



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:05 AM GMT

PDB ID : 3KSD
Title : Crystal Structure of C151S+H178N mutant of Glyceraldehyde-3-phosphate dehydrogenase 1 (GAPDH1) from Staphylococcus aureus MRSA252 complexed with NAD at 2.2 angstrom resolution
Authors : Mukherjee, S.; Dutta, D.; Saha, B.; Das, A.K.
Deposited on : 2009-11-22
Resolution : 2.20 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

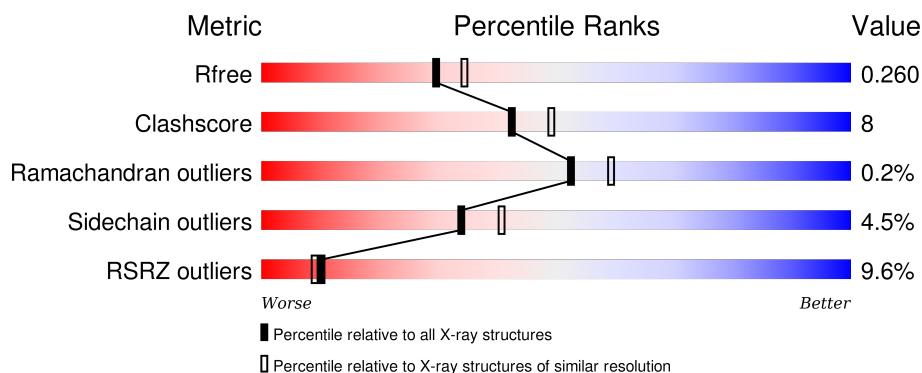
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	O	336	<div> <div>8%</div> <div>76%</div> <div>22%</div> <div>..</div> </div>
1	P	336	<div> <div>16%</div> <div>85%</div> <div>13%</div> <div>..</div> </div>
1	Q	336	<div> <div>7%</div> <div>80%</div> <div>18%</div> <div>..</div> </div>
1	R	336	<div> <div>7%</div> <div>80%</div> <div>17%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10582 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.

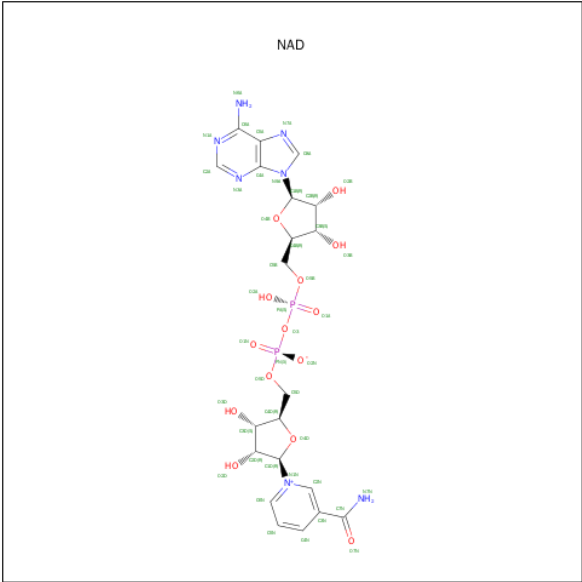
- Molecule 1 is a protein called Glyceraldehyde-3-phosphate dehydrogenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	333	Total	C	N	O	S	0	0	0
			2521	1569	433	511	8			
1	R	333	Total	C	N	O	S	0	1	0
			2524	1570	433	512	9			
1	O	332	Total	C	N	O	S	0	1	0
			2516	1567	429	512	8			
1	P	334	Total	C	N	O	S	0	1	0
			2530	1574	434	514	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	151	SER	CYS	ENGINEERED	UNP Q6GIL8
Q	178	ASN	HIS	ENGINEERED	UNP Q6GIL8
R	151	SER	CYS	ENGINEERED	UNP Q6GIL8
R	178	ASN	HIS	ENGINEERED	UNP Q6GIL8
O	151	SER	CYS	ENGINEERED	UNP Q6GIL8
O	178	ASN	HIS	ENGINEERED	UNP Q6GIL8
P	151	SER	CYS	ENGINEERED	UNP Q6GIL8
P	178	ASN	HIS	ENGINEERED	UNP Q6GIL8

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	Q	1	Total	C	N	O	P	0	0
			43	21	7	13	2		
2	R	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	O	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	P	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

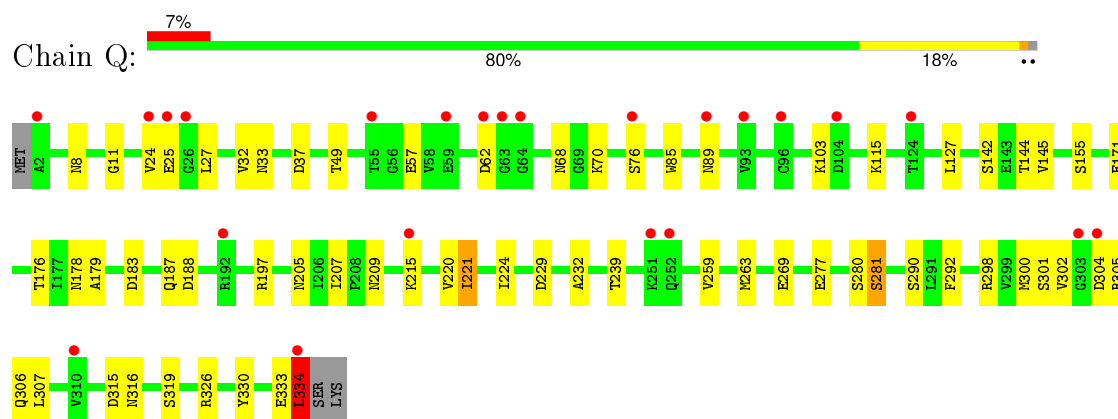
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	Q	102	Total	O	0	0
			102	102		
3	R	83	Total	O	0	0
			83	83		
3	O	65	Total	O	0	0
			65	65		
3	P	66	Total	O	0	0
			66	66		

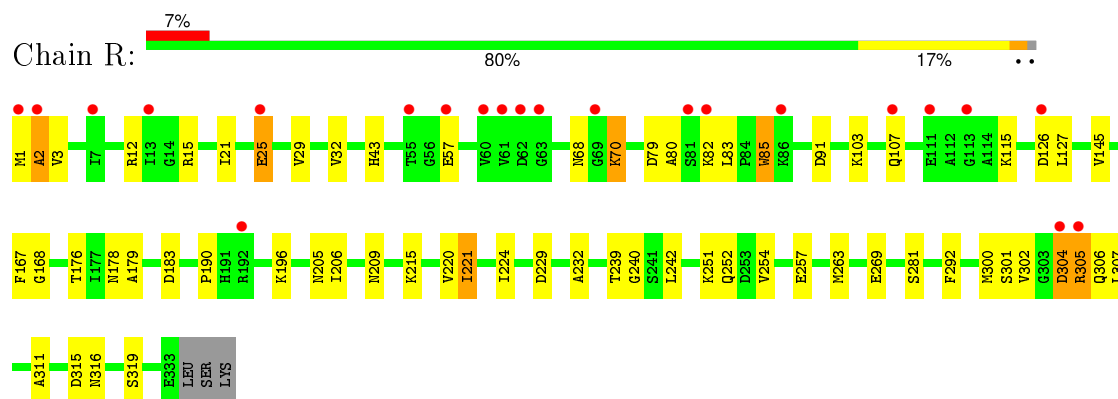
3 Residue-property plots [i](#)

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.

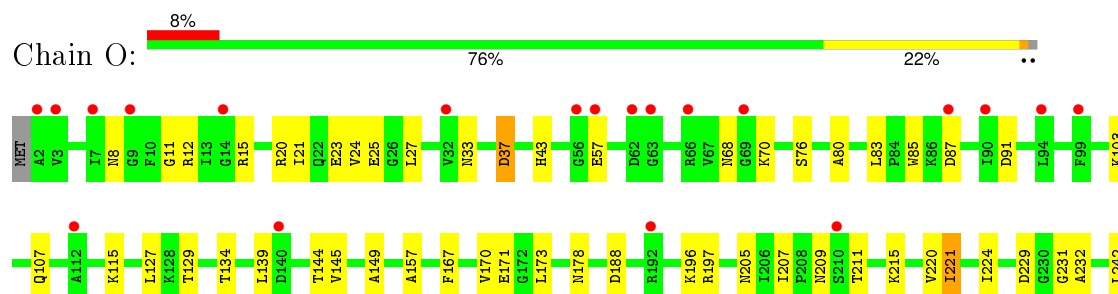
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

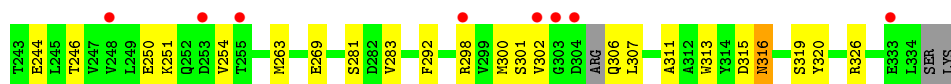


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

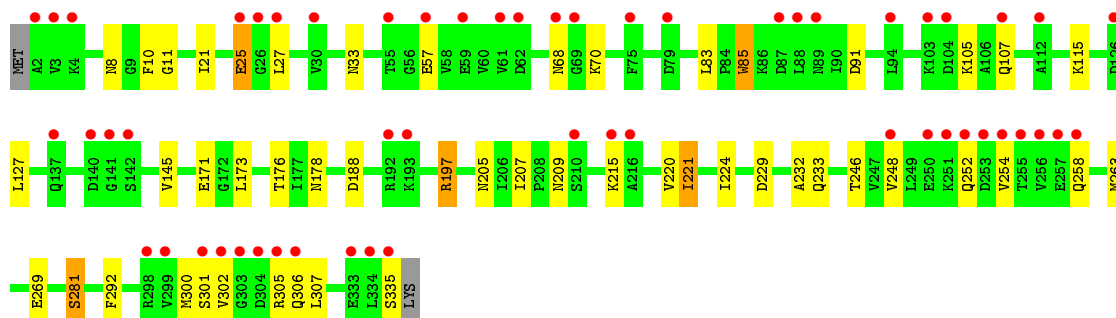
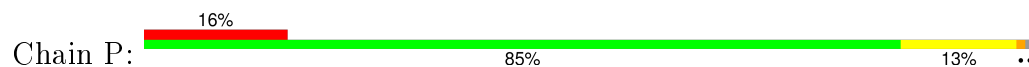


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1





- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.65Å 94.31Å 87.98Å 90.00° 106.02° 90.00°	Depositor
Resolution (Å)	25.84 – 2.20 25.84 – 2.20	Depositor EDS
% Data completeness (in resolution range)	90.0 (25.84-2.20) 90.1 (25.84-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0095	Depositor
R, R_{free}	0.205 , 0.258 0.210 , 0.260	Depositor DCC
R_{free} test set	2439 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	33.7	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 57.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 47902 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10582	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: NAD

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	O	0.77	0/2551	0.80	4/3456 (0.1%)
1	P	0.71	0/2566	0.74	2/3477 (0.1%)
1	Q	0.91	0/2554	0.86	4/3461 (0.1%)
1	R	0.86	0/2560	0.81	0/3468
All	All	0.82	0/10231	0.81	10/13862 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	334	LEU	CA-CB-CG	8.56	134.99	115.30
1	Q	326	ARG	NE-CZ-NH2	-7.93	116.33	120.30
1	O	37	ASP	CB-CG-OD2	6.90	124.51	118.30
1	P	188	ASP	CB-CG-OD1	6.06	123.75	118.30
1	O	188	ASP	CB-CG-OD2	-5.96	112.94	118.30
1	Q	326	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	P	197	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	O	12	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	Q	37	ASP	CB-CG-OD2	5.42	123.18	118.30
1	O	12	ARG	NE-CZ-NH2	-5.33	117.64	120.30

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	O	2516	0	2491	56	7
1	P	2530	0	2509	44	3
1	Q	2521	0	2499	49	5
1	R	2524	0	2505	50	1
2	O	44	0	26	2	0
2	P	44	0	26	1	0
2	Q	43	0	26	1	0
2	R	44	0	26	0	0
3	O	65	0	0	4	0
3	P	66	0	0	8	0
3	Q	102	0	0	5	0
3	R	83	0	0	3	0
All	All	10582	0	10108	169	9

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 8.

All (169) 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:R:304:ASP:OD2	1:R:305:ARG:HG2	1.36	1.25
1:Q:281:SER:HB2	1:R:205:ASN:HD21	1.12	1.08
1:Q:144:THR:HG22	3:Q:414:HOH:O	1.70	0.90
1:O:205:ASN:HD21	1:P:281[A]:SER:HB2	1.36	0.89
1:Q:281:SER:HB2	1:R:205:ASN:ND2	1.88	0.88
1:Q:221:ILE:HG12	1:Q:224:ILE:HG12	1.54	0.87
1:P:171:GLU:HG2	3:P:382:HOH:O	1.73	0.87
1:R:304:ASP:CG	1:R:305:ARG:N	2.30	0.85
1:R:221:ILE:HG12	1:R:224:ILE:HG12	1.59	0.84
1:P:306:GLN:HB2	3:P:341:HOH:O	1.77	0.83
1:P:221:ILE:HG12	1:P:224:ILE:HG12	1.60	0.81
1:O:298:ARG:HD2	3:O:340:HOH:O	1.80	0.80
1:O:221:ILE:HG12	1:O:224:ILE:HG12	1.62	0.79
1:P:252:GLN:HA	3:P:341:HOH:O	1.84	0.77
1:O:37:ASP:HB2	3:O:341:HOH:O	1.84	0.77
1:Q:333:GLU:O	1:Q:334:LEU:HD12	1.85	0.76
1:Q:215:LYS:HE2	1:Q:229:ASP:OD1	1.86	0.76
1:Q:171:GLU:OE2	1:R:305:ARG:HB3	1.88	0.74
1:Q:115:LYS:HG3	3:Q:346:HOH:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:205:ASN:ND2	1:R:281[A]:SER:OG	2.20	0.72
1:P:301:SER:HB3	1:P:306:GLN:HG2	1.71	0.71
1:O:281:SER:HB2	1:P:205:ASN:HD21	1.56	0.69
1:O:307:LEU:HD23	1:P:173:LEU:HG	1.75	0.69
1:O:301:SER:HB3	1:O:306:GLN:HG2	1.75	0.68
1:R:300:MET:CE	1:R:302:VAL:HG23	2.24	0.68
1:Q:300:MET:CE	1:Q:302:VAL:HG23	2.23	0.67
1:O:127:LEU:HD22	1:O:145:VAL:O	1.95	0.67
1:R:304:ASP:OD2	1:R:305:ARG:CG	2.30	0.66
1:R:304:ASP:OD2	1:R:305:ARG:N	2.29	0.66
1:O:300:MET:CE	1:O:302:VAL:HG23	2.25	0.65
1:Q:333:GLU:O	1:Q:334:LEU:HB2	1.97	0.65
1:Q:334:LEU:HD11	3:Q:350:HOH:O	1.97	0.65
1:R:215:LYS:HE2	1:R:229:ASP:OD1	1.97	0.64
1:Q:205:ASN:HD21	1:R:281[B]:SER:HB3	1.60	0.64
1:O:205:ASN:ND2	1:P:281[A]:SER:HB2	2.11	0.63
1:R:301:SER:HB3	1:R:306:GLN:HG2	1.81	0.63
1:P:105:LYS:NZ	3:P:370:HOH:O	2.32	0.63
1:P:300:MET:CE	1:P:302:VAL:HG23	2.29	0.63
1:Q:205:ASN:HD21	1:R:281[A]:SER:HG	1.48	0.62
1:Q:333:GLU:O	1:Q:334:LEU:CB	2.48	0.62
1:R:91:ASP:HA	1:R:115:LYS:HD2	1.82	0.62
1:O:173:LEU:HD12	1:P:246:THR:HG23	1.82	0.61
1:R:215:LYS:NZ	3:R:374:HOH:O	2.33	0.61
1:Q:127:LEU:HD22	1:Q:145:VAL:O	1.99	0.61
1:P:215:LYS:HE2	1:P:229:ASP:OD1	2.02	0.60
1:R:127:LEU:HD22	1:R:145:VAL:O	2.00	0.60
1:R:178:ASN:HD22	1:R:232:ALA:HB1	1.66	0.60
1:R:80:ALA:HA	1:R:83:LEU:HD12	1.82	0.60
1:P:248:VAL:HB	3:P:382:HOH:O	2.01	0.59
1:P:127:LEU:HD22	1:P:145:VAL:O	2.03	0.59
1:P:11:GLY:HA3	2:P:0:NAD:O5B	2.03	0.58
1:O:80:ALA:HA	1:O:83:LEU:HD12	1.85	0.58
1:P:258:GLN:NE2	3:P:397:HOH:O	2.36	0.58
1:R:70:LYS:HB3	3:R:386:HOH:O	2.02	0.58
1:R:315:ASP:O	1:R:319:SER:HB2	2.04	0.57
1:O:281:SER:CB	1:P:205:ASN:HD21	2.17	0.57
1:O:178:ASN:HD22	1:O:232:ALA:HB1	1.71	0.56
1:Q:215:LYS:HE3	3:Q:419:HOH:O	2.05	0.55
1:R:300:MET:HE3	1:R:302:VAL:HG23	1.88	0.54
1:O:215:LYS:HE2	1:O:229:ASP:OD1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:176:THR:HG23	1:Q:178:ASN:ND2	2.23	0.53
1:R:1:MET:O	1:R:3:VAL:N	2.42	0.53
1:R:301:SER:HA	1:R:305:ARG:O	2.09	0.53
1:O:281:SER:HB2	1:P:205:ASN:ND2	2.23	0.53
1:Q:301:SER:HB3	1:Q:306:GLN:HG2	1.92	0.52
1:R:167:PHE:HA	1:R:251:LYS:HE2	1.91	0.52
1:Q:178:ASN:HD22	1:Q:232:ALA:HB1	1.74	0.52
1:Q:49:THR:HG23	1:O:281:SER:OG	2.10	0.52
1:O:91:ASP:HA	1:O:115:LYS:HD2	1.91	0.51
1:O:11:GLY:HA3	2:O:0:NAD:O5B	2.10	0.51
1:Q:197:ARG:HB3	1:Q:207:ILE:HG23	1.93	0.50
1:Q:179:ALA:CB	1:Q:239:THR:HA	2.41	0.50
1:Q:280:SER:HB3	1:R:206:ILE:HB	1.94	0.50
1:O:300:MET:HE2	1:O:302:VAL:HG23	1.92	0.50
1:O:127:LEU:CD2	1:O:145:VAL:O	2.60	0.49
1:Q:304:ASP:OD2	1:Q:305:ARG:HG3	2.11	0.49
1:Q:57:GLU:H	1:Q:68:ASN:ND2	2.10	0.49
1:Q:24:VAL:HG11	1:Q:27:LEU:HD22	1.93	0.49
1:O:315:ASP:O	1:O:319:SER:HB2	2.13	0.49
1:O:316:ASN:H	1:O:316:ASN:ND2	2.10	0.49
1:P:8:ASN:OD1	1:P:33:ASN:HB3	2.13	0.49
1:Q:103:LYS:HB2	1:Q:127:LEU:HG	1.95	0.48
1:P:57:GLU:H	1:P:68:ASN:ND2	2.11	0.48
1:O:320:TYR:CG	2:O:0:NAD:H5N	2.49	0.48
1:Q:300:MET:HE3	1:Q:302:VAL:HG23	1.96	0.48
1:O:197:ARG:HB3	1:O:207:ILE:HG23	1.95	0.47
1:O:144:THR:HG22	3:O:347:HOH:O	2.13	0.47
1:O:263:MET:HG3	1:O:292:PHE:CZ	2.49	0.47
1:R:240:GLY:H	1:R:316:ASN:ND2	2.12	0.47
1:O:20:ARG:NH1	1:O:23:GLU:OE1	2.47	0.47
1:Q:205:ASN:ND2	1:R:281[B]:SER:HB3	2.28	0.47
1:Q:333:GLU:O	1:Q:334:LEU:CD1	2.60	0.47
1:O:205:ASN:HD22	1:P:281[B]:SER:H	1.61	0.47
1:O:205:ASN:HD22	1:P:281[A]:SER:H	1.60	0.47
1:P:197:ARG:HB3	1:P:207:ILE:HG23	1.96	0.47
1:P:21:ILE:HG23	1:P:27:LEU:HB3	1.97	0.46
1:O:246:THR:HG23	1:P:173:LEU:HD12	1.97	0.46
1:R:176:THR:HG23	1:R:178:ASN:ND2	2.30	0.46
1:Q:11:GLY:HA3	2:Q:337:NAD:O5B	2.15	0.46
1:R:57:GLU:H	1:R:68:ASN:ND2	2.13	0.46
1:Q:176:THR:HG23	1:Q:178:ASN:HD21	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:83:LEU:HD13	1:P:85:TRP:CZ2	2.50	0.46
1:Q:315:ASP:O	1:Q:319:SER:HB2	2.16	0.46
1:O:300:MET:SD	1:P:229:ASP:HB2	2.56	0.46
1:P:263:MET:HG3	1:P:292:PHE:CZ	2.51	0.45
1:Q:330:TYR:CE1	1:Q:334:LEU:HD13	2.51	0.45
1:O:306:GLN:N	3:O:343:HOH:O	2.48	0.45
1:R:43:HIS:ND1	1:O:196:LYS:HE2	2.32	0.45
1:O:244:GLU:OE2	1:P:233:GLN:NE2	2.43	0.45
1:P:176:THR:HG23	1:P:178:ASN:HD21	1.81	0.45
1:R:196:LYS:HE2	1:O:43:HIS:ND1	2.31	0.45
1:O:171:GLU:HG3	1:P:307:LEU:HD22	1.98	0.45
1:Q:298:ARG:HG3	1:Q:298:ARG:HH11	1.82	0.45
1:P:207:ILE:HD13	3:P:342:HOH:O	2.17	0.45
1:Q:281:SER:CB	1:R:205:ASN:HD21	2.03	0.45
1:R:196:LYS:HE3	1:O:43:HIS:HB2	1.99	0.44
1:Q:281:SER:H	1:R:205:ASN:HD22	1.64	0.44
1:O:283:VAL:HG11	1:O:313:TRP:HB3	1.99	0.44
1:R:1:MET:O	1:R:2:ALA:C	2.56	0.44
1:O:167:PHE:HA	1:O:251:LYS:HE2	1.99	0.44
1:P:252:GLN:CD	3:P:341:HOH:O	2.56	0.44
1:P:91:ASP:HA	1:P:115:LYS:HD2	1.99	0.43
1:P:263:MET:HG3	1:P:292:PHE:CE1	2.53	0.43
1:O:57:GLU:H	1:O:68:ASN:ND2	2.16	0.43
1:O:205:ASN:HD21	1:P:281[B]:SER:CB	2.27	0.43
1:Q:155:SER:HA	1:Q:290:SER:HB2	2.01	0.43
1:R:242:LEU:HD11	1:R:311:ALA:HB1	2.01	0.43
1:O:21:ILE:HG23	1:O:27:LEU:HB3	2.01	0.43
1:Q:57:GLU:H	1:Q:68:ASN:HD21	1.66	0.43
1:O:24:VAL:HG11	1:O:27:LEU:HD22	2.00	0.43
1:P:176:THR:HG23	1:P:178:ASN:ND2	2.34	0.43
1:P:300:MET:HE2	1:P:302:VAL:HG23	2.01	0.42
1:O:11:GLY:O	1:O:15:ARG:HG3	2.19	0.42
1:O:242:LEU:HD11	1:O:311:ALA:HB1	2.01	0.42
1:Q:263:MET:HG3	1:Q:292:PHE:CE1	2.54	0.42
1:R:12:ARG:HH11	1:R:15:ARG:NH2	2.18	0.42
1:R:183:ASP:N	1:R:183:ASP:OD2	2.52	0.42
1:O:103:LYS:HB2	1:O:127:LEU:HG	2.01	0.42
1:Q:144:THR:CG2	3:Q:414:HOH:O	2.48	0.42
1:Q:316:ASN:H	1:Q:316:ASN:ND2	2.17	0.42
1:O:173:LEU:CD2	1:O:229:ASP:HB3	2.49	0.42
1:O:8:ASN:OD1	1:O:33:ASN:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:127:LEU:HA	1:P:127:LEU:HD23	1.83	0.42
1:O:211:THR:HG22	1:O:231:GLY:HA2	2.02	0.42
1:R:263:MET:HG3	1:R:292:PHE:CZ	2.55	0.42
1:P:57:GLU:H	1:P:68:ASN:HD21	1.68	0.41
1:Q:187:GLN:O	1:Q:188:ASP:C	2.58	0.41
1:P:300:MET:HE3	1:P:302:VAL:HG23	2.01	0.41
1:R:196:LYS:NZ	3:R:399:HOH:O	2.54	0.41
1:R:21:ILE:HG21	1:R:29:VAL:HG23	2.02	0.41
1:Q:307:LEU:HD11	1:R:307:LEU:HD11	2.02	0.41
1:R:103:LYS:HB2	1:R:127:LEU:HG	2.02	0.41
1:R:176:THR:HG23	1:R:178:ASN:HD21	1.86	0.41
1:R:126:ASP:OD1	1:R:126:ASP:N	2.51	0.41
1:Q:259:VAL:O	1:Q:263:MET:HG2	2.20	0.41
1:R:83:LEU:HD13	1:R:85:TRP:CZ2	2.56	0.41
1:P:178:ASN:HD22	1:P:232:ALA:HB1	1.86	0.41
1:R:79:ASP:OD2	1:R:82:LYS:HG2	2.21	0.41
1:R:301:SER:CB	1:R:306:GLN:HG2	2.48	0.41
1:Q:127:LEU:HA	1:Q:127:LEU:HD23	1.76	0.41
1:O:316:ASN:HD22	1:O:316:ASN:H	1.68	0.41
1:O:263:MET:HG3	1:O:292:PHE:CE1	2.55	0.41
1:O:170:VAL:HG11	1:P:305:ARG:NH1	2.36	0.41
1:R:179:ALA:CB	1:R:239:THR:HA	2.51	0.40
1:O:134:THR:HG21	1:O:157:ALA:HB1	2.02	0.40
1:Q:183:ASP:OD2	1:Q:183:ASP:N	2.53	0.40
1:O:173:LEU:HD23	1:O:229:ASP:HB3	2.03	0.40
1:Q:8:ASN:OD1	1:Q:33:ASN:HB3	2.22	0.40
1:O:129:THR:HG22	1:O:149:ALA:HB2	2.03	0.40

All (9) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:62:ASP:OD2	1:O:139:LEU:C[2_546]	1.75	0.45
1:Q:62:ASP:OD2	1:O:139:LEU:O[2_546]	1.83	0.37
1:Q:142:SER:OG	1:Q:277:GLU:OE2[2_545]	1.85	0.35
1:Q:89:ASN:ND2	1:Q:306:GLN:NE2[2_545]	1.95	0.25
1:O:326:ARG:CZ	1:P:25:GLU:OE1[1_455]	1.96	0.24
1:O:326:ARG:NH1	1:P:25:GLU:OE1[1_455]	2.10	0.10
1:Q:305:ARG:NH2	1:O:250:GLU:OE1[1_554]	2.12	0.08
1:O:326:ARG:NE	1:P:25:GLU:OE1[1_455]	2.12	0.08
1:R:25:GLU:OE2	1:O:87:ASP:OD2[1_655]	2.13	0.07

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	O	329/336 (98%)	314 (95%)	15 (5%)	0	100	100
1	P	333/336 (99%)	319 (96%)	14 (4%)	0	100	100
1	Q	331/336 (98%)	316 (96%)	15 (4%)	0	100	100
1	R	332/336 (99%)	318 (96%)	12 (4%)	2 (1%)	30	29
All	All	1325/1344 (99%)	1267 (96%)	56 (4%)	2 (0%)	52	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	R	2	ALA
1	R	168	GLY

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	O	271/274 (99%)	259 (96%)	12 (4%)	35	42
1	P	273/274 (100%)	260 (95%)	13 (5%)	31	37
1	Q	271/274 (99%)	260 (96%)	11 (4%)	37	45
1	R	272/274 (99%)	257 (94%)	15 (6%)	27	30
All	All	1087/1096 (99%)	1036 (95%)	51 (5%)	34	39

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Q	25	GLU
1	Q	32	VAL
1	Q	70	LYS
1	Q	76	SER
1	Q	85	TRP
1	Q	209	ASN
1	Q	220	VAL
1	Q	221	ILE
1	Q	269	GLU
1	Q	281	SER
1	Q	334	LEU
1	R	25	GLU
1	R	32	VAL
1	R	70	LYS
1	R	85	TRP
1	R	107	GLN
1	R	190	PRO
1	R	209	ASN
1	R	220	VAL
1	R	221	ILE
1	R	252	GLN
1	R	254	VAL
1	R	257	GLU
1	R	269	GLU
1	R	304	ASP
1	R	305	ARG
1	O	25	GLU
1	O	70	LYS
1	O	76	SER
1	O	85	TRP
1	O	107	GLN
1	O	209	ASN
1	O	220	VAL
1	O	221	ILE
1	O	254	VAL
1	O	269[A]	GLU
1	O	269[B]	GLU
1	O	316	ASN
1	P	10	PHE
1	P	25	GLU
1	P	70	LYS
1	P	85	TRP
1	P	107	GLN

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Mol	Chain	Res	Type
1	P	209	ASN
1	P	220	VAL
1	P	221	ILE
1	P	254	VAL
1	P	269	GLU
1	P	281[A]	SER
1	P	281[B]	SER
1	P	335	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	Q	68	ASN
1	Q	178	ASN
1	Q	205	ASN
1	Q	258	GLN
1	Q	265	ASN
1	Q	296	GLN
1	Q	316	ASN
1	R	68	ASN
1	R	178	ASN
1	R	205	ASN
1	R	258	GLN
1	R	265	ASN
1	R	296	GLN
1	R	316	ASN
1	O	68	ASN
1	O	178	ASN
1	O	205	ASN
1	O	252	GLN
1	O	258	GLN
1	O	265	ASN
1	P	68	ASN
1	P	178	ASN
1	P	205	ASN
1	P	258	GLN
1	P	265	ASN
1	P	296	GLN
1	P	316	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 ⓘ

4 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	NAD	O	0	-	38,48,48	1.70	3 (7%)	47,73,73	2.07	5 (10%)
2	NAD	P	0	-	38,48,48	1.79	3 (7%)	47,73,73	2.19	11 (23%)
2	NAD	Q	337	-	36,47,48	1.72	4 (11%)	39,70,73	2.81	9 (23%)
2	NAD	R	0	-	38,48,48	1.72	5 (13%)	47,73,73	2.30	8 (17%)

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	NAD	O	0	-	-	0/22/62/62	0/5/5/5
2	NAD	P	0	-	-	0/22/62/62	0/5/5/5
2	NAD	Q	337	-	-	0/16/60/62	0/5/5/5
2	NAD	R	0	-	-	0/22/62/62	0/5/5/5

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	R	0	NAD	PA-O2A	-2.02	1.46	1.54
2	Q	337	NAD	C2A-N1A	2.02	1.37	1.33
2	R	0	NAD	O5B-C5B	2.42	1.54	1.44
2	Q	337	NAD	PN-O3	2.53	1.64	1.60
2	R	0	NAD	C2A-N1A	2.84	1.39	1.33
2	R	0	NAD	C2A-N3A	3.45	1.38	1.32
2	O	0	NAD	C2A-N1A	3.47	1.40	1.33
2	P	0	NAD	C2A-N1A	3.67	1.40	1.33
2	Q	337	NAD	C2A-N3A	4.12	1.39	1.32
2	P	0	NAD	C2A-N3A	4.27	1.39	1.32
2	O	0	NAD	C2A-N3A	4.30	1.39	1.32
2	O	0	NAD	O7N-C7N	7.46	1.40	1.24
2	R	0	NAD	O7N-C7N	7.63	1.40	1.24
2	Q	337	NAD	O7N-C7N	8.03	1.41	1.24
2	P	0	NAD	O7N-C7N	8.24	1.41	1.24

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	337	NAD	N3A-C2A-N1A	-13.19	118.79	128.89
2	R	0	NAD	N3A-C2A-N1A	-12.21	119.54	128.89
2	O	0	NAD	N3A-C2A-N1A	-11.51	120.08	128.89
2	P	0	NAD	N3A-C2A-N1A	-10.14	121.13	128.89
2	R	0	NAD	O7N-C7N-N7N	-4.16	116.74	122.59
2	P	0	NAD	PN-O3-PA	-4.07	121.31	132.73
2	Q	337	NAD	O7N-C7N-C3N	-3.53	115.73	119.59
2	R	0	NAD	O3B-C3B-C4B	-3.26	101.27	111.05
2	Q	337	NAD	C1B-N9A-C4A	-3.10	122.26	126.94
2	R	0	NAD	C4A-C5A-N7A	-3.04	106.68	109.48
2	Q	337	NAD	O3B-C3B-C4B	-2.97	102.15	111.05
2	O	0	NAD	O3-PN-O5D	-2.75	95.65	102.94
2	P	0	NAD	O7N-C7N-C3N	-2.58	116.77	119.59
2	P	0	NAD	C4B-O4B-C1B	-2.51	106.96	109.72
2	Q	337	NAD	O3B-C3B-C2B	-2.48	103.75	111.83
2	R	0	NAD	O5B-C5B-C4B	-2.33	100.52	109.12
2	P	0	NAD	C1B-N9A-C4A	-2.15	123.69	126.94
2	P	0	NAD	O5D-PN-O1N	-2.00	101.84	109.62
2	O	0	NAD	C2N-C3N-C4N	2.00	120.52	118.29
2	O	0	NAD	O2A-PA-O3	2.03	114.29	105.09
2	P	0	NAD	O2N-PN-O1N	2.06	123.70	112.53
2	Q	337	NAD	C2A-N1A-C6A	2.07	122.47	118.77
2	Q	337	NAD	C4B-O4B-C1B	2.08	112.00	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	P	0	NAD	O2A-PA-O3	2.26	115.33	105.09
2	R	0	NAD	O4D-C1D-N1N	2.30	110.66	108.13
2	P	0	NAD	O4D-C1D-N1N	2.31	110.67	108.13
2	R	0	NAD	O3-PA-O5B	2.45	109.43	102.94
2	O	0	NAD	O2N-PN-O1N	2.55	126.32	112.53
2	P	0	NAD	O3-PA-O5B	3.20	111.42	102.94
2	R	0	NAD	C3N-C7N-N7N	3.50	121.65	117.82
2	P	0	NAD	C3N-C7N-N7N	3.77	121.94	117.82
2	Q	337	NAD	C3N-C7N-N7N	5.26	123.57	117.82
2	Q	337	NAD	O4D-C1D-N1N	5.48	114.15	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	O	0	NAD	2	0
2	P	0	NAD	1	0
2	Q	337	NAD	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 ⓘ

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	O	332/336 (98%)	0.54	28 (8%) 14 13	21, 37, 62, 85	0
1	P	334/336 (99%)	0.93	55 (16%) 2 2	21, 46, 78, 106	0
1	Q	333/336 (99%)	0.41	23 (6%) 20 19	17, 31, 51, 58	0
1	R	333/336 (99%)	0.40	22 (6%) 22 21	19, 32, 53, 62	0
All	All	1332/1344 (99%)	0.57	128 (9%) 10 9	17, 36, 66, 106	0

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	P	303	GLY	10.1
1	P	304	ASP	7.8
1	P	25	GLU	7.0
1	R	61	VAL	6.6
1	P	254	VAL	6.5
1	R	62	ASP	6.2
1	P	302	VAL	6.2
1	P	253	ASP	5.7
1	P	2	ALA	5.6
1	P	252	GLN	5.5
1	P	301	SER	5.2
1	O	62	ASP	5.1
1	O	304	ASP	5.0
1	P	335	SER	4.9
1	Q	55	THR	4.9
1	Q	304	ASP	4.9
1	P	126	ASP	4.6
1	Q	62	ASP	4.6
1	O	303	GLY	4.6
1	Q	63	GLY	4.5
1	R	63	GLY	4.5

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Mol	Chain	Res	Type	RSRZ
1	P	61	VAL	4.4
1	O	56	GLY	4.3
1	O	7	ILE	4.2
1	O	69	GLY	4.2
1	R	55	THR	4.1
1	P	299	VAL	4.0
1	P	137	GLN	4.0
1	P	192	ARG	3.9
1	P	298	ARG	3.8
1	R	192	ARG	3.8
1	O	192	ARG	3.8
1	P	62	ASP	3.7
1	R	25	GLU	3.7
1	O	253	ASP	3.6
1	O	302	VAL	3.6
1	P	256	VAL	3.6
1	Q	2	ALA	3.6
1	P	55	THR	3.5
1	R	81	SER	3.5
1	R	126	ASP	3.5
1	P	305	ARG	3.4
1	Q	26	GLY	3.4
1	O	63	GLY	3.4
1	P	210	SER	3.4
1	P	88	LEU	3.4
1	Q	192	ARG	3.3
1	Q	334	LEU	3.3
1	R	86	LYS	3.3
1	P	141	GLY	3.3
1	P	3	VAL	3.2
1	Q	24	VAL	3.2
1	O	57	GLU	3.1
1	R	60	VAL	3.1
1	P	250	GLU	3.1
1	P	255	THR	3.1
1	O	66	ARG	3.1
1	Q	215	LYS	3.0
1	R	69	GLY	3.0
1	R	107	GLN	3.0
1	P	258	GLN	3.0
1	Q	25	GLU	2.9
1	R	7	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	P	27	LEU	2.9
1	O	2	ALA	2.9
1	P	251	LYS	2.9
1	Q	59	GLU	2.8
1	P	248	VAL	2.8
1	O	140	ASP	2.8
1	Q	76	SER	2.7
1	R	111	GLU	2.7
1	P	334	LEU	2.7
1	Q	303	GLY	2.7
1	O	3	VAL	2.7
1	R	2	ALA	2.7
1	P	112	ALA	2.6
1	O	298	ARG	2.6
1	Q	64	GLY	2.6
1	R	113	GLY	2.6
1	P	30	VAL	2.6
1	R	305	ARG	2.5
1	P	89	ASN	2.5
1	P	215	LYS	2.5
1	Q	252	GLN	2.5
1	P	59	GLU	2.5
1	P	26	GLY	2.5
1	O	90	ILE	2.5
1	O	248	VAL	2.4
1	R	13	ILE	2.4
1	P	107	GLN	2.4
1	O	14	GLY	2.4
1	P	87	ASP	2.4
1	O	333	GLU	2.4
1	P	69	GLY	2.4
1	P	306	GLN	2.4
1	Q	124	THR	2.4
1	P	333	GLU	2.4
1	P	75	PHE	2.4
1	P	257	GLU	2.4
1	P	104	ASP	2.3
1	R	1	MET	2.3
1	O	210	SER	2.3
1	O	32	VAL	2.3
1	R	82	LYS	2.3
1	P	140	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	Q	104	ASP	2.3
1	P	103	LYS	2.3
1	P	94	LEU	2.3
1	R	304	ASP	2.3
1	Q	89	ASN	2.2
1	O	112	ALA	2.2
1	P	4	LYS	2.2
1	Q	251	LYS	2.2
1	O	9	GLY	2.2
1	Q	93	VAL	2.2
1	P	193	LYS	2.2
1	R	57	GLU	2.2
1	P	57	GLU	2.2
1	O	99	PHE	2.1
1	P	216	ALA	2.1
1	P	68	ASN	2.1
1	O	87	ASP	2.1
1	Q	96	CYS	2.1
1	Q	310	VAL	2.1
1	O	255	THR	2.0
1	O	94	LEU	2.0
1	P	142	SER	2.0
1	P	79	ASP	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	NAD	P	0	44/44	0.96	0.11	-0.95	20,29,36,38	0
2	NAD	R	0	44/44	0.96	0.11	-1.00	14,24,26,29	0
2	NAD	Q	337	43/44	0.96	0.10	-1.19	14,24,27,29	0
2	NAD	O	0	44/44	0.97	0.10	-1.21	19,27,30,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.