



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:57 PM GMT

PDB ID : 1J4X  
Title : HUMAN VH1-RELATED DUAL-SPECIFICITY PHOSPHATASE C124S  
MUTANT-PEPTIDE COMPLEX  
Authors : Schumacher, M.A.; Todd, J.L.; Tanner, K.G.; Denu, J.M.  
Deposited on : 2001-12-13  
Resolution : 2.75 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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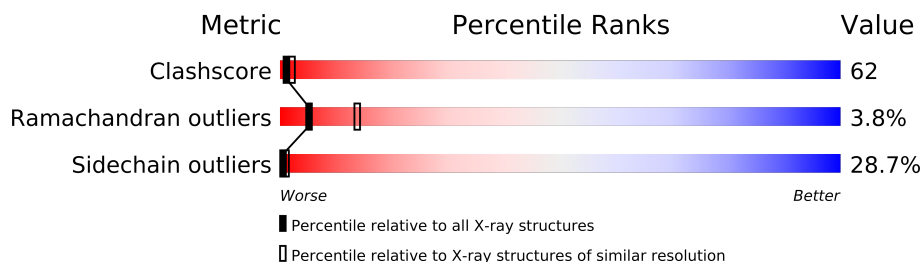
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.75 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	79885	2995 (2.80-2.72)
Ramachandran outliers	78287	2941 (2.80-2.72)
Sidechain outliers	78261	2944 (2.80-2.72)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	184	
2	D	11	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1546 atoms, of which 0 are hydrogen and 0 are deuterium.

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 DUAL SPECIFICITY PROTEIN PHOSPHATASE 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	178	Total	C	N	O	S	0	0	0
			1384	863	248	266	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	SER	CYS	ENGINEERED	UNP P51452

- Molecule 2 is a protein called DDE(AHP)(TPO)G(PTR)VATR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	11	Total	C	N	O	P	0	0	0
			96	53	14	27	2			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total	O	0	0
			57	57		
3	D	9	Total	O	0	0
			9	9		

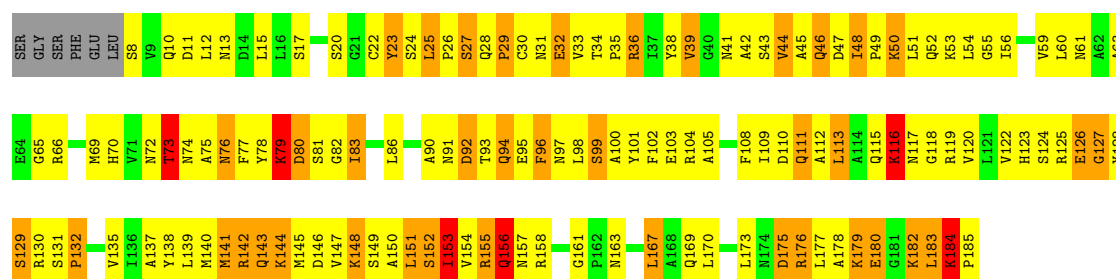
### 3 Residue-property plots

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.

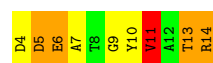
#### • Molecule 1: DUAL SPECIFICITY PROTEIN PHOSPHATASE 3

Chain A: 



#### • Molecule 2: DDE(AHP)(TPO)G(PTR)VATR

Chain D: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	34.39Å 55.77Å 101.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.75	Depositor
% Data completeness (in resolution range)	90.0 (10.00-2.75)	Depositor
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.188 , 0.260	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1546	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	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: TPO, AHP, PTR

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	4.64	4/1406 (0.3%)	1.67	21/1897 (1.1%)
2	D	1.79	1/57 (1.8%)	2.75	5/71 (7.0%)
All	All	4.57	5/1463 (0.3%)	1.72	26/1968 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	3	0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	36	ARG	CB-CG	148.90	5.54	1.52
1	A	36	ARG	CG-CD	81.25	3.55	1.51
1	A	180	GLU	CG-CD	-6.17	1.42	1.51
1	A	184	LYS	CG-CD	-5.94	1.32	1.52
2	D	13	THR	CA-CB	5.15	1.66	1.53

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	ARG	CB-CG-CD	-31.80	28.92	111.60
1	A	36	ARG	CG-CD-NE	15.76	144.89	111.80
1	A	36	ARG	CA-CB-CG	-14.65	81.17	113.40
2	D	11	VAL	N-CA-C	9.49	136.62	111.00
2	D	11	VAL	C-N-CA	6.92	138.99	121.70
2	D	6	GLU	N-CA-C	6.89	129.61	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	ARG	NE-CZ-NH2	-6.87	116.87	120.30
2	D	13	THR	N-CA-CB	-6.85	97.28	110.30
1	A	155	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	A	80	ASP	CA-CB-CG	-6.71	98.63	113.40
1	A	92	ASP	CB-CG-OD2	-6.35	112.59	118.30
2	D	14	ARG	N-CA-CB	6.28	121.90	110.60
1	A	167	LEU	CA-CB-CG	-6.23	100.97	115.30
1	A	29	PRO	N-CA-C	6.16	128.12	112.10
1	A	79	LYS	O-C-N	6.02	132.34	122.70
1	A	80	ASP	N-CA-CB	5.96	121.32	110.60
1	A	32	GLU	N-CA-C	-5.83	95.26	111.00
1	A	153	ILE	CB-CA-C	-5.58	100.44	111.60
1	A	127	GLY	N-CA-C	-5.53	99.28	113.10
1	A	116	LYS	N-CA-C	5.46	125.75	111.00
1	A	113	LEU	CB-CG-CD1	5.45	120.26	111.00
1	A	20	SER	N-CA-CB	5.42	118.63	110.50
1	A	79	LYS	C-N-CA	5.15	134.57	121.70
1	A	156	GLN	CB-CA-C	-5.12	100.16	110.40
1	A	176	ARG	N-CA-C	-5.06	97.35	111.00
1	A	73	THR	N-CA-C	-5.04	97.40	111.00

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	6	GLU	CA
2	D	8	TPO	CA
2	D	11	VAL	CA

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1384	0	1378	171	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	96	0	75	43	0
3	A	57	0	0	15	0
3	D	9	0	0	0	0
All	All	1546	0	1453	183	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 62.

All (183) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:128:TYR:CD2	2:D:11:VAL:HB	1.82	1.15
1:A:33:VAL:HG21	1:A:154:VAL:HG13	1.36	1.04
1:A:178:ALA:HB2	1:A:183:LEU:HD12	1.42	1.01
1:A:163:ASN:HA	2:D:13:THR:HB	1.42	0.99
1:A:44:VAL:HG13	1:A:50:LYS:HD3	1.42	0.99
1:A:126:GLU:HB2	2:D:10:PTR:HE1	1.40	0.98
1:A:39:VAL:HG12	1:A:158:ARG:HD3	1.46	0.98
1:A:128:TYR:HD2	2:D:11:VAL:HB	1.24	0.96
1:A:126:GLU:HG2	2:D:10:PTR:HD1	1.48	0.93
1:A:48:ILE:HA	1:A:51:LEU:HD12	1.48	0.93
2:D:10:PTR:CE2	2:D:11:VAL:HG22	1.98	0.93
1:A:25:LEU:HD22	2:D:9:GLY:HA2	1.50	0.92
1:A:97:ASN:ND2	1:A:100:ALA:HB2	1.85	0.92
1:A:44:VAL:CG1	1:A:50:LYS:HD3	2.03	0.89
1:A:32:GLU:HB2	1:A:38:TYR:CE2	2.09	0.88
1:A:36:ARG:HG2	1:A:119:ARG:NH1	1.92	0.85
1:A:126:GLU:CG	2:D:10:PTR:HD1	2.07	0.84
1:A:38:TYR:CD1	1:A:56:ILE:HD11	2.11	0.84
1:A:128:TYR:CE2	2:D:11:VAL:HB	2.14	0.83
1:A:25:LEU:CD2	2:D:9:GLY:HA2	2.09	0.82
1:A:163:ASN:HA	2:D:13:THR:CB	2.10	0.80
1:A:36:ARG:HB3	1:A:113:LEU:HD21	1.64	0.80
1:A:128:TYR:CD2	2:D:11:VAL:CB	2.64	0.80
1:A:102:PHE:N	3:A:252:HOH:O	2.16	0.79
1:A:36:ARG:CB	1:A:36:ARG:HD2	2.12	0.79
1:A:39:VAL:CG1	1:A:158:ARG:HD3	2.14	0.76
1:A:129:SER:HB2	2:D:11:VAL:HG21	1.68	0.75
1:A:180:GLU:HB2	1:A:182:LYS:HG3	1.69	0.75
1:A:125:ARG:HB3	2:D:10:PTR:O2P	1.88	0.74
1:A:36:ARG:HB3	1:A:113:LEU:CD2	2.17	0.73
1:A:105:ALA:O	1:A:109:ILE:HD12	1.88	0.73
1:A:126:GLU:HB2	2:D:10:PTR:CE1	2.16	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:150:ALA:O	1:A:154:VAL:HG23	1.89	0.72
2:D:10:PTR:CE2	2:D:11:VAL:CG2	2.68	0.71
1:A:128:TYR:HD2	2:D:11:VAL:CB	2.02	0.71
1:A:91:ASN:HB2	1:A:96:PHE:CD1	2.26	0.70
1:A:176:ARG:O	1:A:180:GLU:HG3	1.90	0.70
1:A:116:LYS:CA	3:A:240:HOH:O	2.40	0.69
1:A:28:GLN:HB2	1:A:29:PRO:HD2	1.73	0.69
1:A:161:GLY:O	2:D:11:VAL:CG1	2.41	0.68
1:A:153:ILE:O	1:A:157:ASN:ND2	2.27	0.68
1:A:8:SER:O	1:A:11:ASP:HB2	1.94	0.68
1:A:76:ASN:ND2	1:A:76:ASN:H	1.93	0.67
1:A:41:ASN:HB2	1:A:125:ARG:O	1.94	0.67
1:A:163:ASN:CA	2:D:13:THR:HB	2.23	0.67
1:A:101:TYR:HA	3:A:253:HOH:O	1.95	0.67
1:A:30:CYS:O	1:A:31:ASN:ND2	2.27	0.66
1:A:74:ASN:O	1:A:77:PHE:HB3	1.95	0.66
1:A:101:TYR:HD2	1:A:104:ARG:HD2	1.60	0.66
1:A:122:VAL:HG11	1:A:135:VAL:HG23	1.76	0.66
1:A:103:GLU:HG2	3:A:248:HOH:O	1.95	0.66
1:A:110:ASP:OD2	1:A:142:ARG:NE	2.28	0.66
1:A:98:LEU:O	3:A:252:HOH:O	2.15	0.65
1:A:10:GLN:OE1	1:A:13:ASN:ND2	2.29	0.64
1:A:90:ALA:C	1:A:91:ASN:HD22	2.00	0.64
2:D:11:VAL:HG12	2:D:13:THR:HG22	1.79	0.64
1:A:175:ASP:N	1:A:175:ASP:OD1	2.28	0.64
1:A:111:GLN:HG2	3:A:233:HOH:O	1.97	0.64
1:A:50:LYS:HZ2	1:A:54:LEU:HD21	1.64	0.63
1:A:177:LEU:HD23	1:A:177:LEU:N	2.13	0.63
1:A:94:GLN:NE2	2:D:14:ARG:O	2.30	0.62
1:A:142:ARG:HG2	1:A:142:ARG:HH11	1.63	0.62
1:A:15:LEU:HD12	1:A:151:LEU:HD22	1.82	0.62
1:A:117:ASN:N	3:A:240:HOH:O	2.32	0.62
1:A:108:PHE:HB2	3:A:213:HOH:O	2.00	0.61
1:A:28:GLN:O	1:A:158:ARG:NH1	2.33	0.61
1:A:143:GLN:HB3	1:A:145:MET:HE3	1.83	0.61
1:A:109:ILE:O	1:A:112:ALA:HB3	2.02	0.60
1:A:27:SER:HB2	2:D:5:ASP:OD1	2.01	0.60
1:A:33:VAL:HG23	1:A:34:THR:HG23	1.84	0.60
1:A:101:TYR:N	3:A:252:HOH:O	2.34	0.60
1:A:116:LYS:HE3	3:A:251:HOH:O	2.02	0.60
1:A:99:SER:HB3	1:A:169:GLN:OE1	2.02	0.59
1:A:183:LEU:O	1:A:184:LYS:HB2	2.02	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:175:ASP:O	1:A:179:LYS:HE3	2.03	0.59
1:A:140:MET:HE2	1:A:147:VAL:CB	2.33	0.59
1:A:73:THR:HG22	1:A:77:PHE:CE2	2.38	0.58
1:A:28:GLN:HB2	1:A:29:PRO:CD	2.33	0.58
1:A:155:ARG:O	1:A:155:ARG:HD2	2.04	0.58
1:A:140:MET:HE2	1:A:147:VAL:HA	1.85	0.58
1:A:26:PRO:HD3	1:A:128:TYR:CE1	2.38	0.58
1:A:153:ILE:HD13	1:A:153:ILE:N	2.18	0.57
1:A:70:HIS:HE1	1:A:90:ALA:O	1.88	0.57
1:A:32:GLU:HA	1:A:38:TYR:HA	1.86	0.57
1:A:42:ALA:O	1:A:45:ALA:HB3	2.04	0.57
1:A:36:ARG:HG2	1:A:119:ARG:CZ	2.35	0.56
1:A:138:TYR:O	1:A:142:ARG:HB2	2.05	0.56
1:A:140:MET:CE	1:A:147:VAL:HG23	2.35	0.56
2:D:10:PTR:CZ	2:D:11:VAL:CG2	2.83	0.56
1:A:140:MET:HE2	1:A:147:VAL:HB	1.86	0.56
1:A:116:LYS:N	3:A:240:HOH:O	2.39	0.55
1:A:126:GLU:HG2	2:D:10:PTR:N	2.22	0.55
1:A:32:GLU:HB2	1:A:38:TYR:CD2	2.40	0.54
1:A:143:GLN:HB3	1:A:145:MET:CE	2.37	0.54
1:A:39:VAL:O	1:A:39:VAL:HG12	2.07	0.54
1:A:137:ALA:O	1:A:141:MET:HG2	2.07	0.54
1:A:48:ILE:HG22	1:A:49:PRO:CD	2.38	0.53
1:A:126:GLU:HG2	2:D:9:GLY:C	2.28	0.53
1:A:97:ASN:HD21	1:A:100:ALA:HB2	1.70	0.53
1:A:180:GLU:HB2	1:A:182:LYS:CG	2.38	0.52
1:A:81:SER:OG	1:A:83:ILE:HG13	2.09	0.52
1:A:48:ILE:HG22	1:A:49:PRO:HD3	1.91	0.52
1:A:184:LYS:N	1:A:185:PRO:HD3	2.15	0.52
1:A:63:ALA:HB1	1:A:130:ARG:NH1	2.24	0.52
1:A:69:MET:HG2	3:A:236:HOH:O	2.10	0.52
1:A:25:LEU:HB3	1:A:26:PRO:HD2	1.92	0.52
1:A:128:TYR:CE2	2:D:11:VAL:CA	2.92	0.52
1:A:23:TYR:CD1	1:A:23:TYR:N	2.77	0.52
1:A:140:MET:HE2	1:A:147:VAL:CA	2.40	0.52
1:A:178:ALA:CB	1:A:183:LEU:HD12	2.30	0.52
1:A:105:ALA:C	1:A:109:ILE:HD12	2.30	0.51
2:D:7:AHP:HB2	2:D:7:AHP:HZ3	1.91	0.51
1:A:163:ASN:HA	2:D:13:THR:CA	2.40	0.51
1:A:161:GLY:O	2:D:11:VAL:HG12	2.11	0.51
1:A:52:GLN:O	1:A:55:GLY:N	2.41	0.51
1:A:76:ASN:HA	1:A:79:LYS:HB2	1.93	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:152:SER:O	1:A:156:GLN:HG3	2.12	0.50
1:A:90:ALA:HA	1:A:96:PHE:CZ	2.47	0.50
1:A:140:MET:CE	1:A:147:VAL:HB	2.42	0.50
1:A:26:PRO:HD3	1:A:128:TYR:HE1	1.76	0.49
1:A:92:ASP:HB2	1:A:130:ARG:NE	2.28	0.49
1:A:92:ASP:OD2	2:D:10:PTR:OH	2.30	0.49
1:A:48:ILE:HG22	1:A:49:PRO:N	2.28	0.49
1:A:131:SER:OG	1:A:132:PRO:HD3	2.13	0.49
1:A:60:LEU:HD12	1:A:86:LEU:O	2.12	0.49
1:A:60:LEU:HA	1:A:86:LEU:O	2.13	0.48
1:A:61:ASN:HA	1:A:123:HIS:CD2	2.49	0.48
1:A:142:ARG:CG	1:A:142:ARG:HH11	2.26	0.48
1:A:120:VAL:HG13	1:A:120:VAL:O	2.13	0.48
1:A:76:ASN:ND2	1:A:76:ASN:N	2.58	0.48
1:A:45:ALA:O	1:A:78:TYR:OH	2.27	0.48
1:A:124:SER:OG	1:A:125:ARG:N	2.47	0.47
2:D:4:ASP:OD2	2:D:5:ASP:N	2.45	0.47
1:A:140:MET:HE1	1:A:147:VAL:HG23	1.97	0.47
1:A:12:LEU:HD23	1:A:12:LEU:N	2.30	0.47
1:A:97:ASN:HD22	1:A:100:ALA:HB2	1.71	0.46
1:A:36:ARG:HD3	1:A:118:GLY:O	2.16	0.46
1:A:23:TYR:HD1	1:A:23:TYR:N	2.13	0.46
1:A:127:GLY:HA2	1:A:131:SER:OG	2.16	0.46
1:A:135:VAL:HG12	3:A:265:HOH:O	2.15	0.46
2:D:10:PTR:CZ	2:D:11:VAL:HG23	2.44	0.46
2:D:10:PTR:CG	2:D:11:VAL:N	2.79	0.46
1:A:122:VAL:HG12	1:A:131:SER:HB2	1.97	0.46
1:A:25:LEU:HD22	2:D:9:GLY:CA	2.35	0.46
1:A:63:ALA:CB	1:A:130:ARG:NH1	2.80	0.46
1:A:184:LYS:HB3	1:A:184:LYS:HE2	1.21	0.45
1:A:73:THR:HG22	1:A:77:PHE:CD2	2.51	0.45
1:A:47:ASP:OD1	1:A:49:PRO:HD2	2.17	0.45
1:A:73:THR:O	1:A:74:ASN:HB3	2.16	0.44
1:A:35:PRO:O	1:A:36:ARG:HB2	2.17	0.44
1:A:173:LEU:O	1:A:177:LEU:HG	2.17	0.44
2:D:4:ASP:O	2:D:5:ASP:HB2	2.17	0.44
1:A:167:LEU:HD23	1:A:167:LEU:HA	1.47	0.44
1:A:117:ASN:N	1:A:117:ASN:OD1	2.50	0.43
2:D:10:PTR:CG	2:D:11:VAL:H	2.31	0.43
1:A:139:LEU:CD1	3:A:265:HOH:O	2.66	0.43
1:A:36:ARG:HH11	1:A:118:GLY:N	2.17	0.43
1:A:59:VAL:HG12	1:A:60:LEU:N	2.34	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:MET:CE	1:A:147:VAL:CB	2.97	0.42
1:A:148:LYS:HE2	1:A:148:LYS:HB3	1.22	0.42
1:A:36:ARG:HD3	1:A:36:ARG:HH11	1.78	0.42
2:D:13:THR:HG23	2:D:13:THR:H	1.50	0.42
1:A:42:ALA:O	1:A:46:GLN:HG2	2.19	0.42
1:A:183:LEU:HA	1:A:183:LEU:HD23	1.40	0.42
1:A:93:THR:HA	3:A:221:HOH:O	2.20	0.42
1:A:128:TYR:CE2	2:D:11:VAL:CB	2.93	0.42
1:A:28:GLN:CB	1:A:29:PRO:CD	2.97	0.41
1:A:56:ILE:HD13	1:A:56:ILE:HA	1.78	0.41
1:A:142:ARG:CG	1:A:142:ARG:NH1	2.80	0.41
1:A:126:GLU:CB	2:D:10:PTR:HD1	2.50	0.41
1:A:27:SER:HB2	2:D:5:ASP:CG	2.41	0.41
1:A:56:ILE:HD13	1:A:119:ARG:HB3	2.02	0.41
1:A:61:ASN:ND2	1:A:73:THR:OG1	2.53	0.41
1:A:161:GLY:O	2:D:11:VAL:HB	2.21	0.41
1:A:98:LEU:C	1:A:100:ALA:H	2.23	0.41
1:A:36:ARG:CG	1:A:119:ARG:CZ	2.98	0.41
1:A:140:MET:HE1	1:A:147:VAL:CG2	2.51	0.41
1:A:65:GLY:O	1:A:72:ASN:HB2	2.20	0.41
1:A:150:ALA:HA	1:A:153:ILE:HG12	2.02	0.41
1:A:99:SER:HA	1:A:102:PHE:CE1	2.57	0.40
1:A:170:LEU:O	1:A:173:LEU:HB3	2.21	0.40
2:D:10:PTR:N	2:D:10:PTR:HD1	2.37	0.40
1:A:128:TYR:HE2	2:D:11:VAL:O	2.05	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	176/184 (96%)	142 (81%)	29 (16%)	5 (3%)	8	21
2	D	6/11 (54%)	4 (67%)	0	2 (33%)	0	0
All	All	182/195 (93%)	146 (80%)	29 (16%)	7 (4%)	5	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	LYS
2	D	11	VAL
1	A	82	GLY
1	A	183	LEU
2	D	5	ASP
1	A	75	ALA
1	A	144	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	151/156 (97%)	107 (71%)	44 (29%)	0	1
2	D	6/6 (100%)	5 (83%)	1 (17%)	3	8
All	All	157/162 (97%)	112 (71%)	45 (29%)	0	1

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	SER
1	A	22	CYS
1	A	23	TYR
1	A	24	SER
1	A	25	LEU
1	A	27	SER
1	A	39	VAL
1	A	43	SER
1	A	44	VAL
1	A	46	GLN
1	A	48	ILE
1	A	50	LYS
1	A	53	LYS
1	A	66	ARG
1	A	73	THR
1	A	76	ASN
1	A	79	LYS

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Mol	Chain	Res	Type
1	A	80	ASP
1	A	83	ILE
1	A	94	GLN
1	A	95	GLU
1	A	96	PHE
1	A	99	SER
1	A	111	GLN
1	A	115	GLN
1	A	116	LYS
1	A	126	GLU
1	A	129	SER
1	A	132	PRO
1	A	141	MET
1	A	142	ARG
1	A	143	GLN
1	A	144	LYS
1	A	146	ASP
1	A	148	LYS
1	A	149	SER
1	A	151	LEU
1	A	152	SER
1	A	153	ILE
1	A	156	GLN
1	A	175	ASP
1	A	179	LYS
1	A	182	LYS
1	A	184	LYS
2	D	6	GLU

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

Mol	Chain	Res	Type
1	A	70	HIS
1	A	74	ASN
1	A	76	ASN
1	A	91	ASN
1	A	97	ASN
1	A	143	GLN
1	A	157	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

3 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	PTR	D	10	2	16,16,17	4.56	6 (37%)	20,22,24	1.40	4 (20%)
2	AHP	D	7	2	8,8,9	6.95	3 (37%)	6,8,10	2.09	1 (16%)
2	TPO	D	8	2	10,10,11	6.30	6 (60%)	12,14,16	4.34	5 (41%)

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	PTR	D	10	2	-	0/9/11/13	0/1/1/1
2	AHP	D	7	2	-	0/5/7/9	0/0/0/0
2	TPO	D	8	2	1/1/3/4	0/9/11/13	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	7	AHP	O-C	18.55	1.24	1.11
2	D	8	TPO	O-C	16.21	1.22	1.11
2	D	10	PTR	O-C	14.68	1.21	1.11
2	D	10	PTR	P-OH	8.13	1.72	1.60
2	D	8	TPO	CB-CA	7.24	1.64	1.53
2	D	8	TPO	P-O2P	5.80	1.75	1.54
2	D	8	TPO	CA-C	5.56	1.59	1.48
2	D	7	AHP	CB-CA	5.35	1.57	1.53
2	D	10	PTR	CA-C	4.39	1.56	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	7	AHP	CA-C	3.21	1.54	1.48
2	D	10	PTR	P-O1P	3.14	1.61	1.51
2	D	8	TPO	P-O3P	2.69	1.64	1.54
2	D	10	PTR	CE2-CZ	2.35	1.43	1.38
2	D	8	TPO	P-O1P	2.33	1.59	1.51
2	D	10	PTR	P-O3P	-2.13	1.46	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	8	TPO	C-CA-N	12.73	132.64	111.94
2	D	7	AHP	C-CA-N	4.87	118.69	113.83
2	D	8	TPO	OG1-CB-CG2	-4.75	102.19	110.13
2	D	8	TPO	OG1-CB-CA	4.29	114.75	107.55
2	D	8	TPO	CG2-CB-CA	-3.15	106.73	113.20
2	D	10	PTR	O3P-P-OH	3.05	117.72	105.46
2	D	8	TPO	P-OG1-CB	2.47	131.05	120.17
2	D	10	PTR	OH-CZ-CE2	2.23	126.10	119.23
2	D	10	PTR	CG-CB-CA	-2.19	109.25	114.42
2	D	10	PTR	O3P-P-O2P	-2.05	99.62	107.61

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	8	TPO	CA

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

There are no ligands in this entry.



## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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.