	MPD	C	206	-	-	X	-

2 Entry composition (i)

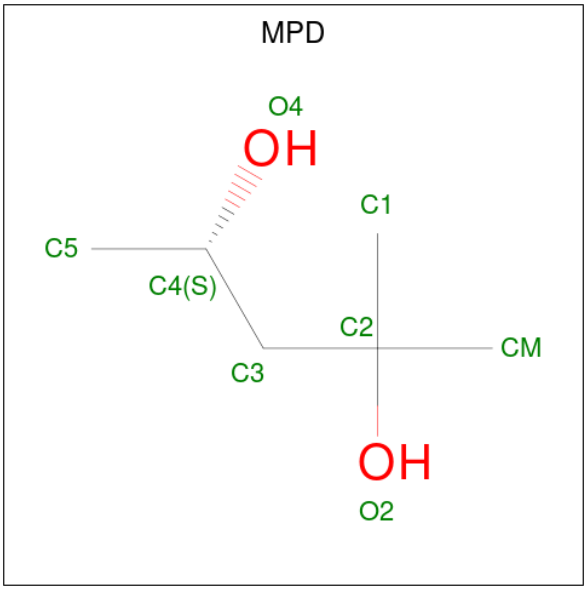
There are 3 unique types of molecules in this entry. The entry contains 4949 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nudix homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	148	1180	764	198	216	2	0	0	0
1	B	147	1175	761	197	215	2	0	0	0
1	C	148	1180	764	198	216	2	0	0	0
1	D	150	1193	773	200	218	2	0	0	0

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	8	6	2	0	0
2	B	1	8	6	2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	B	1	8	6	2	0	0
			Total	C	O		
2	C	1	8	6	2	0	0
			Total	C	O		
2	C	1	8	6	2	0	0
			Total	C	O		
2	C	1	8	6	2	0	0

- Molecule 3 is water.

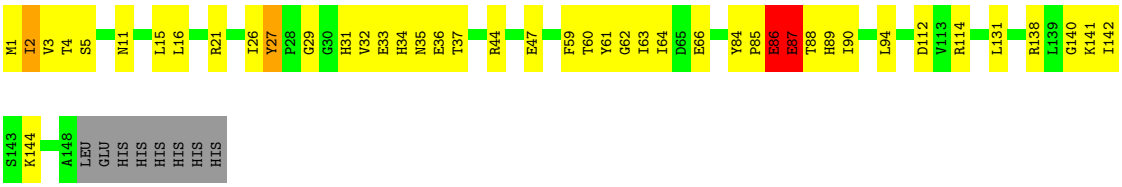
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	42	42	42	0	0
			Total	O		
3	B	39	39	39	0	0
			Total	O		
3	C	54	54	54	0	0
			Total	O		
3	D	38	38	38	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

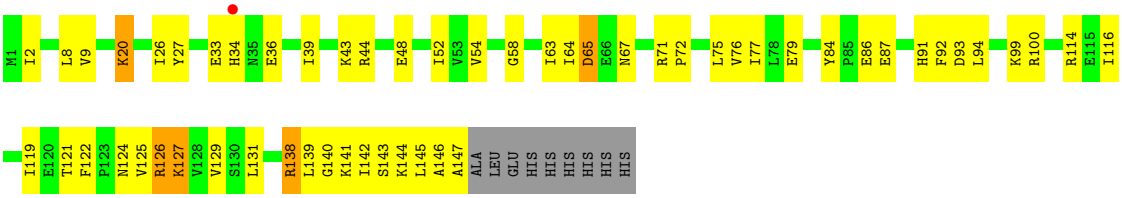
- Molecule 1: Nudix homolog

Chain A:



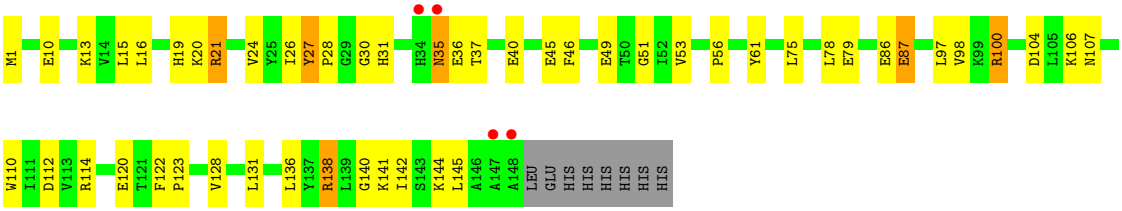
- Molecule 1: Nudix homolog

Chain B:



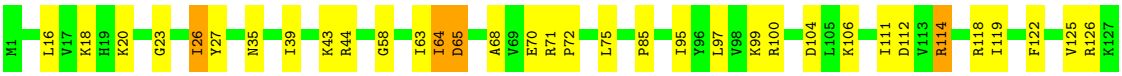
- Molecule 1: Nudix homolog

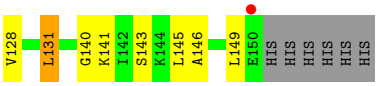
Chain C:



- Molecule 1: Nudix homolog

Chain D:




V128
L131
G140
K141
I142
S143
K144
L145
A146
L149
E150
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	52.47Å 81.46Å 73.89Å 90.00° 99.96° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 72.77 – 2.39	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-2.40) 94.7 (72.77-2.39)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 2.40Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.190 , 0.275 0.185 , 0.268	Depositor DCC
R_{free} test set	1103 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å^2)	29.2	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 60.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4949	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.68 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0622e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MPD

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	0.64	0/1203	0.86	1/1631 (0.1%)
1	B	0.61	0/1198	0.82	0/1624
1	C	0.59	0/1203	0.81	0/1631
1	D	0.60	0/1216	0.79	0/1649
All	All	0.61	0/4820	0.82	1/6535 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	87	GLU	N-CA-C	-5.23	96.89	111.00

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	A	1180	0	1219	44	0
1	B	1175	0	1214	60	0
1	C	1180	0	1219	58	0
1	D	1193	0	1232	39	0
2	A	8	0	14	4	0
2	B	16	0	28	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	24	0	42	13	0
3	A	42	0	0	1	0
3	B	39	0	0	3	0
3	C	54	0	0	1	0
3	D	38	0	0	1	0
All	All	4949	0	4968	187	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 19.

All (187) 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:B:146:ALA:HB3	3:B:241:HOH:O	1.61	0.98
1:A:141:LYS:HD3	1:A:144:LYS:HE2	1.45	0.95
1:A:59:PHE:HB3	2:A:203:MPD:H31	1.48	0.94
1:A:21:ARG:NH2	1:A:86:GLU:OE2	2.04	0.90
1:B:20:LYS:H	1:B:20:LYS:CE	1.86	0.88
1:B:99:LYS:HG3	2:B:202:MPD:H4	1.56	0.86
1:C:141:LYS:HE3	1:C:144:LYS:NZ	1.95	0.82
1:B:144:LYS:HG2	3:B:237:HOH:O	1.80	0.80
1:A:2:ILE:C	1:A:2:ILE:HD13	2.04	0.78
1:D:27:TYR:OH	1:D:128:VAL:HB	1.84	0.77
1:A:87:GLU:HG3	1:A:89:HIS:HE1	1.49	0.77
1:B:2:ILE:HD12	1:B:2:ILE:H	1.49	0.76
1:C:20:LYS:H	2:C:205:MPD:H52	1.50	0.75
1:C:138:ARG:HH11	1:C:138:ARG:HB2	1.51	0.75
1:C:13:LYS:HB2	2:C:206:MPD:C5	2.17	0.74
1:D:141:LYS:HD3	1:D:141:LYS:C	2.08	0.74
1:B:138:ARG:HA	1:B:138:ARG:HE	1.52	0.73
1:A:44:ARG:O	1:A:47:GLU:HB3	1.88	0.73
1:C:45:GLU:O	1:C:49:GLU:HG3	1.89	0.73
3:A:238:HOH:O	1:C:138:ARG:HG3	1.89	0.72
1:B:58:GLY:O	1:D:145:LEU:HD21	1.90	0.71
1:B:20:LYS:H	1:B:20:LYS:HE2	1.54	0.71
1:C:141:LYS:HA	1:C:144:LYS:HB3	1.73	0.70
1:B:20:LYS:H	1:B:20:LYS:HE3	1.55	0.70
1:A:87:GLU:HG3	1:A:89:HIS:CE1	2.26	0.69
1:B:126:ARG:HD2	1:B:127:LYS:N	2.07	0.69
1:C:51:GLY:HA3	1:C:106:LYS:HG2	1.74	0.69
1:B:2:ILE:HD12	1:B:2:ILE:N	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:ARG:HD2	1:B:127:LYS:H	1.58	0.69
1:C:20:LYS:HG3	2:C:205:MPD:H53	1.75	0.68
1:C:13:LYS:HG2	1:C:112:ASP:HA	1.74	0.68
1:D:26:ILE:HG22	1:D:122:PHE:HB2	1.74	0.68
1:B:27:TYR:OH	1:B:129:VAL:HG23	1.94	0.67
1:C:13:LYS:HB2	2:C:206:MPD:H51	1.78	0.66
1:B:140:GLY:O	1:B:144:LYS:HG3	1.94	0.66
1:A:85:PRO:O	1:A:86:GLU:OE1	2.14	0.66
1:C:86:GLU:HG2	1:C:87:GLU:HG2	1.78	0.66
1:A:84:TYR:O	1:A:86:GLU:N	2.28	0.66
1:C:13:LYS:HD2	2:C:206:MPD:H53	1.78	0.65
1:A:60:THR:O	2:A:203:MPD:H53	1.97	0.64
1:B:141:LYS:O	1:B:144:LYS:HB2	1.96	0.64
1:C:138:ARG:HB2	1:C:138:ARG:NH1	2.13	0.64
1:C:141:LYS:HE3	1:C:144:LYS:HZ2	1.61	0.63
1:C:141:LYS:HE3	1:C:144:LYS:HZ1	1.63	0.63
1:D:26:ILE:HG22	1:D:122:PHE:CB	2.29	0.62
1:A:131:LEU:HD11	1:B:63:ILE:HD11	1.82	0.62
1:C:110:TRP:CE3	2:C:206:MPD:H52	2.34	0.62
1:C:19:HIS:HA	2:C:205:MPD:H52	1.80	0.62
1:A:86:GLU:HG3	1:A:87:GLU:OE2	2.00	0.61
1:C:79:GLU:CD	1:D:71:ARG:HH21	2.03	0.61
1:B:52:ILE:HD12	1:B:100:ARG:NH2	2.14	0.61
1:D:26:ILE:HD12	1:D:27:TYR:O	2.01	0.60
1:A:84:TYR:HB3	1:A:86:GLU:CG	2.31	0.60
1:D:26:ILE:CG2	1:D:122:PHE:HB2	2.30	0.60
1:D:104:ASP:O	1:D:106:LYS:HG3	2.00	0.60
1:A:2:ILE:HD11	1:A:32:VAL:HG21	1.82	0.59
1:A:21:ARG:HH22	1:A:86:GLU:HG3	1.67	0.59
1:B:8:LEU:HD12	2:B:202:MPD:HM1	1.84	0.59
1:C:13:LYS:HD2	2:C:206:MPD:C5	2.31	0.59
1:D:118:ARG:O	1:D:118:ARG:HG2	2.02	0.59
1:A:2:ILE:HD13	1:A:2:ILE:O	2.03	0.59
1:C:27:TYR:OH	1:C:128:VAL:HG21	2.03	0.58
1:C:20:LYS:H	2:C:205:MPD:C5	2.16	0.58
1:D:18:LYS:HE3	1:D:23:GLY:O	2.03	0.58
1:A:131:LEU:CD1	1:B:63:ILE:HD11	2.34	0.58
1:A:63:ILE:HD13	1:B:127:LYS:HD2	1.84	0.57
1:A:3:VAL:HG12	1:A:31:HIS:HA	1.86	0.57
1:C:138:ARG:HH11	1:C:138:ARG:CB	2.18	0.57
1:B:44:ARG:O	1:B:48:GLU:HG3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLY:O	1:C:141:LYS:HE2	2.05	0.57
1:B:64:ILE:HD12	1:B:65:ASP:H	1.69	0.57
1:A:112:ASP:OD2	1:A:114:ARG:NH2	2.38	0.56
1:C:75:LEU:HD22	1:C:97:LEU:HD13	1.86	0.56
1:D:85:PRO:HG2	3:D:173:HOH:O	2.05	0.56
1:D:112:ASP:OD1	1:D:114:ARG:HG2	2.06	0.56
1:B:139:LEU:CD1	1:D:149:LEU:HD13	2.37	0.55
1:C:24:VAL:HG12	1:C:120:GLU:HG2	1.88	0.55
1:C:61:TYR:CD1	1:D:131:LEU:HD11	2.42	0.55
1:D:27:TYR:OH	1:D:125:VAL:HA	2.07	0.55
1:C:30:GLY:HA3	1:C:45:GLU:CD	2.28	0.54
1:B:20:LYS:HE3	1:B:20:LYS:N	2.23	0.53
1:A:26:ILE:HG12	1:A:27:TYR:N	2.24	0.53
1:C:86:GLU:O	1:C:87:GLU:HB3	2.09	0.53
1:B:143:SER:HB2	1:D:146:ALA:HB1	1.90	0.52
1:D:111:ILE:HD12	1:D:119:ILE:HD13	1.91	0.52
1:C:10:GLU:OE2	1:C:100:ARG:NH1	2.42	0.52
1:C:26:ILE:HG12	1:C:27:TYR:N	2.24	0.52
1:B:114:ARG:HH11	1:B:114:ARG:HG3	1.75	0.52
1:C:21:ARG:NH2	1:C:86:GLU:OE1	2.42	0.52
1:D:64:ILE:HD12	1:D:65:ASP:H	1.75	0.52
1:B:76:VAL:HG22	1:B:77:ILE:N	2.24	0.52
1:B:52:ILE:HD12	1:B:100:ARG:CZ	2.41	0.51
1:A:84:TYR:O	1:A:87:GLU:O	2.28	0.51
1:C:1:MET:CG	1:C:31:HIS:CD2	2.94	0.50
1:B:141:LYS:O	1:B:144:LYS:CB	2.59	0.50
1:D:114:ARG:HG2	1:D:114:ARG:HH11	1.77	0.50
1:A:2:ILE:CD1	1:A:32:VAL:HG21	2.42	0.49
1:A:140:GLY:O	1:A:144:LYS:HG2	2.13	0.49
1:B:75:LEU:HA	2:B:204:MPD:H12	1.93	0.49
1:B:124:ASN:OD1	1:B:125:VAL:N	2.45	0.49
1:D:63:ILE:HG22	1:D:64:ILE:N	2.28	0.49
1:A:85:PRO:O	1:A:86:GLU:CB	2.60	0.49
1:D:16:LEU:HD12	1:D:111:ILE:HG13	1.93	0.49
1:D:63:ILE:CG2	1:D:64:ILE:N	2.76	0.49
1:B:2:ILE:H	1:B:2:ILE:CD1	2.20	0.49
1:B:84:TYR:HB2	1:B:87:GLU:O	2.13	0.49
1:B:39:ILE:O	1:B:43:LYS:HG3	2.13	0.48
1:B:146:ALA:O	1:B:147:ALA:C	2.51	0.48
1:A:4:THR:O	1:A:29:GLY:HA3	2.13	0.48
1:A:61:TYR:CE2	2:A:203:MPD:HM2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:LYS:HG3	2:B:202:MPD:C4	2.35	0.48
1:C:30:GLY:HA3	1:C:45:GLU:HG3	1.94	0.48
1:C:20:LYS:N	2:C:205:MPD:H52	2.26	0.48
1:C:21:ARG:HH22	1:C:86:GLU:CD	2.17	0.48
1:C:110:TRP:CD2	2:C:206:MPD:H31	2.49	0.47
1:C:86:GLU:O	1:C:87:GLU:CB	2.62	0.47
1:A:84:TYR:HB2	1:A:87:GLU:HG2	1.97	0.47
1:B:52:ILE:CD1	1:B:100:ARG:NH2	2.77	0.47
1:C:141:LYS:HA	1:C:144:LYS:CB	2.42	0.47
1:B:52:ILE:CG2	1:B:100:ARG:HG3	2.45	0.47
1:C:122:PHE:CD1	1:C:123:PRO:HD2	2.49	0.47
1:D:141:LYS:HD3	1:D:141:LYS:O	2.14	0.47
1:B:26:ILE:HG22	1:B:122:PHE:HB2	1.96	0.46
1:D:39:ILE:O	1:D:43:LYS:HG3	2.15	0.46
1:C:1:MET:SD	1:C:1:MET:N	2.79	0.46
1:A:84:TYR:C	1:A:86:GLU:HG2	2.36	0.46
1:C:53:VAL:HA	3:C:252:HOH:O	2.15	0.46
1:C:37:THR:HG23	1:C:40:GLU:OE1	2.15	0.46
1:A:84:TYR:CB	1:A:86:GLU:HG2	2.46	0.46
1:C:30:GLY:HA3	1:C:45:GLU:CG	2.46	0.46
1:B:86:GLU:O	1:B:87:GLU:HB3	2.16	0.46
1:A:2:ILE:CD1	1:A:32:VAL:CG2	2.94	0.46
1:B:143:SER:HB2	1:D:146:ALA:CB	2.46	0.45
1:A:84:TYR:HB3	1:A:86:GLU:HG2	1.98	0.45
1:C:56:PRO:HA	1:C:98:VAL:HG12	1.98	0.45
1:A:59:PHE:HB3	2:A:203:MPD:C3	2.35	0.45
1:B:77:ILE:HA	1:B:93:ASP:O	2.16	0.45
1:D:70:GLU:HG2	1:D:71:ARG:N	2.30	0.45
1:C:15:LEU:O	1:C:16:LEU:HD23	2.16	0.45
1:A:85:PRO:O	1:A:86:GLU:HB3	2.17	0.45
1:C:13:LYS:HB2	2:C:206:MPD:H52	1.97	0.45
1:A:15:LEU:O	1:A:16:LEU:HD23	2.17	0.44
1:C:26:ILE:HG12	1:C:27:TYR:H	1.80	0.44
1:A:138:ARG:O	1:A:142:ILE:HG13	2.18	0.44
1:C:136:LEU:O	1:C:140:GLY:N	2.49	0.44
1:D:44:ARG:HG2	1:D:44:ARG:HH11	1.83	0.44
1:D:140:GLY:O	1:D:143:SER:OG	2.33	0.44
1:A:66:GLU:H	1:A:66:GLU:CD	2.22	0.44
1:C:1:MET:HG2	1:C:31:HIS:CD2	2.53	0.44
1:D:35:ASN:OD1	1:D:35:ASN:N	2.47	0.44
1:A:1:MET:N	1:B:34:HIS:O	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:CG	1:A:90:ILE:O	2.67	0.43
1:B:138:ARG:NH2	1:B:141:LYS:HZ2	2.17	0.43
1:B:144:LYS:O	1:B:146:ALA:N	2.51	0.43
1:C:131:LEU:HA	1:C:131:LEU:HD23	1.68	0.43
1:D:75:LEU:N	1:D:95:ILE:O	2.47	0.43
1:B:54:VAL:HA	1:B:99:LYS:O	2.19	0.43
1:B:139:LEU:HD13	1:D:149:LEU:HD13	2.01	0.43
1:C:1:MET:HG3	1:C:31:HIS:CD2	2.53	0.43
1:D:97:LEU:HD12	1:D:97:LEU:HA	1.87	0.43
1:A:37:THR:HB	1:B:92:PHE:CD1	2.54	0.42
1:D:112:ASP:OD2	1:D:114:ARG:NH1	2.52	0.42
1:C:78:LEU:HD23	1:D:70:GLU:HA	2.00	0.42
1:A:33:GLU:HB2	1:A:36:GLU:OE1	2.20	0.42
1:B:33:GLU:HB2	1:B:36:GLU:HG2	2.02	0.42
1:B:144:LYS:C	1:B:146:ALA:H	2.23	0.42
1:B:67:ASN:HB3	3:B:222:HOH:O	2.20	0.42
1:C:19:HIS:HA	2:C:205:MPD:C5	2.48	0.42
1:C:49:GLU:O	1:C:107:ASN:HB3	2.20	0.42
1:B:116:ILE:CD1	1:B:129:VAL:HG21	2.51	0.41
1:D:99:LYS:HG2	1:D:100:ARG:N	2.34	0.41
1:B:126:ARG:CD	1:B:127:LYS:N	2.81	0.41
1:D:58:GLY:HA3	1:D:72:PRO:HB2	2.03	0.41
1:D:114:ARG:HH11	1:D:114:ARG:CG	2.34	0.41
1:B:9:VAL:O	2:B:202:MPD:HM2	2.20	0.41
1:B:52:ILE:HG23	1:B:100:ARG:HG3	2.02	0.41
1:B:71:ARG:HB3	1:B:72:PRO:HD2	2.03	0.41
1:B:142:ILE:HD12	1:B:142:ILE:N	2.36	0.41
1:C:28:PRO:HB2	1:C:46:PHE:CD1	2.56	0.41
1:B:119:ILE:HG13	1:B:121:THR:HG23	2.03	0.41
1:B:79:GLU:HA	1:B:91:HIS:O	2.21	0.40
1:C:104:ASP:O	1:C:106:LYS:HG2	2.22	0.40
1:A:131:LEU:HD23	1:A:131:LEU:HA	1.87	0.40
1:D:64:ILE:HD12	1:D:68:ALA:O	2.22	0.40
1:B:77:ILE:HD13	1:B:94:LEU:CD2	2.51	0.40
1:A:21:ARG:HH22	1:A:86:GLU:CG	2.32	0.40
1:C:112:ASP:OD1	1:C:114:ARG:HB2	2.21	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	146/156 (94%)	136 (93%)	6 (4%)	4 (3%)	5	5
1	B	145/156 (93%)	135 (93%)	9 (6%)	1 (1%)	22	32
1	C	146/156 (94%)	134 (92%)	9 (6%)	3 (2%)	7	8
1	D	148/156 (95%)	143 (97%)	4 (3%)	1 (1%)	22	32
All	All	585/624 (94%)	548 (94%)	28 (5%)	9 (2%)	10	14

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	86	GLU
1	A	87	GLU
1	C	36	GLU
1	C	35	ASN
1	C	87	GLU
1	A	88	THR
1	B	145	LEU
1	D	64	ILE
1	A	64	ILE

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	130/138 (94%)	121 (93%)	9 (7%)	15	25
1	B	130/138 (94%)	124 (95%)	6 (5%)	27	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	130/138 (94%)	123 (95%)	7 (5%)	22	36
1	D	131/138 (95%)	125 (95%)	6 (5%)	27	43
All	All	521/552 (94%)	493 (95%)	28 (5%)	22	36

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

Mol	Chain	Res	Type
1	A	2	ILE
1	A	5	SER
1	A	11	ASN
1	A	27	TYR
1	A	34	HIS
1	A	35	ASN
1	A	86	GLU
1	A	87	GLU
1	A	94	LEU
1	B	20	LYS
1	B	65	ASP
1	B	126	ARG
1	B	127	LYS
1	B	131	LEU
1	B	138	ARG
1	C	21	ARG
1	C	27	TYR
1	C	35	ASN
1	C	100	ARG
1	C	138	ARG
1	C	142	ILE
1	C	145	LEU
1	D	20	LYS
1	D	26	ILE
1	D	65	ASP
1	D	114	ARG
1	D	126	ARG
1	D	131	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN

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Mol	Chain	Res	Type
1	A	89	HIS
1	B	34	HIS
1	B	67	ASN
1	B	89	HIS
1	B	107	ASN
1	C	107	ASN

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)

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	$\#\|Z\| > 2$
2	MPD	A	203	-	7,7,7	0.63	0	9,10,10	0.47	0
2	MPD	B	202	-	7,7,7	0.70	0	9,10,10	0.51	0
2	MPD	C	205	-	7,7,7	0.49	0	9,10,10	0.48	0
2	MPD	C	206	-	7,7,7	0.46	0	9,10,10	0.59	0
2	MPD	C	201	-	7,7,7	0.45	0	9,10,10	0.79	0
2	MPD	B	204	-	7,7,7	0.38	0	9,10,10	0.70	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
2	MPD	A	203	-	-	0/5/5/5	-
2	MPD	B	202	-	-	0/5/5/5	-
2	MPD	C	205	-	-	1/5/5/5	-
2	MPD	C	206	-	-	3/5/5/5	-
2	MPD	C	201	-	-	3/5/5/5	-
2	MPD	B	204	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	201	MPD	C1-C2-C3-C4
2	C	201	MPD	O2-C2-C3-C4
2	C	201	MPD	CM-C2-C3-C4
2	C	206	MPD	C1-C2-C3-C4
2	B	204	MPD	O2-C2-C3-C4
2	C	206	MPD	O2-C2-C3-C4
2	C	205	MPD	C1-C2-C3-C4
2	C	206	MPD	CM-C2-C3-C4

There are no ring outliers.

5 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	203	MPD	4	0
2	B	202	MPD	4	0
2	C	205	MPD	6	0
2	C	206	MPD	7	0
2	B	204	MPD	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	148/156 (94%)	-0.41	0	100	100	16, 29, 60, 82	0
1	B	147/156 (94%)	-0.49	1 (0%)	87	86	13, 27, 59, 75	0
1	C	148/156 (94%)	-0.41	4 (2%)	54	52	14, 25, 62, 83	0
1	D	150/156 (96%)	-0.52	1 (0%)	87	86	14, 28, 51, 70	0
All	All	593/624 (95%)	-0.46	6 (1%)	82	80	13, 27, 59, 83	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	34	HIS	4.0
1	C	148	ALA	3.1
1	B	34	HIS	2.7
1	C	147	ALA	2.5
1	D	150	GLU	2.1
1	C	35	ASN	2.1

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 monosaccharides 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
2	MPD	A	203	8/8	0.69	0.33	78,79,82,84	0
2	MPD	B	202	8/8	0.83	0.38	62,69,71,73	0
2	MPD	C	201	8/8	0.84	0.21	45,54,63,65	0
2	MPD	C	206	8/8	0.85	0.37	63,69,72,72	0
2	MPD	B	204	8/8	0.88	0.21	38,50,61,64	0
2	MPD	C	205	8/8	0.91	0.26	57,62,68,68	0

6.5 Other polymers (i)

There are no such residues in this entry.