Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:27 PM GMT

PDB ID	:	1AR8
Title	:	P1/MAHONEY POLIOVIRUS, MUTANT P1095S
Authors	:	Wien, M.W.; Curry, S.; Filman, D.J.; Hogle, J.M.
Deposited on	:	1997-08-11
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

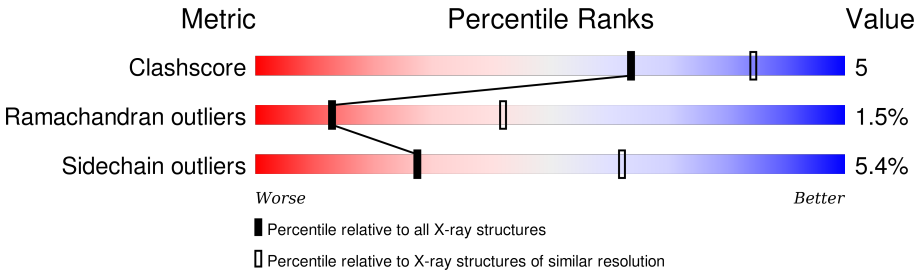
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (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 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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	0	8	25% 75%
2	1	302	77% 15% •• 6%
3	2	272	79% 16% • ••
4	3	238	84% 13% ••
5	4	68	66% 18% • 12%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7184 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 P1/MAHONEY POLIOVIRUS.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	0	8	Total	C	N	O	0	0	0
			43	24	8	11			

- Molecule 2 is a protein called P1/MAHONEY POLIOVIRUS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	1	283	Total	C	N	O	S	0	0	0
			2221	1414	378	424	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	95	SER	PRO	ENGINEERED	UNP P03300

- Molecule 3 is a protein called P1/MAHONEY POLIOVIRUS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	2	268	Total	C	N	O	S	0	0	0
			2085	1317	358	396	14			

- Molecule 4 is a protein called P1/MAHONEY POLIOVIRUS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	3	235	Total	C	N	O	S	0	0	0
			1834	1169	299	349	17			

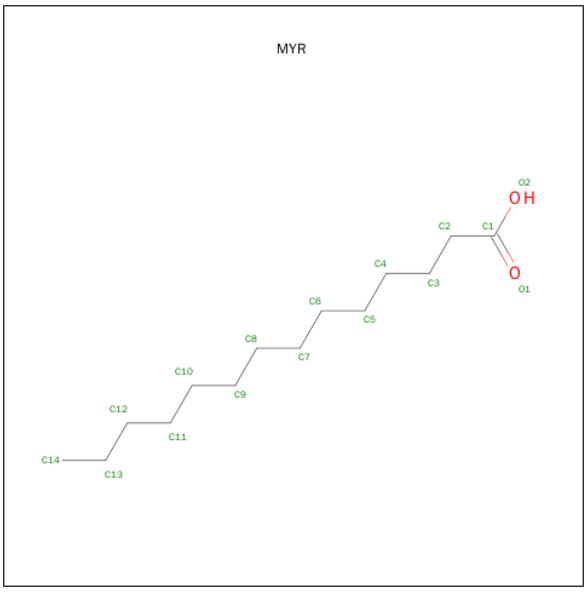
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	123	SER	PHE	CONFLICT	UNP P03300

- Molecule 5 is a protein called P1/MAHONEY POLIOVIRUS.

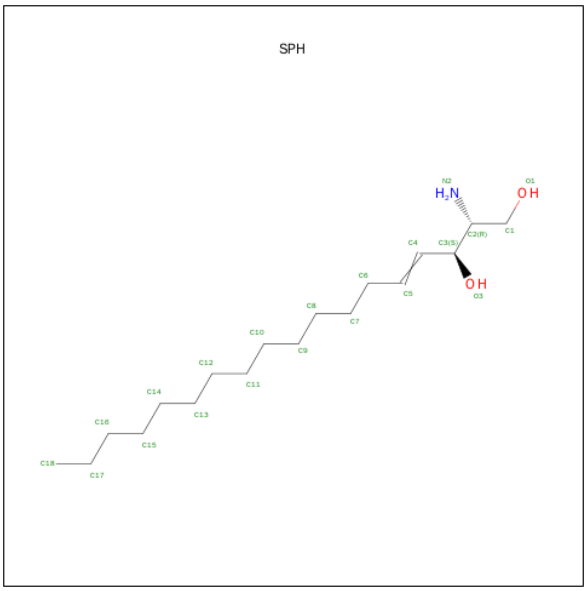
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	4	60	Total	C	N	O	S	0	0	0
			462	286	78	97	1			

- Molecule 6 is MYRISTIC ACID (three-letter code: MYR) (formula: $C_{14}H_{28}O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	4	1	Total	C	O		0	0
			15	14	1			

- Molecule 7 is SPHINGOSINE (three-letter code: SPH) (formula: $C_{18}H_{37}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	1	1	Total	C	N	O	0	0
			21	18	1	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	0	1	Total	O	0	0
			1	1		
8	1	184	Total	O	0	0
			184	184		
8	2	143	Total	O	0	0
			143	143		
8	3	135	Total	O	0	0
			135	135		
8	4	40	Total	O	0	0
			40	40		

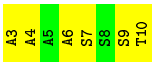
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

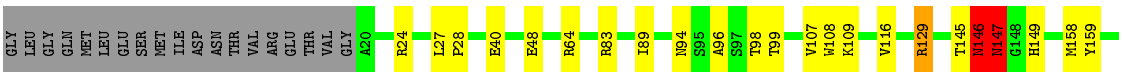
- Molecule 1: P1/MAHONEY POLIOVIRUS

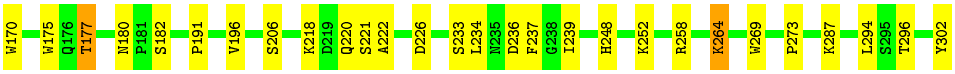
Chain 0:



- Molecule 2: P1/MAHONEY POLIOVIRUS

Chain 1:



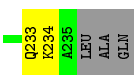


- Molecule 3: P1/MAHONEY POLIOVIRUS

Chain 2:

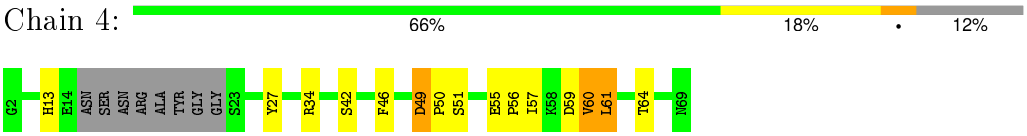
- Molecule 4: P1/MAHONEY POLIOVIRUS

Chain 3:



- Molecule 5: P1/MAHONEY POLIOVIRUS

Chain 4:


66%
18%
12%
G2
H13
E14
ASN
SER
ASN
ARG
ALA
TYR
GLY
GLY
S23
Y27
R34
S42
F46
D49
P50
S51
E55
P56
I57
K58
D59
V60
L61
T64
N69

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	323.04Å 358.22Å 380.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	11.00 – 2.90	Depositor
% Data completeness (in resolution range)	34.0 (11.00-2.90)	Depositor
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.0	Depositor
R, R_{free}	0.270 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7184	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SPH, MYR

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	0	0.47	0/41	1.53	0/53
2	1	0.81	0/2283	1.36	23/3120 (0.7%)
3	2	0.84	1/2142 (0.0%)	1.43	20/2928 (0.7%)
4	3	0.79	0/1881	1.28	12/2562 (0.5%)
5	4	0.75	0/469	1.37	3/632 (0.5%)
All	All	0.81	1/6816 (0.0%)	1.36	58/9295 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	2	7	CYS	CB-SG	5.06	1.90	1.82

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	2	5	GLU	OE1-CD-OE2	-10.41	110.80	123.30
3	2	78	TRP	CD1-CG-CD2	8.50	113.10	106.30
2	1	175	TRP	CD1-CG-CD2	8.48	113.08	106.30
2	1	170	TRP	CD1-CG-CD2	8.36	112.99	106.30
4	3	110	TRP	CD1-CG-CD2	8.33	112.96	106.30
3	2	80	TRP	CD1-CG-CD2	8.14	112.81	106.30
3	2	71	TRP	CE2-CD2-CG	-8.03	100.87	107.30
2	1	175	TRP	CE2-CD2-CG	-8.02	100.89	107.30
4	3	170	TRP	CE2-CD2-CG	-7.95	100.94	107.30
3	2	38	TRP	CD1-CG-CD2	7.93	112.64	106.30
3	2	80	TRP	CE2-CD2-CG	-7.93	100.96	107.30
3	2	227	TRP	CD1-CG-CD2	7.91	112.63	106.30
3	2	227	TRP	CE2-CD2-CG	-7.80	101.06	107.30
4	3	110	TRP	CE2-CD2-CG	-7.78	101.08	107.30
4	3	170	TRP	CD1-CG-CD2	7.75	112.50	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	1	269	TRP	CD1-CG-CD2	7.67	112.44	106.30
3	2	78	TRP	CE2-CD2-CG	-7.63	101.20	107.30
3	2	79	TRP	CD1-CG-CD2	7.59	112.37	106.30
4	3	156	TRP	CE2-CD2-CG	-7.57	101.25	107.30
5	4	34	ARG	NE-CZ-NH1	7.56	124.08	120.30
3	2	79	TRP	CE2-CD2-CG	-7.37	101.40	107.30
2	1	269	TRP	CE2-CD2-CG	-7.35	101.42	107.30
3	2	71	TRP	CD1-CG-CD2	7.34	112.17	106.30
2	1	108	TRP	CE2-CD2-CG	-7.23	101.52	107.30
2	1	108	TRP	CD1-CG-CD2	7.20	112.06	106.30
3	2	38	TRP	CE2-CD2-CG	-7.13	101.60	107.30
4	3	156	TRP	CD1-CG-CD2	6.96	111.87	106.30
2	1	170	TRP	CE2-CD2-CG	-6.94	101.75	107.30
2	1	129	ARG	NE-CZ-NH1	6.71	123.65	120.30
2	1	129	ARG	NE-CZ-NH2	-6.63	116.98	120.30
3	2	270	ARG	NE-CZ-NH1	6.11	123.36	120.30
2	1	24	ARG	NE-CZ-NH2	-6.03	117.28	120.30
2	1	175	TRP	CG-CD2-CE3	6.01	139.31	133.90
3	2	264	ARG	NE-CZ-NH1	5.93	123.27	120.30
5	4	34	ARG	NE-CZ-NH2	-5.84	117.38	120.30
3	2	5	GLU	CG-CD-OE1	5.75	129.80	118.30
5	4	27	TYR	CB-CG-CD2	-5.70	117.58	121.00
3	2	78	TRP	CG-CD1-NE1	-5.69	104.41	110.10
3	2	227	TRP	CG-CD2-CE3	5.57	138.92	133.90
2	1	147	ASN	OD1-CG-ND2	-5.52	109.21	121.90
2	1	170	TRP	CG-CD1-NE1	-5.51	104.58	110.10
4	3	170	TRP	CG-CD2-CE3	5.51	138.86	133.90
2	1	83	ARG	NE-CZ-NH1	5.44	123.02	120.30
2	1	146	ASN	N-CA-CB	-5.44	100.80	110.60
3	2	71	TRP	CG-CD2-CE3	5.37	138.73	133.90
2	1	175	TRP	CB-CG-CD1	-5.30	120.11	127.00
4	3	206	ARG	NE-CZ-NH2	-5.30	117.65	120.30
3	2	168	THR	CA-CB-CG2	5.29	119.80	112.40
4	3	110	TRP	CG-CD1-NE1	-5.25	104.85	110.10
4	3	226	ARG	NE-CZ-NH1	5.22	122.91	120.30
2	1	226	ASP	N-CA-C	5.16	124.94	111.00
2	1	64	ARG	NE-CZ-NH1	5.11	122.86	120.30
4	3	110	TRP	CG-CD2-CE3	5.08	138.47	133.90
2	1	147	ASN	CB-CG-ND2	5.07	128.87	116.70
4	3	65	MET	CG-SD-CE	-5.03	92.16	100.20
2	1	108	TRP	CG-CD2-CE3	5.02	138.42	133.90
2	1	196	VAL	CG1-CB-CG2	-5.01	102.88	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	1	264	LYS	CA-CB-CG	5.01	124.43	113.40

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	0	43	0	39	2	0
2	1	2221	0	2171	29	0
3	2	2085	0	2000	27	0
4	3	1834	0	1816	18	0
5	4	462	0	446	7	0
6	4	15	0	27	1	0
7	1	21	0	37	4	0
8	0	1	0	0	0	0
8	1	184	0	0	3	0
8	2	143	0	0	2	0
8	3	135	0	0	2	0
8	4	40	0	0	0	0
All	All	7184	0	6536	69	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 5.

All (69) 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)
2:1:158:MET:SD	2:1:177:THR:HG23	2.16	0.85
3:2:5:GLU:OE2	3:2:7:CYS:HB3	1.78	0.83
4:3:124:MET:HG3	8:3:277:HOH:O	1.81	0.80
1:0:6:ALA:CA	1:0:7:SER:N	2.48	0.76
2:1:177:THR:HG22	2:1:180:ASN:HB2	1.69	0.73
2:1:40:GLU:HB3	5:4:64:THR:HB	1.71	0.72
2:1:287:LYS:HA	3:2:137:ASN:HD21	1.55	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2:5:GLU:HG2	3:2:9:TYR:HD2	1.57	0.70
2:1:218:LYS:HD2	8:2:405:HOH:O	1.93	0.68
2:1:107:VAL:HG13	2:1:239:ILE:HD13	1.75	0.68
2:1:109:LYS:HA	2:1:239:ILE:HG22	1.78	0.66
2:1:236:ASP:HB3	2:1:237:PHE:CD2	2.32	0.64
3:2:5:GLU:OE2	3:2:7:CYS:CB	2.46	0.63
2:1:218:LYS:HA	3:2:270:ARG:HB2	1.80	0.63
3:2:187:LEU:HD22	4:3:65:MET:CE	2.29	0.61
3:2:143:THR:HG23	3:2:173:ARG:HA	1.80	0.61
3:2:213:VAL:HG22	4:3:37:PRO:HG2	1.86	0.58
2:1:177:THR:HG21	2:1:182:SER:OG	2.04	0.57
5:4:57:ILE:HD11	5:4:61:LEU:HB3	1.87	0.57
7:1:0:SPH:H4	8:1:448:HOH:O	2.04	0.56
3:2:179:TYR:HA	4:3:65:MET:CE	2.36	0.56
2:1:159:TYR:HB2	7:1:0:SPH:H162	1.87	0.56
3:2:146:GLN:NE2	3:2:270:ARG:HD2	2.22	0.55
2:1:48:GLU:HA	3:2:197:ILE:HB	1.89	0.55
2:1:287:LYS:HA	3:2:137:ASN:ND2	2.21	0.54
2:1:89:ILE:HG12	2:1:258:ARG:HG2	1.91	0.53
2:1:191:PRO:HG2	4:3:13:TYR:HB2	1.92	0.52
3:2:48:ASN:HB3	3:2:49:PRO:HD3	1.92	0.52
8:3:317:HOH:O	5:4:46:PHE:HB2	2.11	0.51
2:1:222:ALA:HA	8:1:336:HOH:O	2.09	0.51
2:1:149:HIS:CE1	8:1:303:HOH:O	2.64	0.50
2:1:96:ALA:HB2	2:1:248:HIS:CD2	2.47	0.50
2:1:27:LEU:HB3	2:1:28:PRO:HD2	1.94	0.50
4:3:167:VAL:O	4:3:169:PRO:HD3	2.11	0.49
3:2:30:ASN:HD21	5:4:59:ASP:HB2	1.76	0.49
4:3:55:PHE:HE2	4:3:212:GLY:HA3	1.78	0.49
2:1:302:TYR:CE1	4:3:189:TYR:HB3	2.49	0.48
3:2:60:ALA:O	3:2:255:PRO:HG2	2.15	0.47
4:3:64:THR:O	4:3:67:MET:HG2	2.14	0.47
3:2:73:LYS:HE2	3:2:243:SER:O	2.15	0.47
5:4:13:HIS:H	5:4:13:HIS:CD2	2.32	0.47
4:3:87:LEU:HD13	4:3:190:ILE:HD11	1.97	0.47
2:1:273:PRO:HB3	3:2:189:ASN:HB2	1.97	0.47
2:1:206:SER:O	7:1:0:SPH:H11	2.14	0.46
2:1:294:LEU:HD12	4:3:67:MET:SD	2.56	0.45
4:3:149:MET:CE	4:3:150:LEU:HG	2.46	0.45
2:1:94:ASN:ND2	2:1:248:HIS:HA	2.31	0.45
3:2:34:ALA:HB3	3:2:211:PRO:HD2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:296:THR:HA	4:3:57:LEU:O	2.17	0.45
3:2:5:GLU:CD	3:2:7:CYS:H	2.20	0.44
3:2:63:PHE:CD1	3:2:254:ALA:HB2	2.53	0.43
5:4:49:ASP:HA	5:4:50:PRO:HD3	1.88	0.43
3:2:37:ARG:HG3	4:3:37:PRO:HB3	2.00	0.43
4:3:143:LYS:HD3	4:3:143:LYS:HA	1.75	0.43
3:2:138:THR:HG23	8:2:395:HOH:O	2.19	0.43
2:1:96:ALA:O	2:1:98:THR:HG23	2.19	0.42
2:1:237:PHE:CD2	7:1:0:SPH:H71	2.55	0.42
5:4:55:GLU:N	5:4:56:PRO:HD3	2.34	0.42
6:4:1:MYR:H71	6:4:1:MYR:H101	1.80	0.42
4:3:71:ARG:HB2	4:3:71:ARG:CZ	2.49	0.41
2:1:96:ALA:HB2	2:1:248:HIS:NE2	2.35	0.41
2:1:116:VAL:CG2	4:3:233:GLN:HB2	2.50	0.41
3:2:13:VAL:HG22	3:2:26:GLN:HA	2.03	0.41
1:0:3:ALA:O	1:0:6:ALA:N	2.53	0.41
3:2:37:ARG:O	3:2:211:PRO:HG3	2.20	0.40
3:2:166:ASN:OD1	3:2:168:THR:HG22	2.22	0.40
4:3:88:SER:HB3	4:3:91:SER:OG	2.22	0.40
3:2:146:GLN:HE21	3:2:270:ARG:HD2	1.85	0.40
3:2:239:PHE:O	3:2:241:SER:N	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	0	4/8 (50%)	3 (75%)	0	1 (25%)	0 0
2	1	281/302 (93%)	256 (91%)	21 (8%)	4 (1%)	14 44
3	2	266/272 (98%)	246 (92%)	15 (6%)	5 (2%)	10 35
4	3	233/238 (98%)	222 (95%)	10 (4%)	1 (0%)	39 74

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
5	4	56/68 (82%)	51 (91%)	3 (5%)	2 (4%)	4 18
All	All	840/888 (95%)	778 (93%)	49 (6%)	13 (2%)	13 42

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	1	99	THR
3	2	166	ASN
1	0	4	ALA
2	1	147	ASN
2	1	234	LEU
3	2	240	ALA
5	4	60	VAL
2	1	146	ASN
3	2	48	ASN
4	3	234	LYS
3	2	7	CYS
3	2	114	ALA
5	4	51	SER

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	0	4/4 (100%)	2 (50%)	2 (50%)	0 0
2	1	245/261 (94%)	235 (96%)	10 (4%)	37 73
3	2	228/232 (98%)	213 (93%)	15 (7%)	21 51
4	3	210/212 (99%)	201 (96%)	9 (4%)	35 71
5	4	52/57 (91%)	48 (92%)	4 (8%)	16 42
All	All	739/766 (96%)	699 (95%)	40 (5%)	27 62

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	0	9	SER
1	0	10	THR
2	1	129	ARG
2	1	145	THR
2	1	146	ASN
2	1	147	ASN
2	1	177	THR
2	1	220	GLN
2	1	221	SER
2	1	233	SER
2	1	252	LYS
2	1	264	LYS
3	2	10	SER
3	2	11	ASP
3	2	45	SER
3	2	52	GLN
3	2	73	LYS
3	2	128	PRO
3	2	139	THR
3	2	150	PRO
3	2	152	GLU
3	2	158	THR
3	2	165	ASN
3	2	168	THR
3	2	241	SER
3	2	251	LEU
3	2	264	ARG
4	3	33	PRO
4	3	71	ARG
4	3	85	LEU
4	3	146	LYS
4	3	149	MET
4	3	163	SER
4	3	208	MET
4	3	218	ASN
4	3	224	LEU
5	4	42	SER
5	4	49	ASP
5	4	60	VAL
5	4	61	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	1	69	HIS
3	2	48	ASN
3	2	165	ASN
4	3	218	ASN
5	4	13	HIS
5	4	31	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

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	$\#\|Z\| > 2$
7	SPH	1	0	-	19,20,20	0.42	0	16,21,21	1.98	3 (18%)
6	MYR	4	1	5	14,14,15	0.49	0	12,13,15	0.89	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
7	SPH	1	0	-	-	0/21/21/21	0/0/0/0
6	MYR	4	1	5	-	0/11/12/13	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	1	0	SPH	O3-C3-C2	-4.80	99.47	107.48
7	1	0	SPH	O3-C3-C4	-4.40	98.42	110.78
7	1	0	SPH	O1-C1-C2	3.35	118.57	111.53

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
7	1	0	SPH	4	0
6	4	1	MYR	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.