



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:46 am GMT

PDB ID : 2B4H
Title : Crystal Structure of the Rhesus Rotavirus VP5 Antigen Domain Dimer
Authors : Yoder, J.D.; Dormitzer, P.R.
Deposited on : 2005-09-24
Resolution : 1.60 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

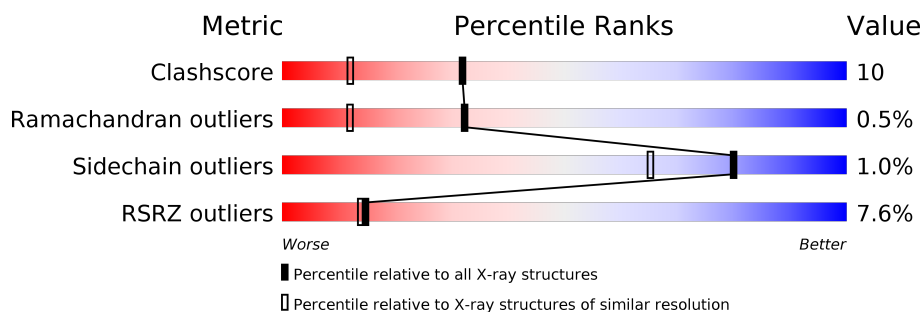
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (1.60-1.60)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	254	
1	B	254	

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	TRS	A	5001	-	-	-	X
3	MPD	A	2001	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MPD	A	3001	-	-	-	X
3	MPD	B	1001	-	-	-	X
3	MPD	B	4001	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3993 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 Outer capsid protein VP4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	0	0
			1801	1148	302	345	6			
1	B	218	Total	C	N	O	S	0	0	0
			1734	1107	291	330	6			

There are 42 discrepancies between the modelled and reference sequences:

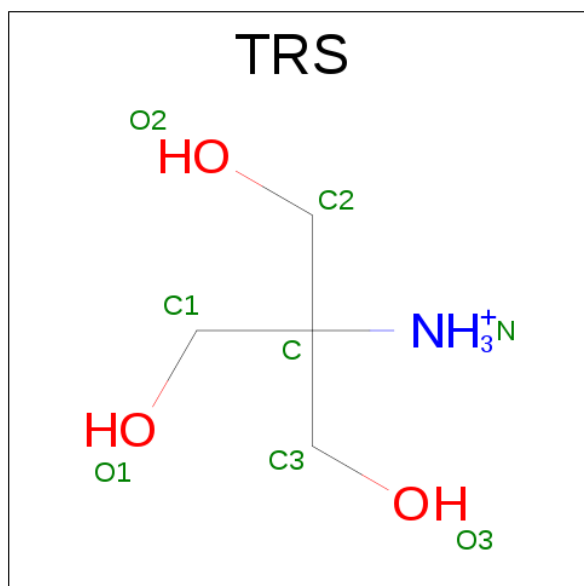
Chain	Residue	Modelled	Actual	Comment	Reference
A	226	MET	-	EXPRESSION TAG	UNP Q91HI9
A	227	GLY	-	EXPRESSION TAG	UNP Q91HI9
A	228	SER	-	EXPRESSION TAG	UNP Q91HI9
A	229	SER	-	EXPRESSION TAG	UNP Q91HI9
A	230	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	231	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	232	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	233	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	234	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	235	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	236	SER	-	EXPRESSION TAG	UNP Q91HI9
A	237	SER	-	EXPRESSION TAG	UNP Q91HI9
A	238	GLY	-	EXPRESSION TAG	UNP Q91HI9
A	239	LEU	-	EXPRESSION TAG	UNP Q91HI9
A	240	VAL	-	EXPRESSION TAG	UNP Q91HI9
A	241	PRO	-	EXPRESSION TAG	UNP Q91HI9
A	242	ARG	-	EXPRESSION TAG	UNP Q91HI9
A	243	GLY	-	EXPRESSION TAG	UNP Q91HI9
A	244	SER	-	EXPRESSION TAG	UNP Q91HI9
A	245	HIS	-	EXPRESSION TAG	UNP Q91HI9
A	246	MET	-	EXPRESSION TAG	UNP Q91HI9
B	226	MET	-	EXPRESSION TAG	UNP Q91HI9
B	227	GLY	-	EXPRESSION TAG	UNP Q91HI9
B	228	SER	-	EXPRESSION TAG	UNP Q91HI9
B	229	SER	-	EXPRESSION TAG	UNP Q91HI9

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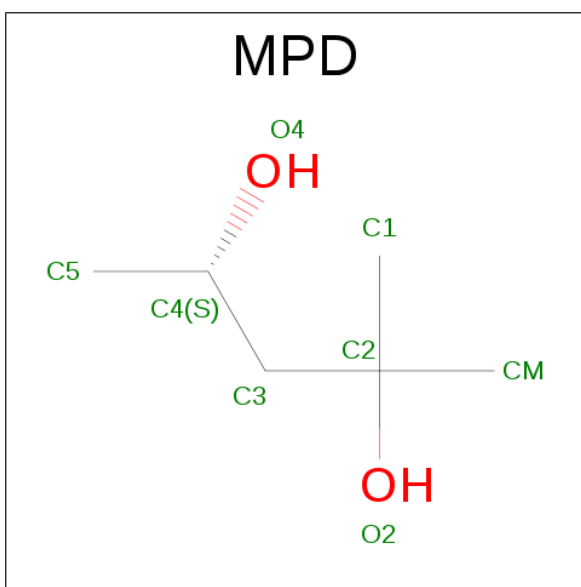
Chain	Residue	Modelled	Actual	Comment	Reference
B	230	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	231	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	232	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	233	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	234	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	235	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	236	SER	-	EXPRESSION TAG	UNP Q91HI9
B	237	SER	-	EXPRESSION TAG	UNP Q91HI9
B	238	GLY	-	EXPRESSION TAG	UNP Q91HI9
B	239	LEU	-	EXPRESSION TAG	UNP Q91HI9
B	240	VAL	-	EXPRESSION TAG	UNP Q91HI9
B	241	PRO	-	EXPRESSION TAG	UNP Q91HI9
B	242	ARG	-	EXPRESSION TAG	UNP Q91HI9
B	243	GLY	-	EXPRESSION TAG	UNP Q91HI9
B	244	SER	-	EXPRESSION TAG	UNP Q91HI9
B	245	HIS	-	EXPRESSION TAG	UNP Q91HI9
B	246	MET	-	EXPRESSION TAG	UNP Q91HI9

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	222	Total	O	0	0
			222	222		
4	B	196	Total	O	0	0
			196	196		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.36Å 54.80Å 65.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.50 – 1.60 45.51 – 1.50	Depositor EDS
% Data completeness (in resolution range)	93.6 (45.50-1.60) 83.6 (45.51-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 1.50Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.203 , 0.225 0.206 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.922	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.6	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	3993	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 26.96 % 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 2.3955e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: TRS, 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	0.59	0/1846	0.77	0/2508
1	B	0.62	0/1779	0.80	0/2417
All	All	0.61	0/3625	0.78	0/4925

There are no bond length outliers.

There are no bond angle outliers.

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	1801	0	1735	40	0
1	B	1734	0	1667	40	0
2	A	8	0	11	0	0
3	A	16	0	28	5	0
3	B	16	0	28	0	0
4	A	222	0	0	2	0
4	B	196	0	0	8	0
All	All	3993	0	3469	71	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (71) 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:419:VAL:HG12	1:A:420:SER:H	1.39	0.88
1:B:419:VAL:HG12	1:B:420:SER:H	1.37	0.87
1:A:337:PHE:H	3:A:3001:MPD:H12	1.47	0.79
1:A:416:THR:HG22	1:A:417:ASP:H	1.52	0.73
1:A:416:THR:HG22	1:A:417:ASP:N	2.05	0.72
1:A:321:ASN:HB2	1:A:352:TYR:HE1	1.55	0.69
1:A:419:VAL:HG12	1:A:420:SER:N	2.08	0.68
1:B:419:VAL:HG12	1:B:420:SER:N	2.09	0.67
1:B:260:SER:HB2	1:B:262:TRP:HE1	1.60	0.67
1:B:416:THR:O	1:B:418:PHE:N	2.29	0.66
1:A:307:ARG:O	1:A:310:GLU:HG2	1.97	0.64
1:A:263:LYS:HD3	1:A:264:GLU:H	1.63	0.62
1:B:381:THR:CG2	4:B:4152:HOH:O	2.48	0.62
1:A:261:LEU:HD23	1:B:261:LEU:HD23	1.81	0.61
1:A:441:ARG:O	3:A:3001:MPD:H51	2.01	0.60
1:A:261:LEU:HB3	1:A:476:SER:HB3	1.85	0.58
1:B:284:SER:HB2	1:B:290:LYS:HB2	1.85	0.57
1:A:315:HIS:ND1	3:A:2001:MPD:H4	2.19	0.57
1:A:416:THR:O	1:A:418:PHE:N	2.38	0.56
1:A:416:THR:CG2	1:A:417:ASP:H	2.17	0.56
1:B:307:ARG:O	1:B:310:GLU:HG2	2.06	0.56
1:A:419:VAL:CG1	1:A:420:SER:H	2.16	0.56
1:A:250:ALA:O	1:B:269:ARG:HA	2.07	0.55
1:A:337:PHE:H	3:A:3001:MPD:C1	2.19	0.54
1:A:250:ALA:O	1:B:269:ARG:CA	2.56	0.54
1:A:269:ARG:HH11	1:A:269:ARG:HG3	1.73	0.54
1:B:366:VAL:HG13	1:B:367:TYR:N	2.24	0.53
1:B:287:LEU:HD22	1:B:289:TYR:CE1	2.44	0.53
1:B:287:LEU:HD23	1:B:287:LEU:O	2.08	0.53
1:A:250:ALA:HB3	1:B:268:ASN:HB3	1.91	0.53
1:B:416:THR:HG22	1:B:417:ASP:N	2.24	0.51
1:A:416:THR:CG2	1:A:417:ASP:N	2.72	0.51
1:A:263:LYS:CD	1:A:264:GLU:H	2.23	0.51
1:A:261:LEU:HD22	1:B:259:THR:CG2	2.41	0.50
1:B:460:GLU:HG2	4:B:4079:HOH:O	2.11	0.49
1:A:467:ARG:HD2	4:A:5087:HOH:O	2.11	0.49
1:A:278:ALA:HB2	1:A:299:ALA:HB2	1.95	0.49
1:B:369:ARG:HH11	1:B:369:ARG:HG2	1.77	0.49
1:A:381:THR:HG21	4:A:5067:HOH:O	2.12	0.49
1:B:371:LEU:HD22	1:B:468:PHE:CZ	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:287:LEU:HD23	4:B:4085:HOH:O	2.13	0.48
1:B:381:THR:HG22	4:B:4152:HOH:O	2.12	0.48
1:B:260:SER:CB	1:B:262:TRP:HE1	2.25	0.48
1:B:366:VAL:HG13	1:B:367:TYR:H	1.77	0.48
1:A:337:PHE:N	3:A:3001:MPD:H12	2.25	0.48
1:B:381:THR:HG23	4:B:4152:HOH:O	2.12	0.47
1:B:287:LEU:HD21	4:B:4052:HOH:O	2.14	0.47
1:B:366:VAL:CG1	4:B:4111:HOH:O	2.63	0.46
1:B:381:THR:O	1:B:457:ASN:HB2	2.15	0.46
1:B:366:VAL:HG12	4:B:4111:HOH:O	2.14	0.46
1:B:287:LEU:HD22	1:B:289:TYR:HE1	1.79	0.46
1:A:470:LEU:HD21	1:A:472:SER:HB3	1.97	0.46
1:B:369:ARG:NH1	1:B:369:ARG:HG2	2.29	0.46
1:A:250:ALA:N	1:B:467:ARG:HG2	2.31	0.45
1:B:419:VAL:CG1	1:B:420:SER:H	2.17	0.45
1:A:419:VAL:CG1	1:A:420:SER:N	2.78	0.45
1:A:369:ARG:HB2	1:B:367:TYR:OH	2.17	0.45
1:A:329:ASN:ND2	1:A:331:GLY:H	2.15	0.45
1:A:256:VAL:O	1:A:256:VAL:HG12	2.17	0.45
1:A:270:ASP:OD1	1:A:467:ARG:HG2	2.16	0.45
1:A:261:LEU:HD22	1:B:259:THR:HG21	1.98	0.44
1:A:371:LEU:HD23	1:A:372:ALA:N	2.32	0.44
1:A:269:ARG:NH1	1:A:269:ARG:HG3	2.32	0.43
1:A:393:GLN:HB3	1:A:441:ARG:CZ	2.48	0.43
1:B:278:ALA:HB2	1:B:299:ALA:HB2	2.00	0.43
1:B:321:ASN:HB2	1:B:352:TYR:HE1	1.84	0.43
1:A:371:LEU:C	1:A:371:LEU:HD23	2.39	0.42
1:B:416:THR:CG2	1:B:417:ASP:N	2.83	0.42
1:B:371:LEU:HD11	1:B:411:LEU:HB2	2.03	0.41
1:A:473:LEU:HD23	1:B:473:LEU:CD2	2.51	0.41
1:B:262:TRP:HB3	1:B:473:LEU:HG	2.03	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	225/254 (89%)	217 (96%)	7 (3%)	1 (0%)	38	16
1	B	216/254 (85%)	210 (97%)	5 (2%)	1 (0%)	32	12
All	All	441/508 (87%)	427 (97%)	12 (3%)	2 (0%)	32	12

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	417	ASP
1	B	417	ASP

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	198/221 (90%)	196 (99%)	2 (1%)	80	65
1	B	190/221 (86%)	188 (99%)	2 (1%)	78	61
All	All	388/442 (88%)	384 (99%)	4 (1%)	80	65

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	321	ASN
1	A	323	MET
1	B	287	LEU
1	B	323	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	300	ASN
1	A	321	ASN

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Mol	Chain	Res	Type
1	A	329	ASN
1	A	414	GLN
1	B	376	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 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	MPD	A	2001	-	7,7,7	0.51	0	9,10,10	0.47	0
3	MPD	A	3001	-	7,7,7	0.39	0	9,10,10	0.65	0
2	TRS	A	5001	-	7,7,7	1.98	2 (28%)	9,9,9	2.37	5 (55%)
3	MPD	B	1001	-	7,7,7	0.44	0	9,10,10	0.61	0
3	MPD	B	4001	-	7,7,7	0.55	0	9,10,10	0.42	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MPD	A	2001	-	-	0/5/5/5	0/0/0/0
3	MPD	A	3001	-	-	0/5/5/5	0/0/0/0
2	TRS	A	5001	-	-	0/9/9/9	0/0/0/0
3	MPD	B	1001	-	-	0/5/5/5	0/0/0/0
3	MPD	B	4001	-	-	0/5/5/5	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	5001	TRS	C2-C	-4.68	1.41	1.52
2	A	5001	TRS	O1-C1	-2.28	1.34	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	5001	TRS	C2-C-C1	-2.38	104.31	111.06
2	A	5001	TRS	C3-C-C1	-2.26	104.65	111.06
2	A	5001	TRS	O2-C2-C	2.10	116.56	110.47
2	A	5001	TRS	C3-C-N	3.76	115.73	107.73
2	A	5001	TRS	O3-C3-C	4.22	122.70	110.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2001	MPD	1	0
3	A	3001	MPD	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	227/254 (89%)	0.39	20 (8%) 11 10	16, 26, 59, 94	0
1	B	218/254 (85%)	0.26	14 (6%) 20 19	14, 27, 55, 96	0
All	All	445/508 (87%)	0.33	34 (7%) 15 14	14, 27, 58, 96	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	418	PHE	13.2
1	B	418	PHE	10.6
1	A	419	VAL	8.0
1	A	417	ASP	7.8
1	A	250	ALA	7.5
1	B	417	ASP	7.4
1	B	419	VAL	6.6
1	A	256	VAL	5.8
1	A	416	THR	5.2
1	A	261	LEU	5.1
1	A	258	LYS	4.5
1	B	476	SER	4.1
1	B	261	LEU	3.9
1	B	259	THR	3.8
1	A	262	TRP	3.5
1	B	262	TRP	3.3
1	A	255	VAL	3.3
1	A	476	SER	3.2
1	B	366	VAL	3.2
1	A	254	ILE	3.2
1	A	259	THR	3.1
1	B	416	THR	2.9
1	A	257	SER	2.9
1	A	474	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	308	ASP	2.7
1	A	265	MET	2.5
1	B	260	SER	2.5
1	A	251	ASN	2.4
1	A	473	LEU	2.4
1	B	265	MET	2.3
1	B	443	ARG	2.3
1	A	267	TYR	2.3
1	B	267	TYR	2.0
1	A	443	ARG	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	TRS	A	5001	8/8	0.68	0.21	23.59	71,71,71,72	0
3	MPD	A	2001	8/8	0.80	0.26	11.99	52,53,53,53	0
3	MPD	A	3001	8/8	0.73	0.28	11.37	60,61,61,62	0
3	MPD	B	4001	8/8	0.47	0.30	7.77	53,54,55,55	0
3	MPD	B	1001	8/8	0.81	0.16	2.33	29,35,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.